



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 20, 2016 – 01:37 AM GMT

PDB ID : 4WQR
Title : Complex of 70S ribosome with tRNA-Phe and mRNA with C-A mismatch in the first position in the A-site.
Authors : Rozov, A.; Demeshkina, N.; Yusupov, M.; Yusupova, G.
Deposited on : 2014-10-22
Resolution : 3.15 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20026982

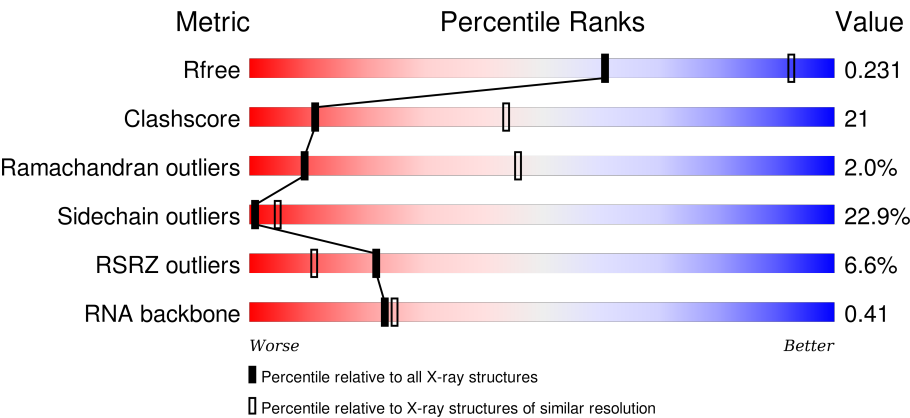
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



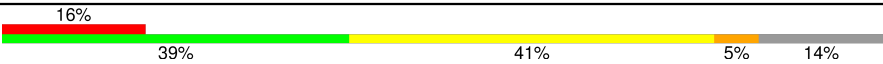


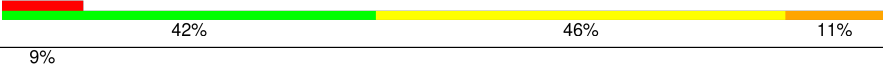
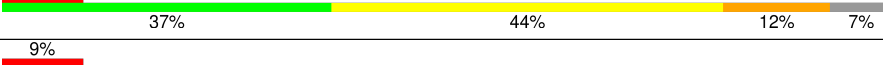



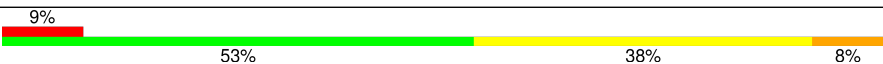

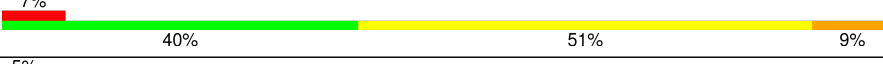
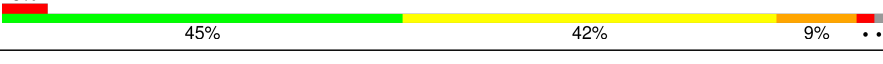



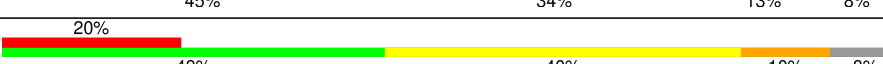

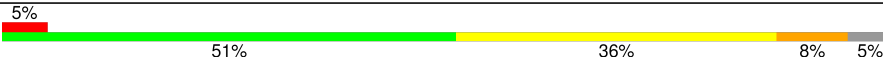
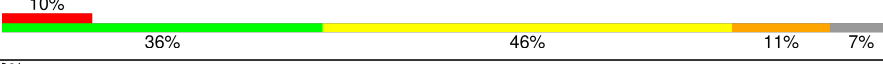



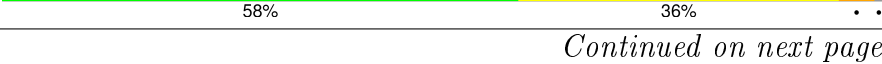


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1112 (3.20-3.12)
Clashscore	102246	1249 (3.20-3.12)
Ramachandran outliers	100387	1222 (3.20-3.12)
Sidechain outliers	100360	1221 (3.20-3.12)
RSRZ outliers	91569	1117 (3.20-3.12)
RNA backbone	2183	1046 (3.62-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	13	1522	<div><div></div><div><div>27%</div><div>44%</div><div>24%</div><div>• •</div></div></div>
1	1G	1522	<div><div></div><div><div>31%</div><div>46%</div><div>19%</div><div>• •</div></div></div>
2	12	256	<div><div>11%</div><div><div>48%</div><div>36%</div><div>8%</div><div>• 7%</div></div></div>
2	1E	256	<div><div>10%</div><div><div>38%</div><div>41%</div><div>13%</div><div>• 7%</div></div></div>

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Mol	Chain	Length	Quality of chain
3	22	239	
3	2E	239	
4	32	209	
4	3E	209	
5	42	162	
5	4E	162	
6	52	101	
6	5E	101	
7	62	156	
7	6E	156	
8	72	138	
8	7E	138	
9	82	128	
9	8E	128	
10	1A	105	
10	1I	105	
11	2A	129	
11	2I	129	
12	3A	132	
12	3I	132	
13	4A	126	
13	4I	126	
14	5A	61	
14	5I	61	
15	6A	89	


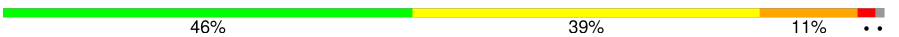



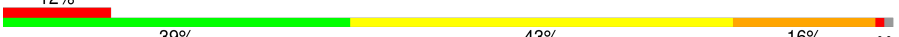

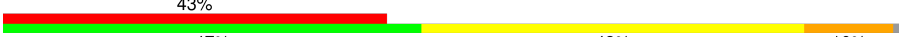
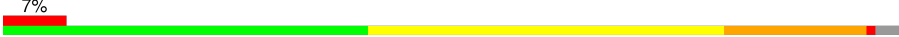

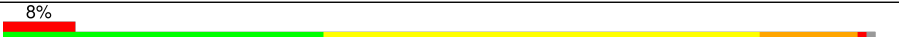
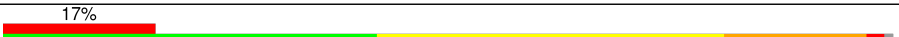

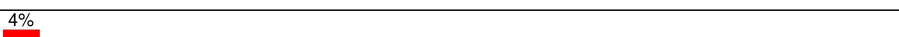
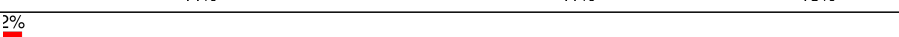
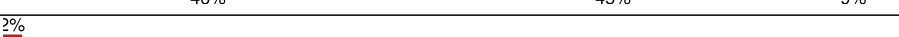



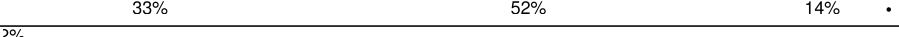


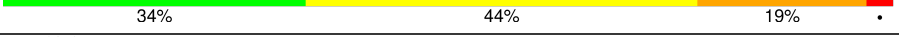


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Mol	Chain	Length	Quality of chain
15	6I	89	
16	7A	88	
16	7I	88	
17	8A	105	
17	8I	105	
18	9A	88	
18	9I	88	
19	AA	93	
19	AI	93	
20	BA	106	
20	BI	106	
21	1B	27	
21	1F	27	
22	1K	76	
22	1L	76	
23	2K	77	
23	2L	77	
24	3K	76	
24	3L	76	
25	4K	30	
25	4L	30	
26	14	2917	
26	1H	2917	
27	16	122	
27	1J	122	

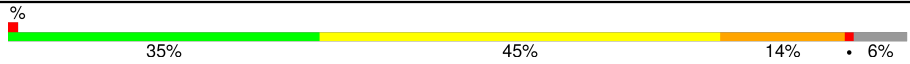

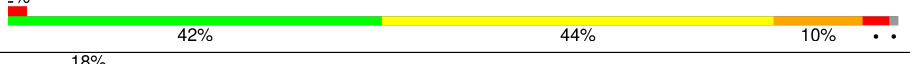
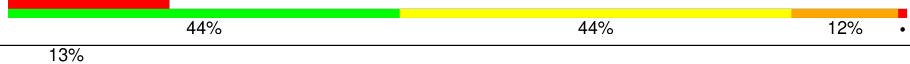
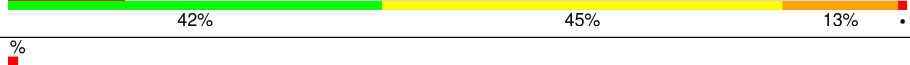
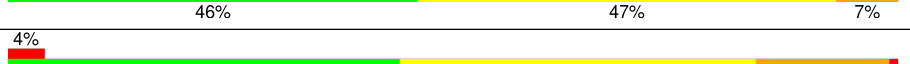
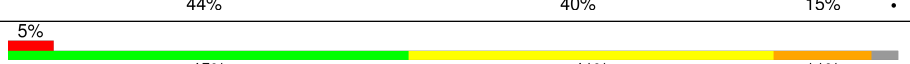
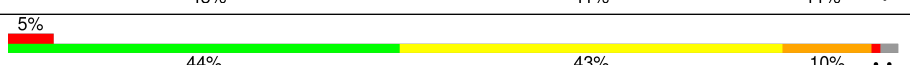
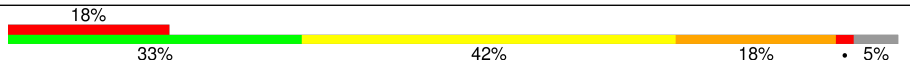
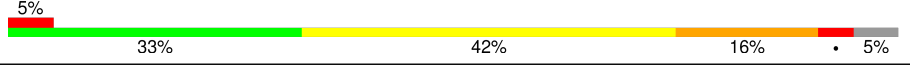
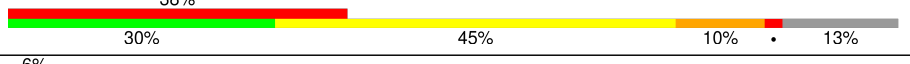
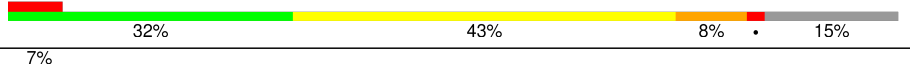
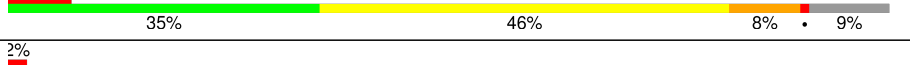


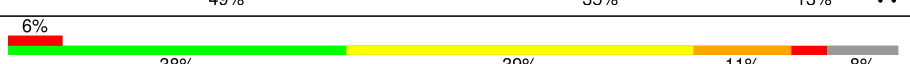
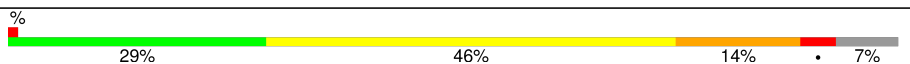
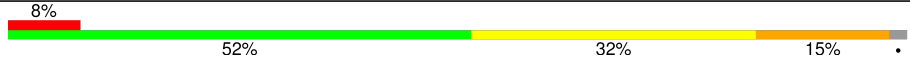
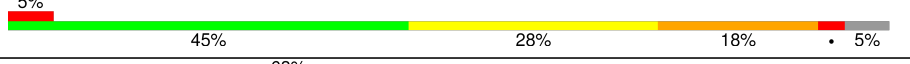
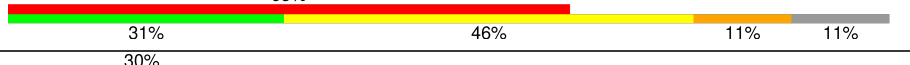
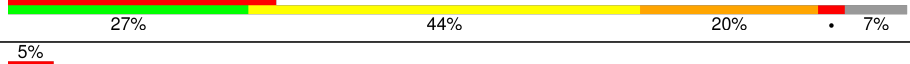




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Mol	Chain	Length	Quality of chain
28	11	276	
28	19	276	
29	21	206	
29	29	206	
30	31	210	
30	39	210	
31	41	182	
31	49	182	
32	51	180	
32	59	180	
33	61	148	
33	69	148	
34	15	140	
34	58	140	
35	25	122	
35	68	122	
36	35	150	
36	78	150	
37	45	141	
37	88	141	
38	55	118	
38	98	118	
39	65	112	
39	A8	112	
40	75	146	

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Mol	Chain	Length	Quality of chain
40	B8	146	
41	85	118	
41	C8	118	
42	95	101	
42	D8	101	
43	A5	113	
43	E8	113	
44	B5	96	
44	F8	96	
45	C5	110	
45	G8	110	
46	D5	206	
46	H8	206	
47	E5	85	
47	I8	85	
48	F5	98	
48	J8	98	
49	G5	72	
49	K8	72	
50	H5	60	
50	L8	60	
51	I5	71	
51	M8	71	
52	J5	60	
52	N8	60	

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Mol	Chain	Length	Quality of chain
53	K5	54	
53	O8	54	
54	L5	49	
54	P8	49	
55	M5	65	
55	Q8	65	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
56	MG	11	301	-	-	-	X
56	MG	11	303	-	-	-	X
56	MG	13	1601	-	-	-	X
56	MG	13	1603	-	-	-	X
56	MG	13	1604	-	-	-	X
56	MG	13	1605	-	-	-	X
56	MG	13	1607	-	-	-	X
56	MG	13	1608	-	-	-	X
56	MG	13	1610	-	-	-	X
56	MG	13	1612	-	-	-	X
56	MG	13	1614	-	-	-	X
56	MG	13	1615	-	-	-	X
56	MG	13	1619	-	-	-	X
56	MG	13	1620	-	-	-	X
56	MG	13	1625	-	-	-	X
56	MG	13	1629	-	-	-	X
56	MG	13	1630	-	-	-	X
56	MG	13	1631	-	-	-	X
56	MG	13	1641	-	-	-	X
56	MG	13	1643	-	-	-	X
56	MG	13	1650	-	-	-	X
56	MG	13	1651	-	-	-	X
56	MG	13	1654	-	-	-	X
56	MG	13	1655	-	-	-	X
56	MG	13	1658	-	-	-	X
56	MG	13	1661	-	-	-	X
56	MG	13	1663	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
56	MG	13	1668	-	-	-	X
56	MG	13	1669	-	-	-	X
56	MG	13	1671	-	-	-	X
56	MG	13	1677	-	-	-	X
56	MG	13	1679	-	-	-	X
56	MG	13	1681	-	-	-	X
56	MG	13	1699	-	-	-	X
56	MG	13	1751	-	-	-	X
56	MG	14	3004	-	-	-	X
56	MG	14	3007	-	-	-	X
56	MG	14	3011	-	-	-	X
56	MG	14	3013	-	-	-	X
56	MG	14	3015	-	-	-	X
56	MG	14	3018	-	-	-	X
56	MG	14	3022	-	-	-	X
56	MG	14	3023	-	-	-	X
56	MG	14	3026	-	-	-	X
56	MG	14	3033	-	-	-	X
56	MG	14	3034	-	-	-	X
56	MG	14	3035	-	-	-	X
56	MG	14	3036	-	-	-	X
56	MG	14	3039	-	-	-	X
56	MG	14	3040	-	-	-	X
56	MG	14	3041	-	-	-	X
56	MG	14	3043	-	-	-	X
56	MG	14	3044	-	-	-	X
56	MG	14	3046	-	-	-	X
56	MG	14	3050	-	-	-	X
56	MG	14	3053	-	-	-	X
56	MG	14	3054	-	-	-	X
56	MG	14	3058	-	-	-	X
56	MG	14	3064	-	-	-	X
56	MG	14	3065	-	-	-	X
56	MG	14	3070	-	-	-	X
56	MG	14	3071	-	-	-	X
56	MG	14	3073	-	-	-	X
56	MG	14	3079	-	-	-	X
56	MG	14	3084	-	-	-	X
56	MG	14	3085	-	-	-	X
56	MG	14	3090	-	-	-	X
56	MG	14	3091	-	-	-	X
56	MG	14	3100	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
56	MG	14	3104	-	-	-	X
56	MG	14	3108	-	-	-	X
56	MG	14	3110	-	-	-	X
56	MG	14	3111	-	-	-	X
56	MG	14	3112	-	-	-	X
56	MG	14	3114	-	-	-	X
56	MG	14	3120	-	-	-	X
56	MG	14	3136	-	-	-	X
56	MG	14	3139	-	-	-	X
56	MG	14	3141	-	-	-	X
56	MG	14	3142	-	-	-	X
56	MG	14	3144	-	-	-	X
56	MG	14	3146	-	-	-	X
56	MG	14	3148	-	-	-	X
56	MG	14	3150	-	-	-	X
56	MG	14	3152	-	-	-	X
56	MG	14	3155	-	-	-	X
56	MG	14	3160	-	-	-	X
56	MG	14	3162	-	-	-	X
56	MG	14	3165	-	-	-	X
56	MG	14	3167	-	-	-	X
56	MG	14	3178	-	-	-	X
56	MG	14	3185	-	-	-	X
56	MG	14	3186	-	-	-	X
56	MG	14	3200	-	-	-	X
56	MG	14	3227	-	-	-	X
56	MG	14	3234	-	-	-	X
56	MG	14	3246	-	-	-	X
56	MG	14	3248	-	-	-	X
56	MG	14	3253	-	-	-	X
56	MG	14	3387	-	-	-	X
56	MG	16	205	-	-	-	X
56	MG	1G	1601	-	-	-	X
56	MG	1G	1602	-	-	-	X
56	MG	1G	1603	-	-	-	X
56	MG	1G	1605	-	-	-	X
56	MG	1G	1611	-	-	-	X
56	MG	1G	1617	-	-	-	X
56	MG	1G	1618	-	-	-	X
56	MG	1G	1632	-	-	-	X
56	MG	1G	1633	-	-	-	X
56	MG	1G	1635	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
56	MG	1G	1639	-	-	-	X
56	MG	1G	1641	-	-	-	X
56	MG	1G	1651	-	-	-	X
56	MG	1G	1654	-	-	-	X
56	MG	1G	1658	-	-	-	X
56	MG	1G	1659	-	-	-	X
56	MG	1G	1663	-	-	-	X
56	MG	1H	3001	-	-	-	X
56	MG	1H	3004	-	-	-	X
56	MG	1H	3007	-	-	-	X
56	MG	1H	3009	-	-	-	X
56	MG	1H	3019	-	-	-	X
56	MG	1H	3020	-	-	-	X
56	MG	1H	3021	-	-	-	X
56	MG	1H	3022	-	-	-	X
56	MG	1H	3023	-	-	-	X
56	MG	1H	3024	-	-	-	X
56	MG	1H	3028	-	-	-	X
56	MG	1H	3029	-	-	-	X
56	MG	1H	3031	-	-	-	X
56	MG	1H	3033	-	-	-	X
56	MG	1H	3036	-	-	-	X
56	MG	1H	3041	-	-	-	X
56	MG	1H	3042	-	-	-	X
56	MG	1H	3046	-	-	-	X
56	MG	1H	3047	-	-	-	X
56	MG	1H	3050	-	-	-	X
56	MG	1H	3053	-	-	-	X
56	MG	1H	3055	-	-	-	X
56	MG	1H	3056	-	-	-	X
56	MG	1H	3057	-	-	-	X
56	MG	1H	3066	-	-	-	X
56	MG	1H	3068	-	-	-	X
56	MG	1H	3071	-	-	-	X
56	MG	1H	3072	-	-	-	X
56	MG	1H	3073	-	-	-	X
56	MG	1H	3075	-	-	-	X
56	MG	1H	3079	-	-	-	X
56	MG	1H	3082	-	-	-	X
56	MG	1H	3083	-	-	-	X
56	MG	1H	3087	-	-	-	X
56	MG	1H	3088	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
56	MG	1H	3090	-	-	-	X
56	MG	1H	3092	-	-	-	X
56	MG	1H	3096	-	-	-	X
56	MG	1H	3097	-	-	-	X
56	MG	1H	3099	-	-	-	X
56	MG	1H	3110	-	-	-	X
56	MG	1H	3113	-	-	-	X
56	MG	1H	3117	-	-	-	X
56	MG	1H	3120	-	-	-	X
56	MG	1H	3122	-	-	-	X
56	MG	1H	3126	-	-	-	X
56	MG	1H	3129	-	-	-	X
56	MG	1H	3130	-	-	-	X
56	MG	1H	3133	-	-	-	X
56	MG	1H	3137	-	-	-	X
56	MG	1H	3148	-	-	-	X
56	MG	1H	3153	-	-	-	X
56	MG	1H	3154	-	-	-	X
56	MG	1H	3164	-	-	-	X
56	MG	1H	3165	-	-	-	X
56	MG	1H	3167	-	-	-	X
56	MG	1H	3174	-	-	-	X
56	MG	1H	3194	-	-	-	X
56	MG	1H	3197	-	-	-	X
56	MG	1H	3201	-	-	-	X
56	MG	1H	3202	-	-	-	X
56	MG	1H	3217	-	-	-	X
56	MG	1H	3234	-	-	-	X
56	MG	1H	3241	-	-	-	X
56	MG	1H	3255	-	-	-	X
56	MG	1H	3258	-	-	-	X
56	MG	1H	3281	-	-	-	X
56	MG	1H	3292	-	-	-	X
56	MG	1H	3297	-	-	-	X
56	MG	1H	3300	-	-	-	X
56	MG	1H	3309	-	-	-	X
56	MG	1H	3312	-	-	-	X
56	MG	1H	3316	-	-	-	X
56	MG	1H	3318	-	-	-	X
56	MG	1H	3326	-	-	-	X
56	MG	1H	3524	-	-	-	X
56	MG	1J	205	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
56	MG	1J	206	-	-	-	X
56	MG	29	302	-	-	-	X
56	MG	2K	103	-	-	-	X
56	MG	2L	101	-	-	-	X
56	MG	31	301	-	-	-	X
56	MG	3I	201	-	-	-	X
56	MG	78	201	-	-	-	X
56	MG	85	201	-	-	-	X
56	MG	98	201	-	-	-	X
56	MG	F5	101	-	-	-	X
57	ZN	3E	303	-	-	-	X
57	ZN	G8	201	-	-	-	X

2 Entry composition [i](#)

There are 58 unique types of molecules in this entry. The entry contains 300537 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called 16S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	13	1498	Total	C	N	O	P	0	0	0
			32207	14334	5973	10402	1498			
1	1G	1497	Total	C	N	O	P	0	0	0
			32182	14324	5968	10394	1496			

- Molecule 2 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	1E	237	Total	C	N	O	S	0	0	0
			1924	1228	344	347	5			
2	12	237	Total	C	N	O	S	0	0	0
			1924	1228	344	347	5			

- Molecule 3 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	2E	205	Total	C	N	O	S	0	0	0
			1605	1011	313	280	1			
3	22	206	Total	C	N	O	S	0	0	0
			1612	1016	314	281	1			

- Molecule 4 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	3E	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			
4	32	208	Total	C	N	O	S	0	0	0
			1703	1066	339	291	7			

- Molecule 5 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	4E	151	Total	C	N	O	S	0	0	0
			1155	729	218	204	4			
5	42	151	Total	C	N	O	S	0	0	0
			1155	729	218	204	4			

- Molecule 6 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	5E	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			
6	52	101	Total	C	N	O	S	0	0	0
			843	531	155	154	3			

- Molecule 7 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	6E	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			
7	62	155	Total	C	N	O	S	0	0	0
			1257	781	252	218	6			

- Molecule 8 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	7E	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			
8	72	138	Total	C	N	O	S	0	0	0
			1116	705	215	193	3			

- Molecule 9 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	8E	127	Total	C	N	O		0	0	0
			1010	639	197	174				
9	82	127	Total	C	N	O		0	0	0
			1010	639	197	174				

- Molecule 10 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	1I	99	Total	C	N	O	S	0	0	0
			801	504	157	139	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	1A	99	Total	C	N	O	S	0	0	0
			801	504	157	139	1			

- Molecule 11 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	2I	119	Total	C	N	O	S	0	0	0
			885	549	168	165	3			
11	2A	119	Total	C	N	O	S	0	0	0
			885	549	168	165	3			

- Molecule 12 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	3I	125	Total	C	N	O	S	0	0	0
			975	614	196	164	1			
12	3A	125	Total	C	N	O	S	0	0	0
			975	614	196	164	1			

- Molecule 13 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	4I	118	Total	C	N	O	S	0	0	0
			938	580	193	163	2			
13	4A	117	Total	C	N	O	S	0	0	0
			933	577	192	162	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
4I	119	ALA	GLY	conflict	UNP P80377
4A	119	ALA	GLY	conflict	UNP P80377

- Molecule 14 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	5I	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			
14	5A	58	Total	C	N	O	S	0	0	0
			476	303	99	70	4			

- Molecule 15 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	6I	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			
15	6A	88	Total	C	N	O	S	0	0	0
			734	459	147	126	2			

- Molecule 16 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	7I	84	Total	C	N	O	S	0	0	0
			705	446	140	118	1			
16	7A	84	Total	C	N	O	S	0	0	0
			705	446	140	118	1			

- Molecule 17 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	8I	100	Total	C	N	O	S	0	0	0
			834	534	155	143	2			
17	8A	100	Total	C	N	O	S	0	0	0
			834	534	155	143	2			

- Molecule 18 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	9I	72	Total	C	N	O	0	0	0
			591	376	117	98			
18	9A	72	Total	C	N	O	0	0	0
			591	376	117	98			

- Molecule 19 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	AI	81	Total	C	N	O	S	0	0	0
			647	413	119	113	2			
19	AA	78	Total	C	N	O	S	0	0	0
			624	398	115	109	2			

- Molecule 20 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	BI	99	Total	C	N	O	S	0	0	0
			763	470	162	129	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	BA	99	Total	C	N	O	S	0	0	0
			763	470	162	129	2			

- Molecule 21 is a protein called 30S ribosomal protein Thx.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	1F	25	Total	C	N	O	0	0	0
			217	134	52	31			
21	1B	25	Total	C	N	O	0	0	0
			217	134	52	31			

- Molecule 22 is a RNA chain called tRNA-Phe.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
22	1K	76	Total	C	N	O	P	S	0	0	0
			1626	729	290	531	75	1			
22	1L	76	Total	C	N	O	P	S	0	0	0
			1626	729	290	531	75	1			

- Molecule 23 is a RNA chain called tRNA-fMet.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
23	2K	77	Total	C	N	O	P	S	0	0	0
			1646	735	298	535	77	1			
23	2L	77	Total	C	N	O	P	S	0	0	0
			1646	735	298	535	77	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
2K	18	C	U	conflict	GB 675817920
2L	18	C	U	conflict	GB 675817920

- Molecule 24 is a RNA chain called tRNA-Phe.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	3K	76	Total	C	N	O	P	0	0	0
			1619	723	290	531	75			
24	3L	76	Total	C	N	O	P	0	0	0
			1619	723	290	531	75			

- Molecule 25 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	4K	14	Total	C	N	O	P	0	0	0
			301	136	60	91	14			
25	4L	6	Total	C	N	O	P	0	0	0
			131	59	27	39	6			

- Molecule 26 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	1H	2912	Total	C	N	O	P	0	0	0
			62707	27911	11722	20163	2911			
26	14	2909	Total	C	N	O	P	0	0	0
			62647	27884	11716	20139	2908			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1H	161	U	-	insertion	GB 48268
1H	493	G	-	insertion	GB 48268
1H	1228	G	-	insertion	GB 48268
14	161	U	-	insertion	GB 48268
14	493	G	-	insertion	GB 48268
14	1228	G	-	insertion	GB 48268

- Molecule 27 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	16	122	Total	C	N	O	P	0	0	0
			2617	1166	486	844	121			
27	1J	122	Total	C	N	O	P	0	0	0
			2617	1166	486	844	121			

- Molecule 28 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	11	273	Total	C	N	O	S	0	0	0
			2120	1338	421	358	3			
28	19	273	Total	C	N	O	S	0	0	0
			2120	1338	421	358	3			

- Molecule 29 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	21	205	Total	C	N	O	S	0	0	0
			1568	991	300	271	6			
29	29	205	Total	C	N	O	S	0	0	0
			1568	991	300	271	6			

- Molecule 30 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	31	202	Total	C	N	O	S	0	0	0
			1585	1011	297	275	2			
30	39	208	Total	C	N	O	S	0	0	0
			1627	1037	304	283	3			

- Molecule 31 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	41	181	Total	C	N	O	S	0	0	0
			1473	942	268	259	4			
31	49	181	Total	C	N	O	S	0	0	0
			1473	942	268	259	4			

- Molecule 32 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	51	174	Total	C	N	O	S	0	0	0
			1336	848	251	236	1			
32	59	170	Total	C	N	O	S	0	0	0
			1307	829	245	232	1			

- Molecule 33 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	61	146	Total	C	N	O	S	0	0	0
			1136	726	201	208	1			
33	69	146	Total	C	N	O	S	0	0	0
			1136	726	201	208	1			

- Molecule 34 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	58	138	Total	C	N	O	S	0	0	0
			1104	712	206	182	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	15	138	Total	C	N	O	S	0	0	0
			1104	712	206	182	4			

- Molecule 35 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	68	122	Total	C	N	O	S	0	0	0
			933	588	171	170	4			
35	25	122	Total	C	N	O	S	0	0	0
			933	588	171	170	4			

- Molecule 36 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	78	150	Total	C	N	O	S	0	0	0
			1145	712	232	198	3			
36	35	150	Total	C	N	O	S	0	0	0
			1145	712	232	198	3			

- Molecule 37 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	88	138	Total	C	N	O	S	0	0	0
			1087	693	208	180	6			
37	45	141	Total	C	N	O	S	0	0	0
			1122	715	212	188	7			

- Molecule 38 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	98	118	Total	C	N	O	S	0	0	0
			968	604	203	160	1			
38	55	117	Total	C	N	O		0	0	0
			960	599	202	159				

- Molecule 39 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
39	A8	111	Total	C	N	O	0	0	0
			882	556	176	150			
39	65	111	Total	C	N	O	0	0	0
			882	556	176	150			

- Molecule 40 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	B8	137	Total	C	N	O	S	0	0	0
			1141	710	234	196	1			
40	75	137	Total	C	N	O	S	0	0	0
			1141	710	234	196	1			

- Molecule 41 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	C8	117	Total	C	N	O	S	0	0	0
			964	610	202	151	1			
41	85	117	Total	C	N	O	S	0	0	0
			964	610	202	151	1			

- Molecule 42 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	D8	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			
42	95	101	Total	C	N	O	S	0	0	0
			779	501	142	135	1			

- Molecule 43 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
43	E8	113	Total	C	N	O	S	0	0	0
			900	566	177	155	2			
43	A5	113	Total	C	N	O	S	0	0	0
			900	566	177	155	2			

- Molecule 44 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	F8	94	Total	C	N	O	S	0	0	0
			742	482	134	125	1			
44	B5	93	Total	C	N	O		0	0	0
			730	474	132	124				

- Molecule 45 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
45	G8	104	Total	C	N	O	S	0	0	0
			791	510	149	127	5			
45	C5	104	Total	C	N	O	S	0	0	0
			794	510	152	127	5			

- Molecule 46 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	H8	175	Total	C	N	O	S	0	0	0
			1397	892	251	251	3			
46	D5	179	Total	C	N	O	S	0	0	0
			1428	911	255	259	3			

- Molecule 47 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	I8	80	Total	C	N	O	S	0	0	0
			627	388	132	106	1			
47	E5	77	Total	C	N	O	S	0	0	0
			613	379	129	104	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I8	6	ALA	GLY	conflict	UNP P60493
I8	8	ALA	GLY	conflict	UNP P60493
E5	6	ALA	GLY	conflict	UNP P60493
E5	8	ALA	GLY	conflict	UNP P60493

- Molecule 48 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
48	J8	97	Total	C	N	O	S	0	0	0
			763	481	150	131	1			
48	F5	97	Total	C	N	O	S	0	0	0
			763	481	150	131	1			

- Molecule 49 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	K8	67	Total	C	N	O	S	0	0	0
			563	349	114	99	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	G5	66	Total	C	N	O	S	0	0	0
			558	346	113	98	1			

- Molecule 50 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	L8	57	Total	C	N	O		0	0	0
			452	288	88	76				
50	H5	59	Total	C	N	O		0	0	0
			469	298	90	81				

- Molecule 51 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
51	M8	66	Total	C	N	O	S	0	0	0
			533	335	96	97	5			
51	I5	63	Total	C	N	O	S	0	0	0
			515	326	93	91	5			

- Molecule 52 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	N8	58	Total	C	N	O	S	0	0	0
			454	285	89	75	5			
52	J5	59	Total	C	N	O	S	0	0	0
			459	288	90	76	5			

- Molecule 53 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	O8	45	Total	C	N	O	S	0	0	0
			389	241	79	65	4			
53	K5	45	Total	C	N	O	S	0	0	0
			389	241	79	65	4			

- Molecule 54 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
54	P8	48	Total	C	N	O	S	0	0	0
			418	257	104	55	2			
54	L5	46	Total	C	N	O	S	0	0	0
			398	245	98	53	2			

- Molecule 55 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
55	Q8	60	Total	C	N	O	S	0	0	0
			480	306	98	74	2			
55	M5	60	Total	C	N	O	S	0	0	0
			477	303	98	74	2			

- Molecule 56 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
56	98	1	Total	Mg	0	0
			1	1		
56	P8	1	Total	Mg	0	0
			1	1		
56	85	1	Total	Mg	0	0
			1	1		
56	C5	1	Total	Mg	0	0
			1	1		
56	13	151	Total	Mg	0	0
			151	151		
56	1J	9	Total	Mg	0	0
			9	9		
56	35	1	Total	Mg	0	0
			1	1		
56	C8	1	Total	Mg	0	0
			1	1		
56	75	1	Total	Mg	0	0
			1	1		
56	16	13	Total	Mg	0	0
			13	13		
56	21	2	Total	Mg	0	0
			2	2		
56	31	1	Total	Mg	0	0
			1	1		
56	Q8	1	Total	Mg	0	0
			1	1		
56	L8	1	Total	Mg	0	0
			1	1		
56	3I	1	Total	Mg	0	0
			1	1		
56	I8	2	Total	Mg	0	0
			2	2		
56	5E	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
56	29	2	Total 2	Mg 2	0	0
56	2K	3	Total 3	Mg 3	0	0
56	J8	2	Total 2	Mg 2	0	0
56	4A	1	Total 1	Mg 1	0	0
56	AI	1	Total 1	Mg 1	0	0
56	1G	103	Total 103	Mg 103	0	0
56	11	4	Total 4	Mg 4	0	0
56	E8	1	Total 1	Mg 1	0	0
56	1H	529	Total 529	Mg 529	0	0
56	F5	1	Total 1	Mg 1	0	0
56	88	2	Total 2	Mg 2	0	0
56	49	1	Total 1	Mg 1	0	0
56	14	389	Total 389	Mg 389	0	0
56	78	2	Total 2	Mg 2	0	0
56	3E	2	Total 2	Mg 2	0	0
56	6A	1	Total 1	Mg 1	0	0
56	1K	2	Total 2	Mg 2	0	0
56	41	2	Total 2	Mg 2	0	0
56	2L	2	Total 2	Mg 2	0	0

- Molecule 57 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
57	32	1	Total Zn 1 1	0	0
57	3E	1	Total Zn 1 1	0	0
57	5I	1	Total Zn 1 1	0	0
57	5A	1	Total Zn 1 1	0	0
57	G8	1	Total Zn 1 1	0	0
57	C5	1	Total Zn 1 1	0	0

- Molecule 58 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
58	13	212	Total O 212 212	0	0
58	3E	1	Total O 1 1	0	0
58	4E	2	Total O 2 2	0	0
58	8E	2	Total O 2 2	0	0
58	1I	2	Total O 2 2	0	0
58	3I	1	Total O 1 1	0	0
58	5I	2	Total O 2 2	0	0
58	1K	1	Total O 1 1	0	0
58	3K	1	Total O 1 1	0	0
58	4K	4	Total O 4 4	0	0
58	1H	1097	Total O 1097 1097	0	0
58	16	16	Total O 16 16	0	0
58	11	9	Total O 9 9	0	0
58	21	5	Total O 5 5	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
58	31	10	Total 10	O 10	0	0
58	58	1	Total 1	O 1	0	0
58	78	4	Total 4	O 4	0	0
58	98	1	Total 1	O 1	0	0
58	C8	2	Total 2	O 2	0	0
58	D8	1	Total 1	O 1	0	0
58	E8	2	Total 2	O 2	0	0
58	G8	3	Total 3	O 3	0	0
58	I8	6	Total 6	O 6	0	0
58	L8	2	Total 2	O 2	0	0
58	P8	2	Total 2	O 2	0	0
58	Q8	2	Total 2	O 2	0	0
58	1G	99	Total 99	O 99	0	0
58	7A	2	Total 2	O 2	0	0
58	BA	1	Total 1	O 1	0	0
58	14	730	Total 730	O 730	0	0
58	1J	12	Total 12	O 12	0	0
58	19	11	Total 11	O 11	0	0
58	29	6	Total 6	O 6	0	0
58	39	7	Total 7	O 7	0	0
58	55	1	Total 1	O 1	0	0

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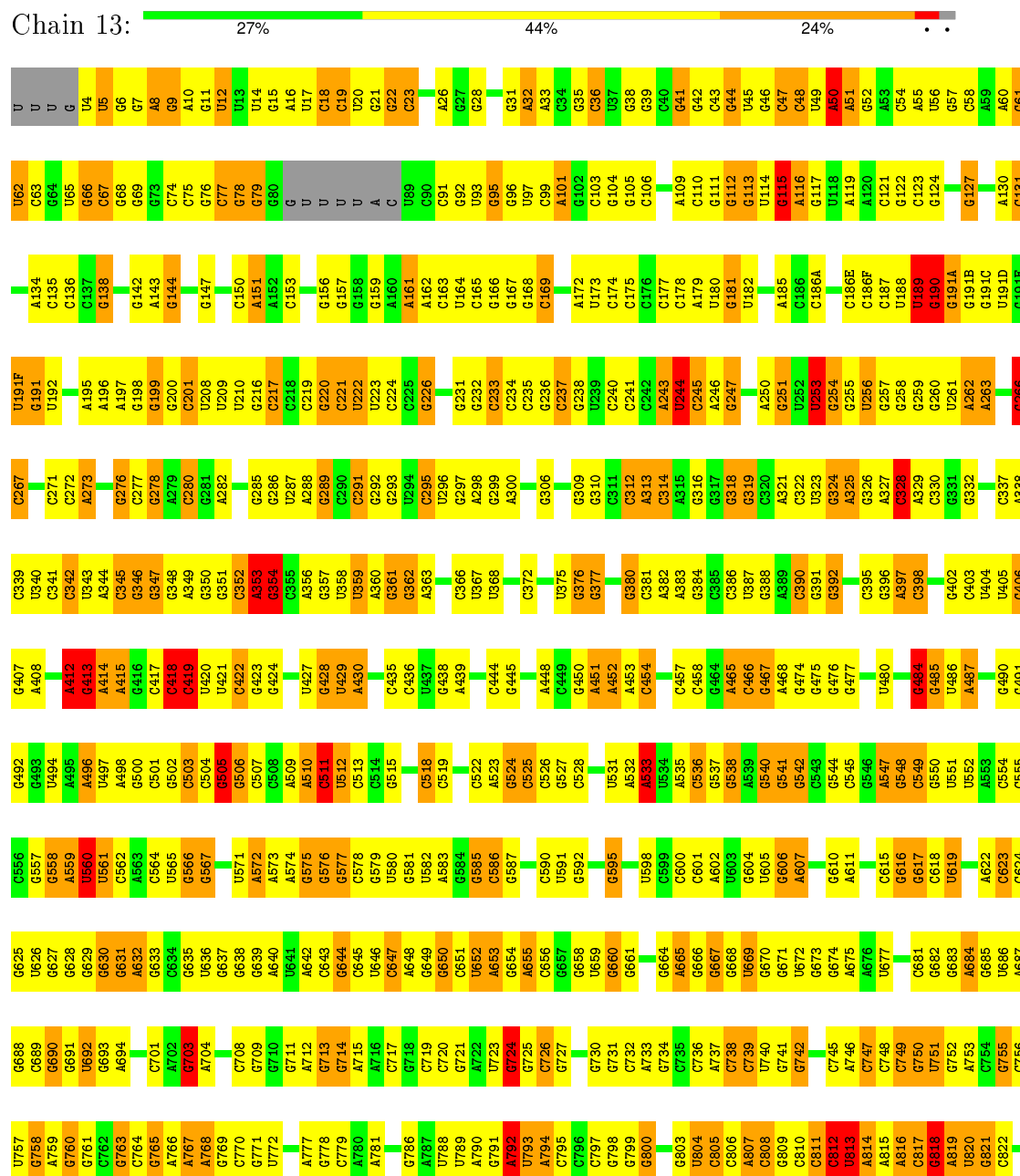
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
58	85	1	Total	O	0	0
			1	1		
58	A5	1	Total	O	0	0
			1	1		
58	H5	1	Total	O	0	0
			1	1		
58	L5	1	Total	O	0	0
			1	1		

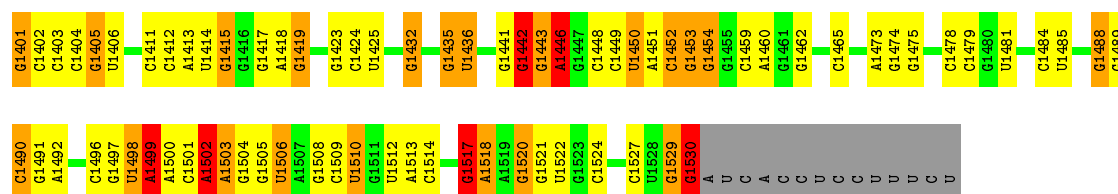
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($\text{RSRZ} > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

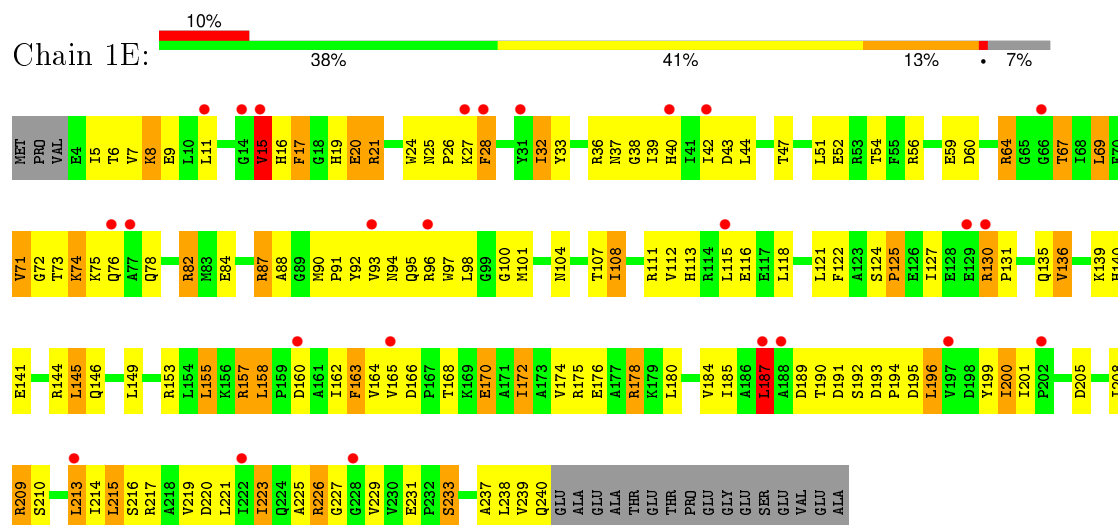
- Molecule 1: 16S ribosomal RNA



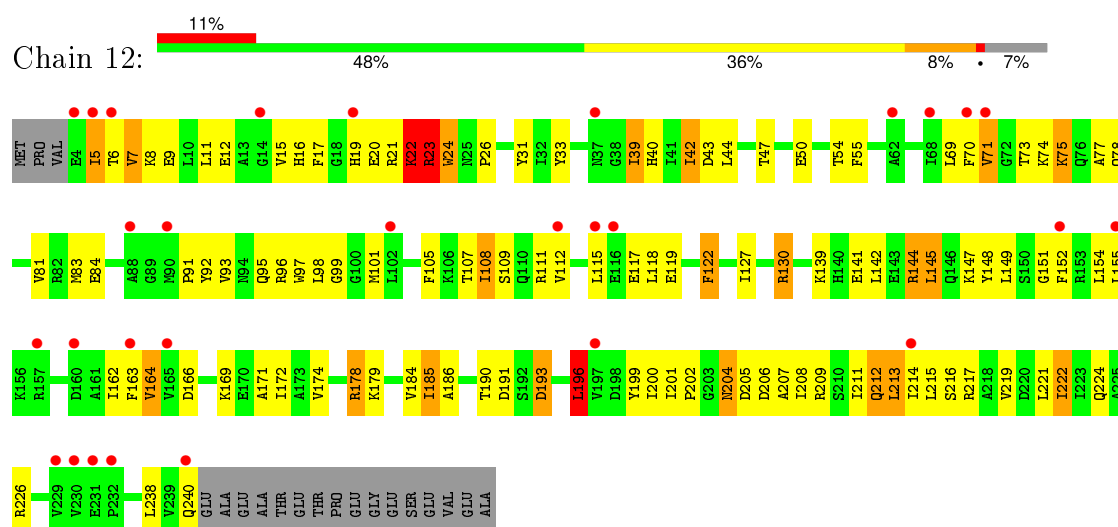
C1336	U1205	G1139	G1009	G947	C982	C808	C736	G662	C589	C525	A374	G305
G1275	G1206	C1140	G1010	C948	C983	C811	A737	A653	C590	C526	A374	G306
G1276	G1207	C1141	G1013	U950	U884	C812	C738	G664	C592	G527	U375	G309
C1208	C1208	G1142	G1074	G985	G1074	C812	C738	A665	C592	C444	G376	G310
U1211	U1211	G1143	G1075	G986	G986	C812	C738	G666	C596	G529	G377	
U1212	U1212	G1144	G1078	U952	G887	C814	G741	G667	C597	G530	A382	A313
A1213	A1213	G1145	U1079	G953	G888	C815	U743	G668	C598	U531	A383	C314
U1281	C1214	A1146	A1080	G954	A889	C816	C744	U669	C599	A532	A384	A315
G1282	G1215	C1147	U1079	U955	G890	C817	C744	G670	C600	U534	G388	G316
G1283	G1216	U1148	G1081	U956	U891	C818	A746	G671	C601	A451	A389	G317
G1284	G1217	U1149	G1082	U957	A892	C819	C747	G672	C602	A452	C390	G318
G1285	G1218	U1150	U1083	A958	C893	C820	C748	G673	G604	A453	G391	G319
G1286	G1219	U1151	G1084	A959	G894	C821	C749	G674	U605	C457	G392	C320
A1287	U1219	A1152	U1085	U960	G895	C822	G750	A675	C606	C458		A321
A1288	G1220	C1153	U1086	U961	C996	C823		U677	C607	C459		
A1289		C1154		C962	C997	C826	C754		A607	C460	G396	C322
A1290	G1224	G1155	A1092	G963	G898	C827	G755	C680	A608	G541	C397	U323
G1291	A1225	G1156	A1093	A964	C899	C828	G756	G681	A609	G544	C398	G324
U1292	C1226	A1157	G1094	A965	A900	C829	U757	G682		C545	C399	A325
G1293	A1227	C1158	U1095	G966	A901	C830	G758	A684	A614	C546	C400	G326
G1294	C1228	U1159	C1096	G967	G902	C831	A759	U685	C615	C547	C401	A327
G1295	G1231	C1160	C1097	A968	C903	U831	G760	A687	C616	A547	C402	C328
A1296	U1232	C1161	C1098	A969	G904	C832	G761	G688	C617	C475	C403	A329
G1297	U1233	C1162	G1099	C970	U905	C833	G762	C689	U618	U552	U404	C330
G1298	G1234	C1163	A1101	G971	G906	C834	U772	C690	U619	A553	U405	A331
A1299	U1235	G1171	A1102	G972	A909	C835	G773	G691	C620	C554	G406	G332
G1300	A1236	C1172	C1103	A974	C910	C836	A776	G692	A621	C555	G407	G333
U1301	C1237	G1173	G1104	A975	U911	C837	A777	U686	C622	C556	A408	C334
G1302	A1238	G1174	A1105	G976	C912	C838	G768	G693	C623	G557	C409	
G1303	C1239	G1175	C1106	A977	A913	C839	G769	A694	C624	G558	G410	G337
G1304	U1240	A1176	G1107	A978	A914	C840	U772	U697	C625	A559	A411	A338
G1305	G1241	G1177	G1108	C979	A915	C841			U626	U560	A412	C339
U1307	C1242	A1178	A1111	U982	G916	C842	G775	C701	C627	C561	G413	U340
G1310	U1247	G1180	C1112	U983	G917	C843	G776	A702	C628	C562	A414	C341
G1311	A1248	G1181	C1113	A984	A918	C844	A777	G703	G630	C563	G416	C345
G1312	C1249	G1182	C1114	C984	A919	C845	G778	A704	G631	U565	C417	G346
U1313	A1250	A1183	C1115	C985	G922	C846	C779	C707	G633	C566	C418	G347
G1314	A1251	G1184	C1116	A986	A923	C847	A780	C708		C567	C419	G348
A1252	C1252	G1185	C1117	G987	C924	C848	A782	G709	U636	G568	U420	A349
G1315	G1253	G1186	C1118	G988	C925	C849	G783	G710	C637	C569	U421	G350
G1316	C1254	G1187	C1119	U991	G926	C850	G784	G713	G638	U571	C422	G351
G1317	U1255	A1188	U1122	U992	G927	C851	G785	G714	C639	C502	G423	C352
A1318	A1256	C1189	A1123	G993	C932	C852		G715	A640	C503	G424	C353
A1319	U1257	G1190	A1124	A994	C933	C853		A716	U641	C504	G425	C354
C1321	G1258	A1191	U1055	C995	G934	C854	G791	C717	A642	A574	G426	C355
G1322	C1259	G1192	U1056	G998	A935	C855	A792	C720	U646	G575	U427	A356
A1323	C1260	G1193	U1126	C998A	A936	C856	U793	G721	C647	G577	G428	G357
A1324	U1261	U1194	G1058	U999	C937	C857	A794	A722	C648	C578	U429	U358
C1325	G1262	C1195	C1059	U999	A937	C858		A723	G649	U512	A430	U359
C1326	C1263	U1196	U1060	A1000	A938	C859	G798	G724	U650	C513	A431	A360
C1327		G1197	G1061	G1001	G939	C860		G725	G651	G579	A432	G361
G1328	C1267	U1198	U1062	G1002	C940	C861	U801	A802	U652	U580	C433	G362
A1329	A1268	U1199	C1063	G1003	G941	C862	A803	G803	U653	C581	U434	A363
U1330	C1270	C1200	G1064	A1004	G942	C863	A804	A729	A653	C582	C435	
G1331	G1271	U1201	U1065	A1005	U943	C864	U804	G730	G585	G584	C436	C366
	U1336	G1202	U1066	C1006	G944	C865	C806	G731	C586	C522	U437	U367
G1334	C1272	C1203	A1067	C1007	G945	C866	C807	C808	G587	A523	G438	G371
C1335	G1273	A1204	G1068	C1008	A946	C867	A807	C735	G588	G524	A440	C372



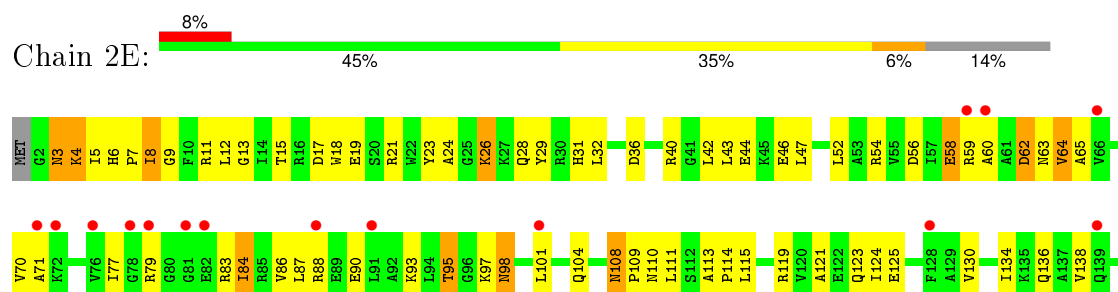
• Molecule 2: 30S ribosomal protein S2

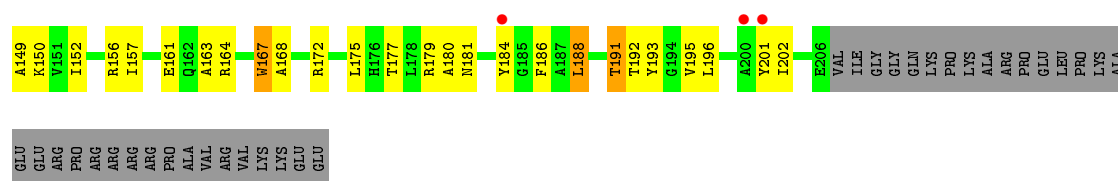


• Molecule 2: 30S ribosomal protein S2

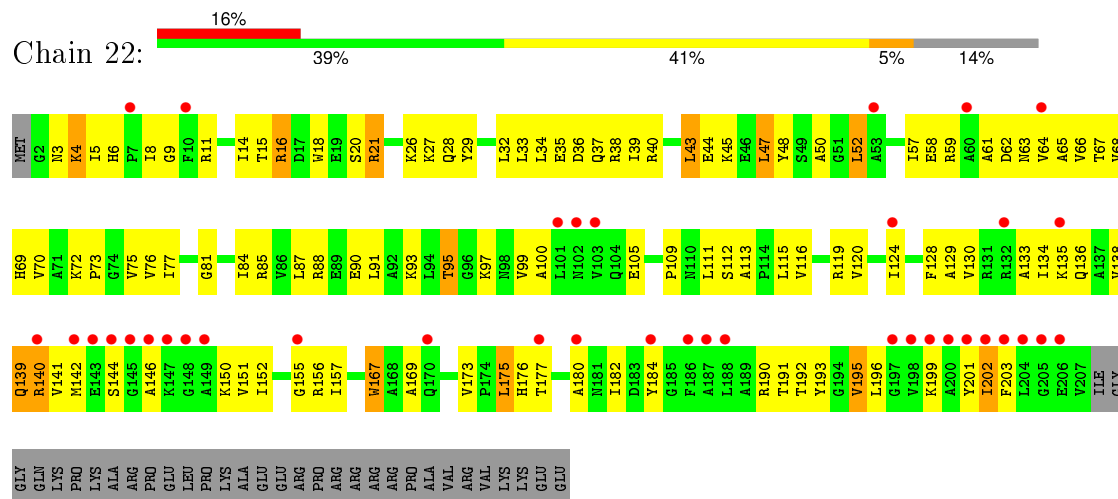


• Molecule 3: 30S ribosomal protein S3

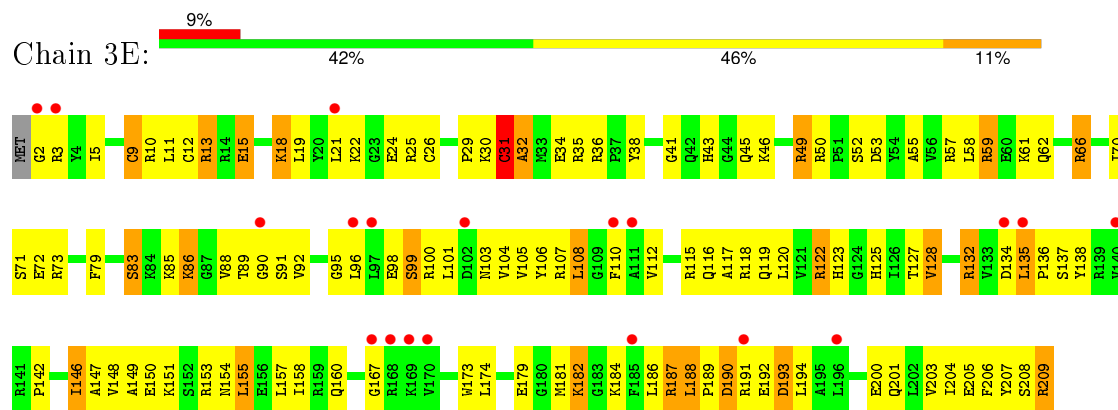




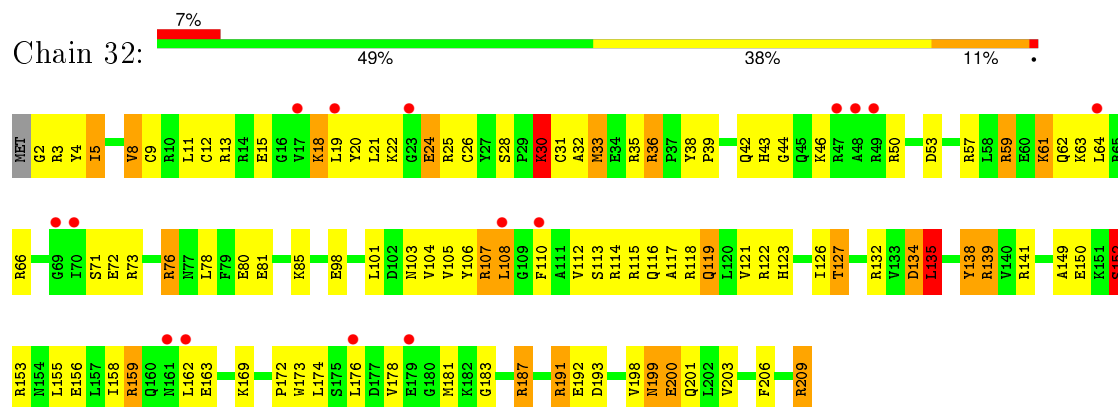
• Molecule 3: 30S ribosomal protein S3



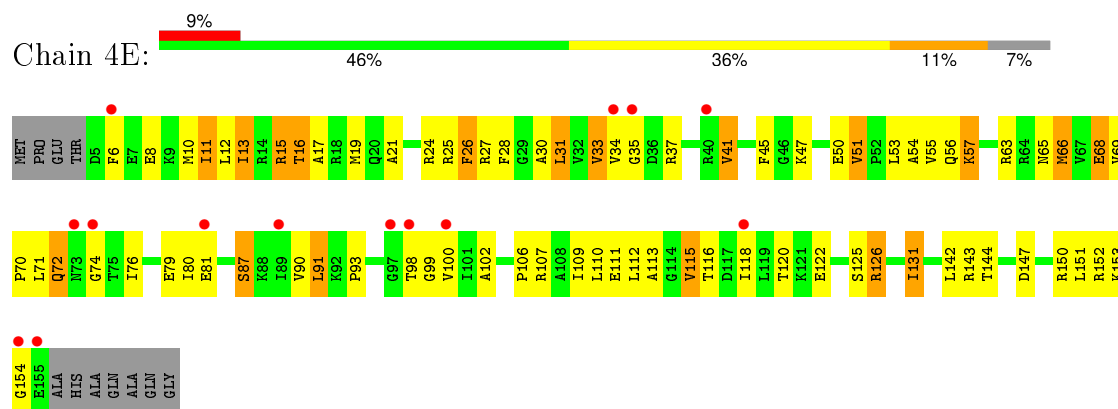
• Molecule 4: 30S ribosomal protein S4



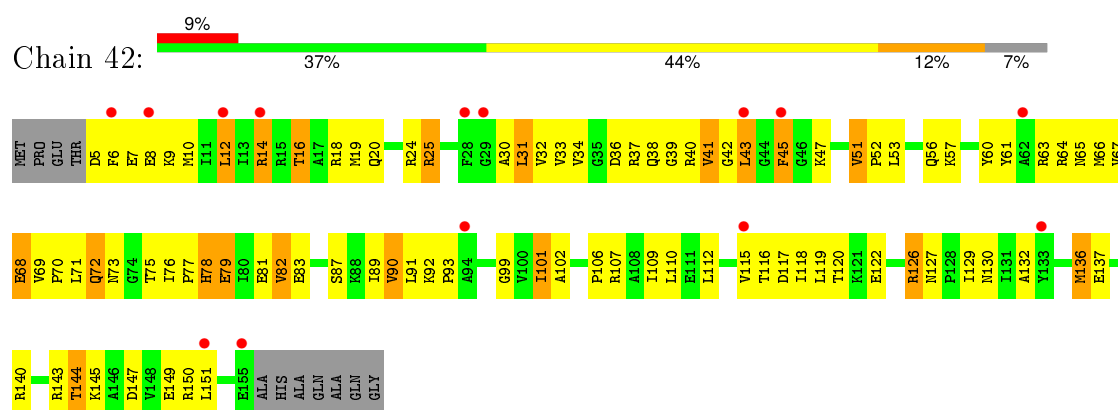
• Molecule 4: 30S ribosomal protein S4



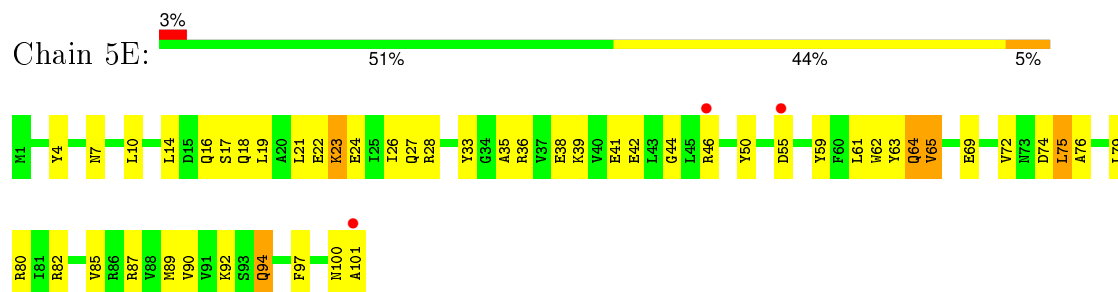
- Molecule 5: 30S ribosomal protein S5



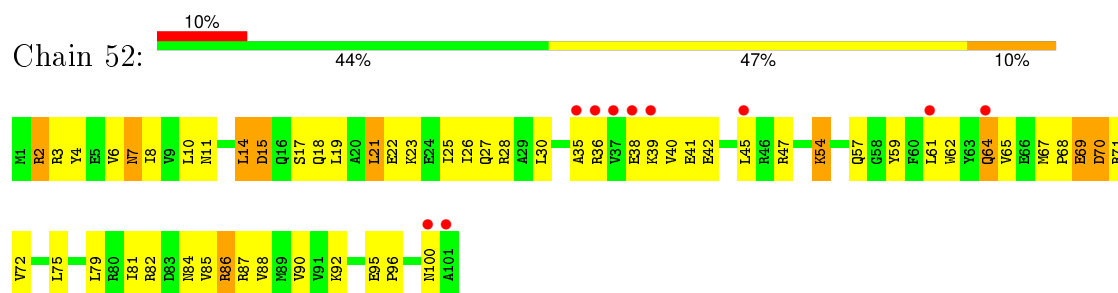
- Molecule 5: 30S ribosomal protein S5



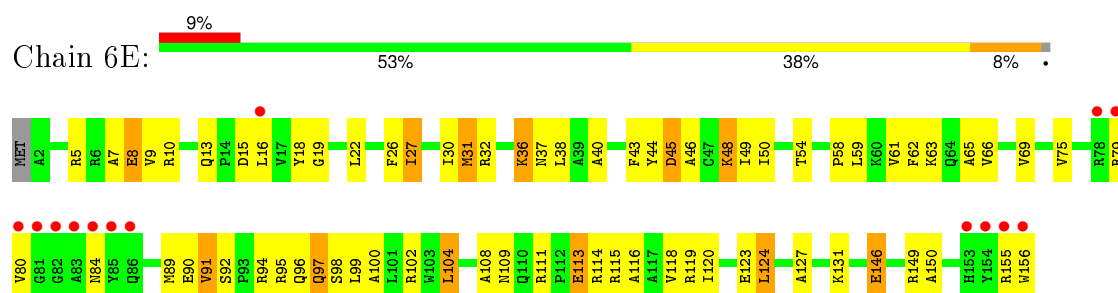
- Molecule 6: 30S ribosomal protein S6



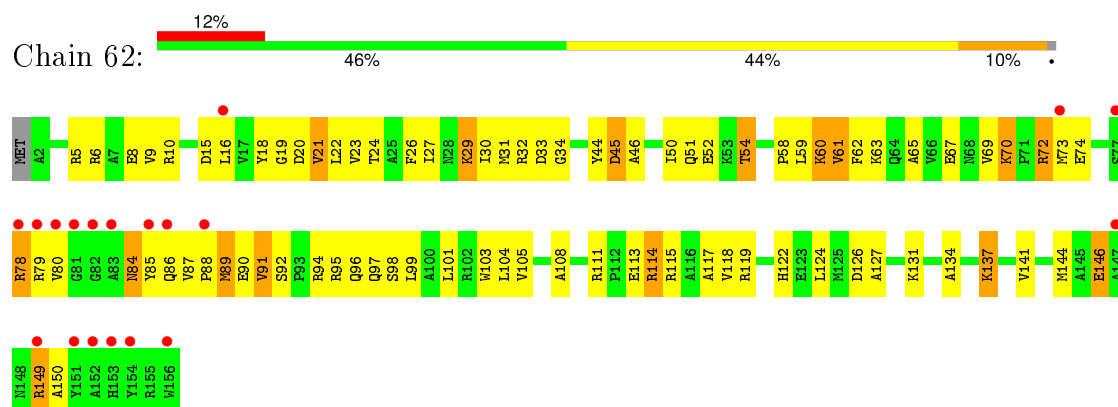
- Molecule 6: 30S ribosomal protein S6



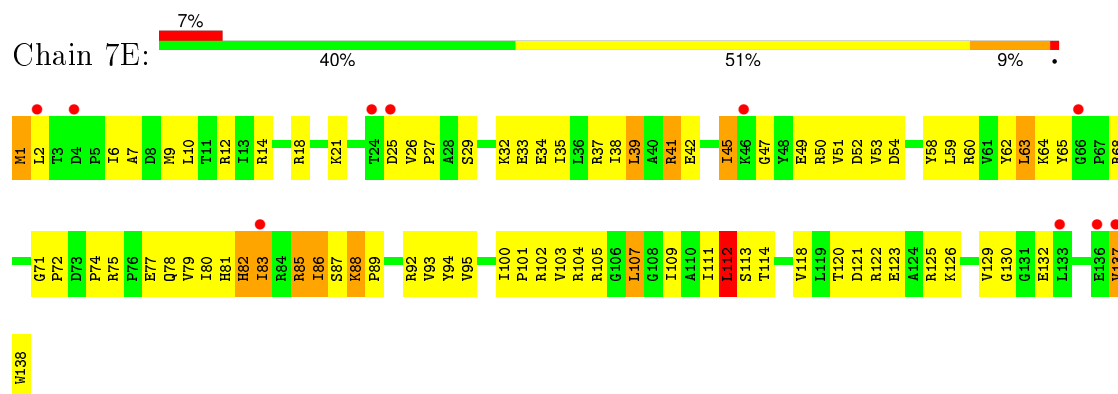
- Molecule 7: 30S ribosomal protein S7



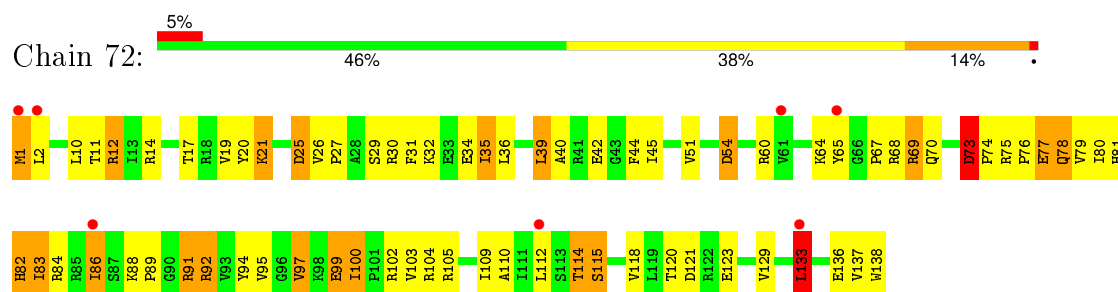
• Molecule 7: 30S ribosomal protein S7



• Molecule 8: 30S ribosomal protein S8

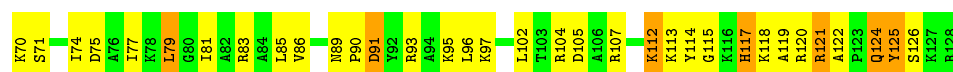


• Molecule 8: 30S ribosomal protein S8

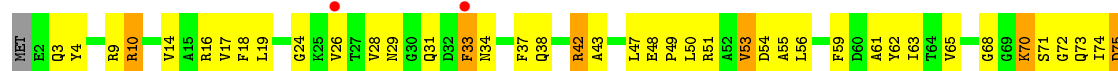


• Molecule 9: 30S ribosomal protein S9

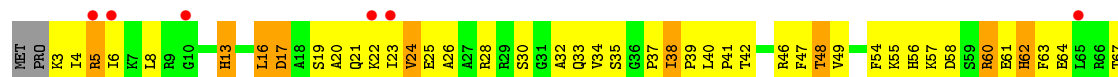




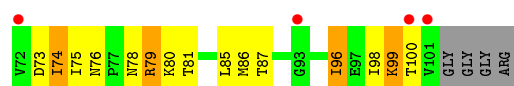
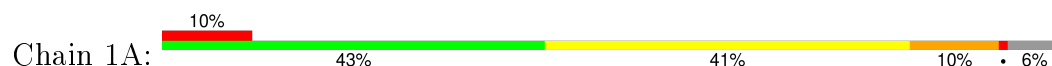
• Molecule 9: 30S ribosomal protein S9



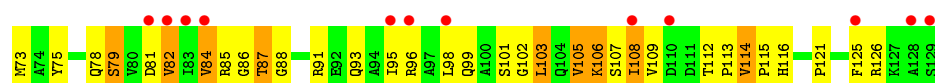
• Molecule 10: 30S ribosomal protein S10



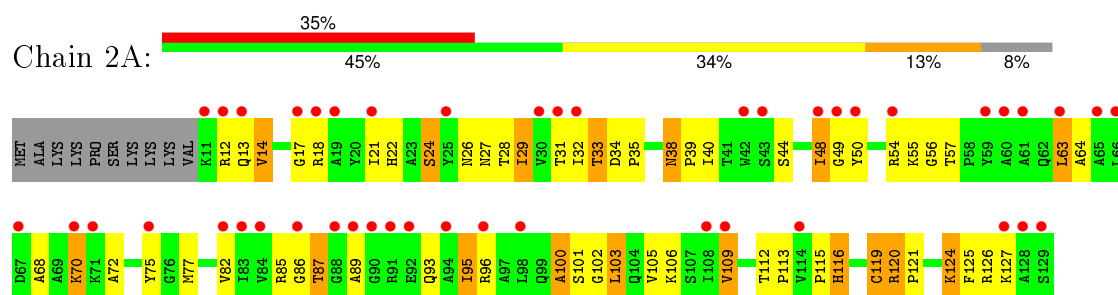
• Molecule 10: 30S ribosomal protein S10



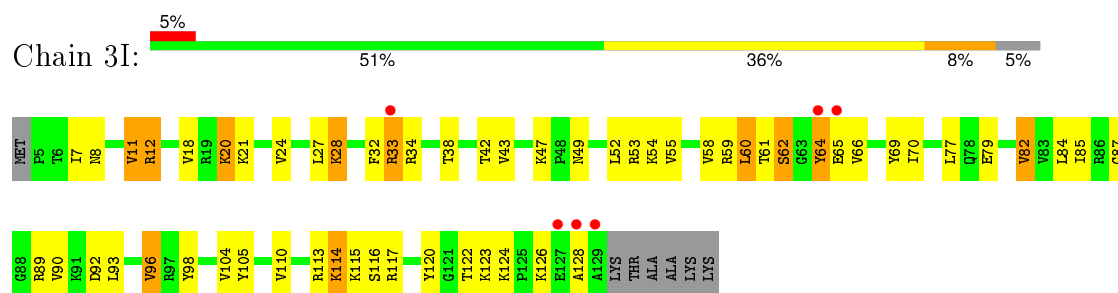
• Molecule 11: 30S ribosomal protein S11



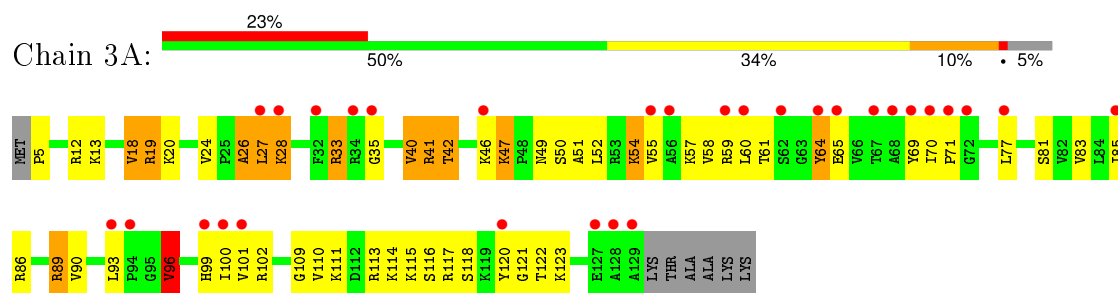
• Molecule 11: 30S ribosomal protein S11



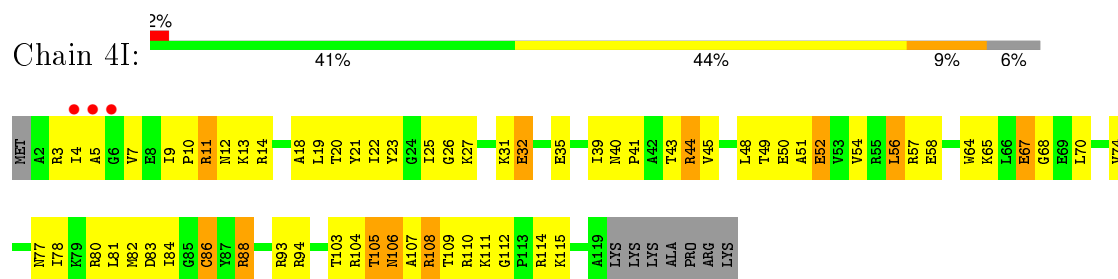
- Molecule 12: 30S ribosomal protein S12



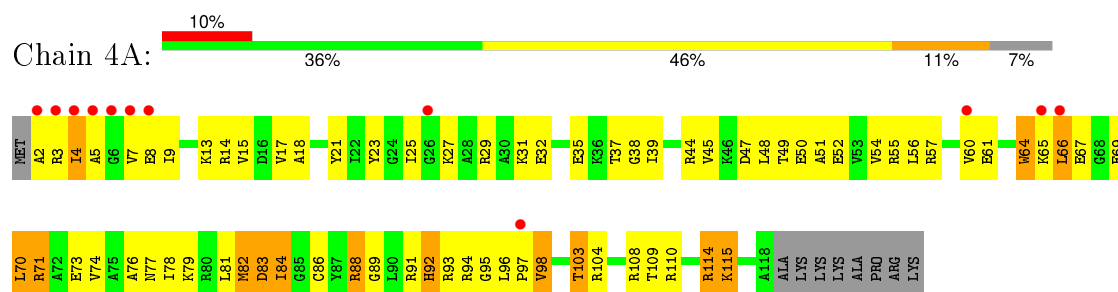
- Molecule 12: 30S ribosomal protein S12



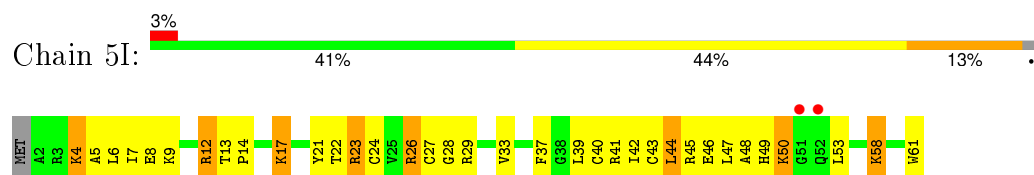
- Molecule 13: 30S ribosomal protein S13



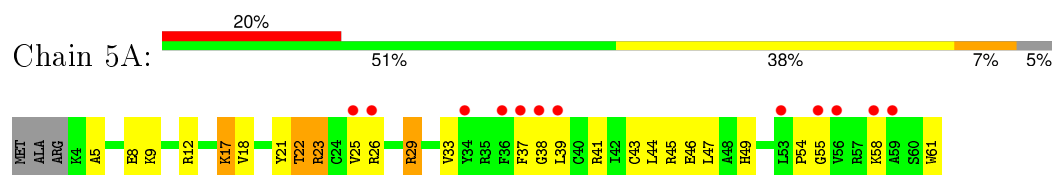
- Molecule 13: 30S ribosomal protein S13



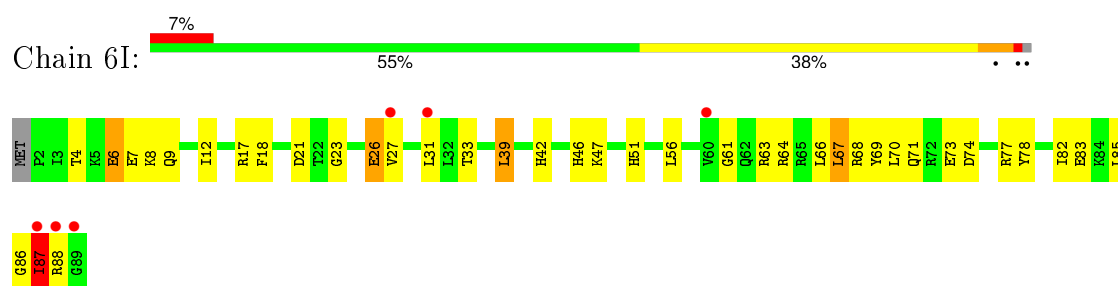
- Molecule 14: 30S ribosomal protein S14 type Z



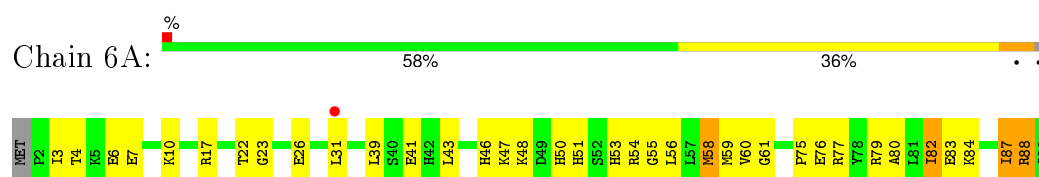
- Molecule 14: 30S ribosomal protein S14 type Z



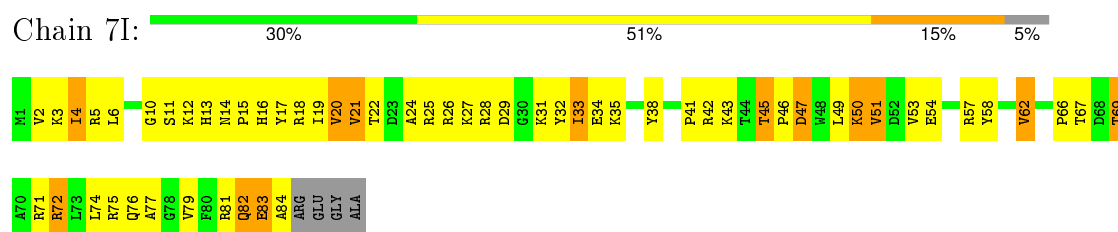
- Molecule 15: 30S ribosomal protein S15



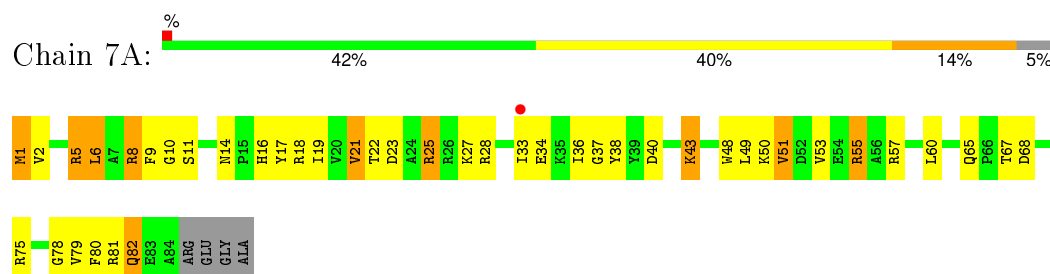
- Molecule 15: 30S ribosomal protein S15



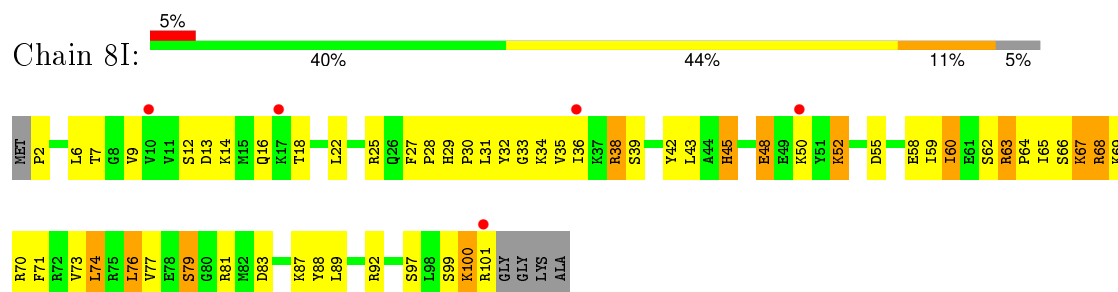
- Molecule 16: 30S ribosomal protein S16



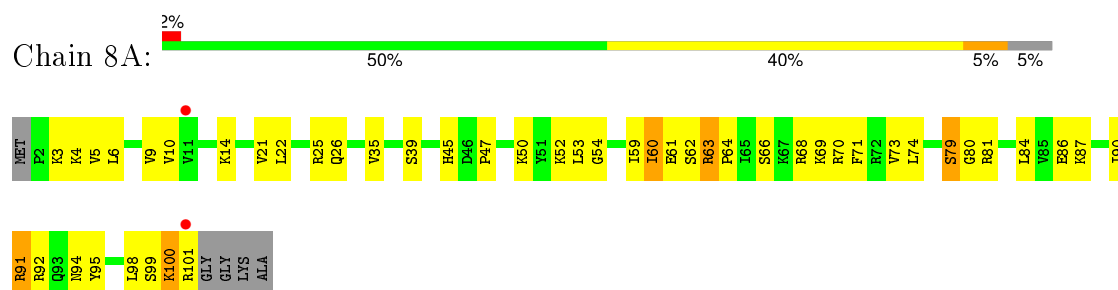
- Molecule 16: 30S ribosomal protein S16



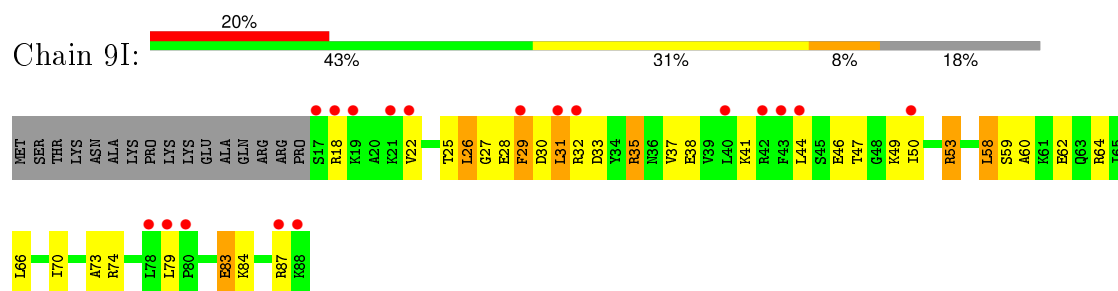
- Molecule 17: 30S ribosomal protein S17



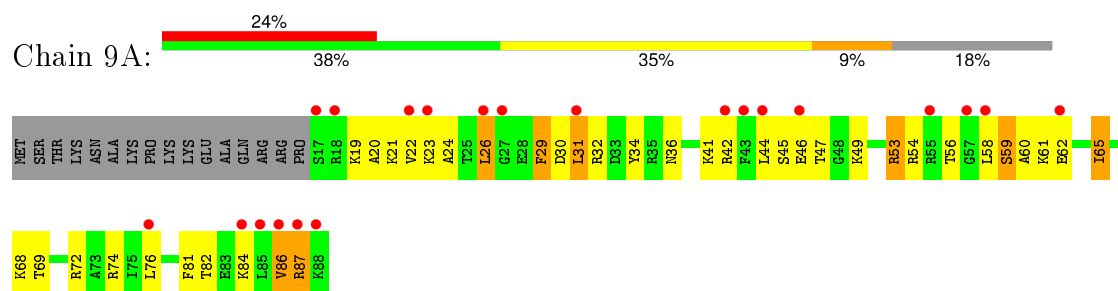
- Molecule 17: 30S ribosomal protein S17



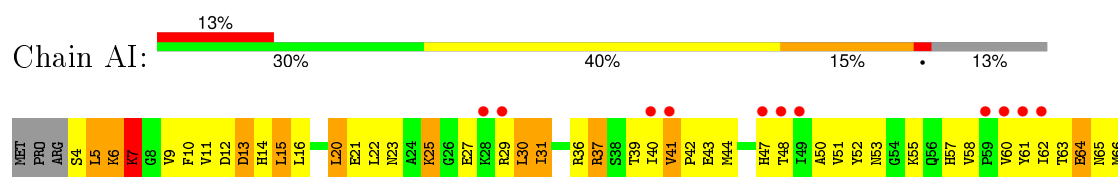
- Molecule 18: 30S ribosomal protein S18

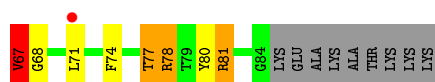


- Molecule 18: 30S ribosomal protein S18

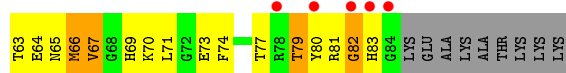
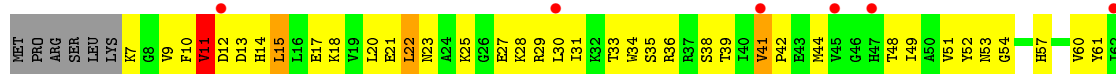


- Molecule 19: 30S ribosomal protein S19





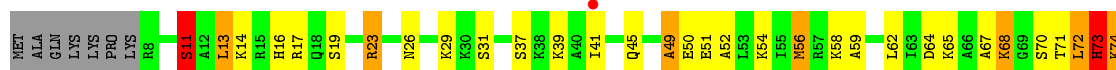
- Molecule 19: 30S ribosomal protein S19



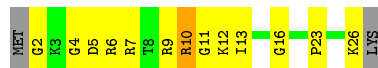
- Molecule 20: 30S ribosomal protein S20



- Molecule 20: 30S ribosomal protein S20



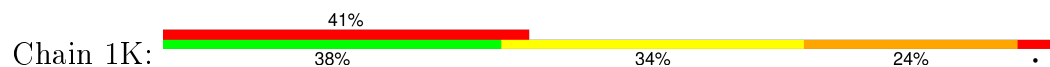
- Molecule 21: 30S ribosomal protein Thx

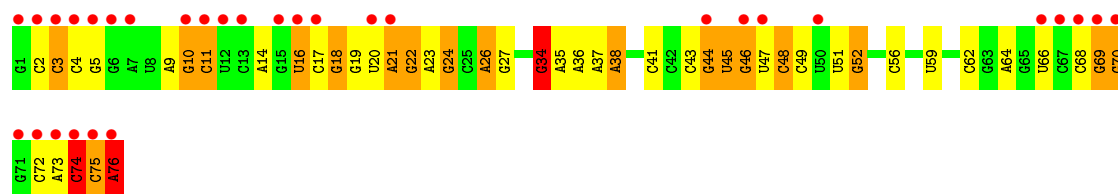


- Molecule 21: 30S ribosomal protein Thx

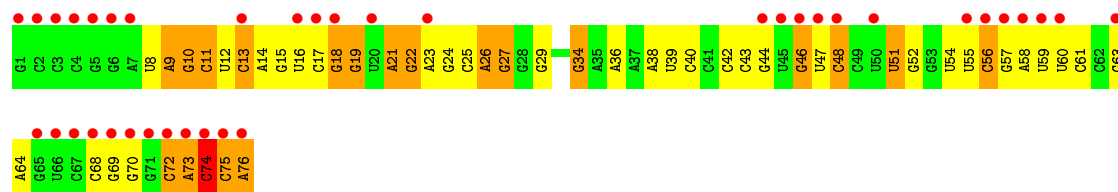


- Molecule 22: tRNA-Phe

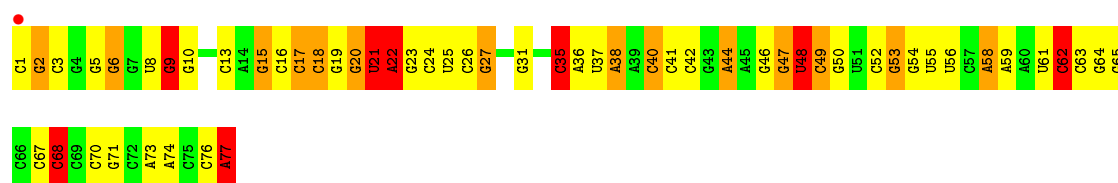




• Molecule 22: tRNA-Phe



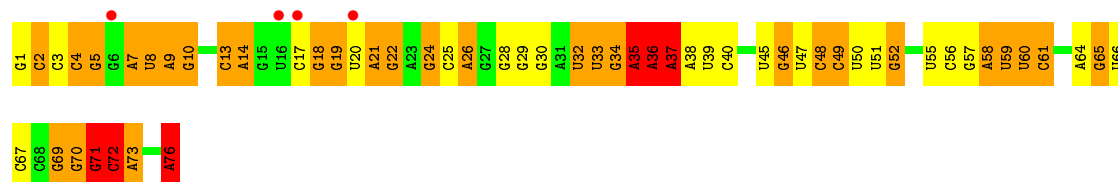
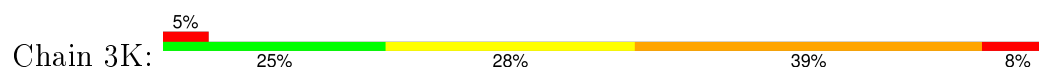
• Molecule 23: tRNA-fMet



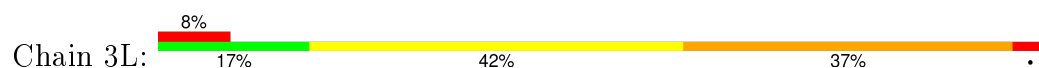
• Molecule 23: tRNA-fMet

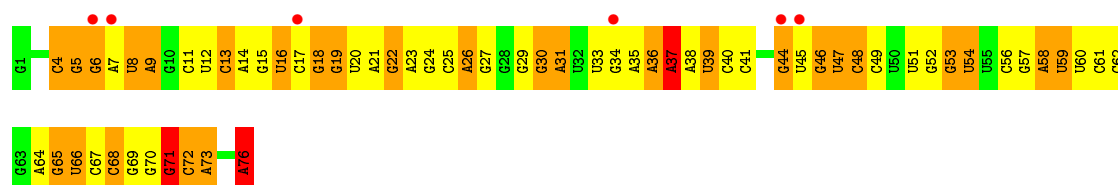


• Molecule 24: tRNA-Phe



• Molecule 24: tRNA-Phe





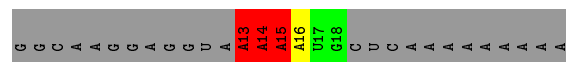
• Molecule 25: mRNA

Chain 4K: . 27% 17% 53%



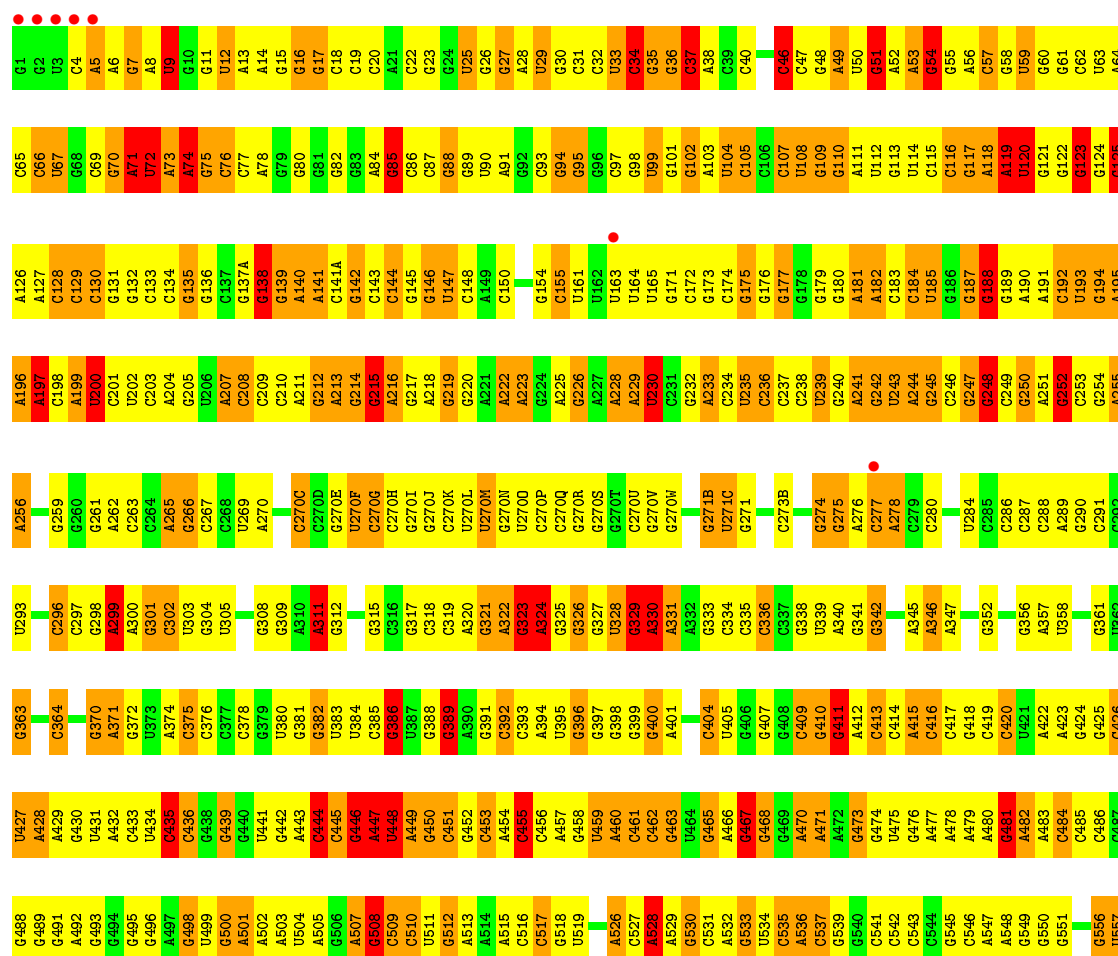
• Molecule 25: mRNA

Chain 4L: 7% 10% 80%



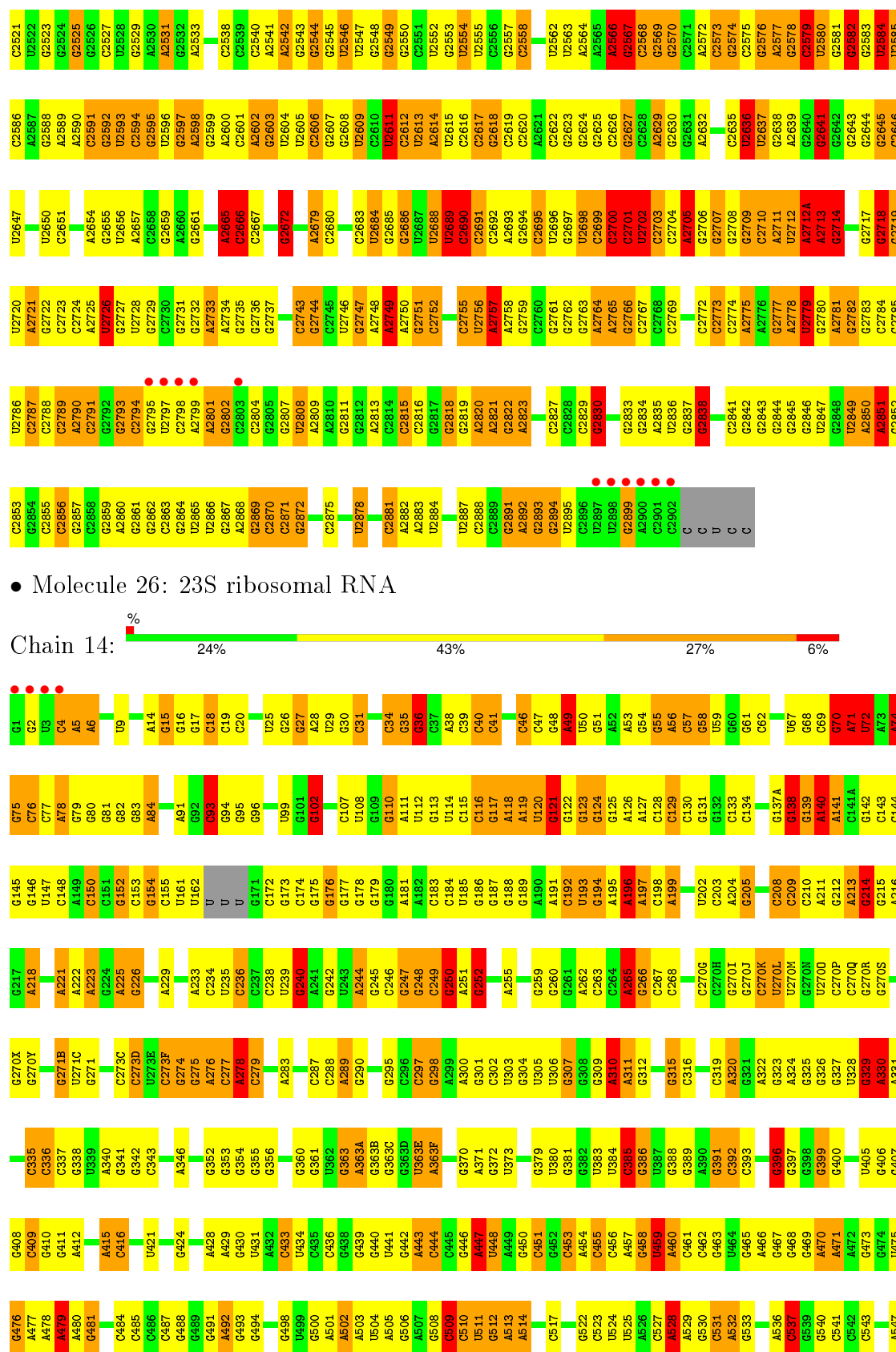
• Molecule 26: 23S ribosomal RNA

Chain 1H: 18% 41% 32% 8%



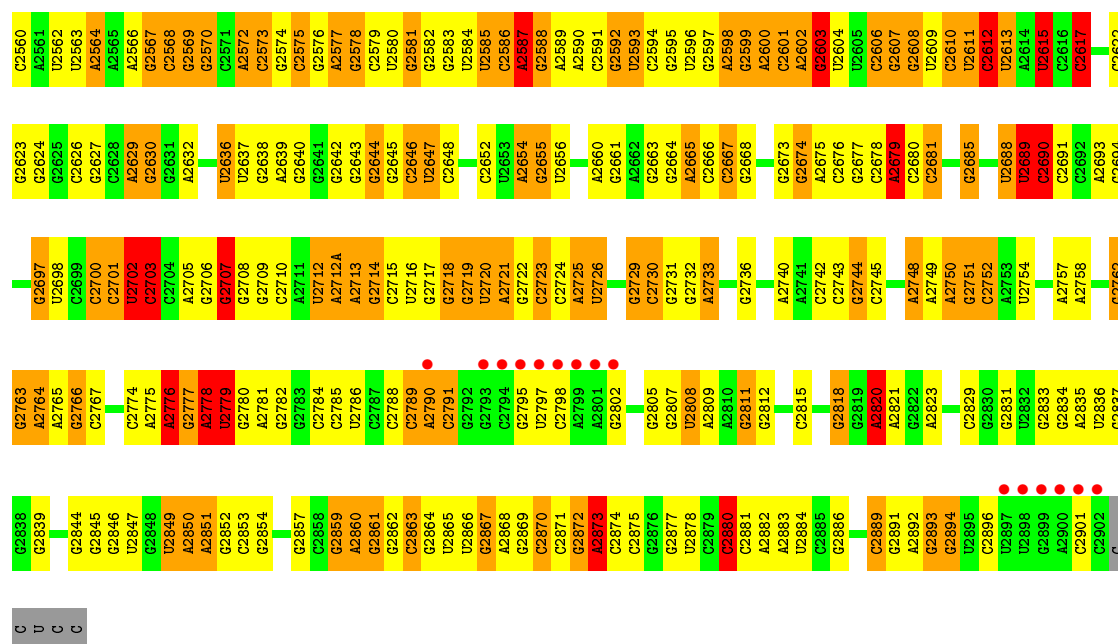
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G1283	A1284	G1285	A1286	G1287	G1288	G1289	G1290	G1291	G1292	G1293	G1294	G1295	G1296	G1297	G1298	G1299	G1300	A1301	G1302	G1303	G1304	G1305	G1306	G1307	G1308	G1309	G1310	G1311	G1312	G1313	G1314	G1315	G1316	G1317	G1318	G1319	G1320	A1321	G1322	G1323	G1324	G1325	G1326	G1327	G1328	G1329	G1330	G1331	G1332	G1333	G1334	A1335	G1336	G1337	G1338	G1339	G1340	G1341	G1342	A1343			
G1223	G1224	G1225	G1226	G1227	G1228	G1229	G1230	G1231	G1232	G1233	G1234	G1235	G1236	G1237	G1238	G1239	G1240	A1241	A1242	G1243	G1244	G1245	G1246	G1247	G1248	G1249	G1250	G1251	G1252	G1253	A1254	G1255	G1256	G1257	G1258	G1259	G1260	G1261	G1262	G1263	A1264	A1265	G1266	G1267	G1268	A1269	G1270	G1271	A1272	G1273	A1274	A1275	A1276	G1277	G1278	G1279	G1280	G1281	G1282				
G1161	G1162	G1163	G1164	G1165	G1166	G1167	G1168	G1169	G1170	G1171	G1172	G1173	G1174	G1175	G1176	G1177	G1178	G1179	G1180	G1181	G1182	G1183	G1184	G1185	G1186	G1187	G1188	G1189	G1190	G1191	G1192	G1193	A1194	G1195	G1196	G1197	G1198	G1199	G1200	G1201	G1202	G1203	A1204	G1205	G1206	G1207	G1208	G1209	A1210	A1211	A1212	G1213	G1214	G1215	G1216	G1217	G1218	G1219	A1220	G1221	G1222		
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C617	C618	C619	C620	C621	C622	C623	C624	C625	C626	C627	C628	C629	C630	C631	C632	C633	C634	C635	C636	C637	C638	C639	C640	C641	C642	C643	C644	C645	C646	C647	C648	C649	C650	C651	C652	C653	C654	C655	C656	C657	C658	C659	C660	C661	C662	C663	C664	C665	C666	C667	C668	C669	C670	C671	C672	C673	C674	C675					
G558	G559	C560	C561	C562	C563	C564	C565	C566	C567	C568	C569	C570	C571	C572	C573	C574	C575	C576	C577	C578	C579	C580	C581	C582	C583	C584	C585	C586	C587	C588	C589	C590	C591	C592	C593	C594	C595	C596	C597	C598	C599	C600	C601	C602	C603	C604	C605	C606	C607	C608	C609	C610	C611	C612	C613	C614	C615	C616	C617				



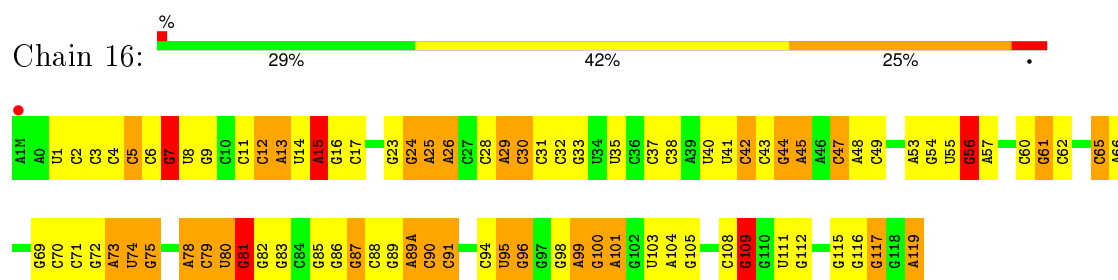




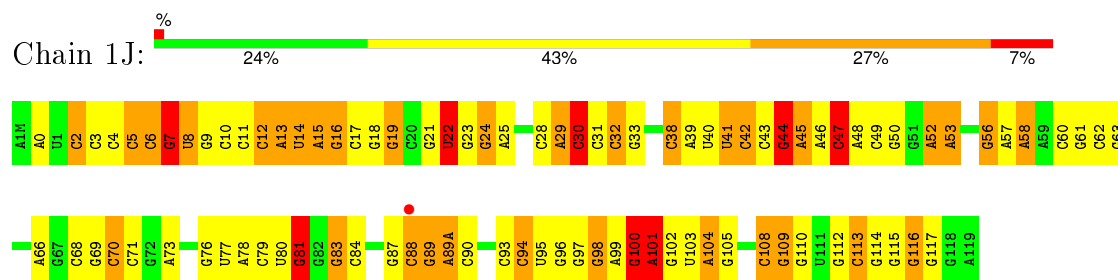




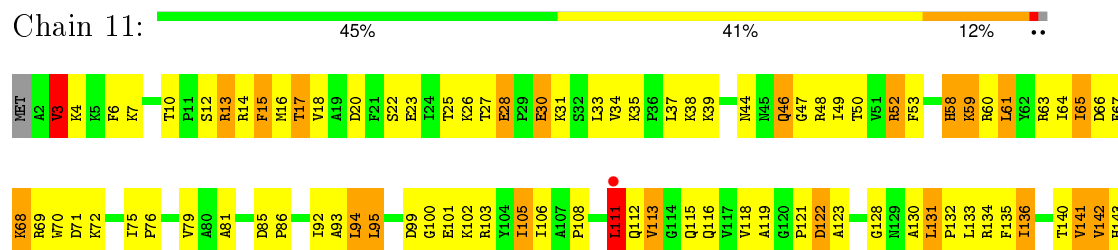
• Molecule 27: 5S ribosomal RNA

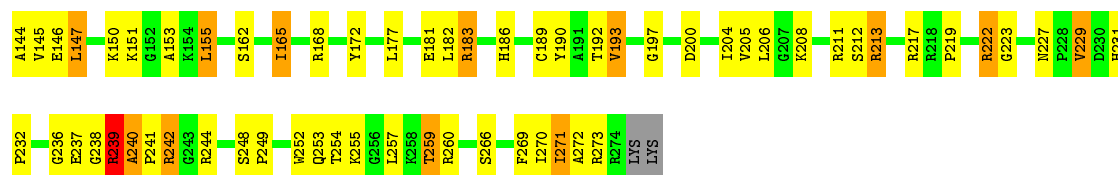


• Molecule 27: 5S ribosomal RNA

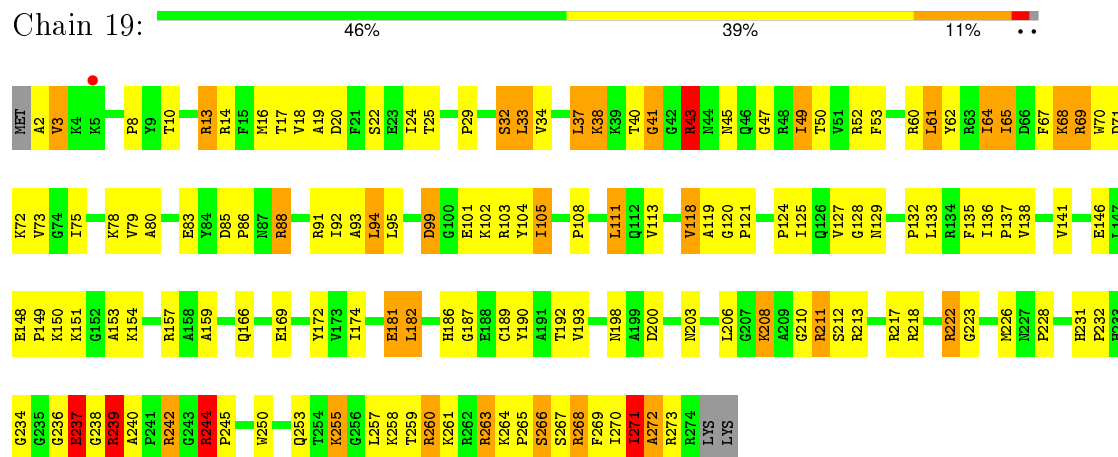


• Molecule 28: 50S ribosomal protein L2

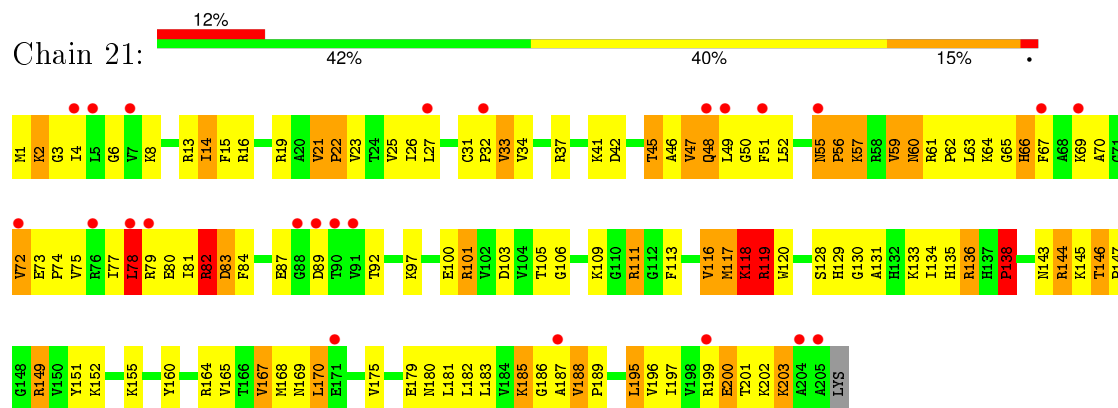




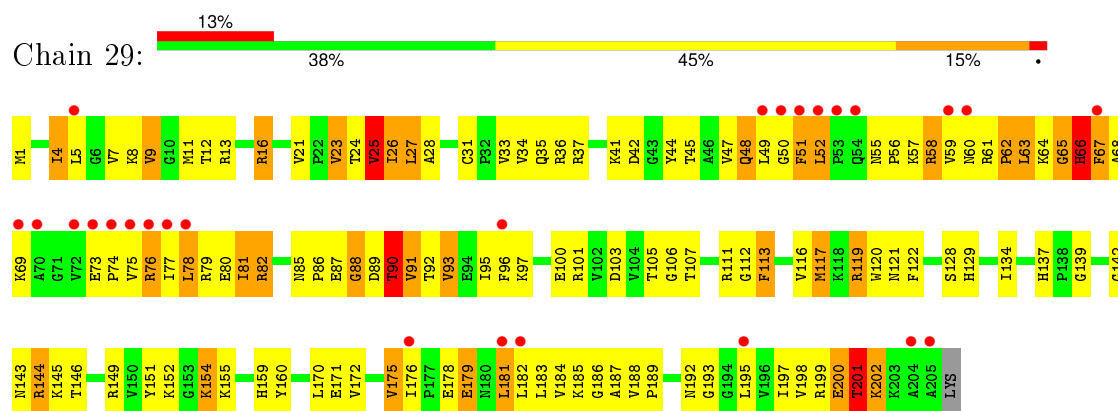
• Molecule 28: 50S ribosomal protein L2



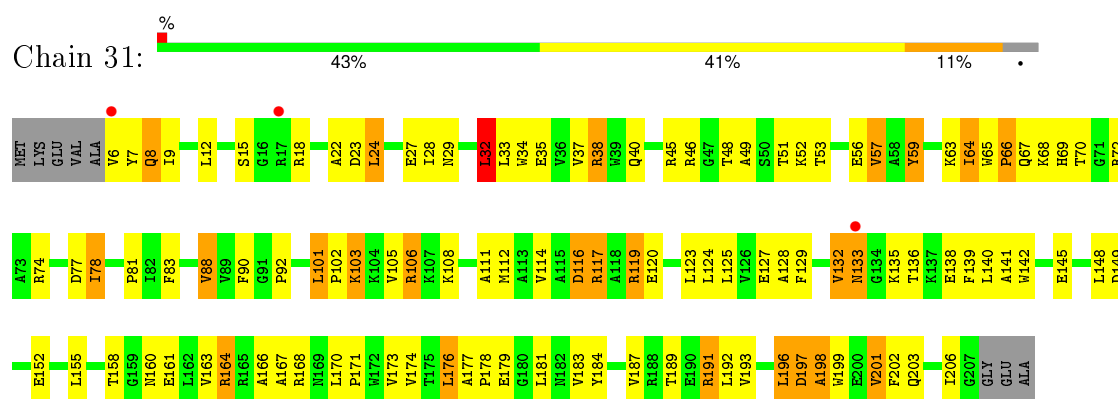
• Molecule 29: 50S ribosomal protein L3



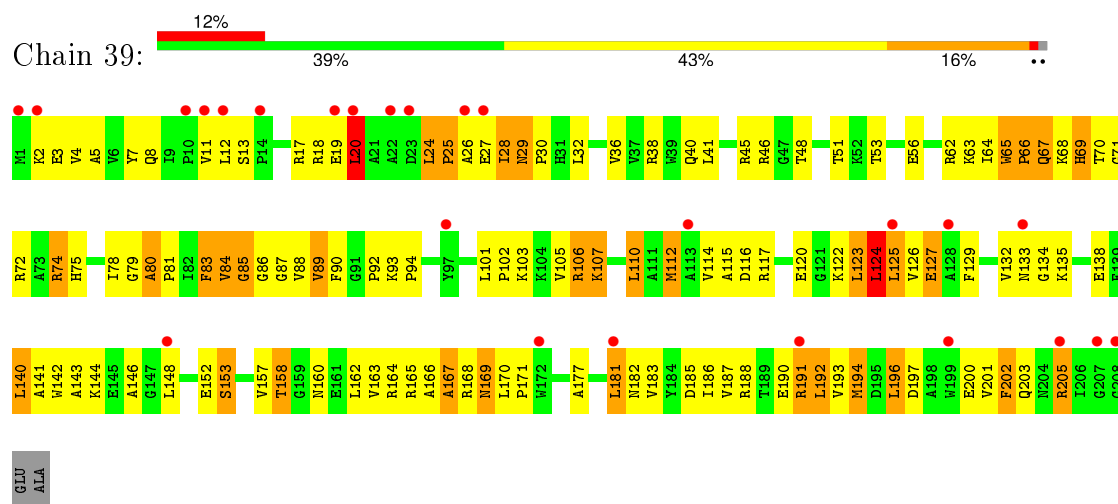
• Molecule 29: 50S ribosomal protein L3



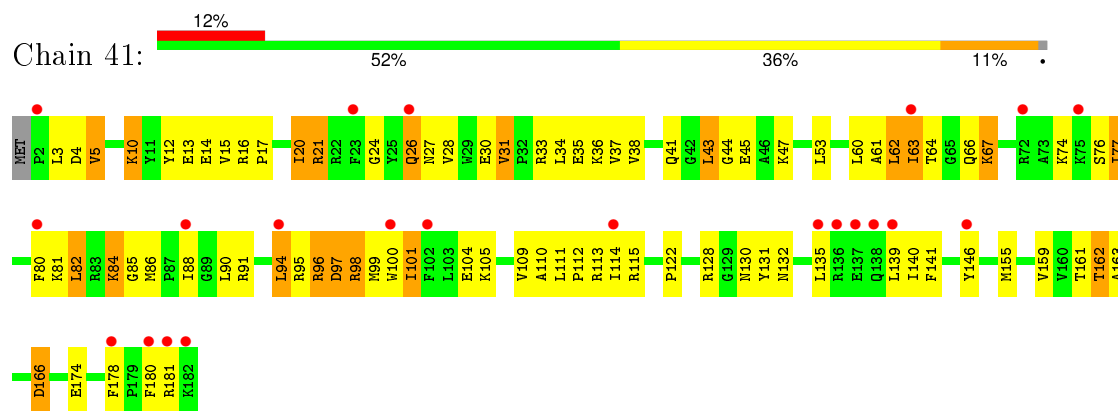
• Molecule 30: 50S ribosomal protein L4



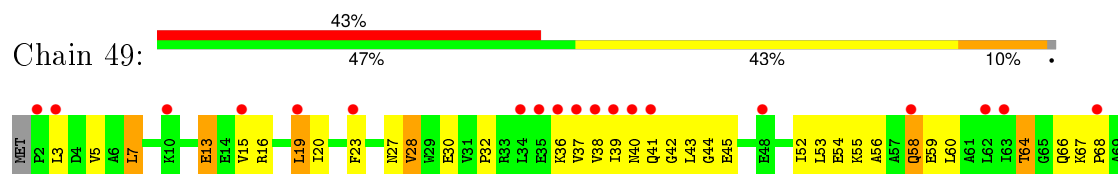
• Molecule 30: 50S ribosomal protein L4

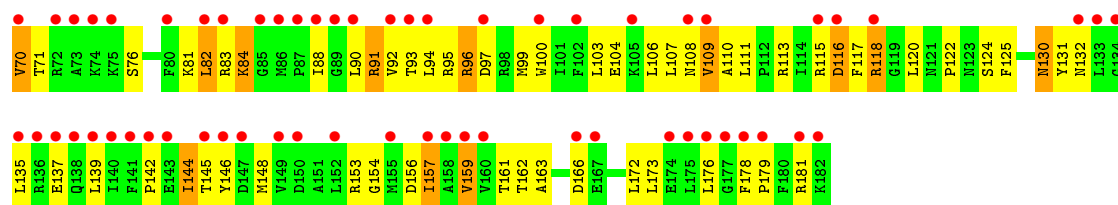


• Molecule 31: 50S ribosomal protein L5

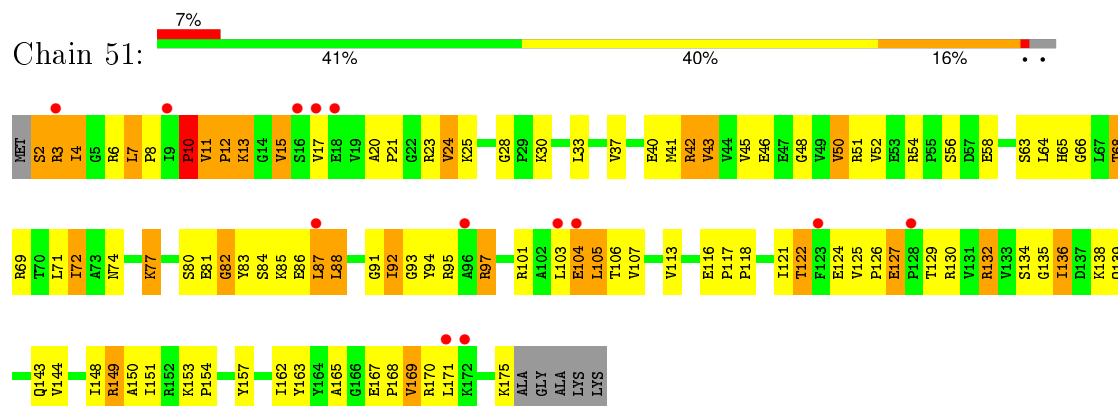


• Molecule 31: 50S ribosomal protein L5

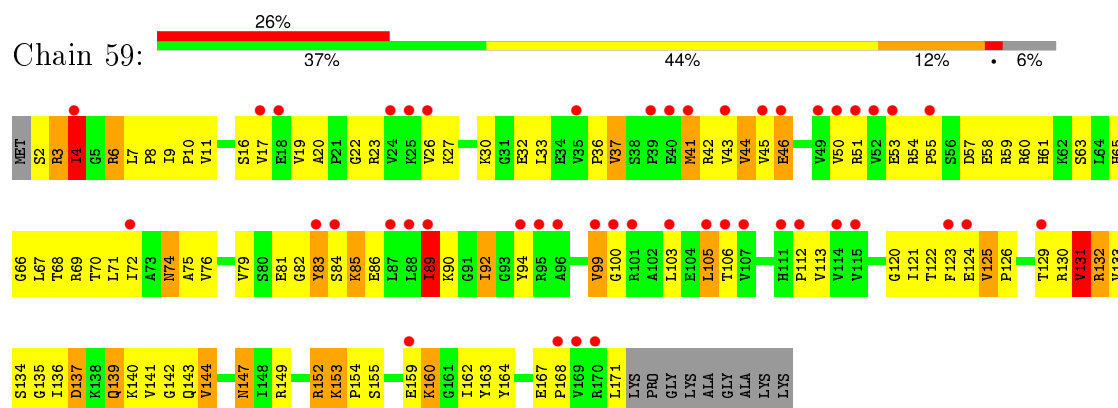




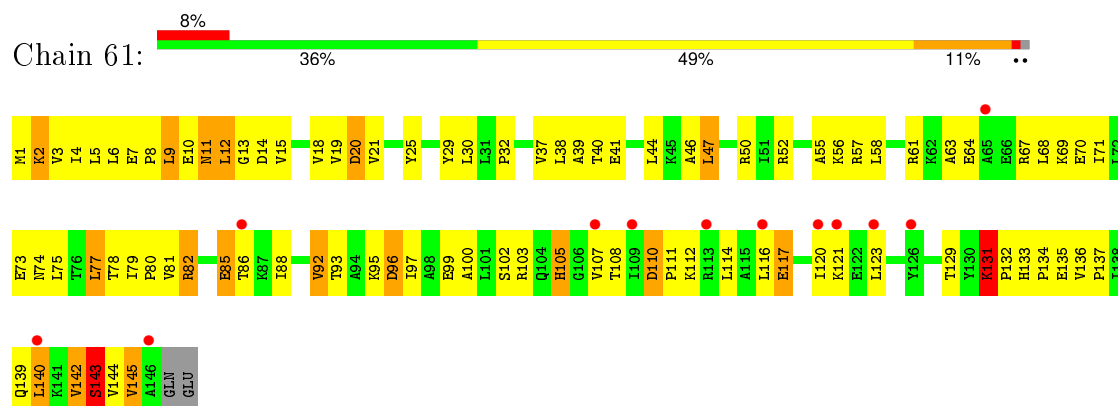
• Molecule 32: 50S ribosomal protein L6



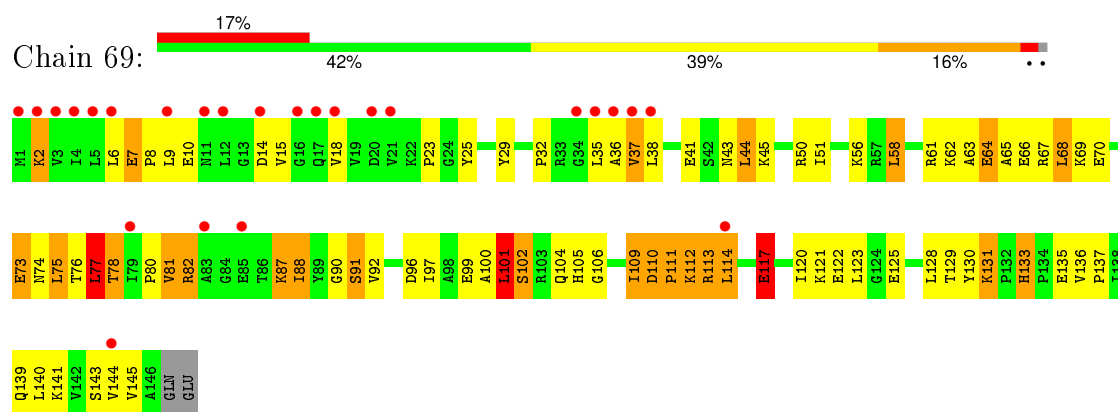
• Molecule 32: 50S ribosomal protein L6



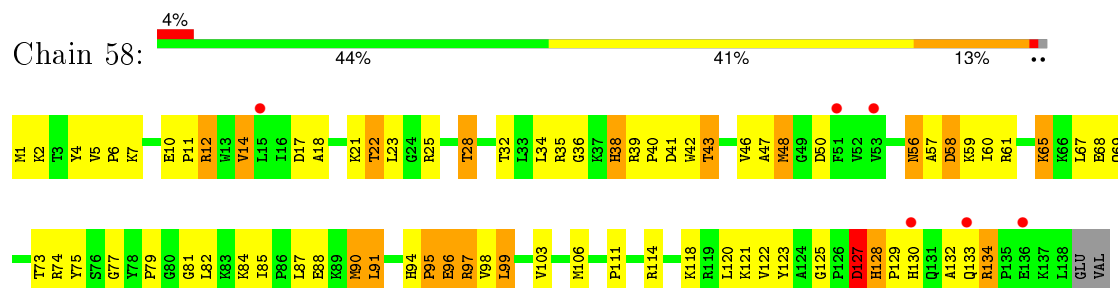
• Molecule 33: 50S ribosomal protein L9



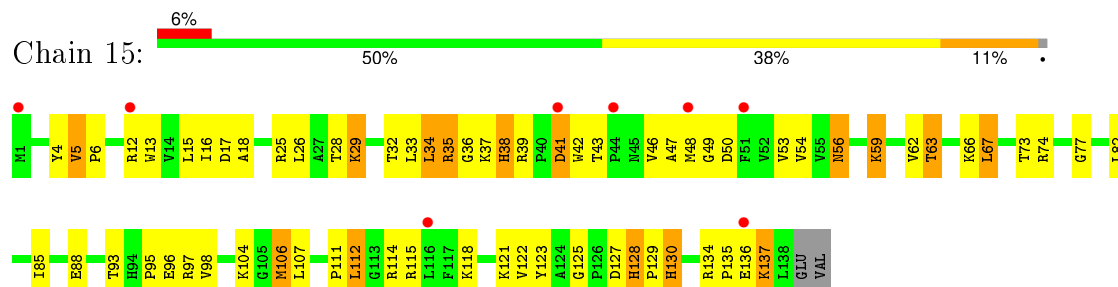
• Molecule 33: 50S ribosomal protein L9



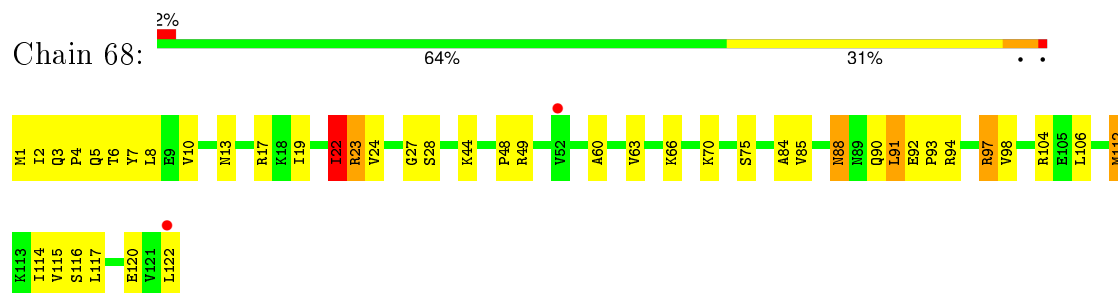
- Molecule 34: 50S ribosomal protein L13



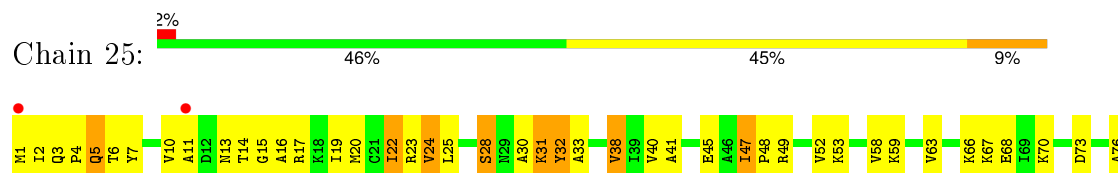
- Molecule 34: 50S ribosomal protein L13



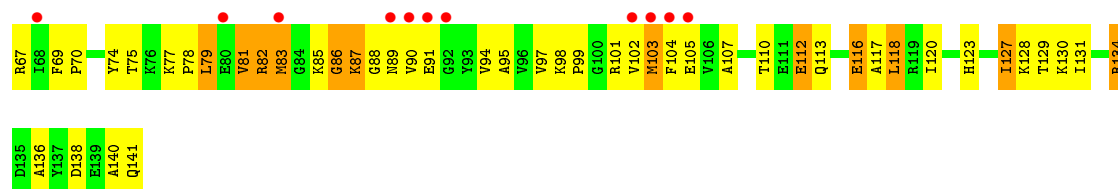
- Molecule 35: 50S ribosomal protein L14



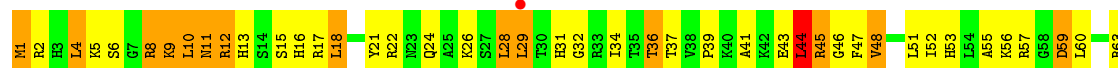
- Molecule 35: 50S ribosomal protein L14



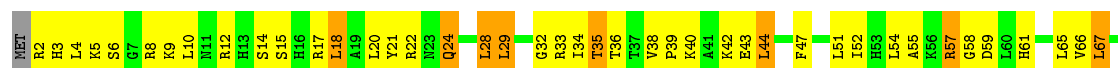




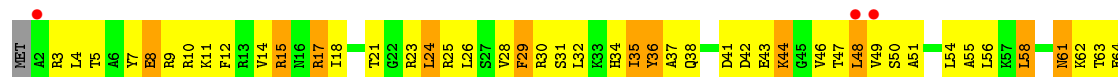
- Molecule 38: 50S ribosomal protein L17



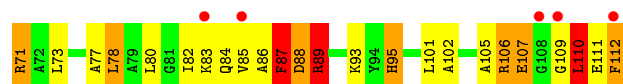
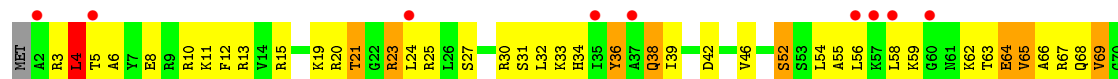
- Molecule 38: 50S ribosomal protein L17



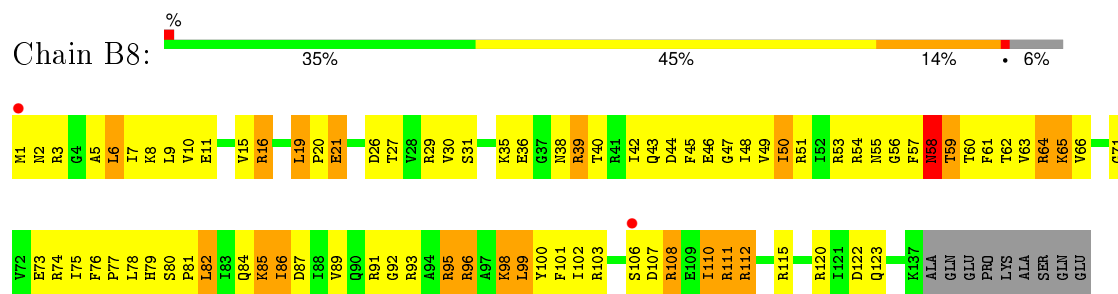
- Molecule 39: 50S ribosomal protein L18



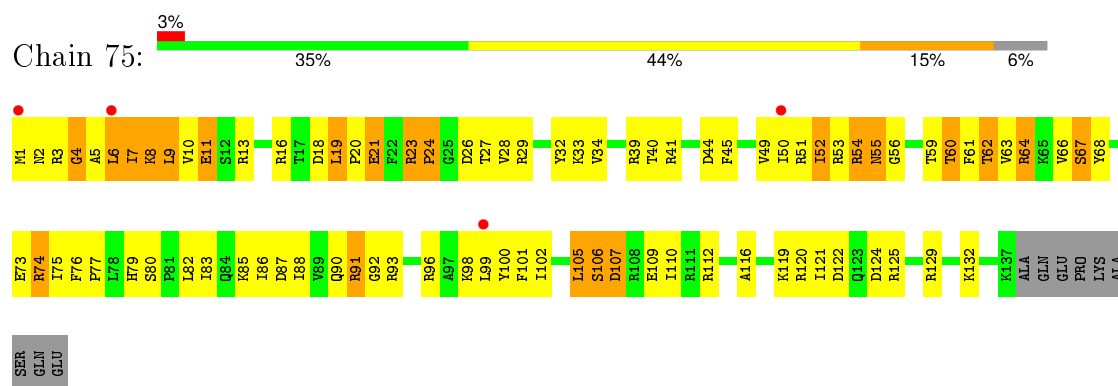
- Molecule 39: 50S ribosomal protein L18



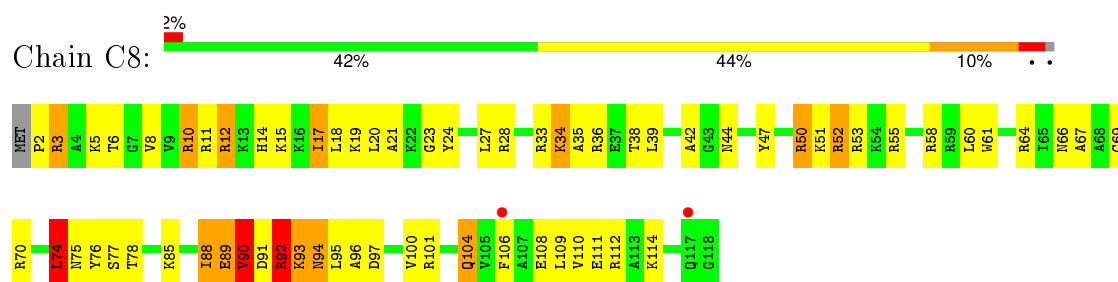
- Molecule 40: 50S ribosomal protein L19



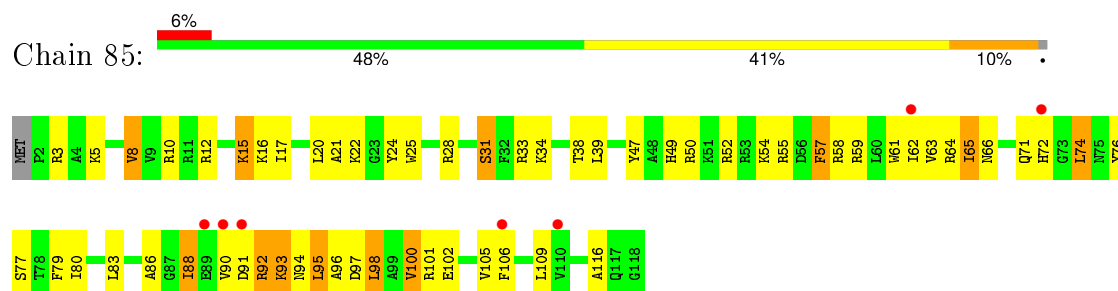
- Molecule 40: 50S ribosomal protein L19



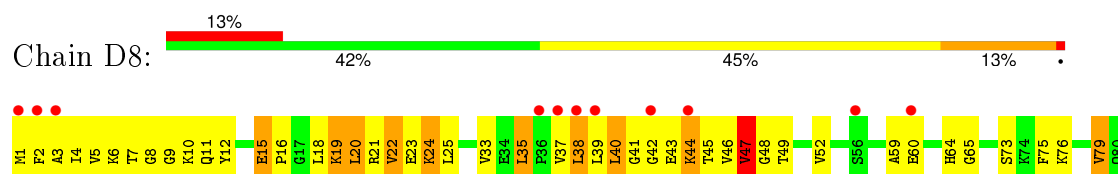
- Molecule 41: 50S ribosomal protein L20

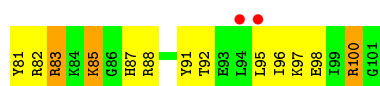


- Molecule 41: 50S ribosomal protein L20

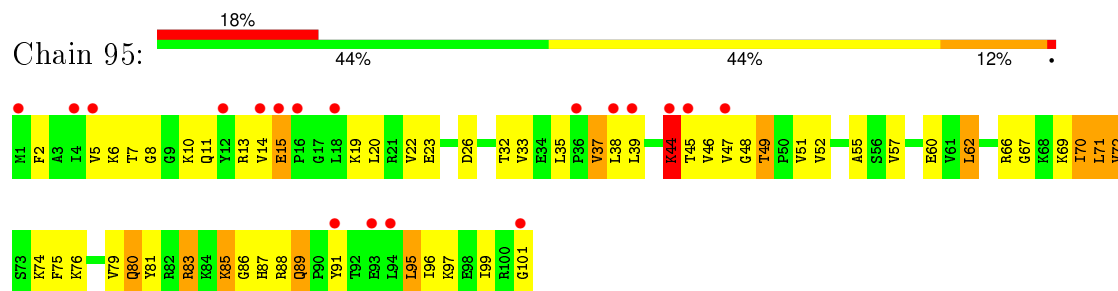


- Molecule 42: 50S ribosomal protein L21

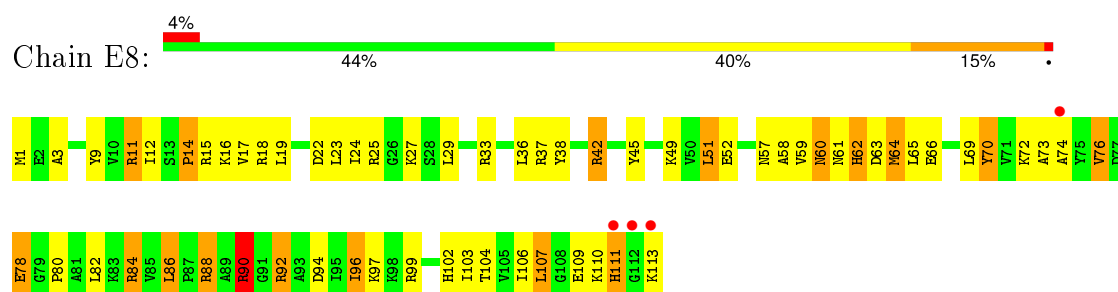




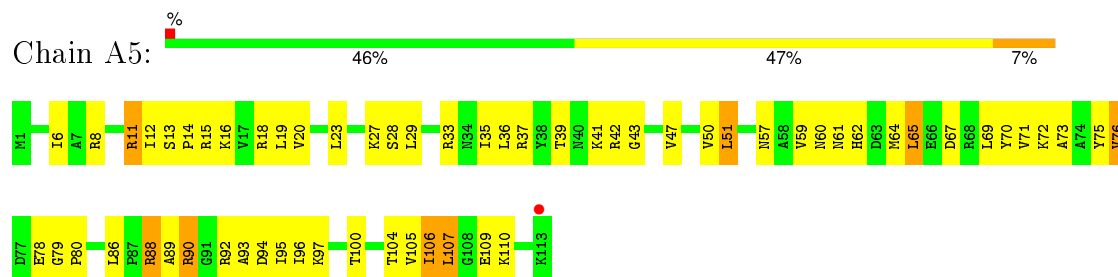
- Molecule 42: 50S ribosomal protein L21



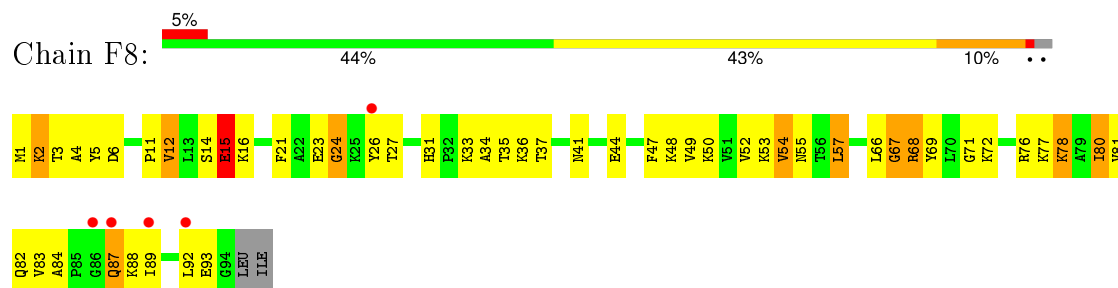
- Molecule 43: 50S ribosomal protein L22



- Molecule 43: 50S ribosomal protein L22

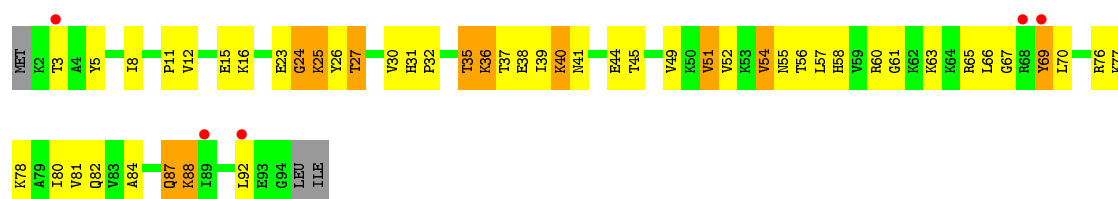


- Molecule 44: 50S ribosomal protein L23

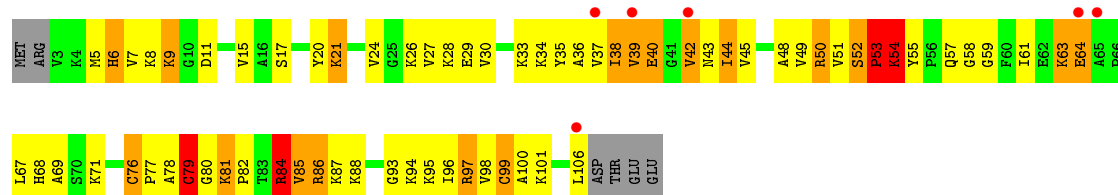


- Molecule 44: 50S ribosomal protein L23

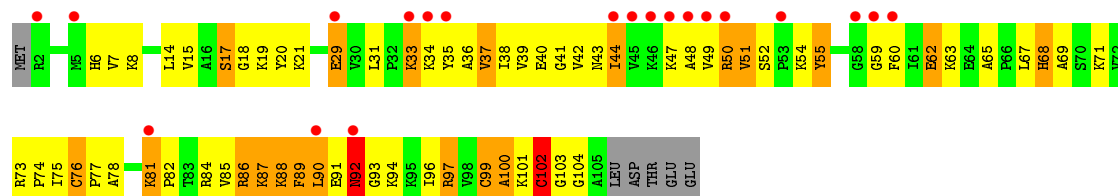




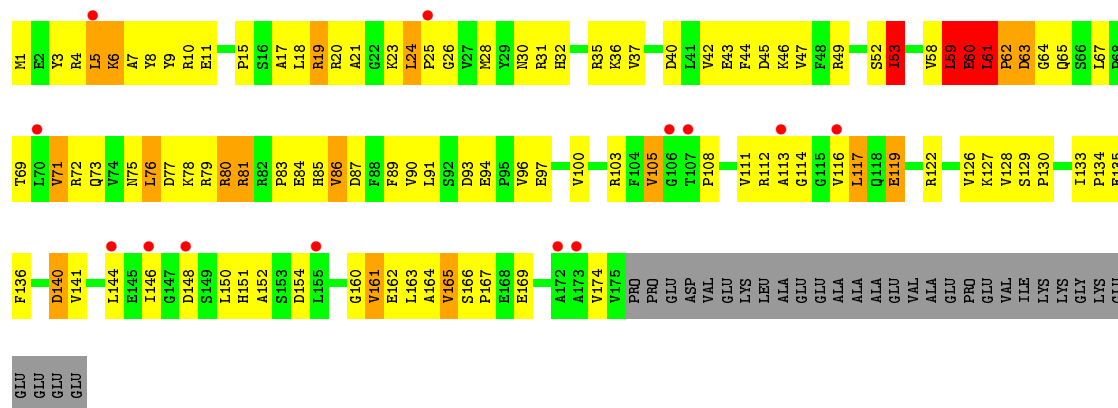
- Molecule 45: 50S ribosomal protein L24



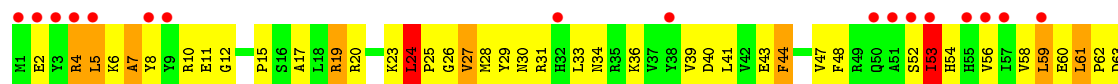
- Molecule 45: 50S ribosomal protein L24



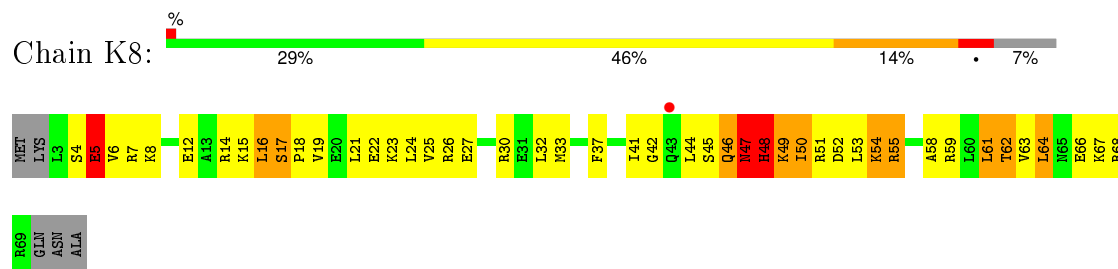
- Molecule 46: 50S ribosomal protein L25



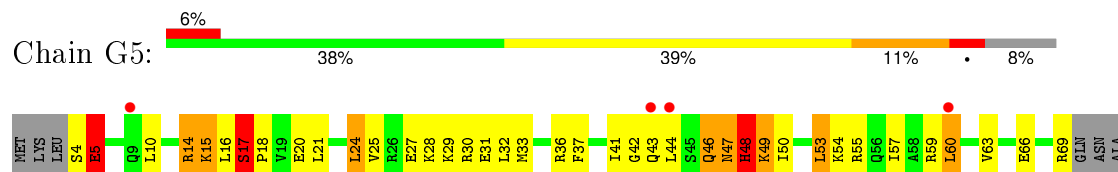
- Molecule 46: 50S ribosomal protein L25



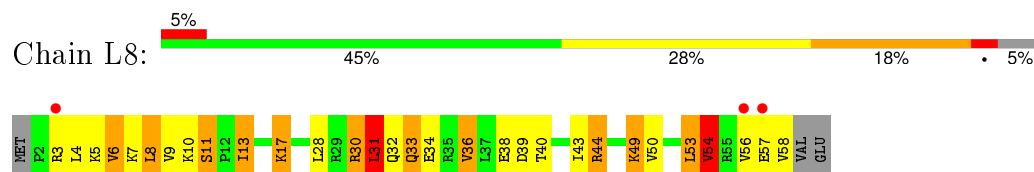
- Molecule 49: 50S ribosomal protein L29



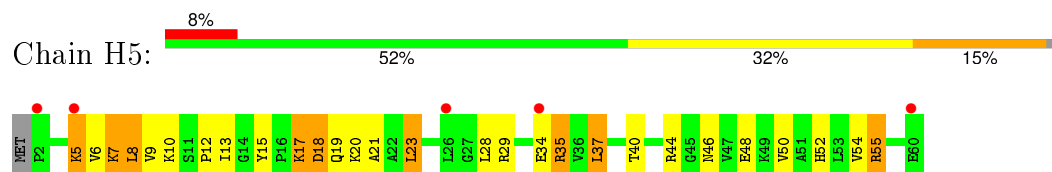
- Molecule 49: 50S ribosomal protein L29



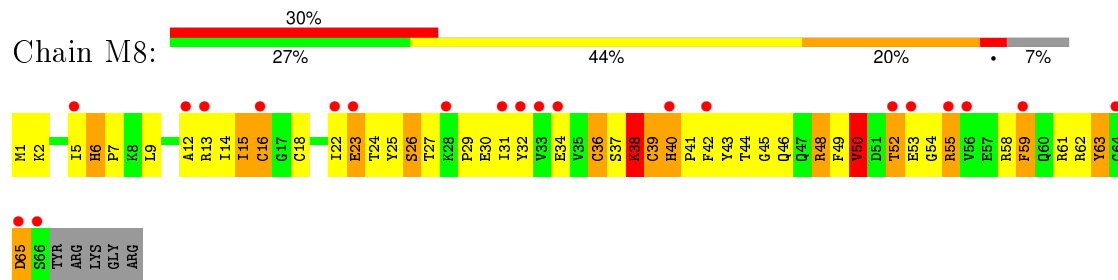
- Molecule 50: 50S ribosomal protein L30



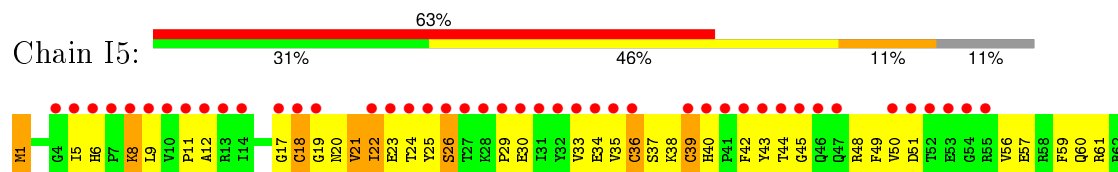
- Molecule 50: 50S ribosomal protein L30



- Molecule 51: 50S ribosomal protein L31

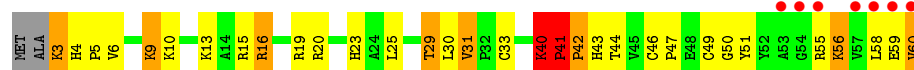
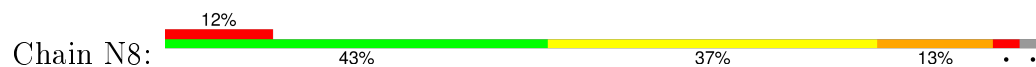


- Molecule 51: 50S ribosomal protein L31

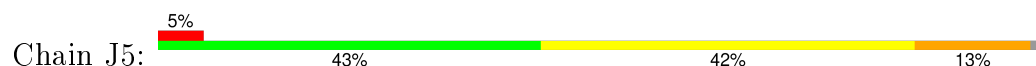




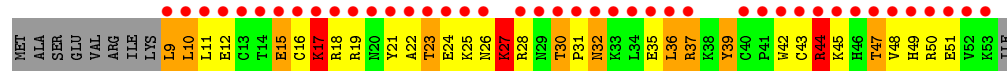
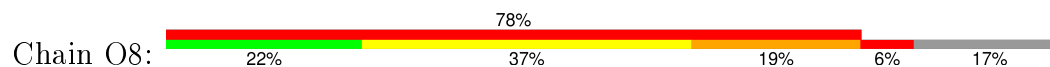
- Molecule 52: 50S ribosomal protein L32



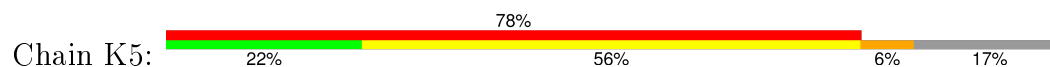
- Molecule 52: 50S ribosomal protein L32



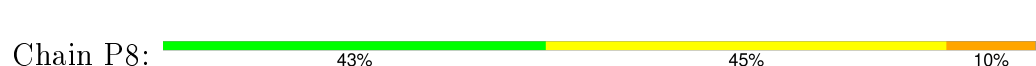
- Molecule 53: 50S ribosomal protein L33



- Molecule 53: 50S ribosomal protein L33



- Molecule 54: 50S ribosomal protein L34

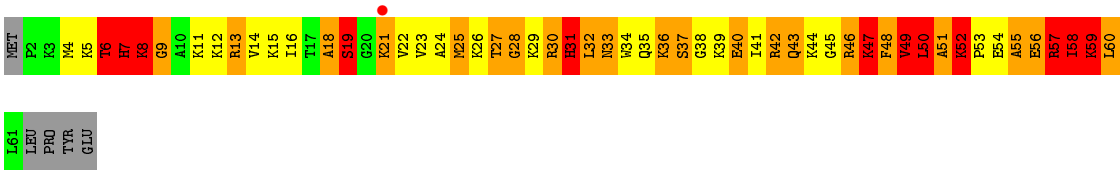


- Molecule 54: 50S ribosomal protein L34

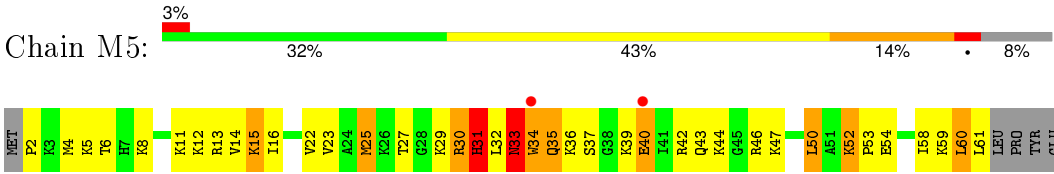


- Molecule 55: 50S ribosomal protein L35





• Molecule 55: 50S ribosomal protein L35



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	210.10Å 448.80Å 621.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	152.19 – 3.15 152.19 – 3.15	Depositor EDS
% Data completeness (in resolution range)	99.9 (152.19-3.15) 93.3 (152.19-3.15)	Depositor EDS
R_{merge}	0.37	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 3.13Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.189 , 0.231 0.192 , 0.231	Depositor DCC
R_{free} test set	2000 reflections (0.21%)	DCC
Wilson B-factor (Å ²)	87.2	Xtriage
Anisotropy	0.287	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 87.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 1000693 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	300537	wwPDB-VP
Average B, all atoms (Å ²)	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: OMC, 5MU, ZN, MIA, MG, 4SU, 7MG, PSU

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	13	0.82	13/36052 (0.0%)	1.55	642/56266 (1.1%)
1	1G	0.71	2/36025 (0.0%)	1.38	354/56227 (0.6%)
2	12	0.37	0/1959	0.64	2/2642 (0.1%)
2	1E	0.45	0/1959	0.73	3/2642 (0.1%)
3	22	0.38	0/1636	0.62	0/2205
3	2E	0.51	0/1629	0.70	0/2195
4	32	0.50	0/1733	0.73	1/2318 (0.0%)
4	3E	0.65	2/1733 (0.1%)	0.78	2/2318 (0.1%)
5	42	0.48	0/1171	0.72	1/1576 (0.1%)
5	4E	0.55	0/1171	0.77	1/1576 (0.1%)
6	52	0.55	0/856	0.70	0/1154
6	5E	0.58	0/856	0.76	0/1154
7	62	0.44	0/1276	0.59	0/1709
7	6E	0.47	0/1276	0.61	0/1709
8	72	0.42	0/1136	0.67	1/1527 (0.1%)
8	7E	0.57	0/1136	0.78	2/1527 (0.1%)
9	82	0.40	0/1029	0.66	1/1379 (0.1%)
9	8E	0.47	0/1029	0.72	0/1379
10	1A	0.39	0/814	0.62	0/1095
10	1I	0.47	0/814	0.67	0/1095
11	2A	0.49	0/900	0.75	0/1213
11	2I	0.57	0/900	0.79	1/1213 (0.1%)
12	3A	0.61	0/991	0.81	1/1327 (0.1%)
12	3I	0.74	0/991	0.94	0/1327
13	4A	0.36	0/943	0.64	0/1265
13	4I	0.51	0/948	0.76	0/1272
14	5A	0.43	0/485	0.73	0/643
14	5I	0.66	1/501 (0.2%)	0.88	3/664 (0.5%)
15	6A	0.49	0/745	0.60	0/992
15	6I	0.60	0/745	0.82	0/992
16	7A	0.51	0/721	0.69	0/970
16	7I	0.52	0/721	0.73	0/970

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
17	8A	0.54	0/847	0.65	0/1131
17	8I	0.62	0/847	0.78	0/1131
18	9A	0.53	0/596	0.72	1/790 (0.1%)
18	9I	0.60	1/596 (0.2%)	0.79	0/790
19	AA	0.40	0/638	0.66	1/860 (0.1%)
19	AI	0.58	0/661	0.84	1/890 (0.1%)
20	BA	0.50	0/765	0.77	0/1007
20	BI	0.42	0/765	0.70	1/1007 (0.1%)
21	1B	0.55	0/221	0.64	0/288
21	1F	0.50	0/221	0.73	0/288
22	1K	0.53	0/1739	1.15	15/2708 (0.6%)
22	1L	0.44	0/1739	0.96	4/2708 (0.1%)
23	2K	1.13	6/1721 (0.3%)	1.67	45/2682 (1.7%)
23	2L	0.99	7/1721 (0.4%)	1.43	18/2682 (0.7%)
24	3K	0.52	0/1809	1.29	25/2819 (0.9%)
24	3L	0.48	2/1809 (0.1%)	1.14	15/2819 (0.5%)
25	4K	1.00	0/338	1.43	5/524 (1.0%)
25	4L	0.79	0/147	1.61	5/227 (2.2%)
26	14	0.99	121/70167 (0.2%)	1.77	2211/109541 (2.0%)
26	1H	1.20	257/70233 (0.4%)	2.06	3865/109643 (3.5%)
27	16	1.01	3/2928 (0.1%)	1.83	103/4568 (2.3%)
27	1J	0.83	1/2928 (0.0%)	1.53	48/4568 (1.1%)
28	11	0.96	3/2170 (0.1%)	1.14	13/2926 (0.4%)
28	19	0.80	0/2170	1.01	6/2926 (0.2%)
29	21	0.75	0/1601	0.99	3/2160 (0.1%)
29	29	0.76	0/1601	0.99	2/2160 (0.1%)
30	31	0.84	0/1620	1.02	5/2194 (0.2%)
30	39	0.67	0/1662	0.95	1/2249 (0.0%)
31	41	0.60	0/1498	0.86	2/2016 (0.1%)
31	49	0.44	0/1498	0.66	0/2016
32	51	0.64	0/1362	0.93	2/1841 (0.1%)
32	59	0.38	0/1332	0.69	1/1802 (0.1%)
33	61	0.55	0/1151	0.83	1/1558 (0.1%)
33	69	0.49	0/1151	0.79	4/1558 (0.3%)
34	15	0.53	0/1131	0.77	0/1525
34	58	0.65	0/1131	0.88	0/1525
35	25	0.66	0/943	0.83	0/1269
35	68	0.75	0/943	0.87	1/1269 (0.1%)
36	35	0.73	0/1162	1.15	6/1544 (0.4%)
36	78	0.81	0/1162	1.12	3/1544 (0.2%)
37	45	0.72	0/1143	0.96	1/1527 (0.1%)
37	88	0.96	3/1107 (0.3%)	1.15	7/1478 (0.5%)
38	55	0.71	0/974	0.93	1/1302 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
38	98	0.67	0/982	1.00	4/1312 (0.3%)
39	65	0.57	0/892	0.96	3/1187 (0.3%)
39	A8	0.78	0/892	1.05	4/1187 (0.3%)
40	75	0.65	0/1155	0.79	0/1542
40	B8	0.70	0/1155	0.89	0/1542
41	85	0.64	0/982	0.81	0/1306
41	C8	0.80	0/982	0.97	5/1306 (0.4%)
42	95	0.67	0/790	0.91	2/1057 (0.2%)
42	D8	0.68	0/790	0.90	1/1057 (0.1%)
43	A5	0.74	0/911	0.88	1/1220 (0.1%)
43	E8	0.72	0/911	1.07	7/1220 (0.6%)
44	B5	0.83	0/744	0.89	0/1000
44	F8	0.96	2/756 (0.3%)	1.03	1/1014 (0.1%)
45	C5	0.78	0/807	0.97	1/1076 (0.1%)
45	G8	0.78	0/804	1.07	2/1073 (0.2%)
46	D5	0.45	0/1460	0.71	1/1982 (0.1%)
46	H8	0.56	0/1427	0.87	2/1935 (0.1%)
47	E5	0.65	0/621	0.94	1/827 (0.1%)
47	I8	0.87	0/635	1.03	0/847
48	F5	0.68	0/770	1.00	4/1022 (0.4%)
48	J8	0.81	0/770	1.00	2/1022 (0.2%)
49	G5	0.65	1/560 (0.2%)	0.84	1/741 (0.1%)
49	K8	0.88	0/565	1.08	1/748 (0.1%)
50	H5	0.54	0/474	0.78	0/635
50	L8	0.84	0/457	1.08	4/613 (0.7%)
51	I5	0.46	0/527	0.76	0/709
51	M8	0.64	0/545	0.84	0/733
52	J5	0.70	0/473	0.91	1/639 (0.2%)
52	N8	0.77	0/468	1.02	2/632 (0.3%)
53	K5	0.61	0/396	0.96	0/529
53	O8	0.83	1/396 (0.3%)	0.89	1/529 (0.2%)
54	L5	0.89	0/406	1.04	2/536 (0.4%)
54	P8	1.04	0/426	1.17	2/561 (0.4%)
55	M5	0.99	1/483 (0.2%)	1.14	1/634 (0.2%)
55	Q8	1.47	5/486 (1.0%)	1.86	11/638 (1.7%)
All	All	0.90	432/322727 (0.1%)	1.57	7504/483212 (1.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	12	0	1
2	1E	0	1
4	32	0	4
4	3E	0	1
9	82	0	2
10	1A	0	2
11	2A	0	2
12	3A	0	1
12	3I	0	1
15	6I	0	1
19	AI	0	2
20	BA	0	3
28	11	0	1
28	19	0	4
29	21	0	3
29	29	0	4
30	31	0	1
30	39	0	6
31	41	0	2
31	49	0	1
32	59	0	3
33	61	0	5
33	69	0	2
34	15	0	2
34	58	0	2
35	68	0	1
36	35	0	4
36	78	0	2
37	45	0	6
37	88	0	2
38	55	0	1
38	98	0	3
39	65	0	2
39	A8	0	2
40	75	0	1
40	B8	0	1
41	85	0	2
41	C8	0	2
42	95	0	1
43	A5	0	3
44	B5	0	2
44	F8	0	1
45	C5	0	3

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Mol	Chain	#Chirality outliers	#Planarity outliers
45	G8	0	3
46	D5	0	2
46	H8	0	2
47	I8	0	2
48	F5	0	1
48	J8	0	2
49	G5	0	3
49	K8	0	3
51	I5	0	2
51	M8	0	2
52	N8	0	2
53	K5	0	3
53	O8	0	3
55	M5	0	3
55	Q8	0	13
All	All	0	142

All (432) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
23	2L	21	U	C5-C6	18.43	1.50	1.34
23	2K	21	U	C5-C6	16.94	1.49	1.34
26	1H	774	A	N9-C4	-14.63	1.29	1.37
26	14	783	A	N9-C4	-13.79	1.29	1.37
26	1H	783	A	N9-C4	-12.87	1.30	1.37
26	1H	2430	A	N9-C4	-12.58	1.30	1.37
26	1H	1614	A	N9-C4	-12.54	1.30	1.37
26	14	774	A	N9-C4	-12.19	1.30	1.37
26	1H	783	A	N3-C4	-11.76	1.27	1.34
26	1H	1698	A	N9-C4	-11.70	1.30	1.37
23	2L	21	U	C2-N3	11.68	1.46	1.37
23	2K	21	U	C2-N3	11.65	1.46	1.37
26	1H	74	A	N9-C4	-10.88	1.31	1.37
26	14	783	A	N3-C4	-10.72	1.28	1.34
26	1H	676	A	N9-C4	-10.41	1.31	1.37
1	13	792	A	C5-C6	-10.34	1.31	1.41
26	1H	2287	A	N9-C4	-10.29	1.31	1.37
26	1H	676	A	N9-C8	10.21	1.46	1.37
26	14	1332	G	N9-C4	-10.07	1.29	1.38
26	1H	1899	G	N9-C4	-10.01	1.29	1.38
26	1H	71	A	N9-C4	-9.88	1.31	1.37
26	14	74	A	N9-C4	-9.82	1.31	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	14	1786	A	N9-C4	-9.74	1.32	1.37
23	2K	21	U	N1-C2	9.56	1.47	1.38
1	13	792	A	N9-C4	-9.44	1.32	1.37
26	14	528	A	N9-C4	-9.43	1.32	1.37
26	1H	1698	A	N3-C4	-9.26	1.29	1.34
26	1H	2442	C	N1-C6	-9.18	1.31	1.37
23	2L	21	U	N1-C2	9.15	1.46	1.38
4	3E	12	CYS	CB-SG	9.12	1.97	1.82
26	1H	197	A	N3-C4	-9.12	1.29	1.34
26	1H	860	U	N1-C2	8.91	1.46	1.38
26	1H	805	G	N7-C5	-8.85	1.33	1.39
23	2L	21	U	C4-C5	8.84	1.51	1.43
23	2K	21	U	C4-C5	8.80	1.51	1.43
26	1H	1332	G	N9-C4	-8.78	1.30	1.38
26	1H	945	A	N7-C5	-8.75	1.34	1.39
26	1H	621	A	N9-C4	-8.40	1.32	1.37
26	1H	138	G	N9-C8	8.39	1.43	1.37
26	1H	676	A	C5-C4	8.35	1.44	1.38
26	1H	1899	G	C2-N3	-8.28	1.26	1.32
28	11	237	GLU	CG-CD	8.21	1.64	1.51
26	1H	2576	G	C8-N7	-8.20	1.26	1.30
26	1H	1786	A	N3-C4	-8.10	1.29	1.34
26	14	1698	A	N9-C4	-8.05	1.33	1.37
26	1H	2062	A	N3-C4	8.00	1.39	1.34
26	1H	1786	A	C5-C6	-7.96	1.33	1.41
26	1H	2392	A	N9-C4	-7.95	1.33	1.37
26	1H	1899	G	N9-C8	7.95	1.43	1.37
26	14	2287	A	N9-C4	-7.94	1.33	1.37
26	14	1698	A	N7-C5	-7.91	1.34	1.39
26	1H	1899	G	N3-C4	-7.91	1.29	1.35
26	1H	2346	A	N3-C4	-7.90	1.30	1.34
26	14	751	A	N9-C4	-7.86	1.33	1.37
28	11	122	ASP	CB-CG	7.85	1.68	1.51
26	1H	774	A	C5-C6	-7.74	1.34	1.41
26	14	2506	U	C2-N3	7.68	1.43	1.37
26	1H	1786	A	N9-C4	-7.64	1.33	1.37
26	1H	2062	A	N7-C5	7.59	1.43	1.39
26	1H	783	A	C5-C6	-7.49	1.34	1.41
1	13	1227	A	N9-C4	-7.49	1.33	1.37
26	1H	2490	G	N9-C8	7.49	1.43	1.37
26	1H	74	A	N3-C4	-7.48	1.30	1.34
1	13	810	C	N1-C6	-7.45	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	14	1678	G	N9-C4	-7.44	1.31	1.38
26	1H	1969	A	C6-N1	-7.39	1.30	1.35
26	1H	829	A	N7-C5	-7.38	1.34	1.39
1	13	815	A	N9-C4	-7.37	1.33	1.37
4	3E	9	CYS	CB-SG	7.37	1.94	1.82
26	1H	783	A	N7-C5	-7.36	1.34	1.39
26	1H	2451	A	N9-C4	-7.34	1.33	1.37
26	14	1698	A	C5-C6	-7.31	1.34	1.41
26	1H	735	A	C5-C4	-7.29	1.33	1.38
26	1H	821	A	N7-C5	-7.27	1.34	1.39
26	1H	2713	A	C5-C4	7.25	1.43	1.38
23	2K	21	U	N3-C4	7.24	1.45	1.38
26	1H	2713	A	N9-C4	-7.22	1.33	1.37
26	1H	777	A	N9-C4	-7.16	1.33	1.37
26	14	1313	U	N1-C2	-7.16	1.32	1.38
26	1H	2346	A	N7-C5	-7.15	1.34	1.39
26	1H	71	A	C5-C6	-7.13	1.34	1.41
26	1H	676	A	N3-C4	-7.08	1.30	1.34
14	5I	27	CYS	CB-SG	-7.06	1.70	1.82
26	1H	1678	G	N9-C8	7.05	1.42	1.37
26	1H	1021	A	N9-C4	-7.04	1.33	1.37
26	1H	245	G	N7-C5	-6.97	1.35	1.39
26	1H	2490	G	N9-C4	-6.94	1.32	1.38
26	1H	735	A	N3-C4	-6.94	1.30	1.34
26	1H	71	A	C6-N6	-6.93	1.28	1.33
26	1H	945	A	C5-C6	-6.83	1.34	1.41
26	1H	698	C	N1-C6	-6.82	1.33	1.37
26	14	783	A	N7-C5	-6.82	1.35	1.39
26	1H	2451	A	C6-N1	-6.79	1.30	1.35
55	Q8	49	VAL	CA-CB	6.78	1.69	1.54
26	14	2506	U	N1-C2	6.71	1.44	1.38
26	1H	1355	G	C6-N1	-6.66	1.34	1.39
23	2K	21	U	N1-C6	6.66	1.44	1.38
26	1H	939	G	N3-C4	-6.65	1.30	1.35
55	Q8	6	THR	CA-CB	6.63	1.70	1.53
26	1H	1966	A	N9-C4	-6.63	1.33	1.37
26	1H	2247	A	N3-C4	-6.62	1.30	1.34
26	1H	845	G	C2-N3	6.61	1.38	1.32
26	14	1332	G	N9-C8	6.58	1.42	1.37
26	14	1289	C	N1-C6	-6.55	1.33	1.37
26	14	1612	C	N1-C6	-6.54	1.33	1.37
26	14	1332	G	N3-C4	-6.54	1.30	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	1142(A)	A	N9-C4	-6.52	1.33	1.37
26	14	780	G	N7-C5	-6.50	1.35	1.39
26	14	1617	C	N1-C6	-6.50	1.33	1.37
26	1H	1564	C	N3-C4	-6.47	1.29	1.33
26	1H	528	A	N9-C4	-6.47	1.33	1.37
26	1H	777	A	N3-C4	-6.47	1.30	1.34
26	14	71	A	N9-C4	-6.46	1.33	1.37
26	1H	621	A	N9-C8	6.45	1.43	1.37
26	14	1142(A)	A	N3-C4	-6.39	1.31	1.34
26	1H	2589	A	C5-C4	-6.37	1.34	1.38
26	1H	945	A	N3-C4	-6.34	1.31	1.34
26	1H	70	G	C6-N1	-6.33	1.35	1.39
44	F8	15	GLU	CG-CD	6.33	1.61	1.51
26	1H	1950	G	N9-C8	6.32	1.42	1.37
26	1H	138	G	C5-C4	6.32	1.42	1.38
26	1H	197	A	N9-C4	-6.30	1.34	1.37
26	14	123	G	C6-N1	-6.29	1.35	1.39
26	1H	679	C	N1-C6	-6.28	1.33	1.37
26	1H	1616	A	C5-C6	-6.27	1.35	1.41
26	1H	2594	C	N1-C6	-6.27	1.33	1.37
26	1H	732	C	N1-C6	-6.26	1.33	1.37
26	1H	774	A	N7-C5	-6.26	1.35	1.39
26	14	2360	A	N9-C4	-6.26	1.34	1.37
26	14	674	G	N7-C5	6.25	1.43	1.39
26	1H	2053	G	C5-C4	-6.24	1.33	1.38
26	1H	140	A	C5-C6	-6.23	1.35	1.41
44	F8	15	GLU	CB-CG	6.22	1.64	1.52
26	1H	805	G	C6-N1	-6.20	1.35	1.39
26	14	1902	C	C4-N4	-6.19	1.28	1.33
26	14	2346	A	N3-C4	-6.16	1.31	1.34
26	1H	1332	G	N3-C4	-6.16	1.31	1.35
26	1H	746	A	N3-C4	-6.16	1.31	1.34
26	14	211	A	N9-C4	-6.15	1.34	1.37
26	1H	2373	G	C2-N3	6.15	1.37	1.32
1	13	1483	A	N9-C4	-6.14	1.34	1.37
26	14	2430	A	N9-C4	-6.14	1.34	1.37
26	1H	1815	A	C5-C4	-6.13	1.34	1.38
37	88	82	ARG	N-CA	6.13	1.58	1.46
26	14	1786	A	C5-C4	6.13	1.43	1.38
26	1H	2249	U	C2-N3	-6.12	1.33	1.37
23	2L	21	U	N3-C4	6.12	1.44	1.38
26	1H	663	G	N7-C5	-6.12	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	2031	A	N9-C4	6.12	1.41	1.37
23	2L	21	U	N1-C6	6.11	1.43	1.38
26	1H	949	C	N1-C6	-6.11	1.33	1.37
26	1H	123	G	C6-O6	-6.10	1.18	1.24
26	1H	1825	A	C6-N1	-6.10	1.31	1.35
26	14	693	C	N3-C4	-6.10	1.29	1.33
26	1H	684	G	C6-N1	-6.09	1.35	1.39
26	1H	2256	G	N1-C2	-6.08	1.32	1.37
26	1H	2518	A	N9-C4	-6.08	1.34	1.37
26	1H	1786	A	N7-C5	-6.07	1.35	1.39
26	1H	2393	A	N7-C5	-6.07	1.35	1.39
26	1H	792	G	N7-C5	-6.07	1.35	1.39
26	1H	1275	A	N7-C5	-6.07	1.35	1.39
26	14	945	A	C5-C6	-6.07	1.35	1.41
26	14	774	A	N3-C4	-6.07	1.31	1.34
26	14	783	A	C5-C6	-6.06	1.35	1.41
55	Q8	49	VAL	CB-CG2	6.05	1.65	1.52
26	1H	1313	U	C4-C5	-6.05	1.38	1.43
26	14	1142(A)	A	N9-C4	-6.04	1.34	1.37
26	1H	960	A	N9-C4	-6.04	1.34	1.37
26	14	782	A	C6-N1	-6.04	1.31	1.35
55	Q8	40	GLU	CB-CG	6.03	1.63	1.52
26	14	1786	A	N3-C4	-6.02	1.31	1.34
26	1H	2310	A	N9-C4	6.01	1.41	1.37
26	1H	2430	A	N1-C2	6.00	1.39	1.34
26	1H	946	G	N9-C4	-6.00	1.33	1.38
26	14	1786	A	C5-C6	-6.00	1.35	1.41
26	14	1950	G	C2-N3	5.97	1.37	1.32
27	1J	89(A)	A	N9-C4	5.97	1.41	1.37
26	1H	49	A	N7-C5	-5.97	1.35	1.39
26	1H	1202	C	N1-C6	-5.97	1.33	1.37
26	1H	2600	A	N7-C5	-5.96	1.35	1.39
27	16	99	A	C5-C4	5.96	1.43	1.38
26	1H	1616	A	N9-C4	-5.94	1.34	1.37
26	1H	71	A	N9-C8	5.94	1.42	1.37
26	1H	2071	A	C6-N6	-5.93	1.29	1.33
26	1H	2062	A	C5-C4	5.92	1.42	1.38
26	14	1332	G	C8-N7	5.92	1.34	1.30
26	14	1904	G	N9-C8	-5.92	1.33	1.37
26	1H	1332	G	N9-C8	5.91	1.42	1.37
26	14	1616	A	C5-C6	-5.91	1.35	1.41
26	14	2015	A	N9-C4	-5.91	1.34	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	14	2247	A	C6-N1	-5.91	1.31	1.35
26	14	1612	C	C2-O2	5.90	1.29	1.24
26	14	1950	G	C5-C4	5.89	1.42	1.38
26	1H	2032	G	C6-N1	-5.88	1.35	1.39
26	14	204	A	N7-C5	-5.88	1.35	1.39
26	14	1616	A	N9-C4	-5.87	1.34	1.37
26	1H	2379	G	C2-N3	5.87	1.37	1.32
26	14	2330	G	C2-N3	5.86	1.37	1.32
26	1H	693	C	N3-C4	-5.84	1.29	1.33
26	1H	2062	A	C5-C6	5.83	1.46	1.41
26	1H	838	C	N1-C6	-5.81	1.33	1.37
26	14	792	G	C6-N1	-5.80	1.35	1.39
26	1H	330	A	N9-C4	-5.80	1.34	1.37
26	1H	682	G	C5-C4	-5.80	1.34	1.38
26	14	2703	C	N1-C6	-5.80	1.33	1.37
26	1H	2448	A	C5-C6	-5.80	1.35	1.41
53	O8	42	TRP	CB-CG	5.80	1.60	1.50
26	1H	805	G	N9-C8	-5.79	1.33	1.37
1	13	1502	A	N9-C4	-5.79	1.34	1.37
26	1H	2082	A	N9-C4	-5.79	1.34	1.37
26	1H	1434	A	N9-C4	-5.79	1.34	1.37
26	1H	774	A	N9-C8	5.77	1.42	1.37
26	14	1678	G	N3-C4	-5.77	1.31	1.35
26	1H	2708	G	C2-N3	5.77	1.37	1.32
26	14	784	A	C6-N1	-5.77	1.31	1.35
26	14	1307	A	C6-N1	-5.76	1.31	1.35
26	1H	265	A	C5-C6	-5.76	1.35	1.41
26	14	1354	A	N9-C4	-5.75	1.34	1.37
26	1H	1616	A	N7-C5	-5.74	1.35	1.39
26	14	567	A	N9-C4	-5.74	1.34	1.37
26	1H	1566	A	C8-N7	5.74	1.35	1.31
26	1H	71	A	C5-C4	5.73	1.42	1.38
26	1H	2451	A	N3-C4	-5.73	1.31	1.34
26	14	945	A	N7-C5	-5.72	1.35	1.39
26	1H	689	A	N9-C4	-5.71	1.34	1.37
26	14	2606	C	N1-C6	-5.71	1.33	1.37
26	14	211	A	N7-C5	-5.71	1.35	1.39
26	1H	138	G	N7-C5	5.70	1.42	1.39
26	1H	2509	G	C5-C4	-5.70	1.34	1.38
26	1H	1799	G	N3-C4	5.69	1.39	1.35
26	1H	2082	A	N3-C4	-5.69	1.31	1.34
26	1H	2072	G	C8-N7	-5.69	1.27	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	14	2612	C	N3-C4	5.68	1.38	1.33
26	1H	1668	A	N3-C4	-5.66	1.31	1.34
55	M5	54	GLU	CG-CD	5.66	1.60	1.51
26	1H	1823	G	C6-N1	-5.65	1.35	1.39
26	14	204	A	C5-C4	-5.65	1.34	1.38
26	14	2725	A	N9-C4	-5.64	1.34	1.37
26	1H	2246	G	N9-C8	-5.64	1.33	1.37
26	14	2359	C	N3-C4	-5.63	1.30	1.33
27	16	119	A	N9-C4	5.63	1.41	1.37
26	1H	201	C	N1-C6	-5.62	1.33	1.37
26	1H	990	A	N7-C5	-5.62	1.35	1.39
26	1H	1556	C	N1-C6	-5.62	1.33	1.37
26	1H	832	G	C2-N3	-5.61	1.28	1.32
26	14	74	A	C5-C6	-5.61	1.35	1.41
26	1H	2502	G	C6-N1	-5.61	1.35	1.39
26	14	2542	A	N7-C5	5.60	1.42	1.39
26	1H	598	G	N7-C5	-5.60	1.35	1.39
37	88	139	GLU	CG-CD	5.60	1.60	1.51
26	1H	2712	U	C2-O2	-5.59	1.17	1.22
26	1H	1815	A	N3-C4	-5.59	1.31	1.34
1	13	792	A	N7-C5	-5.58	1.35	1.39
26	14	2430	A	N7-C5	-5.58	1.35	1.39
26	14	204	A	N3-C4	-5.58	1.31	1.34
26	1H	239	U	C2-N3	-5.58	1.33	1.37
26	1H	1616	A	N9-C8	5.58	1.42	1.37
26	14	330	A	N9-C4	-5.58	1.34	1.37
26	14	204	A	C5-C6	-5.56	1.36	1.41
26	1H	1210	A	C5-C6	-5.56	1.36	1.41
26	14	1997	G	C2-N3	5.56	1.37	1.32
1	13	1408	A	N7-C5	-5.55	1.35	1.39
26	1H	1632	A	C5-C6	-5.54	1.36	1.41
26	14	2392	A	C5-C4	5.54	1.42	1.38
26	1H	1322	A	N9-C4	-5.53	1.34	1.37
26	14	676	A	C5-C4	5.52	1.42	1.38
1	13	808	C	N1-C6	-5.51	1.33	1.37
26	1H	1354	A	C5-C6	-5.51	1.36	1.41
26	1H	1678	G	N9-C4	-5.51	1.33	1.38
1	1G	1473	A	N9-C4	-5.50	1.34	1.37
26	14	211	A	C5-C6	-5.50	1.36	1.41
26	1H	1246	A	N7-C5	-5.50	1.35	1.39
26	1H	2346	A	N9-C4	-5.50	1.34	1.37
26	14	2332	U	N1-C2	5.49	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	739	G	C5-C6	-5.49	1.36	1.42
26	1H	685	A	N7-C5	-5.49	1.35	1.39
26	1H	973	A	N9-C4	-5.48	1.34	1.37
26	1H	2764	A	N9-C4	-5.48	1.34	1.37
26	14	828	U	N3-C4	-5.48	1.33	1.38
26	14	774	A	C5-C6	-5.48	1.36	1.41
26	14	1605	C	N1-C6	-5.48	1.33	1.37
26	14	945	A	N3-C4	-5.48	1.31	1.34
26	14	1308	A	N7-C5	-5.47	1.35	1.39
26	14	2082	A	N7-C5	-5.47	1.35	1.39
26	1H	675	A	C5-C6	-5.47	1.36	1.41
26	1H	1698	A	C5-C6	-5.46	1.36	1.41
26	1H	432	A	C5-C6	-5.45	1.36	1.41
26	14	1313	U	C2-O2	-5.45	1.17	1.22
26	14	1674	G	N7-C5	-5.45	1.35	1.39
26	1H	2577	A	C6-N1	-5.44	1.31	1.35
26	1H	2247	A	N9-C4	-5.44	1.34	1.37
26	1H	2254	C	N1-C2	-5.43	1.34	1.40
26	1H	952	G	C5-C4	-5.43	1.34	1.38
26	1H	188	G	C2-N3	5.42	1.37	1.32
26	1H	1977	A	N9-C8	-5.42	1.33	1.37
26	14	2433	A	N9-C4	-5.42	1.34	1.37
1	13	1498	U	N1-C2	5.41	1.43	1.38
26	14	1342	A	N3-C4	-5.41	1.31	1.34
26	1H	933	A	N3-C4	-5.41	1.31	1.34
26	1H	1599	C	N3-C4	-5.40	1.30	1.33
26	1H	1638	C	N3-C4	-5.40	1.30	1.33
26	1H	1950	G	C5-C4	5.40	1.42	1.38
26	14	784	A	N3-C4	-5.40	1.31	1.34
26	1H	2277	G	N9-C8	-5.40	1.34	1.37
26	14	2506	U	N3-C4	5.40	1.43	1.38
26	1H	2009	G	C5-C4	-5.39	1.34	1.38
26	1H	57	C	N3-C4	-5.39	1.30	1.33
26	14	777	A	N9-C4	-5.39	1.34	1.37
26	14	122	G	N9-C4	-5.38	1.33	1.38
1	1G	354	G	N7-C5	-5.38	1.36	1.39
26	14	664	C	N1-C6	-5.38	1.33	1.37
26	1H	774	A	C2-N3	-5.37	1.28	1.33
37	88	82	ARG	CA-C	5.37	1.67	1.52
26	14	1558	A	N3-C4	-5.37	1.31	1.34
26	14	2688	U	N3-C4	-5.37	1.33	1.38
26	1H	2609	U	N1-C6	-5.37	1.33	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	2361	A	N9-C4	-5.35	1.34	1.37
26	1H	2230	G	C2-N3	-5.34	1.28	1.32
26	1H	2072	G	N3-C4	5.33	1.39	1.35
26	1H	1382	G	C5-C6	-5.32	1.37	1.42
26	1H	1336	A	N1-C2	-5.32	1.29	1.34
26	1H	1681	G	N9-C4	-5.32	1.33	1.38
26	1H	2579	C	N3-C4	-5.31	1.30	1.33
27	16	112	G	N3-C4	5.30	1.39	1.35
26	14	1972	A	N9-C8	-5.29	1.33	1.37
24	3L	37	A	N9-C4	5.29	1.41	1.37
26	14	788	A	N7-C5	-5.29	1.36	1.39
26	1H	608	A	N9-C4	-5.29	1.34	1.37
26	1H	685	A	C5-C6	-5.29	1.36	1.41
26	1H	946	G	C5-C4	-5.29	1.34	1.38
26	1H	1204	A	N7-C5	-5.29	1.36	1.39
26	1H	774	A	N3-C4	-5.28	1.31	1.34
28	11	144	ALA	CA-CB	-5.26	1.41	1.52
26	1H	933	A	C6-N1	-5.26	1.31	1.35
26	14	954	G	C5-C6	5.26	1.47	1.42
26	1H	1219	G	N3-C4	-5.25	1.31	1.35
26	14	1384	A	N3-C4	-5.25	1.31	1.34
55	Q8	40	GLU	CG-CD	5.25	1.59	1.51
26	1H	2318	G	N3-C4	-5.25	1.31	1.35
26	14	1785	A	N7-C5	-5.25	1.36	1.39
26	1H	2392	A	N7-C5	-5.25	1.36	1.39
26	1H	265	A	N9-C4	-5.24	1.34	1.37
24	3L	76	A	C5-C4	5.24	1.42	1.38
1	13	1502	A	P-O5'	-5.24	1.54	1.59
26	1H	123	G	C6-N1	-5.24	1.35	1.39
26	14	2025	C	N1-C6	-5.24	1.34	1.37
26	1H	262	A	N9-C4	-5.24	1.34	1.37
26	1H	581	C	N3-C4	-5.24	1.30	1.33
26	1H	1968	G	C8-N7	-5.24	1.27	1.30
26	1H	452	G	N7-C5	5.23	1.42	1.39
26	1H	1814	G	C6-N1	-5.23	1.35	1.39
26	14	2713	A	C5-C4	5.23	1.42	1.38
26	1H	448	U	N1-C6	-5.23	1.33	1.38
26	1H	1915	U	N1-C2	5.23	1.43	1.38
26	1H	205	G	C2-N3	5.22	1.36	1.32
26	1H	1258	C	N1-C6	-5.22	1.34	1.37
26	1H	1313	U	N1-C2	-5.22	1.33	1.38
26	14	773	U	C2-O2	-5.22	1.17	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	2377	A	N9-C4	-5.22	1.34	1.37
26	1H	1189	A	C5-C6	-5.22	1.36	1.41
26	1H	2432	A	C5-C4	-5.22	1.35	1.38
26	14	1204	A	N9-C4	-5.22	1.34	1.37
26	14	744	G	N7-C5	-5.21	1.36	1.39
26	1H	792	G	C5-C6	-5.19	1.37	1.42
26	1H	1355	G	N1-C2	-5.19	1.33	1.37
26	1H	2048	G	N3-C4	-5.19	1.31	1.35
26	14	509	C	N1-C6	-5.18	1.34	1.37
26	1H	1637	A	C6-N1	-5.18	1.31	1.35
26	14	954	G	N7-C5	5.18	1.42	1.39
26	14	1786	A	N7-C5	-5.17	1.36	1.39
26	1H	140	A	N7-C5	-5.16	1.36	1.39
26	1H	1931	U	C2-N3	-5.16	1.34	1.37
26	1H	663	G	N9-C8	-5.16	1.34	1.37
26	1H	784	A	N3-C4	-5.15	1.31	1.34
26	1H	1367	A	N9-C4	-5.15	1.34	1.37
1	13	535	A	C6-N1	-5.15	1.31	1.35
26	1H	138	G	N3-C4	5.15	1.39	1.35
26	1H	2452	C	N1-C6	-5.14	1.34	1.37
26	1H	746	A	C5-C6	-5.14	1.36	1.41
26	1H	2053	G	C2-N3	-5.14	1.28	1.32
26	14	1950	G	N1-C2	5.14	1.41	1.37
26	1H	1825	A	C6-N6	-5.12	1.29	1.33
26	14	2584	U	N1-C2	5.12	1.43	1.38
26	1H	2230	G	N3-C4	-5.12	1.31	1.35
26	1H	132	G	C5-C4	5.12	1.42	1.38
26	1H	667	U	C2-N3	5.11	1.41	1.37
26	14	2441	C	N3-C4	-5.11	1.30	1.33
26	14	779	U	C2-N3	-5.10	1.34	1.37
26	1H	398	G	N7-C5	-5.10	1.36	1.39
26	1H	681	G	C6-N1	-5.10	1.35	1.39
26	1H	783	A	C6-N1	-5.10	1.31	1.35
26	1H	530	G	N9-C8	5.10	1.41	1.37
26	1H	2713	A	N1-C2	5.10	1.39	1.34
26	1H	805	G	C6-O6	-5.10	1.19	1.24
26	1H	138	G	C6-N1	5.09	1.43	1.39
26	1H	1135	C	N3-C4	-5.09	1.30	1.33
26	1H	613	U	C2-N3	-5.09	1.34	1.37
26	1H	1213	A	C5-C6	-5.08	1.36	1.41
23	2L	77	A	N9-C4	-5.08	1.34	1.37
26	1H	2247	A	C6-N1	-5.08	1.31	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	1H	338	G	N7-C5	-5.08	1.36	1.39
26	14	1845	G	C5-C4	-5.08	1.34	1.38
26	1H	447	A	N3-C4	-5.07	1.31	1.34
26	1H	2233	U	N1-C2	-5.07	1.33	1.38
26	14	832	G	C5-C4	-5.06	1.34	1.38
18	9I	83	GLU	CG-CD	5.06	1.59	1.51
26	1H	682	G	N9-C8	-5.06	1.34	1.37
26	1H	2058	A	N3-C4	-5.06	1.31	1.34
26	14	567	A	N3-C4	-5.06	1.31	1.34
26	1H	1385	G	N9-C4	-5.05	1.33	1.38
26	1H	1349	A	C5-C4	5.05	1.42	1.38
26	1H	2070	G	N9-C8	-5.05	1.34	1.37
26	14	2385	C	N3-C4	5.05	1.37	1.33
26	1H	2271	G	C2-N3	5.05	1.36	1.32
49	G5	5	GLU	CG-CD	5.05	1.59	1.51
26	1H	2713	A	N9-C8	5.05	1.41	1.37
26	1H	793	A	C5-C6	-5.04	1.36	1.41
26	1H	807	U	N1-C6	-5.04	1.33	1.38
26	1H	2430	A	C5-C4	5.04	1.42	1.38
26	1H	1728	G	N3-C4	5.04	1.39	1.35
26	14	698	C	N1-C6	-5.04	1.34	1.37
26	1H	465	G	C6-O6	5.04	1.28	1.24
26	14	138	G	N9-C4	5.04	1.42	1.38
26	1H	695	G	C6-N1	-5.03	1.36	1.39
26	1H	1418	G	N7-C5	5.03	1.42	1.39
26	14	801	G	C6-N1	-5.03	1.36	1.39
26	14	1353	A	N3-C4	-5.03	1.31	1.34
26	1H	2297	C	N3-C4	-5.03	1.30	1.33
26	14	2518	A	N9-C4	-5.03	1.34	1.37
26	1H	2059	A	N9-C4	-5.02	1.34	1.37
26	14	1286	A	N7-C5	-5.02	1.36	1.39
26	1H	1613	G	C6-N1	-5.01	1.36	1.39

All (7504) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1899	G	N3-C4-N9	-26.68	109.99	126.00
26	1H	676	A	C2-N3-C4	-22.63	99.29	110.60
26	1H	2430	A	C2-N3-C4	-22.43	99.39	110.60
26	1H	1899	G	N3-C4-C5	21.40	139.30	128.60
26	1H	783	A	C2-N3-C4	-21.20	100.00	110.60
26	1H	945	A	C6-C5-N7	-20.79	117.75	132.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1786	A	C2-N3-C4	-19.80	100.70	110.60
26	14	1332	G	N3-C4-N9	-19.74	114.15	126.00
26	14	1786	A	C5-N7-C8	-19.43	94.19	103.90
26	14	783	A	C2-N3-C4	-19.31	100.94	110.60
26	1H	74	A	C2-N3-C4	-19.29	100.96	110.60
26	1H	945	A	N1-C6-N6	19.08	130.05	118.60
26	1H	801	G	O5'-P-OP2	-18.79	88.15	110.70
26	14	74	A	C2-N3-C4	-18.33	101.44	110.60
26	14	1786	A	N7-C8-N9	17.88	122.74	113.80
26	1H	917	A	N1-C2-N3	17.78	138.19	129.30
26	14	1332	G	N3-C4-C5	17.72	137.46	128.60
26	1H	2490	G	C5-N7-C8	-17.69	95.45	104.30
26	1H	1332	G	N3-C4-N9	-17.61	115.44	126.00
26	1H	1899	G	C2-N3-C4	-17.39	103.21	111.90
26	14	741	G	O5'-P-OP1	-17.32	89.92	110.70
26	1H	917	A	C2-N3-C4	-17.19	102.00	110.60
26	1H	2598	A	O5'-P-OP1	-17.12	90.15	110.70
26	1H	2448	A	N1-C6-N6	16.89	128.73	118.60
26	1H	945	A	N7-C8-N9	16.80	122.20	113.80
26	1H	945	A	C5-N7-C8	-16.75	95.52	103.90
26	1H	774	A	N3-C4-C5	16.59	138.41	126.80
26	1H	1698	A	C2-N3-C4	-16.52	102.34	110.60
26	14	1332	G	C2-N3-C4	-16.28	103.76	111.90
26	1H	1899	G	N3-C2-N2	-16.28	108.51	119.90
26	1H	1639	U	O5'-P-OP2	-16.23	91.09	105.70
26	1H	1332	G	N3-C4-C5	16.20	136.70	128.60
26	1H	945	A	C4-C5-C6	15.96	124.98	117.00
26	1H	1786	A	N1-C2-N3	15.94	137.27	129.30
26	1H	774	A	C2-N3-C4	-15.94	102.63	110.60
26	1H	1931	U	N3-C2-O2	-15.80	111.14	122.20
26	14	1698	A	N1-C6-N6	15.68	128.00	118.60
26	14	1984	G	O5'-P-OP2	-15.67	91.60	105.70
26	1H	2490	G	C4-C5-N7	15.64	117.06	110.80
26	1H	51	G	O5'-P-OP1	-15.51	91.74	105.70
26	14	2275	C	C6-N1-C2	-15.50	114.10	120.30
26	1H	774	A	N3-C4-N9	-15.45	115.04	127.40
26	1H	71	A	C2-N3-C4	-15.41	102.90	110.60
26	1H	1678	G	C2-N3-C4	-15.40	104.20	111.90
26	14	2430	A	C2-N3-C4	-15.38	102.91	110.60
26	1H	1332	G	C2-N3-C4	-15.36	104.22	111.90
26	1H	1786	A	C5-N7-C8	-15.17	96.32	103.90
26	1H	676	A	C5-N7-C8	-15.06	96.37	103.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	945	A	N1-C6-N6	15.05	127.63	118.60
26	14	1253	A	N1-C6-N6	14.99	127.60	118.60
26	14	774	A	C2-N3-C4	-14.94	103.13	110.60
26	1H	2287	A	C2-N3-C4	-14.92	103.14	110.60
26	14	2436	G	O5'-P-OP1	-14.78	92.40	105.70
26	14	1899	G	C2-N3-C4	-14.73	104.53	111.90
26	14	783	A	C5-N7-C8	-14.66	96.57	103.90
1	13	792	A	C2-N3-C4	-14.66	103.27	110.60
26	1H	676	A	N3-C4-N9	-14.55	115.76	127.40
26	1H	140	A	C5-N7-C8	-14.48	96.66	103.90
26	1H	1829	A	O5'-P-OP1	-14.46	92.68	105.70
26	1H	1616	A	C5-N7-C8	-14.44	96.68	103.90
26	1H	621	A	C5-N7-C8	-14.43	96.68	103.90
1	13	792	A	N1-C6-N6	14.42	127.25	118.60
26	1H	860	U	N3-C2-O2	-14.42	112.11	122.20
26	1H	676	A	N3-C4-C5	14.30	136.81	126.80
26	1H	774	A	C5-N7-C8	-14.29	96.75	103.90
26	1H	1786	A	C6-C5-N7	-14.22	122.35	132.30
26	1H	1982	C	O5'-P-OP2	-14.16	92.96	105.70
26	14	1612	C	C6-N1-C2	14.04	125.92	120.30
26	1H	863	A	O5'-P-OP2	-14.00	93.10	105.70
26	1H	860	U	C4-C5-C6	13.93	128.06	119.70
26	1H	2374	C	C5-C6-N1	-13.92	114.04	121.00
26	1H	49	A	O5'-P-OP2	-13.82	93.26	105.70
26	14	2712	U	C5-C6-N1	-13.80	115.80	122.70
26	1H	783	A	C5-N7-C8	-13.79	97.00	103.90
26	1H	1786	A	N7-C8-N9	13.79	120.69	113.80
26	1H	621	A	C2-N3-C4	-13.66	103.77	110.60
24	3K	76	A	C5-N7-C8	-13.54	97.13	103.90
26	14	774	A	N3-C4-C5	13.51	136.26	126.80
26	14	945	A	C6-C5-N7	-13.49	122.86	132.30
26	14	737	C	N1-C2-O2	-13.45	110.83	118.90
26	14	2273	A	O5'-P-OP2	-13.41	93.63	105.70
26	1H	265	A	C2-N3-C4	-13.37	103.92	110.60
26	1H	2375	G	C8-N9-C4	13.30	111.72	106.40
26	1H	2490	G	N3-C4-C5	13.28	135.24	128.60
26	1H	1786	A	N1-C6-N6	13.28	126.56	118.60
26	14	2430	A	N1-C6-N6	13.27	126.56	118.60
26	1H	1899	G	N9-C4-C5	13.24	110.69	105.40
26	1H	210	C	C6-N1-C2	13.21	125.58	120.30
26	1H	140	A	N7-C8-N9	13.21	120.40	113.80
26	14	330	A	C2-N3-C4	-13.20	104.00	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	945	A	C4-C5-N7	13.14	117.27	110.70
26	1H	2584	U	N3-C2-O2	-13.06	113.06	122.20
26	1H	245	G	N1-C6-O6	13.05	127.73	119.90
26	1H	2430	A	N1-C2-N3	13.03	135.81	129.30
26	1H	2448	A	C5-C6-N6	-13.02	113.28	123.70
26	1H	2346	A	N1-C2-N3	13.00	135.80	129.30
26	1H	1332	G	C5-N7-C8	-12.99	97.80	104.30
26	1H	2053	G	C5-C6-O6	-12.96	120.82	128.60
1	13	690	G	C6-C5-N7	-12.95	122.63	130.40
26	14	783	A	N1-C6-N6	12.94	126.37	118.60
26	14	783	A	C5-C6-N1	-12.93	111.23	117.70
26	14	2042	A	O5'-P-OP2	-12.89	94.10	105.70
26	1H	1678	G	N3-C4-C5	12.88	135.04	128.60
26	1H	2713	A	C5-N7-C8	-12.88	97.46	103.90
26	1H	71	A	C5-N7-C8	-12.88	97.46	103.90
26	1H	698	C	C6-N1-C2	12.85	125.44	120.30
26	1H	1405	U	O5'-P-OP2	-12.84	94.14	105.70
26	1H	1678	G	C4-C5-N7	12.83	115.93	110.80
1	13	792	A	C4-C5-N7	12.81	117.11	110.70
26	1H	945	A	C8-N9-C4	-12.77	100.69	105.80
26	1H	676	A	N7-C8-N9	12.77	120.18	113.80
26	1H	1678	G	C5-N7-C8	-12.75	97.92	104.30
26	1H	2700	C	C6-N1-C2	12.71	125.39	120.30
26	14	1786	A	C4-C5-N7	12.69	117.04	110.70
26	1H	641	C	O5'-P-OP1	-12.67	94.30	105.70
26	1H	1496	A	C8-N9-C4	-12.67	100.73	105.80
26	1H	2430	A	N1-C6-N6	12.65	126.19	118.60
26	1H	2490	G	N7-C8-N9	12.65	119.42	113.10
26	1H	140	A	N1-C6-N6	12.65	126.19	118.60
26	1H	1614	A	C5-N7-C8	-12.64	97.58	103.90
27	16	47	C	C6-N1-C2	12.64	125.36	120.30
26	14	1786	A	C2-N3-C4	-12.62	104.29	110.60
26	1H	783	A	C5-C6-N1	-12.60	111.40	117.70
26	1H	1006	C	O5'-P-OP1	-12.59	94.37	105.70
24	3K	76	A	N1-C6-N6	12.54	126.13	118.60
26	14	510	C	O5'-P-OP2	-12.54	94.41	105.70
26	14	1698	A	C2-N3-C4	-12.51	104.35	110.60
26	1H	808	G	O5'-P-OP2	-12.50	94.45	105.70
26	14	140	A	N7-C8-N9	12.50	120.05	113.80
26	1H	676	A	C5-C6-N1	-12.50	111.45	117.70
26	14	1816	G	O5'-P-OP1	-12.46	94.48	105.70
26	14	1332	G	C5-N7-C8	-12.46	98.07	104.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	774	A	N3-C4-N9	-12.44	117.45	127.40
26	1H	945	A	O4'-C1'-N9	12.41	118.13	108.20
26	14	1698	A	C6-C5-N7	-12.41	123.61	132.30
26	14	2023	G	O5'-P-OP2	-12.41	94.53	105.70
26	14	828	U	C5-C4-O4	12.35	133.31	125.90
26	1H	1698	A	C5-N7-C8	-12.34	97.73	103.90
26	1H	245	G	C6-C5-N7	-12.34	123.00	130.40
26	14	1698	A	C4-C5-N7	12.31	116.85	110.70
26	1H	1496	A	N7-C8-N9	12.30	119.95	113.80
26	1H	2468	G	O4'-C1'-N9	12.29	118.04	108.20
26	1H	1616	A	C4-C5-N7	12.29	116.84	110.70
26	14	189	G	C8-N9-C4	12.28	111.31	106.40
26	1H	728	G	C8-N9-C4	12.24	111.30	106.40
26	14	2430	A	C5-C6-N1	-12.23	111.58	117.70
26	1H	774	A	C5-C6-N1	-12.21	111.59	117.70
26	14	2079	U	O5'-P-OP1	-12.21	94.71	105.70
26	1H	1614	A	C2-N3-C4	-12.19	104.50	110.60
26	14	1698	A	C5-N7-C8	-12.16	97.82	103.90
26	14	74	A	N3-C4-C5	12.16	135.31	126.80
26	1H	34	C	O5'-P-OP1	-12.15	94.76	105.70
26	14	140	A	C5-N7-C8	-12.14	97.83	103.90
26	14	2873	A	N1-C2-N3	12.12	135.36	129.30
26	1H	2490	G	C2-N3-C4	-12.11	105.85	111.90
26	1H	2442	C	N3-C4-N4	12.10	126.47	118.00
26	14	1616	A	C2-N3-C4	-12.09	104.55	110.60
26	14	2700	C	C6-N1-C2	12.06	125.12	120.30
26	1H	120	U	C5-C6-N1	-12.06	116.67	122.70
26	1H	2346	A	C2-N3-C4	-12.04	104.58	110.60
26	1H	2377	A	C8-N9-C4	12.05	110.62	105.80
26	1H	846	C	O5'-P-OP1	-12.04	94.87	105.70
26	1H	74	A	C5-C6-N1	-12.00	111.70	117.70
26	1H	2375	G	C5-C6-O6	-12.00	121.40	128.60
26	14	74	A	C5-C6-N1	-12.00	111.70	117.70
26	14	679	C	N1-C2-O2	-12.00	111.70	118.90
26	1H	814	C	O5'-P-OP2	-11.96	94.93	105.70
26	1H	329	G	O5'-P-OP2	-11.94	94.95	105.70
26	14	528	A	C2-N3-C4	-11.92	104.64	110.60
26	1H	860	U	C5-C6-N1	-11.92	116.74	122.70
26	1H	691	C	C6-N1-C2	11.91	125.06	120.30
26	1H	140	A	C4-C5-N7	11.87	116.63	110.70
26	1H	245	G	C5-C6-O6	-11.83	121.50	128.60
26	1H	2346	A	C6-C5-N7	-11.82	124.03	132.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	783	A	N3-C4-C5	11.78	135.04	126.80
26	1H	201	C	C6-N1-C2	11.77	125.01	120.30
26	1H	2713	A	C2-N3-C4	-11.76	104.72	110.60
1	13	1502	A	C2-N3-C4	-11.75	104.72	110.60
26	1H	2392	A	C5-N7-C8	-11.69	98.06	103.90
26	1H	774	A	C6-N1-C2	11.62	125.57	118.60
1	13	792	A	O4'-C1'-N9	11.61	117.48	108.20
26	1H	2374	C	C6-N1-C2	11.60	124.94	120.30
26	1H	1204	A	O4'-C1'-N9	11.58	117.46	108.20
26	14	991	C	O5'-P-OP1	-11.57	95.29	105.70
26	14	801	G	N1-C6-O6	-11.56	112.96	119.90
24	3K	76	A	N7-C8-N9	11.55	119.58	113.80
26	1H	1257	C	N1-C2-O2	-11.55	111.97	118.90
26	14	676	A	C2-N3-C4	-11.54	104.83	110.60
26	14	1600	C	O5'-P-OP2	-11.53	95.32	105.70
1	1G	108	G	C5-C6-O6	-11.50	121.70	128.60
26	1H	2622	C	C6-N1-C2	11.47	124.89	120.30
26	1H	828	U	C5-C4-O4	11.45	132.77	125.90
1	1G	332	G	C8-N9-C4	11.45	110.98	106.40
26	1H	633	A	N1-C6-N6	11.44	125.47	118.60
26	1H	2346	A	N7-C8-N9	11.44	119.52	113.80
26	1H	1616	A	N7-C8-N9	11.43	119.51	113.80
1	13	1502	A	O5'-P-OP2	-11.42	95.42	105.70
26	1H	683	C	N3-C4-C5	11.39	126.46	121.90
26	1H	758	C	N3-C4-C5	11.39	126.46	121.90
26	14	835	A	O5'-P-OP2	-11.39	95.45	105.70
26	14	2518	A	N1-C6-N6	11.38	125.43	118.60
26	1H	735	A	C8-N9-C4	11.37	110.35	105.80
1	13	792	A	C6-C5-N7	-11.36	124.35	132.30
26	1H	2392	A	C2-N3-C4	-11.35	104.92	110.60
26	1H	1021	A	C2-N3-C4	-11.34	104.93	110.60
26	14	1786	A	C8-N9-C4	-11.34	101.27	105.80
26	1H	578	A	O5'-P-OP2	-11.33	95.50	105.70
26	1H	945	A	C4-N9-C1'	11.32	146.69	126.30
26	1H	676	A	C8-N9-C4	-11.32	101.27	105.80
26	1H	621	A	N1-C6-N6	11.31	125.38	118.60
26	1H	787	U	O5'-P-OP1	11.29	124.25	110.70
26	1H	2430	A	N3-C4-C5	11.27	134.69	126.80
26	1H	141	A	C5-N7-C8	-11.24	98.28	103.90
26	1H	2573	C	C6-N1-C2	-11.20	115.82	120.30
27	1J	30	C	C6-N1-C2	-11.20	115.82	120.30
26	14	1602	U	O5'-P-OP2	11.19	124.13	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1404	C	O5'-P-OP2	-11.17	95.65	105.70
26	14	621	A	C2-N3-C4	-11.17	105.02	110.60
26	14	2392	A	C2-N3-C4	-11.16	105.02	110.60
26	1H	144	C	C6-N1-C2	11.16	124.76	120.30
1	13	899	C	C6-N1-C2	11.12	124.75	120.30
26	1H	1349	A	O5'-P-OP1	-11.11	95.71	105.70
26	1H	814	C	C5-C6-N1	-11.09	115.46	121.00
23	2K	21	U	C5-C4-O4	-11.09	119.25	125.90
26	1H	2584	U	C5-C4-O4	11.08	132.55	125.90
26	1H	2712	U	C2-N1-C1'	11.06	130.97	117.70
26	1H	140	A	C6-C5-N7	-11.06	124.56	132.30
26	1H	1990	C	C6-N1-C2	-11.06	115.88	120.30
26	1H	2375	G	N9-C4-C5	-11.05	100.98	105.40
26	14	1899	G	N1-C2-N3	11.03	130.52	123.90
1	13	525	C	C5-C6-N1	11.01	126.51	121.00
26	1H	679	C	C5-C6-N1	-11.01	115.49	121.00
26	1H	1931	U	N1-C2-N3	11.01	121.51	114.90
26	1H	1786	A	C4-C5-N7	11.01	116.20	110.70
26	1H	1225	C	C6-N1-C2	10.99	124.70	120.30
26	1H	621	A	C4-C5-N7	10.99	116.19	110.70
26	1H	860	U	C2-N1-C1'	10.99	130.88	117.70
26	1H	621	A	N3-C4-C5	10.98	134.49	126.80
1	13	789	U	C5-C4-O4	10.98	132.49	125.90
26	14	320	A	O5'-P-OP2	-10.98	95.82	105.70
26	14	783	A	N3-C4-N9	-10.96	118.63	127.40
26	1H	659	C	C6-N1-C2	10.96	124.68	120.30
26	1H	852	G	O5'-P-OP2	-10.95	95.85	105.70
26	1H	2330	G	C5-C6-O6	-10.94	122.04	128.60
26	14	1825	A	O5'-P-OP2	-10.94	95.85	105.70
26	14	783	A	N7-C8-N9	10.93	119.27	113.80
36	35	147	LEU	CA-CB-CG	10.93	140.44	115.30
26	1H	839	U	O5'-P-OP2	-10.90	95.89	105.70
26	1H	1632	A	N1-C6-N6	10.90	125.14	118.60
26	1H	252	G	O5'-P-OP2	-10.89	95.90	105.70
26	14	782	A	N1-C6-N6	-10.89	112.07	118.60
26	14	793	A	O5'-P-OP2	-10.89	95.90	105.70
24	3K	76	A	C4-C5-N7	10.89	116.14	110.70
26	1H	1382	G	C5-C6-O6	-10.88	122.07	128.60
26	1H	1950	G	N7-C8-N9	10.88	118.54	113.10
26	14	1253	A	C5-C6-N6	-10.88	115.00	123.70
1	1G	108	G	C4-C5-N7	10.88	115.15	110.80
26	14	1029	A	C8-N9-C4	10.85	110.14	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1786	A	C6-C5-N7	-10.85	124.70	132.30
26	1H	1496	A	C5-N7-C8	-10.85	98.47	103.90
26	1H	193	U	C5-C6-N1	-10.84	117.28	122.70
26	1H	2507	C	N3-C2-O2	-10.84	114.31	121.90
26	1H	140	A	C8-N9-C4	-10.83	101.47	105.80
26	1H	111	A	O5'-P-OP2	-10.82	95.96	105.70
26	1H	2502	G	N3-C4-C5	-10.82	123.19	128.60
26	14	945	A	C4-C5-N7	10.81	116.11	110.70
26	1H	2430	A	C5-N7-C8	-10.81	98.50	103.90
26	1H	252	G	O5'-P-OP1	10.80	123.67	110.70
26	1H	1950	G	C2-N3-C4	-10.80	106.50	111.90
26	1H	774	A	C4-C5-N7	10.79	116.09	110.70
26	1H	2607	G	C6-C5-N7	-10.77	123.94	130.40
26	1H	945	A	O5'-P-OP2	-10.77	96.01	105.70
26	1H	783	A	N7-C8-N9	10.75	119.17	113.80
26	1H	330	A	C2-N3-C4	-10.74	105.23	110.60
26	1H	1950	G	C5-N7-C8	-10.73	98.93	104.30
26	14	71	A	N1-C6-N6	10.72	125.03	118.60
26	1H	964	C	O5'-P-OP1	-10.72	96.05	105.70
26	14	2275	C	C5-C6-N1	10.71	126.36	121.00
26	1H	1825	A	N1-C6-N6	-10.70	112.18	118.60
1	13	966	G	C5-C6-O6	-10.70	122.18	128.60
26	1H	2518	A	C5-N7-C8	-10.69	98.56	103.90
26	1H	678	C	C2-N3-C4	-10.69	114.56	119.90
26	1H	2595	G	C4-C5-N7	10.68	115.07	110.80
26	1H	2713	A	N7-C8-N9	10.67	119.14	113.80
26	1H	2085	C	O5'-P-OP2	-10.67	96.10	105.70
1	13	792	A	C5-N7-C8	-10.66	98.57	103.90
26	1H	1558	A	N1-C6-N6	10.66	124.99	118.60
26	1H	74	A	C5-N7-C8	-10.65	98.57	103.90
26	14	2430	A	C6-C5-N7	-10.65	124.84	132.30
26	14	1616	A	C5-N7-C8	-10.65	98.57	103.90
26	1H	2420	C	O5'-P-OP1	-10.64	96.12	105.70
26	1H	815	C	C6-N1-C2	10.63	124.55	120.30
26	1H	2062	A	C8-N9-C4	10.63	110.05	105.80
26	1H	2713	A	N1-C6-N6	10.63	124.98	118.60
26	14	1619	G	O5'-P-OP2	-10.63	96.13	105.70
26	1H	783	A	N3-C4-C5	10.61	134.22	126.80
26	14	856	C	O5'-P-OP1	-10.61	96.16	105.70
26	14	140	A	N1-C6-N6	10.60	124.96	118.60
26	1H	2328	A	N1-C2-N3	10.59	134.59	129.30
24	3L	76	A	N7-C8-N9	10.58	119.09	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1660	C	O5'-P-OP2	-10.57	96.19	105.70
26	14	676	A	C5-N7-C8	-10.53	98.63	103.90
26	14	1678	G	C2-N3-C4	-10.53	106.63	111.90
26	1H	103	A	C8-N9-C4	10.53	110.01	105.80
26	1H	2401	U	C5-C6-N1	10.52	127.96	122.70
26	1H	530	G	N1-C6-O6	-10.52	113.59	119.90
26	1H	783	A	O5'-P-OP2	-10.48	96.26	105.70
26	1H	621	A	C5-C6-N1	-10.47	112.46	117.70
26	14	189	G	N9-C4-C5	-10.46	101.22	105.40
26	14	1678	G	N3-C4-C5	10.46	133.83	128.60
26	1H	2699	C	C6-N1-C2	10.46	124.48	120.30
26	14	1678	G	N3-C4-N9	-10.45	119.73	126.00
26	1H	746	A	O5'-P-OP2	10.45	123.23	110.70
26	1H	461	C	N1-C2-O2	-10.41	112.65	118.90
26	1H	783	A	N3-C4-N9	-10.41	119.07	127.40
26	1H	1379	A	C5-N7-C8	-10.40	98.70	103.90
26	1H	1914	C	N3-C2-O2	-10.40	114.62	121.90
26	1H	1021	A	C5-N7-C8	-10.40	98.70	103.90
26	1H	2346	A	O4'-C1'-N9	10.40	116.52	108.20
24	3K	76	A	C6-C5-N7	-10.39	125.03	132.30
26	1H	1518	C	O5'-P-OP1	-10.38	96.36	105.70
26	1H	2636	U	O5'-P-OP1	-10.38	96.35	105.70
26	1H	1779	U	O5'-P-OP2	-10.38	96.36	105.70
26	1H	1950	G	C8-N9-C4	-10.38	102.25	106.40
26	1H	144	C	C5-C6-N1	-10.38	115.81	121.00
23	2K	21	U	N3-C4-C5	10.36	120.82	114.60
26	1H	741	G	O5'-P-OP1	-10.34	96.39	105.70
26	1H	74	A	N1-C6-N6	10.34	124.80	118.60
26	14	801	G	C5-C6-O6	10.33	134.80	128.60
26	1H	456	C	C5-C6-N1	-10.33	115.84	121.00
26	1H	1566	A	O5'-P-OP2	-10.33	96.40	105.70
26	14	74	A	N1-C6-N6	10.33	124.80	118.60
26	1H	1195	G	C5-C6-O6	-10.31	122.41	128.60
26	14	2056	G	C5-C6-O6	-10.31	122.41	128.60
1	13	738	C	C6-N1-C2	-10.30	116.18	120.30
1	13	781	A	C5-C6-N6	-10.30	115.46	123.70
26	1H	528	A	C6-N1-C2	10.30	124.78	118.60
26	1H	783	A	C6-C5-N7	-10.30	125.09	132.30
1	13	690	G	C4-C5-N7	10.29	114.91	110.80
26	1H	2346	A	C5-N7-C8	-10.27	98.77	103.90
26	14	917	A	O5'-P-OP1	-10.27	96.46	105.70
26	14	2702	U	O4'-C1'-N1	10.27	116.41	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2521	C	C6-N1-C2	10.25	124.40	120.30
1	13	1502	A	C5-N7-C8	-10.25	98.78	103.90
26	1H	774	A	N1-C6-N6	10.25	124.75	118.60
26	1H	2507	C	C6-N1-C2	-10.24	116.20	120.30
26	14	71	A	C5-N7-C8	-10.24	98.78	103.90
26	14	1496	A	N7-C8-N9	10.23	118.92	113.80
26	1H	2544	G	C5-C6-O6	-10.23	122.46	128.60
26	1H	971	C	C6-N1-C2	-10.23	116.21	120.30
26	1H	491	G	O5'-P-OP1	-10.22	96.50	105.70
1	13	789	U	N3-C2-O2	-10.21	115.05	122.20
26	1H	74	A	N3-C4-C5	10.20	133.94	126.80
26	14	2873	A	N7-C8-N9	10.19	118.90	113.80
26	14	2430	A	N1-C2-N3	10.19	134.40	129.30
26	1H	2346	A	C4-C5-C6	10.18	122.09	117.00
26	14	140	A	C4-C5-N7	10.15	115.78	110.70
26	1H	459	U	O5'-P-OP2	-10.15	96.57	105.70
26	1H	2465	C	C5-C6-N1	-10.14	115.93	121.00
26	14	2702	U	N3-C2-O2	-10.12	115.12	122.20
26	14	1930	G	O5'-P-OP1	-10.11	96.60	105.70
26	1H	192	C	C6-N1-C2	10.10	124.34	120.30
26	14	1975	G	C5-C6-O6	-10.09	122.55	128.60
26	1H	51	G	C8-N9-C4	10.09	110.44	106.40
23	2K	74	A	N1-C6-N6	10.08	124.65	118.60
26	1H	265	A	C5-N7-C8	-10.07	98.86	103.90
26	1H	692	C	C2-N3-C4	-10.07	114.86	119.90
26	1H	1332	G	N7-C8-N9	10.07	118.14	113.10
26	1H	1377	G	O5'-P-OP2	-10.07	96.64	105.70
26	14	967	C	O5'-P-OP2	-10.06	96.65	105.70
26	1H	2447	G	N3-C4-C5	-10.05	123.58	128.60
26	1H	945	A	N1-C2-N3	10.04	134.32	129.30
26	14	1204	A	C2-N3-C4	-10.04	105.58	110.60
26	1H	2311	A	N1-C2-N3	10.03	134.31	129.30
27	16	5	C	C6-N1-C2	10.03	124.31	120.30
1	1G	1397	C	C6-N1-C2	-10.03	116.29	120.30
26	1H	686	G	C8-N9-C4	10.02	110.41	106.40
26	14	189	G	C5-C6-O6	-10.02	122.59	128.60
26	14	2017	U	O5'-P-OP1	-10.01	96.69	105.70
26	1H	432	A	N1-C6-N6	9.99	124.60	118.60
26	14	784	A	C5-C6-N6	9.99	131.69	123.70
26	1H	676	A	O4'-C1'-N9	9.98	116.18	108.20
26	14	783	A	C6-C5-N7	-9.98	125.32	132.30
26	1H	120	U	C4-C5-C6	9.97	125.68	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	180	G	C8-N9-C4	9.97	110.39	106.40
26	1H	2436	G	N3-C2-N2	-9.97	112.92	119.90
26	1H	2448	A	N9-C4-C5	-9.97	101.81	105.80
26	1H	1698	A	N1-C6-N6	9.96	124.58	118.60
26	1H	659	C	C5-C6-N1	-9.96	116.02	121.00
26	1H	1558	A	O5'-P-OP1	-9.95	96.74	105.70
26	1H	2490	G	N3-C4-N9	-9.95	120.03	126.00
26	1H	2609	U	C5-C6-N1	-9.95	117.72	122.70
26	1H	236	C	C4-C5-C6	9.95	122.37	117.40
26	1H	2253	G	C5-C6-O6	-9.93	122.64	128.60
26	1H	735	A	N7-C8-N9	-9.92	108.84	113.80
1	13	963	G	N3-C4-N9	9.92	131.95	126.00
26	14	2000	G	C8-N9-C4	9.88	110.35	106.40
26	1H	2346	A	N1-C6-N6	9.87	124.52	118.60
23	2L	21	U	C5-C4-O4	-9.87	119.98	125.90
26	14	945	A	C5-N7-C8	-9.86	98.97	103.90
26	14	453	C	C6-N1-C2	9.86	124.24	120.30
26	1H	787	U	O5'-P-OP2	-9.85	96.83	105.70
26	1H	2424	C	N1-C2-O2	9.85	124.81	118.90
26	1H	788	A	N9-C4-C5	-9.85	101.86	105.80
26	14	2430	A	C4-C5-C6	9.85	121.92	117.00
26	1H	2710	C	C6-N1-C2	9.84	124.24	120.30
26	1H	972	G	N1-C6-O6	-9.83	114.00	119.90
26	1H	1204	A	C2-N3-C4	-9.83	105.69	110.60
26	14	773	U	N1-C2-N3	9.83	120.80	114.90
26	1H	945	A	C5-C6-N6	-9.83	115.84	123.70
26	1H	2639	A	C2-N3-C4	-9.82	105.69	110.60
26	14	528	A	N3-C4-C5	9.82	133.67	126.80
26	1H	860	U	N1-C2-O2	9.80	129.66	122.80
26	1H	2312	U	O5'-P-OP1	-9.80	96.88	105.70
26	1H	859	G	N3-C4-C5	9.80	133.50	128.60
26	1H	2689	U	N3-C4-O4	-9.79	112.55	119.40
26	14	1266	G	C8-N9-C4	9.79	110.32	106.40
26	14	2439	A	P-O3'-C3'	9.79	131.44	119.70
26	1H	2599	G	N1-C6-O6	-9.79	114.03	119.90
26	1H	2070	G	O5'-P-OP2	-9.78	96.90	105.70
26	1H	2379	G	C5-C6-O6	-9.78	122.73	128.60
26	1H	812	C	N1-C2-O2	-9.77	113.04	118.90
26	1H	1324	G	N1-C6-O6	9.77	125.76	119.90
26	1H	621	A	N7-C8-N9	9.76	118.68	113.80
26	1H	1899	G	C8-N9-C1'	9.75	139.68	127.00
26	1H	2442	C	C5-C4-N4	-9.75	113.38	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	783	A	N1-C2-N3	9.75	134.17	129.30
26	14	252	G	C5-N7-C8	9.74	109.17	104.30
1	13	781	A	N1-C6-N6	9.74	124.44	118.60
26	1H	2689	U	N3-C2-O2	-9.74	115.38	122.20
26	1H	1284	A	O5'-P-OP2	-9.73	96.95	105.70
26	14	1614	A	C2-N3-C4	-9.72	105.74	110.60
26	1H	1210	A	C5-N7-C8	-9.72	99.04	103.90
26	1H	2773	C	C6-N1-C2	9.72	124.19	120.30
1	13	974	A	N1-C6-N6	9.70	124.42	118.60
26	1H	2270	G	C8-N9-C4	9.69	110.27	106.40
26	14	252	G	O5'-P-OP2	-9.69	96.98	105.70
26	1H	1977	A	C2-N3-C4	-9.67	105.77	110.60
1	13	1505	G	O5'-P-OP2	9.67	122.30	110.70
26	1H	910	A	N1-C6-N6	9.67	124.40	118.60
26	14	1404	C	O5'-P-OP1	-9.66	97.00	105.70
26	1H	2607	G	C4-C5-N7	9.66	114.66	110.80
1	1G	254	G	O5'-P-OP1	-9.65	97.01	105.70
26	1H	2448	A	C6-C5-N7	-9.65	125.55	132.30
27	16	30	C	C6-N1-C2	-9.65	116.44	120.30
26	1H	2448	A	C4-C5-N7	9.64	115.52	110.70
26	1H	2698	U	C5-C6-N1	-9.63	117.88	122.70
26	14	783	A	C4-C5-N7	9.63	115.52	110.70
1	13	525	C	C6-N1-C2	-9.63	116.45	120.30
26	14	774	A	C5-N7-C8	-9.63	99.09	103.90
26	1H	446	G	N1-C6-O6	9.62	125.67	119.90
26	14	639	U	O5'-P-OP2	-9.62	97.04	105.70
26	14	1342	A	C2-N3-C4	-9.62	105.79	110.60
26	1H	2374	C	C2-N3-C4	-9.61	115.10	119.90
26	1H	2346	A	C8-N9-C4	-9.60	101.96	105.80
26	1H	2380	C	C5-C6-N1	-9.59	116.20	121.00
26	14	252	G	C4-C5-N7	-9.59	106.96	110.80
26	1H	1614	A	N1-C6-N6	9.59	124.35	118.60
26	14	2873	A	C6-C5-N7	-9.59	125.59	132.30
26	14	830	G	C8-N9-C4	9.58	110.23	106.40
26	1H	528	A	N3-C4-N9	-9.57	119.74	127.40
26	1H	1599	C	O5'-P-OP2	-9.57	97.09	105.70
24	3L	71	G	O4'-C1'-N9	9.56	115.85	108.20
26	1H	1255	U	N3-C2-O2	9.55	128.89	122.20
26	1H	1564	C	N3-C4-N4	-9.55	111.31	118.00
26	1H	679	C	C2-N3-C4	-9.55	115.13	119.90
27	16	44	G	C4-N9-C1'	-9.55	114.09	126.50
26	14	2712	U	N3-C4-O4	-9.54	112.72	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1278	A	C8-N9-C4	9.54	109.62	105.80
26	14	2712	U	N3-C2-O2	-9.54	115.52	122.20
26	1H	913	U	N3-C2-O2	-9.53	115.53	122.20
1	13	792	A	N9-C4-C5	-9.53	101.99	105.80
26	1H	783	A	C8-N9-C4	-9.53	101.99	105.80
26	14	687	C	O5'-P-OP1	-9.52	97.13	105.70
26	14	2080	G	O5'-P-OP2	-9.51	97.14	105.70
26	1H	2465	C	C6-N1-C2	9.51	124.10	120.30
26	1H	2598	A	O5'-P-OP2	9.51	122.11	110.70
26	14	2713	A	C2-N3-C4	-9.51	105.85	110.60
1	13	1508	G	O5'-P-OP1	-9.49	97.16	105.70
26	1H	2598	A	C8-N9-C4	9.49	109.59	105.80
26	1H	1528	A	C8-N9-C4	-9.48	102.01	105.80
26	1H	1192	G	C8-N9-C4	9.48	110.19	106.40
26	14	737	C	N3-C4-N4	9.48	124.64	118.00
26	14	140	A	C8-N9-C4	-9.48	102.01	105.80
26	14	114	U	C5-C4-O4	-9.48	120.21	125.90
26	1H	816	C	N3-C4-N4	9.47	124.63	118.00
1	13	525	C	N3-C4-N4	9.47	124.63	118.00
1	13	810	C	O5'-P-OP2	-9.47	97.18	105.70
26	14	133	C	C6-N1-C2	9.46	124.09	120.30
24	3L	76	A	C5-N7-C8	-9.46	99.17	103.90
26	1H	1255	U	N1-C2-O2	-9.46	116.18	122.80
26	14	2374	C	C5-C6-N1	-9.45	116.27	121.00
26	14	2617	C	C6-N1-C2	9.45	124.08	120.30
26	14	2688	U	N1-C2-N3	9.45	120.57	114.90
26	14	2287	A	C2-N3-C4	-9.44	105.88	110.60
26	1H	1616	A	C8-N9-C4	-9.44	102.03	105.80
27	1J	103	U	C5-C6-N1	-9.43	117.99	122.70
26	1H	2449	U	N3-C4-O4	9.42	125.99	119.40
1	1G	862	C	C2-N1-C1'	9.41	129.15	118.80
26	1H	1931	U	C5-C4-O4	9.40	131.54	125.90
26	14	330	A	N1-C6-N6	9.40	124.24	118.60
26	1H	133	C	C6-N1-C2	9.40	124.06	120.30
26	1H	432	A	C5-C6-N6	-9.39	116.19	123.70
26	1H	2392	A	N7-C8-N9	9.39	118.49	113.80
26	14	848	G	N3-C4-C5	-9.38	123.91	128.60
26	14	1786	A	N1-C6-N6	9.38	124.23	118.60
26	1H	860	U	C6-N1-C1'	-9.38	108.06	121.20
1	1G	1499	A	C8-N9-C4	9.37	109.55	105.80
26	1H	783	A	N1-C6-N6	9.37	124.22	118.60
26	1H	1122	G	C8-N9-C4	9.37	110.15	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	692	C	C4-C5-C6	9.37	122.08	117.40
26	14	2374	C	C6-N1-C2	9.37	124.05	120.30
26	14	2873	A	C2-N3-C4	-9.37	105.92	110.60
26	1H	664	C	O5'-P-OP2	-9.36	97.28	105.70
26	1H	678	C	N3-C4-C5	9.36	125.64	121.90
26	1H	71	A	C4-C5-N7	9.35	115.38	110.70
26	1H	1021	A	N7-C8-N9	9.35	118.48	113.80
26	14	1314	C	N1-C2-O2	9.35	124.51	118.90
26	1H	655	A	C8-N9-C4	-9.35	102.06	105.80
26	1H	1332	G	C8-N9-C4	-9.35	102.66	106.40
26	1H	694	U	O5'-P-OP2	-9.34	97.29	105.70
26	1H	265	A	N7-C8-N9	9.34	118.47	113.80
26	14	675	A	N9-C4-C5	-9.34	102.06	105.80
1	13	1504	G	O5'-P-OP1	-9.34	97.30	105.70
26	1H	1614	A	N3-C4-C5	9.34	133.34	126.80
26	1H	2053	G	N1-C6-O6	9.33	125.50	119.90
26	1H	814	C	C2-N3-C4	-9.33	115.23	119.90
1	13	690	G	N7-C8-N9	9.32	117.76	113.10
26	1H	2710	C	C5-C6-N1	-9.32	116.34	121.00
26	1H	71	A	N7-C8-N9	9.32	118.46	113.80
1	1G	108	G	N1-C6-O6	9.32	125.49	119.90
26	1H	835	A	C6-N1-C2	-9.32	113.01	118.60
26	1H	1931	U	C5-C6-N1	-9.31	118.04	122.70
26	1H	930	U	C5-C4-O4	9.31	131.49	125.90
26	1H	782	A	N1-C6-N6	-9.30	113.02	118.60
1	13	690	G	N1-C6-O6	9.30	125.48	119.90
27	16	5	C	N3-C4-C5	9.30	125.62	121.90
26	1H	1885	A	C8-N9-C4	9.29	109.52	105.80
26	14	528	A	N3-C4-N9	-9.30	119.96	127.40
24	3K	71	G	O4'-C1'-N9	9.28	115.63	108.20
26	1H	837	C	O5'-P-OP1	-9.29	97.34	105.70
26	14	1350	C	O5'-P-OP1	-9.29	97.34	105.70
26	1H	676	A	N1-C2-N3	9.28	133.94	129.30
26	1H	788	A	N1-C6-N6	9.27	124.17	118.60
26	14	856	C	C6-N1-C2	-9.26	116.59	120.30
26	1H	74	A	N1-C2-N3	9.26	133.93	129.30
26	1H	667	U	N1-C2-O2	-9.25	116.32	122.80
26	1H	1614	A	C4-C5-N7	9.25	115.33	110.70
26	14	1925	C	N1-C2-O2	-9.25	113.35	118.90
1	13	789	U	N1-C2-N3	9.25	120.45	114.90
26	14	621	A	C5-C6-N1	-9.25	113.08	117.70
26	14	2544	G	C5-C6-O6	-9.24	123.05	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1660	C	N3-C4-C5	9.24	125.60	121.90
26	14	1274	A	N1-C6-N6	9.24	124.14	118.60
26	14	1496	A	C5-N7-C8	-9.23	99.28	103.90
26	1H	2544	G	N1-C6-O6	9.23	125.44	119.90
1	13	903	G	O5'-P-OP2	-9.23	97.39	105.70
26	1H	2053	G	N3-C2-N2	-9.23	113.44	119.90
26	1H	2440	C	C2-N3-C4	9.23	124.51	119.90
26	14	2217	G	N1-C6-O6	9.22	125.43	119.90
26	14	675	A	C8-N9-C4	9.21	109.48	105.80
26	1H	1204	A	C6-C5-N7	-9.21	125.86	132.30
26	1H	2072	G	N9-C4-C5	-9.21	101.72	105.40
22	1K	76	A	N7-C8-N9	9.20	118.40	113.80
26	1H	526	A	N1-C6-N6	-9.20	113.08	118.60
1	13	690	G	C5-N7-C8	-9.20	99.70	104.30
26	1H	2068	U	O5'-P-OP1	-9.20	97.42	105.70
26	1H	2271	G	N3-C4-N9	9.20	131.52	126.00
1	1G	811	C	N1-C2-O2	-9.20	113.38	118.90
26	1H	1700	A	O5'-P-OP2	-9.19	97.43	105.70
26	14	2426	A	N1-C6-N6	9.19	124.11	118.60
26	1H	2544	G	C8-N9-C4	9.18	110.07	106.40
26	1H	655	A	N7-C8-N9	9.18	118.39	113.80
26	1H	695	G	N1-C6-O6	-9.17	114.40	119.90
26	1H	180	G	N9-C4-C5	-9.17	101.73	105.40
26	14	1283	G	N3-C4-C5	-9.17	124.02	128.60
26	14	307	G	O5'-P-OP1	-9.16	97.45	105.70
26	1H	757	U	O5'-P-OP2	-9.16	97.46	105.70
27	16	100	G	C8-N9-C4	9.16	110.06	106.40
26	1H	456	C	C6-N1-C2	9.15	123.96	120.30
26	1H	2712	U	C6-N1-C1'	-9.15	108.39	121.20
26	14	2518	A	C5-N7-C8	-9.15	99.33	103.90
27	16	79	C	C6-N1-C2	-9.14	116.64	120.30
26	1H	456	C	O5'-P-OP2	-9.14	97.47	105.70
26	1H	1338	G	OP1-P-OP2	-9.14	105.89	119.60
26	1H	1189	A	N1-C6-N6	9.13	124.08	118.60
26	1H	107	C	C6-N1-C2	9.13	123.95	120.30
26	1H	575	A	O5'-P-OP1	-9.13	97.48	105.70
27	16	115	G	C5-C6-N1	9.13	116.07	111.50
26	1H	841	A	N1-C6-N6	9.13	124.08	118.60
26	14	530	G	C2-N3-C4	-9.12	107.34	111.90
26	14	265	A	C2-N3-C4	-9.12	106.04	110.60
26	14	676	A	N7-C8-N9	9.12	118.36	113.80
26	1H	658	C	O5'-P-OP2	-9.11	97.50	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1366	A	N1-C6-N6	9.11	124.07	118.60
26	1H	141	A	C4-C5-N7	9.11	115.25	110.70
26	14	2712	U	C2-N3-C4	-9.10	121.54	127.00
23	2L	21	U	N3-C4-C5	9.10	120.06	114.60
26	1H	1021	A	C5-C6-N1	-9.09	113.15	117.70
26	1H	116	C	N1-C2-O2	-9.09	113.45	118.90
26	1H	1789	A	C5-C6-N1	9.09	122.24	117.70
26	1H	1950	G	N3-C4-N9	-9.09	120.55	126.00
26	1H	2330	G	C8-N9-C4	9.09	110.03	106.40
1	13	1502	A	C4-C5-N7	9.09	115.24	110.70
26	14	1564	C	N3-C4-N4	-9.09	111.64	118.00
26	1H	453	C	N1-C2-O2	-9.08	113.45	118.90
26	1H	703	U	C5-C4-O4	9.07	131.34	125.90
26	14	1253	A	N9-C4-C5	-9.07	102.17	105.80
26	14	1342	A	N1-C2-N3	9.07	133.83	129.30
26	1H	560	C	O5'-P-OP2	9.06	121.58	110.70
26	1H	2287	A	C5-C6-N1	-9.06	113.17	117.70
1	1G	449	C	C6-N1-C2	-9.06	116.67	120.30
26	14	945	A	C2-N3-C4	-9.06	106.07	110.60
26	14	1313	U	C6-N1-C2	-9.05	115.57	121.00
26	1H	141	A	N7-C8-N9	9.05	118.33	113.80
26	1H	621	A	C6-N1-C2	9.05	124.03	118.60
26	1H	796	C	O5'-P-OP2	-9.04	97.56	105.70
26	1H	371	A	N1-C6-N6	9.04	124.02	118.60
26	1H	593	G	N1-C2-N3	9.04	129.32	123.90
26	1H	1604	C	C6-N1-C2	-9.04	116.68	120.30
26	1H	1681	G	C4-C5-N7	9.04	114.41	110.80
26	14	1812	A	O5'-P-OP2	-9.03	97.57	105.70
26	1H	2392	A	N3-C4-C5	9.03	133.12	126.80
26	14	140	A	C6-C5-N7	-9.02	125.99	132.30
26	14	1605	C	O5'-P-OP1	-9.02	97.58	105.70
1	13	814	A	O5'-P-OP2	9.02	121.52	110.70
26	1H	1817	G	C8-N9-C4	9.01	110.00	106.40
1	13	865	A	N1-C6-N6	9.01	124.01	118.60
26	14	2392	A	O5'-P-OP1	-9.01	97.59	105.70
26	1H	1304	C	N3-C4-C5	9.01	125.50	121.90
26	1H	1931	U	C4-C5-C6	9.00	125.10	119.70
26	14	2447	G	O5'-P-OP1	-8.99	97.60	105.70
26	1H	1437	C	C6-N1-C2	-8.99	116.70	120.30
26	1H	2430	A	C5-C6-N1	-8.98	113.21	117.70
26	1H	1022	G	N9-C4-C5	8.97	108.99	105.40
55	Q8	25	MET	N-CA-C	8.97	135.21	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	681	G	C2-N3-C4	-8.96	107.42	111.90
26	14	2249	U	N3-C2-O2	-8.96	115.93	122.20
26	1H	2375	G	N7-C8-N9	-8.96	108.62	113.10
26	1H	179	G	N3-C2-N2	-8.96	113.63	119.90
26	1H	1528	A	N7-C8-N9	8.96	118.28	113.80
26	1H	2446	G	N1-C6-O6	8.96	125.28	119.90
1	13	583	A	C8-N9-C4	8.95	109.38	105.80
1	13	990	C	C6-N1-C2	-8.94	116.72	120.30
26	1H	2392	A	C4-C5-N7	8.94	115.17	110.70
1	13	690	G	O4'-C1'-N9	8.94	115.35	108.20
26	1H	2712	U	O4'-C1'-N1	8.94	115.35	108.20
26	14	568	U	C4-C5-C6	-8.94	114.34	119.70
26	1H	2330	G	N9-C4-C5	-8.93	101.83	105.40
26	14	1807	G	C8-N9-C4	8.93	109.97	106.40
26	1H	1899	G	C5-C6-N1	-8.93	107.04	111.50
26	1H	537	C	O5'-P-OP1	8.93	121.41	110.70
26	1H	1698	A	C5-C6-N1	-8.92	113.24	117.70
26	14	2713	A	C5-N7-C8	-8.92	99.44	103.90
26	1H	120	U	C5-C4-O4	8.91	131.25	125.90
26	1H	193	U	C4-C5-C6	8.91	125.05	119.70
26	14	2776	A	C8-N9-C4	-8.91	102.23	105.80
26	14	2851	A	O5'-P-OP1	-8.91	97.68	105.70
26	1H	1266	G	N9-C4-C5	-8.91	101.84	105.40
26	14	2518	A	O4'-C1'-N9	-8.91	101.07	108.20
26	14	2357	U	O5'-P-OP2	-8.91	97.69	105.70
26	14	737	C	N3-C2-O2	8.90	128.13	121.90
26	14	1661	G	C5-C6-O6	-8.90	123.26	128.60
26	14	1970	A	O5'-P-OP2	-8.90	97.69	105.70
26	14	2873	A	N1-C6-N6	8.90	123.94	118.60
26	1H	737	C	C5-C6-N1	-8.89	116.55	121.00
26	14	1647	G	O5'-P-OP1	-8.89	97.70	105.70
26	1H	789	A	O5'-P-OP1	-8.89	97.70	105.70
26	14	2005	A	C8-N9-C4	8.88	109.35	105.80
26	1H	1678	G	N3-C4-N9	-8.88	120.67	126.00
26	1H	2591	C	O5'-P-OP1	-8.88	97.71	105.70
26	1H	686	G	N9-C4-C5	-8.87	101.85	105.40
26	1H	123	G	C5-C6-N1	8.87	115.93	111.50
26	1H	673	C	O5'-P-OP1	8.87	121.34	110.70
1	13	896	C	C6-N1-C2	8.86	123.84	120.30
22	1K	76	A	C8-N9-C4	-8.86	102.25	105.80
26	14	2087	G	C8-N9-C4	8.86	109.94	106.40
26	14	2688	U	N3-C2-O2	-8.85	116.00	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1616	A	C4-C5-N7	8.84	115.12	110.70
26	1H	1394	U	O5'-P-OP1	-8.84	97.75	105.70
26	1H	140	A	C5-C6-N6	-8.84	116.63	123.70
26	1H	2417	C	O5'-P-OP2	-8.84	97.75	105.70
26	1H	2433	A	N1-C2-N3	8.84	133.72	129.30
26	14	1408	C	N1-C2-O2	-8.84	113.60	118.90
26	1H	632	A	O5'-P-OP2	8.83	121.30	110.70
26	1H	1404	C	C6-N1-C2	8.83	123.83	120.30
26	1H	783	A	C4-C5-N7	8.82	115.11	110.70
26	1H	194	G	C5-C6-O6	-8.81	123.31	128.60
26	1H	681	G	N1-C2-N2	-8.81	108.27	116.20
26	14	949	C	N1-C2-O2	-8.81	113.61	118.90
26	14	2690	C	O5'-P-OP2	-8.81	97.77	105.70
26	1H	695	G	C5-C6-O6	8.80	133.88	128.60
26	1H	1204	A	N1-C2-N3	8.80	133.70	129.30
26	1H	203	C	O5'-P-OP2	8.80	121.26	110.70
26	1H	2392	A	C5-C6-N1	-8.80	113.30	117.70
26	14	1204	A	C5-C6-N1	-8.80	113.30	117.70
26	1H	120	U	N3-C2-O2	-8.80	116.04	122.20
26	14	2401	U	C5-C6-N1	8.79	127.10	122.70
26	1H	1255	U	N3-C4-O4	8.79	125.55	119.40
26	14	71	A	C4-C5-N7	8.79	115.10	110.70
26	1H	1786	A	C4-C5-C6	8.79	121.39	117.00
26	14	779	U	N3-C4-O4	-8.79	113.25	119.40
26	1H	226	G	O4'-C1'-N9	8.78	115.22	108.20
26	1H	1201	C	N3-C2-O2	8.77	128.04	121.90
36	35	62	LEU	N-CA-C	8.77	134.69	111.00
26	1H	1613	G	N3-C2-N2	8.77	126.04	119.90
1	13	656	C	C6-N1-C2	-8.77	116.79	120.30
26	14	2591	C	N1-C2-O2	-8.77	113.64	118.90
26	14	974(A)	C	C6-N1-C2	-8.77	116.79	120.30
26	1H	2446	G	C5-C6-O6	-8.76	123.34	128.60
26	14	208	C	N3-C4-C5	8.76	125.41	121.90
26	1H	528	A	C5-C6-N1	-8.76	113.32	117.70
26	14	2339	G	O5'-P-OP2	-8.76	97.82	105.70
26	14	2610	C	O5'-P-OP1	-8.76	97.82	105.70
26	14	871	U	O5'-P-OP1	-8.75	97.83	105.70
26	1H	1604	C	N3-C4-C5	-8.75	118.40	121.90
26	14	2035	G	O4'-C1'-N9	8.74	115.19	108.20
26	1H	1142(A)	A	C2-N3-C4	-8.74	106.23	110.60
26	1H	658	C	N3-C2-O2	-8.74	115.78	121.90
26	1H	1596	A	N1-C6-N6	-8.74	113.36	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	568	U	N3-C4-C5	8.74	119.84	114.60
26	14	2577	A	N1-C6-N6	8.74	123.84	118.60
26	14	2092	U	C5-C4-O4	8.74	131.14	125.90
26	14	1678	G	C5-N7-C8	-8.73	99.93	104.30
26	1H	1314	C	C2-N1-C1'	8.73	128.41	118.80
26	14	783	A	C8-N9-C4	-8.73	102.31	105.80
1	13	542	G	O5'-P-OP1	-8.73	97.84	105.70
26	1H	639	U	N3-C2-O2	-8.72	116.09	122.20
26	1H	957	A	N1-C6-N6	8.72	123.83	118.60
26	1H	2713	A	C4-C5-N7	8.72	115.06	110.70
26	14	2388	A	O4'-C1'-N9	8.72	115.18	108.20
1	1G	904	C	O5'-P-OP1	-8.72	97.85	105.70
26	1H	1376	C	C6-N1-C2	-8.71	116.81	120.30
26	1H	596	G	C5-C6-O6	-8.71	123.38	128.60
26	14	664	C	O5'-P-OP2	-8.71	97.86	105.70
26	14	1790	C	C6-N1-C2	8.70	123.78	120.30
26	1H	2578	G	C8-N9-C4	8.69	109.88	106.40
26	1H	763	G	N3-C4-C5	-8.69	124.25	128.60
26	1H	2518	A	N7-C8-N9	8.69	118.15	113.80
26	1H	796	C	N3-C4-C5	8.69	125.38	121.90
26	1H	1304	C	C6-N1-C2	8.69	123.78	120.30
26	1H	2390	U	O5'-P-OP1	-8.69	97.88	105.70
26	1H	738	G	C5-C6-O6	-8.69	123.39	128.60
26	1H	2609	U	C4-C5-C6	8.69	124.91	119.70
26	1H	204	A	C6-N1-C2	-8.68	113.39	118.60
26	1H	1612	C	N1-C2-O2	-8.68	113.69	118.90
26	1H	1616	A	N1-C6-N6	8.68	123.81	118.60
26	14	676	A	N1-C2-N3	8.68	133.64	129.30
26	1H	2451	A	N1-C6-N6	-8.68	113.39	118.60
26	1H	2755	C	C6-N1-C2	-8.68	116.83	120.30
1	13	963	G	N1-C2-N2	-8.67	108.40	116.20
26	1H	1574	C	C5-C6-N1	-8.67	116.67	121.00
26	1H	1798	U	O5'-P-OP2	-8.67	97.90	105.70
26	1H	2599	G	C6-C5-N7	8.67	135.60	130.40
26	14	1121	C	C6-N1-C2	8.67	123.77	120.30
26	1H	1779	U	C6-N1-C2	8.67	126.20	121.00
26	1H	1636	C	N1-C2-O2	-8.66	113.70	118.90
26	1H	1698	A	C4-C5-N7	8.66	115.03	110.70
26	1H	2627	G	C5-C6-O6	-8.66	123.41	128.60
26	14	2518	A	C6-C5-N7	-8.65	126.24	132.30
26	1H	107	C	C5-C6-N1	-8.65	116.67	121.00
26	1H	1257	C	C2-N3-C4	-8.65	115.57	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	679	C	N1-C2-O2	-8.65	113.71	118.90
26	1H	1162	G	C8-N9-C4	-8.65	102.94	106.40
26	1H	1350	C	C6-N1-C2	8.65	123.76	120.30
26	14	2873	A	C4-C5-C6	8.65	121.32	117.00
26	1H	1969	A	N1-C6-N6	-8.65	113.41	118.60
26	1H	1698	A	C6-C5-N7	-8.64	126.25	132.30
26	1H	1799	G	N3-C4-N9	8.63	131.18	126.00
26	1H	1425	G	C5-C6-N1	8.63	115.82	111.50
26	1H	2330	G	N1-C6-O6	8.63	125.08	119.90
26	1H	2599	G	C4-C5-N7	-8.63	107.35	110.80
26	1H	1678	G	C6-C5-N7	-8.63	125.22	130.40
26	1H	1767	C	O5'-P-OP1	-8.63	97.94	105.70
26	1H	512	G	O4'-C1'-N9	8.62	115.10	108.20
26	1H	2544	G	N9-C4-C5	-8.62	101.95	105.40
1	13	760	G	N1-C6-O6	8.62	125.07	119.90
26	1H	1950	G	N3-C4-C5	8.61	132.91	128.60
26	14	935	C	C6-N1-C2	8.62	123.75	120.30
26	1H	470	A	N1-C2-N3	8.61	133.60	129.30
26	1H	658	C	N1-C2-O2	8.61	124.06	118.90
1	1G	1517	G	O5'-P-OP2	-8.61	97.96	105.70
26	14	2272	U	O5'-P-OP2	-8.60	97.96	105.70
1	13	1322	C	C2-N1-C1'	8.60	128.26	118.80
26	1H	481	G	O5'-P-OP2	-8.60	97.96	105.70
26	1H	2552	U	N1-C2-O2	-8.60	116.78	122.80
26	1H	1376	C	O5'-P-OP1	-8.59	97.97	105.70
24	3L	76	A	N1-C6-N6	8.59	123.75	118.60
26	14	1366	A	N1-C6-N6	8.59	123.75	118.60
26	14	2447	G	P-O3'-C3'	8.59	130.00	119.70
26	1H	2380	C	C2-N3-C4	-8.58	115.61	119.90
26	14	1840	G	N1-C6-O6	8.58	125.05	119.90
26	1H	2779	U	C5-C6-N1	-8.58	118.41	122.70
26	14	971	C	C6-N1-C2	-8.57	116.87	120.30
26	1H	736	C	O5'-P-OP2	8.57	120.99	110.70
26	14	1279	G	N1-C6-O6	-8.57	114.76	119.90
26	14	2070	G	C2-N3-C4	-8.57	107.61	111.90
26	1H	528	A	N3-C4-C5	8.57	132.80	126.80
26	1H	1681	G	N3-C4-C5	8.57	132.88	128.60
26	14	2332	U	N1-C2-O2	8.57	128.80	122.80
26	1H	245	G	C4-C5-N7	8.56	114.22	110.80
1	13	1128	C	C6-N1-C2	-8.56	116.88	120.30
26	1H	1914	C	N1-C2-O2	8.56	124.04	118.90
26	1H	621	A	N3-C4-N9	-8.56	120.55	127.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	945	A	C8-N9-C1'	-8.56	112.29	127.70
26	1H	691	C	C5-C6-N1	-8.56	116.72	121.00
26	1H	409	C	C6-N1-C2	8.55	123.72	120.30
26	1H	1201	C	N1-C2-O2	-8.55	113.77	118.90
1	13	1422	G	O5'-P-OP2	-8.54	98.01	105.70
26	1H	245	G	N9-C4-C5	-8.54	101.98	105.40
26	1H	1786	A	C8-N9-C4	-8.54	102.38	105.80
27	16	115	G	C5-C6-O6	-8.54	123.47	128.60
26	1H	1827	C	C2-N3-C4	-8.54	115.63	119.90
26	1H	1021	A	N1-C6-N6	8.54	123.72	118.60
26	1H	1621	U	C5-C4-O4	-8.54	120.78	125.90
26	14	1902	C	N3-C4-C5	8.54	125.32	121.90
26	14	615	G	O4'-C1'-N9	8.54	115.03	108.20
26	1H	1332	G	N3-C2-N2	-8.54	113.92	119.90
26	1H	745	G	N1-C6-O6	8.53	125.02	119.90
26	1H	945	A	C2-N3-C4	-8.53	106.33	110.60
26	1H	2275	C	C6-N1-C2	-8.53	116.89	120.30
26	1H	2688	U	N1-C2-N3	8.53	120.02	114.90
26	14	2544	G	N1-C6-O6	8.53	125.02	119.90
26	1H	1338	G	C2-N3-C4	8.53	116.17	111.90
26	14	922	U	O5'-P-OP1	-8.53	98.02	105.70
26	1H	1678	G	N7-C8-N9	8.53	117.36	113.10
26	1H	1431	U	C5-C6-N1	8.52	126.96	122.70
26	14	1786	A	N9-C1'-C2'	8.52	125.08	114.00
26	1H	471	A	C2-N3-C4	-8.52	106.34	110.60
1	13	1502	A	C6-C5-N7	-8.52	126.34	132.30
26	1H	1698	A	N1-C2-N3	8.51	133.56	129.30
26	1H	1613	G	N1-C2-N2	-8.51	108.54	116.20
26	1H	821	A	OP1-P-OP2	8.51	132.36	119.60
26	1H	1899	G	C8-N9-C4	-8.51	103.00	106.40
26	1H	321	G	C6-C5-N7	-8.51	125.30	130.40
23	2K	17	C	N1-C2-O2	8.51	124.00	118.90
26	1H	2440	C	C5-C4-N4	8.51	126.16	120.20
26	1H	613	U	N3-C2-O2	-8.50	116.25	122.20
26	1H	948	G	C5-C6-O6	-8.50	123.50	128.60
26	1H	103	A	N9-C4-C5	-8.50	102.40	105.80
26	1H	122	G	C2-N3-C4	-8.50	107.65	111.90
1	13	1502	A	N1-C6-N6	8.49	123.70	118.60
26	1H	1528	A	O4'-C1'-N9	8.49	114.99	108.20
27	16	7	G	C4-C5-N7	8.49	114.20	110.80
26	14	2435	A	C8-N9-C4	-8.49	102.40	105.80
26	14	768	G	O5'-P-OP2	-8.49	98.06	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	827	U	O5'-P-OP2	-8.49	98.06	105.70
26	1H	190	A	C8-N9-C4	8.49	109.19	105.80
26	1H	2778	A	O5'-P-OP2	-8.48	98.06	105.70
26	1H	25	U	C5-C4-O4	-8.48	120.81	125.90
26	1H	679	C	C6-N1-C2	8.48	123.69	120.30
26	1H	1574	C	C2-N3-C4	-8.48	115.66	119.90
27	16	7	G	C5-N7-C8	-8.48	100.06	104.30
26	1H	2385	C	C2-N3-C4	-8.47	115.67	119.90
26	14	1349	A	N1-C6-N6	8.46	123.68	118.60
26	1H	71	A	N1-C2-N3	8.45	133.53	129.30
26	14	312	G	O5'-P-OP1	-8.45	98.10	105.70
26	1H	1202	C	N3-C4-C5	-8.44	118.52	121.90
27	16	112	G	C8-N9-C4	8.44	109.78	106.40
26	1H	2072	G	N3-C2-N2	8.44	125.81	119.90
26	1H	436	C	C6-N1-C2	8.44	123.67	120.30
26	14	1673	U	O5'-P-OP1	-8.44	98.11	105.70
26	14	528	A	C4-N9-C1'	-8.44	111.11	126.30
26	14	1571	A	C6-N1-C2	-8.44	113.54	118.60
26	1H	444	C	O5'-P-OP1	8.43	120.82	110.70
26	1H	507	A	O5'-P-OP2	8.43	120.82	110.70
26	1H	1649	G	N3-C4-C5	-8.43	124.39	128.60
26	14	2036	C	O5'-P-OP2	-8.43	98.11	105.70
26	1H	1698	A	N3-C4-C5	8.42	132.69	126.80
26	14	1253	A	C6-C5-N7	-8.42	126.40	132.30
1	13	963	G	N9-C4-C5	-8.42	102.03	105.40
26	1H	209	C	O5'-P-OP2	-8.42	98.12	105.70
26	1H	1665	A	O5'-P-OP1	-8.41	98.13	105.70
26	1H	1707	G	C5-C6-O6	-8.41	123.55	128.60
26	1H	967	C	C5-C6-N1	-8.40	116.80	121.00
26	1H	201	C	C5-C6-N1	-8.40	116.80	121.00
26	1H	1312	U	O5'-P-OP1	-8.40	98.14	105.70
26	1H	774	A	C8-N9-C1'	8.40	142.82	127.70
26	14	2570	G	N1-C6-O6	8.40	124.94	119.90
26	1H	1826	G	OP1-P-O3'	8.40	123.68	105.20
26	1H	686	G	OP1-P-OP2	8.39	132.19	119.60
26	14	530	G	N3-C4-C5	8.39	132.80	128.60
26	1H	664	C	C6-N1-C2	8.39	123.66	120.30
23	2L	21	U	C2-N3-C4	-8.39	121.97	127.00
26	1H	917	A	N1-C6-N6	8.39	123.63	118.60
26	1H	2429	G	C8-N9-C4	-8.39	103.05	106.40
26	1H	330	A	C5-N7-C8	-8.38	99.71	103.90
26	1H	913	U	O5'-P-OP2	-8.38	98.16	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2618	G	C8-N9-C4	-8.38	103.05	106.40
26	14	558	G	C8-N9-C4	8.38	109.75	106.40
26	1H	2310	A	C8-N9-C4	-8.38	102.45	105.80
26	14	1899	G	N1-C2-N2	-8.37	108.66	116.20
26	1H	85	G	O5'-P-OP2	-8.37	98.16	105.70
26	1H	1818	U	O5'-P-OP2	-8.37	98.17	105.70
26	14	2688	U	C5-C6-N1	-8.36	118.52	122.70
26	1H	1639	U	N3-C2-O2	-8.36	116.35	122.20
26	1H	245	G	N3-C4-N9	8.36	131.01	126.00
26	1H	2430	A	N3-C4-N9	-8.36	120.72	127.40
26	1H	265	A	C6-C5-N7	-8.35	126.45	132.30
26	1H	918	A	O5'-P-OP1	-8.35	98.18	105.70
1	13	690	G	C4-N9-C1'	8.35	137.35	126.50
26	1H	189	G	N1-C6-O6	8.34	124.91	119.90
26	1H	517	C	C5-C4-N4	-8.34	114.36	120.20
1	1G	345	C	C2-N1-C1'	8.34	127.98	118.80
26	1H	2698	U	O5'-P-OP2	-8.34	98.19	105.70
26	1H	674	G	C8-N9-C4	8.33	109.73	106.40
26	1H	2627	G	N9-C4-C5	-8.33	102.07	105.40
1	13	811	C	C2-N3-C4	-8.33	115.74	119.90
1	1G	691	G	N1-C6-O6	8.33	124.90	119.90
1	13	656	C	C5-C6-N1	8.32	125.16	121.00
26	1H	121	G	N3-C4-N9	8.32	130.99	126.00
26	14	1608	A	N1-C6-N6	-8.32	113.61	118.60
26	14	2439	A	N1-C6-N6	8.32	123.59	118.60
26	1H	948	G	N1-C6-O6	8.32	124.89	119.90
26	1H	2287	A	N1-C6-N6	8.32	123.59	118.60
26	1H	1297	C	OP1-P-O3'	8.31	123.49	105.20
26	1H	1422	G	O5'-P-OP1	-8.31	98.22	105.70
26	1H	455	C	C6-N1-C2	8.31	123.62	120.30
26	1H	1501	C	O5'-P-OP2	8.31	120.67	110.70
24	3K	71	G	C4-N9-C1'	-8.31	115.70	126.50
26	1H	814	C	N3-C4-C5	8.31	125.22	121.90
26	1H	1502	C	C6-N1-C2	-8.30	116.98	120.30
26	1H	1811	G	N1-C6-O6	-8.30	114.92	119.90
26	1H	688	U	N1-C2-N3	8.30	119.88	114.90
26	1H	698	C	C5-C4-N4	-8.30	114.39	120.20
26	1H	1313	U	C5-C6-N1	8.29	126.85	122.70
1	13	122	G	N1-C6-O6	8.29	124.87	119.90
26	1H	238	C	C5-C6-N1	-8.29	116.86	121.00
26	1H	2595	G	C5-N7-C8	-8.29	100.16	104.30
26	1H	865	C	O5'-P-OP2	8.28	120.64	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2318	G	N7-C8-N9	8.28	117.24	113.10
26	1H	2713	A	C6-C5-N7	-8.28	126.50	132.30
26	14	1853	A	O5'-P-OP1	-8.28	98.25	105.70
26	14	1939	U	O5'-P-OP1	-8.28	98.25	105.70
26	1H	200	U	O5'-P-OP1	-8.28	98.25	105.70
26	1H	845	G	N3-C2-N2	8.27	125.69	119.90
26	14	913	U	O5'-P-OP2	-8.27	98.26	105.70
26	1H	470	A	N1-C6-N6	8.27	123.56	118.60
1	13	692	U	N3-C4-O4	8.26	125.18	119.40
26	1H	470	A	C2-N3-C4	-8.26	106.47	110.60
26	1H	973	A	C2-N3-C4	-8.26	106.47	110.60
26	1H	2318	G	C5-N7-C8	-8.26	100.17	104.30
1	1G	1346	A	P-O3'-C3'	8.26	129.61	119.70
26	1H	1272	A	C4-C5-C6	-8.25	112.87	117.00
26	14	205	G	C8-N9-C4	8.25	109.70	106.40
1	13	781	A	N9-C4-C5	-8.25	102.50	105.80
26	1H	265	A	C5-C6-N1	-8.25	113.58	117.70
26	1H	1566	A	O5'-P-OP1	8.25	120.60	110.70
26	14	252	G	N1-C6-O6	-8.25	114.95	119.90
26	14	2331	G	C5-C6-O6	-8.25	123.65	128.60
26	14	2392	A	N7-C8-N9	8.25	117.92	113.80
1	13	1526	G	N1-C6-O6	8.24	124.85	119.90
26	1H	196	A	O4'-C1'-N9	8.24	114.79	108.20
26	1H	448	U	C4-C5-C6	8.24	124.64	119.70
26	1H	1900	A	O5'-P-OP1	8.24	120.58	110.70
26	1H	1950	G	O4'-C1'-N9	8.24	114.79	108.20
26	1H	2589	A	OP2-P-O3'	8.23	123.31	105.20
26	1H	418	G	O5'-P-OP1	-8.23	98.30	105.70
26	1H	2568	C	O5'-P-OP1	-8.23	98.30	105.70
1	13	1488	G	N1-C6-O6	8.22	124.83	119.90
26	1H	1899	G	C6-C5-N7	8.22	135.33	130.40
26	1H	2072	G	N1-C2-N3	-8.22	118.97	123.90
26	1H	2827	C	N3-C4-N4	8.22	123.76	118.00
1	13	781	A	C4-C5-N7	8.22	114.81	110.70
26	1H	517	C	N3-C4-N4	8.22	123.75	118.00
26	14	774	A	C5-C6-N1	-8.22	113.59	117.70
26	14	1975	G	N1-C6-O6	8.21	124.83	119.90
26	1H	1379	A	C4-C5-N7	8.21	114.81	110.70
26	1H	2688	U	C5-C4-O4	8.21	130.83	125.90
26	1H	262	A	N1-C6-N6	8.21	123.53	118.60
1	13	789	U	C4-C5-C6	8.21	124.62	119.70
26	1H	575	A	O5'-P-OP2	8.21	120.55	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	681	G	C8-N9-C4	8.21	109.68	106.40
26	14	1997	G	N3-C4-C5	-8.21	124.50	128.60
26	14	2056	G	N3-C2-N2	-8.21	114.16	119.90
26	1H	1379	A	N7-C8-N9	8.20	117.90	113.80
26	14	751	A	OP1-P-OP2	-8.20	107.30	119.60
26	14	189	G	N1-C6-O6	8.20	124.82	119.90
26	14	2346	A	N7-C8-N9	8.20	117.90	113.80
26	1H	1639	U	O5'-P-OP1	8.19	120.53	110.70
26	14	2272	U	O5'-P-OP1	8.20	120.53	110.70
26	1H	199	A	N1-C6-N6	-8.19	113.69	118.60
1	1G	690	G	N3-C4-N9	-8.19	121.09	126.00
26	14	1441	G	C8-N9-C4	8.19	109.68	106.40
1	13	1489	G	C8-N9-C4	8.19	109.67	106.40
1	13	1517	G	C5-C6-O6	-8.18	123.69	128.60
26	14	130	C	C6-N1-C2	8.18	123.57	120.30
26	14	148	C	C6-N1-C2	8.17	123.57	120.30
26	1H	796	C	C2-N3-C4	-8.17	115.81	119.90
26	1H	1128	A	O5'-P-OP1	-8.17	98.35	105.70
26	1H	1922	G	O5'-P-OP2	-8.17	98.35	105.70
26	1H	2232	U	C5-C4-O4	8.17	130.80	125.90
26	1H	1327	C	N1-C2-O2	-8.17	114.00	118.90
26	1H	2374	C	N3-C4-C5	8.16	125.16	121.90
26	1H	62	C	C6-N1-C2	8.16	123.56	120.30
26	1H	860	U	O5'-P-OP1	8.16	120.49	110.70
26	14	676	A	N1-C6-N6	8.16	123.49	118.60
26	14	1283	G	O5'-P-OP2	-8.16	98.36	105.70
1	13	966	G	N1-C6-O6	8.15	124.79	119.90
26	1H	197	A	N1-C2-N3	8.15	133.38	129.30
26	1H	470	A	C5-C6-N6	-8.15	117.18	123.70
26	1H	683	C	C6-N1-C2	8.15	123.56	120.30
26	1H	2639	A	C5-C6-N1	-8.15	113.62	117.70
26	14	798	G	N1-C6-O6	8.15	124.79	119.90
26	14	2688	U	C4-C5-C6	8.15	124.59	119.70
26	14	565	C	C5-C6-N1	-8.15	116.93	121.00
26	14	1950	G	N7-C8-N9	8.15	117.17	113.10
26	14	117	G	O5'-P-OP1	8.14	120.47	110.70
26	1H	2584	U	N3-C4-O4	-8.14	113.70	119.40
26	14	2068	U	O5'-P-OP1	-8.14	98.37	105.70
26	1H	245	G	C8-N9-C1'	-8.14	116.42	127.00
26	14	2375	G	C8-N9-C4	8.14	109.66	106.40
26	14	510	C	C6-N1-C2	-8.14	117.04	120.30
26	1H	208	C	C5-C6-N1	-8.13	116.93	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1210	A	N1-C6-N6	8.13	123.48	118.60
26	1H	141	A	N1-C6-N6	8.13	123.47	118.60
26	1H	2447	G	O4'-C1'-N9	8.12	114.70	108.20
26	1H	2573	C	N3-C4-C5	-8.12	118.65	121.90
26	14	1786	A	C5-C6-N1	-8.13	113.64	117.70
26	14	252	G	N7-C8-N9	-8.12	109.04	113.10
26	1H	699	A	N1-C6-N6	8.12	123.47	118.60
26	1H	1899	G	N1-C2-N3	8.12	128.77	123.90
26	14	1339	G	O5'-P-OP2	8.12	120.45	110.70
26	14	2278	A	N9-C4-C5	8.12	109.05	105.80
26	1H	198	C	N3-C4-C5	8.12	125.15	121.90
26	1H	992	C	C6-N1-C2	-8.12	117.05	120.30
26	1H	1969	A	C5-C6-N6	8.12	130.19	123.70
26	14	2590	A	C8-N9-C4	8.11	109.05	105.80
26	1H	950	G	N1-C2-N2	-8.11	108.90	116.20
26	1H	1950	G	C5-C6-N1	-8.11	107.45	111.50
26	14	1289	C	O5'-P-OP1	-8.11	98.40	105.70
26	14	2392	A	C8-N9-C4	-8.11	102.56	105.80
26	1H	265	A	N1-C6-N6	8.10	123.46	118.60
26	1H	1307	A	O5'-P-OP1	-8.10	98.41	105.70
26	1H	1781	C	C6-N1-C2	8.10	123.54	120.30
26	14	2360	A	C2-N3-C4	-8.10	106.55	110.60
26	1H	606	U	O5'-P-OP2	-8.10	98.42	105.70
26	1H	2058	A	C8-N9-C4	-8.09	102.56	105.80
26	1H	676	A	C4-C5-N7	8.09	114.75	110.70
26	1H	1210	A	C2-N3-C4	-8.09	106.55	110.60
26	1H	1573	G	OP1-P-O3'	-8.09	87.40	105.20
26	14	2689	U	C5-C6-N1	-8.09	118.66	122.70
26	1H	736	C	C2-N3-C4	-8.09	115.86	119.90
26	1H	391	G	C2-N3-C4	-8.08	107.86	111.90
23	2K	21	U	C4-C5-C6	-8.08	114.85	119.70
26	14	1299	G	O5'-P-OP1	-8.08	98.43	105.70
1	13	764	C	C6-N1-C2	8.07	123.53	120.30
23	2K	61	U	O5'-P-OP2	-8.07	98.43	105.70
26	1H	135	G	C8-N9-C4	8.07	109.63	106.40
26	1H	1195	G	N1-C6-O6	8.07	124.74	119.90
26	14	945	A	C4-C5-C6	8.07	121.03	117.00
26	14	2712	U	C5-C4-O4	8.07	130.74	125.90
26	1H	128	C	C6-N1-C2	8.06	123.53	120.30
26	1H	2584	U	N1-C2-N3	8.06	119.74	114.90
26	14	1619	G	C5-C6-O6	-8.06	123.76	128.60
26	14	2778	A	O5'-P-OP2	-8.06	98.44	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	1482	G	O5'-P-OP2	-8.06	98.44	105.70
26	1H	609	A	N1-C6-N6	8.06	123.44	118.60
26	14	752	A	N1-C2-N3	8.06	133.33	129.30
1	13	1530	G	N3-C4-C5	8.06	132.63	128.60
26	1H	814	C	C6-N1-C2	8.06	123.52	120.30
26	1H	190	A	N9-C4-C5	-8.06	102.58	105.80
26	14	1332	G	C8-N9-C1'	8.05	137.47	127.00
26	1H	1363	C	C2-N3-C4	-8.05	115.87	119.90
26	14	1313	U	N1-C2-O2	-8.05	117.16	122.80
1	1G	413	G	C4-C5-N7	-8.05	107.58	110.80
26	1H	567	A	O5'-P-OP1	-8.05	98.46	105.70
26	1H	1496	A	C4-C5-N7	8.05	114.72	110.70
26	1H	2476	A	C8-N9-C4	-8.05	102.58	105.80
26	1H	507	A	N1-C6-N6	8.04	123.43	118.60
26	1H	2822	G	N1-C6-O6	8.04	124.73	119.90
26	14	1332	G	N7-C8-N9	8.04	117.12	113.10
26	14	2005	A	N7-C8-N9	-8.04	109.78	113.80
26	14	2078	C	C6-N1-C2	-8.04	117.08	120.30
1	13	1498	U	C2-N1-C1'	8.04	127.34	117.70
26	1H	1198	U	C5-C6-N1	-8.04	118.68	122.70
26	1H	2576	G	C8-N9-C4	8.04	109.61	106.40
26	1H	2607	G	N1-C2-N2	-8.03	108.97	116.20
26	14	1632	A	N1-C6-N6	8.03	123.42	118.60
26	14	2073	C	N1-C2-O2	-8.03	114.08	118.90
26	14	2592	G	O5'-P-OP2	-8.03	98.47	105.70
26	1H	858	U	O5'-P-OP2	-8.03	98.48	105.70
26	1H	660	G	C5-N7-C8	-8.02	100.29	104.30
26	1H	747	U	OP1-P-OP2	8.02	131.64	119.60
26	1H	2094	G	O5'-P-OP2	-8.02	98.48	105.70
26	1H	2427	C	O5'-P-OP2	8.02	120.33	110.70
26	14	1029	A	N9-C4-C5	-8.02	102.59	105.80
26	14	2323	G	C8-N9-C4	8.02	109.61	106.40
1	1G	1260	C	C6-N1-C2	-8.02	117.09	120.30
54	L5	34	ARG	NE-CZ-NH1	-8.02	116.29	120.30
1	13	813	U	N3-C4-O4	-8.02	113.79	119.40
26	1H	1632	A	C4-C5-N7	8.02	114.71	110.70
26	1H	1241	A	N1-C6-N6	8.02	123.41	118.60
26	14	71	A	C2-N3-C4	-8.02	106.59	110.60
26	14	1396	U	N3-C2-O2	-8.02	116.59	122.20
26	14	1616	A	N7-C8-N9	8.02	117.81	113.80
26	14	1145	C	C6-N1-C2	-8.01	117.09	120.30
26	14	196	A	O4'-C1'-N9	8.01	114.61	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1600	C	O5'-P-OP2	-8.01	98.49	105.70
26	1H	793	A	C5-C6-N6	-8.01	117.29	123.70
26	1H	1129	A	O5'-P-OP2	-8.00	98.50	105.70
26	1H	2272	U	O5'-P-OP1	8.00	120.30	110.70
26	1H	1899	G	C5-C6-O6	8.00	133.40	128.60
26	1H	2346	A	C4-N9-C1'	8.00	140.70	126.30
26	14	1187	G	O5'-P-OP2	-8.00	98.50	105.70
26	1H	1298	C	C5-C6-N1	8.00	125.00	121.00
26	1H	85	G	O5'-P-OP1	7.99	120.29	110.70
26	1H	447	A	O5'-P-OP1	-7.99	98.50	105.70
28	11	52	ARG	NE-CZ-NH1	-7.99	116.31	120.30
24	3L	76	A	C6-C5-N7	-7.99	126.71	132.30
1	13	792	A	C5-C6-N6	-7.98	117.31	123.70
1	13	1260	C	C6-N1-C2	-7.98	117.11	120.30
22	1K	74	C	C2-N1-C1'	7.98	127.58	118.80
26	14	569	U	C5-C6-N1	-7.98	118.71	122.70
26	1H	2689	U	N1-C2-N3	7.98	119.69	114.90
26	14	698	C	OP1-P-OP2	7.97	131.56	119.60
26	14	2237	G	N1-C2-N2	-7.97	109.03	116.20
26	1H	908	C	O5'-P-OP2	-7.97	98.53	105.70
26	1H	1694	C	P-O3'-C3'	7.97	129.26	119.70
26	14	1314	C	C6-N1-C1'	-7.97	111.24	120.80
26	1H	2030	A	N1-C6-N6	7.97	123.38	118.60
26	14	1376	C	O5'-P-OP1	-7.96	98.53	105.70
26	1H	691	C	N3-C4-C5	7.96	125.08	121.90
26	1H	635	C	O5'-P-OP2	-7.96	98.53	105.70
26	14	1189	A	OP1-P-OP2	-7.96	107.66	119.60
26	1H	736	C	N1-C2-O2	-7.96	114.12	118.90
1	13	966	G	C8-N9-C4	7.96	109.58	106.40
26	14	2584	U	C2-N1-C1'	7.96	127.25	117.70
1	13	577	G	N1-C6-O6	7.96	124.67	119.90
26	1H	2440	C	N3-C4-C5	-7.96	118.72	121.90
26	1H	2775	A	N1-C6-N6	7.96	123.37	118.60
26	14	2873	A	C5-N7-C8	-7.96	99.92	103.90
26	14	2374	C	C2-N3-C4	-7.95	115.92	119.90
26	1H	1647	G	O5'-P-OP1	-7.95	98.55	105.70
26	1H	2712	U	N3-C2-O2	-7.95	116.64	122.20
26	14	1942	C	C6-N1-C2	-7.95	117.12	120.30
26	1H	631	A	N7-C8-N9	-7.95	109.83	113.80
26	1H	1252	G	C8-N9-C4	7.94	109.58	106.40
26	1H	681	G	N1-C2-N3	7.94	128.66	123.90
26	1H	784	A	C5-C6-N6	7.94	130.05	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1598	C	O5'-P-OP2	7.94	120.23	110.70
1	13	889	A	O5'-P-OP2	-7.94	98.56	105.70
22	1L	74	C	O4'-C1'-N1	7.94	114.55	108.20
26	1H	123	G	C6-N1-C2	-7.94	120.34	125.10
26	1H	966	G	N1-C6-O6	-7.94	115.14	119.90
1	13	1158	C	N1-C2-O2	7.93	123.66	118.90
26	14	2406	U	O4'-C1'-N1	-7.93	101.85	108.20
24	3K	76	A	C2-N3-C4	-7.93	106.63	110.60
26	14	2513	G	C5-C6-O6	-7.93	123.84	128.60
1	13	811	C	C5-C6-N1	-7.93	117.03	121.00
26	1H	113	G	N1-C6-O6	7.93	124.66	119.90
26	1H	852	G	O5'-P-OP1	7.93	120.22	110.70
26	1H	1335	U	N3-C2-O2	-7.93	116.65	122.20
26	14	1204	A	O4'-C1'-N9	7.93	114.54	108.20
26	1H	1836	C	N3-C4-C5	-7.93	118.73	121.90
26	1H	1842	G	C8-N9-C4	7.93	109.57	106.40
26	14	2307	G	O4'-C1'-N9	7.93	114.54	108.20
26	1H	530	G	C5-C6-O6	7.92	133.35	128.60
26	1H	2271	G	N3-C4-C5	-7.92	124.64	128.60
1	1G	1395	C	O5'-P-OP1	-7.92	98.57	105.70
26	14	664	C	C5-C6-N1	-7.92	117.04	121.00
26	14	1956	U	N1-C2-O2	7.92	128.35	122.80
26	14	189	G	C4-C5-N7	7.92	113.97	110.80
26	14	676	A	O4'-C1'-N9	7.92	114.54	108.20
1	1G	882	C	O5'-P-OP1	-7.92	98.57	105.70
26	14	1600	C	C5-C6-N1	-7.92	117.04	121.00
23	2K	6	G	C8-N9-C4	7.92	109.57	106.40
26	1H	641	C	O5'-P-OP2	7.92	120.20	110.70
26	14	2253	G	C5-C6-O6	-7.91	123.85	128.60
26	1H	1314	C	C6-N1-C2	-7.91	117.14	120.30
26	14	2330	G	C8-N9-C4	7.91	109.56	106.40
26	14	568	U	N1-C2-N3	-7.91	110.16	114.90
26	14	2313	C	C6-N1-C2	-7.90	117.14	120.30
1	13	812	C	O5'-P-OP2	7.90	120.18	110.70
26	1H	731	C	C5-C4-N4	-7.90	114.67	120.20
26	14	459	U	N3-C2-O2	-7.90	116.67	122.20
26	1H	765	G	O5'-P-OP1	-7.90	98.59	105.70
27	16	81	G	C4-C5-N7	7.90	113.96	110.80
26	1H	2627	G	C4-C5-N7	7.89	113.96	110.80
26	14	2392	A	C5-C6-N1	-7.89	113.75	117.70
26	1H	2439	A	C5-N7-C8	-7.89	99.95	103.90
24	3K	32	U	C2-N1-C1'	7.89	127.16	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1189	A	C5-C6-N6	-7.89	117.39	123.70
26	1H	2032	G	C2-N3-C4	-7.89	107.96	111.90
26	1H	73	A	C2-N3-C4	7.88	114.54	110.60
26	1H	447	A	C8-N9-C4	-7.88	102.65	105.80
26	1H	2466	C	N3-C4-C5	7.88	125.05	121.90
26	14	621	A	C5-N7-C8	-7.88	99.96	103.90
26	14	2374	C	N3-C4-C5	7.88	125.05	121.90
26	14	1956	U	O5'-P-OP2	-7.88	98.61	105.70
26	14	681	G	N9-C4-C5	-7.88	102.25	105.40
1	13	1526	G	C4-C5-N7	7.88	113.95	110.80
26	1H	1438	U	N3-C4-C5	-7.88	109.87	114.60
26	14	2351	G	N3-C4-N9	7.88	130.72	126.00
26	1H	815	C	N3-C4-C5	7.87	125.05	121.90
26	1H	2054	A	OP2-P-O3'	7.87	122.52	105.20
26	1H	197	A	OP2-P-O3'	7.87	122.52	105.20
26	1H	622	G	N3-C2-N2	7.87	125.41	119.90
26	1H	2028	U	O5'-P-OP1	-7.87	98.62	105.70
1	1G	197	A	N7-C8-N9	7.87	117.74	113.80
26	1H	913	U	N1-C2-O2	7.87	128.31	122.80
26	1H	965	C	O5'-P-OP1	-7.87	98.62	105.70
26	14	1496	A	N1-C6-N6	7.87	123.32	118.60
26	1H	144	C	C2-N3-C4	-7.86	115.97	119.90
26	1H	1305	C	C6-N1-C2	7.86	123.44	120.30
24	3K	76	A	C8-N9-C4	-7.85	102.66	105.80
26	14	1314	C	C2-N1-C1'	7.85	127.44	118.80
26	1H	1622	G	N1-C6-O6	-7.85	115.19	119.90
26	1H	932	G	N1-C2-N2	-7.84	109.14	116.20
26	14	74	A	C5-N7-C8	-7.84	99.98	103.90
26	14	2046	G	N1-C6-O6	-7.84	115.19	119.90
26	14	2452	C	N3-C4-N4	7.84	123.49	118.00
26	1H	203	C	O5'-P-OP1	-7.84	98.64	105.70
26	1H	1210	A	C4-C5-N7	7.84	114.62	110.70
26	1H	630	G	C8-N9-C4	7.84	109.53	106.40
26	1H	189	G	C8-N9-C4	7.83	109.53	106.40
1	13	1205	U	N3-C4-C5	-7.83	109.90	114.60
26	1H	213	A	O5'-P-OP1	7.83	120.10	110.70
26	1H	1678	G	N1-C6-O6	7.83	124.60	119.90
1	13	1195	C	C6-N1-C2	-7.83	117.17	120.30
26	1H	728	G	N7-C8-N9	-7.83	109.19	113.10
1	13	880	C	C6-N1-C2	7.83	123.43	120.30
26	1H	769	G	N1-C2-N2	-7.83	109.15	116.20
26	14	2217	G	C6-C5-N7	-7.83	125.70	130.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	252	G	O5'-P-OP1	7.83	120.09	110.70
26	1H	845	G	C4-C5-N7	7.83	113.93	110.80
26	1H	1204	A	N1-C6-N6	7.82	123.29	118.60
26	1H	2261	C	OP2-P-O3'	7.82	122.41	105.20
26	1H	2036	C	C6-N1-C2	-7.82	117.17	120.30
26	14	2501	C	N3-C4-C5	7.82	125.03	121.90
26	1H	793	A	N1-C6-N6	7.82	123.29	118.60
26	1H	1382	G	C4-C5-N7	7.82	113.93	110.80
43	E8	42	ARG	NE-CZ-NH1	-7.82	116.39	120.30
26	14	74	A	N3-C4-N9	-7.82	121.14	127.40
26	14	1121	C	C5-C6-N1	-7.82	117.09	121.00
1	13	522	C	O5'-P-OP2	-7.81	98.67	105.70
26	1H	148	C	C5-C6-N1	-7.81	117.09	121.00
26	1H	1200	C	C2-N3-C4	-7.81	115.99	119.90
26	1H	1639	U	C5-C6-N1	-7.81	118.80	122.70
1	1G	1442	G	N3-C4-C5	7.81	132.50	128.60
26	14	801	G	O5'-P-OP2	-7.81	98.67	105.70
26	1H	2261	C	O5'-P-OP2	-7.80	98.68	105.70
26	1H	2609	U	C2-N3-C4	-7.80	122.32	127.00
1	1G	1112	C	C6-N1-C2	-7.80	117.18	120.30
26	1H	2552	U	N1-C2-N3	7.79	119.58	114.90
26	1H	631	A	C8-N9-C4	7.79	108.92	105.80
1	1G	1527	C	C5-C6-N1	-7.79	117.10	121.00
26	14	569	U	C2-N3-C4	-7.79	122.33	127.00
26	14	2829	C	N1-C2-O2	-7.79	114.22	118.90
26	14	735	A	C8-N9-C4	7.79	108.92	105.80
26	1H	839	U	N1-C2-N3	7.79	119.57	114.90
26	14	2244	U	N1-C2-O2	-7.79	117.35	122.80
26	1H	2375	G	C4-C5-N7	7.78	113.91	110.80
26	14	1609	A	C6-N1-C2	-7.78	113.93	118.60
26	1H	22	C	O5'-P-OP1	-7.78	98.70	105.70
26	1H	728	G	N9-C4-C5	-7.78	102.29	105.40
26	1H	1915	U	N3-C2-O2	-7.78	116.76	122.20
26	1H	2508	G	N9-C4-C5	7.78	108.51	105.40
26	1H	1204	A	N7-C8-N9	7.78	117.69	113.80
36	78	20	GLY	N-CA-C	7.77	132.53	113.10
23	2L	21	U	N1-C2-N3	7.77	119.56	114.90
26	14	1661	G	C8-N9-C4	7.77	109.51	106.40
26	1H	1933	G	C8-N9-C4	-7.77	103.29	106.40
1	13	644	G	O5'-P-OP2	-7.76	98.71	105.70
26	14	1309	G	C8-N9-C1'	-7.76	116.91	127.00
26	14	49	A	P-O3'-C3'	7.76	129.02	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	138	G	N7-C8-N9	7.76	116.98	113.10
26	1H	2377	A	N9-C4-C5	-7.76	102.69	105.80
26	1H	2444	G	C8-N9-C4	-7.76	103.30	106.40
27	16	81	G	C5-N7-C8	-7.76	100.42	104.30
26	14	974(A)	C	C5-C4-N4	7.76	125.63	120.20
26	1H	2062	A	N9-C4-C5	-7.76	102.70	105.80
26	14	2688	U	C5-C4-O4	7.76	130.55	125.90
26	1H	860	U	C2-N3-C4	-7.75	122.35	127.00
26	1H	214	G	C2-N3-C4	7.75	115.78	111.90
26	14	2044	C	N1-C2-O2	-7.75	114.25	118.90
26	1H	121	G	C5-C6-N1	7.75	115.37	111.50
26	1H	1355	G	N1-C6-O6	-7.75	115.25	119.90
26	14	2346	A	C8-N9-C4	-7.75	102.70	105.80
26	14	2463	C	C6-N1-C2	7.75	123.40	120.30
26	14	2502	G	C8-N9-C4	-7.75	103.30	106.40
26	1H	2473	U	C2-N1-C1'	7.75	126.99	117.70
26	1H	16	G	N3-C2-N2	-7.74	114.48	119.90
26	1H	526	A	N9-C4-C5	7.74	108.90	105.80
26	1H	2866	U	C5-C4-O4	7.74	130.54	125.90
26	1H	189	G	C5-C6-O6	-7.74	123.96	128.60
27	1J	101	A	C5-N7-C8	-7.74	100.03	103.90
26	1H	1928	A	O5'-P-OP1	-7.74	98.74	105.70
26	1H	2614	A	C2-N3-C4	7.74	114.47	110.60
26	14	2598	A	C8-N9-C4	7.74	108.89	105.80
26	1H	1437	C	C2-N1-C1'	7.73	127.31	118.80
26	1H	2031	A	C5-C6-N6	-7.73	117.52	123.70
26	14	1934	C	C6-N1-C2	7.73	123.39	120.30
26	1H	788	A	C8-N9-C4	7.73	108.89	105.80
26	14	2429	G	OP2-P-O3'	7.73	122.21	105.20
26	1H	141	A	O4'-C1'-N9	7.73	114.38	108.20
1	1G	893	C	C6-N1-C2	7.72	123.39	120.30
1	13	974	A	C6-C5-N7	-7.72	126.90	132.30
26	14	1612	C	N3-C4-C5	7.72	124.99	121.90
26	14	784	A	C5-C6-N1	-7.72	113.84	117.70
26	1H	482	A	C8-N9-C4	-7.72	102.71	105.80
26	1H	763	G	C6-N1-C2	-7.72	120.47	125.10
26	1H	2508	G	C8-N9-C4	-7.71	103.31	106.40
26	14	783	A	N1-C2-N3	7.71	133.16	129.30
26	1H	633	A	C6-C5-N7	-7.71	126.91	132.30
26	1H	2503	A	N1-C2-N3	-7.71	125.45	129.30
26	14	2301	C	C6-N1-C2	-7.71	117.22	120.30
26	14	127	A	C5-C6-N6	-7.70	117.54	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1612	C	C5-C6-N1	-7.70	117.15	121.00
26	1H	698	C	C5-C6-N1	-7.70	117.15	121.00
26	14	679	C	N3-C2-O2	7.70	127.29	121.90
1	13	1158	C	C2-N1-C1'	7.70	127.27	118.80
26	1H	667	U	N3-C2-O2	7.69	127.59	122.20
26	1H	2607	G	N9-C4-C5	-7.69	102.32	105.40
26	1H	2318	G	O4'-C1'-N9	7.69	114.35	108.20
4	3E	12	CYS	CA-CB-SG	7.69	127.84	114.00
26	1H	148	C	C6-N1-C2	7.69	123.38	120.30
26	1H	2548	G	N1-C6-O6	-7.69	115.28	119.90
1	13	1526	G	C5-C6-O6	-7.69	123.99	128.60
26	1H	1297	C	OP2-P-O3'	-7.69	88.29	105.20
26	1H	138	G	C8-N9-C4	-7.69	103.33	106.40
48	F5	82	LEU	CA-CB-CG	7.68	132.97	115.30
1	13	1227	A	C2-N3-C4	-7.68	106.76	110.60
1	1G	812	C	P-O3'-C3'	7.68	128.92	119.70
26	14	1304	C	N3-C2-O2	-7.68	116.52	121.90
26	1H	190	A	C5-C6-N6	-7.68	117.56	123.70
26	14	1698	A	C5-C6-N1	-7.68	113.86	117.70
26	1H	616	A	N1-C6-N6	7.67	123.20	118.60
39	65	110	LEU	CA-CB-CG	7.67	132.95	115.30
26	1H	71	A	O4'-C1'-N9	-7.67	102.06	108.20
1	1G	1286	A	C8-N9-C4	-7.67	102.73	105.80
26	14	468	G	OP1-P-OP2	-7.67	108.10	119.60
26	14	561	G	N3-C4-N9	-7.67	121.40	126.00
26	1H	599	G	N3-C2-N2	7.66	125.26	119.90
26	1H	103	A	N1-C6-N6	7.66	123.19	118.60
26	14	683	C	N3-C4-C5	7.66	124.96	121.90
26	14	2679	A	O5'-P-OP2	-7.66	98.81	105.70
26	1H	1232	G	C5-C6-O6	-7.66	124.01	128.60
1	1G	1432	G	C5-C6-N1	-7.66	107.67	111.50
26	1H	2030	A	C5-C6-N6	-7.65	117.58	123.70
26	1H	1266	G	C8-N9-C4	7.65	109.46	106.40
26	1H	2403	C	C6-N1-C2	-7.65	117.24	120.30
26	1H	2424	C	C2-N3-C4	7.65	123.72	119.90
26	14	134	C	C2-N3-C4	-7.65	116.08	119.90
26	1H	2287	A	N3-C4-C5	7.65	132.15	126.80
26	1H	675	A	C4-C5-N7	7.64	114.52	110.70
26	1H	2318	G	C8-N9-C4	-7.64	103.34	106.40
28	11	131	LEU	CA-CB-CG	7.64	132.88	115.30
26	14	1142	U	N1-C2-O2	7.64	128.15	122.80
26	14	2426	A	N9-C4-C5	-7.64	102.74	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	821	A	C4-C5-C6	7.64	120.82	117.00
1	13	1491	G	OP2-P-O3'	7.64	122.00	105.20
26	1H	1204	A	C5-N7-C8	-7.64	100.08	103.90
26	1H	271(B)	G	P-O3'-C3'	7.63	128.86	119.70
26	1H	1123	C	C2-N3-C4	-7.63	116.08	119.90
26	1H	2490	G	C8-N9-C4	-7.63	103.35	106.40
1	13	910	C	N3-C4-C5	7.63	124.95	121.90
26	1H	1616	A	C6-C5-N7	-7.62	126.96	132.30
26	1H	1616	A	O4'-C1'-N9	7.62	114.30	108.20
26	1H	2230	G	N3-C4-N9	-7.62	121.43	126.00
26	1H	2586	C	C5-C4-N4	-7.62	114.86	120.20
26	1H	2401	U	C6-N1-C2	-7.62	116.43	121.00
26	1H	2485	G	N9-C4-C5	-7.62	102.35	105.40
26	14	1394	U	O5'-P-OP2	7.62	119.84	110.70
26	1H	245	G	C4-N9-C1'	7.62	136.40	126.50
26	1H	2576	G	N9-C4-C5	-7.62	102.35	105.40
26	1H	88	G	N3-C4-C5	-7.61	124.79	128.60
26	1H	2411	A	N1-C6-N6	-7.61	114.03	118.60
26	1H	735	A	N1-C6-N6	7.61	123.17	118.60
26	14	330	A	C5-N7-C8	-7.61	100.09	103.90
26	14	470	A	C5-C6-N6	-7.61	117.61	123.70
1	13	186(A)	C	C6-N1-C2	-7.61	117.26	120.30
26	1H	1310	G	C5-C6-O6	-7.61	124.03	128.60
26	1H	1622	G	C6-C5-N7	7.61	134.97	130.40
26	14	2512	C	O5'-P-OP1	-7.61	98.85	105.70
26	1H	607	U	C6-N1-C2	7.61	125.56	121.00
26	1H	2271	G	C6-N1-C2	-7.60	120.54	125.10
26	1H	930	U	N3-C4-O4	-7.60	114.08	119.40
26	1H	2731	G	C6-C5-N7	-7.60	125.84	130.40
26	14	1379	A	C5-N7-C8	-7.60	100.10	103.90
26	14	2587	A	N1-C6-N6	7.60	123.16	118.60
26	1H	874	G	O5'-P-OP2	-7.60	98.86	105.70
26	14	780	G	O5'-P-OP1	-7.59	98.87	105.70
26	1H	1564	C	C5-C4-N4	7.59	125.51	120.20
26	1H	2451	A	N3-C4-N9	-7.59	121.33	127.40
26	14	1396	U	O5'-P-OP1	-7.59	98.87	105.70
26	1H	739	G	O5'-P-OP2	-7.59	98.87	105.70
26	1H	1914	C	C6-N1-C2	-7.58	117.27	120.30
26	14	2247	A	N1-C2-N3	7.58	133.09	129.30
1	13	1512	U	C5-C6-N1	-7.58	118.91	122.70
26	1H	560	C	O5'-P-OP1	-7.58	98.88	105.70
26	14	2592	G	N3-C4-N9	7.58	130.55	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2699	C	C5-C4-N4	-7.58	114.90	120.20
26	14	330	A	C4-C5-N7	7.58	114.49	110.70
26	1H	738	G	C6-N1-C2	-7.57	120.56	125.10
26	14	2331	G	N9-C4-C5	-7.57	102.37	105.40
26	1H	78	A	C8-N9-C4	7.57	108.83	105.80
26	14	199	A	O5'-P-OP2	-7.57	98.89	105.70
1	1G	898	G	C8-N9-C4	7.57	109.43	106.40
26	1H	1702	G	O5'-P-OP1	-7.57	98.89	105.70
26	1H	2616	C	N1-C2-O2	-7.57	114.36	118.90
23	2L	21	U	N3-C2-O2	-7.57	116.90	122.20
26	14	1313	U	C5-C6-N1	7.57	126.48	122.70
26	1H	2448	A	C5-N7-C8	-7.57	100.12	103.90
27	16	112	G	N9-C4-C5	-7.57	102.37	105.40
1	13	1128	C	C5-C6-N1	7.56	124.78	121.00
26	14	2278	A	C8-N9-C4	-7.56	102.78	105.80
1	13	1488	G	C6-C5-N7	-7.56	125.86	130.40
26	1H	139	G	N3-C4-C5	-7.56	124.82	128.60
1	1G	690	G	N3-C4-C5	7.56	132.38	128.60
26	14	1572	A	N1-C6-N6	7.56	123.14	118.60
26	1H	825	C	N3-C4-N4	7.56	123.29	118.00
26	1H	1977	A	O5'-P-OP2	-7.56	98.89	105.70
26	1H	1602	U	O5'-P-OP2	7.56	119.77	110.70
26	1H	2466	C	C5-C4-N4	-7.56	114.91	120.20
26	1H	2627	G	N1-C6-O6	7.56	124.44	119.90
1	1G	576	G	C6-C5-N7	-7.56	125.86	130.40
26	1H	74	A	N3-C4-N9	-7.56	121.35	127.40
26	1H	1220	A	N1-C6-N6	-7.56	114.07	118.60
26	1H	2053	G	N1-C2-N2	7.56	123.00	116.20
26	1H	2578	G	N7-C8-N9	-7.56	109.32	113.10
26	14	848	G	C4-N9-C1'	7.55	136.32	126.50
23	2K	21	U	C2-N1-C1'	7.55	126.76	117.70
1	13	1519	A	C5-C6-N6	7.55	129.74	123.70
26	1H	203	C	N1-C2-O2	-7.55	114.37	118.90
26	1H	1368	G	O5'-P-OP2	-7.55	98.91	105.70
26	14	1564	C	N3-C2-O2	-7.55	116.62	121.90
26	1H	792	G	C6-C5-N7	-7.55	125.87	130.40
1	1G	197	A	C8-N9-C4	-7.55	102.78	105.80
26	14	1681	G	C5-N7-C8	-7.55	100.53	104.30
26	1H	679	C	C4-C5-C6	7.54	121.17	117.40
26	1H	692	C	C5-C6-N1	-7.54	117.23	121.00
26	1H	733	G	N9-C4-C5	-7.54	102.38	105.40
26	1H	74	A	O4'-C1'-N9	-7.54	102.17	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1229(A)	G	O5'-P-OP2	-7.54	98.91	105.70
26	1H	2552	U	C5-C6-N1	-7.54	118.93	122.70
26	1H	2688	U	C4-C5-C6	7.54	124.23	119.70
26	1H	128	C	O5'-P-OP2	-7.54	98.91	105.70
26	1H	733	G	C6-C5-N7	-7.54	125.88	130.40
26	1H	1266	G	N1-C6-O6	7.54	124.42	119.90
26	1H	1614	A	O4'-C1'-N9	7.54	114.23	108.20
26	1H	2274	A	OP1-P-OP2	-7.54	108.29	119.60
1	1G	1501	C	C6-N1-C2	7.54	123.32	120.30
26	14	2435	A	N7-C8-N9	7.54	117.57	113.80
26	1H	932	G	N3-C2-N2	7.54	125.18	119.90
26	1H	1360	A	N1-C6-N6	7.54	123.12	118.60
26	14	466	A	C5-C6-N6	-7.54	117.67	123.70
26	14	1278	A	N7-C8-N9	-7.54	110.03	113.80
26	14	193	U	C2-N3-C4	-7.54	122.48	127.00
26	1H	1700	A	OP1-P-OP2	7.54	130.90	119.60
26	1H	2540	C	C6-N1-C2	7.53	123.31	120.30
26	14	1306	C	O5'-P-OP1	-7.53	98.92	105.70
26	1H	2232	U	N3-C2-O2	-7.53	116.93	122.20
26	14	2377	A	C8-N9-C4	7.53	108.81	105.80
26	14	2712	U	N1-C2-N3	7.53	119.42	114.90
26	1H	2447	G	N1-C6-O6	-7.52	115.39	119.90
26	14	2473	U	C2-N1-C1'	7.52	126.73	117.70
1	13	1526	G	C6-C5-N7	-7.52	125.89	130.40
26	1H	2430	A	C4-C5-N7	7.52	114.46	110.70
26	1H	635	C	O5'-P-OP1	7.52	119.72	110.70
26	1H	2070	G	N3-C4-N9	7.52	130.51	126.00
27	16	98	G	OP1-P-OP2	7.52	130.88	119.60
26	14	2506	U	C2-N1-C1'	7.52	126.72	117.70
26	14	2592	G	N3-C4-C5	-7.52	124.84	128.60
26	1H	791	C	OP1-P-O3'	-7.52	88.66	105.20
1	13	575	G	O4'-C1'-N9	-7.51	102.19	108.20
26	1H	663	G	C4-C5-C6	7.51	123.31	118.80
26	14	1462	C	C6-N1-C2	-7.51	117.29	120.30
1	13	817	C	C6-N1-C2	7.51	123.31	120.30
26	1H	1278	A	C5-C6-N6	-7.51	117.69	123.70
26	14	110	G	N1-C6-O6	7.51	124.41	119.90
1	13	967	C	N1-C2-O2	-7.51	114.39	118.90
26	1H	1428	C	C5-C6-N1	-7.51	117.25	121.00
26	14	1366	A	C4-C5-N7	7.51	114.45	110.70
26	1H	271(C)	U	C2-N1-C1'	7.51	126.71	117.70
26	1H	311	A	O5'-P-OP1	-7.50	98.95	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1671	U	OP1-P-OP2	7.50	130.85	119.60
26	1H	1022	G	C8-N9-C4	-7.50	103.40	106.40
26	1H	2406	U	O4'-C1'-N1	-7.50	102.20	108.20
26	14	681	G	N3-C4-N9	7.50	130.50	126.00
26	1H	1589	C	O5'-P-OP2	7.50	119.70	110.70
26	1H	1698	A	N7-C8-N9	7.50	117.55	113.80
27	16	44	G	C8-N9-C1'	7.50	136.75	127.00
28	19	272	ALA	N-CA-C	7.50	131.25	111.00
26	1H	839	U	C4-C5-C6	7.50	124.20	119.70
26	1H	1213	A	O5'-P-OP2	7.50	119.69	110.70
26	1H	1614	A	N7-C8-N9	7.50	117.55	113.80
26	1H	1253	A	C8-N9-C4	7.50	108.80	105.80
26	14	528	A	C8-N9-C1'	7.50	141.19	127.70
1	13	233	C	C6-N1-C2	-7.49	117.30	120.30
26	1H	593	G	C2-N3-C4	-7.49	108.16	111.90
1	13	185	A	C8-N9-C4	-7.49	102.80	105.80
26	1H	575	A	C6-N1-C2	-7.49	114.11	118.60
26	1H	1022	G	P-O3'-C3'	7.49	128.69	119.70
26	1H	1269	A	C2-N3-C4	-7.49	106.86	110.60
26	1H	1797	C	C5-C6-N1	-7.49	117.26	121.00
23	2K	74	A	C5-C6-N6	-7.48	117.71	123.70
26	1H	784	A	N9-C4-C5	7.48	108.79	105.80
26	1H	146	G	N1-C6-O6	7.48	124.39	119.90
26	1H	575	A	N1-C2-N3	7.48	133.04	129.30
26	1H	2286	A	C8-N9-C4	-7.48	102.81	105.80
26	14	670	A	C8-N9-C4	7.48	108.79	105.80
26	14	974(A)	C	N3-C2-O2	-7.48	116.66	121.90
26	14	2000	G	N7-C8-N9	-7.48	109.36	113.10
26	1H	198	C	C2-N3-C4	-7.48	116.16	119.90
26	14	252	G	C6-C5-N7	7.48	134.89	130.40
26	1H	2672	G	C6-C5-N7	-7.48	125.92	130.40
26	1H	2622	C	N3-C2-O2	7.47	127.13	121.90
26	1H	692	C	C5-C4-N4	-7.47	114.97	120.20
26	1H	1623	G	N1-C6-O6	-7.47	115.42	119.90
1	1G	197	A	P-O3'-C3'	7.47	128.66	119.70
1	1G	1442	G	N3-C4-N9	-7.47	121.52	126.00
26	1H	1356	G	C5-C6-O6	-7.47	124.12	128.60
26	1H	1468	C	C6-N1-C2	-7.47	117.31	120.30
26	1H	2385	C	O5'-P-OP2	-7.47	98.98	105.70
26	14	1333	C	N3-C4-C5	7.47	124.89	121.90
26	14	2880	C	C6-N1-C2	-7.46	117.31	120.30
1	13	1502	A	N7-C8-N9	7.46	117.53	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	936	C	C6-N1-C2	7.46	123.28	120.30
26	1H	948	G	C5-N7-C8	-7.46	100.57	104.30
26	1H	1632	A	N9-C4-C5	-7.46	102.81	105.80
26	14	2449	U	N3-C4-O4	7.46	124.62	119.40
26	14	2542	A	C8-N9-C4	7.46	108.78	105.80
1	13	892	A	C2-N3-C4	-7.46	106.87	110.60
1	13	966	G	N9-C4-C5	-7.46	102.42	105.40
23	2L	40	C	C6-N1-C2	-7.46	117.32	120.30
26	1H	733	G	N1-C2-N2	-7.46	109.49	116.20
1	13	1197	G	OP1-P-O3'	7.45	121.60	105.20
26	1H	566	U	C5-C4-O4	-7.45	121.43	125.90
26	1H	1279	G	N1-C6-O6	-7.45	115.43	119.90
26	1H	2485	G	C8-N9-C4	7.45	109.38	106.40
26	1H	2392	A	C6-N1-C2	7.45	123.07	118.60
26	1H	2509	G	C5-C6-N1	7.45	115.23	111.50
26	14	134	C	N3-C4-C5	7.45	124.88	121.90
26	14	786	C	C6-N1-C2	7.45	123.28	120.30
26	1H	837	C	O5'-P-OP2	7.45	119.63	110.70
26	1H	2371	G	N1-C6-O6	7.45	124.37	119.90
26	14	737	C	C5-C4-N4	-7.45	114.99	120.20
1	13	778	G	O5'-P-OP2	-7.44	99.00	105.70
26	1H	2392	A	N3-C4-N9	-7.44	121.44	127.40
26	14	1681	G	C4-C5-N7	7.44	113.78	110.80
26	14	2577	A	C5-C6-N1	-7.44	113.98	117.70
29	29	78	LEU	CA-CB-CG	7.44	132.42	115.30
26	1H	2595	G	C5-C6-O6	-7.44	124.13	128.60
26	1H	1305	C	N3-C4-C5	7.44	124.88	121.90
26	1H	2212	A	O4'-C1'-N9	7.44	114.15	108.20
26	14	134	C	C5-C6-N1	-7.44	117.28	121.00
26	14	1314	C	N1-C2-N3	-7.44	113.99	119.20
26	1H	2028	U	O5'-P-OP2	7.44	119.63	110.70
26	14	693	C	N3-C4-N4	-7.44	112.79	118.00
26	14	1332	G	N9-C4-C5	7.44	108.38	105.40
26	14	1776	G	O5'-P-OP1	7.44	119.63	110.70
26	1H	840	C	C6-N1-C2	7.44	123.28	120.30
26	1H	1332	G	N1-C6-O6	7.44	124.36	119.90
26	1H	2453	A	N1-C6-N6	-7.44	114.14	118.60
1	1G	862	C	C6-N1-C2	-7.44	117.33	120.30
26	14	856	C	C2-N1-C1'	7.44	126.98	118.80
24	3K	76	A	C5-C6-N6	-7.43	117.75	123.70
26	14	2542	A	N7-C8-N9	-7.43	110.08	113.80
26	1H	446	G	N9-C4-C5	-7.43	102.43	105.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	337	C	C5-C6-N1	7.43	124.72	121.00
26	14	2042	A	O5'-P-OP1	7.43	119.62	110.70
1	13	967	C	N3-C2-O2	7.43	127.10	121.90
26	1H	699	A	C2-N3-C4	7.43	114.31	110.60
26	1H	1632	A	C5-C6-N6	-7.43	117.76	123.70
26	1H	2392	A	C8-N9-C4	-7.43	102.83	105.80
26	14	912	C	C6-N1-C2	-7.43	117.33	120.30
26	1H	782	A	N9-C4-C5	7.43	108.77	105.80
26	1H	1729	A	O4'-C1'-N9	7.43	114.14	108.20
26	1H	2287	A	O5'-P-OP2	-7.43	99.02	105.70
26	1H	2822	G	C6-C5-N7	-7.43	125.94	130.40
26	1H	1394	U	O5'-P-OP2	7.42	119.61	110.70
26	1H	1602	U	N1-C2-N3	7.42	119.36	114.90
26	1H	1626	G	C8-N9-C4	-7.42	103.43	106.40
1	13	963	G	N3-C2-N2	7.42	125.10	119.90
1	13	576	G	N3-C4-N9	7.42	130.45	126.00
26	1H	2822	G	N9-C4-C5	-7.42	102.43	105.40
26	1H	1990	C	N3-C2-O2	-7.42	116.71	121.90
26	1H	2689	U	C5-C4-O4	7.42	130.35	125.90
26	14	488	G	C8-N9-C4	7.42	109.37	106.40
26	14	769	G	OP1-P-O3'	7.42	121.52	105.20
1	13	910	C	C6-N1-C2	7.42	123.27	120.30
26	14	1471	A	C8-N9-C4	-7.41	102.83	105.80
26	1H	512	G	O5'-P-OP1	-7.41	99.03	105.70
26	1H	1336	A	C2-N3-C4	7.41	114.30	110.60
26	1H	2430	A	C6-C5-N7	-7.41	127.12	132.30
26	1H	775	G	O4'-C1'-N9	7.40	114.12	108.20
26	14	2601	C	C6-N1-C2	-7.40	117.34	120.30
26	1H	1596	A	OP2-P-O3'	7.39	121.47	105.20
26	14	209	C	C2-N3-C4	-7.39	116.20	119.90
26	14	441	U	O5'-P-OP1	-7.39	99.05	105.70
26	14	502	A	C2-N3-C4	-7.39	106.90	110.60
1	13	353	A	C8-N9-C4	-7.39	102.84	105.80
26	1H	796	C	C5-C6-N1	-7.39	117.30	121.00
26	1H	1381	G	N3-C2-N2	-7.39	114.73	119.90
1	13	792	A	N3-C4-C5	7.39	131.97	126.80
26	14	1022	G	N9-C4-C5	7.39	108.36	105.40
26	1H	986	C	OP1-P-OP2	-7.38	108.52	119.60
26	1H	1248	G	N3-C2-N2	-7.38	114.73	119.90
26	1H	2192	G	C6-C5-N7	-7.38	125.97	130.40
1	1G	690	G	C5-N7-C8	-7.38	100.61	104.30
26	14	1340	U	C6-N1-C2	7.38	125.43	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1956	U	N3-C2-O2	-7.38	117.03	122.20
26	1H	2830	G	C8-N9-C4	-7.38	103.45	106.40
26	1H	2689	U	C2-N3-C4	-7.38	122.57	127.00
26	14	1400	G	O5'-P-OP1	7.38	119.56	110.70
26	1H	510	C	O5'-P-OP2	-7.38	99.06	105.70
26	1H	2056	G	OP1-P-O3'	7.38	121.44	105.20
26	1H	40	C	C5-C6-N1	-7.37	117.31	121.00
26	1H	1707	G	N1-C6-O6	7.37	124.32	119.90
26	14	114	U	N3-C4-O4	7.37	124.56	119.40
26	14	2386	C	C6-N1-C2	7.37	123.25	120.30
26	1H	2599	G	C5-C6-O6	7.37	133.02	128.60
23	2K	77	A	O5'-P-OP1	-7.37	99.07	105.70
39	A8	54	LEU	CA-CB-CG	7.36	132.24	115.30
26	14	574	C	C6-N1-C2	7.36	123.25	120.30
26	14	779	U	N1-C2-O2	7.36	127.95	122.80
1	13	690	G	C2-N3-C4	-7.36	108.22	111.90
26	1H	1950	G	C6-C5-N7	-7.36	125.98	130.40
26	14	2375	G	N7-C8-N9	-7.36	109.42	113.10
26	1H	121	G	C5-C6-O6	-7.36	124.19	128.60
26	1H	2518	A	C8-N9-C4	-7.36	102.86	105.80
26	14	1585	C	N1-C2-O2	7.36	123.31	118.90
26	1H	1332	G	C5-C6-N1	-7.35	107.82	111.50
1	13	1371	G	O5'-P-OP2	7.35	119.52	110.70
26	14	1402	C	C6-N1-C2	-7.35	117.36	120.30
26	1H	586	A	O5'-P-OP2	-7.35	99.08	105.70
26	1H	2699	C	C2-N3-C4	-7.35	116.22	119.90
1	1G	328	C	C6-N1-C2	-7.35	117.36	120.30
26	14	856	C	C5-C6-N1	7.35	124.67	121.00
26	1H	66	C	C6-N1-C2	-7.35	117.36	120.30
26	14	741	G	O5'-P-OP2	7.35	119.52	110.70
26	14	774	A	O5'-P-OP2	-7.35	99.09	105.70
26	14	2867	G	O4'-C1'-N9	7.35	114.08	108.20
1	13	525	C	C5-C4-N4	-7.34	115.06	120.20
26	14	1359	A	C8-N9-C4	7.34	108.74	105.80
26	14	2779	U	N3-C2-O2	-7.34	117.06	122.20
1	1G	108	G	C5-N7-C8	-7.34	100.63	104.30
26	14	15	G	O5'-P-OP1	-7.34	99.09	105.70
26	14	1671	U	O5'-P-OP1	-7.34	99.09	105.70
26	14	1762	A	O4'-C1'-N9	7.34	114.07	108.20
26	1H	208	C	C2-N3-C4	-7.34	116.23	119.90
26	1H	1187	G	N1-C6-O6	7.34	124.30	119.90
26	1H	1269	A	C5-N7-C8	-7.34	100.23	103.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1950	G	C4-C5-N7	7.33	113.73	110.80
26	14	1602	U	N3-C4-C5	-7.33	110.20	114.60
26	1H	1210	A	C6-C5-N7	-7.32	127.17	132.30
26	1H	2311	A	C2-N3-C4	-7.32	106.94	110.60
26	1H	2439	A	C2-N3-C4	-7.32	106.94	110.60
26	14	463	G	OP1-P-O3'	7.32	121.31	105.20
26	1H	737	C	C4-C5-C6	7.32	121.06	117.40
26	1H	265	A	C4-C5-N7	7.32	114.36	110.70
26	14	683	C	C2-N3-C4	-7.32	116.24	119.90
26	14	1391	U	O5'-P-OP1	-7.32	99.11	105.70
26	1H	116	C	C5-C6-N1	-7.32	117.34	121.00
26	1H	2059	A	N1-C2-N3	7.32	132.96	129.30
37	88	86	GLY	N-CA-C	-7.32	94.81	113.10
26	14	808	G	O5'-P-OP2	-7.32	99.11	105.70
26	1H	1681	G	C5-N7-C8	-7.32	100.64	104.30
26	1H	2330	G	C2-N3-C4	-7.32	108.24	111.90
26	1H	970	C	O5'-P-OP1	-7.31	99.12	105.70
26	1H	1606	G	C5-C6-O6	-7.31	124.21	128.60
26	1H	1797	C	C4-C5-C6	7.31	121.06	117.40
26	1H	1284	A	N1-C6-N6	7.31	122.99	118.60
27	16	7	G	C6-C5-N7	-7.31	126.01	130.40
26	1H	667	U	N3-C4-O4	7.31	124.52	119.40
26	1H	1189	A	C4-C5-N7	7.31	114.36	110.70
26	1H	1640	C	O5'-P-OP1	7.31	119.47	110.70
26	14	2237	G	N3-C2-N2	7.31	125.02	119.90
26	1H	950	G	C6-C5-N7	-7.31	126.02	130.40
26	1H	2295	C	C6-N1-C2	-7.31	117.38	120.30
26	14	1379	A	C4-C5-N7	7.31	114.35	110.70
26	1H	1198	U	C2-N3-C4	-7.30	122.62	127.00
26	14	1632	A	C5-C6-N6	-7.30	117.86	123.70
26	14	2324	C	C6-N1-C2	7.30	123.22	120.30
26	1H	664	C	C5-C6-N1	-7.30	117.35	121.00
26	14	487	C	O5'-P-OP1	-7.30	99.13	105.70
26	14	2331	G	C8-N9-C4	7.30	109.32	106.40
26	1H	2499	C	C6-N1-C2	-7.30	117.38	120.30
26	14	675	A	N1-C6-N6	7.29	122.98	118.60
26	1H	508	G	C4-C5-N7	7.28	113.71	110.80
26	1H	870	A	C5-C6-N1	7.28	121.34	117.70
26	1H	949	C	C5-C6-N1	-7.28	117.36	121.00
26	14	2038	G	C8-N9-C4	7.28	109.31	106.40
1	13	1310	G	C8-N9-C4	7.28	109.31	106.40
26	1H	2437	U	N3-C4-C5	-7.28	110.23	114.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
27	16	53	A	C8-N9-C4	-7.28	102.89	105.80
26	14	2045	C	C6-N1-C2	7.28	123.21	120.30
1	1G	449	C	N3-C2-O2	-7.27	116.81	121.90
26	14	2502	G	O5'-P-OP1	-7.27	99.15	105.70
1	1G	1297	C	P-O3'-C3'	7.27	128.43	119.70
26	1H	965	C	C5-C6-N1	7.27	124.64	121.00
26	1H	1611	C	C2-N3-C4	-7.27	116.27	119.90
26	1H	2598	A	N9-C4-C5	-7.27	102.89	105.80
24	3K	32	U	N3-C2-O2	-7.26	117.11	122.20
26	1H	2346	A	C5-C6-N1	-7.26	114.07	117.70
26	1H	1210	A	N7-C8-N9	7.26	117.43	113.80
26	1H	1573	G	OP2-P-O3'	7.26	121.17	105.20
26	14	1840	G	N3-C2-N2	-7.26	114.82	119.90
26	14	2503	A	O5'-P-OP2	-7.26	99.17	105.70
26	14	1029	A	N7-C8-N9	-7.26	110.17	113.80
26	14	2518	A	C2-N3-C4	-7.26	106.97	110.60
26	1H	1393	A	O5'-P-OP2	-7.26	99.17	105.70
26	14	565	C	C4-C5-C6	7.26	121.03	117.40
26	1H	738	G	N1-C2-N3	7.25	128.25	123.90
26	1H	845	G	OP1-P-O3'	7.25	121.16	105.20
27	1J	70	C	C6-N1-C2	-7.25	117.40	120.30
26	1H	2731	G	N1-C6-O6	7.25	124.25	119.90
26	14	2247	A	N1-C6-N6	-7.25	114.25	118.60
26	14	2713	A	N7-C8-N9	7.25	117.43	113.80
26	1H	630	G	O5'-P-OP1	7.25	119.40	110.70
26	1H	2004	G	O5'-P-OP2	-7.25	99.18	105.70
27	16	23	G	C8-N9-C4	7.25	109.30	106.40
26	1H	1546	C	O5'-P-OP1	-7.25	99.18	105.70
26	14	1698	A	N7-C8-N9	7.25	117.42	113.80
26	14	2518	A	C4-C5-N7	7.25	114.32	110.70
26	14	759	G	C5-C6-O6	-7.24	124.25	128.60
26	1H	1899	G	N1-C2-N2	7.24	122.72	116.20
26	14	310	A	O5'-P-OP1	-7.24	99.18	105.70
26	14	415	A	O5'-P-OP2	-7.24	99.18	105.70
26	1H	702	G	O5'-P-OP2	-7.24	99.18	105.70
28	11	52	ARG	NE-CZ-NH2	7.24	123.92	120.30
1	1G	332	G	C5-C6-O6	-7.24	124.25	128.60
1	13	1198	G	O5'-P-OP1	-7.24	99.19	105.70
26	1H	74	A	C6-C5-N7	-7.24	127.23	132.30
26	1H	99	U	N3-C2-O2	-7.24	117.13	122.20
26	1H	1899	G	C4-N9-C1'	-7.24	117.09	126.50
26	14	1325	G	N3-C4-N9	7.24	130.34	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2873	A	C8-N9-C4	-7.24	102.91	105.80
26	1H	199	A	C6-C5-N7	7.24	137.37	132.30
26	1H	762	U	N1-C2-O2	7.23	127.86	122.80
26	1H	2070	G	N1-C2-N2	-7.23	109.69	116.20
26	14	2689	U	P-O3'-C3'	7.23	128.38	119.70
26	1H	129	C	C4-C5-C6	7.23	121.01	117.40
26	14	74	A	C6-N1-C2	7.23	122.94	118.60
26	1H	1157	G	N1-C2-N3	7.23	128.24	123.90
1	1G	315	A	N1-C6-N6	7.23	122.94	118.60
26	14	2829	C	N3-C2-O2	7.23	126.96	121.90
26	1H	1894	C	C6-N1-C2	-7.22	117.41	120.30
1	13	576	G	C8-N9-C1'	-7.22	117.61	127.00
26	1H	785	G	C4-C5-N7	-7.22	107.91	110.80
26	1H	1204	A	C4-C5-C6	7.22	120.61	117.00
26	1H	1672	C	N1-C2-O2	-7.22	114.57	118.90
26	1H	2548	G	C5-C6-N1	7.22	115.11	111.50
1	1G	811	C	C6-N1-C2	-7.22	117.41	120.30
26	14	1309	G	N1-C6-O6	7.22	124.23	119.90
26	1H	2279	G	N3-C2-N2	7.22	124.95	119.90
26	1H	1241	A	C2-N3-C4	-7.22	106.99	110.60
1	13	827	U	C2-N1-C1'	7.21	126.36	117.70
26	1H	2552	U	C4-C5-C6	7.21	124.03	119.70
26	1H	636	G	O5'-P-OP2	7.21	119.36	110.70
26	1H	2577	A	N1-C6-N6	-7.21	114.27	118.60
26	1H	1401	G	C8-N9-C4	-7.21	103.52	106.40
26	14	2870	C	C6-N1-C2	-7.21	117.42	120.30
26	14	215	G	N1-C6-O6	7.21	124.23	119.90
26	1H	675	A	N9-C4-C5	-7.21	102.92	105.80
26	1H	906	G	N1-C6-O6	-7.21	115.57	119.90
26	1H	2599	G	C5-N7-C8	7.21	107.91	104.30
1	13	767	A	O5'-P-OP1	-7.21	99.21	105.70
26	1H	1604	C	N1-C2-O2	-7.21	114.58	118.90
26	1H	1955	U	C5-C4-O4	7.21	130.22	125.90
26	14	561	G	N9-C4-C5	7.21	108.28	105.40
26	1H	1938	A	N1-C6-N6	7.20	122.92	118.60
26	1H	1977	A	C8-N9-C4	7.20	108.68	105.80
26	14	951	C	N1-C2-O2	7.20	123.22	118.90
1	1G	1405	G	O5'-P-OP2	-7.20	99.22	105.70
26	14	2392	A	C5-N7-C8	-7.20	100.30	103.90
26	1H	2867	G	O5'-P-OP1	-7.20	99.22	105.70
26	1H	2688	U	N3-C2-O2	-7.20	117.16	122.20
26	1H	568	U	N3-C2-O2	7.19	127.23	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1142	U	C2-N1-C1'	7.19	126.33	117.70
26	14	1835	G	N1-C6-O6	-7.19	115.58	119.90
26	14	247	G	N1-C6-O6	7.19	124.21	119.90
26	1H	945	A	C5-C6-N1	-7.19	114.11	117.70
26	1H	2871	C	O5'-P-OP2	-7.19	99.23	105.70
26	1H	141	A	O5'-P-OP2	-7.18	99.23	105.70
26	1H	508	G	C6-C5-N7	-7.18	126.09	130.40
26	1H	510	C	N3-C4-C5	-7.18	119.03	121.90
26	14	1272	A	N1-C6-N6	7.18	122.91	118.60
26	1H	330	A	N1-C2-N3	7.18	132.89	129.30
26	1H	1885	A	N7-C8-N9	-7.18	110.21	113.80
26	1H	840	C	O5'-P-OP2	-7.17	99.24	105.70
26	14	133	C	C5-C6-N1	-7.17	117.41	121.00
26	1H	2318	G	N3-C4-N9	-7.17	121.70	126.00
26	1H	2439	A	O5'-P-OP2	-7.17	99.25	105.70
26	14	1773	A	O5'-P-OP1	7.17	119.31	110.70
26	1H	78	A	C2-N3-C4	-7.17	107.02	110.60
26	1H	146	G	C8-N9-C4	7.17	109.27	106.40
26	1H	150	C	O5'-P-OP2	-7.17	99.25	105.70
26	1H	752	A	P-O3'-C3'	7.17	128.30	119.70
26	14	1342	A	N1-C6-N6	7.17	122.90	118.60
26	1H	823	G	C8-N9-C4	7.17	109.27	106.40
26	1H	2311	A	O4'-C1'-N9	7.17	113.93	108.20
26	1H	2430	A	O5'-P-OP2	-7.17	99.25	105.70
26	1H	2546	U	C5-C6-N1	-7.17	119.12	122.70
27	16	41	U	C5-C6-N1	-7.17	119.12	122.70
26	14	1609	A	N1-C2-N3	7.16	132.88	129.30
26	1H	259	G	N1-C6-O6	7.16	124.20	119.90
1	1G	1442	G	C4-N9-C1'	-7.16	117.19	126.50
26	1H	1340	U	N3-C4-O4	7.16	124.41	119.40
26	14	1950	G	C5-N7-C8	-7.16	100.72	104.30
26	1H	587	C	N3-C2-O2	-7.16	116.89	121.90
26	14	2555	U	O5'-P-OP1	-7.16	99.26	105.70
27	1J	30	C	C2-N1-C1'	7.16	126.67	118.80
26	1H	2378	A	N1-C6-N6	7.16	122.89	118.60
26	14	1339	G	O5'-P-OP1	-7.16	99.26	105.70
26	14	2492	U	O5'-P-OP2	7.16	119.29	110.70
26	1H	1204	A	C4-N9-C1'	7.15	139.18	126.30
26	1H	1278	A	N1-C6-N6	7.15	122.89	118.60
26	14	2685	G	C5-C6-O6	-7.15	124.31	128.60
26	14	933	A	N1-C6-N6	7.15	122.89	118.60
26	14	2433	A	N7-C8-N9	7.15	117.38	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	459	U	N3-C2-O2	-7.15	117.19	122.20
26	1H	2429	G	N9-C4-C5	7.15	108.26	105.40
1	13	714	G	O5'-P-OP1	-7.15	99.27	105.70
26	1H	693	C	OP2-P-O3'	7.15	120.92	105.20
26	1H	2018	G	C6-C5-N7	-7.15	126.11	130.40
25	4L	13	A	P-O3'-C3'	7.15	128.28	119.70
1	13	690	G	C8-N9-C4	-7.15	103.54	106.40
26	1H	2507	C	N1-C2-O2	7.15	123.19	118.90
26	14	2703	C	C4-C5-C6	7.15	120.97	117.40
26	1H	2509	G	C8-N9-C4	7.14	109.26	106.40
1	1G	817	C	C6-N1-C2	7.14	123.16	120.30
26	1H	435	C	N1-C2-O2	7.14	123.19	118.90
26	14	790	C	O5'-P-OP2	-7.14	99.27	105.70
26	1H	599	G	N3-C4-N9	7.14	130.28	126.00
26	1H	2617	C	C6-N1-C2	7.14	123.16	120.30
26	1H	1827	C	C5-C6-N1	-7.14	117.43	121.00
26	14	795	C	O5'-P-OP2	-7.14	99.28	105.70
26	14	2872	G	C8-N9-C4	-7.14	103.55	106.40
26	1H	2004	G	OP1-P-OP2	7.14	130.31	119.60
26	14	574	C	N3-C4-N4	-7.14	113.00	118.00
26	14	672	C	O5'-P-OP2	-7.14	99.28	105.70
26	1H	2721	A	N1-C6-N6	7.13	122.88	118.60
26	14	584	C	N3-C4-C5	7.13	124.75	121.90
26	1H	1969	A	N9-C4-C5	7.13	108.65	105.80
1	13	768	A	C6-N1-C2	-7.13	114.32	118.60
26	1H	836	G	C2-N3-C4	7.13	115.47	111.90
26	1H	2031	A	C5-C6-N1	7.13	121.27	117.70
26	14	391	G	N1-C6-O6	7.13	124.18	119.90
26	1H	1602	U	C4-C5-C6	7.13	123.98	119.70
26	1H	448	U	N1-C2-N3	7.13	119.18	114.90
26	1H	908	C	C5-C6-N1	-7.13	117.44	121.00
26	1H	2438	U	C5-C6-N1	-7.13	119.14	122.70
26	14	982	C	C6-N1-C2	-7.12	117.45	120.30
26	1H	2726	U	C5-C6-N1	-7.12	119.14	122.70
1	1G	1417	G	N1-C6-O6	7.12	124.17	119.90
26	1H	825	C	N1-C2-O2	-7.12	114.63	118.90
28	11	122	ASP	CB-CG-OD2	7.12	124.71	118.30
1	13	1354	C	C6-N1-C2	-7.12	117.45	120.30
26	1H	213	A	C8-N9-C4	7.12	108.65	105.80
26	14	1558	A	C2-N3-C4	-7.12	107.04	110.60
1	13	901	A	N1-C2-N3	7.11	132.86	129.30
26	1H	2437	U	C5-C4-O4	7.11	130.17	125.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1379	A	N1-C6-N6	7.11	122.87	118.60
26	14	1520	U	C5-C4-O4	7.11	130.17	125.90
26	14	2258	C	C5-C4-N4	-7.11	115.22	120.20
26	14	2700	C	C5-C6-N1	-7.11	117.44	121.00
26	14	492	A	O5'-P-OP2	-7.11	99.30	105.70
26	14	1564	C	C5-C4-N4	7.11	125.18	120.20
26	14	1948	G	O5'-P-OP1	-7.11	99.30	105.70
26	1H	859	G	N3-C4-N9	-7.11	121.74	126.00
26	14	1330	C	O5'-P-OP1	-7.11	99.30	105.70
26	14	1261	C	O5'-P-OP1	-7.10	99.31	105.70
26	14	1373	A	C8-N9-C4	7.10	108.64	105.80
26	1H	248	G	O5'-P-OP2	-7.10	99.31	105.70
26	1H	188	G	C6-C5-N7	-7.10	126.14	130.40
26	14	770	G	OP1-P-OP2	-7.10	108.95	119.60
26	14	1247	A	O5'-P-OP2	-7.10	99.31	105.70
26	1H	660	G	N7-C8-N9	7.10	116.65	113.10
26	14	916	G	O5'-P-OP1	-7.10	99.31	105.70
26	1H	2328	A	C6-N1-C2	-7.09	114.34	118.60
1	1G	413	G	C6-C5-N7	7.09	134.66	130.40
26	1H	446	G	C6-C5-N7	-7.09	126.14	130.40
26	14	673	C	O5'-P-OP1	7.09	119.21	110.70
26	14	2707	G	O4'-C1'-N9	7.09	113.87	108.20
1	13	1322	C	C5-C6-N1	7.09	124.54	121.00
1	1G	1485	U	O5'-P-OP2	-7.09	99.32	105.70
26	1H	262	A	C5-C6-N6	-7.08	118.03	123.70
26	1H	1326	U	O5'-P-OP2	-7.08	99.32	105.70
26	1H	950	G	N1-C2-N3	7.08	128.15	123.90
1	1G	442	C	C6-N1-C2	-7.08	117.47	120.30
26	14	2331	G	N1-C6-O6	7.08	124.15	119.90
26	1H	1298	C	C6-N1-C2	-7.08	117.47	120.30
26	1H	2439	A	O4'-C1'-N9	-7.08	102.54	108.20
26	14	1332	G	C8-N9-C4	-7.08	103.57	106.40
26	1H	2010	G	N1-C6-O6	7.08	124.15	119.90
26	14	613	U	N3-C2-O2	-7.08	117.25	122.20
1	1G	332	G	N9-C4-C5	-7.08	102.57	105.40
26	1H	1612	C	C4-C5-C6	7.07	120.94	117.40
26	14	2056	G	C6-N1-C2	-7.07	120.86	125.10
26	1H	628	G	C8-N9-C4	7.07	109.23	106.40
26	1H	1325	G	O5'-P-OP2	7.07	119.19	110.70
26	1H	1595	G	O5'-P-OP1	-7.07	99.33	105.70
26	1H	1616	A	C5-C6-N6	-7.07	118.04	123.70
26	14	74	A	C4-C5-N7	7.07	114.23	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1249	U	N1-C2-O2	-7.07	117.85	122.80
26	1H	919	G	C4-C5-C6	7.07	123.04	118.80
26	14	1241	A	C2-N3-C4	-7.07	107.07	110.60
26	14	945	A	C5-C6-N6	-7.06	118.05	123.70
26	14	1698	A	N9-C4-C5	-7.06	102.97	105.80
26	14	2335	A	O4'-C1'-N9	7.06	113.85	108.20
26	1H	1943	U	O5'-P-OP2	-7.06	99.34	105.70
1	13	580	U	C5-C6-N1	-7.06	119.17	122.70
1	13	1322	C	C6-N1-C1'	-7.06	112.33	120.80
23	2K	21	U	C2-N3-C4	-7.06	122.76	127.00
26	1H	51	G	N7-C8-N9	-7.06	109.57	113.10
26	1H	577	G	OP1-P-OP2	-7.06	109.01	119.60
26	1H	2451	A	N9-C4-C5	7.06	108.62	105.80
26	14	127	A	OP1-P-O3'	7.06	120.73	105.20
26	14	2074	U	O5'-P-OP1	-7.06	99.35	105.70
26	14	861	A	O5'-P-OP1	-7.06	99.35	105.70
26	1H	1658	C	N3-C4-N4	7.06	122.94	118.00
26	1H	432	A	N9-C4-C5	-7.05	102.98	105.80
26	1H	866	A	N7-C8-N9	7.05	117.33	113.80
26	1H	841	A	C5-C6-N6	-7.05	118.06	123.70
26	1H	2351	G	N3-C4-N9	7.05	130.23	126.00
26	14	2584	U	C6-N1-C1'	-7.05	111.33	121.20
26	1H	121	G	C6-N1-C2	-7.05	120.87	125.10
26	1H	78	A	N9-C4-C5	-7.05	102.98	105.80
26	1H	2056	G	N3-C4-N9	7.05	130.23	126.00
26	1H	1361	G	C8-N9-C4	7.05	109.22	106.40
26	1H	2379	G	N1-C6-O6	7.05	124.13	119.90
26	1H	2607	G	C2-N3-C4	-7.04	108.38	111.90
1	13	1189	C	O5'-P-OP2	7.04	119.15	110.70
26	14	2501	C	C2-N1-C1'	-7.04	111.05	118.80
26	1H	598	G	O5'-P-OP2	-7.04	99.36	105.70
26	1H	2031	A	C2-N3-C4	7.04	114.12	110.60
26	14	130	C	N3-C4-C5	7.04	124.72	121.90
26	14	755	C	N1-C2-O2	-7.04	114.67	118.90
26	1H	2701	C	N3-C2-O2	-7.04	116.97	121.90
26	14	1280	G	OP2-P-O3'	7.04	120.69	105.20
26	1H	1699	G	C8-N9-C4	-7.04	103.58	106.40
1	13	910	C	C5-C6-N1	-7.04	117.48	121.00
26	1H	127	A	O5'-P-OP2	-7.04	99.37	105.70
26	1H	950	G	C2-N3-C4	-7.04	108.38	111.90
26	14	1782	C	N3-C2-O2	7.04	126.83	121.90
26	1H	146	G	C4-C5-N7	7.03	113.61	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2331	G	C8-N9-C4	7.03	109.21	106.40
26	14	2378	A	N1-C6-N6	7.03	122.82	118.60
26	1H	2272	U	O5'-P-OP2	-7.03	99.37	105.70
26	1H	214	G	N3-C4-C5	-7.03	125.08	128.60
26	14	2712(A)	A	C8-N9-C4	7.03	108.61	105.80
26	1H	598	G	C6-C5-N7	-7.03	126.18	130.40
26	1H	2048	G	N9-C4-C5	7.03	108.21	105.40
26	14	2361	A	C2-N3-C4	-7.03	107.09	110.60
26	1H	211	A	C8-N9-C4	7.03	108.61	105.80
27	16	99	A	OP1-P-OP2	7.03	130.14	119.60
1	13	865	A	C4-C5-N7	7.02	114.21	110.70
26	14	1623	G	N9-C4-C5	7.02	108.21	105.40
26	1H	631	A	C5-N7-C8	7.02	107.41	103.90
26	1H	1006	C	O5'-P-OP2	7.02	119.13	110.70
26	14	385	C	OP1-P-OP2	7.02	130.13	119.60
26	1H	1187	G	OP2-P-O3'	7.02	120.64	105.20
26	1H	1602	U	C5-C6-N1	-7.02	119.19	122.70
26	1H	1010	A	C8-N9-C4	7.02	108.61	105.80
26	1H	71	A	C6-C5-N7	-7.02	127.39	132.30
26	1H	1940	U	N3-C4-O4	7.02	124.31	119.40
26	1H	1277	G	C2-N3-C4	-7.02	108.39	111.90
26	1H	2871	C	O5'-P-OP1	7.02	119.12	110.70
1	1G	11	G	O5'-P-OP1	-7.02	99.39	105.70
26	1H	816	C	N3-C4-C5	-7.01	119.09	121.90
26	1H	1402	C	C6-N1-C2	-7.01	117.49	120.30
26	14	752	A	C6-N1-C2	-7.01	114.39	118.60
26	14	2702	U	O5'-P-OP2	-7.01	99.39	105.70
26	14	774	A	C4-C5-N7	7.01	114.21	110.70
26	1H	2199	A	N1-C6-N6	-7.01	114.39	118.60
26	1H	2368	C	O5'-P-OP1	-7.01	99.39	105.70
26	14	2307	G	C4-N9-C1'	7.01	135.61	126.50
26	1H	40	C	O5'-P-OP2	-7.01	99.39	105.70
26	1H	948	G	C4-C5-N7	7.01	113.60	110.80
26	14	1673	U	O5'-P-OP2	7.01	119.11	110.70
26	1H	2497	A	C4-C5-C6	7.01	120.50	117.00
26	14	736	C	N3-C2-O2	7.01	126.80	121.90
26	14	1307	A	C2-N3-C4	-7.01	107.10	110.60
1	13	5	U	P-O3'-C3'	7.00	128.11	119.70
26	1H	1191	G	OP1-P-OP2	7.00	130.11	119.60
26	14	1934	C	O5'-P-OP2	-7.00	99.40	105.70
26	1H	2025	C	C6-N1-C2	-7.00	117.50	120.30
26	14	193	U	C5-C6-N1	-7.00	119.20	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	946	G	O5'-P-OP1	-7.00	99.40	105.70
26	1H	131	G	C5-C6-O6	-7.00	124.40	128.60
26	1H	1604	C	O5'-P-OP1	-7.00	99.40	105.70
26	1H	2703	C	C6-N1-C2	-7.00	117.50	120.30
27	16	65	C	N1-C2-O2	7.00	123.10	118.90
26	14	2518	A	C5-C6-N6	-7.00	118.10	123.70
1	1G	1286	A	N7-C8-N9	7.00	117.30	113.80
26	1H	1681	G	N1-C6-O6	6.99	124.10	119.90
26	1H	52	A	N1-C2-N3	-6.99	125.80	129.30
1	13	1128	C	C2-N1-C1'	6.99	126.49	118.80
26	1H	263	C	O5'-P-OP2	-6.99	99.41	105.70
26	1H	507	A	C5-C6-N6	-6.99	118.11	123.70
26	1H	989	G	C5-C6-O6	-6.99	124.41	128.60
26	1H	1496	A	C6-C5-N7	-6.99	127.41	132.30
26	14	2346	A	O4'-C1'-N9	6.99	113.79	108.20
26	14	1332	G	N3-C2-N2	-6.99	115.01	119.90
26	1H	1312	U	C5-C4-O4	6.99	130.09	125.90
26	1H	1363	C	O5'-P-OP2	-6.99	99.41	105.70
26	14	841	A	C2-N3-C4	-6.99	107.11	110.60
26	1H	686	G	N7-C8-N9	-6.98	109.61	113.10
1	13	1049	U	C2-N1-C1'	-6.98	109.32	117.70
11	2I	102	GLY	N-CA-C	-6.98	95.65	113.10
1	13	47	C	N1-C2-O2	-6.98	114.71	118.90
1	13	982	U	N3-C2-O2	-6.98	117.31	122.20
26	1H	785	G	N3-C2-N2	-6.98	115.02	119.90
26	1H	915	C	N1-C2-O2	6.98	123.09	118.90
26	1H	2259	G	OP1-P-OP2	-6.97	109.14	119.60
26	14	451	C	O5'-P-OP1	-6.97	99.42	105.70
26	14	456	C	N1-C2-O2	-6.97	114.72	118.90
26	1H	1300	U	N1-C2-N3	6.97	119.08	114.90
1	1G	266	G	P-O3'-C3'	6.97	128.07	119.70
26	1H	699	A	N1-C2-N3	-6.97	125.81	129.30
27	16	101	A	C2-N3-C4	-6.97	107.11	110.60
26	1H	265	A	C8-N9-C4	-6.97	103.01	105.80
26	1H	2375	G	N1-C6-O6	6.97	124.08	119.90
26	14	133	C	N3-C4-C5	6.97	124.69	121.90
26	1H	2330	G	C4-C5-N7	6.97	113.59	110.80
26	1H	2689	U	C5-C6-N1	-6.97	119.22	122.70
26	1H	74	A	C4-C5-N7	6.96	114.18	110.70
26	1H	530	G	N3-C4-N9	-6.96	121.82	126.00
26	1H	599	G	N1-C2-N2	-6.96	109.93	116.20
26	14	48	G	OP2-P-O3'	6.96	120.52	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	165	U	C2-N1-C1'	6.96	126.06	117.70
26	14	471	A	C2-N3-C4	-6.96	107.12	110.60
1	13	1435	G	C8-N9-C4	6.96	109.19	106.40
26	1H	784	A	N1-C6-N6	-6.96	114.42	118.60
26	1H	1613	G	C5-C6-O6	6.96	132.78	128.60
26	1H	2506	U	N1-C2-O2	6.96	127.67	122.80
26	14	1698	A	C5-C6-N6	-6.96	118.13	123.70
26	1H	1314	C	O5'-P-OP2	-6.96	99.44	105.70
26	1H	2310	A	N7-C8-N9	6.96	117.28	113.80
26	14	2351	G	N3-C4-C5	-6.96	125.12	128.60
1	13	19	C	C6-N1-C2	-6.96	117.52	120.30
26	1H	2521	C	C5-C6-N1	-6.96	117.52	121.00
26	14	935	C	C5-C6-N1	-6.96	117.52	121.00
26	14	1989	G	N3-C2-N2	-6.96	115.03	119.90
26	1H	528	A	C8-N9-C1'	6.96	140.22	127.70
26	1H	774	A	N7-C8-N9	6.96	117.28	113.80
26	14	265	A	C6-C5-N7	-6.96	127.43	132.30
26	14	2249	U	C6-N1-C2	-6.96	116.83	121.00
26	1H	2513	G	O5'-P-OP2	-6.95	99.44	105.70
1	1G	904	C	N1-C2-O2	-6.95	114.73	118.90
26	1H	1437	C	N3-C2-O2	-6.95	117.03	121.90
26	1H	839	U	OP1-P-OP2	6.95	130.02	119.60
26	1H	432	A	C4-C5-N7	6.95	114.17	110.70
1	1G	266	G	C8-N9-C4	-6.95	103.62	106.40
26	14	1929	G	OP1-P-OP2	6.95	130.02	119.60
26	14	2272	U	N3-C4-O4	-6.95	114.54	119.40
26	14	1954	G	C8-N9-C1'	6.94	136.03	127.00
26	1H	134	C	C5-C6-N1	-6.94	117.53	121.00
26	1H	1624	G	O5'-P-OP2	-6.94	99.45	105.70
26	1H	2585	U	N3-C4-O4	-6.94	114.54	119.40
26	1H	2638	G	N3-C2-N2	6.94	124.76	119.90
26	14	774	A	N1-C6-N6	6.94	122.77	118.60
26	14	2554	U	O5'-P-OP1	-6.94	99.45	105.70
26	1H	239	U	N3-C4-O4	-6.94	114.54	119.40
26	1H	1153	C	O5'-P-OP2	-6.94	99.46	105.70
26	14	1441	G	N7-C8-N9	-6.94	109.63	113.10
26	1H	1296	G	O5'-P-OP2	-6.94	99.46	105.70
26	1H	1598	C	OP1-P-O3'	6.94	120.46	105.20
26	1H	236	C	C5-C6-N1	-6.93	117.53	121.00
26	1H	2006	C	C6-N1-C2	6.93	123.07	120.30
26	14	197	A	C6-N1-C2	-6.93	114.44	118.60
26	14	330	A	C6-C5-N7	-6.93	127.44	132.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	784	A	N1-C6-N6	-6.93	114.44	118.60
26	1H	583	G	OP1-P-O3'	6.93	120.44	105.20
26	1H	1928	A	C5-C6-N1	6.93	121.17	117.70
26	14	2697	G	C2-N3-C4	-6.93	108.44	111.90
26	1H	1229(A)	G	C6-C5-N7	-6.93	126.24	130.40
26	1H	2068	U	C5-C4-O4	6.93	130.06	125.90
26	1H	918	A	O5'-P-OP2	6.92	119.01	110.70
26	1H	2371	G	C5-C6-N1	-6.92	108.04	111.50
26	1H	2827	C	C5-C4-N4	-6.92	115.35	120.20
26	14	776	G	N1-C6-O6	6.92	124.06	119.90
26	1H	613	U	N1-C2-O2	6.92	127.65	122.80
26	1H	208	C	OP2-P-O3'	6.92	120.43	105.20
26	1H	2253	G	O5'-P-OP2	-6.92	99.47	105.70
27	16	7	G	N7-C8-N9	6.92	116.56	113.10
26	1H	1594	G	C8-N9-C4	-6.92	103.63	106.40
26	14	1616	A	C6-C5-N7	-6.92	127.46	132.30
26	1H	1225	C	C5-C6-N1	-6.92	117.54	121.00
26	1H	52	A	O5'-P-OP2	-6.92	99.48	105.70
26	1H	1834	U	N3-C2-O2	-6.92	117.36	122.20
26	1H	2277	G	C4-C5-N7	-6.92	108.03	110.80
26	14	807	U	N3-C4-O4	6.92	124.24	119.40
26	1H	774	A	C4-N9-C1'	-6.92	113.85	126.30
26	14	2272	U	N3-C2-O2	-6.92	117.36	122.20
1	13	974	A	C4-C5-N7	6.91	114.16	110.70
26	1H	786	C	C5-C6-N1	-6.91	117.54	121.00
26	14	140	A	C2-N3-C4	-6.91	107.14	110.60
26	1H	1336	A	N1-C6-N6	-6.91	114.45	118.60
26	14	2278	A	N1-C6-N6	-6.91	114.45	118.60
26	1H	1241	A	C6-C5-N7	-6.91	127.46	132.30
26	1H	1673	U	C5-C6-N1	-6.91	119.25	122.70
26	14	2707	G	C5-C6-N1	6.91	114.95	111.50
26	1H	1275	A	N1-C6-N6	6.91	122.74	118.60
26	1H	1332	G	C4-C5-N7	6.91	113.56	110.80
26	1H	1237	A	OP1-P-OP2	-6.90	109.24	119.60
26	1H	2032	G	N1-C2-N3	6.90	128.04	123.90
26	14	1482	U	C5-C4-O4	6.90	130.04	125.90
26	1H	2697	G	OP1-P-OP2	6.90	129.95	119.60
1	13	1525	G	N3-C4-C5	6.90	132.05	128.60
26	14	676	A	C4-C5-N7	6.90	114.15	110.70
26	1H	671	C	C2-N1-C1'	-6.90	111.21	118.80
26	1H	645	C	C5-C6-N1	6.90	124.45	121.00
26	14	2464	C	C6-N1-C2	6.90	123.06	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	827	U	N3-C2-O2	-6.89	117.37	122.20
26	14	512	G	O4'-C1'-N9	6.89	113.72	108.20
26	14	1616	A	O4'-C1'-N9	6.89	113.72	108.20
26	1H	122	G	C6-C5-N7	-6.89	126.26	130.40
26	1H	284	U	O5'-P-OP1	-6.89	99.50	105.70
39	A8	24	LEU	CA-CB-CG	6.89	131.15	115.30
52	N8	41	PRO	C-N-CD	-6.89	105.44	120.60
26	1H	966	G	C8-N9-C4	6.89	109.16	106.40
26	1H	1353	A	N1-C2-N3	6.89	132.75	129.30
26	1H	468	G	C5-C6-O6	-6.89	124.47	128.60
26	1H	1249	U	N3-C2-O2	6.89	127.02	122.20
26	14	735	A	N7-C8-N9	-6.89	110.36	113.80
26	14	787	U	O5'-P-OP1	6.89	118.97	110.70
26	1H	2713	A	C5-C6-N1	-6.89	114.26	117.70
27	16	5	C	C2-N3-C4	-6.89	116.46	119.90
26	1H	1340	U	C5-C4-O4	-6.88	121.77	125.90
26	14	2286	A	N7-C8-N9	6.88	117.24	113.80
27	16	24	G	N3-C4-C5	-6.88	125.16	128.60
26	1H	1303	G	O5'-P-OP2	-6.88	99.51	105.70
26	14	71	A	C5-C6-N6	-6.88	118.20	123.70
1	13	1395	C	C6-N1-C2	-6.88	117.55	120.30
26	1H	982	C	O5'-P-OP2	-6.88	99.51	105.70
1	13	896	C	C5-C6-N1	-6.87	117.56	121.00
26	1H	265	A	O4'-C1'-N9	6.87	113.70	108.20
26	1H	470	A	C4-C5-N7	6.87	114.14	110.70
26	1H	1299	G	C5-C6-O6	-6.87	124.48	128.60
26	1H	1626	G	N3-C2-N2	-6.87	115.09	119.90
26	1H	2287	A	N1-C2-N3	6.87	132.74	129.30
26	14	991	C	C6-N1-C2	-6.87	117.55	120.30
26	14	2326	C	N3-C4-C5	-6.87	119.15	121.90
1	13	586	C	C6-N1-C2	6.87	123.05	120.30
26	1H	802	A	N1-C6-N6	6.87	122.72	118.60
26	1H	1428	C	N3-C4-N4	-6.87	113.19	118.00
26	14	1332	G	OP1-P-O3'	6.87	120.32	105.20
26	14	1304	C	N1-C2-O2	6.87	123.02	118.90
26	14	2463	C	N3-C2-O2	6.87	126.71	121.90
26	1H	596	G	N1-C6-O6	6.87	124.02	119.90
26	1H	1931	U	C2-N3-C4	-6.87	122.88	127.00
1	1G	1260	C	C5-C6-N1	6.87	124.43	121.00
26	14	1314	C	C5-C6-N1	6.87	124.43	121.00
26	1H	246	C	C5-C6-N1	-6.86	117.57	121.00
26	1H	684	G	N9-C4-C5	6.86	108.14	105.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1604	C	O5'-P-OP1	-6.86	99.52	105.70
32	59	153	LYS	C-N-CD	6.86	142.81	128.40
26	1H	946	G	N7-C8-N9	-6.86	109.67	113.10
26	1H	2444	G	N3-C2-N2	-6.86	115.10	119.90
26	14	1760	A	C8-N9-C4	6.86	108.54	105.80
1	1G	1200	C	C2-N1-C1'	6.86	126.34	118.80
26	14	1276	A	N9-C4-C5	-6.86	103.06	105.80
26	14	1312	U	O5'-P-OP1	-6.86	99.53	105.70
26	1H	71	A	N3-C4-C5	6.86	131.60	126.80
26	1H	838	C	C4-C5-C6	6.86	120.83	117.40
26	14	965	C	C6-N1-C2	-6.86	117.56	120.30
26	14	1786	A	N3-C4-C5	6.86	131.60	126.80
26	1H	330	A	N1-C6-N6	6.85	122.71	118.60
26	1H	1247	A	C6-N1-C2	-6.85	114.49	118.60
26	14	1261	C	N3-C4-C5	6.85	124.64	121.90
36	35	65	ARG	N-CA-C	-6.85	92.50	111.00
1	13	899	C	N3-C2-O2	6.85	126.69	121.90
26	1H	449	A	N1-C6-N6	-6.85	114.49	118.60
26	1H	2501	C	C6-N1-C2	6.85	123.04	120.30
1	1G	690	G	N7-C8-N9	6.85	116.52	113.10
1	13	12	U	C6-N1-C2	-6.85	116.89	121.00
26	1H	1213	A	N1-C6-N6	6.85	122.71	118.60
26	1H	1372	U	C4-C5-C6	6.85	123.81	119.70
26	14	574	C	C2-N1-C1'	-6.85	111.27	118.80
26	14	2307	G	N7-C8-N9	6.85	116.52	113.10
1	1G	884	U	N3-C2-O2	-6.84	117.41	122.20
26	14	1386	C	N3-C4-C5	-6.84	119.16	121.90
1	1G	913	A	P-O3'-C3'	6.84	127.91	119.70
26	14	855	G	C8-N9-C4	-6.84	103.66	106.40
26	1H	335	C	C5-C6-N1	6.84	124.42	121.00
26	1H	486	C	O5'-P-OP2	6.84	118.91	110.70
26	1H	793	A	O5'-P-OP2	-6.84	99.54	105.70
26	1H	942	G	N3-C2-N2	-6.84	115.11	119.90
26	1H	2447	G	C4-C5-N7	-6.84	108.06	110.80
26	14	990	A	O5'-P-OP1	-6.84	99.55	105.70
27	16	60	C	C6-N1-C2	-6.84	117.56	120.30
26	1H	130	C	C2-N3-C4	-6.84	116.48	119.90
26	1H	593	G	O5'-P-OP2	-6.84	99.55	105.70
26	1H	624	C	O5'-P-OP2	6.84	118.90	110.70
26	1H	946	G	C8-N9-C4	6.84	109.14	106.40
26	1H	1789	A	C6-N1-C2	-6.84	114.50	118.60
27	16	47	C	C5-C6-N1	-6.84	117.58	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	188	G	N1-C6-O6	6.84	124.00	119.90
26	14	1318	C	C6-N1-C2	-6.84	117.56	120.30
26	14	2701	C	P-O3'-C3'	6.84	127.90	119.70
26	14	127	A	C5-C6-N1	6.83	121.12	117.70
1	13	813	U	C4-C5-C6	-6.83	115.60	119.70
26	1H	598	G	C4-C5-C6	6.83	122.90	118.80
26	1H	1215	G	C2-N3-C4	6.83	115.32	111.90
26	1H	1648	C	N1-C2-O2	-6.83	114.80	118.90
1	1G	121	C	N1-C2-O2	6.83	123.00	118.90
26	14	2281	C	C6-N1-C2	-6.83	117.57	120.30
26	1H	34	C	O5'-P-OP2	6.83	118.90	110.70
26	14	530	G	N1-C6-O6	6.83	124.00	119.90
26	1H	834	C	OP2-P-O3'	6.83	120.21	105.20
26	14	56	A	N1-C6-N6	-6.83	114.50	118.60
26	14	621	A	N7-C8-N9	6.83	117.21	113.80
26	1H	2286	A	N7-C8-N9	6.82	117.21	113.80
26	1H	2246	G	C5-N7-C8	6.82	107.71	104.30
26	1H	2342	C	C5-C6-N1	6.82	124.41	121.00
26	14	2307	G	C8-N9-C4	-6.82	103.67	106.40
1	13	115	G	N3-C4-C5	-6.82	125.19	128.60
26	1H	2708	G	O5'-P-OP2	-6.82	99.56	105.70
26	14	916	G	O5'-P-OP2	6.82	118.88	110.70
26	1H	398	G	O5'-P-OP2	-6.82	99.57	105.70
26	1H	761	A	O5'-P-OP2	-6.82	99.57	105.70
26	1H	1022	G	N3-C2-N2	-6.82	115.13	119.90
26	1H	1754	C	N3-C4-C5	-6.82	119.17	121.90
26	1H	2502	G	O5'-P-OP1	-6.82	99.57	105.70
26	1H	2830	G	N7-C8-N9	6.82	116.51	113.10
26	1H	391	G	N1-C2-N3	6.81	127.99	123.90
1	1G	862	C	N3-C2-O2	-6.81	117.13	121.90
26	14	2359	C	N3-C4-N4	-6.81	113.23	118.00
26	14	2607	G	N3-C4-N9	6.81	130.09	126.00
26	1H	1410	G	C4-N9-C1'	-6.81	117.65	126.50
43	E8	51	LEU	CA-CB-CG	6.81	130.97	115.30
26	14	848	G	C8-N9-C4	-6.81	103.68	106.40
26	1H	1266	G	C5-C6-O6	-6.81	124.52	128.60
26	14	71	A	O4'-C1'-N9	-6.81	102.75	108.20
26	14	265	A	N1-C6-N6	6.81	122.69	118.60
26	14	1782	C	N3-C4-N4	6.81	122.77	118.00
26	14	2314	C	C6-N1-C2	-6.81	117.58	120.30
26	14	2873	A	C5-C6-N1	-6.81	114.30	117.70
27	1J	14	U	O5'-P-OP2	-6.81	99.57	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1302	A	OP1-P-OP2	6.81	129.81	119.60
1	13	1322	C	N1-C2-O2	6.81	122.98	118.90
1	13	507	C	C6-N1-C2	-6.80	117.58	120.30
26	1H	2070	G	O5'-P-OP1	6.80	118.86	110.70
1	13	897	C	C5-C6-N1	-6.80	117.60	121.00
26	1H	1322	A	C5-N7-C8	-6.80	100.50	103.90
26	1H	2638	G	N3-C4-N9	6.80	130.08	126.00
1	1G	534	U	N3-C2-O2	-6.80	117.44	122.20
26	14	2005	A	O5'-P-OP2	-6.80	99.58	105.70
26	14	2685	G	N1-C6-O6	6.80	123.98	119.90
24	3K	35	A	O5'-P-OP2	-6.80	99.58	105.70
26	1H	1971	A	C6-N1-C2	-6.80	114.52	118.60
26	14	1616	A	N3-C4-C5	6.80	131.56	126.80
26	14	1496	A	C6-C5-N7	-6.79	127.54	132.30
1	13	1158	C	N3-C2-O2	-6.79	117.14	121.90
26	1H	845	G	P-O3'-C3'	6.79	127.85	119.70
26	14	676	A	OP1-P-OP2	6.79	129.79	119.60
26	14	953	A	N1-C6-N6	6.79	122.68	118.60
26	14	1251	C	OP1-P-OP2	6.79	129.79	119.60
26	14	2606	C	OP1-P-OP2	6.79	129.79	119.60
26	1H	1021	A	C8-N9-C4	-6.79	103.08	105.80
26	14	2059	A	O4'-C1'-N9	6.79	113.63	108.20
26	1H	1201	C	C5-C4-N4	-6.79	115.45	120.20
26	14	954	G	C4-C5-N7	-6.79	108.08	110.80
26	14	1379	A	N9-C1'-C2'	6.79	122.83	114.00
1	13	422	C	P-O3'-C3'	6.79	127.84	119.70
26	1H	682	G	C8-N9-C4	6.79	109.11	106.40
26	14	2332	U	N3-C4-O4	-6.79	114.65	119.40
26	14	2061	G	O5'-P-OP2	-6.79	99.59	105.70
1	13	496	A	C8-N9-C4	-6.78	103.09	105.80
26	1H	1817	G	N7-C8-N9	-6.78	109.71	113.10
26	1H	375	C	O5'-P-OP2	-6.78	99.60	105.70
26	1H	1444	G	N1-C6-O6	-6.78	115.83	119.90
45	C5	103	GLY	N-CA-C	6.78	130.06	113.10
26	1H	940	G	C6-N1-C2	-6.78	121.03	125.10
26	1H	1558	A	C5-C6-N6	-6.78	118.28	123.70
1	1G	885	G	C8-N9-C4	6.78	109.11	106.40
26	14	684	G	C8-N9-C4	-6.78	103.69	106.40
26	1H	508	G	C5-N7-C8	-6.78	100.91	104.30
26	1H	598	G	N1-C6-O6	6.78	123.97	119.90
26	1H	684	G	C8-N9-C4	-6.78	103.69	106.40
26	1H	1699	G	O4'-C1'-N9	6.78	113.62	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	798	G	C5-C6-O6	-6.78	124.53	128.60
26	14	1241	A	C5-C6-N1	-6.78	114.31	117.70
1	13	1505	G	OP1-P-OP2	-6.77	109.44	119.60
26	1H	199	A	C4-C5-C6	-6.77	113.61	117.00
26	14	945	A	C5-C6-N1	-6.77	114.31	117.70
26	1H	267	C	N3-C4-C5	6.77	124.61	121.90
26	14	209	C	N3-C4-C5	6.77	124.61	121.90
26	1H	2591	C	O5'-P-OP2	6.77	118.82	110.70
26	14	1623	G	C8-N9-C1'	6.77	135.80	127.00
26	1H	142	G	N3-C4-N9	-6.77	121.94	126.00
26	1H	232	G	O5'-P-OP2	-6.77	99.61	105.70
5	42	31	LEU	CA-CB-CG	6.77	130.86	115.30
26	1H	2466	C	C6-N1-C2	6.77	123.01	120.30
26	14	775	G	N3-C4-C5	-6.77	125.22	128.60
26	1H	448	U	C5-C6-N1	-6.76	119.32	122.70
1	1G	697	U	O5'-P-OP2	-6.76	99.61	105.70
26	1H	1780	A	N1-C6-N6	-6.76	114.54	118.60
26	1H	2439	A	N1-C6-N6	6.76	122.66	118.60
26	1H	676	A	C6-N1-C2	6.76	122.66	118.60
26	14	1349	A	C5-N7-C8	-6.76	100.52	103.90
1	13	328	C	N1-C2-O2	6.76	122.95	118.90
26	1H	1299	G	C4-C5-N7	6.76	113.50	110.80
26	1H	1621	U	N3-C4-O4	6.76	124.13	119.40
26	1H	2441	C	N3-C4-N4	-6.76	113.27	118.00
26	14	2429	G	O5'-P-OP1	6.76	118.81	110.70
26	1H	465	G	C5-C6-N1	-6.75	108.12	111.50
26	1H	931	G	N3-C4-C5	-6.75	125.22	128.60
26	1H	1162	G	N7-C8-N9	6.75	116.48	113.10
1	1G	1314	C	C6-N1-C2	-6.75	117.60	120.30
26	14	786	C	N3-C4-C5	6.75	124.60	121.90
26	14	2681	C	N3-C2-O2	-6.75	117.17	121.90
26	1H	452	G	C2-N3-C4	6.75	115.28	111.90
26	1H	2856	C	C6-N1-C2	-6.75	117.60	120.30
26	14	2329	G	C5-C6-N1	6.75	114.88	111.50
26	1H	828	U	N3-C4-O4	-6.75	114.67	119.40
26	1H	1924	C	N3-C2-O2	-6.75	117.17	121.90
26	14	1600	C	C6-N1-C2	6.75	123.00	120.30
26	1H	1157	G	N1-C6-O6	6.75	123.95	119.90
26	14	1397	U	C5-C4-O4	6.75	129.95	125.90
26	1H	2246	G	C4-C5-N7	-6.75	108.10	110.80
27	16	44	G	C6-C5-N7	6.75	134.45	130.40
26	1H	673	C	OP1-P-OP2	-6.75	109.48	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1972	A	OP1-P-OP2	-6.75	109.48	119.60
26	1H	1249	U	C2-N3-C4	-6.75	122.95	127.00
26	1H	210	C	N3-C4-C5	6.74	124.60	121.90
26	1H	1849	G	O5'-P-OP1	-6.74	99.63	105.70
26	14	768	G	OP1-P-OP2	6.74	129.72	119.60
26	14	79	G	C5-C6-O6	-6.74	124.56	128.60
26	1H	1673	U	C2-N1-C1'	-6.74	109.61	117.70
26	14	467	G	O5'-P-OP2	-6.74	99.64	105.70
26	1H	1225	C	OP1-P-OP2	6.74	129.71	119.60
26	1H	2435	A	N7-C8-N9	6.73	117.17	113.80
26	1H	94	G	C5-C6-O6	-6.73	124.56	128.60
26	1H	1607	C	C2-N1-C1'	6.73	126.21	118.80
26	1H	2433	A	C4-C5-C6	6.73	120.37	117.00
26	1H	2083	G	C2-N3-C4	-6.73	108.54	111.90
26	1H	675	A	C5-N7-C8	-6.73	100.54	103.90
26	1H	1257	C	N1-C2-N3	6.73	123.91	119.20
26	1H	1940	U	N1-C2-N3	6.73	118.94	114.90
26	1H	2620	C	N1-C2-O2	-6.73	114.86	118.90
26	14	827	U	N1-C2-O2	-6.73	118.09	122.80
1	13	912	C	C6-N1-C2	6.73	122.99	120.30
14	5I	27	CYS	CA-CB-SG	-6.73	101.89	114.00
26	1H	978	G	OP1-P-O3'	6.73	120.00	105.20
26	1H	205	G	C8-N9-C4	6.72	109.09	106.40
26	1H	2283	C	O5'-P-OP2	-6.72	99.65	105.70
26	14	1366	A	C5-C6-N6	-6.72	118.32	123.70
26	14	1696	G	O5'-P-OP2	-6.72	99.65	105.70
26	1H	1888	G	N3-C4-C5	-6.72	125.24	128.60
26	1H	2441	C	C5-C4-N4	6.72	124.91	120.20
26	1H	2779	U	C2-N3-C4	-6.72	122.97	127.00
26	14	1496	A	C8-N9-C4	-6.72	103.11	105.80
26	14	1826	G	N7-C8-N9	-6.72	109.74	113.10
23	2K	21	U	N3-C2-O2	-6.72	117.50	122.20
26	1H	2298	A	O5'-P-OP2	-6.72	99.65	105.70
26	14	530	G	C4-C5-N7	6.72	113.49	110.80
26	1H	1337	G	OP1-P-O3'	6.72	119.98	105.20
27	16	87	G	C8-N9-C4	6.72	109.09	106.40
26	1H	971	C	N3-C4-C5	-6.72	119.21	121.90
26	14	802	A	C6-N1-C2	-6.72	114.57	118.60
26	14	2029	G	O5'-P-OP1	-6.72	99.66	105.70
26	1H	778	G	C2-N3-C4	-6.71	108.54	111.90
26	1H	2690	C	N3-C4-C5	-6.71	119.21	121.90
26	14	2217	G	C4-C5-C6	6.71	122.83	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2324	C	N3-C4-C5	6.71	124.59	121.90
26	14	2042	A	C8-N9-C4	6.71	108.48	105.80
26	1H	862	G	O5'-P-OP2	-6.71	99.66	105.70
26	1H	2296	U	N3-C4-O4	6.71	124.10	119.40
26	1H	2709	G	O5'-P-OP1	6.71	118.75	110.70
26	1H	2011	U	N3-C2-O2	6.71	126.90	122.20
26	1H	2082	A	C6-N1-C2	-6.71	114.57	118.60
26	14	566	U	C5-C6-N1	-6.71	119.35	122.70
26	14	2512	C	C6-N1-C2	6.71	122.98	120.30
26	1H	1604	C	O5'-P-OP2	6.71	118.75	110.70
55	Q8	28	GLY	N-CA-C	6.70	129.86	113.10
26	1H	2363	C	C6-N1-C2	6.70	122.98	120.30
26	14	2490	G	C8-N9-C4	-6.70	103.72	106.40
26	1H	2502	G	C8-N9-C4	-6.70	103.72	106.40
26	14	1694	C	N1-C2-O2	6.70	122.92	118.90
26	1H	146	G	N9-C4-C5	-6.70	102.72	105.40
26	1H	188	G	C4-C5-N7	6.70	113.48	110.80
26	1H	415	A	N1-C6-N6	6.70	122.62	118.60
26	1H	1347	G	N1-C6-O6	6.70	123.92	119.90
26	1H	1356	G	N1-C6-O6	6.70	123.92	119.90
1	1G	1518	A	O5'-P-OP1	-6.70	99.67	105.70
26	14	1141	U	P-O3'-C3'	6.70	127.74	119.70
26	14	1625	C	O5'-P-OP2	-6.70	99.67	105.70
26	14	2702	U	N1-C2-O2	6.70	127.49	122.80
26	1H	732	C	N1-C2-O2	-6.70	114.88	118.90
26	1H	777	A	N1-C2-N3	6.70	132.65	129.30
26	1H	1006	C	C6-N1-C2	6.70	122.98	120.30
27	16	101	A	C8-N9-C4	6.70	108.48	105.80
26	14	2593	U	OP1-P-OP2	-6.70	109.56	119.60
1	1G	894	G	C4-C5-N7	6.69	113.48	110.80
26	1H	609	A	C8-N9-C4	6.69	108.48	105.80
26	1H	1348	G	OP1-P-O3'	6.69	119.92	105.20
26	1H	2823	A	C8-N9-C4	-6.69	103.12	105.80
1	13	932	C	C6-N1-C2	-6.69	117.62	120.30
1	13	974	A	O4'-C1'-N9	6.69	113.55	108.20
26	1H	2541	A	O5'-P-OP1	-6.69	99.68	105.70
26	14	918	A	O5'-P-OP1	-6.69	99.68	105.70
26	1H	1253	A	N7-C8-N9	-6.69	110.46	113.80
26	1H	2243	U	OP2-P-O3'	6.69	119.91	105.20
26	14	2286	A	C6-C5-N7	-6.69	127.62	132.30
26	14	2591	C	O5'-P-OP2	-6.69	99.68	105.70
26	1H	1619	G	N3-C4-C5	-6.69	125.26	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	51	G	N9-C4-C5	-6.68	102.73	105.40
26	1H	729	G	C8-N9-C4	-6.68	103.73	106.40
1	1G	32	A	C8-N9-C4	-6.68	103.13	105.80
26	14	639	U	C5-C4-O4	6.68	129.91	125.90
26	14	681	G	C8-N9-C1'	-6.68	118.31	127.00
26	14	2547	U	O5'-P-OP1	6.68	118.72	110.70
26	1H	1819	A	O5'-P-OP1	-6.68	99.69	105.70
26	1H	2782	G	N1-C6-O6	6.68	123.91	119.90
26	14	1571	A	C5-C6-N1	6.68	121.04	117.70
1	13	582	U	N3-C4-C5	6.68	118.61	114.60
26	14	2585	U	C2-N1-C1'	6.68	125.72	117.70
26	1H	371	A	C5-C6-N6	-6.68	118.36	123.70
26	1H	1159	U	N3-C2-O2	-6.68	117.53	122.20
26	14	134	C	C6-N1-C2	6.68	122.97	120.30
26	1H	2485	G	C2-N3-C4	-6.68	108.56	111.90
26	1H	2620	C	C6-N1-C2	6.68	122.97	120.30
22	1K	74	C	C6-N1-C1'	-6.68	112.79	120.80
26	1H	72	U	O5'-P-OP2	-6.67	99.69	105.70
26	1H	928	G	N3-C4-C5	6.67	131.94	128.60
26	1H	1759	A	O5'-P-OP1	-6.67	99.69	105.70
28	11	111	LEU	CA-CB-CG	6.67	130.65	115.30
26	14	1815	A	OP1-P-O3'	6.67	119.88	105.20
26	1H	1378	A	O5'-P-OP1	-6.67	99.69	105.70
26	1H	2566	A	C8-N9-C4	-6.67	103.13	105.80
26	14	2087	G	N7-C8-N9	-6.67	109.76	113.10
1	13	892	A	N1-C6-N6	6.67	122.60	118.60
26	1H	680	G	C5-C6-O6	-6.67	124.60	128.60
26	14	2495	G	C5-C6-N1	-6.67	108.17	111.50
26	1H	1528	A	C5-N7-C8	-6.67	100.56	103.90
26	1H	62	C	C5-C6-N1	-6.67	117.67	121.00
26	1H	396	G	N1-C6-O6	6.67	123.90	119.90
26	14	2267	A	OP1-P-OP2	6.67	129.60	119.60
26	1H	2330	G	C6-C5-N7	-6.67	126.40	130.40
26	1H	2665	A	N1-C2-N3	6.67	132.63	129.30
26	1H	2672	G	C8-N9-C4	-6.67	103.73	106.40
26	1H	2595	G	N9-C4-C5	-6.67	102.73	105.40
26	14	1344	G	N1-C6-O6	6.67	123.90	119.90
26	1H	470	A	C5-N7-C8	-6.66	100.57	103.90
26	1H	691	C	N3-C2-O2	6.66	126.56	121.90
26	1H	1385	G	N3-C4-N9	-6.66	122.00	126.00
26	1H	2373	G	N1-C2-N3	6.66	127.90	123.90
26	1H	2775	A	N9-C4-C5	-6.66	103.13	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1955	U	N1-C2-N3	6.66	118.90	114.90
26	14	2518	A	N7-C8-N9	6.66	117.13	113.80
1	13	1322	C	O5'-P-OP2	-6.66	99.70	105.70
26	1H	2503	A	C2-N3-C4	6.66	113.93	110.60
26	1H	462	C	O5'-P-OP2	-6.66	99.71	105.70
26	1H	528	A	C5-N7-C8	-6.66	100.57	103.90
26	1H	763	G	N3-C4-N9	6.66	130.00	126.00
26	1H	2275	C	N3-C4-C5	-6.66	119.24	121.90
26	1H	205	G	N9-C4-C5	-6.66	102.74	105.40
26	1H	2409	G	C5-C6-O6	-6.66	124.61	128.60
26	14	330	A	N1-C2-N3	6.66	132.63	129.30
26	1H	2276	G	N3-C2-N2	-6.65	115.24	119.90
26	14	2286	A	N1-C6-N6	6.65	122.59	118.60
26	1H	814	C	N3-C4-N4	-6.65	113.34	118.00
26	1H	1622	G	C4-C5-N7	-6.65	108.14	110.80
26	1H	1665	A	N1-C2-N3	6.65	132.63	129.30
26	14	2554	U	N1-C2-O2	-6.65	118.14	122.80
26	14	2712	U	C4-C5-C6	6.65	123.69	119.70
26	1H	627	A	C8-N9-C4	6.65	108.46	105.80
26	1H	1196	C	N3-C4-N4	6.65	122.66	118.00
26	1H	2342	C	C6-N1-C2	-6.65	117.64	120.30
26	14	502	A	N1-C2-N3	6.65	132.62	129.30
26	1H	610	C	C2-N3-C4	-6.65	116.58	119.90
26	1H	1429	G	O5'-P-OP2	-6.65	99.72	105.70
1	13	961	U	O5'-P-OP2	-6.65	99.72	105.70
26	1H	1326	U	N3-C2-O2	-6.65	117.55	122.20
26	14	1471	A	N7-C8-N9	6.65	117.12	113.80
24	3K	71	G	C8-N9-C1'	6.64	135.64	127.00
26	14	307	G	OP1-P-OP2	6.64	129.57	119.60
26	1H	2379	G	N9-C4-C5	-6.64	102.74	105.40
26	14	2374	C	N3-C4-N4	-6.64	113.35	118.00
26	14	2422	A	O5'-P-OP2	-6.64	99.72	105.70
26	1H	1193	G	O5'-P-OP1	6.64	118.67	110.70
26	14	681	G	N1-C6-O6	6.64	123.89	119.90
26	1H	1471	A	N1-C6-N6	6.64	122.58	118.60
26	14	2030	A	O5'-P-OP2	-6.64	99.72	105.70
1	13	548	G	O5'-P-OP2	-6.64	99.73	105.70
26	1H	639	U	C5-C4-O4	6.64	129.88	125.90
26	1H	1786	A	N9-C1'-C2'	6.64	122.63	114.00
26	1H	59	U	N3-C4-C5	-6.63	110.62	114.60
26	14	475	U	C6-N1-C2	-6.63	117.02	121.00
26	14	1599	C	O5'-P-OP1	-6.63	99.73	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	901	A	C2-N3-C4	-6.63	107.28	110.60
26	1H	1970	A	C2-N3-C4	6.63	113.92	110.60
26	14	775	G	N3-C4-N9	6.63	129.98	126.00
1	13	1517	G	O5'-P-OP2	-6.63	99.73	105.70
26	1H	463	G	N3-C2-N2	6.63	124.54	119.90
26	1H	580	C	N3-C4-C5	-6.63	119.25	121.90
26	1H	1649	G	N1-C6-O6	-6.63	115.92	119.90
26	1H	2265	U	N3-C4-O4	6.63	124.04	119.40
26	14	1253	A	C4-C5-N7	6.63	114.02	110.70
26	14	2730	C	C6-N1-C2	-6.63	117.65	120.30
26	1H	1124	C	C2-N3-C4	-6.63	116.59	119.90
26	1H	1391	U	C5-C6-N1	6.63	126.02	122.70
26	1H	245	G	C4-C5-C6	6.63	122.78	118.80
26	1H	1027	A	C2-N3-C4	-6.63	107.29	110.60
26	1H	2324	C	O5'-P-OP2	-6.63	99.74	105.70
26	1H	2449	U	O5'-P-OP1	-6.63	99.73	105.70
26	14	2866	U	O5'-P-OP2	-6.63	99.74	105.70
1	13	314	C	O5'-P-OP2	-6.62	99.74	105.70
1	13	1113	C	C6-N1-C2	-6.62	117.65	120.30
26	1H	657	U	C5-C6-N1	-6.62	119.39	122.70
26	1H	1558	A	C6-C5-N7	-6.62	127.66	132.30
26	1H	2084	C	C6-N1-C2	6.62	122.95	120.30
26	1H	2261	C	OP1-P-O3'	-6.62	90.63	105.20
26	14	186	G	N1-C2-N3	6.62	127.87	123.90
26	14	775	G	N1-C2-N2	-6.62	110.24	116.20
26	14	945	A	N7-C8-N9	6.62	117.11	113.80
26	14	1272	A	C5-C6-N6	-6.62	118.40	123.70
26	1H	179	G	N1-C6-O6	6.62	123.87	119.90
26	1H	943	U	O5'-P-OP1	-6.62	99.74	105.70
26	1H	1641	A	O5'-P-OP2	6.62	118.65	110.70
27	16	29	A	OP1-P-OP2	-6.62	109.67	119.60
26	14	581	C	C6-N1-C2	-6.62	117.65	120.30
26	14	779	U	N3-C4-C5	6.62	118.57	114.60
26	14	2598	A	N7-C8-N9	-6.62	110.49	113.80
1	13	800	G	N1-C6-O6	6.62	123.87	119.90
26	1H	146	G	C5-C6-O6	-6.62	124.63	128.60
26	1H	2409	G	C4-C5-N7	6.62	113.45	110.80
26	14	693	C	C5-C6-N1	-6.62	117.69	121.00
26	14	1337	G	OP1-P-O3'	6.62	119.76	105.20
26	1H	1158	C	C5-C6-N1	-6.62	117.69	121.00
26	1H	2708	G	N3-C4-N9	6.62	129.97	126.00
26	1H	528	A	C2-N3-C4	-6.62	107.29	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1395	A	O4'-C1'-N9	6.62	113.49	108.20
26	1H	2326	C	N1-C2-O2	6.62	122.87	118.90
26	1H	2419	U	OP1-P-O3'	6.62	119.75	105.20
26	1H	2491	U	N3-C2-O2	6.62	126.83	122.20
26	14	1437	C	C6-N1-C2	-6.62	117.65	120.30
26	1H	1324	G	N3-C2-N2	-6.61	115.27	119.90
26	1H	1516	U	C5-C4-O4	6.61	129.87	125.90
1	13	577	G	C5-C6-O6	-6.61	124.63	128.60
1	13	770	C	N3-C2-O2	6.61	126.53	121.90
1	1G	558	G	C5-C6-N1	-6.61	108.19	111.50
24	3L	76	A	O4'-C1'-N9	6.61	113.49	108.20
26	1H	1857	G	C6-C5-N7	-6.61	126.43	130.40
26	1H	1902	C	C4-C5-C6	6.61	120.70	117.40
26	1H	2585	U	C5-C6-N1	-6.61	119.39	122.70
26	14	784	A	N9-C4-C5	6.61	108.44	105.80
26	14	1463	C	C6-N1-C2	-6.61	117.66	120.30
26	1H	528	A	O4'-C1'-N9	-6.61	102.91	108.20
26	1H	378	C	N3-C4-N4	6.61	122.62	118.00
26	1H	782	A	C5-C6-N6	6.61	128.99	123.70
26	14	1678	G	N7-C8-N9	6.61	116.40	113.10
26	1H	71	A	N1-C6-N6	6.61	122.56	118.60
26	1H	99	U	N1-C2-O2	6.61	127.42	122.80
26	1H	2712(A)	A	C8-N9-C4	6.61	108.44	105.80
26	14	372	G	O4'-C1'-N9	6.60	113.48	108.20
26	14	1394	U	O5'-P-OP1	-6.60	99.76	105.70
26	14	2440	C	N3-C4-C5	-6.60	119.26	121.90
1	1G	1490	C	O5'-P-OP2	-6.60	99.76	105.70
26	1H	2036	C	OP2-P-O3'	6.60	119.72	105.20
26	1H	2084	C	C5-C6-N1	-6.60	117.70	121.00
26	14	852	G	N1-C6-O6	-6.60	115.94	119.90
26	1H	2439	A	OP2-P-O3'	-6.60	90.68	105.20
26	14	1210	A	C5-N7-C8	-6.60	100.60	103.90
36	35	85	LEU	CA-CB-CG	6.60	130.47	115.30
26	1H	305	U	C6-N1-C2	-6.60	117.04	121.00
1	13	1151	A	O4'-C1'-N9	6.59	113.47	108.20
23	2K	21	U	C6-N1-C1'	-6.59	111.97	121.20
26	1H	811	U	C5-C6-N1	-6.59	119.40	122.70
1	1G	631	G	C8-N9-C4	-6.59	103.76	106.40
26	14	330	A	C5-C6-N1	-6.59	114.40	117.70
26	14	1021	A	C2-N3-C4	-6.59	107.30	110.60
23	2K	6	G	C5-C6-O6	-6.59	124.65	128.60
26	1H	410	G	O5'-P-OP2	6.59	118.61	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	910	A	C4-C5-C6	6.59	120.30	117.00
26	1H	1433	U	O5'-P-OP2	-6.59	99.77	105.70
55	Q8	59	LYS	CD-CE-NZ	6.59	126.86	111.70
26	1H	675	A	O5'-P-OP2	-6.59	99.77	105.70
26	1H	2468	G	C6-C5-N7	-6.59	126.45	130.40
26	14	1968	G	C4-C5-N7	6.59	113.44	110.80
54	L5	34	ARG	NE-CZ-NH2	6.59	123.59	120.30
26	1H	2331	G	C2-N3-C4	-6.59	108.61	111.90
26	1H	2639	A	N1-C6-N6	6.59	122.55	118.60
26	14	363(F)	A	C8-N9-C4	6.59	108.43	105.80
26	14	475	U	C2-N1-C1'	6.59	125.61	117.70
26	14	698	C	O5'-P-OP1	-6.58	99.77	105.70
26	14	921	G	C8-N9-C4	-6.58	103.77	106.40
26	1H	1639	U	N3-C4-O4	-6.58	114.79	119.40
26	1H	594	U	C5-C6-N1	-6.58	119.41	122.70
26	1H	2731	G	C5-C6-O6	-6.58	124.65	128.60
26	14	2436	G	N3-C2-N2	-6.58	115.29	119.90
26	1H	671	C	N3-C4-N4	-6.58	113.39	118.00
44	F8	67	GLY	N-CA-C	-6.58	96.65	113.10
1	13	902	G	O5'-P-OP2	-6.58	99.78	105.70
1	13	1446	A	O4'-C1'-N9	6.58	113.46	108.20
27	1J	101	A	N1-C6-N6	6.58	122.55	118.60
26	1H	1968	G	C2-N3-C4	6.57	115.19	111.90
26	1H	1195	G	N3-C2-N2	-6.57	115.30	119.90
1	13	181	G	N3-C4-C5	-6.57	125.31	128.60
1	13	795	C	N1-C2-O2	-6.57	114.96	118.90
26	1H	51	G	O4'-C1'-N9	-6.57	102.94	108.20
26	1H	1330	C	O5'-P-OP1	-6.57	99.79	105.70
26	14	703	U	C5-C4-O4	6.57	129.84	125.90
26	14	819	A	C8-N9-C4	-6.57	103.17	105.80
26	14	940	G	C2-N3-C4	6.57	115.19	111.90
26	14	2392	A	N1-C2-N3	6.57	132.59	129.30
26	14	2595	G	O5'-P-OP1	-6.57	99.79	105.70
26	1H	247	G	C8-N9-C4	6.57	109.03	106.40
26	1H	828	U	N3-C2-O2	-6.57	117.60	122.20
26	14	110	G	C5-C6-O6	-6.57	124.66	128.60
26	1H	807	U	N3-C4-O4	6.57	124.00	119.40
26	1H	2446	G	C4-C5-N7	6.57	113.43	110.80
26	14	244	A	N1-C6-N6	6.57	122.54	118.60
26	1H	80	G	O5'-P-OP1	-6.57	99.79	105.70
26	1H	693	C	O5'-P-OP2	-6.57	99.79	105.70
26	1H	1021	A	C4-C5-N7	6.57	113.98	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	792	A	C3'-C2'-C1'	-6.56	96.25	101.50
26	1H	1830	C	N1-C2-O2	-6.56	114.96	118.90
23	2K	74	A	N9-C4-C5	-6.56	103.17	105.80
26	1H	874	G	O5'-P-OP1	6.56	118.58	110.70
26	1H	1300	U	OP1-P-O3'	6.56	119.64	105.20
26	1H	1358	G	C6-C5-N7	-6.56	126.46	130.40
26	14	2496	C	C5-C4-N4	-6.56	115.61	120.20
1	13	770	C	O5'-P-OP2	6.56	118.57	110.70
26	14	1899	G	C5-C6-N1	-6.56	108.22	111.50
1	13	285	G	N1-C6-O6	6.56	123.84	119.90
26	1H	2731	G	N3-C4-N9	6.56	129.94	126.00
26	1H	1373	A	C8-N9-C4	6.56	108.42	105.80
22	1K	38	A	N1-C6-N6	6.56	122.53	118.60
26	1H	138	G	C5-N7-C8	-6.56	101.02	104.30
26	14	2607	G	C8-N9-C1'	-6.56	118.48	127.00
26	1H	1203	G	O5'-P-OP2	-6.55	99.80	105.70
26	1H	1228	G	C2-N3-C4	-6.55	108.62	111.90
1	1G	566	G	O4'-C1'-N9	-6.55	102.96	108.20
1	13	1348	U	C5-C4-O4	6.55	129.83	125.90
26	1H	2604	U	N3-C2-O2	-6.55	117.61	122.20
26	14	2359	C	C5-C4-N4	6.55	124.79	120.20
26	1H	37	C	N3-C4-C5	-6.55	119.28	121.90
26	1H	922	U	O5'-P-OP2	6.55	118.56	110.70
26	1H	1347	G	C5-C6-O6	-6.55	124.67	128.60
26	1H	396	G	C5-C6-O6	-6.55	124.67	128.60
26	1H	2607	G	N3-C2-N2	6.55	124.48	119.90
26	14	2449	U	C5-C4-O4	-6.55	121.97	125.90
26	14	2581	G	OP1-P-OP2	6.55	129.42	119.60
1	13	560	U	C5-C6-N1	6.55	125.97	122.70
26	1H	1416	G	O4'-C1'-N9	6.55	113.44	108.20
26	1H	2586	C	N3-C4-C5	6.55	124.52	121.90
26	14	1786	A	C4-N9-C1'	6.55	138.08	126.30
1	13	763	G	C8-N9-C4	-6.54	103.78	106.40
26	1H	984	A	O5'-P-OP2	-6.54	99.81	105.70
26	14	1470	G	C5-C6-N1	-6.54	108.23	111.50
26	1H	197	A	C2-N3-C4	-6.54	107.33	110.60
26	14	1644	C	O5'-P-OP2	-6.54	99.81	105.70
26	1H	66	C	C5-C6-N1	6.54	124.27	121.00
26	14	1558	A	N1-C2-N3	6.54	132.57	129.30
26	14	2270	G	O5'-P-OP1	-6.54	99.81	105.70
26	1H	2546	U	N1-C2-O2	-6.54	118.22	122.80
26	1H	849	A	N1-C6-N6	6.54	122.52	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2210	G	C4-N9-C1'	6.54	135.00	126.50
26	14	1276	A	C8-N9-C4	6.54	108.42	105.80
26	1H	1839	G	C4-N9-C1'	6.54	135.00	126.50
26	1H	2307	G	N1-C6-O6	6.54	123.82	119.90
26	1H	2009	G	OP1-P-O3'	6.54	119.58	105.20
26	1H	2434	A	O5'-P-OP1	-6.54	99.82	105.70
1	1G	576	G	C4-N9-C1'	6.54	135.00	126.50
1	1G	690	G	O5'-P-OP2	-6.54	99.82	105.70
24	3L	76	A	C4-C5-N7	6.54	113.97	110.70
26	14	1543	A	O5'-P-OP1	6.54	118.54	110.70
26	14	2338	G	O5'-P-OP1	-6.54	99.82	105.70
26	1H	335	C	C6-N1-C2	-6.53	117.69	120.30
26	1H	1280	G	OP1-P-OP2	-6.53	109.80	119.60
26	1H	1931	U	N3-C4-O4	-6.53	114.83	119.40
26	14	188	G	C5-C6-O6	-6.53	124.68	128.60
26	14	1597	A	C8-N9-C4	6.53	108.41	105.80
26	14	2330	G	N1-C6-O6	6.53	123.82	119.90
26	1H	455	C	N3-C4-C5	6.53	124.51	121.90
26	1H	2390	U	O5'-P-OP2	6.53	118.54	110.70
1	1G	894	G	N1-C6-O6	6.53	123.82	119.90
26	14	2607	G	N9-C4-C5	-6.53	102.79	105.40
26	1H	500	G	C4-C5-N7	-6.53	108.19	110.80
26	1H	203	C	N3-C2-O2	6.53	126.47	121.90
26	1H	507	A	C4-C5-N7	6.53	113.96	110.70
26	1H	1278	A	C6-N1-C2	-6.53	114.68	118.60
1	1G	689	C	C6-N1-C2	-6.53	117.69	120.30
26	1H	613	U	C5-C6-N1	-6.53	119.44	122.70
26	1H	698	C	OP1-P-OP2	6.53	129.39	119.60
26	1H	2271	G	C8-N9-C1'	-6.53	118.52	127.00
1	1G	1436	U	C5-C6-N1	6.53	125.96	122.70
26	14	210	C	C2-N3-C4	-6.53	116.64	119.90
26	14	2275	C	P-O3'-C3'	6.53	127.53	119.70
1	13	418	C	N1-C2-O2	6.52	122.81	118.90
26	1H	2069	G	C8-N9-C4	6.52	109.01	106.40
26	1H	446	G	C5-C6-O6	-6.52	124.69	128.60
26	1H	1348	G	C5-C6-O6	-6.52	124.69	128.60
26	1H	1895	C	O5'-P-OP1	-6.52	99.83	105.70
26	1H	146	G	N3-C4-C5	6.52	131.86	128.60
26	1H	1022	G	C4-C5-N7	-6.52	108.19	110.80
26	1H	2011	U	OP1-P-O3'	6.52	119.54	105.20
1	1G	108	G	C6-C5-N7	-6.52	126.49	130.40
26	14	116	C	N1-C2-O2	-6.52	114.99	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	383	U	O5'-P-OP2	6.52	118.52	110.70
1	1G	576	G	N1-C6-O6	6.52	123.81	119.90
22	1K	74	C	N1-C2-O2	6.51	122.81	118.90
26	1H	2570	G	N9-C4-C5	6.51	108.00	105.40
26	1H	209	C	C2-N3-C4	-6.51	116.64	119.90
26	1H	1216	G	O5'-P-OP1	-6.51	99.84	105.70
26	1H	1831	G	O5'-P-OP2	6.51	118.51	110.70
1	1G	354	G	C6-C5-N7	-6.51	126.50	130.40
26	1H	866	A	C4-N9-C1'	6.51	138.01	126.30
26	14	1396	U	C5-C6-N1	-6.51	119.45	122.70
26	1H	1123	C	C5-C6-N1	-6.51	117.75	121.00
26	1H	2335	A	O5'-P-OP2	-6.51	99.84	105.70
26	14	1332	G	N1-C2-N3	6.51	127.80	123.90
26	1H	576	U	C4-C5-C6	6.50	123.60	119.70
26	1H	1694	C	OP2-P-O3'	6.50	119.51	105.20
26	1H	195	A	C5-N7-C8	-6.50	100.65	103.90
1	1G	692	U	N3-C4-O4	6.50	123.95	119.40
26	14	1958	C	C6-N1-C2	6.50	122.90	120.30
26	1H	134	C	C2-N3-C4	-6.50	116.65	119.90
26	14	655	A	N1-C6-N6	-6.50	114.70	118.60
26	1H	1931	U	N1-C2-O2	6.50	127.35	122.80
26	1H	202	U	N3-C4-C5	6.50	118.50	114.60
26	1H	783	A	OP1-P-OP2	6.50	129.35	119.60
26	1H	1572	A	N1-C6-N6	6.50	122.50	118.60
26	1H	1967	C	N3-C2-O2	-6.50	117.35	121.90
26	1H	2447	G	C6-N1-C2	-6.50	121.20	125.10
1	1G	1442	G	C8-N9-C1'	6.50	135.45	127.00
26	14	210	C	C5-C6-N1	-6.50	117.75	121.00
26	1H	32	C	C6-N1-C2	-6.50	117.70	120.30
26	1H	182	A	C8-N9-C4	6.50	108.40	105.80
26	1H	1142(A)	A	C5-C6-N1	-6.50	114.45	117.70
26	1H	1145	C	O5'-P-OP1	-6.50	99.85	105.70
26	1H	1763	G	O5'-P-OP2	-6.50	99.85	105.70
26	1H	1973	G	N1-C2-N2	-6.50	110.35	116.20
26	14	102	G	O4'-C1'-N9	6.50	113.40	108.20
26	14	735	A	O5'-P-OP2	-6.50	99.85	105.70
26	14	2512	C	N3-C4-C5	6.50	124.50	121.90
26	14	530	G	N9-C4-C5	-6.50	102.80	105.40
26	1H	781	A	O5'-P-OP1	-6.49	99.86	105.70
26	1H	2699	C	C5-C6-N1	-6.49	117.75	121.00
26	14	681	G	C5-C6-O6	-6.49	124.70	128.60
26	14	1586	A	C8-N9-C4	-6.49	103.20	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	780	G	C6-C5-N7	-6.49	126.51	130.40
26	1H	71	A	C8-N9-C4	-6.49	103.20	105.80
26	14	121	G	C6-N1-C2	-6.49	121.21	125.10
26	14	1283	G	N3-C4-N9	6.49	129.89	126.00
26	1H	593	G	C6-C5-N7	-6.49	126.51	130.40
26	14	2503	A	C2-N3-C4	6.49	113.84	110.60
1	13	760	G	C6-C5-N7	-6.49	126.51	130.40
26	1H	330	A	C4-C5-N7	6.49	113.94	110.70
26	1H	910	A	C6-C5-N7	-6.49	127.76	132.30
26	1H	265	A	N1-C2-N3	6.48	132.54	129.30
26	1H	2582	G	N3-C4-N9	6.48	129.89	126.00
26	1H	2881	C	O5'-P-OP1	-6.48	99.86	105.70
1	1G	1465	C	C6-N1-C2	-6.48	117.71	120.30
26	14	2580	U	N3-C4-O4	6.48	123.94	119.40
26	14	2880	C	C2-N1-C1'	6.48	125.93	118.80
26	1H	971	C	C4-C5-C6	6.48	120.64	117.40
26	1H	1332	G	C8-N9-C1'	6.48	135.43	127.00
26	1H	2439	A	OP1-P-O3'	6.48	119.46	105.20
26	1H	758	C	C4-C5-C6	-6.48	114.16	117.40
26	1H	2264	C	N3-C2-O2	-6.48	117.36	121.90
26	1H	507	A	C6-C5-N7	-6.48	127.77	132.30
26	1H	2563	U	C5-C4-O4	6.48	129.79	125.90
26	14	1367	A	C2-N3-C4	-6.48	107.36	110.60
26	14	2447	G	N1-C6-O6	6.48	123.79	119.90
26	14	1022	G	N3-C4-N9	-6.48	122.11	126.00
26	1H	839	U	N3-C4-C5	-6.48	110.72	114.60
26	1H	971	C	N1-C2-O2	-6.48	115.01	118.90
26	14	391	G	N9-C4-C5	-6.48	102.81	105.40
26	1H	2468	G	C4-N9-C1'	6.47	134.92	126.50
27	16	115	G	C6-N1-C2	-6.47	121.22	125.10
26	14	2731	G	N7-C8-N9	6.47	116.34	113.10
26	1H	734	A	C5-N7-C8	-6.47	100.66	103.90
26	1H	2541	A	O5'-P-OP2	6.47	118.47	110.70
26	14	1978	A	OP2-P-O3'	6.47	119.44	105.20
26	14	780	G	N1-C6-O6	6.47	123.78	119.90
1	13	285	G	C2-N3-C4	-6.47	108.67	111.90
1	13	1483	A	C5-N7-C8	-6.47	100.67	103.90
26	1H	745	G	C6-C5-N7	-6.47	126.52	130.40
26	1H	1622	G	C5-C6-O6	6.47	132.48	128.60
26	1H	2283	C	N1-C2-O2	-6.47	115.02	118.90
26	1H	277	C	N1-C2-O2	6.47	122.78	118.90
26	1H	1314	C	N3-C2-O2	-6.47	117.37	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
37	45	82	ARG	N-CA-C	6.47	128.46	111.00
26	1H	2269	A	C2-N3-C4	-6.47	107.37	110.60
1	13	684	A	C8-N9-C4	-6.46	103.21	105.80
26	1H	691	C	C2-N3-C4	-6.46	116.67	119.90
26	1H	944	G	N7-C8-N9	6.46	116.33	113.10
26	14	2287	A	N3-C4-C5	6.46	131.32	126.80
26	14	2390	U	O5'-P-OP1	-6.46	99.88	105.70
26	1H	1340	U	C5-C6-N1	-6.46	119.47	122.70
26	1H	1205	U	N1-C2-N3	6.46	118.78	114.90
26	1H	1627	G	O5'-P-OP2	-6.46	99.88	105.70
1	13	982	U	N1-C2-O2	6.46	127.32	122.80
1	13	1412	C	C6-N1-C2	6.46	122.88	120.30
26	1H	121	G	C8-N9-C1'	-6.46	118.60	127.00
26	1H	1597	A	O4'-C1'-N9	6.46	113.37	108.20
26	14	211	A	C2-N3-C4	-6.46	107.37	110.60
26	1H	1780	A	N1-C2-N3	6.46	132.53	129.30
26	1H	1789	A	C5-C6-N6	-6.46	118.53	123.70
26	1H	2232	U	O5'-P-OP2	-6.46	99.89	105.70
26	14	470	A	N1-C6-N6	6.46	122.47	118.60
26	14	2258	C	N3-C4-N4	6.46	122.52	118.00
1	13	865	A	C5-N7-C8	-6.46	100.67	103.90
26	1H	194	G	C8-N9-C4	6.46	108.98	106.40
26	1H	748	G	N3-C4-N9	-6.46	122.13	126.00
26	1H	928	G	C5-C6-O6	-6.46	124.73	128.60
26	1H	2598	A	OP2-P-O3'	6.46	119.41	105.20
43	E8	90	ARG	NE-CZ-NH1	-6.46	117.07	120.30
26	14	1309	G	C6-C5-N7	-6.46	126.53	130.40
26	14	1931	U	OP1-P-OP2	-6.46	109.92	119.60
26	1H	2550	G	N1-C6-O6	6.46	123.77	119.90
26	1H	330	A	C6-C5-N7	-6.45	127.78	132.30
26	1H	1197	G	OP2-P-O3'	6.45	119.40	105.20
26	1H	2377	A	N7-C8-N9	-6.45	110.57	113.80
26	14	113	G	N3-C4-C5	6.45	131.83	128.60
26	14	670	A	N9-C4-C5	-6.45	103.22	105.80
26	14	997	G	N1-C6-O6	-6.45	116.03	119.90
26	1H	613	U	C6-N1-C1'	-6.45	112.17	121.20
26	1H	1202	C	C4-C5-C6	6.45	120.63	117.40
26	14	864	G	OP1-P-OP2	-6.45	109.92	119.60
1	13	12	U	N3-C4-C5	-6.45	110.73	114.60
1	13	1227	A	N3-C4-C5	6.45	131.32	126.80
26	1H	1624	G	C5-C6-N1	6.45	114.72	111.50
26	14	148	C	C5-C6-N1	-6.45	117.78	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1562	A	N1-C6-N6	6.45	122.47	118.60
1	13	1310	G	O5'-P-OP2	-6.45	99.90	105.70
26	1H	2058	A	N9-C4-C5	6.45	108.38	105.80
26	14	130	C	C2-N3-C4	-6.45	116.68	119.90
26	1H	791	C	P-O3'-C3'	6.45	127.44	119.70
26	1H	1609	A	C8-N9-C4	6.45	108.38	105.80
26	1H	1915	U	N3-C4-C5	6.45	118.47	114.60
26	1H	676	A	OP1-P-OP2	6.45	129.27	119.60
26	1H	966	G	N1-C2-N2	-6.45	110.40	116.20
26	1H	1367	A	C2-N3-C4	-6.45	107.38	110.60
26	1H	1899	G	C4-C5-C6	-6.45	114.93	118.80
26	1H	2079	U	C4-C5-C6	6.45	123.57	119.70
1	1G	1124	G	O4'-C1'-N9	6.44	113.36	108.20
1	13	904	C	N3-C4-C5	6.44	124.48	121.90
1	13	1378	C	C6-N1-C2	-6.44	117.72	120.30
26	1H	965	C	N3-C4-N4	6.44	122.51	118.00
26	1H	2253	G	C4-C5-N7	6.44	113.38	110.80
26	1H	2666	C	N3-C4-N4	6.44	122.51	118.00
1	1G	925	G	C8-N9-C4	6.44	108.98	106.40
26	14	265	A	C5-C6-N1	-6.44	114.48	117.70
26	14	1349	A	C4-C5-N7	6.44	113.92	110.70
26	1H	99	U	C2-N1-C1'	6.44	125.43	117.70
26	1H	1916	A	C8-N9-C4	-6.44	103.22	105.80
26	14	2075	U	C5-C6-N1	-6.44	119.48	122.70
26	14	1345	C	C6-N1-C2	-6.44	117.72	120.30
26	1H	633	A	C5-C6-N1	-6.44	114.48	117.70
26	14	330	A	N3-C4-C5	6.44	131.31	126.80
26	1H	1382	G	C4-N9-C1'	-6.44	118.13	126.50
2	1E	158	LEU	CA-CB-CG	6.43	130.10	115.30
26	1H	1336	A	C5-C6-N1	6.43	120.92	117.70
26	1H	1972	A	C5-C6-N6	-6.43	118.55	123.70
26	14	215	G	C5-C6-O6	-6.43	124.74	128.60
26	14	1821	A	C6-N1-C2	-6.43	114.74	118.60
26	14	2430	A	OP1-P-OP2	-6.43	109.95	119.60
26	14	2547	U	OP2-P-O3'	6.43	119.36	105.20
26	1H	60	G	O5'-P-OP2	-6.43	99.91	105.70
26	1H	610	C	C5-C6-N1	-6.43	117.78	121.00
27	16	9	G	OP2-P-O3'	6.43	119.35	105.20
26	14	773	U	N3-C2-O2	-6.43	117.70	122.20
26	1H	1843	C	C5-C6-N1	-6.43	117.78	121.00
26	1H	51	G	OP2-P-O3'	6.43	119.34	105.20
26	1H	119	A	C4-C5-N7	-6.43	107.48	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2703	C	N3-C2-O2	-6.43	117.40	121.90
26	14	1982	C	C5-C6-N1	6.43	124.22	121.00
26	14	1999	C	C5-C4-N4	-6.43	115.70	120.20
1	1G	1417	G	C5-C6-N1	-6.43	108.29	111.50
26	14	2588	G	O5'-P-OP1	6.43	118.41	110.70
26	14	2873	A	C4-N9-C1'	6.43	137.87	126.30
26	1H	1815	A	C6-N1-C2	-6.43	114.74	118.60
26	14	1236	G	C8-N9-C4	6.43	108.97	106.40
26	14	1654	A	N1-C6-N6	-6.43	114.74	118.60
1	13	226	G	N3-C4-N9	6.42	129.85	126.00
26	1H	2307	G	C2-N3-C4	-6.42	108.69	111.90
26	14	1640	C	O5'-P-OP2	-6.42	99.92	105.70
26	14	2078	C	N3-C4-C5	-6.42	119.33	121.90
26	14	2249	U	N1-C2-O2	6.42	127.30	122.80
26	1H	1215	G	C8-N9-C4	-6.42	103.83	106.40
26	1H	2060	A	C2-N3-C4	6.42	113.81	110.60
26	1H	809	G	C5-C6-O6	-6.42	124.75	128.60
26	14	138	G	C5-C6-O6	-6.42	124.75	128.60
26	14	1904	G	O5'-P-OP2	-6.42	99.92	105.70
26	1H	51	G	N3-C2-N2	6.42	124.39	119.90
26	1H	919	G	N3-C4-C5	-6.42	125.39	128.60
26	1H	1021	A	C6-C5-N7	-6.42	127.81	132.30
26	14	740	U	C5-C4-O4	6.42	129.75	125.90
26	14	2224	G	N1-C6-O6	6.42	123.75	119.90
22	1K	34	G	C8-N9-C4	-6.42	103.83	106.40
26	1H	2585	U	N3-C2-O2	-6.42	117.71	122.20
27	16	79	C	OP2-P-O3'	6.42	119.31	105.20
1	13	484	G	N3-C4-C5	-6.41	125.39	128.60
1	13	1190	G	N1-C6-O6	6.41	123.75	119.90
26	1H	132	G	N7-C8-N9	6.41	116.31	113.10
26	1H	328	U	O5'-P-OP1	-6.41	99.93	105.70
26	1H	1805	U	OP1-P-O3'	-6.41	91.09	105.20
26	1H	2248	C	N3-C4-C5	6.41	124.47	121.90
26	1H	2419	U	OP1-P-OP2	-6.41	109.98	119.60
26	1H	2436	G	N1-C2-N2	6.41	121.97	116.20
1	13	1335	C	C6-N1-C2	6.41	122.86	120.30
23	2K	17	C	C2-N1-C1'	6.41	125.85	118.80
26	1H	1558	A	P-O3'-C3'	6.41	127.39	119.70
26	1H	1950	G	C5-C6-O6	6.41	132.45	128.60
26	1H	1968	G	C5-C6-O6	-6.41	124.75	128.60
26	1H	1982	C	C2-N1-C1'	6.41	125.85	118.80
26	14	664	C	C4-C5-C6	6.41	120.61	117.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	974	A	C5-C6-N6	-6.41	118.57	123.70
26	1H	194	G	N3-C4-N9	6.41	129.84	126.00
26	1H	2827	C	N1-C2-O2	-6.41	115.06	118.90
26	14	1804	C	C5-C4-N4	-6.41	115.71	120.20
26	1H	190	A	N1-C6-N6	6.41	122.44	118.60
26	1H	2027	G	C4-C5-N7	-6.41	108.24	110.80
26	1H	2048	G	C8-N9-C4	-6.41	103.84	106.40
26	1H	2599	G	N7-C8-N9	-6.41	109.90	113.10
1	1G	632	A	P-O3'-C3'	6.41	127.39	119.70
26	14	830	G	N9-C4-C5	-6.41	102.84	105.40
26	14	1252	G	C8-N9-C4	6.41	108.96	106.40
26	14	1663	C	C5-C4-N4	-6.41	115.72	120.20
26	14	1689	A	O5'-P-OP2	-6.41	99.94	105.70
1	13	767	A	C8-N9-C4	6.40	108.36	105.80
1	13	1234	C	N3-C2-O2	-6.40	117.42	121.90
26	1H	77	C	C5-C4-N4	-6.40	115.72	120.20
26	1H	1655	A	N1-C6-N6	6.40	122.44	118.60
26	1H	1700	A	C8-N9-C4	6.40	108.36	105.80
26	1H	2300	G	C6-C5-N7	-6.40	126.56	130.40
1	1G	537	G	O5'-P-OP1	-6.40	99.94	105.70
1	13	1519	A	C5-C6-N1	-6.40	114.50	117.70
26	1H	2627	G	C6-C5-N7	-6.40	126.56	130.40
26	1H	2708	G	N9-C4-C5	-6.40	102.84	105.40
26	1H	2827	C	N3-C2-O2	6.40	126.38	121.90
26	14	792	G	C5-C6-N1	6.40	114.70	111.50
26	14	2292	C	O5'-P-OP2	-6.40	99.94	105.70
26	1H	675	A	C5-C6-N6	-6.40	118.58	123.70
26	14	1131	G	O4'-C1'-N9	6.40	113.32	108.20
26	14	2237	G	N3-C4-N9	6.40	129.84	126.00
26	1H	809	G	N3-C4-N9	6.40	129.84	126.00
26	14	945	A	N9-C4-C5	-6.40	103.24	105.80
23	2K	68	C	C2-N1-C1'	6.40	125.83	118.80
26	1H	458	G	O4'-C1'-N9	6.40	113.32	108.20
26	1H	2252	G	OP1-P-OP2	6.40	129.19	119.60
26	14	1597	A	N7-C8-N9	-6.40	110.60	113.80
26	1H	238	C	C6-N1-C2	6.39	122.86	120.30
26	1H	2442	C	C4-C5-C6	6.39	120.60	117.40
24	3L	76	A	C8-N9-C4	-6.39	103.24	105.80
26	1H	1310	G	O5'-P-OP2	6.39	118.37	110.70
27	16	100	G	N7-C8-N9	-6.39	109.90	113.10
1	1G	1281	U	C2-N1-C1'	6.39	125.37	117.70
26	14	660	G	O5'-P-OP2	-6.39	99.95	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1607	C	C6-N1-C1'	-6.39	113.13	120.80
27	16	112	G	N3-C4-N9	6.39	129.83	126.00
26	14	1393	A	O5'-P-OP2	-6.39	99.95	105.70
26	1H	1123	C	N1-C2-O2	-6.39	115.07	118.90
26	14	1434	A	C8-N9-C4	6.39	108.36	105.80
36	35	62	LEU	CB-CG-CD2	-6.39	100.14	111.00
1	13	792	A	N9-C1'-C2'	6.39	122.31	114.00
26	14	1826	G	C5-N7-C8	6.39	107.49	104.30
1	13	1502	A	N1-C2-N3	6.38	132.49	129.30
26	14	1324	G	O5'-P-OP2	-6.38	99.95	105.70
1	13	963	G	C8-N9-C1'	-6.38	118.70	127.00
26	14	1266	G	N7-C8-N9	-6.38	109.91	113.10
26	14	1760	A	N7-C8-N9	-6.38	110.61	113.80
26	1H	1019	U	N3-C2-O2	-6.38	117.73	122.20
27	16	23	G	N3-C4-C5	6.38	131.79	128.60
8	72	133	LEU	CA-CB-CG	6.38	129.97	115.30
1	13	681	C	C6-N1-C2	-6.38	117.75	120.30
26	1H	1606	G	C5-C6-N1	6.38	114.69	111.50
26	1H	2737	G	N1-C6-O6	6.38	123.73	119.90
26	14	2346	A	N1-C2-N3	6.38	132.49	129.30
26	1H	645	C	C6-N1-C2	-6.38	117.75	120.30
1	13	1205	U	N1-C2-O2	-6.37	118.34	122.80
26	1H	1122	G	N9-C4-C5	-6.37	102.85	105.40
26	1H	1799	G	N9-C4-C5	-6.37	102.85	105.40
26	1H	2028	U	C5-C6-N1	-6.37	119.51	122.70
26	1H	2346	A	C4-C5-N7	6.37	113.89	110.70
1	1G	1502	A	C6-C5-N7	-6.37	127.84	132.30
22	1K	74	C	O4'-C1'-N1	6.37	113.30	108.20
26	1H	974	G	O4'-C1'-N9	-6.37	103.10	108.20
26	1H	1336	A	N9-C4-C5	6.37	108.35	105.80
26	1H	1436	G	P-O3'-C3'	6.37	127.35	119.70
26	1H	2853	C	OP2-P-O3'	6.37	119.22	105.20
26	14	683	C	N1-C2-O2	-6.37	115.08	118.90
26	14	802	A	C5-C6-N1	6.37	120.89	117.70
26	14	1801	G	O5'-P-OP1	-6.37	99.97	105.70
26	14	2606	C	O5'-P-OP1	-6.37	99.97	105.70
27	16	26	A	O5'-P-OP2	6.37	118.34	110.70
26	14	2210	G	C4-N9-C1'	6.37	134.78	126.50
26	14	2689	U	C2-N3-C4	-6.37	123.18	127.00
1	13	623	C	C6-N1-C2	-6.37	117.75	120.30
26	1H	371	A	C6-C5-N7	-6.37	127.84	132.30
26	1H	587	C	N1-C2-O2	6.37	122.72	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2271	G	C5-C6-N1	6.37	114.68	111.50
26	1H	2375	G	N9-C1'-C2'	-6.37	105.00	112.00
26	14	806	C	C5-C4-N4	-6.37	115.74	120.20
26	1H	1425	G	C6-N1-C2	-6.36	121.28	125.10
26	1H	1579	A	O4'-C1'-N9	6.36	113.29	108.20
26	1H	1680	U	C5-C4-O4	-6.36	122.08	125.90
26	14	1142(A)	A	C2-N3-C4	-6.36	107.42	110.60
26	1H	122	G	N1-C2-N3	6.36	127.72	123.90
26	1H	1264	G	OP1-P-O3'	6.36	119.19	105.20
26	14	681	G	C8-N9-C4	6.36	108.94	106.40
26	1H	1842	G	N7-C8-N9	-6.36	109.92	113.10
37	88	2	LEU	CA-CB-CG	6.36	129.92	115.30
55	Q8	50	LEU	CA-CB-CG	-6.36	100.68	115.30
26	14	202	U	N1-C2-O2	6.36	127.25	122.80
26	14	811	U	N1-C2-N3	6.36	118.71	114.90
1	13	667	G	C8-N9-C4	-6.36	103.86	106.40
26	1H	383	U	O4'-C1'-N1	6.36	113.28	108.20
26	1H	675	A	N1-C6-N6	6.36	122.41	118.60
26	1H	2445	G	C8-N9-C4	-6.36	103.86	106.40
26	1H	1303	G	OP2-P-O3'	6.35	119.18	105.20
26	1H	1604	C	N3-C4-N4	6.35	122.45	118.00
28	11	213	ARG	NE-CZ-NH2	-6.35	117.12	120.30
26	14	2247	A	N9-C4-C5	6.35	108.34	105.80
26	1H	2212	A	N1-C6-N6	6.35	122.41	118.60
26	1H	262	A	C8-N9-C4	6.35	108.34	105.80
26	1H	1334	G	C6-C5-N7	-6.35	126.59	130.40
26	1H	1733	G	N1-C6-O6	6.35	123.71	119.90
26	14	649	G	N1-C6-O6	6.35	123.71	119.90
26	14	2546	U	OP1-P-OP2	6.35	129.13	119.60
26	1H	127	A	N1-C6-N6	6.35	122.41	118.60
26	1H	129	C	C5-C6-N1	-6.35	117.83	121.00
26	1H	829	A	OP1-P-OP2	6.35	129.12	119.60
26	1H	1839	G	C6-C5-N7	-6.35	126.59	130.40
26	14	129	C	N1-C2-O2	6.35	122.71	118.90
26	1H	633	A	C5-N7-C8	-6.35	100.73	103.90
26	1H	792	G	C5-C6-O6	-6.35	124.79	128.60
26	1H	1225	C	C2-N1-C1'	-6.35	111.82	118.80
26	1H	1496	A	N1-C6-N6	6.35	122.41	118.60
26	1H	2018	G	C8-N9-C4	-6.35	103.86	106.40
26	14	1528	A	N1-C6-N6	6.35	122.41	118.60
26	1H	748	G	C8-N9-C1'	6.35	135.25	127.00
26	1H	842	G	C8-N9-C4	6.34	108.94	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1501	C	O5'-P-OP1	-6.34	99.99	105.70
26	1H	1592	C	C6-N1-C2	6.34	122.84	120.30
26	14	265	A	C5-N7-C8	-6.34	100.73	103.90
26	14	2575	C	C5-C4-N4	6.34	124.64	120.20
26	1H	934	G	O5'-P-OP1	-6.34	99.99	105.70
26	1H	2502	G	C6-N1-C2	-6.34	121.29	125.10
27	16	81	G	N7-C8-N9	6.34	116.27	113.10
32	51	171	LEU	CA-CB-CG	6.34	129.89	115.30
1	13	328	C	O5'-P-OP1	-6.34	99.99	105.70
55	Q8	13	ARG	NE-CZ-NH1	-6.34	117.13	120.30
26	14	466	A	N1-C6-N6	6.34	122.41	118.60
1	13	361	G	N1-C2-N3	6.34	127.70	123.90
26	14	204	A	C6-N1-C2	-6.34	114.80	118.60
26	14	1997	G	N3-C4-N9	6.34	129.80	126.00
27	16	78	A	C2-N3-C4	-6.34	107.43	110.60
1	1G	332	G	N7-C8-N9	-6.34	109.93	113.10
26	1H	1248	G	O5'-P-OP1	6.34	118.30	110.70
26	14	741	G	C5-C6-O6	-6.34	124.80	128.60
26	14	1843	C	C6-N1-C2	6.34	122.83	120.30
33	69	131	LYS	C-N-CD	-6.34	106.66	120.60
26	1H	378	C	C5-C4-N4	-6.33	115.77	120.20
26	1H	691	C	C5-C4-N4	-6.33	115.77	120.20
26	14	2277	G	C5-C6-O6	6.33	132.40	128.60
26	1H	1305	C	C5-C6-N1	-6.33	117.83	121.00
26	1H	1672	C	N3-C4-C5	6.33	124.43	121.90
26	14	1826	G	C8-N9-C4	6.33	108.93	106.40
26	1H	2232	U	N1-C2-N3	6.33	118.70	114.90
26	14	1274	A	C5-C6-N6	-6.33	118.64	123.70
26	1H	640	C	OP1-P-O3'	6.33	119.12	105.20
26	1H	741	G	C5-C6-O6	-6.33	124.80	128.60
26	1H	2264	C	C6-N1-C2	-6.33	117.77	120.30
1	1G	14	U	C5-C6-N1	6.33	125.86	122.70
1	1G	345	C	N1-C2-O2	6.33	122.70	118.90
26	14	297	C	C6-N1-C2	-6.33	117.77	120.30
26	14	641	C	O5'-P-OP2	6.33	118.30	110.70
26	14	1270	C	OP2-P-O3'	6.33	119.12	105.20
26	1H	120	U	N1-C2-N3	6.33	118.70	114.90
23	2K	22	A	O5'-P-OP1	-6.33	100.01	105.70
26	1H	140	A	C2-N3-C4	-6.33	107.44	110.60
26	1H	536	A	C8-N9-C4	-6.33	103.27	105.80
26	1H	1617	C	O5'-P-OP1	-6.33	100.01	105.70
27	16	53	A	N7-C8-N9	6.33	116.96	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	138	G	C8-N9-C4	-6.33	103.87	106.40
26	14	1781	C	C6-N1-C2	6.33	122.83	120.30
26	1H	204	A	C5-C6-N1	6.32	120.86	117.70
26	1H	866	A	C5-N7-C8	-6.32	100.74	103.90
26	14	786	C	C5-C6-N1	-6.32	117.84	121.00
26	14	1608	A	N9-C4-C5	6.32	108.33	105.80
26	1H	1347	G	N3-C2-N2	-6.32	115.47	119.90
26	1H	501	A	OP1-P-OP2	6.32	129.08	119.60
26	1H	728	G	O5'-P-OP2	-6.32	100.01	105.70
26	1H	137(A)	G	C5-C6-O6	-6.32	124.81	128.60
26	1H	467	G	C8-N9-C4	6.32	108.93	106.40
26	14	1251	C	N3-C4-N4	6.32	122.42	118.00
26	14	2477	C	C2-N1-C1'	6.32	125.75	118.80
26	1H	2695	C	N1-C2-O2	-6.32	115.11	118.90
1	1G	896	C	N3-C4-C5	6.32	124.43	121.90
26	14	122	G	C8-N9-C4	6.32	108.93	106.40
26	14	804	A	C8-N9-C4	6.32	108.33	105.80
26	1H	1373	A	N7-C8-N9	-6.32	110.64	113.80
26	1H	2281	C	C5-C4-N4	-6.32	115.78	120.20
26	1H	2581	G	N3-C2-N2	6.32	124.32	119.90
26	14	210	C	N3-C4-C5	6.32	124.43	121.90
26	14	681	G	C6-C5-N7	-6.32	126.61	130.40
26	14	827	U	C5-C4-O4	-6.32	122.11	125.90
26	14	2764	A	C8-N9-C4	6.32	108.33	105.80
48	F5	36	GLY	N-CA-C	6.32	128.89	113.10
26	14	566	U	C6-N1-C2	6.31	124.79	121.00
26	14	1216	G	N1-C6-O6	6.31	123.69	119.90
26	1H	36	G	O5'-P-OP2	-6.31	100.02	105.70
26	1H	541	C	N3-C2-O2	-6.31	117.48	121.90
26	1H	568	U	C5-C4-O4	-6.31	122.11	125.90
26	1H	1809	A	C5-C6-N1	6.31	120.86	117.70
27	1J	56	G	N3-C4-N9	6.31	129.79	126.00
26	1H	827	U	C5-C6-N1	-6.31	119.54	122.70
26	1H	1618	A	C5-N7-C8	-6.31	100.74	103.90
27	16	15	A	O4'-C1'-N9	6.31	113.25	108.20
1	13	913	A	C8-N9-C4	6.31	108.32	105.80
26	1H	1034	G	C5-C6-O6	-6.31	124.81	128.60
26	1H	2773	C	C5-C6-N1	-6.31	117.84	121.00
26	14	1325	G	N3-C4-C5	-6.31	125.45	128.60
1	13	576	G	C4-N9-C1'	6.31	134.70	126.50
26	1H	2708	G	C8-N9-C4	6.31	108.92	106.40
26	14	2065	C	C2-N1-C1'	6.31	125.74	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2427	C	N1-C2-O2	-6.31	115.12	118.90
26	14	2607	G	C6-C5-N7	-6.31	126.62	130.40
1	13	804	U	C5-C6-N1	-6.31	119.55	122.70
27	16	100	G	N9-C4-C5	-6.31	102.88	105.40
1	13	280	C	C6-N1-C2	6.30	122.82	120.30
26	1H	781	A	OP1-P-OP2	6.30	129.06	119.60
26	1H	1200	C	C5-C6-N1	-6.30	117.85	121.00
26	1H	1265	A	O5'-P-OP2	-6.30	100.03	105.70
26	1H	1308	A	C8-N9-C4	-6.30	103.28	105.80
26	1H	1936	A	C5-C6-N6	-6.30	118.66	123.70
26	1H	2026	C	C4-C5-C6	6.30	120.55	117.40
26	14	57	C	N3-C2-O2	6.30	126.31	121.90
26	14	1407	C	O5'-P-OP1	-6.30	100.03	105.70
26	14	1411	C	O5'-P-OP1	6.30	118.27	110.70
26	14	2433	A	C8-N9-C4	-6.30	103.28	105.80
1	13	821	G	C8-N9-C1'	-6.30	118.81	127.00
26	14	1349	A	C6-C5-N7	-6.30	127.89	132.30
1	13	583	A	N9-C4-C5	-6.30	103.28	105.80
26	1H	138	G	C4-C5-C6	-6.30	115.02	118.80
26	1H	142	G	C4-N9-C1'	-6.30	118.31	126.50
26	1H	748	G	C4-N9-C1'	-6.30	118.31	126.50
26	1H	1521	G	C8-N9-C4	-6.30	103.88	106.40
26	1H	2518	A	C4-C5-N7	6.30	113.85	110.70
26	14	1024	G	N1-C6-O6	6.30	123.68	119.90
26	14	1950	G	O4'-C1'-N9	6.30	113.24	108.20
26	1H	2580	U	OP2-P-O3'	6.30	119.06	105.20
26	1H	2620	C	N3-C2-O2	6.30	126.31	121.90
26	14	2346	A	C4-N9-C1'	6.30	137.64	126.30
26	1H	110	G	OP1-P-OP2	6.30	129.05	119.60
26	1H	994	C	C6-N1-C2	-6.30	117.78	120.30
26	1H	1385	G	N3-C4-C5	6.30	131.75	128.60
26	14	76	C	C6-N1-C2	-6.30	117.78	120.30
26	1H	1836	C	C6-N1-C2	-6.30	117.78	120.30
1	1G	1502	A	N7-C8-N9	6.30	116.95	113.80
26	14	2731	G	C6-C5-N7	-6.30	126.62	130.40
1	13	1495	U	C5-C4-O4	6.29	129.68	125.90
26	1H	1138	G	C8-N9-C1'	-6.29	118.82	127.00
1	13	1299	A	N1-C6-N6	6.29	122.38	118.60
26	1H	1413	G	C8-N9-C4	-6.29	103.88	106.40
26	14	690	G	N1-C6-O6	6.29	123.68	119.90
26	14	2490	G	N7-C8-N9	6.29	116.25	113.10
1	13	169	C	C2-N1-C1'	6.29	125.72	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	842	G	C5-C6-O6	-6.29	124.83	128.60
26	1H	1299	G	N9-C4-C5	-6.29	102.88	105.40
26	14	1614	A	C5-C6-N1	-6.29	114.55	117.70
26	1H	1229(A)	G	C2-N3-C4	-6.29	108.75	111.90
26	1H	1349	A	C2-N3-C4	-6.29	107.45	110.60
26	1H	1657	C	C6-N1-C2	-6.29	117.78	120.30
26	1H	2603	G	OP1-P-O3'	6.29	119.04	105.20
26	1H	790	C	N3-C2-O2	6.29	126.30	121.90
26	1H	1209	G	N1-C6-O6	6.29	123.67	119.90
26	1H	1613	G	N1-C6-O6	-6.29	116.13	119.90
26	1H	1698	A	N3-C4-N9	-6.29	122.37	127.40
1	1G	963	G	C4-N9-C1'	6.29	134.67	126.50
26	14	568	U	N1-C2-O2	6.29	127.20	122.80
26	1H	140	A	O4'-C1'-N9	6.29	113.23	108.20
26	1H	2761	G	C8-N9-C4	6.29	108.92	106.40
1	1G	121	C	C6-N1-C1'	-6.29	113.25	120.80
26	1H	2584	U	N1-C2-O2	6.29	127.20	122.80
1	1G	913	A	OP2-P-O3'	6.29	119.03	105.20
26	14	2076	U	O5'-P-OP2	-6.29	100.04	105.70
26	1H	266	G	O5'-P-OP2	-6.28	100.05	105.70
26	1H	1471	A	C4-C5-N7	6.28	113.84	110.70
26	1H	1598	C	N3-C4-C5	-6.28	119.39	121.90
26	1H	1695	G	N3-C2-N2	6.28	124.30	119.90
26	1H	2698	U	C4-C5-C6	6.28	123.47	119.70
25	4K	17	U	C5-C6-N1	-6.28	119.56	122.70
26	1H	823	G	N9-C4-C5	-6.28	102.89	105.40
1	13	541	G	C5-C6-O6	-6.28	124.83	128.60
26	1H	630	G	N7-C8-N9	-6.28	109.96	113.10
26	1H	1402	C	C5-C6-N1	6.28	124.14	121.00
28	19	111	LEU	CA-CB-CG	6.28	129.74	115.30
26	1H	657	U	OP2-P-O3'	6.28	119.01	105.20
26	1H	1424	G	N1-C2-N3	6.28	127.67	123.90
26	1H	2822	G	C4-C5-N7	6.28	113.31	110.80
1	13	764	C	N3-C4-C5	6.28	124.41	121.90
26	1H	468	G	C8-N9-C4	6.28	108.91	106.40
26	1H	811	U	N1-C2-O2	-6.28	118.41	122.80
26	1H	1379	A	N1-C6-N6	6.28	122.37	118.60
26	1H	1764	G	N1-C2-N3	6.28	127.67	123.90
26	14	1643	G	OP2-P-O3'	6.28	119.01	105.20
26	1H	966	G	C5-C6-O6	6.28	132.37	128.60
1	1G	353	A	N1-C6-N6	6.27	122.36	118.60
26	14	189	G	N7-C8-N9	-6.27	109.96	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	3E	32	ALA	N-CA-C	-6.27	94.06	111.00
26	14	2713	A	N1-C6-N6	6.27	122.36	118.60
26	1H	966	G	N7-C8-N9	-6.27	109.96	113.10
26	14	252	G	C5-C6-O6	6.27	132.36	128.60
26	1H	2389	G	OP1-P-O3'	6.27	118.99	105.20
26	14	196	A	O5'-P-OP2	-6.27	100.06	105.70
1	13	1517	G	C5-C6-N1	6.27	114.63	111.50
26	1H	845	G	N9-C4-C5	-6.27	102.89	105.40
26	1H	1786	A	C5-C6-N1	-6.27	114.57	117.70
26	14	798	G	N3-C2-N2	-6.27	115.51	119.90
26	14	2713	A	C4-C5-N7	6.27	113.83	110.70
26	1H	606	U	N3-C2-O2	-6.27	117.81	122.20
26	1H	906	G	C6-C5-N7	6.27	134.16	130.40
19	AI	25	LYS	N-CA-C	-6.26	94.09	111.00
26	1H	1193	G	N7-C8-N9	-6.26	109.97	113.10
26	1H	2192	G	C4-N9-C1'	6.26	134.64	126.50
26	1H	2247	A	O5'-P-OP1	-6.26	100.06	105.70
19	AA	82	GLY	N-CA-C	6.26	128.76	113.10
26	1H	2427	C	N1-C2-O2	-6.26	115.14	118.90
26	1H	2450	A	N1-C2-N3	6.26	132.43	129.30
26	14	848	G	N3-C4-N9	6.26	129.76	126.00
26	14	1804	C	C5-C6-N1	6.26	124.13	121.00
26	1H	1224	G	O5'-P-OP1	-6.26	100.07	105.70
1	1G	576	G	C4-C5-C6	6.26	122.56	118.80
26	14	1142	U	N3-C2-O2	-6.26	117.82	122.20
1	13	579	G	N1-C6-O6	6.26	123.65	119.90
26	14	2247	A	C5-C6-N6	6.26	128.70	123.70
22	1K	48	C	N1-C2-O2	6.25	122.65	118.90
1	13	1313	U	C5-C6-N1	6.25	125.83	122.70
26	1H	61	G	C8-N9-C4	6.25	108.90	106.40
26	1H	1520	U	C5-C4-O4	6.25	129.65	125.90
26	1H	1806	C	OP1-P-OP2	6.25	128.98	119.60
26	1H	2282	G	O5'-P-OP1	-6.25	100.07	105.70
26	1H	2625	G	C5-C6-O6	-6.25	124.85	128.60
12	3A	27	LEU	CA-CB-CG	6.25	129.68	115.30
26	1H	671	C	C6-N1-C2	6.25	122.80	120.30
26	1H	850	C	C5-C4-N4	-6.25	115.82	120.20
26	1H	2502	G	N3-C4-N9	6.25	129.75	126.00
26	14	1801	G	C5-C6-O6	-6.25	124.85	128.60
1	13	888	G	C8-N9-C4	6.25	108.90	106.40
26	1H	698	C	N3-C4-C5	6.25	124.40	121.90
26	1H	2314	C	O5'-P-OP2	-6.25	100.08	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	2L	27	G	C8-N9-C4	6.25	108.90	106.40
26	14	1313	U	N3-C4-C5	-6.25	110.85	114.60
26	1H	1349	A	N1-C6-N6	6.25	122.35	118.60
26	14	703	U	N3-C4-O4	-6.25	115.03	119.40
26	14	1908	C	C6-N1-C2	-6.25	117.80	120.30
26	1H	179	G	N1-C2-N3	6.24	127.65	123.90
26	1H	762	U	C2-N1-C1'	6.24	125.19	117.70
26	1H	1607	C	N3-C4-N4	6.24	122.37	118.00
26	1H	2378	A	N9-C4-C5	-6.24	103.30	105.80
26	14	138	G	C2-N3-C4	6.24	115.02	111.90
26	14	1378	A	N1-C2-N3	-6.24	126.18	129.30
26	14	2499	C	N1-C2-O2	-6.24	115.15	118.90
26	1H	1366	A	N9-C4-C5	-6.24	103.30	105.80
26	14	1281	G	C4-C5-N7	6.24	113.30	110.80
26	1H	1379	A	N9-C1'-C2'	6.24	122.11	114.00
26	1H	52	A	C2-N3-C4	6.24	113.72	110.60
26	1H	263	C	O5'-P-OP1	6.24	118.18	110.70
26	1H	763	G	C4-C5-C6	6.24	122.54	118.80
26	1H	1594	G	N7-C8-N9	6.24	116.22	113.10
26	14	826	U	N3-C4-C5	-6.24	110.86	114.60
26	14	2707	G	N1-C6-O6	-6.24	116.16	119.90
26	1H	784	A	N3-C4-N9	-6.24	122.41	127.40
26	1H	2036	C	C5-C6-N1	6.24	124.12	121.00
1	13	1227	A	O5'-P-OP2	-6.23	100.09	105.70
26	1H	470	A	C6-N1-C2	-6.23	114.86	118.60
27	16	45	A	OP1-P-OP2	6.23	128.95	119.60
26	1H	427	U	N3-C2-O2	6.23	126.56	122.20
26	1H	526	A	C5-C6-N6	6.23	128.69	123.70
26	1H	2762	G	C4-C5-N7	6.23	113.29	110.80
26	14	187	G	C6-N1-C2	-6.23	121.36	125.10
26	14	2240	C	C6-N1-C2	-6.23	117.81	120.30
26	1H	336	C	N3-C4-N4	6.23	122.36	118.00
26	1H	655	A	C4-C5-C6	6.23	120.12	117.00
26	1H	1256	G	C8-N9-C1'	-6.23	118.90	127.00
26	1H	1300	U	N1-C2-O2	-6.23	118.44	122.80
1	1G	121	C	C2-N1-C1'	6.23	125.65	118.80
22	1K	34	G	N7-C8-N9	6.23	116.22	113.10
26	1H	1914	C	C5-C4-N4	6.23	124.56	120.20
1	13	22	G	O5'-P-OP1	-6.23	100.10	105.70
1	13	291	C	N1-C2-O2	-6.23	115.16	118.90
1	13	518	C	C6-N1-C2	6.23	122.79	120.30
26	1H	835	A	C5-C6-N1	6.23	120.81	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	949	C	C4-C5-C6	6.23	120.51	117.40
26	14	1030	G	C4-C5-N7	6.23	113.29	110.80
26	14	2084	C	N1-C2-O2	-6.23	115.16	118.90
26	14	1427	A	N1-C6-N6	-6.23	114.86	118.60
1	13	505	G	O5'-P-OP2	6.22	118.17	110.70
26	1H	733	G	N3-C2-N2	6.22	124.26	119.90
26	1H	2359	C	C5-C6-N1	-6.22	117.89	121.00
26	14	1125	G	N1-C6-O6	-6.22	116.17	119.90
26	14	2427	C	O5'-P-OP2	6.22	118.17	110.70
26	1H	305	U	C5-C6-N1	6.22	125.81	122.70
26	1H	823	G	N1-C2-N3	6.22	127.63	123.90
26	14	2720	U	C5-C4-O4	6.22	129.63	125.90
1	13	1260	C	C5-C6-N1	6.22	124.11	121.00
26	1H	689	A	O5'-P-OP2	-6.22	100.10	105.70
1	13	972	C	C6-N1-C2	-6.22	117.81	120.30
1	13	1498	U	P-O3'-C3'	6.22	127.16	119.70
26	1H	188	G	OP1-P-OP2	6.22	128.93	119.60
26	1H	1287	A	OP1-P-OP2	-6.22	110.27	119.60
26	1H	1655	A	C5-C6-N6	-6.22	118.72	123.70
1	1G	598	U	N1-C2-O2	-6.22	118.45	122.80
26	1H	2449	U	N3-C4-C5	-6.22	110.87	114.60
26	1H	827	U	N1-C2-O2	-6.22	118.45	122.80
26	14	203	C	C2-N3-C4	-6.22	116.79	119.90
26	14	1988	C	C5-C4-N4	-6.22	115.85	120.20
1	13	585	G	C8-N9-C4	6.21	108.89	106.40
23	2K	9	G	C2-N3-C4	6.21	115.01	111.90
26	1H	2451	A	C5-C6-N6	6.21	128.67	123.70
26	14	1394	U	C5-C4-O4	6.21	129.63	125.90
26	14	2681	C	C5-C4-N4	6.21	124.55	120.20
26	1H	1416	G	C4-N9-C1'	-6.21	118.42	126.50
26	1H	1942	C	C6-N1-C2	-6.21	117.81	120.30
26	1H	1021	A	N3-C4-C5	6.21	131.15	126.80
26	1H	1363	C	C5-C6-N1	-6.21	117.89	121.00
26	1H	2554	U	C5-C6-N1	6.21	125.81	122.70
26	1H	1641	A	O5'-P-OP1	-6.21	100.11	105.70
26	1H	2010	G	O5'-P-OP2	6.21	118.15	110.70
26	1H	736	C	C6-N1-C2	6.21	122.78	120.30
1	13	770	C	N1-C2-O2	-6.21	115.18	118.90
1	13	789	U	C6-N1-C2	-6.21	117.28	121.00
26	1H	59	U	OP2-P-O3'	6.21	118.85	105.20
26	1H	241	A	O5'-P-OP2	-6.21	100.11	105.70
26	1H	1269	A	N7-C8-N9	6.21	116.90	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	817	C	C5-C6-N1	-6.21	117.90	121.00
26	1H	561	G	OP1-P-OP2	6.20	128.91	119.60
26	1H	967	C	C6-N1-C2	6.20	122.78	120.30
26	1H	249	C	N3-C2-O2	-6.20	117.56	121.90
26	1H	112	U	N3-C2-O2	6.20	126.54	122.20
26	1H	2054	A	C5-C6-N6	-6.20	118.74	123.70
26	14	2406	U	OP1-P-O3'	6.20	118.84	105.20
1	13	893	C	N1-C2-O2	6.20	122.62	118.90
26	1H	585	G	N1-C6-O6	6.20	123.62	119.90
26	1H	826	U	N1-C2-O2	-6.20	118.46	122.80
26	14	1388	G	O5'-P-OP2	-6.20	100.12	105.70
26	1H	121	G	N3-C4-C5	-6.20	125.50	128.60
26	1H	786	C	C2-N3-C4	-6.20	116.80	119.90
27	1J	22	U	C5-C6-N1	6.20	125.80	122.70
26	1H	1574	C	O5'-P-OP1	6.20	118.13	110.70
26	1H	2254	C	N1-C2-O2	-6.20	115.18	118.90
26	1H	78	A	N1-C6-N6	6.19	122.32	118.60
26	1H	238	C	C2-N3-C4	-6.19	116.80	119.90
26	1H	779	U	C6-N1-C2	6.19	124.72	121.00
1	13	23	C	C6-N1-C2	-6.19	117.82	120.30
26	1H	124	G	C5-C6-N1	6.19	114.60	111.50
26	1H	774	A	C8-N9-C4	-6.19	103.32	105.80
26	1H	1915	U	C2-N3-C4	-6.19	123.28	127.00
27	16	7	G	N1-C6-O6	6.19	123.62	119.90
27	16	60	C	C5-C6-N1	6.19	124.10	121.00
26	14	330	A	N9-C4-C5	-6.19	103.32	105.80
26	14	597	U	N1-C2-O2	-6.19	118.47	122.80
26	14	2375	G	C5-N7-C8	6.19	107.40	104.30
26	14	2776	A	N7-C8-N9	6.19	116.90	113.80
26	1H	1228	G	N1-C2-N3	6.19	127.61	123.90
26	1H	1298	C	N1-C2-O2	6.19	122.61	118.90
26	1H	1647	G	C4-C5-N7	-6.19	108.32	110.80
1	1G	249	U	O5'-P-OP2	-6.19	100.13	105.70
26	14	2276	G	N3-C2-N2	-6.19	115.57	119.90
26	1H	175	G	N1-C6-O6	-6.19	116.19	119.90
26	1H	566	U	C6-N1-C2	6.19	124.71	121.00
26	1H	1537	C	C6-N1-C2	-6.19	117.83	120.30
27	16	61	G	C8-N9-C4	-6.19	103.92	106.40
26	14	40	C	C6-N1-C2	-6.19	117.83	120.30
26	1H	996	A	C8-N9-C4	6.18	108.27	105.80
26	1H	108	U	O5'-P-OP1	-6.18	100.14	105.70
1	1G	1508	G	C8-N9-C4	6.18	108.87	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1528	A	N7-C8-N9	6.18	116.89	113.80
1	13	1158	C	C6-N1-C2	-6.18	117.83	120.30
26	1H	130	C	C5-C4-N4	-6.18	115.88	120.20
26	1H	1249	U	C5-C4-O4	-6.18	122.19	125.90
26	1H	1614	A	C5-C6-N1	-6.18	114.61	117.70
26	14	2584	U	OP1-P-OP2	-6.18	110.33	119.60
26	1H	1790	C	N1-C2-O2	-6.18	115.19	118.90
35	68	22	ILE	CG1-CB-CG2	-6.18	97.81	111.40
26	1H	1266	G	C4-C5-N7	6.18	113.27	110.80
26	1H	1614	A	N3-C4-N9	-6.18	122.46	127.40
26	1H	2061	G	OP1-P-O3'	6.18	118.79	105.20
26	1H	2585	U	N1-C2-O2	6.18	127.12	122.80
26	1H	2713	A	N3-C4-C5	6.18	131.12	126.80
43	E8	11	ARG	NE-CZ-NH1	6.18	123.39	120.30
1	13	1405	G	C5-C6-O6	-6.17	124.89	128.60
26	1H	74	A	N7-C8-N9	6.17	116.89	113.80
37	88	82	ARG	N-CA-C	6.17	127.67	111.00
26	14	1407	C	N3-C2-O2	6.17	126.22	121.90
26	14	1661	G	N1-C6-O6	6.17	123.60	119.90
27	16	49	C	N3-C2-O2	6.17	126.22	121.90
26	14	2590	A	N7-C8-N9	-6.17	110.71	113.80
1	13	507	C	N3-C4-C5	-6.17	119.43	121.90
26	1H	216	A	N1-C2-N3	6.17	132.38	129.30
26	1H	651	G	O5'-P-OP2	6.17	118.11	110.70
26	14	1129	A	O5'-P-OP2	-6.17	100.15	105.70
26	1H	16	G	N1-C2-N3	6.17	127.60	123.90
26	1H	770	G	OP1-P-OP2	-6.17	110.35	119.60
26	1H	817	C	OP2-P-O3'	6.17	118.77	105.20
26	1H	1257	C	C5-C4-N4	-6.17	115.88	120.20
26	1H	2018	G	N7-C8-N9	6.17	116.18	113.10
26	1H	2550	G	O5'-P-OP2	-6.17	100.15	105.70
26	14	1474	C	N1-C2-O2	6.17	122.60	118.90
26	1H	2497	A	N3-C4-N9	6.17	132.33	127.40
1	1G	345	C	C5-C6-N1	6.17	124.08	121.00
26	14	2700	C	C5-C4-N4	-6.17	115.88	120.20
26	1H	2271	G	C6-C5-N7	-6.17	126.70	130.40
26	1H	2308	G	N1-C6-O6	6.17	123.60	119.90
26	14	621	A	N1-C6-N6	6.17	122.30	118.60
26	1H	1391	U	N3-C4-O4	6.16	123.72	119.40
1	1G	334	C	C6-N1-C2	6.16	122.77	120.30
1	1G	862	C	C6-N1-C1'	-6.16	113.40	120.80
26	14	776	G	C5-C6-O6	-6.16	124.90	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	1332	A	C8-N9-C4	-6.16	103.34	105.80
26	14	57	C	N1-C2-O2	-6.16	115.20	118.90
26	1H	594	U	C6-N1-C2	6.16	124.69	121.00
26	14	558	G	N7-C8-N9	-6.16	110.02	113.10
26	14	2346	A	C5-N7-C8	-6.16	100.82	103.90
26	1H	1623	G	OP2-P-O3'	6.16	118.75	105.20
1	1G	1190	G	C4-C5-N7	-6.16	108.34	110.80
26	1H	455	C	N3-C4-N4	-6.16	113.69	118.00
48	J8	80	LEU	CA-CB-CG	6.16	129.46	115.30
26	14	1571	A	C5-C6-N6	-6.16	118.78	123.70
26	14	1572	A	C5-C6-N6	-6.16	118.78	123.70
26	14	2681	C	N3-C4-N4	-6.16	113.69	118.00
26	1H	1324	G	C5-C6-O6	-6.15	124.91	128.60
5	4E	12	LEU	CA-CB-CG	6.15	129.45	115.30
26	1H	1807	G	C5-C6-O6	-6.15	124.91	128.60
26	14	1470	G	N1-C6-O6	6.15	123.59	119.90
26	14	1992	G	C5-C6-N1	6.15	114.58	111.50
26	14	2554	U	O5'-P-OP2	6.15	118.08	110.70
26	1H	1285	G	C5-C6-N1	6.15	114.58	111.50
26	1H	1334	G	N7-C8-N9	6.15	116.17	113.10
26	1H	1559	G	N1-C6-O6	6.15	123.59	119.90
26	14	777	A	N1-C6-N6	-6.15	114.91	118.60
26	14	1342	A	O4'-C1'-N9	6.15	113.12	108.20
1	1G	875	C	N3-C2-O2	-6.15	117.60	121.90
26	14	1340	U	C5-C6-N1	-6.15	119.62	122.70
26	1H	593	G	C6-N1-C2	-6.15	121.41	125.10
26	1H	917	A	C6-C5-N7	-6.15	128.00	132.30
1	1G	898	G	N1-C6-O6	6.15	123.59	119.90
26	14	1253	A	N3-C4-N9	6.15	132.32	127.40
26	14	1407	C	C4-C5-C6	-6.15	114.33	117.40
26	14	2301	C	C5-C6-N1	6.15	124.07	121.00
1	13	23	C	C5-C6-N1	6.15	124.07	121.00
26	1H	2082	A	N1-C2-N3	6.15	132.37	129.30
26	14	801	G	C4-C5-N7	-6.15	108.34	110.80
26	1H	23	G	N3-C2-N2	-6.14	115.60	119.90
26	1H	2051	A	C5-N7-C8	-6.14	100.83	103.90
26	14	2570	G	N3-C2-N2	-6.14	115.60	119.90
26	1H	1264	G	C6-C5-N7	-6.14	126.72	130.40
26	1H	1624	G	C8-N9-C4	6.14	108.86	106.40
1	13	1412	C	C5-C6-N1	-6.14	117.93	121.00
26	1H	247	G	N9-C4-C5	-6.14	102.94	105.40
1	1G	481	G	N3-C4-N9	6.14	129.68	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1813	G	N1-C6-O6	-6.14	116.22	119.90
26	1H	738	G	C4-C5-N7	6.14	113.26	110.80
26	1H	1147	C	C6-N1-C2	6.14	122.76	120.30
26	1H	2439	A	C4-C5-N7	6.14	113.77	110.70
26	14	509	C	C4-C5-C6	6.14	120.47	117.40
26	14	1914	C	C6-N1-C2	-6.14	117.84	120.30
26	14	2304	G	C8-N9-C4	-6.14	103.94	106.40
26	1H	800	A	OP1-P-OP2	6.14	128.81	119.60
26	1H	1786	A	C4-N9-C1'	6.14	137.35	126.30
26	1H	1594	G	OP1-P-O3'	6.14	118.70	105.20
26	14	1795	C	C6-N1-C2	-6.14	117.85	120.30
26	1H	321	G	N1-C6-O6	6.13	123.58	119.90
26	1H	418	G	C8-N9-C4	6.13	108.85	106.40
26	1H	790	C	N1-C2-O2	-6.13	115.22	118.90
26	1H	1677	A	C8-N9-C4	6.13	108.25	105.80
23	2L	35	C	C2-N1-C1'	6.13	125.55	118.80
26	14	1279	G	O5'-P-OP2	-6.13	100.18	105.70
26	1H	2318	G	C2-N3-C4	-6.13	108.83	111.90
26	1H	1353	A	C6-N1-C2	-6.13	114.92	118.60
26	14	1355	G	N7-C8-N9	6.13	116.17	113.10
26	1H	346	A	C8-N9-C4	-6.13	103.35	105.80
26	1H	935	C	C6-N1-C2	6.13	122.75	120.30
26	1H	1929	G	O5'-P-OP2	-6.13	100.18	105.70
26	1H	2424	C	C5-C6-N1	6.13	124.06	121.00
1	1G	353	A	C4-C5-N7	6.13	113.76	110.70
26	14	1355	G	C8-N9-C4	-6.13	103.95	106.40
26	14	1366	A	C5-N7-C8	-6.13	100.84	103.90
26	14	2354	G	C5-C6-O6	-6.13	124.92	128.60
26	1H	130	C	C5-C6-N1	-6.13	117.94	121.00
26	1H	769	G	N3-C2-N2	6.13	124.19	119.90
26	1H	2488	A	N1-C6-N6	6.13	122.28	118.60
26	14	2224	G	C6-C5-N7	-6.13	126.72	130.40
26	14	2473	U	N1-C2-O2	6.13	127.09	122.80
27	1J	6	C	C6-N1-C2	6.13	122.75	120.30
26	1H	239	U	C5-C4-O4	6.12	129.57	125.90
26	1H	1192	G	O5'-P-OP2	-6.12	100.19	105.70
26	1H	2574	G	C5-C6-N1	6.12	114.56	111.50
1	1G	110	C	C6-N1-C2	6.12	122.75	120.30
26	14	1406	U	C5-C6-N1	6.12	125.76	122.70
1	13	807	A	C8-N9-C4	-6.12	103.35	105.80
26	1H	1364	G	N3-C2-N2	6.12	124.19	119.90
1	1G	905	U	C5-C6-N1	-6.12	119.64	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	1470	G	N3-C2-N2	-6.12	115.61	119.90
26	1H	238	C	C4-C5-C6	6.12	120.46	117.40
26	1H	917	A	C5-C6-N1	-6.12	114.64	117.70
1	13	1519	A	C4-C5-N7	-6.12	107.64	110.70
26	1H	2380	C	C6-N1-C2	6.12	122.75	120.30
26	14	639	U	N3-C4-C5	-6.12	110.93	114.60
26	14	1309	G	C4-N9-C1'	6.12	134.46	126.50
26	14	748	G	C8-N9-C1'	6.12	134.96	127.00
1	13	812	C	P-O3'-C3'	6.12	127.04	119.70
26	1H	1574	C	C6-N1-C2	6.12	122.75	120.30
26	1H	1614	A	C6-C5-N7	-6.12	128.02	132.30
26	1H	2045	C	C5-C6-N1	-6.12	117.94	121.00
26	1H	2324	C	C6-N1-C1'	-6.12	113.46	120.80
26	1H	197	A	P-O3'-C3'	6.11	127.04	119.70
26	1H	1386	C	C5-C6-N1	6.11	124.06	121.00
26	1H	2263	C	N3-C4-C5	-6.11	119.45	121.90
1	1G	1157	A	P-O3'-C3'	6.11	127.04	119.70
1	13	690	G	C8-N9-C1'	-6.11	119.06	127.00
26	14	621	A	N3-C4-C5	6.11	131.08	126.80
1	13	582	U	N3-C4-O4	-6.11	115.12	119.40
1	13	1431	C	C6-N1-C2	6.11	122.74	120.30
26	1H	2412	A	C6-N1-C2	-6.11	114.93	118.60
1	13	1432	G	C8-N9-C1'	-6.11	119.06	127.00
26	1H	205	G	C5-C6-O6	-6.11	124.94	128.60
26	14	36	G	C8-N9-C4	-6.11	103.96	106.40
26	14	1313	U	N3-C4-O4	6.11	123.67	119.40
26	14	1939	U	OP2-P-O3'	6.11	118.64	105.20
1	13	1310	G	N9-C4-C5	-6.11	102.96	105.40
26	1H	1158	C	N3-C4-C5	6.11	124.34	121.90
26	1H	1790	C	C5-C4-N4	-6.11	115.93	120.20
26	14	1954	G	C4-N9-C1'	-6.11	118.56	126.50
26	14	2045	C	O5'-P-OP2	-6.11	100.20	105.70
26	14	2329	G	N1-C6-O6	-6.11	116.24	119.90
27	1J	78	A	C8-N9-C4	6.11	108.24	105.80
26	1H	1294	U	O5'-P-OP2	6.10	118.03	110.70
26	1H	447	A	OP2-P-O3'	6.10	118.63	105.20
26	1H	564	C	OP1-P-O3'	6.10	118.63	105.20
26	1H	2245	U	O5'-P-OP1	-6.10	100.21	105.70
26	1H	2419	U	O5'-P-OP2	6.10	118.02	110.70
26	1H	2439	A	N7-C8-N9	6.10	116.85	113.80
26	1H	2496	C	N3-C2-O2	6.10	126.17	121.90
1	1G	690	G	O4'-C1'-N9	6.10	113.08	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2346	A	C4-C5-C6	6.10	120.05	117.00
26	1H	651	G	N1-C6-O6	-6.10	116.24	119.90
26	1H	1775	U	OP1-P-O3'	6.10	118.62	105.20
1	1G	41	G	C8-N9-C4	6.10	108.84	106.40
26	14	613	U	C5-C4-O4	6.10	129.56	125.90
26	14	828	U	N3-C4-C5	-6.10	110.94	114.60
26	14	1939	U	C5-C6-N1	-6.10	119.65	122.70
1	13	690	G	C4-C5-C6	6.10	122.46	118.80
26	1H	470	A	N9-C4-C5	-6.10	103.36	105.80
26	1H	733	G	O5'-P-OP2	-6.10	100.21	105.70
26	1H	738	G	C6-C5-N7	-6.10	126.74	130.40
26	1H	1558	A	C4-C5-N7	6.10	113.75	110.70
26	1H	1660	C	C2-N3-C4	-6.10	116.85	119.90
26	14	462	C	O5'-P-OP2	-6.10	100.21	105.70
26	14	1977	A	OP2-P-O3'	6.10	118.61	105.20
1	13	797	C	N1-C2-O2	-6.10	115.24	118.90
26	1H	461	C	N3-C2-O2	6.10	126.17	121.90
26	1H	2355	C	O5'-P-OP2	6.10	118.02	110.70
26	14	124	G	N1-C6-O6	-6.10	116.24	119.90
26	1H	777	A	C2-N3-C4	-6.09	107.55	110.60
26	1H	1424	G	O5'-P-OP2	-6.09	100.22	105.70
26	1H	1936	A	C5-C6-N1	6.09	120.75	117.70
26	1H	1972	A	N1-C6-N6	6.09	122.26	118.60
1	1G	1499	A	N7-C8-N9	-6.09	110.75	113.80
26	14	1528	A	C5-N7-C8	-6.09	100.85	103.90
26	14	2584	U	N1-C2-O2	6.09	127.07	122.80
1	13	266	G	C5-N7-C8	-6.09	101.25	104.30
1	13	834	C	O5'-P-OP2	-6.09	100.22	105.70
1	13	1323	G	O5'-P-OP1	-6.09	100.22	105.70
26	1H	729	G	C6-C5-N7	-6.09	126.75	130.40
1	1G	729	A	C8-N9-C4	-6.09	103.36	105.80
26	1H	1332	G	N9-C4-C5	6.09	107.84	105.40
26	1H	1647	G	O4'-C1'-N9	-6.09	103.33	108.20
26	1H	1799	G	N3-C2-N2	6.09	124.16	119.90
1	1G	111	G	N1-C6-O6	6.09	123.55	119.90
1	1G	481	G	N3-C4-C5	-6.09	125.56	128.60
1	1G	1502	A	C5-N7-C8	-6.09	100.86	103.90
26	14	1968	G	C5-C6-O6	-6.09	124.95	128.60
26	1H	739	G	C5-C6-O6	-6.09	124.95	128.60
1	1G	738	C	C6-N1-C2	-6.09	117.86	120.30
26	14	752	A	N7-C8-N9	6.09	116.84	113.80
26	14	830	G	N7-C8-N9	-6.09	110.06	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	933	A	C5-N7-C8	-6.09	100.86	103.90
26	14	2581	G	O5'-P-OP2	-6.09	100.22	105.70
26	14	2075	U	OP2-P-O3'	6.08	118.59	105.20
26	1H	422	A	OP1-P-OP2	6.08	128.72	119.60
26	1H	731	C	N3-C4-N4	6.08	122.26	118.00
1	1G	1147	C	C2-N1-C1'	6.08	125.49	118.80
26	14	211	A	N1-C6-N6	6.08	122.25	118.60
26	1H	2075	U	N1-C2-O2	-6.08	118.54	122.80
26	14	1203	G	O5'-P-OP2	-6.08	100.23	105.70
26	14	2689	U	OP2-P-O3'	6.08	118.58	105.20
1	13	781	A	C5-N7-C8	-6.08	100.86	103.90
1	13	926	G	C8-N9-C4	-6.08	103.97	106.40
1	1G	1490	C	O5'-P-OP1	6.08	118.00	110.70
26	14	2275	C	C5'-C4'-O4'	-6.08	101.80	109.10
26	1H	1451	C	N1-C2-O2	6.08	122.55	118.90
26	1H	1857	G	N1-C6-O6	6.08	123.55	119.90
26	14	2056	G	C5-C6-N1	6.08	114.54	111.50
26	1H	789	A	OP2-P-O3'	6.08	118.57	105.20
26	1H	830	G	N1-C2-N3	6.08	127.55	123.90
26	1H	141	A	C6-C5-N7	-6.07	128.05	132.30
26	1H	1382	G	C5-C6-N1	6.07	114.54	111.50
26	1H	1691	C	OP1-P-O3'	6.07	118.56	105.20
26	1H	1786	A	O4'-C1'-N9	6.07	113.06	108.20
26	1H	1956	U	OP1-P-O3'	6.07	118.56	105.20
26	1H	1987	G	N3-C2-N2	-6.07	115.65	119.90
26	1H	453	C	N3-C2-O2	6.07	126.15	121.90
26	1H	1834	U	N1-C2-O2	6.07	127.05	122.80
26	14	1920	C	O5'-P-OP2	-6.07	100.23	105.70
1	13	377	G	O5'-P-OP2	-6.07	100.24	105.70
1	13	1500	A	C5-C6-N1	6.07	120.73	117.70
26	1H	1518	C	O5'-P-OP2	6.07	117.98	110.70
26	1H	2325	G	O5'-P-OP1	-6.07	100.24	105.70
26	1H	2527	C	N3-C4-N4	6.07	122.25	118.00
27	16	44	G	N7-C8-N9	-6.07	110.06	113.10
26	14	2210	G	C8-N9-C1'	-6.07	119.11	127.00
26	14	2586	C	C5-C6-N1	6.07	124.03	121.00
26	1H	543	C	C6-N1-C2	6.07	122.73	120.30
1	1G	884	U	C6-N1-C2	-6.07	117.36	121.00
2	1E	187	LEU	CA-CB-CG	6.07	129.25	115.30
26	1H	107	C	C4-C5-C6	6.07	120.43	117.40
26	1H	1673	U	C6-N1-C2	6.07	124.64	121.00
26	1H	2269	A	O5'-P-OP1	6.07	117.98	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2477	C	C2-N1-C1'	6.07	125.47	118.80
1	1G	898	G	C5-C6-O6	-6.07	124.96	128.60
26	14	698	C	C4-C5-C6	6.07	120.43	117.40
1	13	567	G	O5'-P-OP1	-6.07	100.24	105.70
26	1H	213	A	O5'-P-OP2	-6.07	100.24	105.70
1	1G	230	G	N1-C6-O6	6.07	123.54	119.90
26	14	689	A	C8-N9-C4	6.06	108.23	105.80
26	14	784	A	C4-C5-N7	-6.06	107.67	110.70
1	13	190	G	C4-N9-C1'	6.06	134.38	126.50
1	13	758	G	N1-C6-O6	6.06	123.54	119.90
1	1G	811	C	C5-C6-N1	6.06	124.03	121.00
26	14	840	C	C6-N1-C2	6.06	122.72	120.30
26	1H	2241	A	N1-C2-N3	6.06	132.33	129.30
1	13	122	G	C5-C6-O6	-6.06	124.97	128.60
26	1H	1520	U	C6-N1-C2	-6.06	117.36	121.00
26	1H	1941	C	N3-C4-C5	-6.06	119.48	121.90
26	1H	2684	U	O5'-P-OP1	-6.06	100.25	105.70
26	14	664	C	OP1-P-OP2	6.06	128.69	119.60
26	14	1783	A	C8-N9-C4	-6.06	103.38	105.80
26	1H	557	U	C5-C6-N1	-6.06	119.67	122.70
26	1H	1786	A	C5-C6-N6	-6.06	118.85	123.70
26	1H	1825	A	C5-C6-N6	6.06	128.54	123.70
1	1G	904	C	O5'-P-OP2	6.06	117.97	110.70
26	14	270(X)	G	C5-C6-N1	-6.06	108.47	111.50
27	1J	47	C	OP1-P-O3'	6.06	118.53	105.20
26	14	71	A	C6-C5-N7	-6.06	128.06	132.30
26	14	543	C	N1-C2-O2	6.06	122.53	118.90
26	1H	2017	U	N3-C4-O4	6.05	123.64	119.40
1	1G	400	C	C6-N1-C2	6.05	122.72	120.30
1	1G	576	G	C8-N9-C1'	-6.05	119.13	127.00
26	14	1302	A	N1-C6-N6	-6.05	114.97	118.60
26	14	1780	A	N1-C2-N3	6.05	132.33	129.30
26	14	2439	A	C5-C6-N6	-6.05	118.86	123.70
1	1G	108	G	N9-C4-C5	-6.05	102.98	105.40
26	14	759	G	N3-C2-N2	-6.05	115.66	119.90
26	14	1795	C	N3-C2-O2	-6.05	117.66	121.90
26	14	2071	A	C6-N1-C2	-6.05	114.97	118.60
27	1J	102	G	C8-N9-C4	6.05	108.82	106.40
1	13	528	C	N3-C4-C5	6.05	124.32	121.90
1	13	821	G	N3-C4-N9	6.05	129.63	126.00
26	1H	82	G	OP1-P-O3'	6.05	118.51	105.20
26	1H	2582	G	C5-C6-O6	-6.05	124.97	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	600	G	O5'-P-OP2	-6.05	100.26	105.70
26	1H	1187	G	C5-C6-N1	-6.05	108.47	111.50
26	1H	1836	C	OP1-P-O3'	6.05	118.51	105.20
26	1H	2449	U	C2-N1-C1'	6.05	124.96	117.70
1	1G	652	U	C2-N1-C1'	6.05	124.96	117.70
26	14	1356	G	O5'-P-OP1	-6.05	100.25	105.70
26	1H	830	G	C8-N9-C4	-6.05	103.98	106.40
26	1H	1661	G	N7-C8-N9	-6.05	110.08	113.10
26	14	2261	C	N3-C4-N4	-6.05	113.77	118.00
26	1H	2286	A	C4-C5-C6	6.05	120.02	117.00
26	14	53	A	OP1-P-O3'	6.05	118.50	105.20
26	14	2037	G	N1-C6-O6	-6.05	116.27	119.90
26	14	2353	G	N1-C6-O6	-6.05	116.27	119.90
1	13	1520	G	C5-C6-O6	-6.04	124.97	128.60
26	1H	835	A	N1-C2-N3	6.04	132.32	129.30
26	1H	2623	G	N3-C4-C5	-6.04	125.58	128.60
26	14	1528	A	C4-C5-N7	6.04	113.72	110.70
1	13	1227	A	N3-C4-N9	-6.04	122.56	127.40
1	13	1234	C	C6-N1-C2	-6.04	117.88	120.30
26	1H	222	A	P-O3'-C3'	6.04	126.95	119.70
26	1H	928	G	N3-C2-N2	-6.04	115.67	119.90
26	1H	2672	G	N7-C8-N9	6.04	116.12	113.10
27	16	56	G	C4-N9-C1'	6.04	134.36	126.50
26	14	1322	A	OP2-P-O3'	6.04	118.50	105.20
26	14	2431	U	N3-C2-O2	6.04	126.43	122.20
26	14	2441	C	N3-C4-N4	-6.04	113.77	118.00
1	13	1299	A	C6-C5-N7	-6.04	128.07	132.30
26	1H	203	C	N3-C4-C5	6.04	124.32	121.90
26	1H	210	C	C5-C6-N1	-6.04	117.98	121.00
26	1H	1283	G	O5'-P-OP2	-6.04	100.26	105.70
26	1H	2485	G	N1-C2-N2	-6.04	110.76	116.20
27	16	109	G	C8-N9-C4	-6.04	103.98	106.40
26	14	379	G	O5'-P-OP1	-6.04	100.26	105.70
26	14	2061	G	C8-N9-C4	6.04	108.82	106.40
26	14	2346	A	C6-C5-N7	-6.04	128.07	132.30
1	13	253	U	OP2-P-O3'	6.04	118.49	105.20
26	1H	706	A	O5'-P-OP1	-6.04	100.26	105.70
26	1H	1475	G	N3-C2-N2	-6.04	115.67	119.90
26	1H	784	A	C8-N9-C4	-6.04	103.39	105.80
26	1H	2431	U	OP2-P-O3'	-6.04	91.92	105.20
26	14	2779	U	C2-N3-C4	-6.04	123.38	127.00
26	1H	733	G	C8-N9-C1'	-6.04	119.15	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	629	G	O5'-P-OP2	-6.04	100.27	105.70
26	14	1324	G	N1-C6-O6	6.04	123.52	119.90
26	1H	1122	G	N7-C8-N9	-6.04	110.08	113.10
26	1H	1900	A	C5'-C4'-O4'	-6.04	101.86	109.10
26	1H	2690	C	N1-C2-O2	-6.04	115.28	118.90
26	1H	2838	G	O5'-P-OP1	-6.04	100.27	105.70
26	14	784	A	N3-C4-N9	-6.04	122.57	127.40
26	14	2712	U	O4'-C1'-N1	6.04	113.03	108.20
1	13	1416	G	O5'-P-OP1	-6.03	100.27	105.70
26	1H	46	C	N1-C2-O2	-6.03	115.28	118.90
26	1H	920	G	N7-C8-N9	-6.03	110.08	113.10
26	14	945	A	O4'-C1'-N9	6.03	113.03	108.20
26	14	1527	G	N3-C4-N9	-6.03	122.38	126.00
26	14	2622	C	C2-N1-C1'	-6.03	112.16	118.80
26	1H	698	C	C2-N3-C4	-6.03	116.88	119.90
26	1H	2230	G	N1-C2-N2	6.03	121.63	116.20
1	1G	275	G	N1-C6-O6	6.03	123.52	119.90
26	14	1204	A	C5-N7-C8	-6.03	100.88	103.90
26	1H	131	G	N1-C6-O6	6.03	123.52	119.90
26	1H	528	A	C8-N9-C4	-6.03	103.39	105.80
26	1H	2604	U	N1-C2-O2	6.03	127.02	122.80
26	14	833	U	C4-C5-C6	6.03	123.32	119.70
26	14	2044	C	N3-C2-O2	6.03	126.12	121.90
26	1H	512	G	OP1-P-O3'	6.03	118.47	105.20
26	1H	1265	A	N1-C2-N3	6.03	132.31	129.30
26	1H	1977	A	C5-C6-N1	-6.03	114.69	117.70
26	1H	2028	U	N3-C2-O2	-6.03	117.98	122.20
26	1H	2375	G	C5-C6-N1	6.03	114.51	111.50
26	14	2584	U	O5'-P-OP2	6.03	117.94	110.70
1	13	742	G	C8-N9-C4	6.03	108.81	106.40
26	1H	447	A	N7-C8-N9	6.03	116.81	113.80
26	1H	2275	C	OP1-P-O3'	6.03	118.46	105.20
26	1H	2567	G	O5'-P-OP1	-6.03	100.28	105.70
26	14	250	G	O5'-P-OP1	-6.03	100.28	105.70
26	14	1383	C	O4'-C1'-N1	6.03	113.02	108.20
26	14	2423	U	C6-N1-C2	6.03	124.62	121.00
1	13	564	C	O5'-P-OP1	-6.03	100.28	105.70
26	1H	420	C	C6-N1-C2	6.03	122.71	120.30
26	1H	2830	G	O5'-P-OP2	-6.03	100.28	105.70
26	1H	694	U	O5'-P-OP1	6.02	117.93	110.70
1	13	1307	U	C6-N1-C2	-6.02	117.39	121.00
26	1H	67	U	OP1-P-O3'	6.02	118.45	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1228	G	N3-C2-N2	-6.02	115.68	119.90
26	1H	2490	G	C6-C5-N7	-6.02	126.79	130.40
26	14	2023	G	OP1-P-OP2	6.02	128.63	119.60
26	14	2244	U	N1-C2-N3	6.02	118.51	114.90
23	2K	6	G	N7-C8-N9	-6.02	110.09	113.10
1	1G	690	G	C8-N9-C4	-6.02	103.99	106.40
26	14	2542	A	C5-N7-C8	6.02	106.91	103.90
26	14	2572	A	O5'-P-OP1	-6.02	100.28	105.70
1	13	1103	C	C6-N1-C2	-6.02	117.89	120.30
26	1H	450	G	C5-C6-N1	-6.02	108.49	111.50
26	1H	838	C	C2-N3-C4	-6.02	116.89	119.90
27	16	5	C	C5-C6-N1	-6.02	117.99	121.00
26	14	559	G	C5-C6-N1	-6.02	108.49	111.50
26	1H	1814	G	N1-C2-N2	-6.02	110.78	116.20
26	1H	2433	A	O5'-P-OP2	6.02	117.92	110.70
26	1H	2489	G	N1-C2-N3	6.02	127.51	123.90
27	16	79	C	C5-C6-N1	6.02	124.01	121.00
28	11	271	ILE	N-CA-C	6.02	127.25	111.00
26	14	759	G	C6-N1-C2	-6.02	121.49	125.10
1	13	123	C	O5'-P-OP2	-6.01	100.29	105.70
1	13	1500	A	N1-C6-N6	-6.01	114.99	118.60
26	1H	990	A	C2-N3-C4	-6.01	107.59	110.60
26	1H	1658	C	N1-C2-O2	-6.01	115.29	118.90
26	1H	1969	A	C4-C5-N7	-6.01	107.69	110.70
26	1H	2285	C	C4-C5-C6	-6.01	114.39	117.40
26	1H	2819	G	O5'-P-OP2	-6.01	100.29	105.70
1	1G	586	C	C5-C6-N1	6.01	124.01	121.00
26	1H	610	C	N3-C2-O2	-6.01	117.69	121.90
26	14	674	G	C5-C6-N1	6.01	114.51	111.50
1	13	1520	G	C8-N9-C4	6.01	108.80	106.40
26	1H	118	A	O5'-P-OP1	-6.01	100.29	105.70
26	1H	1987	G	N1-C6-O6	6.01	123.51	119.90
1	1G	945	G	N1-C6-O6	6.01	123.51	119.90
26	14	197	A	OP2-P-O3'	6.01	118.43	105.20
26	14	2423	U	C5-C6-N1	-6.01	119.69	122.70
47	E5	14	ARG	NE-CZ-NH1	-6.01	117.29	120.30
26	1H	25	U	N3-C4-C5	6.01	118.21	114.60
26	1H	768	G	O5'-P-OP2	-6.01	100.29	105.70
26	1H	655	A	C6-C5-N7	-6.01	128.09	132.30
26	1H	2616	C	O5'-P-OP2	6.01	117.91	110.70
1	13	1187	G	C4-N9-C1'	6.01	134.31	126.50
26	1H	1141	U	O4'-C1'-N1	6.01	113.00	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2422	A	O4'-C1'-N9	6.01	113.00	108.20
26	14	693	C	C2-N1-C1'	-6.01	112.19	118.80
26	14	1420	U	N3-C2-O2	-6.01	118.00	122.20
26	14	2502	G	N7-C8-N9	6.01	116.10	113.10
26	14	2523	G	N9-C4-C5	-6.01	103.00	105.40
26	1H	686	G	N3-C4-N9	6.00	129.60	126.00
1	13	285	G	N9-C4-C5	-6.00	103.00	105.40
26	1H	729	G	OP2-P-O3'	6.00	118.41	105.20
26	1H	792	G	C8-N9-C1'	-6.00	119.20	127.00
26	1H	2328	A	C2-N3-C4	-6.00	107.60	110.60
26	1H	2594	C	C6-N1-C2	-6.00	117.90	120.30
26	14	951	C	OP1-P-O3'	6.00	118.41	105.20
26	14	2277	G	O5'-P-OP2	-6.00	100.30	105.70
1	13	964	A	C8-N9-C4	6.00	108.20	105.80
26	1H	456	C	C2-N1-C1'	-6.00	112.20	118.80
26	1H	732	C	C4-C5-C6	6.00	120.40	117.40
26	14	2319	G	N1-C6-O6	-6.00	116.30	119.90
26	1H	429	A	O5'-P-OP1	-6.00	100.30	105.70
26	1H	688	U	OP2-P-O3'	6.00	118.40	105.20
26	1H	470	A	O5'-P-OP1	-6.00	100.30	105.70
26	14	1283	G	OP1-P-OP2	6.00	128.60	119.60
26	14	1402	C	N3-C4-C5	-6.00	119.50	121.90
23	2K	17	C	C6-N1-C1'	-6.00	113.60	120.80
26	1H	331	A	OP1-P-O3'	6.00	118.39	105.20
26	1H	840	C	C5-C6-N1	-6.00	118.00	121.00
26	14	1623	G	C6-C5-N7	6.00	134.00	130.40
27	1J	103	U	O5'-P-OP2	-6.00	100.30	105.70
26	1H	1252	G	N7-C8-N9	-6.00	110.10	113.10
26	1H	2318	G	N3-C4-C5	6.00	131.60	128.60
27	16	12	C	C5-C6-N1	-6.00	118.00	121.00
27	16	81	G	C6-C5-N7	-6.00	126.80	130.40
26	14	636	G	O5'-P-OP1	-6.00	100.30	105.70
26	14	827	U	N3-C2-O2	6.00	126.40	122.20
26	14	1619	G	C5-C6-N1	6.00	114.50	111.50
26	1H	299	A	OP2-P-O3'	5.99	118.39	105.20
26	1H	447	A	N9-C4-C5	5.99	108.20	105.80
26	1H	2743	C	C2-N3-C4	-5.99	116.90	119.90
1	1G	1285	A	P-O3'-C3'	5.99	126.89	119.70
26	14	1930	G	C4-C5-N7	-5.99	108.40	110.80
26	14	2267	A	O5'-P-OP2	-5.99	100.31	105.70
26	14	2623	G	O5'-P-OP1	5.99	117.89	110.70
26	1H	451	C	N1-C2-O2	-5.99	115.31	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	803	U	C5-C6-N1	-5.99	119.70	122.70
26	1H	1990	C	N3-C4-C5	-5.99	119.50	121.90
26	14	1975	G	C4-C5-N7	5.99	113.20	110.80
26	14	2297	C	O5'-P-OP1	-5.99	100.31	105.70
26	1H	568	U	C6-N1-C2	5.99	124.59	121.00
26	1H	662	G	OP1-P-OP2	5.99	128.59	119.60
26	1H	856	C	O5'-P-OP1	-5.99	100.31	105.70
26	14	1371	G	C5-C6-O6	-5.99	125.01	128.60
26	14	1379	A	C5-C6-N6	-5.99	118.91	123.70
26	1H	639	U	O5'-P-OP2	-5.99	100.31	105.70
26	14	244	A	O5'-P-OP1	5.99	117.89	110.70
26	14	511	U	C6-N1-C2	-5.99	117.41	121.00
26	14	568	U	C6-N1-C2	5.99	124.59	121.00
26	14	1616	A	N1-C6-N6	5.99	122.19	118.60
26	14	1678	G	C4-C5-N7	5.99	113.19	110.80
26	1H	2582	G	C4-C5-N7	5.99	113.19	110.80
1	13	970	C	C5-C6-N1	5.99	123.99	121.00
1	13	1371	G	O5'-P-OP1	-5.99	100.31	105.70
26	1H	209	C	C5-C6-N1	-5.99	118.01	121.00
26	1H	1035	U	C5-C4-O4	5.99	129.49	125.90
26	1H	1636	C	N3-C2-O2	5.99	126.09	121.90
26	1H	1671	U	N3-C4-O4	5.99	123.59	119.40
26	1H	2568	C	OP2-P-O3'	5.99	118.37	105.20
26	14	1021	A	C8-N9-C4	-5.99	103.41	105.80
26	1H	127	A	C5-C6-N6	-5.98	118.91	123.70
1	13	889	A	OP1-P-OP2	5.98	128.57	119.60
26	1H	753	C	OP1-P-OP2	5.98	128.57	119.60
26	1H	792	G	N1-C6-O6	5.98	123.49	119.90
26	1H	799	G	N3-C4-C5	5.98	131.59	128.60
26	1H	1334	G	C4-N9-C1'	5.98	134.28	126.50
50	L8	54	VAL	N-CA-C	5.98	127.15	111.00
26	14	1648	C	C6-N1-C2	-5.98	117.91	120.30
26	14	1930	G	N9-C4-C5	5.98	107.79	105.40
26	14	1972	A	OP2-P-O3'	5.98	118.36	105.20
26	1H	1649	G	N3-C4-N9	5.98	129.59	126.00
26	1H	2276	G	C4-C5-N7	-5.98	108.41	110.80
26	1H	2280	G	C2-N3-C4	5.98	114.89	111.90
26	14	801	G	C6-C5-N7	5.98	133.99	130.40
26	1H	1380	G	N3-C2-N2	5.98	124.08	119.90
26	1H	1437	C	N1-C2-O2	5.98	122.49	118.90
26	1H	1971	A	C2-N3-C4	5.98	113.59	110.60
26	1H	2424	C	C4-C5-C6	-5.98	114.41	117.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2726	U	N3-C2-O2	-5.98	118.02	122.20
26	1H	2731	G	C4-C5-N7	5.98	113.19	110.80
26	14	488	G	N7-C8-N9	-5.98	110.11	113.10
26	14	929	G	N3-C4-C5	-5.98	125.61	128.60
26	14	2433	A	C5-N7-C8	-5.98	100.91	103.90
26	1H	730	C	N3-C4-C5	5.98	124.29	121.90
26	1H	2712	U	N1-C2-N3	5.98	118.49	114.90
26	1H	62	C	C2-N3-C4	-5.97	116.91	119.90
26	1H	1777	U	O5'-P-OP1	-5.97	100.32	105.70
26	1H	1780	A	N9-C4-C5	5.97	108.19	105.80
26	14	939	G	N1-C6-O6	5.97	123.48	119.90
26	14	1391	U	O5'-P-OP2	5.97	117.87	110.70
26	14	1728	G	N3-C4-C5	-5.97	125.61	128.60
26	14	2023	G	C6-C5-N7	-5.97	126.81	130.40
27	1J	7	G	N9-C4-C5	-5.97	103.01	105.40
1	13	985	C	O5'-P-OP2	5.97	117.87	110.70
26	1H	633	A	C4-C5-N7	5.97	113.69	110.70
26	1H	1245	G	C6-N1-C2	-5.97	121.52	125.10
26	1H	2376	A	C8-N9-C4	5.97	108.19	105.80
26	1H	2577	A	N9-C4-C5	5.97	108.19	105.80
26	14	2332	U	N3-C2-O2	-5.97	118.02	122.20
26	14	2401	U	C2-N1-C1'	5.97	124.87	117.70
26	14	2560	C	O5'-P-OP1	-5.97	100.32	105.70
26	1H	2310	A	N3-C4-C5	-5.97	122.62	126.80
26	1H	2468	G	C8-N9-C1'	-5.97	119.24	127.00
26	1H	2574	G	C5-C6-O6	-5.97	125.02	128.60
55	Q8	49	VAL	CA-CB-CG2	5.97	119.86	110.90
26	14	1938	A	N1-C2-N3	5.97	132.29	129.30
26	14	2452	C	C5-C4-N4	-5.97	116.02	120.20
1	13	899	C	C5-C6-N1	-5.97	118.02	121.00
1	13	1518	A	C8-N9-C4	5.97	108.19	105.80
26	1H	586	A	N1-C6-N6	-5.97	115.02	118.60
26	1H	1757	U	OP1-P-O3'	5.97	118.33	105.20
33	61	131	LYS	C-N-CD	-5.97	107.47	120.60
1	1G	1500	A	N1-C6-N6	5.97	122.18	118.60
26	14	576	U	OP2-P-O3'	5.97	118.33	105.20
26	14	1954	G	N3-C4-N9	-5.97	122.42	126.00
1	13	1025	U	C5-C6-N1	5.97	125.68	122.70
26	1H	1142(A)	A	C5-N7-C8	-5.97	100.92	103.90
26	1H	2582	G	O5'-P-OP1	5.97	117.86	110.70
26	14	1402	C	N3-C4-N4	5.97	122.18	118.00
26	14	2702	U	C6-N1-C2	-5.97	117.42	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	293	U	N3-C4-O4	5.97	123.58	119.40
26	1H	698	C	N3-C2-O2	5.97	126.08	121.90
26	1H	808	G	OP1-P-OP2	5.97	128.55	119.60
26	1H	1380	G	N1-C2-N2	-5.97	110.83	116.20
26	14	756	C	O5'-P-OP1	-5.97	100.33	105.70
26	1H	114	U	OP1-P-O3'	5.96	118.32	105.20
26	1H	840	C	C2-N3-C4	-5.96	116.92	119.90
26	1H	1142(A)	A	N3-C4-N9	-5.96	122.63	127.40
26	1H	1199	U	C2-N3-C4	-5.96	123.42	127.00
26	1H	1380	G	C6-C5-N7	-5.96	126.82	130.40
26	14	2491	U	OP1-P-O3'	5.96	118.32	105.20
27	1J	81	G	C4-C5-N7	5.96	113.19	110.80
1	13	740	U	O5'-P-OP2	-5.96	100.33	105.70
26	1H	973	A	N3-C4-C5	5.96	130.97	126.80
26	1H	1776	G	N9-C4-C5	-5.96	103.02	105.40
26	1H	1973	G	N7-C8-N9	5.96	116.08	113.10
26	1H	2707	G	C4-N9-C1'	-5.96	118.75	126.50
26	14	694	U	N3-C2-O2	-5.96	118.03	122.20
26	1H	593	G	N1-C6-O6	5.96	123.48	119.90
26	1H	815	C	C5-C6-N1	-5.96	118.02	121.00
26	1H	2782	G	C6-C5-N7	-5.96	126.82	130.40
26	1H	128	C	O5'-P-OP1	5.96	117.85	110.70
26	1H	141	A	C2-N3-C4	-5.96	107.62	110.60
26	1H	733	G	N3-C4-N9	5.96	129.58	126.00
26	1H	1573	G	C8-N9-C4	5.96	108.78	106.40
26	1H	1975	G	N1-C6-O6	-5.96	116.32	119.90
26	1H	2256	G	C5-C6-O6	5.96	132.18	128.60
26	14	1313	U	N1-C2-N3	5.96	118.48	114.90
1	13	889	A	C4-C5-C6	5.96	119.98	117.00
1	13	1113	C	N1-C2-O2	5.96	122.47	118.90
26	14	193	U	C5-C4-O4	-5.96	122.33	125.90
26	14	245	G	O5'-P-OP1	-5.96	100.34	105.70
26	1H	194	G	N9-C4-C5	-5.96	103.02	105.40
26	1H	391	G	C5-C6-N1	-5.96	108.52	111.50
26	1H	1193	G	N1-C6-O6	-5.96	116.33	119.90
26	14	1994	C	O5'-P-OP2	-5.96	100.34	105.70
1	13	328	C	C2-N1-C1'	5.96	125.35	118.80
26	1H	583	G	C8-N9-C4	-5.96	104.02	106.40
26	1H	774	A	C4-C5-C6	-5.96	114.02	117.00
27	1J	100	G	N1-C2-N2	-5.96	110.84	116.20
1	13	1530	G	N3-C4-N9	-5.95	122.43	126.00
26	1H	851	U	OP2-P-O3'	5.95	118.30	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1413	G	N7-C8-N9	5.95	116.08	113.10
38	98	67	LEU	CA-CB-CG	5.95	128.99	115.30
26	14	268	C	C5-C6-N1	5.95	123.98	121.00
26	14	1852	C	O5'-P-OP1	-5.95	100.34	105.70
24	3K	32	U	N1-C2-O2	5.95	126.97	122.80
26	1H	1779	U	OP1-P-OP2	5.95	128.53	119.60
26	1H	1229(A)	G	N1-C6-O6	5.95	123.47	119.90
26	1H	1733	G	C5-C6-N1	-5.95	108.53	111.50
26	1H	2052	G	O5'-P-OP2	-5.95	100.34	105.70
26	1H	2609	U	N1-C2-O2	-5.95	118.63	122.80
27	16	78	A	C8-N9-C4	5.95	108.18	105.80
1	1G	1499	A	O5'-P-OP1	-5.95	100.34	105.70
26	14	2424	C	O5'-P-OP1	-5.95	100.34	105.70
1	13	811	C	N1-C2-O2	-5.95	115.33	118.90
1	13	1498	U	N3-C2-O2	-5.95	118.03	122.20
26	1H	743	G	OP1-P-OP2	5.95	128.52	119.60
26	14	1432	C	C6-N1-C2	5.95	122.68	120.30
24	3K	71	G	C5'-C4'-O4'	5.95	116.24	109.10
26	1H	571	A	C2-N3-C4	-5.95	107.63	110.60
26	1H	835	A	N9-C4-C5	5.95	108.18	105.80
26	1H	1365	A	C2-N3-C4	-5.95	107.63	110.60
26	1H	649	G	C6-C5-N7	-5.95	126.83	130.40
26	1H	736	C	N3-C4-C5	5.95	124.28	121.90
26	1H	812	C	N3-C2-O2	5.95	126.06	121.90
26	1H	1303	G	OP1-P-O3'	-5.95	92.12	105.20
26	1H	1792	G	N1-C6-O6	-5.95	116.33	119.90
26	1H	2061	G	N1-C6-O6	-5.95	116.33	119.90
26	14	2321	G	C8-N9-C4	-5.95	104.02	106.40
26	1H	256	A	N1-C6-N6	5.94	122.17	118.60
26	1H	740	U	O5'-P-OP1	5.94	117.83	110.70
26	1H	2550	G	C6-C5-N7	-5.94	126.83	130.40
26	14	2253	G	O5'-P-OP1	5.94	117.83	110.70
1	13	266	G	N7-C8-N9	5.94	116.07	113.10
1	13	1525	G	C2-N3-C4	-5.94	108.93	111.90
26	1H	1790	C	N3-C2-O2	5.94	126.06	121.90
26	14	2426	A	C6-C5-N7	-5.94	128.14	132.30
1	13	623	C	C5-C6-N1	5.94	123.97	121.00
23	2K	74	A	C6-C5-N7	-5.94	128.14	132.30
26	1H	1325	G	N3-C4-C5	-5.94	125.63	128.60
1	1G	1435	G	O5'-P-OP2	-5.94	100.35	105.70
1	1G	1488	G	N3-C4-N9	5.94	129.56	126.00
26	14	827	U	C2-N3-C4	-5.94	123.44	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1128	A	O5'-P-OP1	-5.94	100.35	105.70
25	4K	16	A	C8-N9-C4	5.94	108.18	105.80
26	1H	2435	A	C5-N7-C8	-5.94	100.93	103.90
1	1G	306	G	N3-C4-C5	5.94	131.57	128.60
1	13	169	C	C6-N1-C2	-5.94	117.92	120.30
14	5I	44	LEU	CA-CB-CG	5.94	128.96	115.30
26	1H	328	U	C6-N1-C2	-5.94	117.44	121.00
26	1H	1382	G	OP2-P-O3'	5.94	118.27	105.20
26	1H	1425	G	C2-N3-C4	5.94	114.87	111.90
26	1H	2666	C	C6-N1-C2	-5.94	117.92	120.30
1	1G	862	C	N1-C2-O2	5.94	122.46	118.90
26	1H	1957	C	C5-C6-N1	-5.94	118.03	121.00
26	14	72	U	N3-C2-O2	-5.94	118.05	122.20
26	1H	481	G	O4'-C1'-N9	5.93	112.95	108.20
26	1H	1966	A	O5'-P-OP2	-5.93	100.36	105.70
27	16	111	U	C5-C4-O4	5.93	129.46	125.90
26	14	479	A	N1-C6-N6	-5.93	115.04	118.60
1	13	1331	G	O5'-P-OP2	-5.93	100.36	105.70
26	14	2577	A	C6-C5-N7	-5.93	128.15	132.30
26	14	2731	G	C8-N9-C4	-5.93	104.03	106.40
26	1H	205	G	N3-C4-N9	5.93	129.56	126.00
26	14	664	C	C2-N3-C4	-5.93	116.93	119.90
26	1H	1274	A	O5'-P-OP1	-5.93	100.36	105.70
26	1H	1698	A	O4'-C1'-N9	5.93	112.94	108.20
26	14	130	C	C5-C4-N4	-5.93	116.05	120.20
26	14	565	C	C6-N1-C2	5.93	122.67	120.30
26	14	2539	C	C6-N1-C2	5.93	122.67	120.30
26	14	2612	C	N1-C2-O2	5.93	122.46	118.90
26	1H	1779	U	C5-C6-N1	-5.93	119.74	122.70
26	14	1328	G	N9-C4-C5	-5.93	103.03	105.40
26	14	1678	G	C8-N9-C4	-5.93	104.03	106.40
26	14	2087	G	N9-C4-C5	-5.93	103.03	105.40
1	13	1489	G	N7-C8-N9	-5.93	110.14	113.10
1	13	1519	A	C8-N9-C4	-5.93	103.43	105.80
26	1H	103	A	N7-C8-N9	-5.93	110.84	113.80
26	1H	787	U	OP1-P-OP2	-5.93	110.71	119.60
26	1H	1446	C	C6-N1-C2	-5.93	117.93	120.30
26	1H	2523	G	N9-C4-C5	5.93	107.77	105.40
26	14	247	G	N3-C4-C5	5.93	131.56	128.60
26	1H	142	G	N3-C4-C5	5.92	131.56	128.60
26	1H	852	G	OP2-P-O3'	5.92	118.23	105.20
26	1H	1204	A	C5-C6-N1	-5.92	114.74	117.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1368	G	C2-N3-C4	5.92	114.86	111.90
26	1H	2707	G	N3-C2-N2	-5.92	115.75	119.90
1	1G	586	C	C6-N1-C2	-5.92	117.93	120.30
1	1G	662	G	N3-C4-C5	-5.92	125.64	128.60
26	14	1496	A	C4-C5-N7	5.92	113.66	110.70
26	14	1961	C	C2-N1-C1'	-5.92	112.28	118.80
26	14	2604	U	C5-C6-N1	-5.92	119.74	122.70
1	13	738	C	N3-C4-C5	-5.92	119.53	121.90
26	1H	239	U	C5-C6-N1	-5.92	119.74	122.70
26	1H	139	G	C2-N3-C4	5.92	114.86	111.90
26	1H	291	C	C6-N1-C2	5.92	122.67	120.30
26	1H	690	G	C2-N3-C4	-5.92	108.94	111.90
26	1H	731	C	C2-N1-C1'	5.92	125.31	118.80
26	1H	1232	G	N1-C6-O6	5.92	123.45	119.90
26	1H	2068	U	C2-N3-C4	5.92	130.55	127.00
26	1H	2373	G	C6-N1-C2	-5.92	121.55	125.10
26	1H	417	C	O5'-P-OP2	5.92	117.80	110.70
26	1H	609	A	C5-C6-N6	-5.92	118.96	123.70
1	13	582	U	C5-C6-N1	-5.92	119.74	122.70
26	1H	1624	G	N7-C8-N9	-5.92	110.14	113.10
26	1H	2351	G	C4-N9-C1'	5.92	134.19	126.50
26	14	698	C	N3-C4-N4	5.92	122.14	118.00
26	14	2592	G	O5'-P-OP1	5.92	117.80	110.70
26	1H	461	C	N3-C4-N4	5.92	122.14	118.00
26	1H	775	G	N3-C4-C5	-5.92	125.64	128.60
26	1H	1300	U	C6-N1-C1'	5.92	129.48	121.20
26	1H	2518	A	C2-N3-C4	-5.92	107.64	110.60
1	1G	1346	A	OP2-P-O3'	5.92	118.22	105.20
1	1G	1475	G	OP1-P-OP2	5.92	128.47	119.60
26	14	380	U	O5'-P-OP2	-5.92	100.38	105.70
26	14	1342	A	N9-C1'-C2'	5.92	121.69	114.00
26	14	1409	C	O5'-P-OP2	-5.92	100.38	105.70
26	1H	1950	G	N1-C2-N2	-5.92	110.88	116.20
1	1G	528	C	O4'-C1'-N1	5.92	112.93	108.20
27	1J	22	U	C2-N1-C1'	5.92	124.80	117.70
1	13	1224	G	O5'-P-OP1	5.91	117.80	110.70
1	13	1417	G	N3-C4-N9	5.91	129.55	126.00
26	1H	259	G	C5-C6-O6	-5.91	125.05	128.60
26	1H	690	G	C8-N9-C4	5.91	108.77	106.40
26	1H	692	C	N3-C4-N4	5.91	122.14	118.00
26	1H	827	U	O5'-P-OP1	5.91	117.80	110.70
26	1H	1621	U	N3-C2-O2	5.91	126.34	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	972	C	C6-N1-C2	-5.91	117.93	120.30
26	14	186	G	N1-C6-O6	5.91	123.45	119.90
26	14	2363	C	C6-N1-C2	5.91	122.67	120.30
26	14	2406	U	P-O3'-C3'	5.91	126.80	119.70
26	14	2584	U	O4'-C1'-N1	5.91	112.93	108.20
26	1H	126	A	OP2-P-O3'	5.91	118.21	105.20
26	14	1320	C	N3-C4-C5	-5.91	119.53	121.90
26	14	1678	G	N3-C2-N2	-5.91	115.76	119.90
26	1H	398	G	OP1-P-OP2	5.91	128.47	119.60
26	1H	699	A	C5-C6-N6	-5.91	118.97	123.70
26	1H	1555	G	O5'-P-OP1	-5.91	100.38	105.70
26	1H	2269	A	C8-N9-C4	5.91	108.16	105.80
26	1H	2351	G	N3-C4-C5	-5.91	125.64	128.60
26	1H	119	A	N1-C6-N6	-5.91	115.06	118.60
26	1H	1159	U	O5'-P-OP2	-5.91	100.38	105.70
26	1H	1520	U	N1-C2-N3	5.91	118.44	114.90
26	1H	2269	A	N9-C4-C5	-5.91	103.44	105.80
26	1H	2618	G	N7-C8-N9	5.91	116.06	113.10
1	1G	345	C	C6-N1-C1'	-5.91	113.71	120.80
26	14	827	U	C5-C6-N1	-5.91	119.75	122.70
26	1H	458	G	OP1-P-OP2	5.91	128.46	119.60
26	1H	2328	A	C4-C5-C6	5.91	119.95	117.00
1	1G	150	C	C6-N1-C2	-5.91	117.94	120.30
1	1G	898	G	N9-C4-C5	-5.91	103.04	105.40
26	14	1268	A	C2-N3-C4	-5.91	107.65	110.60
26	14	2070	G	N1-C2-N3	5.91	127.44	123.90
26	14	2277	G	N1-C6-O6	-5.91	116.36	119.90
26	1H	752	A	C6-N1-C2	-5.90	115.06	118.60
26	1H	1825	A	N9-C4-C5	5.90	108.16	105.80
26	14	2377	A	N3-C4-C5	5.90	130.93	126.80
1	13	1436	U	O5'-P-OP2	5.90	117.78	110.70
26	1H	2231	C	C6-N1-C2	-5.90	117.94	120.30
26	1H	2335	A	O4'-C1'-N9	5.90	112.92	108.20
26	1H	2731	G	C8-N9-C1'	-5.90	119.33	127.00
1	1G	995	C	C6-N1-C2	-5.90	117.94	120.30
26	14	141	A	N7-C8-N9	5.90	116.75	113.80
1	13	1523	G	O5'-P-OP2	-5.90	100.39	105.70
26	1H	2244	U	C2-N3-C4	-5.90	123.46	127.00
26	14	278	A	OP1-P-O3'	5.90	118.18	105.20
26	1H	845	G	C5-N7-C8	-5.90	101.35	104.30
26	1H	1432	C	C4-C5-C6	5.90	120.35	117.40
26	1H	1626	G	N3-C4-N9	-5.90	122.46	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1993	U	C2-N3-C4	-5.90	123.46	127.00
26	1H	2581	G	O4'-C1'-N9	5.90	112.92	108.20
26	14	1807	G	N7-C8-N9	-5.90	110.15	113.10
26	1H	193	U	N1-C2-O2	-5.90	118.67	122.80
26	1H	370	G	O4'-C1'-N9	-5.90	103.48	108.20
26	1H	966	G	N3-C2-N2	5.90	124.03	119.90
26	1H	2483	C	O5'-P-OP1	-5.90	100.39	105.70
26	1H	607	U	N1-C2-N3	-5.90	111.36	114.90
26	1H	802	A	OP2-P-O3'	5.90	118.17	105.20
26	1H	824	A	C2-N3-C4	-5.90	107.65	110.60
26	1H	1204	A	C8-N9-C1'	-5.90	117.09	127.70
26	1H	1318	C	O5'-P-OP1	-5.90	100.39	105.70
26	14	537	C	C5-C6-N1	5.90	123.95	121.00
8	7E	71	GLY	N-CA-C	-5.89	98.36	113.10
26	1H	734	A	C5-C6-N1	-5.89	114.75	117.70
26	1H	734	A	C6-N1-C2	5.89	122.14	118.60
26	14	954	G	N3-C2-N2	-5.89	115.78	119.90
26	1H	88	G	C8-N9-C4	-5.89	104.04	106.40
26	1H	2383	G	C6-C5-N7	-5.89	126.86	130.40
26	1H	2550	G	C4-C5-C6	5.89	122.33	118.80
26	14	1383	C	O5'-P-OP2	-5.89	100.40	105.70
26	14	2388	A	O5'-P-OP1	5.89	117.77	110.70
26	1H	1194	A	C5-C6-N6	-5.89	118.99	123.70
26	14	827	U	O5'-P-OP1	5.89	117.77	110.70
26	14	1000	A	C8-N9-C4	-5.89	103.44	105.80
1	13	1281	U	N3-C2-O2	-5.89	118.08	122.20
26	1H	138	G	C5-C6-N1	5.89	114.44	111.50
26	1H	141	A	C8-N9-C4	-5.89	103.44	105.80
26	1H	148	C	C2-N3-C4	-5.89	116.95	119.90
26	14	1451	C	C6-N1-C2	5.89	122.66	120.30
26	14	1608	A	C5-C6-N6	5.89	128.41	123.70
26	14	1614	A	C6-C5-N7	-5.89	128.18	132.30
26	14	1771	C	C2-N3-C4	-5.89	116.95	119.90
1	13	1502	A	C5-C6-N1	-5.89	114.76	117.70
26	1H	435	C	C5-C6-N1	5.89	123.94	121.00
26	1H	733	G	C2-N3-C4	-5.89	108.96	111.90
26	1H	1341	U	C5-C4-O4	-5.89	122.37	125.90
26	14	1394	U	C6-N1-C1'	5.89	129.44	121.20
1	13	724	G	OP1-P-O3'	5.89	118.15	105.20
8	7E	112	LEU	CA-CB-CG	5.89	128.84	115.30
26	1H	94	G	C8-N9-C4	5.89	108.75	106.40
26	1H	763	G	C4-N9-C1'	5.89	134.15	126.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2312	U	C5-C4-O4	-5.89	122.37	125.90
26	1H	2361	A	C2-N3-C4	-5.89	107.66	110.60
1	13	291	C	C2-N1-C1'	-5.88	112.33	118.80
26	1H	1682	G	C8-N9-C1'	-5.88	119.35	127.00
1	1G	1396	A	O5'-P-OP1	-5.88	100.40	105.70
26	14	866	A	C8-N9-C1'	-5.88	117.11	127.70
26	14	939	G	C4-C5-C6	5.88	122.33	118.80
26	14	1284	A	C5-N7-C8	-5.88	100.96	103.90
26	1H	1215	G	OP1-P-O3'	5.88	118.14	105.20
26	1H	507	A	C5-N7-C8	-5.88	100.96	103.90
26	1H	536	A	N9-C4-C5	5.88	108.15	105.80
26	1H	583	G	C6-N1-C2	-5.88	121.57	125.10
26	1H	630	G	C5-C6-O6	-5.88	125.07	128.60
26	1H	2413	G	C5-C6-O6	-5.88	125.07	128.60
1	13	190	G	N3-C4-C5	-5.88	125.66	128.60
26	1H	103	A	C5-C6-N6	-5.88	119.00	123.70
26	1H	663	G	OP1-P-OP2	5.88	128.42	119.60
26	1H	2473	U	N1-C2-O2	5.88	126.92	122.80
39	A8	110	LEU	CA-CB-CG	5.88	128.82	115.30
26	1H	435	C	C6-N1-C2	-5.88	117.95	120.30
26	1H	1188	U	O5'-P-OP2	-5.88	100.41	105.70
26	1H	1769	G	O5'-P-OP2	-5.88	100.41	105.70
26	1H	2497	A	N3-C4-C5	-5.88	122.69	126.80
26	14	122	G	C5-C6-O6	-5.88	125.07	128.60
26	14	2439	A	C6-C5-N7	-5.88	128.19	132.30
26	1H	40	C	N1-C2-O2	-5.88	115.37	118.90
26	1H	54	G	C6-C5-N7	-5.88	126.87	130.40
26	1H	1291	C	N3-C4-N4	-5.88	113.89	118.00
26	1H	1771	C	N1-C2-O2	-5.88	115.37	118.90
26	1H	2409	G	C6-C5-N7	-5.88	126.87	130.40
26	14	2426	A	C5-C6-N6	-5.88	119.00	123.70
1	13	669	U	N1-C2-N3	5.88	118.42	114.90
26	1H	659	C	C2-N3-C4	-5.87	116.96	119.90
26	1H	996	A	N7-C8-N9	-5.87	110.86	113.80
26	1H	1293	C	OP1-P-O3'	5.87	118.12	105.20
26	1H	1780	A	O5'-P-OP2	5.87	117.75	110.70
26	1H	2067	G	N9-C4-C5	5.87	107.75	105.40
26	1H	2451	A	C4-C5-C6	-5.87	114.06	117.00
26	14	1274	A	C6-C5-N7	-5.87	128.19	132.30
26	14	1976	U	OP1-P-OP2	5.87	128.41	119.60
26	1H	1785	A	OP2-P-O3'	5.87	118.12	105.20
26	1H	2584	U	C4-C5-C6	5.87	123.22	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	774	A	C6-N1-C2	5.87	122.12	118.60
26	14	2013	A	C2-N3-C4	-5.87	107.66	110.60
26	14	2595	G	C4-C5-N7	5.87	113.15	110.80
26	14	2712(A)	A	N9-C4-C5	-5.87	103.45	105.80
1	1G	427	U	N1-C2-O2	5.87	126.91	122.80
23	2L	21	U	C4-C5-C6	-5.87	116.18	119.70
26	1H	786	C	OP2-P-O3'	5.87	118.11	105.20
26	1H	1265	A	C8-N9-C4	-5.87	103.45	105.80
26	1H	2749	A	C8-N9-C4	5.87	108.15	105.80
26	14	1934	C	N1-C2-O2	5.87	122.42	118.90
26	1H	215	G	C4-C5-N7	5.87	113.15	110.80
26	1H	1379	A	C5-C6-N6	-5.87	119.01	123.70
26	1H	1692	U	C5-C4-O4	-5.87	122.38	125.90
26	14	866	A	C4-N9-C1'	5.87	136.86	126.30
26	14	2542	A	O5'-P-OP2	-5.87	100.42	105.70
26	1H	1374	G	O5'-P-OP2	5.87	117.74	110.70
26	1H	1781	C	O5'-P-OP1	-5.87	100.42	105.70
26	1H	1848	A	N1-C6-N6	5.87	122.12	118.60
26	1H	2445	G	N9-C4-C5	5.87	107.75	105.40
1	1G	1234	C	N3-C2-O2	-5.87	117.79	121.90
26	1H	1942	C	N3-C4-C5	5.86	124.25	121.90
1	13	507	C	C4-C5-C6	5.86	120.33	117.40
26	1H	586	A	C5-C6-N6	5.86	128.39	123.70
26	1H	2449	U	OP2-P-O3'	5.86	118.10	105.20
25	4K	20	U	N3-C4-O4	5.86	123.50	119.40
26	1H	46	C	C6-N1-C2	-5.86	117.96	120.30
26	1H	48	G	OP2-P-O3'	5.86	118.09	105.20
26	1H	1355	G	N1-C2-N2	-5.86	110.92	116.20
26	1H	1381	G	C4-C5-N7	-5.86	108.46	110.80
26	14	1021	A	N7-C8-N9	5.86	116.73	113.80
26	14	1782	C	N1-C2-O2	-5.86	115.39	118.90
26	1H	1661	G	C8-N9-C4	5.86	108.74	106.40
26	1H	1998	G	C8-N9-C4	5.86	108.74	106.40
1	13	191	G	N1-C6-O6	5.86	123.41	119.90
26	1H	517	C	C6-N1-C2	-5.86	117.96	120.30
26	1H	2458	G	N3-C4-N9	5.86	129.51	126.00
26	14	1246	A	C8-N9-C4	5.86	108.14	105.80
26	14	1813	G	C4-C5-N7	-5.86	108.46	110.80
26	14	2260	C	C6-N1-C2	5.86	122.64	120.30
26	1H	2714	G	C5-C6-O6	-5.85	125.09	128.60
26	14	213	A	C8-N9-C4	5.85	108.14	105.80
26	14	2490	G	C5-C6-O6	5.85	132.11	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	452	G	N1-C6-O6	-5.85	116.39	119.90
26	1H	657	U	O5'-P-OP2	-5.85	100.43	105.70
26	1H	1671	U	C5-C4-O4	-5.85	122.39	125.90
26	1H	1687	G	C5-C6-N1	-5.85	108.57	111.50
41	C8	12	ARG	NE-CZ-NH2	-5.85	117.37	120.30
55	Q8	52	LYS	C-N-CD	-5.85	107.72	120.60
26	14	658	C	O5'-P-OP2	-5.85	100.43	105.70
26	14	743	G	C8-N9-C4	5.85	108.74	106.40
26	14	1614	A	O4'-C1'-N9	5.85	112.88	108.20
26	14	1728	G	N3-C4-N9	5.85	129.51	126.00
1	13	760	G	C5-C6-O6	-5.85	125.09	128.60
26	1H	1045	A	O4'-C1'-N9	5.85	112.88	108.20
26	1H	1403	C	C5-C4-N4	5.85	124.30	120.20
26	1H	2286	A	C6-C5-N7	-5.85	128.21	132.30
26	1H	2379	G	C6-C5-N7	-5.85	126.89	130.40
26	1H	2416	C	OP2-P-O3'	5.85	118.07	105.20
26	1H	2606	C	OP1-P-OP2	5.85	128.37	119.60
55	Q8	60	LEU	CA-CB-CG	5.85	128.75	115.30
26	14	632	A	O5'-P-OP2	5.85	117.72	110.70
26	14	1614	A	N1-C2-N3	5.85	132.22	129.30
27	1J	102	G	C6-C5-N7	5.85	133.91	130.40
26	1H	508	G	N7-C8-N9	5.85	116.02	113.10
26	1H	1037	G	C5-C6-O6	-5.85	125.09	128.60
26	1H	1259	G	OP2-P-O3'	5.85	118.06	105.20
26	1H	1561	G	OP1-P-O3'	5.85	118.06	105.20
26	1H	2438	U	C2-N3-C4	-5.85	123.49	127.00
1	1G	963	G	N3-C4-N9	5.85	129.51	126.00
1	1G	729	A	N1-C6-N6	-5.85	115.09	118.60
26	1H	1672	C	C2-N1-C1'	-5.84	112.37	118.80
26	1H	1807	G	C5-C6-N1	5.84	114.42	111.50
26	1H	2815	C	O5'-P-OP2	-5.84	100.44	105.70
26	1H	2857	G	O5'-P-OP1	-5.84	100.44	105.70
27	16	74	U	C5-C4-O4	5.84	129.41	125.90
26	14	141	A	C8-N9-C4	-5.84	103.46	105.80
26	14	1728	G	C2-N3-C4	5.84	114.82	111.90
26	1H	113	G	C2-N3-C4	-5.84	108.98	111.90
26	1H	336	C	N1-C2-O2	-5.84	115.39	118.90
26	1H	992	C	OP1-P-O3'	5.84	118.06	105.20
26	1H	1777	U	C4-C5-C6	5.84	123.21	119.70
26	1H	2236	C	N3-C2-O2	5.84	125.99	121.90
26	14	479	A	C4-C5-N7	-5.84	107.78	110.70
26	1H	2390	U	N3-C4-O4	5.84	123.49	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
30	31	176	LEU	CA-CB-CG	5.84	128.74	115.30
26	14	2415	G	N3-C2-N2	-5.84	115.81	119.90
26	14	2624	G	C5-C6-O6	-5.84	125.09	128.60
26	14	2880	C	C5-C6-N1	5.84	123.92	121.00
23	2K	9	G	C8-N9-C4	-5.84	104.06	106.40
26	1H	866	A	C8-N9-C1'	-5.84	117.19	127.70
26	1H	1471	A	C5-N7-C8	-5.84	100.98	103.90
27	16	74	U	N3-C4-O4	-5.84	115.31	119.40
1	1G	1462	G	N3-C4-N9	-5.84	122.50	126.00
26	14	31	C	O5'-P-OP1	-5.84	100.44	105.70
26	1H	1376	C	N3-C4-C5	-5.84	119.56	121.90
26	14	772	C	N1-C2-O2	-5.84	115.40	118.90
26	1H	1513	C	C5-C6-N1	5.84	123.92	121.00
26	14	828	U	N3-C2-O2	-5.84	118.11	122.20
26	14	1414	G	C4-N9-C1'	5.84	134.09	126.50
26	1H	1355	G	OP2-P-O3'	5.83	118.04	105.20
26	14	1686	C	C6-N1-C2	5.83	122.63	120.30
26	1H	296	C	C6-N1-C2	5.83	122.63	120.30
26	1H	1933	G	N9-C4-C5	5.83	107.73	105.40
26	14	2084	C	N3-C4-C5	-5.83	119.57	121.90
26	1H	330	A	N7-C8-N9	5.83	116.72	113.80
26	1H	1195	G	C6-N1-C2	-5.83	121.60	125.10
25	4L	16	A	C8-N9-C4	5.83	108.13	105.80
26	14	1437	C	C5-C6-N1	5.83	123.92	121.00
26	14	657	U	C5-C4-O4	5.83	129.40	125.90
1	13	769	G	N1-C6-O6	5.83	123.40	119.90
1	13	897	C	C2-N3-C4	-5.83	116.99	119.90
24	3K	35	A	N7-C8-N9	5.83	116.72	113.80
26	1H	695	G	N3-C2-N2	5.83	123.98	119.90
26	1H	834	C	N3-C4-N4	5.83	122.08	118.00
26	1H	2502	G	C4-C5-C6	5.83	122.30	118.80
26	1H	2871	C	N3-C4-N4	-5.83	113.92	118.00
1	1G	128	G	N9-C4-C5	5.83	107.73	105.40
26	14	2063	C	O5'-P-OP2	-5.83	100.45	105.70
26	14	2742	C	C6-N1-C2	5.83	122.63	120.30
26	1H	1544	C	N1-C2-O2	5.83	122.40	118.90
1	1G	388	G	C5-C6-O6	5.83	132.10	128.60
26	14	1827	C	OP1-P-O3'	5.83	118.02	105.20
27	1J	101	A	C4-C5-N7	5.83	113.61	110.70
1	13	1064	G	N3-C4-N9	-5.83	122.50	126.00
1	13	1381	U	O4'-C1'-N1	5.83	112.86	108.20
26	1H	2089	U	C6-N1-C2	5.83	124.50	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	769	G	C6-C5-N7	-5.82	126.91	130.40
26	1H	144	C	N3-C4-C5	5.82	124.23	121.90
26	1H	321	G	C4-C5-N7	5.82	113.13	110.80
26	1H	323	G	N1-C6-O6	-5.82	116.41	119.90
26	1H	398	G	N1-C2-N3	5.82	127.39	123.90
26	1H	2684	U	OP1-P-OP2	5.82	128.34	119.60
26	14	806	C	N3-C4-N4	5.82	122.08	118.00
26	14	1651	G	OP1-P-O3'	5.82	118.01	105.20
1	13	1096	C	C6-N1-C2	-5.82	117.97	120.30
26	1H	2010	G	O5'-P-OP1	-5.82	100.46	105.70
26	1H	1415	U	C5-C4-O4	5.82	129.39	125.90
26	14	1790	C	C5-C4-N4	-5.82	116.13	120.20
26	1H	656	G	OP2-P-O3'	5.82	118.00	105.20
26	1H	729	G	N7-C8-N9	5.82	116.01	113.10
1	1G	337	C	N3-C4-N4	5.82	122.07	118.00
26	14	2247	A	C8-N9-C4	-5.82	103.47	105.80
26	14	2431	U	N1-C2-O2	-5.82	118.73	122.80
26	1H	416	C	N3-C4-C5	5.82	124.23	121.90
26	1H	2078	C	N3-C4-C5	5.82	124.23	121.90
26	1H	2300	G	N7-C8-N9	5.82	116.01	113.10
26	1H	2718	G	N1-C6-O6	5.82	123.39	119.90
26	14	1022	G	C4-C5-N7	-5.82	108.47	110.80
26	14	1813	G	C6-C5-N7	5.82	133.89	130.40
1	13	860	A	N1-C6-N6	5.81	122.09	118.60
26	1H	815	C	C2-N3-C4	-5.81	116.99	119.90
26	1H	2083	G	N1-C6-O6	5.81	123.39	119.90
1	1G	1498	U	P-O3'-C3'	5.81	126.68	119.70
1	13	191(F)	U	C5-C6-N1	5.81	125.61	122.70
26	1H	613	U	C2-N1-C1'	5.81	124.67	117.70
26	1H	2286	A	C4-N9-C1'	5.81	136.76	126.30
26	1H	2713	A	C8-N9-C4	-5.81	103.47	105.80
1	1G	334	C	N3-C4-C5	5.81	124.22	121.90
26	14	807	U	C5-C4-O4	-5.81	122.41	125.90
1	13	692	U	C4-C5-C6	5.81	123.19	119.70
26	1H	40	C	C2-N3-C4	-5.81	117.00	119.90
26	1H	2387	U	O5'-P-OP2	5.81	117.67	110.70
26	1H	849	A	N9-C4-C5	-5.81	103.48	105.80
26	1H	1249	U	C6-N1-C2	5.81	124.48	121.00
26	1H	2368	C	C6-N1-C2	-5.81	117.98	120.30
27	16	78	A	O5'-P-OP1	5.81	117.67	110.70
26	14	1791	A	OP1-P-OP2	-5.81	110.88	119.60
26	14	2596	U	OP1-P-OP2	5.81	128.31	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	2K	1	C	C6-N1-C2	-5.81	117.98	120.30
23	2K	68	C	C5-C6-N1	5.81	123.90	121.00
26	1H	741	G	C6-N1-C2	-5.81	121.62	125.10
26	1H	1783	A	O5'-P-OP2	-5.81	100.47	105.70
26	1H	1938	A	O4'-C1'-N9	5.81	112.85	108.20
26	1H	2060	A	C6-C5-N7	5.81	136.37	132.30
26	14	113	G	C8-N9-C4	5.81	108.72	106.40
26	1H	825	C	C5-C4-N4	-5.81	116.14	120.20
26	1H	2435	A	C8-N9-C4	-5.81	103.48	105.80
1	13	892	A	N9-C4-C5	-5.80	103.48	105.80
1	13	1103	C	O5'-P-OP2	-5.80	100.48	105.70
26	1H	29	U	OP1-P-O3'	5.80	117.97	105.20
26	1H	2238	G	OP1-P-OP2	5.80	128.31	119.60
26	1H	2466	C	N3-C2-O2	5.80	125.96	121.90
1	1G	197	A	C4-C5-C6	5.80	119.90	117.00
26	14	575	A	O4'-C1'-N9	5.80	112.84	108.20
26	14	740	U	C2-N3-C4	5.80	130.48	127.00
26	14	2362	G	O5'-P-OP2	5.80	117.66	110.70
1	13	739	C	N1-C2-O2	-5.80	115.42	118.90
26	1H	252	G	C6-N1-C2	-5.80	121.62	125.10
26	1H	806	C	N1-C2-O2	5.80	122.38	118.90
26	1H	1550	C	N1-C2-O2	-5.80	115.42	118.90
26	1H	1807	G	O5'-P-OP2	-5.80	100.48	105.70
26	1H	2555	U	N1-C2-O2	-5.80	118.74	122.80
26	1H	2586	C	N3-C2-O2	5.80	125.96	121.90
26	1H	2777	G	O4'-C1'-N9	-5.80	103.56	108.20
26	14	2020	A	C8-N9-C4	-5.80	103.48	105.80
26	14	2573	C	C2-N1-C1'	5.80	125.18	118.80
1	13	318	G	N1-C6-O6	5.80	123.38	119.90
23	2K	58	A	N1-C2-N3	5.80	132.20	129.30
26	1H	726	G	C8-N9-C4	5.80	108.72	106.40
26	1H	1279	G	O5'-P-OP1	5.80	117.66	110.70
1	1G	1506	U	C5-C4-O4	-5.80	122.42	125.90
26	14	1252	G	O4'-C1'-N9	-5.80	103.56	108.20
26	14	1264	G	C5-C6-O6	5.80	132.08	128.60
26	14	1332	G	C4-N9-C1'	-5.80	118.96	126.50
26	14	2578	G	C5-C6-O6	5.80	132.08	128.60
26	1H	2228	G	C6-C5-N7	-5.80	126.92	130.40
26	14	391	G	C8-N9-C1'	-5.80	119.46	127.00
26	1H	138	G	N9-C1'-C2'	5.80	121.53	114.00
26	1H	602	G	O4'-C1'-N9	5.80	112.84	108.20
26	1H	2708	G	C5-C6-O6	-5.80	125.12	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1543	A	C2-N3-C4	-5.80	107.70	110.60
22	1K	48	C	C2-N1-C1'	5.79	125.17	118.80
26	1H	920	G	C8-N9-C4	5.79	108.72	106.40
26	1H	1602	U	C5-C4-O4	5.79	129.38	125.90
26	14	1241	A	C6-N1-C2	5.79	122.08	118.60
26	1H	667	U	C5-C4-O4	-5.79	122.43	125.90
26	1H	1187	G	C6-C5-N7	-5.79	126.92	130.40
26	1H	1905	C	OP1-P-OP2	-5.79	110.91	119.60
26	1H	2331	G	C5-C6-O6	-5.79	125.12	128.60
26	1H	2413	G	N1-C6-O6	5.79	123.38	119.90
1	1G	353	A	C5-N7-C8	-5.79	101.00	103.90
26	14	1646	C	OP1-P-O3'	5.79	117.94	105.20
1	1G	567	G	N1-C6-O6	-5.79	116.43	119.90
26	14	657	U	C5-C6-N1	-5.79	119.81	122.70
26	1H	2073	C	C5-C4-N4	-5.79	116.15	120.20
26	1H	2552	U	C2-N3-C4	-5.79	123.53	127.00
26	14	661	C	N3-C4-C5	-5.79	119.58	121.90
26	1H	1192	G	N9-C4-C5	-5.79	103.08	105.40
26	1H	1619	G	C6-N1-C2	-5.79	121.63	125.10
26	1H	2387	U	OP2-P-O3'	5.79	117.93	105.20
23	2K	40	C	C5-C4-N4	-5.79	116.15	120.20
26	1H	187	G	C8-N9-C1'	-5.79	119.48	127.00
26	1H	859	G	C8-N9-C4	5.79	108.71	106.40
26	1H	1632	A	C5-N7-C8	-5.79	101.01	103.90
26	1H	2490	G	O4'-C1'-N9	5.79	112.83	108.20
26	14	1366	A	N1-C2-N3	-5.79	126.41	129.30
26	14	2473	U	N3-C2-O2	-5.79	118.15	122.20
26	1H	1897	G	C2-N3-C4	-5.78	109.01	111.90
26	1H	2245	U	C5-C6-N1	5.78	125.59	122.70
26	14	676	A	C6-C5-N7	-5.78	128.25	132.30
26	14	1210	A	C2-N3-C4	-5.78	107.71	110.60
26	14	1681	G	N7-C8-N9	5.78	115.99	113.10
26	14	1948	G	C5-C6-N1	5.78	114.39	111.50
26	14	2274	A	OP2-P-O3'	5.78	117.92	105.20
26	14	2597	G	C8-N9-C1'	-5.78	119.48	127.00
26	1H	265	A	N3-C4-C5	5.78	130.85	126.80
1	1G	729	A	N9-C4-C5	5.78	108.11	105.80
26	14	1202	C	C6-N1-C2	5.78	122.61	120.30
26	14	1307	A	N1-C2-N3	5.78	132.19	129.30
1	13	1201	A	O5'-P-OP2	-5.78	100.50	105.70
26	1H	256	A	N9-C4-C5	-5.78	103.49	105.80
26	1H	1839	G	C8-N9-C1'	-5.78	119.48	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2573	C	N1-C2-N3	5.78	123.25	119.20
26	14	121	G	C5-C6-N1	5.78	114.39	111.50
27	1J	101	A	N7-C8-N9	5.78	116.69	113.80
26	1H	1224	G	C8-N9-C4	5.78	108.71	106.40
26	1H	2624	G	C8-N9-C4	5.78	108.71	106.40
26	1H	1295	C	N1-C2-O2	-5.78	115.43	118.90
26	1H	1670	C	C5-C6-N1	-5.78	118.11	121.00
26	1H	1957	C	OP1-P-OP2	-5.78	110.93	119.60
26	1H	1971	A	C5-C6-N1	5.78	120.59	117.70
1	1G	258	G	N1-C6-O6	5.78	123.37	119.90
26	14	733	G	O5'-P-OP2	-5.78	100.50	105.70
26	1H	633	A	C2-N3-C4	-5.78	107.71	110.60
26	1H	792	G	C4-N9-C1'	5.78	134.01	126.50
26	1H	962	G	OP1-P-OP2	-5.78	110.94	119.60
1	1G	1200	C	N1-C2-O2	5.78	122.36	118.90
29	29	88	GLY	N-CA-C	5.78	127.54	113.10
26	14	1702	G	N1-C6-O6	5.77	123.36	119.90
24	3K	71	G	N7-C8-N9	-5.77	110.21	113.10
26	1H	739	G	N1-C6-O6	5.77	123.36	119.90
26	1H	1382	G	N1-C6-O6	5.77	123.36	119.90
26	1H	1784	A	O5'-P-OP1	5.77	117.63	110.70
26	14	1955	U	N3-C2-O2	-5.77	118.16	122.20
26	1H	2258	C	C5-C4-N4	-5.77	116.16	120.20
1	13	770	C	OP1-P-OP2	-5.77	110.95	119.60
1	13	821	G	C4-N9-C1'	5.77	134.00	126.50
26	1H	973	A	C8-N9-C4	5.77	108.11	105.80
26	1H	1359	A	N1-C2-N3	5.77	132.19	129.30
1	1G	579	G	C4-N9-C1'	5.77	134.00	126.50
1	1G	992	U	P-O3'-C3'	5.77	126.62	119.70
1	1G	1281	U	C5-C6-N1	5.77	125.58	122.70
26	14	1903	G	OP2-P-O3'	5.77	117.89	105.20
26	14	2447	G	N3-C2-N2	-5.77	115.86	119.90
26	14	2587	A	C5-C6-N6	-5.77	119.08	123.70
1	13	500	G	C8-N9-C4	5.77	108.71	106.40
1	13	1057	G	N1-C6-O6	-5.77	116.44	119.90
26	14	77	C	OP2-P-O3'	5.77	117.89	105.20
26	14	676	A	C8-N9-C4	-5.77	103.49	105.80
26	14	2443	C	C5-C4-N4	-5.77	116.16	120.20
26	1H	115	C	C4-C5-C6	5.77	120.28	117.40
26	1H	599	G	N3-C4-C5	-5.77	125.72	128.60
26	14	2510	C	OP1-P-OP2	5.77	128.25	119.60
26	1H	2324	C	C5-C4-N4	-5.76	116.17	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2327	A	C2-N3-C4	5.76	113.48	110.60
26	1H	2336	A	C2-N3-C4	5.76	113.48	110.60
30	31	32	LEU	CA-CB-CG	5.76	128.56	115.30
1	1G	1524	C	C4-C5-C6	5.76	120.28	117.40
26	14	214	G	C8-N9-C4	-5.76	104.09	106.40
26	14	922	U	C5-C6-N1	5.76	125.58	122.70
26	14	1605	C	OP1-P-OP2	5.76	128.25	119.60
26	14	1698	A	N3-C4-C5	5.76	130.84	126.80
26	14	1771	C	N3-C4-C5	5.76	124.21	121.90
26	1H	2763	G	C4-C5-N7	5.76	113.11	110.80
26	14	671	C	OP2-P-O3'	5.76	117.88	105.20
26	1H	119	A	N1-C2-N3	5.76	132.18	129.30
26	1H	463	G	OP1-P-O3'	5.76	117.88	105.20
26	1H	2401	U	C2-N1-C1'	5.76	124.61	117.70
26	14	2355	C	C6-N1-C2	5.76	122.61	120.30
1	13	1433	A	O5'-P-OP1	-5.76	100.52	105.70
23	2K	68	C	C6-N1-C2	-5.76	118.00	120.30
26	1H	1303	G	OP1-P-OP2	5.76	128.24	119.60
26	1H	2346	A	C8-N9-C1'	-5.76	117.33	127.70
1	1G	1145	C	N1-C2-O2	5.76	122.36	118.90
26	14	205	G	N7-C8-N9	-5.76	110.22	113.10
26	14	774	A	C8-N9-C1'	5.76	138.07	127.70
26	14	786	C	C2-N1-C1'	-5.76	112.47	118.80
26	14	948	G	N3-C2-N2	-5.76	115.87	119.90
26	14	1281	G	C5-C6-O6	-5.76	125.14	128.60
26	1H	115	C	C5-C6-N1	-5.76	118.12	121.00
26	1H	1969	A	C8-N9-C4	-5.76	103.50	105.80
26	1H	2297	C	C5-C6-N1	-5.76	118.12	121.00
1	1G	579	G	C6-C5-N7	-5.76	126.94	130.40
26	14	1663	C	N3-C2-O2	5.76	125.93	121.90
26	14	2463	C	N1-C2-O2	-5.76	115.44	118.90
26	1H	670	A	O4'-C1'-N9	-5.76	103.59	108.20
26	1H	2496	C	N1-C2-O2	-5.76	115.45	118.90
26	1H	2869	G	C8-N9-C4	-5.76	104.10	106.40
26	14	740	U	N3-C4-C5	-5.76	111.15	114.60
26	14	2375	G	N9-C1'-C2'	-5.76	105.67	112.00
26	14	262	A	O5'-P-OP2	-5.75	100.52	105.70
26	14	2386	C	C5-C4-N4	-5.75	116.17	120.20
1	13	1327	C	N3-C4-C5	5.75	124.20	121.90
26	1H	208	C	C6-N1-C2	5.75	122.60	120.30
26	1H	2269	A	O5'-P-OP2	-5.75	100.52	105.70
26	1H	609	A	N9-C4-C5	-5.75	103.50	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2295	C	N3-C4-C5	-5.75	119.60	121.90
55	Q8	60	LEU	CB-CG-CD2	5.75	120.78	111.00
26	14	1782	C	P-O3'-C3'	5.75	126.60	119.70
26	14	1315	C	N3-C4-N4	-5.75	113.97	118.00
26	14	2352	A	C2-N3-C4	-5.75	107.72	110.60
26	14	2489	G	OP2-P-O3'	5.75	117.85	105.20
1	13	522	C	C4-C5-C6	5.75	120.27	117.40
26	1H	1287	A	C8-N9-C4	-5.75	103.50	105.80
26	1H	1312	U	O5'-P-OP2	5.75	117.60	110.70
26	1H	2246	G	OP1-P-O3'	5.75	117.84	105.20
26	1H	2782	G	C8-N9-C1'	-5.75	119.53	127.00
26	1H	1340	U	C4-C5-C6	5.75	123.15	119.70
26	1H	1811	G	C5-C6-O6	5.75	132.05	128.60
26	1H	2672	G	C4-C5-C6	5.75	122.25	118.80
26	14	150	C	C5-C4-N4	5.75	124.22	120.20
26	14	441	U	OP1-P-OP2	5.75	128.22	119.60
26	14	1899	G	C5-C6-O6	5.75	132.05	128.60
26	14	2363	C	C2-N3-C4	-5.75	117.03	119.90
1	1G	1488	G	N9-C4-C5	-5.75	103.10	105.40
26	1H	587	C	C4-C5-C6	5.74	120.27	117.40
26	1H	864	G	C2-N3-C4	5.74	114.77	111.90
26	1H	1568	G	OP1-P-OP2	-5.74	110.98	119.60
1	1G	811	C	N3-C2-O2	5.74	125.92	121.90
26	14	809	G	OP2-P-O3'	-5.74	92.56	105.20
26	14	2547	U	O5'-P-OP2	-5.74	100.53	105.70
1	13	418	C	C2-N1-C1'	5.74	125.12	118.80
1	13	1065	U	P-O3'-C3'	5.74	126.59	119.70
26	1H	1693	U	OP1-P-OP2	5.74	128.21	119.60
26	1H	2731	G	N9-C4-C5	-5.74	103.10	105.40
26	14	2325	G	OP1-P-OP2	5.74	128.21	119.60
26	1H	686	G	N1-C2-N2	-5.74	111.03	116.20
26	1H	751	A	N1-C6-N6	-5.74	115.16	118.60
26	1H	1939	U	N3-C4-C5	5.74	118.04	114.60
26	1H	2318	G	C5-C6-N1	-5.74	108.63	111.50
26	14	47	C	C6-N1-C2	5.74	122.60	120.30
26	14	465	G	OP1-P-OP2	-5.74	110.99	119.60
26	14	1616	A	C8-N9-C4	-5.74	103.50	105.80
1	1G	1119	C	C6-N1-C2	-5.74	118.00	120.30
26	14	620	G	C8-N9-C4	-5.74	104.10	106.40
26	14	1343	G	N3-C4-N9	5.74	129.44	126.00
26	14	1396	U	N1-C2-O2	5.74	126.82	122.80
49	G5	42	GLY	N-CA-C	-5.74	98.76	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	974	A	C5-N7-C8	-5.74	101.03	103.90
26	14	948	G	N1-C6-O6	5.74	123.34	119.90
1	13	770	C	N3-C4-N4	5.74	122.02	118.00
26	1H	1321	A	C8-N9-C4	5.74	108.09	105.80
26	1H	1761	C	C5-C4-N4	-5.74	116.19	120.20
26	1H	1957	C	C4-C5-C6	5.74	120.27	117.40
26	1H	2257	U	N3-C4-C5	-5.74	111.16	114.60
26	14	391	G	C6-C5-N7	-5.74	126.96	130.40
26	14	1623	G	N3-C4-N9	-5.74	122.56	126.00
26	14	2517	C	O4'-C1'-N1	5.74	112.79	108.20
33	69	77	LEU	CA-CB-CG	5.74	128.49	115.30
26	1H	195	A	N1-C6-N6	5.73	122.04	118.60
26	1H	853	G	O5'-P-OP2	-5.73	100.54	105.70
26	1H	1241	A	P-O3'-C3'	5.73	126.58	119.70
1	13	313	A	O5'-P-OP2	-5.73	100.54	105.70
26	1H	1267	U	C5-C6-N1	-5.73	119.83	122.70
26	1H	1698	A	O5'-P-OP2	-5.73	100.54	105.70
26	1H	2569	G	O5'-P-OP2	-5.73	100.54	105.70
26	1H	695	G	OP1-P-OP2	-5.73	111.00	119.60
26	1H	2492	U	N3-C2-O2	-5.73	118.19	122.20
26	14	2513	G	N1-C6-O6	5.73	123.34	119.90
1	13	816	A	C8-N9-C4	-5.73	103.51	105.80
26	1H	54	G	O5'-P-OP1	-5.73	100.54	105.70
26	1H	180	G	C5-C6-O6	-5.73	125.16	128.60
1	13	692	U	N3-C4-C5	-5.73	111.16	114.60
26	1H	191	A	N1-C6-N6	5.73	122.04	118.60
26	1H	436	C	N1-C2-N3	-5.73	115.19	119.20
26	1H	847	U	N3-C4-O4	-5.73	115.39	119.40
26	1H	1326	U	C5-C4-O4	5.73	129.34	125.90
26	1H	1768	U	OP2-P-O3'	5.73	117.80	105.20
26	1H	2762	G	C5-N7-C8	-5.73	101.44	104.30
26	14	179	G	OP1-P-OP2	5.73	128.19	119.60
26	1H	686	G	C2-N3-C4	-5.73	109.04	111.90
26	1H	932	G	N1-C6-O6	-5.73	116.47	119.90
26	1H	1598	C	O5'-P-OP1	-5.73	100.55	105.70
26	1H	2607	G	N3-C4-N9	5.73	129.44	126.00
1	1G	345	C	C6-N1-C2	-5.73	118.01	120.30
1	1G	720	C	C5-C6-N1	5.73	123.86	121.00
26	14	416	C	N1-C2-O2	5.73	122.34	118.90
1	13	1284	C	C6-N1-C2	5.72	122.59	120.30
26	1H	404	C	P-O3'-C3'	5.72	126.57	119.70
26	1H	565	C	O5'-P-OP1	-5.72	100.55	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	664	C	C2-N3-C4	-5.72	117.04	119.90
26	1H	664	C	OP1-P-OP2	5.72	128.19	119.60
26	1H	1197	G	C8-N9-C4	5.72	108.69	106.40
1	13	127	G	OP2-P-O3'	5.72	117.79	105.20
1	13	356	A	O4'-C1'-N9	5.72	112.78	108.20
24	3K	37	A	O4'-C1'-N9	5.72	112.78	108.20
26	1H	1833	U	C6-N1-C2	-5.72	117.57	121.00
26	1H	2731	G	C4-N9-C1'	5.72	133.94	126.50
26	1H	195	A	P-O3'-C3'	5.72	126.56	119.70
26	1H	2051	A	N7-C8-N9	5.72	116.66	113.80
26	14	2689	U	N1-C2-N3	5.72	118.33	114.90
1	13	1417	G	C8-N9-C1'	-5.72	119.56	127.00
26	1H	441	U	OP1-P-O3'	5.72	117.78	105.20
26	1H	967	C	N3-C4-C5	5.72	124.19	121.90
26	1H	1632	A	C6-C5-N7	-5.72	128.30	132.30
26	1H	1971	A	OP1-P-O3'	5.72	117.78	105.20
1	1G	995	C	C5-C6-N1	5.72	123.86	121.00
26	14	2443	C	O5'-P-OP2	5.72	117.56	110.70
26	1H	1307	A	C8-N9-C4	5.72	108.09	105.80
26	14	186	G	C8-N9-C4	5.72	108.69	106.40
26	14	1797	C	N1-C2-O2	-5.72	115.47	118.90
1	13	1412	C	C2-N3-C4	-5.72	117.04	119.90
26	1H	983	A	C5-C6-N6	-5.72	119.13	123.70
26	1H	1193	G	OP1-P-OP2	-5.72	111.02	119.60
26	1H	1386	C	C6-N1-C2	-5.72	118.01	120.30
26	1H	2265	U	C6-N1-C2	-5.72	117.57	121.00
26	1H	2467	C	C6-N1-C2	5.72	122.59	120.30
1	13	1064	G	N9-C4-C5	5.71	107.69	105.40
24	3K	76	A	O4'-C1'-N9	5.71	112.77	108.20
26	1H	1368	G	N3-C2-N2	-5.71	115.90	119.90
26	1H	2078	C	O5'-P-OP2	5.71	117.56	110.70
26	14	329	G	C5-C6-N1	5.71	114.36	111.50
26	14	1780	A	N9-C4-C5	5.71	108.09	105.80
1	13	1412	C	N1-C2-O2	-5.71	115.47	118.90
26	1H	2024	G	N1-C6-O6	-5.71	116.47	119.90
26	1H	2690	C	C6-N1-C2	-5.71	118.02	120.30
43	E8	11	ARG	NE-CZ-NH2	-5.71	117.44	120.30
26	14	746	A	O5'-P-OP2	5.71	117.56	110.70
27	1J	103	U	C6-N1-C2	5.71	124.43	121.00
1	13	363	A	N1-C6-N6	-5.71	115.17	118.60
26	1H	967	C	O5'-P-OP2	-5.71	100.56	105.70
26	1H	1775	U	C5-C6-N1	-5.71	119.84	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	568	U	O5'-P-OP1	-5.71	100.56	105.70
26	14	1315	C	OP2-P-O3'	5.71	117.76	105.20
26	14	1582	C	N1-C2-O2	5.71	122.33	118.90
26	1H	805	G	C4-C5-C6	5.71	122.22	118.80
26	1H	1213	A	C4-C5-N7	5.71	113.55	110.70
26	1H	2866	U	N3-C4-C5	-5.71	111.17	114.60
23	2L	35	C	N1-C2-O2	5.71	122.33	118.90
26	14	188	G	C6-C5-N7	-5.71	126.97	130.40
26	14	510	C	N3-C4-C5	-5.71	119.62	121.90
26	1H	835	A	C8-N9-C4	-5.71	103.52	105.80
26	1H	848	G	O5'-P-OP1	5.71	117.55	110.70
26	1H	1019	U	C5-C4-O4	5.71	129.32	125.90
26	1H	2409	G	O5'-P-OP2	-5.71	100.56	105.70
26	14	204	A	C5-C6-N6	-5.71	119.13	123.70
26	14	1127	A	O5'-P-OP1	-5.71	100.56	105.70
26	14	2070	G	N1-C2-N2	-5.71	111.06	116.20
26	14	2293	C	N3-C4-N4	-5.71	114.00	118.00
1	13	865	A	C6-C5-N7	-5.71	128.31	132.30
26	1H	681	G	N7-C8-N9	-5.71	110.25	113.10
26	14	1928	A	C8-N9-C4	5.71	108.08	105.80
26	14	2235	G	C5-C6-O6	-5.71	125.18	128.60
26	14	2277	G	N9-C4-C5	5.71	107.68	105.40
26	1H	125	G	C6-C5-N7	-5.70	126.98	130.40
26	1H	194	G	N1-C6-O6	5.70	123.32	119.90
26	1H	829	A	C2-N3-C4	-5.70	107.75	110.60
26	1H	1366	A	C5-C6-N6	-5.70	119.14	123.70
26	1H	1427	A	C6-N1-C2	-5.70	115.18	118.60
26	1H	1520	U	N3-C2-O2	-5.70	118.21	122.20
26	1H	1830	C	N3-C4-C5	5.70	124.18	121.90
27	16	80	U	OP2-P-O3'	5.70	117.75	105.20
1	1G	1115	C	N3-C4-C5	-5.70	119.62	121.90
23	2L	48	U	P-O3'-C3'	5.70	126.54	119.70
26	14	2021	C	O5'-P-OP1	-5.70	100.57	105.70
26	14	2092	U	N3-C2-O2	-5.70	118.21	122.20
26	14	2615	U	C5-C6-N1	5.70	125.55	122.70
26	1H	2330	G	N1-C2-N3	5.70	127.32	123.90
26	1H	2448	A	N3-C4-N9	5.70	131.96	127.40
26	14	36	G	N7-C8-N9	5.70	115.95	113.10
26	14	93	C	C6-N1-C2	-5.70	118.02	120.30
1	13	1502	A	N9-C4-C5	-5.70	103.52	105.80
26	1H	791	C	OP2-P-O3'	5.70	117.74	105.20
26	1H	1413	G	O5'-P-OP2	5.70	117.54	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1690	A	C6-N1-C2	-5.70	115.18	118.60
26	1H	2255	G	C5-C6-N1	-5.70	108.65	111.50
26	1H	2458	G	C8-N9-C1'	-5.70	119.59	127.00
26	1H	621	A	C8-N9-C4	-5.70	103.52	105.80
26	1H	1197	G	OP1-P-OP2	5.70	128.15	119.60
26	1H	1707	G	C4-C5-N7	5.70	113.08	110.80
26	14	138	G	N3-C4-C5	-5.70	125.75	128.60
26	1H	188	G	C5-N7-C8	-5.70	101.45	104.30
26	1H	2538	C	C6-N1-C2	5.70	122.58	120.30
39	65	73	LEU	CA-CB-CG	5.70	128.40	115.30
1	13	779	C	C6-N1-C2	-5.70	118.02	120.30
26	1H	117	G	C5-C6-O6	-5.70	125.18	128.60
26	1H	426	C	O5'-P-OP1	-5.70	100.58	105.70
26	1H	2545	G	N3-C4-C5	-5.70	125.75	128.60
26	1H	2573	C	C4-C5-C6	5.70	120.25	117.40
26	14	2379	G	O4'-C1'-N9	5.70	112.76	108.20
26	14	2776	A	P-O3'-C3'	5.70	126.53	119.70
26	1H	2212	A	C4-C5-N7	5.69	113.55	110.70
26	1H	2747	G	C6-C5-N7	-5.69	126.98	130.40
26	14	1938	A	C6-N1-C2	-5.69	115.18	118.60
26	14	1949	G	O5'-P-OP2	-5.69	100.58	105.70
26	1H	972	G	O5'-P-OP1	5.69	117.53	110.70
26	1H	1670	C	C6-N1-C2	5.69	122.58	120.30
1	1G	538	G	C6-C5-N7	-5.69	126.98	130.40
1	1G	1059	C	C6-N1-C2	-5.69	118.02	120.30
26	1H	98	G	OP1-P-OP2	5.69	128.14	119.60
26	1H	624	C	O5'-P-OP1	-5.69	100.58	105.70
26	1H	1284	A	C5-N7-C8	-5.69	101.05	103.90
26	14	530	G	C8-N9-C4	5.69	108.68	106.40
26	14	1909	C	N1-C2-O2	5.69	122.31	118.90
26	14	2597	G	O5'-P-OP1	5.69	117.53	110.70
42	95	80	GLN	N-CA-C	5.69	126.36	111.00
26	1H	416	C	C6-N1-C2	5.69	122.58	120.30
26	1H	847	U	C5-C6-N1	-5.69	119.86	122.70
26	1H	2523	G	C8-N9-C4	-5.69	104.12	106.40
1	1G	962	C	N1-C2-O2	5.69	122.31	118.90
26	14	1342	A	C6-C5-N7	-5.69	128.32	132.30
26	14	1585	C	N3-C2-O2	-5.69	117.92	121.90
26	14	2000	G	N9-C4-C5	-5.69	103.12	105.40
26	14	2501	C	C4-C5-C6	-5.69	114.56	117.40
26	1H	76	C	C4-C5-C6	5.69	120.24	117.40
26	1H	2509	G	N7-C8-N9	-5.69	110.26	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	983	A	OP2-P-O3'	5.69	117.71	105.20
26	14	2506	U	N3-C4-O4	5.69	123.38	119.40
1	13	792	A	C1'-O4'-C4'	-5.68	105.35	109.90
26	1H	1639	U	N1-C2-O2	5.68	126.78	122.80
27	16	6	C	N1-C2-O2	-5.68	115.49	118.90
1	1G	361	G	C8-N9-C4	5.68	108.67	106.40
1	13	809	G	O5'-P-OP1	-5.68	100.58	105.70
26	1H	113	G	C4-C5-C6	5.68	122.21	118.80
26	1H	330	A	O5'-P-OP2	-5.68	100.59	105.70
26	1H	671	C	C5-C6-N1	-5.68	118.16	121.00
26	1H	2330	G	C6-N1-C2	-5.68	121.69	125.10
26	1H	2546	U	C2-N1-C1'	-5.68	110.88	117.70
26	14	47	C	C5-C6-N1	-5.68	118.16	121.00
26	14	268	C	C6-N1-C2	-5.68	118.03	120.30
26	14	2721	A	OP1-P-O3'	5.68	117.70	105.20
26	1H	1246	A	C4-C5-C6	5.68	119.84	117.00
26	14	1397	U	N3-C4-O4	-5.68	115.42	119.40
1	13	484	G	C4-N9-C1'	5.68	133.88	126.50
26	1H	259	G	C8-N9-C4	5.68	108.67	106.40
26	1H	957	A	C6-C5-N7	-5.68	128.32	132.30
26	14	752	A	P-O3'-C3'	5.68	126.52	119.70
26	1H	1778	U	OP2-P-O3'	5.68	117.69	105.20
1	1G	330	C	N1-C2-O2	5.68	122.31	118.90
1	13	647	C	C6-N1-C2	-5.68	118.03	120.30
26	1H	53	A	OP1-P-O3'	5.68	117.69	105.20
26	1H	788	A	C5-C6-N1	-5.68	114.86	117.70
26	1H	1374	G	N1-C6-O6	5.68	123.31	119.90
26	1H	1393	A	OP1-P-O3'	5.68	117.69	105.20
26	1H	1935	G	N9-C4-C5	5.68	107.67	105.40
26	14	945	A	N1-C2-N3	5.68	132.14	129.30
1	13	897	C	C6-N1-C2	5.67	122.57	120.30
26	1H	324	A	O5'-P-OP2	5.67	117.51	110.70
26	1H	965	C	OP1-P-OP2	5.67	128.11	119.60
26	1H	1392	A	O5'-P-OP1	-5.67	100.59	105.70
26	14	2432	A	N1-C6-N6	5.67	122.00	118.60
1	13	1496	C	C4-C5-C6	5.67	120.24	117.40
26	14	62	C	C6-N1-C2	5.67	122.57	120.30
26	14	2224	G	C8-N9-C1'	-5.67	119.63	127.00
26	1H	763	G	C6-C5-N7	-5.67	127.00	130.40
26	1H	952	G	C2-N3-C4	5.67	114.73	111.90
26	1H	2345	G	C8-N9-C4	5.67	108.67	106.40
26	14	391	G	C5-C6-O6	-5.67	125.20	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	756	C	C6-N1-C2	-5.67	118.03	120.30
26	14	1298	C	N1-C2-O2	5.67	122.30	118.90
26	14	2595	G	C5-N7-C8	-5.67	101.46	104.30
27	1J	60	C	C6-N1-C2	-5.67	118.03	120.30
26	14	791	C	P-O3'-C3'	5.67	126.50	119.70
26	1H	125	G	C4-C5-N7	5.67	113.07	110.80
26	1H	866	A	O4'-C1'-N9	-5.67	103.67	108.20
26	1H	1204	A	C4-C5-N7	5.67	113.53	110.70
26	1H	1339	G	O5'-P-OP2	5.67	117.50	110.70
1	1G	1436	U	C2-N1-C1'	5.67	124.50	117.70
26	1H	467	G	O5'-P-OP2	-5.67	100.60	105.70
26	1H	975	G	N3-C2-N2	-5.67	115.93	119.90
26	1H	2769	C	C6-N1-C2	-5.67	118.03	120.30
20	BI	13	LEU	CA-CB-CG	5.67	128.33	115.30
26	14	1792	G	C8-N9-C4	5.67	108.67	106.40
1	13	484	G	C8-N9-C4	-5.66	104.14	106.40
1	13	819	A	O5'-P-OP1	-5.66	100.60	105.70
26	1H	104	U	C5-C4-O4	-5.66	122.50	125.90
26	1H	745	G	C4-C5-C6	5.66	122.20	118.80
26	1H	832	G	C8-N9-C4	-5.66	104.13	106.40
26	1H	1244	G	C5-C6-O6	-5.66	125.20	128.60
26	1H	1382	G	C8-N9-C1'	5.66	134.36	127.00
26	1H	1558	A	C2-N3-C4	-5.66	107.77	110.60
26	14	511	U	N3-C4-C5	-5.66	111.20	114.60
26	14	2613	U	C4-C5-C6	5.66	123.10	119.70
26	1H	2059	A	C6-N1-C2	-5.66	115.20	118.60
26	1H	2325	G	OP1-P-OP2	5.66	128.09	119.60
26	1H	2506	U	O5'-P-OP2	5.66	117.49	110.70
26	14	769	G	C5-C6-O6	-5.66	125.20	128.60
26	14	1904	G	C4-C5-N7	-5.66	108.53	110.80
26	1H	649	G	N1-C6-O6	5.66	123.30	119.90
26	14	939	G	C6-C5-N7	-5.66	127.00	130.40
26	14	2351	G	C6-C5-N7	-5.66	127.00	130.40
26	14	2518	A	N1-C2-N3	5.66	132.13	129.30
26	1H	864	G	C5-C6-N1	5.66	114.33	111.50
26	1H	1226	G	OP1-P-OP2	5.66	128.09	119.60
26	1H	2070	G	N3-C2-N2	5.66	123.86	119.90
26	1H	2271	G	C4-N9-C1'	5.66	133.86	126.50
26	14	4	C	C2-N1-C1'	5.66	125.02	118.80
26	1H	2244	U	OP1-P-OP2	-5.66	111.11	119.60
26	1H	2300	G	N1-C6-O6	5.66	123.29	119.90
26	14	827	U	C6-N1-C2	5.66	124.39	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2332	U	O5'-P-OP1	5.66	117.49	110.70
1	13	295	C	O5'-P-OP2	-5.66	100.61	105.70
1	13	887	G	N1-C2-N2	-5.66	111.11	116.20
26	1H	115	C	N3-C4-N4	5.66	121.96	118.00
26	1H	783	A	N9-C1'-C2'	-5.66	105.78	112.00
26	1H	1200	C	N1-C2-O2	-5.66	115.51	118.90
26	1H	1968	G	N9-C1'-C2'	-5.66	105.78	112.00
26	1H	2721	A	C8-N9-C4	5.66	108.06	105.80
1	1G	547	A	C8-N9-C4	5.66	108.06	105.80
26	14	2685	G	C8-N9-C4	5.66	108.66	106.40
24	3K	71	G	C4-C5-N7	-5.65	108.54	110.80
26	1H	504	U	C2-N1-C1'	5.65	124.48	117.70
26	1H	2192	G	C8-N9-C1'	-5.65	119.65	127.00
27	16	14	U	OP1-P-OP2	5.65	128.08	119.60
26	1H	576	U	OP2-P-O3'	5.65	117.64	105.20
26	1H	920	G	C5-N7-C8	5.65	107.13	104.30
26	1H	1124	C	C4-C5-C6	5.65	120.23	117.40
26	1H	1626	G	N7-C8-N9	5.65	115.93	113.10
26	14	1807	G	N9-C4-C5	-5.65	103.14	105.40
1	13	827	U	C6-N1-C1'	-5.65	113.29	121.20
1	13	1436	U	OP1-P-OP2	-5.65	111.12	119.60
26	1H	184	C	C6-N1-C2	5.65	122.56	120.30
26	1H	684	G	N1-C6-O6	-5.65	116.51	119.90
26	1H	1969	A	C5-N7-C8	5.65	106.73	103.90
26	1H	1976	U	N3-C2-O2	-5.65	118.25	122.20
26	1H	2702	U	C5-C6-N1	5.65	125.53	122.70
26	14	393	C	C5-C4-N4	5.65	124.15	120.20
26	14	736	C	O5'-P-OP2	5.65	117.48	110.70
26	1H	246	C	C6-N1-C2	5.65	122.56	120.30
26	1H	912	C	C4-C5-C6	5.65	120.22	117.40
26	1H	1496	A	O4'-C1'-N9	5.65	112.72	108.20
26	14	2079	U	N1-C2-O2	5.65	126.75	122.80
26	14	2722	G	O5'-P-OP2	5.65	117.48	110.70
26	1H	2452	C	N3-C4-C5	5.65	124.16	121.90
26	14	408	G	O5'-P-OP2	-5.65	100.62	105.70
26	1H	692	C	N1-C2-O2	-5.64	115.51	118.90
26	1H	1389	G	C8-N9-C4	-5.64	104.14	106.40
26	1H	1424	G	C6-N1-C2	-5.64	121.71	125.10
26	1H	2465	C	C4-C5-C6	5.64	120.22	117.40
1	1G	221	C	C2-N1-C1'	5.64	125.01	118.80
26	14	74	A	N9-C4-C5	-5.64	103.54	105.80
26	1H	214	G	N3-C4-N9	5.64	129.38	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2501	C	C5-C4-N4	-5.64	116.25	120.20
26	14	728	G	C8-N9-C4	5.64	108.66	106.40
26	14	2049	G	N3-C2-N2	-5.64	115.95	119.90
1	13	566	G	O5'-P-OP2	-5.64	100.62	105.70
26	14	770	G	C8-N9-C4	5.64	108.66	106.40
26	14	2523	G	C5-C6-O6	-5.64	125.22	128.60
26	1H	414	C	C2-N3-C4	-5.64	117.08	119.90
26	1H	944	G	C4-N9-C1'	5.64	133.83	126.50
26	1H	994	C	N3-C4-C5	-5.64	119.64	121.90
26	1H	1195	G	N1-C2-N3	5.64	127.28	123.90
26	14	1162	G	O5'-P-OP1	-5.64	100.62	105.70
26	14	1785	A	C4-C5-C6	5.64	119.82	117.00
1	13	1407	C	N3-C4-N4	-5.64	114.05	118.00
26	1H	2700	C	N3-C4-C5	5.64	124.16	121.90
26	1H	182	A	N7-C8-N9	-5.64	110.98	113.80
26	1H	1728	G	N3-C4-N9	5.64	129.38	126.00
46	H8	24	LEU	CA-CB-CG	5.64	128.27	115.30
1	1G	831	U	C6-N1-C2	-5.64	117.62	121.00
26	14	759	G	OP1-P-OP2	-5.64	111.14	119.60
26	14	1965	C	O5'-P-OP1	-5.64	100.63	105.70
26	14	2072	G	OP1-P-OP2	-5.64	111.14	119.60
27	1J	104	A	C8-N9-C4	5.64	108.06	105.80
26	1H	138	G	C2-N3-C4	5.63	114.72	111.90
26	1H	302	C	O5'-P-OP2	-5.63	100.63	105.70
1	1G	662	G	N3-C4-N9	5.63	129.38	126.00
1	13	808	C	N3-C4-C5	-5.63	119.65	121.90
1	13	1322	C	C2-N3-C4	5.63	122.72	119.90
26	14	2238	G	N1-C2-N2	5.63	121.27	116.20
26	1H	803	U	C2-N3-C4	-5.63	123.62	127.00
26	1H	2247	A	C2-N3-C4	-5.63	107.78	110.60
1	1G	784	C	C6-N1-C2	5.63	122.55	120.30
26	14	809	G	OP1-P-O3'	5.63	117.59	105.20
1	13	881	G	C8-N9-C4	5.63	108.65	106.40
26	1H	2503	A	C5-C6-N6	-5.63	119.20	123.70
26	14	811	U	C4-C5-C6	5.63	123.08	119.70
26	1H	124	G	N3-C4-C5	5.63	131.41	128.60
26	1H	397	G	N3-C2-N2	-5.63	115.96	119.90
26	1H	508	G	C4-N9-C1'	5.63	133.82	126.50
26	1H	560	C	C5-C4-N4	-5.63	116.26	120.20
26	1H	1888	G	N3-C4-N9	5.63	129.38	126.00
26	1H	2061	G	N3-C4-C5	-5.63	125.79	128.60
26	14	265	A	N7-C8-N9	5.63	116.61	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	992	U	P-O3'-C3'	5.63	126.45	119.70
26	1H	2056	G	N9-C4-C5	-5.63	103.15	105.40
26	1H	2082	A	C8-N9-C4	5.63	108.05	105.80
26	14	1262	A	C5-C6-N6	-5.63	119.20	123.70
1	13	1374	A	C2-N3-C4	-5.62	107.79	110.60
1	1G	197	A	C6-C5-N7	-5.62	128.36	132.30
26	14	1965	C	N1-C2-O2	-5.62	115.53	118.90
26	14	2392	A	O4'-C1'-N9	5.62	112.70	108.20
1	13	1149	C	C6-N1-C2	-5.62	118.05	120.30
26	1H	746	A	N1-C6-N6	5.62	121.97	118.60
1	1G	723	U	C2-N1-C1'	5.62	124.45	117.70
26	14	784	A	C6-N1-C2	5.62	121.97	118.60
26	14	2060	A	N9-C4-C5	5.62	108.05	105.80
1	13	1526	G	C5-N7-C8	-5.62	101.49	104.30
26	1H	415	A	C5-C6-N6	-5.62	119.20	123.70
26	1H	1410	G	N3-C4-C5	5.62	131.41	128.60
1	1G	1527	C	N3-C2-O2	-5.62	117.97	121.90
1	13	1329	A	N1-C6-N6	5.62	121.97	118.60
26	1H	312	G	OP2-P-O3'	5.62	117.56	105.20
26	1H	939	G	C5-C6-O6	5.62	131.97	128.60
26	1H	1127	A	N1-C6-N6	5.62	121.97	118.60
26	1H	1294	U	O5'-P-OP1	-5.62	100.64	105.70
26	1H	2446	G	N9-C4-C5	-5.62	103.15	105.40
26	1H	1198	U	N1-C2-N3	5.62	118.27	114.90
26	1H	1291	C	OP1-P-O3'	5.62	117.56	105.20
26	1H	2354	G	OP1-P-O3'	5.62	117.56	105.20
55	Q8	7	HIS	N-CA-C	5.62	126.17	111.00
26	14	864	G	N3-C4-C5	-5.62	125.79	128.60
26	14	1984	G	OP1-P-OP2	5.62	128.03	119.60
26	14	2679	A	C8-N9-C4	5.62	108.05	105.80
26	1H	681	G	N9-C4-C5	-5.62	103.15	105.40
1	13	606	G	N3-C4-N9	5.62	129.37	126.00
1	13	1290	G	C4-N9-C1'	5.62	133.80	126.50
26	1H	113	G	C6-C5-N7	-5.62	127.03	130.40
26	1H	682	G	N7-C8-N9	-5.62	110.29	113.10
26	1H	738	G	N1-C6-O6	5.62	123.27	119.90
26	1H	1246	A	C6-N1-C2	-5.62	115.23	118.60
26	1H	2040	C	C5-C4-N4	-5.62	116.27	120.20
26	1H	2054	A	N1-C6-N6	5.62	121.97	118.60
26	14	2252	G	OP1-P-O3'	5.62	117.55	105.20
1	13	1336	C	P-O3'-C3'	5.61	126.44	119.70
26	1H	624	C	N1-C2-O2	-5.61	115.53	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1124	C	N1-C2-N3	5.61	123.13	119.20
26	14	1768	U	C5-C4-O4	5.61	129.27	125.90
26	14	1834	U	N3-C4-O4	5.61	123.33	119.40
1	13	538	G	N3-C4-N9	5.61	129.37	126.00
26	1H	816	C	C5-C4-N4	-5.61	116.27	120.20
26	1H	967	C	C2-N3-C4	-5.61	117.09	119.90
26	14	1965	C	C2-N1-C1'	-5.61	112.63	118.80
26	14	822	U	O5'-P-OP2	5.61	117.43	110.70
26	14	1348	G	C4-C5-N7	5.61	113.05	110.80
27	1J	49	C	C6-N1-C2	-5.61	118.06	120.30
26	1H	202	U	O5'-P-OP2	5.61	117.43	110.70
26	1H	736	C	C5-C6-N1	-5.61	118.19	121.00
26	1H	2351	G	OP1-P-OP2	5.61	128.01	119.60
26	1H	2549	G	N1-C6-O6	5.61	123.27	119.90
26	14	1899	G	C8-N9-C4	-5.61	104.16	106.40
26	1H	184	C	O5'-P-OP1	-5.61	100.65	105.70
26	1H	847	U	OP1-P-OP2	5.61	128.01	119.60
26	1H	1190	G	OP1-P-OP2	5.61	128.01	119.60
26	1H	1649	G	N1-C2-N2	-5.61	111.15	116.20
26	1H	1900	A	OP1-P-OP2	-5.61	111.19	119.60
26	14	2354	G	N1-C6-O6	5.61	123.27	119.90
26	1H	779	U	N3-C4-C5	5.61	117.96	114.60
26	1H	967	C	O5'-P-OP1	5.61	117.43	110.70
26	1H	1306	C	N3-C4-C5	5.61	124.14	121.90
26	1H	2547	U	N3-C2-O2	5.61	126.12	122.20
26	14	801	G	N9-C4-C5	5.61	107.64	105.40
26	14	2509	G	OP1-P-OP2	5.61	128.01	119.60
26	14	2863	C	C6-N1-C2	5.61	122.54	120.30
26	1H	1335	U	N1-C2-O2	5.60	126.72	122.80
26	14	988	A	O5'-P-OP1	-5.60	100.66	105.70
26	1H	622	G	N3-C4-N9	5.60	129.36	126.00
26	1H	1201	C	N3-C4-N4	5.60	121.92	118.00
26	1H	1899	G	C5-N7-C8	-5.60	101.50	104.30
26	14	2551	C	C5-C6-N1	-5.60	118.20	121.00
1	13	276	G	N3-C2-N2	-5.60	115.98	119.90
26	1H	798	G	C4-C5-N7	5.60	113.04	110.80
26	14	1262	A	N9-C4-C5	-5.60	103.56	105.80
26	14	2291	U	C5-C4-O4	5.60	129.26	125.90
1	13	353	A	N7-C8-N9	5.60	116.60	113.80
26	1H	2249	U	O5'-P-OP1	-5.60	100.66	105.70
26	1H	2351	G	C8-N9-C1'	-5.60	119.72	127.00
26	1H	2604	U	O5'-P-OP1	-5.60	100.66	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
27	16	14	U	N3-C2-O2	-5.60	118.28	122.20
26	14	193	U	C4-C5-C6	5.60	123.06	119.70
26	14	1678	G	C5-C6-N1	-5.60	108.70	111.50
26	1H	831	G	C8-N9-C4	5.60	108.64	106.40
26	1H	1142(A)	A	N3-C4-C5	5.60	130.72	126.80
26	1H	1654	A	C6-N1-C2	-5.60	115.24	118.60
26	14	297	C	N3-C2-O2	-5.60	117.98	121.90
26	14	1209	G	OP1-P-OP2	5.60	128.00	119.60
26	14	1266	G	C5-C6-N1	5.60	114.30	111.50
26	14	2676	C	C5-C6-N1	-5.60	118.20	121.00
27	1J	56	G	C4-N9-C1'	5.60	133.78	126.50
26	1H	1401	G	N7-C8-N9	5.60	115.90	113.10
26	1H	1427	A	N1-C2-N3	5.60	132.10	129.30
26	1H	423	A	OP1-P-OP2	5.59	127.99	119.60
26	14	1409	C	OP1-P-OP2	5.59	127.99	119.60
26	1H	620	G	O5'-P-OP2	-5.59	100.67	105.70
26	14	2079	U	N3-C2-O2	-5.59	118.28	122.20
26	1H	783	A	C4-C5-C6	5.59	119.80	117.00
26	1H	2059	A	O4'-C1'-N9	5.59	112.67	108.20
26	1H	2617	C	N3-C2-O2	5.59	125.81	121.90
26	14	1925	C	C2-N1-C1'	-5.59	112.65	118.80
26	14	2513	G	C4-C5-N7	5.59	113.04	110.80
1	13	936	C	N3-C2-O2	-5.59	117.99	121.90
1	13	963	G	C6-C5-N7	-5.59	127.05	130.40
26	1H	116	C	C2-N1-C1'	-5.59	112.65	118.80
26	1H	631	A	O5'-P-OP1	-5.59	100.67	105.70
26	1H	2294	C	N3-C2-O2	-5.59	117.99	121.90
26	1H	2607	G	C5-N7-C8	-5.59	101.50	104.30
26	14	1949	G	OP1-P-OP2	5.59	127.98	119.60
26	14	2577	A	C4-C5-C6	5.59	119.80	117.00
33	69	102	SER	N-CA-C	-5.59	95.91	111.00
26	1H	616	A	OP2-P-O3'	5.59	117.49	105.20
1	1G	128	G	C4-C5-N7	-5.59	108.56	110.80
26	14	1342	A	C5-C6-N1	-5.59	114.91	117.70
1	13	1373	G	C8-N9-C4	-5.59	104.17	106.40
26	1H	124	G	C4-C5-N7	5.59	113.03	110.80
26	1H	1817	G	N9-C4-C5	-5.59	103.17	105.40
26	1H	2351	G	C6-C5-N7	-5.59	127.05	130.40
26	1H	2467	C	N3-C4-C5	5.59	124.14	121.90
1	1G	111	G	C5-C6-N1	-5.59	108.71	111.50
26	14	1285	G	C5-C6-N1	-5.59	108.71	111.50
26	14	2217	G	C4-N9-C1'	5.59	133.76	126.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2392	A	N1-C6-N6	5.58	121.95	118.60
26	1H	2438	U	C5-C4-O4	-5.58	122.55	125.90
1	1G	251	G	O4'-C1'-N9	-5.58	103.73	108.20
1	1G	924	C	OP1-P-OP2	5.58	127.98	119.60
1	1G	1232	U	O5'-P-OP2	-5.58	100.67	105.70
1	13	1432	G	C6-C5-N7	-5.58	127.05	130.40
26	1H	462	C	OP1-P-OP2	5.58	127.98	119.60
26	1H	1312	U	N1-C2-N3	5.58	118.25	114.90
26	1H	1633	G	OP2-P-O3'	5.58	117.48	105.20
26	1H	2067	G	C8-N9-C4	-5.58	104.17	106.40
43	E8	64	MET	N-CA-C	5.58	126.08	111.00
1	1G	449	C	C5-C4-N4	5.58	124.11	120.20
26	14	660	G	C5-C6-O6	5.58	131.95	128.60
26	14	750	A	OP1-P-O3'	5.58	117.49	105.20
26	1H	404	C	C6-N1-C2	5.58	122.53	120.30
26	1H	1010	A	N7-C8-N9	-5.58	111.01	113.80
26	1H	1349	A	N9-C4-C5	-5.58	103.57	105.80
26	1H	2379	G	C8-N9-C4	5.58	108.63	106.40
1	1G	970	C	O5'-P-OP1	-5.58	100.68	105.70
26	1H	1670	C	C2-N3-C4	-5.58	117.11	119.90
1	13	50	A	OP2-P-O3'	5.58	117.47	105.20
1	13	50	A	P-O3'-C3'	5.58	126.39	119.70
1	13	191	G	N7-C8-N9	5.58	115.89	113.10
26	1H	1272	A	C4-N9-C1'	-5.58	116.26	126.30
26	1H	1658	C	C5-C6-N1	5.58	123.79	121.00
26	1H	2374	C	N3-C4-N4	-5.58	114.10	118.00
27	16	89(A)	A	N1-C6-N6	5.58	121.95	118.60
26	14	632	A	OP1-P-OP2	-5.58	111.23	119.60
1	13	1513	A	C5-C6-N6	-5.58	119.24	123.70
26	1H	468	G	C6-N1-C2	-5.58	121.75	125.10
26	1H	626	U	OP1-P-O3'	5.58	117.47	105.20
26	1H	2477	C	C5-C6-N1	5.58	123.79	121.00
26	14	675	A	C4-C5-N7	5.58	113.49	110.70
26	14	740	U	C6-N1-C2	-5.58	117.65	121.00
26	14	1470	G	OP2-P-O3'	5.58	117.47	105.20
26	14	2703	C	C5-C6-N1	-5.58	118.21	121.00
1	13	36	C	C6-N1-C2	-5.58	118.07	120.30
26	1H	71	A	N3-C4-N9	-5.58	122.94	127.40
26	1H	198	C	C6-N1-C2	5.58	122.53	120.30
26	1H	572	A	OP1-P-OP2	-5.58	111.24	119.60
26	1H	2544	G	C4-C5-N7	5.58	113.03	110.80
26	1H	2549	G	C5-C6-O6	-5.58	125.25	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2718	G	C5-C6-O6	-5.58	125.25	128.60
26	14	150	C	N3-C4-N4	-5.58	114.10	118.00
26	14	597	U	N3-C4-O4	5.58	123.30	119.40
26	14	1612	C	N3-C2-O2	5.58	125.80	121.90
1	13	975	A	O4'-C1'-N9	-5.57	103.74	108.20
26	1H	813	U	OP1-P-OP2	5.57	127.96	119.60
26	14	848	G	C8-N9-C1'	-5.57	119.75	127.00
26	14	2054	A	C8-N9-C4	-5.57	103.57	105.80
26	14	2332	U	C5-C4-O4	5.57	129.24	125.90
26	14	2880	C	N3-C4-C5	-5.57	119.67	121.90
26	14	179	G	C8-N9-C4	5.57	108.63	106.40
26	1H	581	C	C6-N1-C2	-5.57	118.07	120.30
26	1H	944	G	C5-C6-O6	5.57	131.94	128.60
26	1H	954	G	O5'-P-OP2	5.57	117.39	110.70
26	1H	1196	C	N1-C2-O2	-5.57	115.56	118.90
26	1H	1222	C	C6-N1-C2	-5.57	118.07	120.30
26	14	2544	G	N9-C4-C5	-5.57	103.17	105.40
1	13	1468	A	C8-N9-C4	5.57	108.03	105.80
26	1H	424	G	N9-C4-C5	5.57	107.63	105.40
26	1H	249	C	N1-C2-O2	5.57	122.24	118.90
26	1H	972	G	C5-C6-N1	5.57	114.28	111.50
26	1H	1394	U	C2-N3-C4	5.57	130.34	127.00
26	1H	1749	A	C8-N9-C4	-5.57	103.57	105.80
1	1G	115	G	P-O3'-C3'	5.57	126.38	119.70
26	14	1308	A	C4-C5-C6	5.57	119.78	117.00
26	1H	866	A	C4-C5-N7	5.57	113.48	110.70
54	P8	9	ARG	NE-CZ-NH1	-5.57	117.52	120.30
26	14	1954	G	N9-C4-C5	5.57	107.63	105.40
55	M5	33	ASN	N-CA-C	5.57	126.03	111.00
26	1H	134	C	N3-C4-C5	5.56	124.13	121.90
1	13	856	C	C6-N1-C2	-5.56	118.08	120.30
26	1H	696	G	O5'-P-OP2	5.56	117.38	110.70
26	1H	762	U	C6-N1-C1'	-5.56	113.41	121.20
26	1H	802	A	OP1-P-O3'	-5.56	92.96	105.20
26	1H	2045	C	C6-N1-C2	5.56	122.53	120.30
27	16	47	C	N3-C4-C5	5.56	124.12	121.90
1	1G	1147	C	C6-N1-C1'	-5.56	114.13	120.80
26	14	1419	A	O4'-C1'-N9	5.56	112.65	108.20
26	14	1569	A	C8-N9-C4	-5.56	103.58	105.80
26	14	1667	G	C5-C6-N1	-5.56	108.72	111.50
26	1H	1035	U	C5-C6-N1	-5.56	119.92	122.70
26	1H	1144	G	C8-N9-C4	-5.56	104.17	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1352	U	C2-N3-C4	-5.56	123.66	127.00
26	1H	1559	G	N3-C4-C5	5.56	131.38	128.60
26	1H	2018	G	N1-C6-O6	5.56	123.24	119.90
26	1H	448	U	N3-C2-O2	-5.56	118.31	122.20
26	1H	866	A	C6-C5-N7	-5.56	128.41	132.30
26	1H	2672	G	N1-C2-N3	5.56	127.23	123.90
26	14	453	C	C5-C4-N4	-5.56	116.31	120.20
26	14	871	U	OP1-P-O3'	5.56	117.43	105.20
1	13	892	A	C8-N9-C4	5.56	108.02	105.80
26	1H	1284	A	C6-C5-N7	-5.56	128.41	132.30
26	1H	1287	A	O5'-P-OP1	5.56	117.37	110.70
26	1H	1292	U	C5-C6-N1	-5.56	119.92	122.70
26	1H	1611	C	N3-C4-C5	5.56	124.12	121.90
26	1H	2379	G	C4-C5-N7	5.56	113.02	110.80
45	G8	106	LEU	CA-CB-CG	5.56	128.08	115.30
26	14	2732	G	C6-N1-C2	-5.56	121.77	125.10
1	13	789	U	N3-C4-C5	-5.56	111.27	114.60
26	1H	301	G	OP1-P-OP2	5.56	127.93	119.60
26	1H	611	C	C5-C6-N1	-5.56	118.22	121.00
26	1H	1190	G	O5'-P-OP1	-5.56	100.70	105.70
1	1G	963	G	N3-C4-C5	-5.56	125.82	128.60
26	1H	917	A	O4'-C1'-N9	5.55	112.64	108.20
26	1H	1314	C	C6-N1-C1'	-5.55	114.13	120.80
1	1G	306	G	N3-C4-N9	-5.55	122.67	126.00
26	1H	1051	G	C6-C5-N7	-5.55	127.07	130.40
26	14	117	G	C5-C6-N1	5.55	114.28	111.50
26	14	570	G	C8-N9-C4	-5.55	104.18	106.40
26	1H	199	A	C5-C6-N6	5.55	128.14	123.70
26	1H	639	U	N1-C2-O2	5.55	126.69	122.80
26	1H	860	U	N1-C2-N3	5.55	118.23	114.90
26	1H	919	G	N1-C2-N3	5.55	127.23	123.90
26	1H	971	C	N1-C2-N3	5.55	123.09	119.20
26	1H	1272	A	N1-C6-N6	-5.55	115.27	118.60
26	14	1601	G	OP1-P-O3'	5.55	117.41	105.20
26	14	2225	A	P-O3'-C3'	5.55	126.36	119.70
1	13	595	G	C5-C6-O6	5.55	131.93	128.60
1	13	948	C	OP1-P-O3'	5.55	117.41	105.20
1	13	1279	A	N7-C8-N9	5.55	116.57	113.80
26	1H	371	A	N9-C4-C5	-5.55	103.58	105.80
26	1H	955	C	OP1-P-OP2	5.55	127.92	119.60
26	1H	975	G	N1-C2-N2	5.55	121.19	116.20
25	4L	16	A	N7-C8-N9	-5.55	111.03	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	833	U	C5-C6-N1	-5.55	119.92	122.70
26	14	974(A)	C	N3-C4-N4	-5.55	114.11	118.00
26	14	1480	G	N1-C6-O6	5.55	123.23	119.90
26	1H	659	C	N3-C4-C5	5.55	124.12	121.90
26	1H	2686	G	N3-C4-C5	-5.55	125.83	128.60
26	14	614	U	N3-C2-O2	-5.55	118.32	122.20
26	14	2084	C	N3-C2-O2	5.55	125.78	121.90
26	1H	113	G	C5-C6-N1	-5.55	108.73	111.50
26	1H	911	A	N1-C6-N6	5.55	121.93	118.60
26	1H	1487	G	N3-C4-C5	5.55	131.37	128.60
26	1H	1784	A	O4'-C1'-N9	-5.55	103.76	108.20
26	1H	1791	A	C2-N3-C4	5.55	113.37	110.60
26	1H	1982	C	N3-C4-N4	5.55	121.88	118.00
26	1H	2334	G	C5-C6-N1	5.55	114.27	111.50
26	14	409	C	C6-N1-C2	5.55	122.52	120.30
26	1H	109	G	C4-C5-N7	-5.54	108.58	110.80
26	1H	551	G	N7-C8-N9	-5.54	110.33	113.10
26	1H	1158	C	C2-N3-C4	-5.54	117.13	119.90
26	14	767	U	C5-C4-O4	5.54	129.23	125.90
1	13	1195	C	C5-C6-N1	5.54	123.77	121.00
26	1H	830	G	C2-N3-C4	-5.54	109.13	111.90
26	1H	1968	G	N3-C4-N9	5.54	129.33	126.00
26	1H	2368	C	N3-C4-C5	-5.54	119.68	121.90
26	14	1863	G	O5'-P-OP2	-5.54	100.71	105.70
1	13	122	G	N3-C2-N2	-5.54	116.02	119.90
1	13	1378	C	C5-C6-N1	5.54	123.77	121.00
26	1H	945	A	O5'-P-OP1	5.54	117.35	110.70
26	1H	2499	C	N3-C4-N4	5.54	121.88	118.00
26	14	1668	A	N9-C4-C5	-5.54	103.58	105.80
26	14	1979	C	C5-C4-N4	-5.54	116.32	120.20
26	14	2261	C	N3-C4-C5	5.54	124.12	121.90
27	1J	7	G	C6-C5-N7	-5.54	127.08	130.40
1	13	1203	C	C6-N1-C2	-5.54	118.08	120.30
26	1H	825	C	C4-C5-C6	5.54	120.17	117.40
26	1H	2237	G	N3-C2-N2	5.54	123.78	119.90
26	14	2589	A	N7-C8-N9	-5.54	111.03	113.80
27	1J	60	C	C5-C6-N1	5.54	123.77	121.00
1	13	1128	C	N1-C2-O2	5.54	122.22	118.90
26	1H	535	C	C5-C6-N1	-5.54	118.23	121.00
26	1H	1605	C	N3-C4-C5	5.54	124.12	121.90
26	14	187	G	N3-C4-N9	5.54	129.32	126.00
26	14	1277	G	C8-N9-C4	5.54	108.62	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2574	G	C6-N1-C2	-5.54	121.78	125.10
26	14	1937	A	C5-C6-N1	-5.54	114.93	117.70
26	14	1997	G	C4-N9-C1'	5.54	133.70	126.50
27	1J	102	G	N3-C4-C5	5.54	131.37	128.60
1	13	238	G	N7-C8-N9	-5.54	110.33	113.10
26	1H	1200	C	O5'-P-OP1	5.54	117.34	110.70
26	1H	1660	C	N3-C2-O2	-5.54	118.03	121.90
26	1H	2265	U	N3-C4-C5	-5.54	111.28	114.60
26	1H	2708	G	N1-C2-N2	-5.54	111.22	116.20
26	14	205	G	C5-C6-O6	-5.54	125.28	128.60
26	14	2624	G	C2-N3-C4	5.54	114.67	111.90
1	13	740	U	N3-C4-C5	-5.53	111.28	114.60
26	1H	13	A	N1-C6-N6	-5.53	115.28	118.60
26	14	410	G	O5'-P-OP2	5.53	117.34	110.70
26	14	1700	A	O5'-P-OP2	5.53	117.34	110.70
26	14	1844	C	C5-C4-N4	-5.53	116.33	120.20
1	13	506	G	N3-C2-N2	-5.53	116.03	119.90
26	1H	1805	U	OP2-P-O3'	5.53	117.37	105.20
26	1H	2391	G	C5-C6-O6	5.53	131.92	128.60
26	1H	2538	C	C5-C6-N1	-5.53	118.23	121.00
1	1G	337	C	C6-N1-C2	-5.53	118.09	120.30
1	1G	555	C	C2-N1-C1'	5.53	124.88	118.80
26	14	130	C	C5-C6-N1	-5.53	118.23	121.00
26	14	1299	G	C4-C5-N7	5.53	113.01	110.80
26	14	1725	G	C4-N9-C1'	5.53	133.69	126.50
26	14	2392	A	O5'-P-OP2	5.53	117.34	110.70
1	13	667	G	N7-C8-N9	5.53	115.86	113.10
26	1H	639	U	C6-N1-C2	-5.53	117.68	121.00
26	1H	1345	C	OP2-P-O3'	5.53	117.36	105.20
26	1H	1606	G	O5'-P-OP2	-5.53	100.72	105.70
1	1G	687	A	P-O3'-C3'	5.53	126.33	119.70
26	1H	49	A	N1-C6-N6	5.53	121.92	118.60
26	1H	321	G	N3-C4-N9	5.53	129.32	126.00
26	1H	528	A	C5-C6-N6	5.53	128.12	123.70
26	1H	970	C	N1-C2-O2	-5.53	115.58	118.90
26	1H	1621	U	N1-C2-O2	-5.53	118.93	122.80
1	13	972	C	OP2-P-O3'	5.53	117.36	105.20
26	1H	142	G	C8-N9-C1'	5.53	134.18	127.00
26	1H	271(C)	U	C6-N1-C1'	-5.53	113.46	121.20
26	1H	1361	G	N7-C8-N9	-5.53	110.34	113.10
27	16	6	C	C5-C6-N1	-5.53	118.24	121.00
26	14	223	A	C8-N9-C4	-5.53	103.59	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2374	C	C4-C5-C6	5.52	120.16	117.40
26	1H	2468	G	N7-C8-N9	5.52	115.86	113.10
26	14	2254	C	OP2-P-O3'	5.52	117.35	105.20
1	13	788	U	N1-C2-N3	5.52	118.21	114.90
26	1H	805	G	OP1-P-O3'	5.52	117.35	105.20
26	1H	1301	A	N1-C6-N6	5.52	121.91	118.60
26	1H	1916	A	N1-C2-N3	5.52	132.06	129.30
26	1H	2755	C	C5-C6-N1	5.52	123.76	121.00
26	14	1334	G	N3-C2-N2	-5.52	116.03	119.90
26	14	2477	C	C6-N1-C2	-5.52	118.09	120.30
26	1H	468	G	N1-C6-O6	5.52	123.21	119.90
26	1H	1255	U	C5-C4-O4	-5.52	122.59	125.90
26	14	2877	G	O5'-P-OP1	5.52	117.33	110.70
1	13	1369	C	O5'-P-OP2	-5.52	100.73	105.70
26	1H	821	A	O5'-P-OP2	-5.52	100.73	105.70
26	1H	1780	A	C5-C6-N6	5.52	128.12	123.70
26	1H	2300	G	C8-N9-C4	-5.52	104.19	106.40
26	1H	2310	A	C6-N1-C2	-5.52	115.29	118.60
26	1H	2540	C	N3-C4-C5	5.52	124.11	121.90
30	31	176	LEU	CB-CG-CD2	-5.52	101.62	111.00
26	14	999	U	O5'-P-OP2	5.52	117.32	110.70
26	14	1982	C	C6-N1-C2	-5.52	118.09	120.30
1	13	190	G	C8-N9-C4	-5.52	104.19	106.40
26	1H	743	G	N3-C4-C5	-5.52	125.84	128.60
26	1H	1949	G	N3-C2-N2	-5.52	116.04	119.90
26	1H	2501	C	N3-C4-C5	5.52	124.11	121.90
26	1H	2593	U	C5-C6-N1	5.52	125.46	122.70
1	1G	701	C	N1-C2-O2	5.52	122.21	118.90
26	14	1806	C	O5'-P-OP2	-5.52	100.73	105.70
26	14	2592	G	N1-C2-N2	-5.52	111.23	116.20
26	1H	20	C	O5'-P-OP2	-5.52	100.74	105.70
26	1H	33	U	OP1-P-O3'	5.52	117.33	105.20
26	1H	528	A	C4-N9-C1'	-5.52	116.37	126.30
26	1H	2509	G	N9-C4-C5	-5.52	103.19	105.40
26	14	1377	G	N3-C4-C5	-5.52	125.84	128.60
26	14	1519	G	C5-C6-N1	-5.52	108.74	111.50
26	14	2011	U	C6-N1-C2	5.52	124.31	121.00
26	14	2708	G	N1-C2-N2	-5.52	111.23	116.20
1	13	1504	G	P-O3'-C3'	5.51	126.32	119.70
26	1H	463	G	N1-C2-N2	-5.51	111.24	116.20
26	1H	864	G	N3-C4-C5	-5.51	125.84	128.60
26	1H	2736	G	N3-C4-C5	5.51	131.36	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	6	G	N1-C6-O6	5.51	123.21	119.90
26	14	210	C	C6-N1-C2	5.51	122.50	120.30
26	14	602	G	N9-C4-C5	-5.51	103.19	105.40
26	14	2314	C	N3-C2-O2	-5.51	118.04	121.90
26	14	385	C	O5'-P-OP1	-5.51	100.74	105.70
26	1H	775	G	N3-C4-N9	5.51	129.31	126.00
26	1H	805	G	C8-N9-C4	-5.51	104.19	106.40
26	1H	2465	C	C2-N3-C4	-5.51	117.14	119.90
26	1H	2586	C	C6-N1-C2	5.51	122.50	120.30
26	14	2512	C	N3-C2-O2	5.51	125.76	121.90
26	1H	695	G	N1-C2-N2	-5.51	111.24	116.20
26	1H	1605	C	O5'-P-OP1	-5.51	100.74	105.70
27	16	44	G	C8-N9-C4	5.51	108.60	106.40
1	1G	906	G	C8-N9-C4	5.51	108.60	106.40
26	14	684	G	N7-C8-N9	5.51	115.86	113.10
26	14	1653	G	C5-N7-C8	5.51	107.05	104.30
48	F5	29	GLY	N-CA-C	5.51	126.88	113.10
26	1H	2440	C	OP1-P-OP2	5.51	127.86	119.60
26	1H	2871	C	N3-C2-O2	-5.51	118.04	121.90
26	14	1333	C	C4-C5-C6	-5.51	114.65	117.40
23	2K	35	C	C2-N1-C1'	5.51	124.86	118.80
26	1H	809	G	N3-C4-C5	-5.51	125.85	128.60
26	1H	868	U	N3-C2-O2	-5.51	118.35	122.20
1	1G	1061	G	C6-C5-N7	-5.51	127.10	130.40
26	14	479	A	C5-N7-C8	5.51	106.65	103.90
26	1H	1021	A	O4'-C1'-N9	-5.50	103.80	108.20
28	11	147	LEU	CB-CG-CD2	-5.50	101.64	111.00
26	14	665	C	C6-N1-C1'	-5.50	114.19	120.80
26	14	1777	U	N1-C2-N3	5.50	118.20	114.90
26	14	2508	G	C5-C6-N1	5.50	114.25	111.50
26	1H	684	G	N3-C4-C5	-5.50	125.85	128.60
26	1H	1949	G	N3-C4-N9	-5.50	122.70	126.00
26	1H	2073	C	N1-C2-O2	-5.50	115.60	118.90
26	1H	2076	U	N1-C2-O2	-5.50	118.95	122.80
26	1H	2775	A	C5-C6-N6	-5.50	119.30	123.70
39	A8	101	LEU	CA-CB-CG	5.50	127.96	115.30
26	14	475	U	N3-C2-O2	-5.50	118.35	122.20
26	14	748	G	N9-C4-C5	5.50	107.60	105.40
26	14	1277	G	OP1-P-OP2	5.50	127.86	119.60
26	1H	25	U	C6-N1-C2	5.50	124.30	121.00
26	1H	230	U	O5'-P-OP2	-5.50	100.75	105.70
26	1H	1192	G	N7-C8-N9	-5.50	110.35	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2080	G	N1-C6-O6	-5.50	116.60	119.90
1	1G	538	G	C4-N9-C1'	5.50	133.65	126.50
26	1H	1138	G	C4-N9-C1'	5.50	133.65	126.50
26	1H	1334	G	C4-C5-N7	5.50	113.00	110.80
1	1G	932	C	N1-C2-O2	5.50	122.20	118.90
1	13	1335	C	C5-C6-N1	-5.50	118.25	121.00
26	1H	1997	G	N1-C2-N3	5.50	127.20	123.90
26	1H	2611	U	N3-C2-O2	-5.50	118.35	122.20
1	1G	1530	G	N7-C8-N9	5.50	115.85	113.10
26	14	1567	A	C8-N9-C4	-5.50	103.60	105.80
26	14	1796	U	O5'-P-OP2	5.50	117.30	110.70
27	1J	44	G	C4-N9-C1'	-5.50	119.35	126.50
26	1H	51	G	C8-N9-C1'	-5.50	119.86	127.00
26	1H	133	C	N3-C4-C5	5.50	124.10	121.90
26	1H	670	A	N9-C4-C5	-5.50	103.60	105.80
26	1H	864	G	N3-C4-N9	5.50	129.30	126.00
26	1H	1343	G	N3-C4-N9	5.50	129.30	126.00
26	1H	1424	G	N3-C4-N9	5.50	129.30	126.00
26	1H	1685	C	OP1-P-O3'	5.50	117.29	105.20
27	16	8	U	C5-C4-O4	5.50	129.20	125.90
26	14	203	C	N3-C4-C5	5.50	124.10	121.90
26	14	1932	A	O5'-P-OP1	-5.50	100.75	105.70
1	13	138	G	O5'-P-OP1	5.50	117.29	110.70
26	1H	389	G	C8-N9-C4	5.50	108.60	106.40
26	1H	1222	C	C2-N1-C1'	5.50	124.84	118.80
26	1H	2032	G	C5-N7-C8	5.50	107.05	104.30
24	3L	71	G	C4-C5-N7	-5.50	108.60	110.80
26	14	2031	A	C5-C6-N6	-5.50	119.30	123.70
26	14	2258	C	N3-C2-O2	5.50	125.75	121.90
26	1H	719	C	C6-N1-C2	-5.49	118.10	120.30
26	1H	793	A	C4-C5-C6	5.49	119.75	117.00
26	1H	2243	U	O5'-P-OP2	-5.49	100.76	105.70
1	1G	912	C	C5-C6-N1	-5.49	118.25	121.00
26	14	1343	G	N3-C4-C5	-5.49	125.85	128.60
26	14	1820	U	O5'-P-OP2	5.49	117.29	110.70
26	1H	385	C	O5'-P-OP1	-5.49	100.76	105.70
26	1H	1394	U	N3-C4-C5	-5.49	111.31	114.60
1	13	238	G	C8-N9-C4	5.49	108.60	106.40
23	2K	10	G	O5'-P-OP1	-5.49	100.76	105.70
26	1H	1935	G	C5-C6-O6	5.49	131.90	128.60
24	3L	76	A	C2-N3-C4	-5.49	107.86	110.60
26	14	1734	C	C6-N1-C2	-5.49	118.10	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2042	A	C2-N3-C4	-5.49	107.86	110.60
26	14	2248	C	O5'-P-OP1	-5.49	100.76	105.70
26	1H	1951	U	O5'-P-OP2	-5.49	100.76	105.70
26	1H	2592	G	C6-C5-N7	-5.49	127.11	130.40
23	2L	76	C	N1-C2-O2	-5.49	115.61	118.90
26	14	508	G	OP1-P-OP2	5.49	127.83	119.60
26	14	1022	G	C6-C5-N7	5.49	133.69	130.40
26	14	1203	G	O5'-P-OP1	5.49	117.29	110.70
23	2K	74	A	C4-C5-N7	5.49	113.44	110.70
26	1H	848	G	O5'-P-OP2	-5.49	100.76	105.70
26	14	211	A	C5-C6-N6	-5.49	119.31	123.70
26	14	2426	A	C4-C5-N7	5.49	113.44	110.70
26	1H	2253	G	N1-C6-O6	5.49	123.19	119.90
53	O8	42	TRP	CA-CB-CG	5.49	124.12	113.70
1	1G	972	C	O5'-P-OP2	-5.49	100.76	105.70
26	14	915	C	N1-C2-O2	5.49	122.19	118.90
26	14	2442	C	C4-C5-C6	5.49	120.14	117.40
1	13	191(F)	U	C6-N1-C2	-5.48	117.71	121.00
1	13	910	C	C2-N3-C4	-5.48	117.16	119.90
23	2K	68	C	N3-C4-N4	5.48	121.84	118.00
26	1H	2453	A	N9-C4-C5	5.48	107.99	105.80
26	14	780	G	C4-C5-C6	5.48	122.09	118.80
26	14	2263	C	O5'-P-OP2	5.48	117.28	110.70
26	14	2502	G	N3-C4-C5	-5.48	125.86	128.60
26	1H	607	U	C5-C6-N1	-5.48	119.96	122.70
26	1H	687	C	O5'-P-OP1	-5.48	100.77	105.70
26	1H	793	A	C6-N1-C2	-5.48	115.31	118.60
26	1H	1363	C	N3-C4-C5	5.48	124.09	121.90
26	1H	1432	C	N3-C4-N4	5.48	121.84	118.00
26	1H	2061	G	N3-C2-N2	5.48	123.74	119.90
1	1G	320	C	C6-N1-C2	5.48	122.49	120.30
26	14	2330	G	N7-C8-N9	-5.48	110.36	113.10
26	14	2606	C	C2-N3-C4	-5.48	117.16	119.90
26	1H	627	A	N7-C8-N9	-5.48	111.06	113.80
26	1H	1129	A	OP1-P-OP2	5.48	127.82	119.60
26	1H	2324	C	OP1-P-OP2	5.48	127.82	119.60
26	14	263	C	N1-C2-O2	5.48	122.19	118.90
1	13	900	A	C5-C6-N1	5.48	120.44	117.70
26	1H	1051	G	N1-C6-O6	5.48	123.19	119.90
26	1H	1283	G	C8-N9-C1'	-5.48	119.88	127.00
26	14	1703	G	C4-C5-N7	5.48	112.99	110.80
26	1H	54	G	N1-C6-O6	5.48	123.19	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	768	G	C6-C5-N7	-5.48	127.11	130.40
26	14	1309	G	N9-C4-C5	-5.48	103.21	105.40
26	14	2435	A	OP1-P-O3'	5.48	117.25	105.20
26	1H	120	U	N3-C4-O4	-5.47	115.57	119.40
26	1H	622	G	N1-C2-N2	-5.47	111.27	116.20
26	1H	2439	A	C6-C5-N7	-5.47	128.47	132.30
26	1H	2473	U	C6-N1-C1'	-5.47	113.54	121.20
43	E8	107	LEU	CA-CB-CG	5.47	127.89	115.30
1	1G	245	C	N1-C2-O2	-5.47	115.61	118.90
26	14	954	G	N3-C4-C5	-5.47	125.86	128.60
26	14	2231	C	C5-C4-N4	-5.47	116.37	120.20
26	14	2443	C	N3-C4-N4	5.47	121.83	118.00
26	1H	1817	G	C5-N7-C8	5.47	107.04	104.30
26	1H	2359	C	C4-C5-C6	5.47	120.14	117.40
1	1G	1061	G	C4-N9-C1'	5.47	133.62	126.50
1	1G	1502	A	C2-N3-C4	-5.47	107.86	110.60
1	13	651	C	N1-C2-O2	5.47	122.18	118.90
26	1H	1428	C	C5-C4-N4	5.47	124.03	120.20
26	1H	1517	G	OP1-P-O3'	5.47	117.24	105.20
26	1H	371	A	C4-C5-N7	5.47	113.44	110.70
26	1H	634	C	O5'-P-OP2	-5.47	100.78	105.70
26	1H	1624	G	N1-C6-O6	-5.47	116.62	119.90
26	1H	1973	G	N3-C2-N2	5.47	123.73	119.90
26	1H	2782	G	N3-C4-N9	5.47	129.28	126.00
1	1G	328	C	P-O3'-C3'	5.47	126.26	119.70
1	1G	1527	C	C4-C5-C6	5.47	120.14	117.40
26	14	1569	A	O4'-C1'-N9	5.47	112.58	108.20
26	14	2272	U	N1-C2-O2	5.47	126.63	122.80
26	14	2523	G	C8-N9-C4	5.47	108.59	106.40
26	14	2597	G	C8-N9-C4	5.47	108.59	106.40
1	13	1529	G	OP2-P-O3'	5.47	117.23	105.20
1	1G	108	G	N1-C2-N3	-5.47	120.62	123.90
26	14	140	A	C5-C6-N6	-5.47	119.33	123.70
26	14	1827	C	N3-C2-O2	-5.47	118.07	121.90
1	13	226	G	C8-N9-C1'	-5.47	119.89	127.00
1	13	512	U	O5'-P-OP2	-5.47	100.78	105.70
1	13	578	C	N3-C4-N4	5.47	121.83	118.00
26	1H	1610	A	N1-C6-N6	5.47	121.88	118.60
26	1H	2241	A	C2-N3-C4	-5.47	107.87	110.60
26	14	2387	U	C5-C6-N1	-5.47	119.97	122.70
27	1J	103	U	C2-N1-C1'	-5.47	111.14	117.70
26	1H	380	U	N3-C2-O2	-5.46	118.38	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1312	U	P-O3'-C3'	5.46	126.26	119.70
26	1H	1791	A	N1-C2-N3	-5.46	126.57	129.30
26	1H	2438	U	C6-N1-C1'	-5.46	113.55	121.20
1	1G	392	G	C8-N9-C4	5.46	108.58	106.40
26	14	1313	U	C2-N1-C1'	5.46	124.26	117.70
26	14	2420	C	C6-N1-C1'	-5.46	114.24	120.80
26	14	2443	C	O5'-P-OP1	-5.46	100.78	105.70
26	1H	2458	G	C4-N9-C1'	5.46	133.60	126.50
26	1H	2467	C	C5-C6-N1	-5.46	118.27	121.00
26	14	2444	G	O5'-P-OP2	-5.46	100.78	105.70
1	13	237	C	N1-C2-O2	-5.46	115.62	118.90
26	1H	795	C	C6-N1-C2	5.46	122.48	120.30
26	1H	1311	G	C5-C6-O6	-5.46	125.32	128.60
26	1H	1332	G	N1-C2-N3	5.46	127.18	123.90
26	1H	1438	U	N3-C4-O4	5.46	123.22	119.40
26	1H	2299	G	N1-C6-O6	5.46	123.18	119.90
26	1H	2641	G	N1-C6-O6	5.46	123.18	119.90
26	14	782	A	N9-C4-C5	5.46	107.98	105.80
26	14	2003	G	N1-C6-O6	5.46	123.18	119.90
30	39	80	ALA	C-N-CD	5.46	139.87	128.40
26	14	1265	A	C5-C6-N6	-5.46	119.33	123.70
1	13	254	G	O5'-P-OP1	-5.46	100.79	105.70
26	1H	1626	G	N9-C4-C5	5.46	107.58	105.40
1	1G	576	G	C5-C6-N1	-5.46	108.77	111.50
26	14	2286	A	C8-N9-C4	-5.46	103.62	105.80
26	14	2679	A	N7-C8-N9	-5.46	111.07	113.80
1	13	1099	G	N1-C6-O6	-5.46	116.63	119.90
1	13	1475	G	OP1-P-OP2	5.46	127.78	119.60
26	1H	737	C	C2-N1-C1'	-5.46	112.80	118.80
26	1H	818	G	C5-C6-O6	5.46	131.87	128.60
26	1H	1471	A	C6-C5-N7	-5.46	128.48	132.30
26	14	71	A	N7-C8-N9	5.46	116.53	113.80
26	14	530	G	C5-C6-N1	-5.46	108.77	111.50
26	14	784	A	OP1-P-O3'	5.46	117.20	105.20
26	14	831	G	C8-N9-C4	5.46	108.58	106.40
26	14	1623	G	C4-N9-C1'	-5.46	119.41	126.50
26	14	2287	A	C5-C6-N1	-5.46	114.97	117.70
1	13	1416	G	O5'-P-OP2	5.46	117.25	110.70
1	13	1417	G	C4-N9-C1'	5.46	133.59	126.50
26	1H	777	A	C6-N1-C2	-5.46	115.33	118.60
26	1H	2579	C	C4-C5-C6	5.46	120.13	117.40
1	13	751	U	O5'-P-OP1	-5.45	100.79	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	204	A	C5-C6-N6	-5.45	119.34	123.70
26	1H	587	C	C6-N1-C2	-5.45	118.12	120.30
26	1H	2582	G	OP1-P-OP2	-5.45	111.42	119.60
26	14	140	A	OP2-P-O3'	5.45	117.20	105.20
26	14	559	G	C4-C5-N7	-5.45	108.62	110.80
26	14	775	G	N1-C6-O6	-5.45	116.63	119.90
26	14	2328	A	C8-N9-C4	5.45	107.98	105.80
1	13	263	A	O5'-P-OP2	5.45	117.24	110.70
26	1H	49	A	C4-C5-C6	5.45	119.73	117.00
26	14	2007	C	N1-C2-O2	-5.45	115.63	118.90
26	1H	382	G	OP1-P-O3'	5.45	117.19	105.20
26	1H	427	U	N1-C2-N3	-5.45	111.63	114.90
26	1H	1272	A	C6-C5-N7	5.45	136.12	132.30
26	1H	1423	G	N9-C4-C5	-5.45	103.22	105.40
26	1H	1644	C	C6-N1-C1'	-5.45	114.26	120.80
26	14	2624	G	C5-C6-N1	5.45	114.23	111.50
24	3K	32	U	C6-N1-C2	-5.45	117.73	121.00
26	1H	1491	G	N1-C6-O6	5.45	123.17	119.90
26	1H	2280	G	OP1-P-O3'	5.45	117.19	105.20
49	K8	61	LEU	CA-CB-CG	-5.45	102.77	115.30
26	14	2316	C	O5'-P-OP1	-5.45	100.80	105.70
1	13	1222	G	O5'-P-OP1	5.45	117.24	110.70
26	14	678	C	C6-N1-C2	5.45	122.48	120.30
1	13	1521	G	O5'-P-OP1	-5.45	100.80	105.70
26	1H	70	G	N1-C2-N2	-5.45	111.30	116.20
26	1H	2392	A	O4'-C1'-N9	5.45	112.56	108.20
26	1H	2401	U	O5'-P-OP1	-5.45	100.80	105.70
26	1H	2502	G	OP2-P-O3'	5.45	117.18	105.20
27	16	65	C	C2-N1-C1'	5.45	124.79	118.80
50	L8	31	LEU	CA-CB-CG	5.45	127.82	115.30
26	14	1302	A	OP1-P-OP2	5.45	127.77	119.60
26	14	1780	A	C8-N9-C4	-5.45	103.62	105.80
26	14	2249	U	C5-C4-O4	5.45	129.17	125.90
26	1H	1516	U	N3-C2-O2	-5.44	118.39	122.20
26	14	1821	A	N1-C2-N3	5.44	132.02	129.30
1	13	586	C	C5-C6-N1	-5.44	118.28	121.00
26	1H	593	G	C4-C5-C6	5.44	122.07	118.80
26	1H	798	G	N3-C4-C5	5.44	131.32	128.60
26	1H	1193	G	C4-C5-N7	-5.44	108.62	110.80
1	1G	377	G	N3-C4-N9	5.44	129.27	126.00
26	14	1347	G	OP1-P-O3'	5.44	117.17	105.20
26	14	2336	A	O4'-C1'-N9	-5.44	103.85	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2713	A	N3-C4-C5	5.44	130.61	126.80
27	1J	38	C	O5'-P-OP2	-5.44	100.80	105.70
1	13	1497	G	O5'-P-OP2	-5.44	100.80	105.70
26	1H	509	C	O5'-P-OP2	-5.44	100.80	105.70
26	14	1605	C	N1-C2-O2	-5.44	115.64	118.90
26	14	2464	C	C5-C6-N1	-5.44	118.28	121.00
26	1H	473	G	N1-C2-N2	-5.44	111.31	116.20
1	13	715	A	O5'-P-OP2	-5.44	100.81	105.70
1	13	1214	C	C6-N1-C2	5.44	122.47	120.30
26	1H	323	G	O5'-P-OP1	-5.44	100.81	105.70
26	1H	600	G	N3-C4-C5	5.44	131.32	128.60
26	1H	1189	A	N9-C4-C5	-5.44	103.62	105.80
26	1H	2030	A	OP1-P-O3'	5.44	117.16	105.20
26	1H	2238	G	OP2-P-O3'	5.44	117.16	105.20
1	1G	413	G	C5-N7-C8	5.44	107.02	104.30
26	14	2430	A	N9-C4-C5	-5.44	103.62	105.80
27	1J	81	G	C6-C5-N7	-5.44	127.14	130.40
37	88	2	LEU	N-CA-C	-5.44	96.32	111.00
26	14	970	C	N1-C2-O2	-5.44	115.64	118.90
26	14	1904	G	C8-N9-C4	5.44	108.57	106.40
26	1H	585	G	OP1-P-OP2	-5.43	111.45	119.60
26	1H	758	C	C5-C4-N4	-5.43	116.40	120.20
26	1H	973	A	N9-C4-C5	-5.43	103.63	105.80
28	11	236	GLY	N-CA-C	5.43	126.69	113.10
26	14	776	G	O4'-C1'-N9	-5.43	103.85	108.20
26	14	1254	A	O5'-P-OP2	-5.43	100.81	105.70
26	14	1954	G	C6-C5-N7	5.43	133.66	130.40
1	13	18	C	C5-C6-N1	5.43	123.72	121.00
26	1H	1663	C	C5-C4-N4	-5.43	116.40	120.20
26	1H	1827	C	C4-C5-C6	5.43	120.12	117.40
26	1H	2688	U	C5-C6-N1	-5.43	119.98	122.70
26	14	385	C	P-O3'-C3'	5.43	126.22	119.70
26	14	1980	G	N1-C6-O6	5.43	123.16	119.90
26	14	2051	A	C8-N9-C4	-5.43	103.63	105.80
26	14	2378	A	N9-C4-C5	-5.43	103.63	105.80
27	1J	89(A)	A	C2-N3-C4	5.43	113.32	110.60
26	1H	780	G	C4-C5-N7	5.43	112.97	110.80
26	1H	802	A	N9-C4-C5	-5.43	103.63	105.80
26	1H	1157	G	C6-C5-N7	-5.43	127.14	130.40
26	1H	1625	C	N1-C2-O2	5.43	122.16	118.90
26	1H	2072	G	C6-N1-C2	5.43	128.36	125.10
26	1H	743	G	C4-C5-N7	-5.43	108.63	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1193	G	C6-N1-C2	-5.43	121.84	125.10
26	1H	2721	A	N9-C4-C5	-5.43	103.63	105.80
26	14	807	U	C4-C5-C6	5.43	122.96	119.70
26	14	1382	G	C4-N9-C1'	-5.43	119.44	126.50
26	14	1760	A	O5'-P-OP2	-5.43	100.81	105.70
1	13	22	G	N3-C2-N2	-5.43	116.10	119.90
1	13	916	G	C4-N9-C1'	5.43	133.56	126.50
1	13	917	G	O5'-P-OP1	-5.43	100.81	105.70
26	1H	731	C	C2-N3-C4	-5.43	117.19	119.90
26	14	2286	A	C5-N7-C8	-5.43	101.19	103.90
26	14	2351	G	C4-N9-C1'	5.43	133.56	126.50
26	14	2702	U	N1-C1'-C2'	5.43	121.06	114.00
26	14	2730	C	N3-C2-O2	-5.43	118.10	121.90
1	13	1356	G	C8-N9-C4	-5.43	104.23	106.40
26	1H	1157	G	C8-N9-C1'	-5.43	119.95	127.00
26	1H	2388	A	C2-N3-C4	-5.43	107.89	110.60
26	14	1142(A)	A	N1-C2-N3	5.43	132.01	129.30
1	13	1307	U	OP1-P-O3'	5.42	117.13	105.20
26	1H	863	A	OP1-P-OP2	5.42	127.74	119.60
26	1H	2307	G	C4-C5-N7	5.42	112.97	110.80
26	14	2092	U	C6-N1-C2	-5.42	117.75	121.00
26	14	2327	A	N9-C4-C5	5.42	107.97	105.80
26	1H	763	G	OP2-P-O3'	5.42	117.13	105.20
26	1H	1674	G	O4'-C1'-N9	-5.42	103.86	108.20
1	1G	979	C	C4-C5-C6	5.42	120.11	117.40
26	14	215	G	N9-C4-C5	-5.42	103.23	105.40
1	13	1498	U	C6-N1-C1'	-5.42	113.61	121.20
26	1H	135	G	N7-C8-N9	-5.42	110.39	113.10
26	1H	680	G	C8-N9-C4	5.42	108.57	106.40
26	1H	956	G	N1-C6-O6	5.42	123.15	119.90
26	1H	1365	A	C5-C6-N1	-5.42	114.99	117.70
26	1H	2404	C	O5'-P-OP1	-5.42	100.82	105.70
26	1H	2686	G	N3-C4-N9	5.42	129.25	126.00
48	J8	85	LEU	CA-CB-CG	-5.42	102.83	115.30
26	14	1251	C	O5'-P-OP1	-5.42	100.82	105.70
26	14	1404	C	N1-C2-O2	5.42	122.15	118.90
26	1H	660	G	C4-C5-N7	5.42	112.97	110.80
26	1H	1429	G	C5-C6-O6	5.42	131.85	128.60
26	14	1154	G	C4-C5-N7	5.42	112.97	110.80
1	13	640	A	O5'-P-OP1	-5.42	100.82	105.70
26	1H	177	G	C8-N9-C4	-5.42	104.23	106.40
26	1H	2311	A	N7-C8-N9	5.42	116.51	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2329	G	OP1-P-OP2	5.42	127.73	119.60
26	1H	2476	A	N7-C8-N9	5.42	116.51	113.80
1	1G	501	C	C6-N1-C2	-5.42	118.13	120.30
26	14	948	G	C8-N9-C4	-5.42	104.23	106.40
26	14	2712(A)	A	C4-C5-C6	-5.42	114.29	117.00
38	55	79	LEU	CA-CB-CG	5.42	127.76	115.30
26	1H	119	A	C5-N7-C8	5.42	106.61	103.90
26	1H	386	G	O5'-P-OP1	-5.42	100.83	105.70
26	1H	1192	G	N3-C4-N9	5.42	129.25	126.00
26	1H	1293	C	OP1-P-OP2	-5.42	111.48	119.60
26	1H	1428	C	O5'-P-OP1	-5.42	100.83	105.70
26	1H	1520	U	OP2-P-O3'	5.42	117.12	105.20
26	1H	2035	G	O5'-P-OP1	-5.42	100.83	105.70
26	14	59	U	OP2-P-O3'	5.42	117.11	105.20
26	1H	215	G	C2-N3-C4	-5.42	109.19	111.90
26	1H	732	C	C5-C6-N1	-5.42	118.29	121.00
27	16	61	G	N9-C4-C5	5.42	107.57	105.40
26	1H	946	G	C6-C5-N7	5.41	133.65	130.40
26	1H	1899	G	OP2-P-O3'	5.41	117.11	105.20
26	1H	2384	G	C8-N9-C4	-5.41	104.23	106.40
22	1L	74	C	C2-N1-C1'	-5.41	112.84	118.80
26	14	53	A	O5'-P-OP1	-5.41	100.83	105.70
26	14	102	G	O5'-P-OP1	-5.41	100.83	105.70
26	14	804	A	N7-C8-N9	-5.41	111.09	113.80
26	14	1653	G	N7-C8-N9	-5.41	110.39	113.10
1	13	1432	G	N3-C4-N9	5.41	129.25	126.00
26	1H	779	U	N3-C4-O4	-5.41	115.61	119.40
26	1H	844	C	N3-C4-C5	-5.41	119.73	121.90
26	1H	1607	C	C5-C4-N4	-5.41	116.41	120.20
26	14	788	A	N7-C8-N9	5.41	116.50	113.80
26	14	1309	G	C8-N9-C4	5.41	108.56	106.40
1	13	896	C	C2-N3-C4	-5.41	117.20	119.90
26	1H	498	G	O5'-P-OP2	5.41	117.19	110.70
26	1H	818	G	N9-C4-C5	5.41	107.56	105.40
26	1H	1265	A	N9-C4-C5	5.41	107.96	105.80
26	14	1586	A	N7-C8-N9	5.41	116.50	113.80
26	1H	811	U	N1-C2-N3	5.41	118.14	114.90
36	78	88	LEU	CA-CB-CG	5.41	127.73	115.30
1	1G	720	C	C6-N1-C2	-5.41	118.14	120.30
1	1G	1374	A	O4'-C1'-N9	5.41	112.53	108.20
1	13	1525	G	C8-N9-C4	5.41	108.56	106.40
26	1H	202	U	C6-N1-C1'	-5.41	113.63	121.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	583	G	C5-C6-N1	5.41	114.20	111.50
26	1H	2071	A	C5-C6-N6	-5.41	119.38	123.70
26	14	1630(A)	C	C6-N1-C2	-5.41	118.14	120.30
26	14	1971	A	C5-C6-N1	5.41	120.40	117.70
26	14	2329	G	C6-N1-C2	-5.41	121.86	125.10
26	1H	841	A	C4-C5-N7	5.40	113.40	110.70
26	1H	1406	U	OP1-P-O3'	5.40	117.09	105.20
26	14	1975	G	N9-C4-C5	-5.40	103.24	105.40
26	14	1664	A	O4'-C1'-N9	-5.40	103.88	108.20
26	14	1756	G	N1-C6-O6	5.40	123.14	119.90
26	1H	2244	U	C4-C5-C6	5.40	122.94	119.70
26	14	141	A	OP2-P-O3'	5.40	117.08	105.20
46	D5	24	LEU	CA-CB-CG	5.40	127.72	115.30
26	14	476	G	OP1-P-OP2	5.40	127.70	119.60
26	14	2217	G	C5-C6-O6	-5.40	125.36	128.60
26	1H	124	G	C5-C6-O6	-5.40	125.36	128.60
26	1H	1019	U	C6-N1-C2	-5.40	117.76	121.00
26	1H	2613	U	C4-C5-C6	5.40	122.94	119.70
26	1H	2779	U	N3-C2-O2	-5.40	118.42	122.20
1	1G	1498	U	C2-N1-C1'	5.40	124.18	117.70
26	14	1768	U	C2-N1-C1'	-5.40	111.22	117.70
1	13	325	A	C4-C5-C6	-5.40	114.30	117.00
26	1H	424	G	N1-C6-O6	-5.40	116.66	119.90
26	1H	2028	U	N1-C2-N3	5.40	118.14	114.90
26	14	736	C	C4-C5-C6	-5.40	114.70	117.40
1	13	402	G	O5'-P-OP1	5.39	117.17	110.70
26	1H	871	U	O5'-P-OP2	5.39	117.17	110.70
27	16	101	A	N1-C2-N3	5.39	132.00	129.30
1	1G	175	C	C6-N1-C2	-5.39	118.14	120.30
26	14	138	G	OP1-P-O3'	5.39	117.07	105.20
26	14	1784	A	OP1-P-O3'	5.39	117.07	105.20
26	1H	528	A	OP1-P-O3'	5.39	117.06	105.20
26	1H	735	A	C4-C5-C6	5.39	119.70	117.00
26	1H	989	G	N1-C6-O6	5.39	123.14	119.90
26	1H	1357	U	OP1-P-OP2	5.39	127.69	119.60
26	1H	1470	G	OP2-P-O3'	5.39	117.06	105.20
26	1H	1658	C	N3-C4-C5	-5.39	119.74	121.90
26	1H	2362	G	N1-C6-O6	-5.39	116.67	119.90
27	16	85	G	C5-C6-N1	5.39	114.20	111.50
36	78	71	VAL	C-N-CD	5.39	139.72	128.40
26	14	679	C	C2-N3-C4	-5.39	117.20	119.90
26	14	1281	G	N1-C6-O6	5.39	123.14	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1939	U	C2-N1-C1'	-5.39	111.23	117.70
26	14	2724	C	N1-C2-O2	-5.39	115.66	118.90
26	1H	133	C	O5'-P-OP1	5.39	117.17	110.70
26	14	2281	C	O5'-P-OP2	5.39	117.17	110.70
1	13	362	G	O5'-P-OP1	-5.39	100.85	105.70
26	1H	784	A	P-O3'-C3'	5.39	126.17	119.70
26	1H	1249	U	C5-C6-N1	-5.39	120.00	122.70
26	1H	1782	C	O5'-P-OP2	5.39	117.17	110.70
26	1H	2244	U	OP2-P-O3'	5.39	117.06	105.20
37	88	82	ARG	NE-CZ-NH2	5.39	123.00	120.30
1	1G	1048	G	C8-N9-C4	-5.39	104.24	106.40
26	14	2074	U	OP1-P-OP2	5.39	127.69	119.60
2	12	23	ARG	N-CA-C	-5.39	96.45	111.00
26	14	2607	G	C4-C5-C6	5.39	122.03	118.80
1	13	798	G	C2-N3-C4	-5.39	109.21	111.90
1	13	888	G	N1-C2-N2	-5.39	111.35	116.20
26	1H	818	G	O5'-P-OP1	-5.39	100.85	105.70
26	1H	928	G	N1-C2-N2	5.39	121.05	116.20
26	1H	1799	G	P-O3'-C3'	5.39	126.16	119.70
26	1H	2331	G	N1-C2-N3	5.39	127.13	123.90
26	1H	2501	C	N3-C2-O2	5.39	125.67	121.90
26	1H	2584	U	C5-C6-N1	-5.39	120.01	122.70
26	1H	2639	A	C6-C5-N7	-5.39	128.53	132.30
26	1H	2641	G	C5-C6-O6	-5.39	125.37	128.60
26	14	1537	C	C6-N1-C2	-5.39	118.14	120.30
26	14	1979	C	N3-C4-N4	5.39	121.77	118.00
26	1H	139	G	C8-N9-C4	-5.38	104.25	106.40
26	1H	446	G	C2-N3-C4	-5.38	109.21	111.90
26	1H	755	C	N1-C2-O2	-5.38	115.67	118.90
26	1H	963	U	OP1-P-OP2	-5.38	111.53	119.60
26	1H	1202	C	N1-C2-O2	-5.38	115.67	118.90
26	1H	2397	G	C4-C5-N7	5.38	112.95	110.80
26	1H	2437	U	C4-C5-C6	5.38	122.93	119.70
26	14	246	C	N3-C4-C5	5.38	124.05	121.90
26	14	954	G	C6-C5-N7	5.38	133.63	130.40
26	14	1899	G	N3-C4-N9	-5.38	122.77	126.00
26	1H	1901	A	C6-N1-C2	-5.38	115.37	118.60
26	1H	1993	U	C5-C6-N1	-5.38	120.01	122.70
26	14	1660	C	OP1-P-OP2	5.38	127.67	119.60
26	1H	187	G	C4-N9-C1'	5.38	133.50	126.50
26	1H	430	G	OP1-P-O3'	5.38	117.04	105.20
26	1H	620	G	OP1-P-OP2	5.38	127.67	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1417	C	C6-N1-C2	-5.38	118.15	120.30
26	1H	2585	U	C5-C4-O4	5.38	129.13	125.90
27	16	78	A	N3-C4-C5	5.38	130.57	126.80
1	1G	12	U	N3-C2-O2	-5.38	118.43	122.20
26	14	141	A	C5-N7-C8	-5.38	101.21	103.90
26	14	2674	G	O5'-P-OP2	-5.38	100.86	105.70
26	1H	747	U	O5'-P-OP1	-5.38	100.86	105.70
1	1G	1474	G	N1-C6-O6	5.38	123.13	119.90
26	14	696	G	O5'-P-OP2	5.38	117.16	110.70
22	1K	76	A	C5-N7-C8	-5.38	101.21	103.90
26	1H	1003	G	OP1-P-OP2	5.38	127.67	119.60
26	1H	2518	A	N3-C4-N9	-5.38	123.10	127.40
26	14	948	G	N1-C2-N2	5.38	121.04	116.20
26	14	982	C	OP1-P-O3'	5.38	117.03	105.20
26	14	1776	G	N3-C4-N9	5.38	129.23	126.00
1	13	1518	A	C5-N7-C8	5.38	106.59	103.90
26	1H	760	G	N1-C6-O6	5.38	123.13	119.90
26	1H	2030	A	O4'-C1'-N9	-5.38	103.90	108.20
26	1H	2597	G	OP2-P-O3'	5.38	117.03	105.20
55	Q8	43	GLN	N-CA-C	5.38	125.52	111.00
27	1J	102	G	N7-C8-N9	-5.38	110.41	113.10
26	14	676	A	N3-C4-C5	5.38	130.56	126.80
26	14	1125	G	C4-C5-N7	-5.38	108.65	110.80
26	14	2445	G	C5-C6-N1	5.38	114.19	111.50
26	14	2446	G	O5'-P-OP2	-5.38	100.86	105.70
1	13	1187	G	C8-N9-C1'	-5.37	120.01	127.00
26	1H	94	G	N1-C6-O6	5.37	123.12	119.90
26	1H	859	G	C4-N9-C1'	-5.37	119.52	126.50
26	1H	1129	A	C5-C6-N6	-5.37	119.40	123.70
26	1H	2447	G	C5-C6-N1	5.37	114.19	111.50
1	1G	230	G	C5-C6-N1	-5.37	108.81	111.50
26	14	186	G	C5-C6-O6	-5.37	125.38	128.60
26	14	2330	G	C2-N3-C4	-5.37	109.21	111.90
26	14	2681	C	C5-C6-N1	-5.37	118.31	121.00
1	13	900	A	OP1-P-OP2	-5.37	111.54	119.60
26	1H	616	A	OP1-P-OP2	5.37	127.66	119.60
26	1H	752	A	N1-C2-N3	5.37	131.99	129.30
26	1H	782	A	C2-N3-C4	-5.37	107.91	110.60
26	1H	1854	A	N1-C6-N6	-5.37	115.38	118.60
26	1H	2064	C	N3-C4-C5	-5.37	119.75	121.90
1	1G	662	G	C4-N9-C1'	5.37	133.48	126.50
26	14	252	G	C2-N3-C4	5.37	114.59	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	987	G	C6-C5-N7	5.37	133.62	130.40
1	13	1497	G	O4'-C1'-N9	5.37	112.50	108.20
26	1H	215	G	N1-C6-O6	5.37	123.12	119.90
26	1H	645	C	C2-N1-C1'	5.37	124.71	118.80
26	1H	792	G	OP1-P-OP2	5.37	127.66	119.60
26	1H	1534	G	N3-C4-C5	-5.37	125.92	128.60
26	1H	2449	U	C4-C5-C6	5.37	122.92	119.70
26	14	1313	U	O4'-C1'-N1	5.37	112.50	108.20
1	13	974	A	N9-C4-C5	-5.37	103.65	105.80
26	1H	2010	G	OP1-P-O3'	5.37	117.01	105.20
26	1H	2578	G	N1-C6-O6	-5.37	116.68	119.90
38	98	18	LEU	CA-CB-CG	5.37	127.65	115.30
26	14	1704	G	N9-C4-C5	-5.37	103.25	105.40
26	14	2044	C	OP1-P-OP2	5.37	127.65	119.60
1	13	903	G	C6-N1-C2	-5.37	121.88	125.10
26	1H	1248	G	C4-C5-N7	-5.37	108.65	110.80
26	1H	2705	A	OP1-P-OP2	-5.37	111.55	119.60
26	14	2239	G	OP2-P-O3'	5.37	117.01	105.20
1	13	1496	C	N1-C2-O2	-5.37	115.68	118.90
26	1H	116	C	C4-C5-C6	5.37	120.08	117.40
26	1H	1294	U	C5-C6-N1	-5.37	120.02	122.70
26	1H	2607	G	N1-C6-O6	5.37	123.12	119.90
26	14	1570	A	C4-C5-N7	5.37	113.38	110.70
26	14	1797	C	C2-N3-C4	-5.37	117.22	119.90
26	14	2044	C	N3-C4-N4	5.37	121.75	118.00
26	1H	1855	G	N3-C4-N9	5.36	129.22	126.00
26	1H	1982	C	OP1-P-OP2	5.36	127.65	119.60
26	14	116	C	OP1-P-OP2	5.36	127.64	119.60
26	14	2443	C	C6-N1-C2	-5.36	118.15	120.30
26	1H	793	A	C6-C5-N7	-5.36	128.55	132.30
1	13	878	G	N3-C4-N9	5.36	129.22	126.00
26	1H	447	A	O5'-P-OP2	5.36	117.13	110.70
26	1H	1237	A	OP2-P-O3'	5.36	116.99	105.20
26	1H	2655	G	C4-N9-C1'	-5.36	119.53	126.50
27	16	24	G	N3-C4-N9	5.36	129.22	126.00
1	1G	402	G	C8-N9-C4	5.36	108.54	106.40
26	14	2096	U	C5-C6-N1	5.36	125.38	122.70
26	14	2454	G	O5'-P-OP2	-5.36	100.88	105.70
23	2K	9	G	C5-C6-N1	5.36	114.18	111.50
26	1H	530	G	N9-C4-C5	5.36	107.54	105.40
26	1H	1937	A	O5'-P-OP1	5.36	117.13	110.70
26	1H	2079	U	N1-C2-N3	5.36	118.11	114.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	41	98	ARG	NE-CZ-NH1	5.36	122.98	120.30
26	14	2501	C	C6-N1-C1'	5.36	127.23	120.80
26	1H	40	C	C4-C5-C6	5.36	120.08	117.40
26	1H	1236	G	OP1-P-O3'	5.36	116.99	105.20
1	1G	564	C	N3-C4-C5	-5.36	119.76	121.90
26	14	187	G	C8-N9-C1'	-5.36	120.04	127.00
1	13	226	G	C6-C5-N7	-5.36	127.19	130.40
26	1H	574	C	C2-N3-C4	5.36	122.58	119.90
26	1H	866	A	N9-C4-C5	-5.36	103.66	105.80
26	1H	970	C	OP1-P-O3'	-5.36	93.42	105.20
26	1H	1489	U	O5'-P-OP1	-5.36	100.88	105.70
26	1H	2689	U	P-O3'-C3'	5.36	126.13	119.70
1	1G	538	G	N1-C6-O6	5.36	123.11	119.90
26	14	945	A	C4-N9-C1'	5.36	135.94	126.30
26	14	2447	G	C5-C6-O6	-5.36	125.39	128.60
26	1H	67	U	C5-C6-N1	5.35	125.38	122.70
26	1H	397	G	N3-C4-N9	-5.35	122.79	126.00
26	1H	1021	A	N3-C4-N9	-5.35	123.12	127.40
26	1H	1612	C	N3-C4-C5	-5.35	119.76	121.90
26	1H	2266	A	OP1-P-OP2	-5.35	111.57	119.60
26	14	2603	G	OP1-P-O3'	5.35	116.98	105.20
26	14	1029	A	N1-C6-N6	5.35	121.81	118.60
26	14	1352	U	O5'-P-OP2	-5.35	100.88	105.70
1	13	1487	G	O5'-P-OP2	-5.35	100.88	105.70
26	14	2237	G	C8-N9-C1'	-5.35	120.04	127.00
1	13	804	U	OP2-P-O3'	5.35	116.97	105.20
1	13	818	G	C4-C5-N7	-5.35	108.66	110.80
26	1H	1646	C	O5'-P-OP1	-5.35	100.89	105.70
26	1H	2212	A	C6-C5-N7	-5.35	128.56	132.30
26	14	113	G	C5-C6-O6	-5.35	125.39	128.60
26	14	1145	C	C5-C6-N1	5.35	123.67	121.00
26	14	2377	A	C2-N3-C4	-5.35	107.92	110.60
1	13	1519	A	N9-C4-C5	5.35	107.94	105.80
24	3K	35	A	C8-N9-C4	-5.35	103.66	105.80
26	1H	1616	A	C2-N3-C4	-5.35	107.93	110.60
26	1H	2527	C	C5-C6-N1	5.35	123.67	121.00
27	16	33	G	OP1-P-O3'	5.35	116.97	105.20
41	C8	10	ARG	NE-CZ-NH1	-5.35	117.63	120.30
1	1G	337	C	C5-C4-N4	-5.35	116.46	120.20
26	14	179	G	O5'-P-OP2	-5.35	100.89	105.70
26	14	2232	U	C5-C4-O4	5.35	129.11	125.90
26	14	2426	A	N3-C4-N9	5.35	131.68	127.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2726	U	N3-C2-O2	-5.35	118.46	122.20
26	1H	467	G	N7-C8-N9	-5.35	110.43	113.10
26	1H	1935	G	C4-C5-N7	-5.35	108.66	110.80
27	16	86	G	C5-C6-N1	5.35	114.17	111.50
26	1H	1187	G	P-O3'-C3'	5.34	126.11	119.70
26	1H	2431	U	OP1-P-O3'	5.34	116.96	105.20
26	14	862	G	C5-C6-O6	5.34	131.81	128.60
26	14	1519	G	C4-C5-N7	-5.34	108.66	110.80
1	13	812	C	N3-C2-O2	-5.34	118.16	121.90
26	1H	1938	A	C5-C6-N6	-5.34	119.43	123.70
26	1H	1888	G	N1-C6-O6	-5.34	116.69	119.90
26	1H	2212	A	C5-N7-C8	-5.34	101.23	103.90
1	1G	43	C	C6-N1-C2	5.34	122.44	120.30
26	14	1026	U	C5-C6-N1	5.34	125.37	122.70
1	13	557	G	O5'-P-OP2	-5.34	100.89	105.70
1	13	923	A	O5'-P-OP1	-5.34	100.89	105.70
26	1H	856	C	N3-C4-N4	5.34	121.74	118.00
26	1H	1378	A	C5-N7-C8	-5.34	101.23	103.90
26	1H	1410	G	C8-N9-C1'	5.34	133.94	127.00
26	1H	2253	G	C5-C6-N1	5.34	114.17	111.50
1	1G	632	A	OP2-P-O3'	5.34	116.95	105.20
26	14	915	C	N3-C2-O2	-5.34	118.16	121.90
26	14	2356	C	C6-N1-C2	5.34	122.44	120.30
1	13	578	C	N1-C2-O2	-5.34	115.70	118.90
26	1H	742	G	C2-N3-C4	-5.34	109.23	111.90
26	1H	2069	G	N9-C4-C5	-5.34	103.27	105.40
26	1H	2614	A	N1-C2-N3	-5.34	126.63	129.30
30	31	38	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	13	726	C	O5'-P-OP1	-5.34	100.90	105.70
26	1H	200	U	N1-C2-N3	5.34	118.10	114.90
26	1H	1600	C	OP1-P-O3'	5.34	116.94	105.20
26	1H	1830	C	N3-C2-O2	5.34	125.64	121.90
26	1H	2403	C	C5-C6-N1	5.34	123.67	121.00
26	14	2420	C	O5'-P-OP1	-5.34	100.90	105.70
1	13	820	U	O5'-P-OP1	-5.33	100.90	105.70
26	1H	482	A	N7-C8-N9	5.33	116.47	113.80
26	1H	1456	G	N1-C6-O6	5.33	123.10	119.90
26	1H	2040	C	N3-C4-N4	5.33	121.73	118.00
26	14	991	C	OP1-P-OP2	5.33	127.60	119.60
1	13	484	G	P-O3'-C3'	5.33	126.10	119.70
26	1H	470	A	C6-C5-N7	-5.33	128.57	132.30
26	1H	563	G	C8-N9-C4	5.33	108.53	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1157	G	C4-N9-C1'	5.33	133.43	126.50
26	1H	2570	G	N3-C4-N9	-5.33	122.80	126.00
26	1H	2710	C	O5'-P-OP2	5.33	117.10	110.70
26	1H	2765	A	OP1-P-OP2	5.33	127.60	119.60
26	14	1588	C	C6-N1-C2	-5.33	118.17	120.30
26	14	2487	G	N1-C2-N2	-5.33	111.40	116.20
26	1H	621	A	C6-C5-N7	-5.33	128.57	132.30
26	14	779	U	C5-C6-N1	-5.33	120.03	122.70
26	14	912	C	C5-C6-N1	5.33	123.67	121.00
26	14	1904	G	N7-C8-N9	-5.33	110.43	113.10
26	14	2535	G	O5'-P-OP2	-5.33	100.90	105.70
48	F5	11	ARG	NE-CZ-NH1	-5.33	117.63	120.30
26	1H	1373	A	O5'-P-OP1	5.33	117.10	110.70
1	1G	1415	G	C8-N9-C4	5.33	108.53	106.40
26	14	1853	A	OP1-P-OP2	5.33	127.59	119.60
26	14	2330	G	N9-C4-C5	-5.33	103.27	105.40
1	13	1432	G	C4-N9-C1'	5.33	133.43	126.50
26	1H	138	G	O4'-C1'-N9	5.33	112.46	108.20
26	1H	746	A	C4-C5-N7	5.33	113.36	110.70
26	1H	1272	A	OP1-P-O3'	5.33	116.92	105.20
26	1H	1965	C	C4-C5-C6	-5.33	114.73	117.40
26	1H	2287	A	C5-N7-C8	-5.33	101.24	103.90
26	1H	2446	G	C6-C5-N7	-5.33	127.20	130.40
1	1G	413	G	C4-N9-C1'	-5.33	119.57	126.50
26	14	204	A	N1-C6-N6	5.33	121.80	118.60
26	14	1259	G	OP2-P-O3'	5.33	116.92	105.20
26	14	1336	A	O5'-P-OP2	-5.33	100.91	105.70
26	14	2234	G	N3-C4-C5	-5.33	125.94	128.60
1	13	989	C	N1-C2-O2	5.33	122.09	118.90
26	1H	415	A	N9-C4-C5	-5.33	103.67	105.80
26	1H	576	U	N1-C2-O2	-5.33	119.07	122.80
26	1H	1270	C	C4-C5-C6	5.33	120.06	117.40
26	1H	1644	C	C2-N1-C1'	5.33	124.66	118.80
26	14	150	C	C6-N1-C1'	5.33	127.19	120.80
26	14	689	A	O5'-P-OP2	-5.33	100.91	105.70
26	14	725	G	C4-C5-N7	-5.33	108.67	110.80
26	1H	125	G	N3-C2-N2	5.32	123.63	119.90
26	1H	271(C)	U	C5-C6-N1	5.32	125.36	122.70
26	1H	2345	G	N9-C4-C5	-5.32	103.27	105.40
1	1G	1446	A	O4'-C1'-N9	5.32	112.46	108.20
26	14	1142	U	C6-N1-C1'	-5.32	113.75	121.20
26	1H	198	C	C5-C4-N4	-5.32	116.47	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1286	A	OP2-P-O3'	5.32	116.91	105.20
27	16	80	U	N1-C2-O2	5.32	126.53	122.80
1	1G	261	U	N3-C2-O2	-5.32	118.47	122.20
1	1G	1400	C	N3-C4-C5	5.32	124.03	121.90
1	13	1498	U	N1-C2-O2	5.32	126.53	122.80
1	13	1498	U	O4'-C1'-N1	-5.32	103.94	108.20
26	1H	255	A	OP2-P-O3'	5.32	116.91	105.20
26	1H	807	U	C4-C5-C6	5.32	122.89	119.70
26	1H	2485	G	N1-C2-N3	5.32	127.09	123.90
27	16	12	C	C4-C5-C6	5.32	120.06	117.40
26	14	189	G	N3-C4-C5	5.32	131.26	128.60
26	14	677	A	O5'-P-OP2	-5.32	100.91	105.70
26	14	1755	A	C2-N3-C4	-5.32	107.94	110.60
26	14	1840	G	C5-C6-N1	-5.32	108.84	111.50
1	13	582	U	C2-N3-C4	-5.32	123.81	127.00
26	1H	1970	A	C8-N9-C4	-5.32	103.67	105.80
26	14	1394	U	C2-N1-C1'	-5.32	111.32	117.70
42	95	49	THR	C-N-CD	5.32	139.57	128.40
1	13	541	G	N1-C6-O6	5.32	123.09	119.90
1	13	899	C	C2-N1-C1'	-5.32	112.95	118.80
1	13	900	A	C8-N9-C4	5.32	107.93	105.80
26	1H	417	C	C5-C4-N4	-5.32	116.48	120.20
26	1H	768	G	OP1-P-OP2	5.32	127.58	119.60
26	1H	1894	C	C5-C6-N1	5.32	123.66	121.00
26	14	139	G	C4-C5-N7	-5.32	108.67	110.80
26	14	247	G	C5-C6-N1	-5.32	108.84	111.50
26	14	2000	G	C5-N7-C8	5.32	106.96	104.30
1	13	880	C	N3-C4-C5	5.32	124.03	121.90
23	2K	27	G	C5-C6-O6	-5.32	125.41	128.60
26	1H	1301	A	N9-C4-C5	-5.32	103.67	105.80
26	1H	1313	U	C2-N1-C1'	5.32	124.08	117.70
26	1H	2033	A	C2-N3-C4	-5.32	107.94	110.60
26	14	187	G	N3-C4-C5	-5.32	125.94	128.60
26	14	1786	A	OP1-P-O3'	5.32	116.89	105.20
26	14	1840	G	C2-N3-C4	-5.32	109.24	111.90
26	14	2427	C	OP2-P-O3'	5.32	116.89	105.20
26	1H	2244	U	N1-C2-N3	5.31	118.09	114.90
1	13	1530	G	O5'-P-OP2	-5.31	100.92	105.70
26	1H	115	C	N1-C2-O2	-5.31	115.71	118.90
26	1H	910	A	N9-C4-C5	-5.31	103.67	105.80
26	1H	960	A	C8-N9-C4	5.31	107.92	105.80
26	1H	2280	G	C4-C5-N7	-5.31	108.67	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2468	G	O5'-P-OP1	5.31	117.07	110.70
26	1H	2717	G	C8-N9-C4	5.31	108.53	106.40
26	14	1282	U	C5-C6-N1	-5.31	120.04	122.70
27	1J	47	C	N3-C4-C5	5.31	124.03	121.90
26	1H	277	C	C2-N1-C1'	5.31	124.64	118.80
26	1H	1403	C	N3-C4-N4	-5.31	114.28	118.00
26	1H	2312	U	N3-C4-O4	5.31	123.12	119.40
26	1H	2427	C	O5'-P-OP1	-5.31	100.92	105.70
26	14	197	A	C5-C6-N6	-5.31	119.45	123.70
26	14	1436	G	N3-C4-N9	5.31	129.19	126.00
26	14	1663	C	N1-C2-O2	-5.31	115.71	118.90
1	13	771	G	N1-C2-N3	5.31	127.08	123.90
1	13	1276	G	C8-N9-C4	-5.31	104.28	106.40
26	1H	439	G	OP1-P-O3'	5.31	116.88	105.20
26	1H	1839	G	O4'-C1'-N9	-5.31	103.95	108.20
26	1H	2038	G	N3-C2-N2	5.31	123.62	119.90
26	1H	2329	G	C5-C6-O6	-5.31	125.41	128.60
1	1G	772	U	N3-C2-O2	-5.31	118.48	122.20
26	14	583	G	C8-N9-C4	-5.31	104.28	106.40
1	13	1475	G	O5'-P-OP1	-5.31	100.92	105.70
26	1H	564	C	C5-C6-N1	5.31	123.65	121.00
26	1H	1318	C	N3-C4-C5	5.31	124.02	121.90
26	1H	1645	G	C5-C6-N1	5.31	114.15	111.50
26	1H	1982	C	C5-C4-N4	-5.31	116.48	120.20
1	1G	481	G	C4-N9-C1'	5.31	133.40	126.50
26	14	226	G	O4'-C1'-N9	5.31	112.45	108.20
26	1H	1764	G	N1-C2-N2	-5.31	111.42	116.20
27	16	115	G	C4-C5-N7	5.31	112.92	110.80
25	4L	15	A	N1-C6-N6	5.31	121.78	118.60
26	14	115	C	C5-C6-N1	-5.30	118.35	121.00
26	14	594	U	OP2-P-O3'	5.30	116.87	105.20
26	14	621	A	C4-C5-N7	5.30	113.35	110.70
26	14	665	C	N1-C2-O2	5.30	122.08	118.90
26	14	1474	C	C6-N1-C2	-5.30	118.18	120.30
26	14	1981	A	C2-N3-C4	-5.30	107.95	110.60
22	1K	34	G	C4-N9-C1'	5.30	133.39	126.50
26	1H	1786	A	OP1-P-O3'	5.30	116.87	105.20
26	1H	2617	C	N1-C2-O2	-5.30	115.72	118.90
26	14	929	G	N3-C4-N9	5.30	129.18	126.00
26	1H	400	G	C5-C6-O6	-5.30	125.42	128.60
26	1H	2230	G	N3-C4-C5	5.30	131.25	128.60
1	1G	317	G	N7-C8-N9	5.30	115.75	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	436	C	C6-N1-C2	5.30	122.42	120.30
26	14	1204	A	N1-C2-N3	5.30	131.95	129.30
26	14	1836	C	OP1-P-O3'	5.30	116.86	105.20
26	1H	177	G	N1-C6-O6	-5.30	116.72	119.90
26	1H	500	G	OP1-P-OP2	5.30	127.55	119.60
26	1H	796	C	O5'-P-OP1	5.30	117.06	110.70
26	1H	1465	G	O5'-P-OP2	-5.30	100.93	105.70
26	14	829	A	OP1-P-OP2	5.30	127.55	119.60
26	14	1216	G	C4-N9-C1'	5.30	133.39	126.50
26	14	1291	C	O5'-P-OP2	-5.30	100.93	105.70
26	14	1323	U	N1-C2-O2	-5.30	119.09	122.80
26	14	1342	A	OP1-P-OP2	-5.30	111.65	119.60
26	14	2086	U	N1-C2-O2	5.30	126.51	122.80
1	13	1395	C	P-O3'-C3'	5.30	126.06	119.70
26	1H	136	G	N9-C4-C5	-5.30	103.28	105.40
26	1H	794	G	C4-C5-N7	-5.30	108.68	110.80
26	1H	1225	C	N3-C4-N4	-5.30	114.29	118.00
26	1H	1658	C	N3-C2-O2	5.30	125.61	121.90
32	51	82	GLY	N-CA-C	5.30	126.34	113.10
26	14	1373	A	N7-C8-N9	-5.30	111.15	113.80
26	1H	634	C	N3-C4-N4	-5.29	114.29	118.00
26	1H	2082	A	N7-C8-N9	-5.29	111.15	113.80
1	1G	1481	U	O5'-P-OP1	-5.29	100.93	105.70
26	14	2332	U	O5'-P-OP2	-5.29	100.94	105.70
26	14	2713	A	C5-C6-N1	-5.29	115.05	117.70
1	13	112	G	C8-N9-C4	-5.29	104.28	106.40
1	13	328	C	C6-N1-C1'	-5.29	114.45	120.80
26	1H	1197	G	N7-C8-N9	-5.29	110.45	113.10
26	1H	2040	C	N3-C2-O2	5.29	125.61	121.90
26	14	298	G	N1-C6-O6	5.29	123.08	119.90
26	14	1198	U	N3-C2-O2	-5.29	118.49	122.20
26	14	2514	U	O5'-P-OP2	5.29	117.05	110.70
26	1H	1639	U	C5-C4-O4	5.29	129.07	125.90
1	1G	666	G	C6-C5-N7	-5.29	127.22	130.40
26	14	1204	A	N7-C8-N9	5.29	116.45	113.80
26	14	2360	A	C8-N9-C4	5.29	107.92	105.80
26	1H	737	C	C2-N3-C4	-5.29	117.25	119.90
26	1H	125	G	O4'-C1'-N9	-5.29	103.97	108.20
26	1H	746	A	O4'-C1'-N9	5.29	112.43	108.20
26	1H	2593	U	N3-C2-O2	-5.29	118.50	122.20
26	1H	2622	C	C5-C6-N1	-5.29	118.36	121.00
29	21	186	GLY	N-CA-C	5.29	126.32	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	1498	U	O4'-C1'-N1	-5.29	103.97	108.20
26	14	810	U	OP1-P-O3'	5.29	116.83	105.20
1	1G	1488	G	C8-N9-C1'	-5.29	120.13	127.00
26	14	447	A	O4'-C1'-N9	-5.29	103.97	108.20
26	14	584	C	C6-N1-C2	5.29	122.42	120.30
26	14	2675	A	C2-N3-C4	-5.29	107.96	110.60
1	13	572	A	N1-C6-N6	-5.29	115.43	118.60
1	13	836	G	N9-C4-C5	-5.29	103.29	105.40
26	1H	1451	C	N3-C2-O2	-5.29	118.20	121.90
26	1H	1782	C	C4-C5-C6	5.29	120.04	117.40
27	16	7	G	C5-C6-O6	-5.29	125.43	128.60
1	1G	353	A	N7-C8-N9	5.29	116.44	113.80
26	14	267	C	C6-N1-C2	-5.29	118.19	120.30
26	14	1297	C	C2-N1-C1'	-5.29	112.99	118.80
27	1J	30	C	N3-C2-O2	-5.29	118.20	121.90
26	1H	1791	A	OP1-P-OP2	-5.28	111.68	119.60
26	1H	1805	U	N3-C4-O4	5.28	123.10	119.40
26	1H	2287	A	C6-C5-N7	-5.28	128.60	132.30
45	G8	79	CYS	N-CA-C	5.28	125.27	111.00
1	1G	518	C	O5'-P-OP1	5.28	117.04	110.70
26	14	1857	G	C6-C5-N7	-5.28	127.23	130.40
28	19	43	ARG	CG-CD-NE	5.28	122.89	111.80
1	13	1495	U	N3-C4-O4	-5.28	115.70	119.40
26	1H	1562	A	N9-C4-C5	-5.28	103.69	105.80
26	1H	2071	A	C6-N1-C2	-5.28	115.43	118.60
26	14	1235	G	N1-C6-O6	-5.28	116.73	119.90
1	13	936	C	C6-N1-C2	-5.28	118.19	120.30
26	1H	388	G	OP2-P-O3'	5.28	116.82	105.20
28	11	229	VAL	CB-CA-C	-5.28	101.37	111.40
1	1G	631	G	N7-C8-N9	5.28	115.74	113.10
1	1G	1062	U	O5'-P-OP2	-5.28	100.95	105.70
26	14	614	U	C6-N1-C2	-5.28	117.83	121.00
26	14	1240	U	O5'-P-OP2	-5.28	100.95	105.70
26	14	2044	C	C5-C4-N4	-5.28	116.50	120.20
26	14	2578	G	N1-C2-N2	-5.28	111.45	116.20
1	13	354	G	C6-C5-N7	-5.28	127.23	130.40
1	13	380	G	N3-C4-N9	-5.28	122.83	126.00
1	13	1508	G	N3-C4-N9	-5.28	122.83	126.00
26	1H	430	G	C6-C5-N7	-5.28	127.23	130.40
26	1H	1124	C	C5-C6-N1	-5.28	118.36	121.00
26	1H	1416	G	C8-N9-C4	5.28	108.51	106.40
26	1H	2620	C	C5-C4-N4	-5.28	116.50	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	778	G	OP1-P-O3'	5.28	116.81	105.20
26	14	738	G	N1-C2-N2	-5.28	111.45	116.20
26	14	1673	U	C2-N1-C1'	-5.28	111.37	117.70
26	14	2012	G	C5-C6-O6	-5.28	125.43	128.60
26	1H	563	G	C6-C5-N7	5.28	133.57	130.40
26	1H	1824	G	C5-N7-C8	5.28	106.94	104.30
26	14	429	A	N1-C6-N6	5.28	121.77	118.60
26	14	602	G	N3-C4-N9	5.28	129.17	126.00
26	14	613	U	N1-C2-O2	5.28	126.50	122.80
26	14	1414	G	C6-C5-N7	-5.28	127.23	130.40
26	14	1961	C	N3-C2-O2	5.28	125.59	121.90
1	13	511	C	N1-C2-O2	-5.28	115.73	118.90
26	1H	239	U	O5'-P-OP2	-5.28	100.95	105.70
26	1H	741	G	N1-C2-N3	5.28	127.07	123.90
26	1H	1528	A	C2-N3-C4	-5.28	107.96	110.60
26	1H	1609	A	N7-C8-N9	-5.28	111.16	113.80
26	1H	1799	G	N3-C4-C5	-5.28	125.96	128.60
26	1H	1942	C	C5-C6-N1	5.28	123.64	121.00
26	1H	2269	A	N1-C6-N6	5.28	121.77	118.60
26	14	2644	G	C2-N3-C4	-5.28	109.26	111.90
39	65	4	LEU	CA-CB-CG	-5.28	103.17	115.30
26	1H	872	A	C2-N3-C4	-5.27	107.96	110.60
26	1H	914	C	C5-C4-N4	5.27	123.89	120.20
26	1H	1849	G	O5'-P-OP2	5.27	117.03	110.70
1	1G	697	U	C5-C6-N1	-5.27	120.06	122.70
26	14	781	A	OP1-P-OP2	5.27	127.51	119.60
1	13	220	G	C4-N9-C1'	5.27	133.35	126.50
1	13	533	A	O4'-C1'-N9	5.27	112.42	108.20
1	13	811	C	C4-C5-C6	5.27	120.04	117.40
26	1H	911	A	C5-C6-N1	-5.27	115.06	117.70
26	1H	1184	G	OP2-P-O3'	5.27	116.80	105.20
26	1H	2743	C	N1-C2-O2	-5.27	115.74	118.90
41	C8	3	ARG	NE-CZ-NH1	5.27	122.94	120.30
23	2L	40	C	C5-C6-N1	5.27	123.64	121.00
26	14	1267	U	P-O3'-C3'	5.27	126.03	119.70
26	14	1679	U	OP1-P-O3'	5.27	116.80	105.20
1	13	652	U	C5-C6-N1	5.27	125.34	122.70
26	1H	392	C	OP1-P-O3'	5.27	116.80	105.20
26	1H	2265	U	C5-C6-N1	5.27	125.33	122.70
26	14	866	A	O4'-C1'-N9	-5.27	103.98	108.20
26	14	1992	G	C8-N9-C4	5.27	108.51	106.40
1	13	814	A	O5'-P-OP1	-5.27	100.96	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	815	A	C8-N9-C4	5.27	107.91	105.80
1	13	1488	G	N9-C4-C5	-5.27	103.29	105.40
23	2K	38	A	O5'-P-OP2	-5.27	100.96	105.70
26	1H	456	C	C2-N3-C4	-5.27	117.27	119.90
26	1H	992	C	C2-N1-C1'	5.27	124.60	118.80
27	16	95	U	C2-N1-C1'	-5.27	111.38	117.70
26	14	709	U	C5-C4-O4	5.27	129.06	125.90
26	14	771	G	N7-C8-N9	-5.27	110.47	113.10
26	14	1314	C	C2-N3-C4	5.27	122.53	119.90
26	14	1601	G	O5'-P-OP2	-5.27	100.96	105.70
26	1H	1573	G	N1-C6-O6	5.27	123.06	119.90
26	1H	2035	G	C5-C6-N1	-5.27	108.87	111.50
26	14	199	A	C2-N3-C4	5.27	113.23	110.60
26	14	265	A	C4-C5-N7	5.27	113.33	110.70
26	14	1352	U	C5-C6-N1	-5.27	120.07	122.70
26	14	2568	C	C6-N1-C2	-5.27	118.19	120.30
23	2K	15	G	C2-N3-C4	-5.27	109.27	111.90
26	1H	954	G	N3-C2-N2	-5.27	116.21	119.90
26	1H	2341	G	OP1-P-O3'	5.27	116.78	105.20
26	14	114	U	C2-N1-C1'	5.27	124.02	117.70
26	14	2351	G	C5-C6-O6	-5.27	125.44	128.60
1	13	465	A	N7-C8-N9	-5.26	111.17	113.80
26	1H	141	A	C5-C6-N6	-5.26	119.49	123.70
26	1H	586	A	C4-C5-N7	-5.26	108.07	110.70
26	1H	803	U	C5-C4-O4	-5.26	122.74	125.90
26	1H	928	G	C5-N7-C8	-5.26	101.67	104.30
26	1H	1662	C	C4-C5-C6	5.26	120.03	117.40
26	1H	2060	A	C4-C5-C6	-5.26	114.37	117.00
26	1H	2443	C	N3-C4-N4	5.26	121.69	118.00
26	1H	2751	G	C4-N9-C1'	5.26	133.34	126.50
1	1G	427	U	N3-C2-O2	-5.26	118.51	122.20
1	1G	546	G	C8-N9-C4	-5.26	104.29	106.40
26	14	742	G	N1-C2-N3	5.26	127.06	123.90
26	14	980	A	C6-C5-N7	-5.26	128.62	132.30
26	14	2430	A	C5-N7-C8	-5.26	101.27	103.90
26	1H	35	G	C5-C6-O6	5.26	131.76	128.60
1	13	263	A	O5'-P-OP1	-5.26	100.97	105.70
1	13	800	G	C5-C6-O6	-5.26	125.44	128.60
1	13	1306	A	O5'-P-OP2	-5.26	100.96	105.70
1	13	1513	A	OP2-P-O3'	5.26	116.78	105.20
26	1H	1925	C	C6-N1-C2	-5.26	118.19	120.30
37	88	19	GLY	N-CA-C	5.26	126.25	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
24	3L	71	G	C5'-C4'-O4'	5.26	115.41	109.10
26	14	773	U	OP1-P-OP2	-5.26	111.71	119.60
26	14	788	A	C8-N9-C4	-5.26	103.70	105.80
27	1J	98	G	C8-N9-C1'	-5.26	120.16	127.00
1	13	62	U	O5'-P-OP2	-5.26	100.97	105.70
1	13	538	G	N9-C4-C5	-5.26	103.30	105.40
26	1H	528	A	N7-C8-N9	5.26	116.43	113.80
26	1H	660	G	C8-N9-C4	-5.26	104.30	106.40
26	1H	1620	G	O5'-P-OP2	5.26	117.01	110.70
26	1H	2010	G	C5-C6-O6	-5.26	125.44	128.60
26	1H	2666	C	N3-C4-C5	-5.26	119.80	121.90
27	16	78	A	N7-C8-N9	-5.26	111.17	113.80
26	14	236	C	N1-C2-O2	-5.26	115.75	118.90
26	1H	128	C	C5-C6-N1	-5.26	118.37	121.00
26	1H	2072	G	N3-C4-N9	5.26	129.16	126.00
26	1H	2503	A	N1-C6-N6	5.26	121.75	118.60
26	14	1264	G	N1-C6-O6	-5.26	116.75	119.90
1	13	376	G	N3-C4-C5	-5.26	125.97	128.60
26	1H	398	G	C2-N3-C4	-5.26	109.27	111.90
26	1H	2018	G	C5-N7-C8	-5.26	101.67	104.30
26	1H	2239	G	OP1-P-OP2	-5.26	111.72	119.60
26	1H	2246	G	N3-C4-C5	-5.26	125.97	128.60
26	1H	2447	G	C2-N3-C4	5.26	114.53	111.90
26	14	381	G	C8-N9-C4	5.26	108.50	106.40
26	14	769	G	N1-C6-O6	5.26	123.05	119.90
26	14	1271	G	O5'-P-OP2	-5.26	100.97	105.70
26	14	1566	A	C4-C5-C6	-5.26	114.37	117.00
26	1H	1368	G	C6-N1-C2	-5.25	121.95	125.10
26	14	663	G	OP1-P-OP2	5.25	127.48	119.60
26	14	1396	U	C4-C5-C6	5.25	122.85	119.70
1	13	115	G	P-O3'-C3'	5.25	126.00	119.70
26	1H	1254	A	C8-N9-C4	5.25	107.90	105.80
26	1H	1933	G	N7-C8-N9	5.25	115.73	113.10
26	1H	2069	G	OP2-P-O3'	5.25	116.76	105.20
26	1H	2598	A	C4-C5-C6	-5.25	114.37	117.00
26	1H	2782	G	C4-N9-C1'	5.25	133.33	126.50
23	2L	21	U	C2-N1-C1'	5.25	124.00	117.70
26	1H	579	G	C2-N3-C4	5.25	114.53	111.90
26	1H	1265	A	C5'-C4'-C3'	-5.25	107.60	116.00
26	1H	2061	G	C2-N3-C4	5.25	114.53	111.90
26	1H	2550	G	O5'-P-OP1	5.25	117.00	110.70
26	14	192	C	N1-C2-O2	-5.25	115.75	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1605	C	C2-N3-C4	-5.25	117.27	119.90
26	14	1709	U	C6-N1-C2	-5.25	117.85	121.00
26	14	2253	G	N1-C6-O6	5.25	123.05	119.90
26	14	2314	C	N1-C2-O2	5.25	122.05	118.90
26	1H	318	C	O5'-P-OP1	-5.25	100.97	105.70
26	1H	735	A	C5-N7-C8	5.25	106.53	103.90
26	1H	2427	C	C2-N1-C1'	-5.25	113.03	118.80
1	1G	916	G	C2-N3-C4	5.25	114.53	111.90
26	14	93	C	C2-N1-C1'	5.25	124.58	118.80
26	14	870	A	OP1-P-O3'	5.25	116.75	105.20
26	14	2056	G	N1-C6-O6	5.25	123.05	119.90
26	1H	729	G	C4-N9-C1'	5.25	133.32	126.50
26	1H	1125	G	C5-C6-N1	5.25	114.12	111.50
26	1H	2007	C	O5'-P-OP1	5.25	117.00	110.70
26	1H	2373	G	C4-C5-C6	5.25	121.95	118.80
26	14	965	C	O5'-P-OP1	-5.25	100.98	105.70
1	13	963	G	C8-N9-C4	5.25	108.50	106.40
1	13	1395	C	C5-C6-N1	5.25	123.62	121.00
26	1H	33	U	OP2-P-O3'	-5.25	93.66	105.20
26	1H	1225	C	C5-C4-N4	5.25	123.87	120.20
26	1H	1598	C	C6-N1-C2	-5.25	118.20	120.30
27	16	103	U	OP2-P-O3'	5.25	116.74	105.20
26	14	204	A	N1-C2-N3	5.25	131.92	129.30
26	14	690	G	OP1-P-O3'	5.25	116.74	105.20
26	14	1828	G	C4-C5-C6	5.25	121.95	118.80
26	14	2820	A	C2-N3-C4	-5.25	107.98	110.60
26	1H	672	C	OP2-P-O3'	5.25	116.74	105.20
26	1H	1314	C	N1-C2-O2	5.25	122.05	118.90
26	1H	1543	A	C5-C6-N1	-5.25	115.08	117.70
26	1H	2290	G	O5'-P-OP2	5.25	116.99	110.70
26	1H	2434	A	N1-C6-N6	-5.25	115.45	118.60
26	1H	2546	U	N3-C2-O2	5.25	125.87	122.20
26	1H	2878	U	C2-N1-C1'	5.25	123.99	117.70
27	16	87	G	N9-C4-C5	-5.25	103.30	105.40
1	1G	1502	A	N1-C2-N3	5.25	131.92	129.30
26	14	1950	G	C4-N9-C1'	5.25	133.32	126.50
26	14	1977	A	C8-N9-C4	5.25	107.90	105.80
26	1H	795	C	N3-C4-C5	5.24	124.00	121.90
26	1H	1305	C	C2-N3-C4	-5.24	117.28	119.90
26	1H	1391	U	C2-N3-C4	5.24	130.15	127.00
26	1H	1438	U	C6-N1-C2	-5.24	117.85	121.00
26	1H	2451	A	C6-C5-N7	5.24	135.97	132.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2672	G	C4-N9-C1'	5.24	133.32	126.50
26	1H	2822	G	C5-C6-O6	-5.24	125.45	128.60
26	14	598	G	OP1-P-OP2	5.24	127.47	119.60
26	14	1899	G	C6-C5-N7	-5.24	127.25	130.40
26	14	2522	U	N3-C2-O2	-5.24	118.53	122.20
26	1H	121	G	C8-N9-C4	5.24	108.50	106.40
26	1H	663	G	C5-C6-N1	-5.24	108.88	111.50
26	1H	787	U	N3-C4-C5	5.24	117.75	114.60
26	1H	1224	G	C4-N9-C1'	-5.24	119.69	126.50
26	14	458	G	O5'-P-OP2	-5.24	100.98	105.70
26	14	1496	A	O4'-C1'-N9	5.24	112.39	108.20
1	13	418	C	C5-C6-N1	5.24	123.62	121.00
1	13	813	U	O5'-P-OP2	-5.24	100.98	105.70
26	1H	484	C	OP1-P-O3'	5.24	116.73	105.20
26	1H	758	C	C2-N3-C4	-5.24	117.28	119.90
26	1H	775	G	C6-C5-N7	-5.24	127.26	130.40
26	1H	1268	A	N7-C8-N9	-5.24	111.18	113.80
26	1H	1424	G	C8-N9-C1'	-5.24	120.19	127.00
26	1H	2372	G	C5-C6-O6	-5.24	125.45	128.60
30	31	140	LEU	CA-CB-CG	5.24	127.35	115.30
1	1G	1348	U	N3-C2-O2	-5.24	118.53	122.20
26	14	1918	A	N9-C4-C5	-5.24	103.70	105.80
26	14	2330	G	C6-C5-N7	-5.24	127.26	130.40
1	13	359	U	C5-C4-O4	5.24	129.04	125.90
1	13	759	A	OP2-P-O3'	5.24	116.73	105.20
26	1H	1941	C	C6-N1-C2	-5.24	118.20	120.30
26	1H	2690	C	O4'-C1'-N1	5.24	112.39	108.20
29	21	65	GLY	N-CA-C	-5.24	100.00	113.10
26	14	754	C	N3-C4-C5	-5.24	119.81	121.90
26	14	1558	A	P-O3'-C3'	5.24	125.99	119.70
26	14	1703	G	C8-N9-C4	5.24	108.50	106.40
26	14	2224	G	N3-C4-N9	5.24	129.14	126.00
26	1H	2550	G	C4-N9-C1'	5.24	133.31	126.50
1	1G	963	G	C8-N9-C1'	-5.24	120.19	127.00
26	1H	87	C	C5-C4-N4	-5.24	116.53	120.20
26	1H	1191	G	N1-C6-O6	-5.24	116.76	119.90
26	1H	1684	C	O5'-P-OP2	-5.24	100.99	105.70
26	1H	1955	U	O5'-P-OP2	-5.24	100.99	105.70
26	1H	2055	C	OP2-P-O3'	5.24	116.72	105.20
26	1H	2235	G	O5'-P-OP1	5.24	116.98	110.70
26	1H	2495	G	C5-C6-N1	-5.24	108.88	111.50
26	1H	2525	G	C5-C6-O6	-5.24	125.46	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
27	16	17	C	N1-C2-O2	5.24	122.04	118.90
26	14	2681	C	C4-C5-C6	5.24	120.02	117.40
27	1J	56	G	N3-C4-C5	-5.24	125.98	128.60
14	5I	44	LEU	CB-CG-CD1	5.23	119.90	111.00
26	1H	837	C	C5-C4-N4	-5.23	116.54	120.20
26	1H	2737	G	C5-C6-O6	-5.23	125.46	128.60
1	13	1469	G	C5-C6-O6	-5.23	125.46	128.60
26	1H	199	A	N1-C2-N3	-5.23	126.68	129.30
26	1H	835	A	N3-C4-C5	-5.23	123.14	126.80
26	1H	1166	C	C5-C6-N1	5.23	123.62	121.00
26	1H	1564	C	N3-C2-O2	-5.23	118.24	121.90
26	1H	2307	G	C6-C5-N7	-5.23	127.26	130.40
26	1H	2699	C	N3-C4-C5	5.23	123.99	121.90
1	1G	1297	C	OP2-P-O3'	5.23	116.71	105.20
26	14	752	A	C8-N9-C4	-5.23	103.71	105.80
26	14	803	U	O5'-P-OP1	5.23	116.98	110.70
26	14	1279	G	C5-C6-O6	5.23	131.74	128.60
26	14	2426	A	N7-C8-N9	5.23	116.42	113.80
26	14	2599	G	N1-C6-O6	-5.23	116.76	119.90
26	1H	760	G	OP1-P-OP2	-5.23	111.75	119.60
26	1H	827	U	C2-N1-C1'	-5.23	111.42	117.70
26	1H	1334	G	N3-C4-N9	5.23	129.14	126.00
26	1H	1689	A	N1-C6-N6	-5.23	115.46	118.60
26	1H	2626	C	N3-C4-C5	5.23	123.99	121.90
26	1H	674	G	N9-C4-C5	-5.23	103.31	105.40
26	1H	799	G	N3-C4-N9	-5.23	122.86	126.00
1	1G	525	C	C5-C6-N1	5.23	123.61	121.00
26	14	1325	G	C2-N3-C4	5.23	114.51	111.90
1	13	1488	G	C5-C6-O6	-5.23	125.46	128.60
26	1H	2035	G	O4'-C1'-N9	5.23	112.38	108.20
26	1H	2072	G	C4-C5-N7	5.23	112.89	110.80
26	14	688	U	C5-C6-N1	-5.23	120.09	122.70
26	14	863	A	C8-N9-C4	-5.23	103.71	105.80
26	14	2491	U	N1-C2-O2	-5.23	119.14	122.80
26	1H	748	G	N3-C4-C5	5.23	131.21	128.60
26	1H	974	G	O5'-P-OP2	-5.23	101.00	105.70
26	1H	1301	A	C6-C5-N7	-5.23	128.64	132.30
26	1H	1653	G	O5'-P-OP2	-5.23	101.00	105.70
26	1H	2597	G	P-O3'-C3'	5.23	125.97	119.70
26	14	396	G	OP1-P-O3'	5.23	116.70	105.20
1	13	244	U	C4-C5-C6	5.22	122.83	119.70
1	13	1178	G	C8-N9-C4	-5.22	104.31	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	980	A	N3-C4-C5	-5.22	123.14	126.80
26	1H	1238	G	N9-C1'-C2'	-5.22	106.25	112.00
26	1H	1915	U	N1-C2-O2	5.22	126.46	122.80
27	16	75	G	C8-N9-C4	5.22	108.49	106.40
26	14	2004	G	O5'-P-OP1	-5.22	101.00	105.70
26	14	2585	U	O4'-C1'-N1	5.22	112.38	108.20
28	19	244	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	13	1111	A	N1-C6-N6	-5.22	115.47	118.60
1	13	1209	C	N3-C4-N4	5.22	121.66	118.00
23	2K	62	C	C6-N1-C2	-5.22	118.21	120.30
26	1H	97	C	OP1-P-OP2	5.22	127.44	119.60
26	1H	188	G	N1-C2-N2	-5.22	111.50	116.20
26	1H	569	U	N1-C2-O2	-5.22	119.14	122.80
26	1H	792	G	N3-C4-N9	5.22	129.13	126.00
26	1H	1339	G	C5-C6-O6	-5.22	125.47	128.60
26	1H	2323	G	C4-C5-N7	5.22	112.89	110.80
26	1H	2851	A	OP2-P-O3'	5.22	116.69	105.20
26	1H	2870	C	OP2-P-O3'	5.22	116.69	105.20
26	14	315	G	O5'-P-OP2	-5.22	101.00	105.70
26	14	1216	G	C6-C5-N7	-5.22	127.27	130.40
26	14	1351	C	C4-C5-C6	5.22	120.01	117.40
26	14	2032	G	N1-C6-O6	5.22	123.03	119.90
26	1H	455	C	C4-C5-C6	-5.22	114.79	117.40
1	13	181	G	N3-C4-N9	5.22	129.13	126.00
1	13	889	A	C5-C6-N1	-5.22	115.09	117.70
26	1H	667	U	OP2-P-O3'	5.22	116.68	105.20
26	1H	827	U	C2-N3-C4	-5.22	123.87	127.00
26	1H	833	U	C5-C6-N1	-5.22	120.09	122.70
26	1H	1423	G	OP1-P-OP2	5.22	127.43	119.60
1	1G	798	G	N1-C6-O6	5.22	123.03	119.90
26	14	561	G	C4-C5-N7	-5.22	108.71	110.80
26	14	2030	A	C5-C6-N6	-5.22	119.52	123.70
26	14	2511	U	OP1-P-O3'	5.22	116.68	105.20
26	14	2617	C	N3-C2-O2	5.22	125.55	121.90
26	1H	1476	C	O5'-P-OP1	-5.22	101.00	105.70
26	14	2063	C	O5'-P-OP1	5.22	116.96	110.70
26	14	2724	C	OP2-P-O3'	5.22	116.68	105.20
26	1H	322	A	O5'-P-OP1	-5.22	101.00	105.70
26	1H	1356	G	O5'-P-OP1	-5.22	101.01	105.70
26	1H	1386	C	N1-C2-O2	-5.22	115.77	118.90
26	1H	1437	C	C5-C6-N1	5.22	123.61	121.00
27	16	7	G	C2-N3-C4	-5.22	109.29	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	32	135	LEU	CA-CB-CG	5.22	127.30	115.30
26	14	1290	C	OP1-P-OP2	5.22	127.42	119.60
43	A5	92	ARG	NE-CZ-NH2	-5.22	117.69	120.30
25	4K	18	G	N3-C4-N9	-5.21	122.87	126.00
26	1H	2210	G	N7-C8-N9	5.21	115.71	113.10
26	1H	2347	C	C5-C6-N1	5.21	123.61	121.00
26	1H	2702	U	C5'-C4'-O4'	5.21	115.36	109.10
1	1G	1190	G	N9-C4-C5	5.21	107.49	105.40
26	14	186	G	O5'-P-OP1	-5.21	101.01	105.70
26	14	1022	G	C8-N9-C1'	5.21	133.78	127.00
1	13	703	G	C4-N9-C1'	5.21	133.28	126.50
26	1H	203	C	C6-N1-C2	5.21	122.39	120.30
26	1H	751	A	N7-C8-N9	-5.21	111.19	113.80
26	1H	828	U	C4-C5-C6	5.21	122.83	119.70
26	1H	1258	C	OP2-P-O3'	5.21	116.67	105.20
26	1H	1798	U	N3-C4-C5	5.21	117.73	114.60
1	1G	772	U	O5'-P-OP2	-5.21	101.01	105.70
22	1K	74	C	C5-C6-N1	5.21	123.61	121.00
26	1H	1011	G	N3-C4-C5	-5.21	125.99	128.60
26	1H	1250	G	C5-C6-O6	-5.21	125.47	128.60
28	11	239	ARG	NE-CZ-NH1	-5.21	117.69	120.30
26	14	828	U	N3-C4-O4	-5.21	115.75	119.40
26	14	1265	A	N1-C6-N6	5.21	121.73	118.60
26	14	1678	G	N1-C6-O6	5.21	123.03	119.90
26	14	1914	C	C2-N1-C1'	5.21	124.53	118.80
26	14	2713	A	C6-C5-N7	-5.21	128.65	132.30
26	1H	75	G	C5-C6-N1	5.21	114.11	111.50
26	1H	632	A	OP1-P-OP2	-5.21	111.78	119.60
26	1H	889	C	C2-N1-C1'	5.21	124.53	118.80
1	13	1113	C	N3-C2-O2	-5.21	118.25	121.90
26	1H	1775	U	C4-C5-C6	5.21	122.83	119.70
26	1H	2215	G	C8-N9-C4	5.21	108.48	106.40
27	16	95	U	C5-C4-O4	5.21	129.03	125.90
26	14	470	A	OP1-P-OP2	5.21	127.41	119.60
26	14	479	A	N9-C4-C5	5.21	107.88	105.80
26	14	980	A	N1-C6-N6	5.21	121.72	118.60
26	14	2032	G	C6-C5-N7	-5.21	127.28	130.40
26	1H	1824	G	C5-C6-O6	-5.21	125.48	128.60
26	1H	1942	C	C4-C5-C6	-5.21	114.80	117.40
1	1G	1529	G	C4-N9-C1'	5.21	133.27	126.50
26	14	669	G	N3-C4-C5	-5.21	126.00	128.60
26	14	1023	U	C5-C4-O4	5.21	129.02	125.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1798	U	C2-N3-C4	-5.21	123.88	127.00
1	13	1266	G	N3-C4-N9	-5.21	122.88	126.00
26	1H	119	A	N9-C4-C5	5.21	107.88	105.80
26	1H	738	G	N3-C4-N9	5.21	129.12	126.00
26	1H	910	A	C5-C6-N1	-5.21	115.10	117.70
26	1H	2264	C	N1-C2-O2	5.21	122.02	118.90
27	16	14	U	O5'-P-OP2	-5.21	101.02	105.70
27	16	17	C	N3-C2-O2	-5.21	118.26	121.90
25	4K	20	U	C2-N1-C1'	5.20	123.94	117.70
26	1H	219	G	N9-C4-C5	5.20	107.48	105.40
26	1H	1950	G	N3-C2-N2	5.20	123.54	119.90
26	1H	2542	A	C8-N9-C4	5.20	107.88	105.80
26	1H	2712	U	N3-C4-O4	5.20	123.04	119.40
26	14	1769	G	N3-C4-N9	5.20	129.12	126.00
26	14	2335	A	O5'-P-OP1	-5.20	101.02	105.70
27	1J	103	U	N3-C4-O4	-5.20	115.76	119.40
26	1H	1204	A	C8-N9-C4	-5.20	103.72	105.80
26	1H	1560	G	OP1-P-O3'	5.20	116.64	105.20
26	1H	2238	G	O5'-P-OP1	-5.20	101.02	105.70
26	14	2405	G	N1-C6-O6	5.20	123.02	119.90
1	13	123	C	C5-C6-N1	-5.20	118.40	121.00
1	13	412	A	P-O3'-C3'	5.20	125.94	119.70
23	2K	6	G	N9-C4-C5	-5.20	103.32	105.40
26	1H	660	G	C2-N3-C4	-5.20	109.30	111.90
26	1H	834	C	C5-C4-N4	-5.20	116.56	120.20
26	1H	971	C	OP2-P-O3'	5.20	116.64	105.20
26	1H	1559	G	C2-N3-C4	-5.20	109.30	111.90
26	1H	2054	A	OP1-P-O3'	-5.20	93.76	105.20
26	1H	2210	G	C8-N9-C4	-5.20	104.32	106.40
38	98	75	LEU	CA-CB-CG	5.20	127.26	115.30
1	1G	498	A	C8-N9-C4	5.20	107.88	105.80
26	14	1279	G	N3-C4-C5	-5.20	126.00	128.60
26	14	1991	U	C5-C4-O4	5.20	129.02	125.90
26	14	2356	C	C5-C6-N1	-5.20	118.40	121.00
26	14	2593	U	O5'-P-OP1	5.20	116.94	110.70
26	1H	404	C	N3-C2-O2	5.20	125.54	121.90
26	1H	763	G	N1-C2-N3	5.20	127.02	123.90
26	1H	1193	G	C8-N9-C4	5.20	108.48	106.40
26	1H	1307	A	OP1-P-OP2	5.20	127.40	119.60
26	1H	2606	C	OP1-P-O3'	5.20	116.64	105.20
1	1G	244	U	C5-C4-O4	-5.20	122.78	125.90
26	14	863	A	OP2-P-O3'	5.20	116.64	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	974(A)	C	N1-C2-N3	5.20	122.84	119.20
26	14	1573	G	O5'-P-OP1	5.20	116.94	110.70
1	13	557	G	N3-C4-N9	5.20	129.12	126.00
1	13	1417	G	N3-C4-C5	-5.20	126.00	128.60
26	1H	512	G	N1-C6-O6	-5.20	116.78	119.90
26	1H	609	A	N7-C8-N9	-5.20	111.20	113.80
26	1H	931	G	C6-N1-C2	-5.20	121.98	125.10
26	1H	1310	G	N1-C6-O6	5.20	123.02	119.90
26	1H	2397	G	N1-C6-O6	5.20	123.02	119.90
26	14	1284	A	N1-C6-N6	5.20	121.72	118.60
26	14	2540	C	O5'-P-OP2	-5.20	101.02	105.70
26	1H	28	A	C2-N3-C4	5.20	113.20	110.60
26	1H	1021	A	C6-N1-C2	5.20	121.72	118.60
26	1H	1562	A	N1-C6-N6	5.20	121.72	118.60
50	L8	54	VAL	CB-CA-C	-5.20	101.53	111.40
26	14	2298	A	O5'-P-OP2	-5.20	101.02	105.70
27	1J	116	G	N3-C4-C5	5.20	131.20	128.60
26	1H	105	C	C6-N1-C2	-5.19	118.22	120.30
26	1H	2449	U	OP1-P-OP2	5.19	127.39	119.60
26	14	1969	A	OP1-P-OP2	-5.19	111.81	119.60
26	14	2622	C	C5-C6-N1	-5.19	118.40	121.00
26	1H	1186	G	N3-C2-N2	-5.19	116.27	119.90
26	1H	1210	A	C8-N9-C4	-5.19	103.72	105.80
26	1H	2396	G	C4-N9-C1'	-5.19	119.75	126.50
26	1H	2688	U	C6-N1-C2	-5.19	117.89	121.00
1	1G	1339	A	O4'-C1'-N9	5.19	112.35	108.20
26	14	760	G	O5'-P-OP2	5.19	116.93	110.70
26	14	933	A	C4-C5-N7	5.19	113.30	110.70
26	14	1414	G	C8-N9-C1'	-5.19	120.25	127.00
26	14	1664	A	N1-C2-N3	5.19	131.90	129.30
26	14	2607	G	C8-N9-C4	5.19	108.48	106.40
1	13	115	G	C2-N3-C4	5.19	114.50	111.90
1	13	221	C	C6-N1-C2	-5.19	118.22	120.30
1	13	1469	G	N7-C8-N9	5.19	115.69	113.10
22	1K	76	A	O4'-C1'-N9	5.19	112.35	108.20
26	1H	670	A	N1-C6-N6	5.19	121.71	118.60
26	1H	736	C	N3-C2-O2	5.19	125.53	121.90
26	1H	849	A	C5-C6-N6	-5.19	119.55	123.70
26	1H	1367	A	C8-N9-C4	5.19	107.88	105.80
26	1H	1817	G	N3-C2-N2	5.19	123.53	119.90
26	1H	2329	G	C8-N9-C4	5.19	108.48	106.40
26	1H	2645	G	N3-C4-C5	5.19	131.19	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
27	16	91	C	C5-C6-N1	5.19	123.59	121.00
26	14	2456	C	N3-C4-C5	-5.19	119.82	121.90
1	13	1327	C	C6-N1-C2	5.19	122.38	120.30
26	1H	681	G	O5'-P-OP2	5.19	116.93	110.70
26	1H	2558	C	OP2-P-O3'	5.19	116.61	105.20
26	14	1633	G	C8-N9-C4	-5.19	104.32	106.40
1	13	538	G	C6-C5-N7	-5.19	127.29	130.40
26	1H	782	A	N1-C2-N3	5.19	131.89	129.30
26	1H	1311	G	N1-C6-O6	5.19	123.01	119.90
26	1H	1817	G	N3-C4-N9	5.19	129.11	126.00
26	1H	2778	A	C2-N3-C4	-5.19	108.01	110.60
26	14	666	G	C2-N3-C4	-5.19	109.31	111.90
26	14	773	U	C4-C5-C6	5.19	122.81	119.70
26	14	1155	A	OP1-P-O3'	5.19	116.61	105.20
26	14	1528	A	C8-N9-C4	-5.19	103.72	105.80
26	14	1961	C	N1-C2-O2	-5.19	115.79	118.90
26	14	1971	A	OP1-P-O3'	5.19	116.61	105.20
26	1H	431	U	OP1-P-OP2	-5.19	111.82	119.60
26	1H	1348	G	N1-C6-O6	5.19	123.01	119.90
26	1H	2277	G	C5-C6-O6	5.19	131.71	128.60
27	16	49	C	N1-C2-O2	-5.19	115.79	118.90
1	1G	413	G	N1-C6-O6	-5.19	116.79	119.90
1	13	185	A	N7-C8-N9	5.18	116.39	113.80
1	13	576	G	N3-C4-C5	-5.18	126.01	128.60
26	1H	256	A	C8-N9-C4	5.18	107.87	105.80
26	1H	1189	A	C5-N7-C8	-5.18	101.31	103.90
26	14	451	C	O4'-C1'-N1	5.18	112.35	108.20
26	14	583	G	N7-C8-N9	5.18	115.69	113.10
26	14	1319	G	O5'-P-OP2	-5.18	101.03	105.70
26	14	2548	G	C8-N9-C4	-5.18	104.33	106.40
26	1H	616	A	C6-C5-N7	-5.18	128.67	132.30
26	1H	1272	A	C8-N9-C4	5.18	107.87	105.80
26	1H	1471	A	N7-C8-N9	5.18	116.39	113.80
26	1H	1766	U	C5-C4-O4	-5.18	122.79	125.90
26	1H	2599	G	OP2-P-O3'	5.18	116.60	105.20
28	11	95	LEU	CA-CB-CG	5.18	127.22	115.30
46	H8	161	VAL	CB-CA-C	-5.18	101.55	111.40
26	14	774	A	C4-N9-C1'	-5.18	116.97	126.30
26	14	2333	A	C5-N7-C8	5.18	106.49	103.90
26	14	2572	A	N9-C4-C5	5.18	107.87	105.80
1	13	219	C	C6-N1-C2	-5.18	118.23	120.30
26	1H	1726	G	C4-N9-C1'	-5.18	119.77	126.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1G	301	G	O5'-P-OP2	-5.18	101.04	105.70
26	14	1619	G	C4-C5-N7	5.18	112.87	110.80
1	13	266	G	C4-C5-N7	5.18	112.87	110.80
1	13	966	G	N3-C4-N9	5.18	129.11	126.00
26	1H	389	G	N9-C4-C5	-5.18	103.33	105.40
26	1H	417	C	N3-C2-O2	5.18	125.53	121.90
26	1H	568	U	O5'-P-OP2	-5.18	101.04	105.70
26	1H	628	G	N7-C8-N9	-5.18	110.51	113.10
26	1H	683	C	C2-N3-C4	-5.18	117.31	119.90
26	1H	2468	G	N1-C2-N2	-5.18	111.54	116.20
26	14	2376	A	C8-N9-C4	5.18	107.87	105.80
26	14	2601	C	N3-C2-O2	-5.18	118.27	121.90
1	13	1518	A	N7-C8-N9	-5.18	111.21	113.80
26	1H	1241	A	C5-C6-N1	-5.18	115.11	117.70
26	1H	1299	G	N3-C4-N9	5.18	129.11	126.00
26	14	783	A	O5'-P-OP2	-5.18	101.04	105.70
26	14	1208	C	O5'-P-OP1	-5.18	101.04	105.70
27	1J	103	U	C2-N3-C4	-5.18	123.89	127.00
1	13	891	U	OP2-P-O3'	5.18	116.59	105.20
26	1H	580	C	C6-N1-C2	-5.18	118.23	120.30
26	1H	972	G	C5-C6-O6	5.18	131.71	128.60
26	1H	1889	A	C5-N7-C8	-5.18	101.31	103.90
26	14	600	G	C8-N9-C4	5.18	108.47	106.40
2	1E	196	LEU	CA-CB-CG	5.17	127.20	115.30
26	1H	477	A	OP1-P-OP2	-5.17	111.84	119.60
26	1H	2586	C	N1-C2-O2	-5.17	115.80	118.90
26	1H	2751	G	C5-N7-C8	-5.17	101.71	104.30
1	1G	35	G	C8-N9-C4	-5.17	104.33	106.40
1	1G	413	G	C5-C6-O6	5.17	131.71	128.60
1	1G	583	A	C8-N9-C4	5.17	107.87	105.80
1	1G	1200	C	C6-N1-C1'	-5.17	114.59	120.80
1	13	413	G	C4-C5-N7	-5.17	108.73	110.80
23	2K	76	C	C5-C4-N4	-5.17	116.58	120.20
26	1H	1022	G	C6-N1-C2	-5.17	122.00	125.10
50	L8	49	LYS	CD-CE-NZ	-5.17	99.80	111.70
1	13	1290	G	C6-C5-N7	-5.17	127.30	130.40
26	1H	207	A	C2-N3-C4	-5.17	108.01	110.60
26	1H	835	A	OP2-P-O3'	5.17	116.58	105.20
26	1H	1249	U	N3-C4-C5	5.17	117.70	114.60
26	1H	2380	C	N3-C4-C5	5.17	123.97	121.90
1	13	888	G	C2-N3-C4	-5.17	109.31	111.90
26	1H	197	A	C5-N7-C8	-5.17	101.31	103.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	1138	G	C6-C5-N7	-5.17	127.30	130.40
26	14	2039	C	N1-C2-O2	5.17	122.00	118.90
1	13	963	G	N3-C4-C5	-5.17	126.02	128.60
26	1H	613	U	C2-N3-C4	-5.17	123.90	127.00
26	1H	2396	G	C8-N9-C1'	5.17	133.72	127.00
26	14	2319	G	N3-C4-C5	-5.17	126.02	128.60
26	14	2430	A	O5'-P-OP1	5.17	116.90	110.70
26	1H	863	A	N1-C6-N6	-5.17	115.50	118.60
26	1H	1502	C	C5-C6-N1	5.17	123.58	121.00
27	16	90	C	N3-C4-C5	5.17	123.97	121.90
26	14	388	G	N3-C4-C5	5.17	131.18	128.60
26	14	1431	U	N3-C4-C5	5.17	117.70	114.60
26	14	1702	G	C5-C6-O6	-5.17	125.50	128.60
26	14	1772	G	N9-C1'-C2'	-5.17	106.32	112.00
26	14	2013	A	C8-N9-C4	5.17	107.87	105.80
1	13	1455	G	C8-N9-C4	5.17	108.47	106.40
26	1H	2636	U	C6-N1-C2	5.17	124.10	121.00
26	1H	2757	A	N1-C6-N6	5.17	121.70	118.60
26	14	18	C	O5'-P-OP1	-5.17	101.05	105.70
26	14	1527	G	N3-C4-C5	5.17	131.18	128.60
1	13	328	C	N3-C2-O2	-5.16	118.28	121.90
26	1H	474	G	N9-C4-C5	5.16	107.47	105.40
26	1H	948	G	N3-C4-C5	5.16	131.18	128.60
26	1H	1355	G	C5-C6-N1	5.16	114.08	111.50
26	1H	1611	C	C5-C4-N4	-5.16	116.58	120.20
26	1H	1776	G	C4-C5-N7	5.16	112.87	110.80
26	1H	2246	G	N7-C8-N9	-5.16	110.52	113.10
26	14	1797	C	C5-C4-N4	-5.16	116.58	120.20
26	14	2068	U	C5-C6-N1	-5.16	120.12	122.70
26	14	2763	G	N3-C2-N2	5.16	123.51	119.90
26	1H	841	A	C6-C5-N7	-5.16	128.69	132.30
26	1H	1799	G	N1-C2-N2	-5.16	111.55	116.20
26	14	726	G	C5-C6-O6	5.16	131.70	128.60
26	14	982	C	N3-C4-C5	-5.16	119.83	121.90
26	14	2086	U	O5'-P-OP2	-5.16	101.05	105.70
1	13	908	A	C2-N3-C4	-5.16	108.02	110.60
26	1H	1656	C	OP2-P-O3'	5.16	116.55	105.20
31	41	44	GLY	N-CA-C	-5.16	100.20	113.10
26	14	194	G	O5'-P-OP2	5.16	116.89	110.70
26	14	955	C	O5'-P-OP2	-5.16	101.06	105.70
26	14	1253	A	C4-C5-C6	5.16	119.58	117.00
26	14	1311	G	C5-C6-N1	-5.16	108.92	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	928	G	O5'-P-OP1	-5.16	101.06	105.70
26	1H	865	C	OP1-P-OP2	-5.16	111.86	119.60
26	1H	1364	G	OP2-P-O3'	5.16	116.55	105.20
26	1H	1630(A)	C	O5'-P-OP1	-5.16	101.06	105.70
26	1H	2029	G	O5'-P-OP1	-5.16	101.06	105.70
26	1H	2429	G	OP1-P-OP2	-5.16	111.86	119.60
1	1G	552	U	OP2-P-O3'	5.16	116.55	105.20
26	14	2627	G	C8-N9-C1'	-5.16	120.30	127.00
26	1H	825	C	N3-C2-O2	5.16	125.51	121.90
26	1H	1283	G	N3-C4-N9	5.16	129.09	126.00
26	1H	1695	G	N1-C2-N2	-5.16	111.56	116.20
26	1H	2575	C	C5-C6-N1	-5.16	118.42	121.00
26	1H	2697	G	C6-C5-N7	-5.16	127.31	130.40
26	14	2057	A	C2-N3-C4	-5.16	108.02	110.60
26	1H	598	G	N3-C4-C5	-5.16	126.02	128.60
26	1H	1649	G	N3-C2-N2	5.16	123.51	119.90
26	1H	1728	G	N3-C2-N2	5.16	123.51	119.90
26	1H	2275	C	C5'-C4'-O4'	-5.16	102.91	109.10
26	14	2377	A	C4-N9-C1'	-5.16	117.02	126.30
26	1H	1369	G	O5'-P-OP2	-5.15	101.06	105.70
26	1H	1698	A	N9-C1'-C2'	5.15	120.70	114.00
26	1H	1930	G	OP2-P-O3'	5.15	116.54	105.20
26	14	141	A	O4'-C1'-N9	5.15	112.32	108.20
26	1H	236	C	N3-C4-C5	-5.15	119.84	121.90
26	1H	772	C	N3-C4-N4	5.15	121.61	118.00
26	1H	805	G	C6-C5-N7	-5.15	127.31	130.40
26	1H	1214	A	OP2-P-O3'	5.15	116.53	105.20
26	1H	1271	G	N9-C4-C5	-5.15	103.34	105.40
26	1H	2028	U	C4-C5-C6	5.15	122.79	119.70
41	C8	50	ARG	NE-CZ-NH2	-5.15	117.72	120.30
26	14	929	G	C4-N9-C1'	5.15	133.20	126.50
26	14	1277	G	N3-C4-C5	5.15	131.18	128.60
26	14	2043	C	C6-N1-C2	-5.15	118.24	120.30
26	1H	193	U	C2-N3-C4	-5.15	123.91	127.00
26	1H	321	G	C5-C6-O6	-5.15	125.51	128.60
26	1H	411	G	C4-C5-N7	-5.15	108.74	110.80
26	1H	571	A	C8-N9-C4	5.15	107.86	105.80
26	1H	911	A	C5-N7-C8	-5.15	101.33	103.90
26	1H	1257	C	N3-C4-N4	5.15	121.61	118.00
1	1G	509	A	N9-C1'-C2'	-5.15	106.33	112.00
26	14	1121	C	C2-N3-C4	-5.15	117.33	119.90
26	14	1971	A	C8-N9-C4	5.15	107.86	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1293	C	C2-N1-C1'	5.15	124.46	118.80
26	1H	1298	C	C2-N3-C4	5.15	122.47	119.90
26	14	982	C	C5-C6-N1	5.15	123.58	121.00
26	14	1936	A	O4'-C1'-N9	5.15	112.32	108.20
1	13	1292	U	C5-C6-N1	-5.15	120.13	122.70
26	1H	517	C	N1-C2-O2	-5.15	115.81	118.90
26	1H	770	G	C4-C5-N7	5.15	112.86	110.80
26	1H	937	U	C6-N1-C2	5.15	124.09	121.00
26	1H	1692	U	C2-N1-C1'	5.15	123.88	117.70
26	1H	2368	C	O5'-P-OP2	5.15	116.88	110.70
27	16	85	G	C5-C6-O6	-5.15	125.51	128.60
1	1G	397	A	C2-N3-C4	5.15	113.17	110.60
26	14	27	G	OP1-P-O3'	5.15	116.52	105.20
26	14	1030	G	C6-C5-N7	-5.15	127.31	130.40
26	14	1350	C	O5'-P-OP2	5.15	116.88	110.70
1	13	795	C	O5'-P-OP2	-5.15	101.07	105.70
26	1H	508	G	N9-C1'-C2'	5.15	120.69	114.00
1	1G	128	G	N3-C2-N2	-5.15	116.30	119.90
1	1G	176	C	C6-N1-C2	-5.15	118.24	120.30
24	3L	37	A	C2-N3-C4	5.15	113.17	110.60
26	14	757	U	C6-N1-C2	-5.15	117.91	121.00
26	14	1407	C	O5'-P-OP2	5.15	116.88	110.70
26	14	1652	A	O5'-P-OP2	-5.15	101.07	105.70
28	19	43	ARG	NE-CZ-NH2	-5.15	117.73	120.30
1	13	266	G	O4'-C1'-N9	-5.14	104.08	108.20
26	1H	73	A	N3-C4-C5	-5.14	123.20	126.80
26	1H	678	C	C5-C6-N1	-5.14	118.43	121.00
26	1H	838	C	N3-C4-N4	5.14	121.60	118.00
26	1H	1623	G	C4-C5-N7	-5.14	108.74	110.80
26	1H	1800	C	OP1-P-O3'	5.14	116.52	105.20
26	14	932	G	N3-C4-N9	-5.14	122.91	126.00
26	14	1433	U	O5'-P-OP2	-5.14	101.07	105.70
26	14	2437	U	OP1-P-OP2	5.14	127.32	119.60
1	13	124	G	O5'-P-OP1	5.14	116.87	110.70
1	13	653	A	O4'-C1'-N9	5.14	112.31	108.20
1	13	1432	G	C4-C5-C6	5.14	121.89	118.80
26	1H	140	A	OP2-P-O3'	5.14	116.51	105.20
26	1H	542	C	N3-C4-C5	-5.14	119.84	121.90
26	1H	655	A	C5-N7-C8	-5.14	101.33	103.90
26	1H	944	G	C8-N9-C4	-5.14	104.34	106.40
26	1H	972	G	C5-N7-C8	5.14	106.87	104.30
26	1H	1514	U	O5'-P-OP1	-5.14	101.07	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2252	G	N7-C8-N9	-5.14	110.53	113.10
26	1H	2503	A	OP1-P-OP2	-5.14	111.89	119.60
1	1G	345	C	P-O3'-C3'	5.14	125.87	119.70
1	1G	1128	C	P-O3'-C3'	5.14	125.87	119.70
1	1G	1237	C	C6-N1-C2	-5.14	118.24	120.30
23	2L	72	C	OP2-P-O3'	5.14	116.51	105.20
26	14	1276	A	C4-C5-N7	5.14	113.27	110.70
26	14	1332	G	C4-C5-N7	5.14	112.86	110.80
26	14	1439	A	N1-C2-N3	5.14	131.87	129.30
26	14	2265	U	N3-C4-O4	5.14	123.00	119.40
26	14	2627	G	C6-C5-N7	-5.14	127.31	130.40
1	13	1486	G	O5'-P-OP2	-5.14	101.07	105.70
1	13	1514	C	OP2-P-O3'	5.14	116.51	105.20
23	2K	9	G	N1-C6-O6	-5.14	116.82	119.90
26	1H	602	G	C6-C5-N7	-5.14	127.31	130.40
26	14	1826	G	N3-C4-N9	5.14	129.08	126.00
1	13	12	U	C5-C6-N1	5.14	125.27	122.70
1	13	1408	A	N1-C6-N6	5.14	121.68	118.60
26	1H	563	G	N7-C8-N9	-5.14	110.53	113.10
26	1H	1667	G	C4-C5-N7	5.14	112.86	110.80
26	1H	2042	A	N3-C4-N9	-5.14	123.29	127.40
26	1H	2051	A	N1-C6-N6	5.14	121.68	118.60
1	1G	266	G	N3-C4-C5	-5.14	126.03	128.60
26	14	1632	A	C4-C5-N7	5.14	113.27	110.70
1	13	862	C	N3-C4-C5	-5.14	119.84	121.90
26	1H	1142(A)	A	N7-C8-N9	5.14	116.37	113.80
26	1H	1647	G	C5-N7-C8	5.14	106.87	104.30
26	1H	2751	G	N7-C8-N9	5.14	115.67	113.10
1	13	1227	A	C5-C6-N1	-5.14	115.13	117.70
24	3K	72	C	C6-N1-C2	-5.14	118.25	120.30
26	1H	1936	A	C5-N7-C8	-5.14	101.33	103.90
26	14	246	C	C6-N1-C2	5.14	122.36	120.30
26	14	2547	U	OP1-P-O3'	-5.14	93.90	105.20
1	13	577	G	C6-C5-N7	-5.13	127.32	130.40
26	1H	1814	G	C5-C6-O6	5.13	131.68	128.60
26	1H	2246	G	C5-C6-N1	5.13	114.07	111.50
37	88	106	VAL	CB-CA-C	-5.13	101.64	111.40
26	14	1427	A	C4-C5-N7	-5.13	108.13	110.70
26	14	1599	C	OP2-P-O3'	5.13	116.50	105.20
26	14	1628	G	O5'-P-OP2	-5.13	101.08	105.70
26	1H	321	G	C4-N9-C1'	5.13	133.17	126.50
26	1H	1215	G	N3-C4-C5	-5.13	126.03	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1395	A	C2-N3-C4	-5.13	108.03	110.60
26	14	391	G	N3-C4-N9	5.13	129.08	126.00
26	14	684	G	C5-C6-N1	5.13	114.07	111.50
1	13	419	C	C2-N1-C1'	5.13	124.45	118.80
26	1H	73	A	C5-C6-N1	5.13	120.27	117.70
26	1H	680	G	N1-C6-O6	5.13	122.98	119.90
26	1H	2278	A	C6-N1-C2	-5.13	115.52	118.60
27	16	80	U	N3-C2-O2	-5.13	118.61	122.20
26	14	114	U	C6-N1-C1'	-5.13	114.02	121.20
26	14	1331	A	C6-N1-C2	-5.13	115.52	118.60
26	1H	185	U	N1-C2-N3	5.13	117.98	114.90
26	1H	1601	G	OP1-P-O3'	5.13	116.49	105.20
26	1H	1709	U	C6-N1-C2	-5.13	117.92	121.00
26	1H	2363	C	OP2-P-O3'	5.13	116.49	105.20
26	1H	2488	A	C6-C5-N7	-5.13	128.71	132.30
26	14	828	U	C4-C5-C6	5.13	122.78	119.70
26	14	1643	G	O5'-P-OP2	-5.13	101.08	105.70
26	1H	742	G	N3-C2-N2	-5.13	116.31	119.90
26	1H	1448	G	O5'-P-OP1	-5.13	101.08	105.70
26	1H	1835	G	C2-N3-C4	5.13	114.46	111.90
26	1H	2578	G	C5-N7-C8	5.13	106.86	104.30
26	14	189	G	O5'-P-OP1	-5.13	101.08	105.70
26	14	433	C	C6-N1-C2	-5.13	118.25	120.30
26	14	1308	A	N1-C2-N3	5.13	131.86	129.30
26	14	1636	C	O5'-P-OP2	5.13	116.86	110.70
26	14	2329	G	N7-C8-N9	-5.13	110.53	113.10
1	13	809	G	N1-C6-O6	-5.13	116.82	119.90
26	1H	259	G	N9-C4-C5	-5.13	103.35	105.40
26	1H	1060	U	P-O3'-C3'	5.13	125.85	119.70
26	1H	1597	A	N1-C2-N3	5.13	131.86	129.30
26	1H	1688	U	OP2-P-O3'	5.13	116.48	105.20
26	1H	1758	G	N3-C2-N2	-5.13	116.31	119.90
26	1H	2271	G	N1-C2-N2	-5.13	111.59	116.20
26	1H	2555	U	N1-C2-N3	5.13	117.97	114.90
1	1G	579	G	N7-C8-N9	5.13	115.66	113.10
26	14	70	G	N3-C4-C5	-5.13	126.04	128.60
26	14	693	C	OP2-P-O3'	5.13	116.48	105.20
26	14	1520	U	N3-C4-O4	-5.13	115.81	119.40
26	14	2763	G	N1-C6-O6	-5.13	116.83	119.90
26	1H	165	U	N1-C2-O2	5.12	126.39	122.80
26	1H	864	G	C5-C6-O6	-5.12	125.53	128.60
26	1H	2444	G	N9-C4-C5	5.12	107.45	105.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2477	C	C6-N1-C2	-5.12	118.25	120.30
26	1H	2762	G	N1-C6-O6	5.12	122.97	119.90
27	16	8	U	O5'-P-OP1	5.12	116.85	110.70
26	14	113	G	N1-C6-O6	5.12	122.97	119.90
26	14	191	A	C6-N1-C2	5.12	121.67	118.60
26	14	392	C	N3-C4-C5	5.12	123.95	121.90
26	14	954	G	N1-C6-O6	-5.12	116.83	119.90
33	69	131	LYS	C-N-CA	5.12	143.53	122.00
23	2K	44	A	C8-N9-C4	-5.12	103.75	105.80
26	1H	988	A	P-O3'-C3'	5.12	125.85	119.70
26	1H	1759	A	OP1-P-OP2	5.12	127.28	119.60
26	1H	2073	C	OP2-P-O3'	5.12	116.47	105.20
26	1H	2428	G	P-O3'-C3'	5.12	125.85	119.70
23	2L	6	G	C8-N9-C4	5.12	108.45	106.40
23	2L	21	U	O4'-C1'-N1	5.12	112.30	108.20
26	14	694	U	N1-C2-O2	5.12	126.39	122.80
26	14	933	A	C5-C6-N6	-5.12	119.60	123.70
1	13	1335	C	C2-N1-C1'	-5.12	113.17	118.80
26	1H	701	G	C6-C5-N7	-5.12	127.33	130.40
26	1H	2295	C	C4-C5-C6	5.12	119.96	117.40
26	1H	2322	A	N9-C4-C5	5.12	107.85	105.80
26	1H	2497	A	N1-C6-N6	5.12	121.67	118.60
26	14	768	G	C4-C5-C6	5.12	121.87	118.80
26	14	782	A	C8-N9-C4	-5.12	103.75	105.80
26	14	949	C	C2-N1-C1'	-5.12	113.17	118.80
26	1H	574	C	N1-C2-N3	-5.12	115.62	119.20
1	1G	915	A	OP1-P-O3'	5.12	116.47	105.20
26	14	863	A	O5'-P-OP2	-5.12	101.09	105.70
26	14	1633	G	N7-C8-N9	5.12	115.66	113.10
1	13	756	C	OP2-P-O3'	5.12	116.46	105.20
26	1H	194	G	N7-C8-N9	-5.12	110.54	113.10
26	1H	957	A	C5-C6-N6	-5.12	119.61	123.70
26	1H	1654	A	C8-N9-C4	-5.12	103.75	105.80
26	1H	1790	C	C6-N1-C2	5.12	122.35	120.30
26	1H	2875	C	C6-N1-C2	5.12	122.35	120.30
1	1G	401	C	O5'-P-OP1	5.12	116.84	110.70
1	1G	570	G	C8-N9-C4	-5.12	104.35	106.40
26	14	278	A	P-O3'-C3'	5.12	125.84	119.70
26	14	1128	A	C5-C6-N1	5.12	120.26	117.70
26	14	1600	C	N3-C4-N4	-5.12	114.42	118.00
26	14	1968	G	C5-N7-C8	-5.12	101.74	104.30
26	14	2196	C	N1-C2-O2	-5.12	115.83	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2279	G	O5'-P-OP2	-5.12	101.09	105.70
26	14	2381	C	N3-C4-C5	5.12	123.95	121.90
52	J5	51	TYR	CA-CB-CG	5.12	123.12	113.40
26	1H	737	C	N1-C2-O2	-5.12	115.83	118.90
26	1H	811	U	OP1-P-OP2	5.12	127.28	119.60
26	1H	1566	A	C6-N1-C2	-5.12	115.53	118.60
1	1G	1488	G	C8-N9-C4	5.12	108.45	106.40
23	2K	48	U	C2-N1-C1'	5.12	123.84	117.70
26	1H	20	C	OP2-P-O3'	5.12	116.45	105.20
26	1H	1144	G	OP1-P-O3'	5.12	116.46	105.20
26	1H	1192	G	C5-N7-C8	5.12	106.86	104.30
26	1H	1793	C	C6-N1-C2	5.12	122.35	120.30
1	1G	250	A	P-O3'-C3'	5.12	125.84	119.70
1	1G	538	G	C8-N9-C1'	-5.12	120.35	127.00
1	1G	1499	A	N9-C4-C5	-5.12	103.75	105.80
26	14	2323	G	N7-C8-N9	-5.12	110.54	113.10
26	14	2626	C	C5-C4-N4	-5.12	116.62	120.20
1	13	67	C	N3-C2-O2	-5.11	118.32	121.90
26	1H	762	U	C5-C4-O4	-5.11	122.83	125.90
26	1H	2237	G	N9-C4-C5	-5.11	103.36	105.40
26	1H	2346	A	P-O3'-C3'	5.11	125.84	119.70
26	1H	2438	U	C4-C5-C6	5.11	122.77	119.70
26	14	771	G	OP1-P-O3'	5.11	116.45	105.20
26	14	2441	C	C5-C4-N4	5.11	123.78	120.20
26	14	2487	G	N3-C4-N9	5.11	129.07	126.00
26	1H	180	G	N3-C4-N9	5.11	129.07	126.00
26	1H	1430	C	OP1-P-O3'	5.11	116.45	105.20
26	1H	1934	C	OP1-P-O3'	5.11	116.45	105.20
26	1H	585	G	C5-C6-O6	-5.11	125.53	128.60
26	1H	617	G	OP1-P-OP2	-5.11	111.94	119.60
26	1H	736	C	O5'-P-OP1	-5.11	101.10	105.70
26	1H	1804	C	C5-C6-N1	-5.11	118.44	121.00
26	1H	2748	A	O5'-P-OP1	-5.11	101.10	105.70
1	1G	1055	A	N9-C4-C5	-5.11	103.76	105.80
1	1G	1529	G	C8-N9-C4	-5.11	104.36	106.40
26	14	2277	G	C4-C5-N7	-5.11	108.75	110.80
26	1H	1430	C	N3-C4-C5	-5.11	119.86	121.90
26	1H	1939	U	C4-C5-C6	-5.11	116.63	119.70
26	1H	2014	A	C6-N1-C2	-5.11	115.53	118.60
1	1G	306	G	C4-N9-C1'	-5.11	119.86	126.50
26	14	852	G	O5'-P-OP2	-5.11	101.10	105.70
26	14	1267	U	OP2-P-O3'	5.11	116.44	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	2877	G	C5-C6-N1	-5.11	108.94	111.50
1	13	226	G	C4-N9-C1'	5.11	133.14	126.50
26	1H	779	U	N1-C2-N3	-5.11	111.84	114.90
27	16	96	G	C2-N3-C4	5.11	114.45	111.90
1	1G	1435	G	C6-C5-N7	-5.11	127.33	130.40
26	14	748	G	C4-N9-C1'	-5.11	119.86	126.50
26	14	1621	U	N1-C2-O2	-5.11	119.22	122.80
26	14	1941	C	O5'-P-OP1	-5.11	101.10	105.70
26	14	2419	U	C5-C6-N1	5.11	125.25	122.70
26	14	2432	A	N1-C2-N3	5.11	131.85	129.30
26	14	2584	U	N3-C2-O2	-5.11	118.62	122.20
1	13	291	C	O5'-P-OP2	-5.11	101.11	105.70
1	13	690	G	C5-C6-N1	-5.11	108.95	111.50
1	13	1091	U	N3-C2-O2	-5.11	118.63	122.20
1	13	1500	A	C2-N3-C4	5.11	113.15	110.60
26	1H	1161	C	OP1-P-OP2	-5.11	111.94	119.60
26	1H	1308	A	N1-C2-N3	5.11	131.85	129.30
26	1H	1936	A	C4-C5-N7	5.11	113.25	110.70
26	1H	2089	U	N3-C2-O2	5.11	125.77	122.20
26	1H	2488	A	C2-N3-C4	-5.11	108.05	110.60
1	1G	1441	G	O5'-P-OP1	-5.11	101.11	105.70
26	14	2308	G	N3-C4-N9	-5.11	122.94	126.00
26	1H	202	U	N1-C2-O2	5.10	126.37	122.80
26	1H	1368	G	C5-C6-N1	5.10	114.05	111.50
26	1H	2447	G	N3-C4-N9	5.10	129.06	126.00
26	14	1498	C	OP1-P-O3'	5.10	116.43	105.20
1	13	1139	G	C8-N9-C4	5.10	108.44	106.40
1	13	1502	A	N3-C4-C5	5.10	130.37	126.80
26	1H	188	G	N3-C2-N2	5.10	123.47	119.90
26	1H	966	G	O5'-P-OP2	-5.10	101.11	105.70
26	1H	1281	G	C4-C5-C6	-5.10	115.74	118.80
26	1H	1393	A	O4'-C1'-N9	5.10	112.28	108.20
26	1H	1671	U	C2-N1-C1'	5.10	123.82	117.70
26	1H	2056	G	N1-C2-N3	5.10	126.96	123.90
26	1H	2233	U	N1-C2-O2	-5.10	119.23	122.80
26	1H	2509	G	N3-C2-N2	5.10	123.47	119.90
26	14	608	A	C2-N3-C4	-5.10	108.05	110.60
26	14	779	U	N3-C2-O2	-5.10	118.63	122.20
26	14	1663	C	N3-C4-N4	5.10	121.57	118.00
26	1H	189	G	N7-C8-N9	-5.10	110.55	113.10
26	1H	1513	C	C6-N1-C2	-5.10	118.26	120.30
1	13	580	U	N3-C4-O4	-5.10	115.83	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	1128	C	N3-C2-O2	-5.10	118.33	121.90
26	1H	386	G	C2-N3-C4	5.10	114.45	111.90
26	1H	701	G	N1-C2-N3	5.10	126.96	123.90
26	1H	1360	A	C6-C5-N7	-5.10	128.73	132.30
26	1H	1366	A	C4-C5-N7	5.10	113.25	110.70
26	1H	1786	A	C1'-O4'-C4'	-5.10	105.82	109.90
26	1H	2270	G	N9-C4-C5	-5.10	103.36	105.40
1	1G	112	G	N1-C6-O6	5.10	122.96	119.90
26	14	2433	A	O5'-P-OP2	5.10	116.82	110.70
1	13	872	A	O4'-C1'-N9	5.10	112.28	108.20
1	13	1286	A	C8-N9-C4	-5.10	103.76	105.80
26	1H	77	C	OP1-P-OP2	-5.10	111.95	119.60
26	1H	496	G	O5'-P-OP2	-5.10	101.11	105.70
26	1H	762	U	N3-C4-C5	5.10	117.66	114.60
26	1H	1767	C	N3-C2-O2	-5.10	118.33	121.90
26	1H	2070	G	C5-C6-N1	5.10	114.05	111.50
26	1H	2444	G	N1-C2-N2	5.10	120.79	116.20
1	1G	297	G	C8-N9-C4	5.10	108.44	106.40
26	14	1563	G	OP2-P-O3'	5.10	116.42	105.20
26	14	2237	G	OP1-P-OP2	5.10	127.25	119.60
1	13	798	G	N1-C2-N2	-5.10	111.61	116.20
1	13	803	G	C5-C6-O6	5.10	131.66	128.60
26	1H	1241	A	C4-C5-C6	5.10	119.55	117.00
26	1H	2721	A	C4-C5-N7	5.10	113.25	110.70
26	14	2249	U	N3-C4-C5	-5.10	111.54	114.60
26	14	2286	A	C4-N9-C1'	5.10	135.47	126.30
26	1H	617	G	C8-N9-C4	5.09	108.44	106.40
26	1H	1190	G	C5-C6-O6	-5.09	125.54	128.60
26	1H	1215	G	C5-C6-N1	5.09	114.05	111.50
26	1H	1218	C	C5-C4-N4	-5.09	116.63	120.20
26	1H	2065	C	N3-C4-C5	5.09	123.94	121.90
26	1H	2779	U	C4-C5-C6	5.09	122.76	119.70
26	14	1801	G	C4-C5-N7	5.09	112.84	110.80
26	1H	801	G	N3-C2-N2	-5.09	116.33	119.90
26	1H	928	G	N1-C6-O6	5.09	122.96	119.90
26	1H	1948	G	C5-C6-O6	5.09	131.66	128.60
26	1H	2373	G	C8-N9-C1'	-5.09	120.38	127.00
26	1H	2592	G	N3-C4-C5	-5.09	126.05	128.60
26	14	782	A	C5-C6-N1	5.09	120.25	117.70
26	14	1608	A	O5'-P-OP1	-5.09	101.12	105.70
1	13	1407	C	O5'-P-OP2	-5.09	101.12	105.70
26	1H	526	A	C6-C5-N7	5.09	135.86	132.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	961	C	N3-C2-O2	-5.09	118.33	121.90
26	1H	1265	A	C4-C5-C6	5.09	119.55	117.00
26	1H	1879	C	C6-N1-C2	-5.09	118.26	120.30
26	1H	1904	G	N9-C1'-C2'	-5.09	106.40	112.00
26	14	191	A	N1-C6-N6	5.09	121.66	118.60
26	14	215	G	C6-C5-N7	-5.09	127.34	130.40
26	14	1320	C	N3-C4-N4	5.09	121.56	118.00
26	1H	1321	A	N1-C2-N3	5.09	131.84	129.30
26	1H	1692	U	C6-N1-C1'	-5.09	114.07	121.20
26	1H	2026	C	N3-C4-C5	-5.09	119.86	121.90
26	1H	2757	A	O5'-P-OP2	-5.09	101.12	105.70
26	14	610	C	C5-C6-N1	-5.09	118.45	121.00
26	14	677	A	OP1-P-OP2	5.09	127.23	119.60
26	14	2287	A	N1-C6-N6	5.09	121.65	118.60
1	13	189	U	C2-N1-C1'	5.09	123.81	117.70
26	1H	196	A	C6-N1-C2	5.09	121.65	118.60
26	1H	1376	C	N1-C2-N3	5.09	122.76	119.20
26	1H	2392	A	C6-C5-N7	-5.09	128.74	132.30
26	14	194	G	N1-C6-O6	5.09	122.95	119.90
26	14	1826	G	N3-C4-C5	-5.09	126.06	128.60
26	14	1904	G	C5-N7-C8	5.09	106.84	104.30
26	14	2269	A	OP1-P-O3'	5.09	116.39	105.20
26	1H	787	U	C2-N3-C4	-5.09	123.95	127.00
26	1H	1340	U	O4'-C1'-N1	5.09	112.27	108.20
26	1H	1400	G	C8-N9-C4	-5.09	104.36	106.40
26	14	733	G	OP1-P-OP2	5.09	127.23	119.60
26	14	992	C	C6-N1-C2	-5.09	118.27	120.30
26	14	2083	G	N1-C2-N3	5.09	126.95	123.90
26	14	2168	G	C4-N9-C1'	5.09	133.11	126.50
26	14	2564	A	N1-C6-N6	5.09	121.65	118.60
26	1H	508	G	C8-N9-C4	-5.08	104.37	106.40
26	1H	2445	G	N3-C2-N2	-5.08	116.34	119.90
26	1H	2512	C	N1-C2-O2	-5.08	115.85	118.90
26	1H	2751	G	C4-C5-N7	5.08	112.83	110.80
1	1G	555	C	C6-N1-C2	-5.08	118.27	120.30
26	14	725	G	O5'-P-OP1	-5.08	101.12	105.70
26	14	848	G	C2-N3-C4	5.08	114.44	111.90
26	14	2850	A	OP1-P-O3'	5.08	116.39	105.20
26	1H	432	A	N1-C2-N3	-5.08	126.76	129.30
26	1H	940	G	C5-C6-N1	5.08	114.04	111.50
26	1H	1695	G	C4-N9-C1'	5.08	133.11	126.50
26	1H	1784	A	OP1-P-O3'	5.08	116.38	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2056	G	C6-N1-C2	-5.08	122.05	125.10
23	2L	48	U	OP2-P-O3'	5.08	116.38	105.20
26	14	74	A	C6-C5-N7	-5.08	128.74	132.30
26	14	122	G	N7-C8-N9	-5.08	110.56	113.10
26	14	603	A	C4-N9-C1'	5.08	135.45	126.30
26	14	729	G	C2-N3-C4	5.08	114.44	111.90
26	14	2523	G	N1-C6-O6	5.08	122.95	119.90
26	14	2607	G	O5'-P-OP2	-5.08	101.12	105.70
1	13	35	G	C8-N9-C4	-5.08	104.37	106.40
1	13	278	G	N3-C4-N9	5.08	129.05	126.00
23	2K	31	G	C5-C6-N1	-5.08	108.96	111.50
26	1H	62	C	OP2-P-O3'	5.08	116.38	105.20
26	1H	1162	G	N9-C4-C5	5.08	107.43	105.40
26	1H	1400	G	N9-C4-C5	5.08	107.43	105.40
26	1H	2007	C	O5'-P-OP2	-5.08	101.13	105.70
54	P8	11	LYS	CD-CE-NZ	5.08	123.39	111.70
1	1G	815	A	OP2-P-O3'	5.08	116.38	105.20
26	14	49	A	O5'-P-OP2	-5.08	101.13	105.70
26	14	79	G	N1-C6-O6	5.08	122.95	119.90
26	14	805	G	OP1-P-O3'	5.08	116.38	105.20
26	14	871	U	O5'-P-OP2	5.08	116.80	110.70
26	1H	577	G	O5'-P-OP1	5.08	116.80	110.70
26	1H	2440	C	C6-N1-C1'	5.08	126.90	120.80
26	14	1414	G	N3-C4-N9	5.08	129.05	126.00
26	14	1779	U	C6-N1-C1'	-5.08	114.09	121.20
26	14	2429	G	OP1-P-OP2	-5.08	111.98	119.60
1	13	1205	U	C6-N1-C2	-5.08	117.95	121.00
26	1H	2486	G	O5'-P-OP1	5.08	116.79	110.70
26	1H	2701	C	C6-N1-C2	-5.08	118.27	120.30
29	21	119	ARG	NE-CZ-NH1	5.08	122.84	120.30
26	14	138	G	N1-C6-O6	5.08	122.95	119.90
26	14	477	A	C2-N3-C4	-5.08	108.06	110.60
26	14	1695	G	C6-C5-N7	-5.08	127.35	130.40
26	14	2423	U	C2-N3-C4	-5.08	123.95	127.00
26	14	2491	U	N3-C2-O2	5.08	125.75	122.20
27	1J	56	G	C8-N9-C1'	-5.08	120.40	127.00
26	1H	2260	C	C5-C4-N4	-5.08	116.65	120.20
22	1L	40	C	C6-N1-C2	-5.08	118.27	120.30
1	13	574	A	OP1-P-OP2	-5.08	111.99	119.60
26	1H	650	C	OP1-P-O3'	5.08	116.37	105.20
26	1H	1203	G	OP1-P-O3'	5.08	116.37	105.20
26	1H	1264	G	C4-C5-N7	5.08	112.83	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1388	G	O5'-P-OP2	-5.08	101.13	105.70
26	1H	1787	A	O4'-C1'-N9	-5.08	104.14	108.20
26	1H	2582	G	N9-C4-C5	-5.08	103.37	105.40
26	1H	2711	A	OP2-P-O3'	-5.08	94.03	105.20
1	1G	150	C	C5-C6-N1	5.08	123.54	121.00
1	1G	972	C	C5-C6-N1	5.08	123.54	121.00
26	14	947	G	N3-C4-C5	5.08	131.14	128.60
26	14	1623	G	N3-C2-N2	-5.08	116.35	119.90
1	13	765	G	C6-C5-N7	-5.07	127.36	130.40
26	1H	315	G	N1-C6-O6	5.07	122.94	119.90
26	1H	1318	C	N1-C2-O2	5.07	121.94	118.90
26	1H	1325	G	C8-N9-C4	-5.07	104.37	106.40
26	1H	1888	G	C2-N3-C4	5.07	114.44	111.90
26	1H	2497	A	C6-N1-C2	-5.07	115.56	118.60
26	14	1257	C	N1-C2-O2	-5.07	115.86	118.90
26	14	1909	C	N3-C2-O2	-5.07	118.35	121.90
26	14	1926	U	N1-C2-N3	5.07	117.94	114.90
26	1H	468	G	OP1-P-OP2	-5.07	111.99	119.60
26	1H	916	G	O5'-P-OP2	5.07	116.79	110.70
26	1H	1213	A	C2-N3-C4	-5.07	108.06	110.60
26	14	510	C	OP1-P-OP2	5.07	127.21	119.60
26	1H	940	G	C5-C6-O6	-5.07	125.56	128.60
26	1H	1352	U	N1-C2-O2	-5.07	119.25	122.80
26	1H	1491	G	C5-C6-N1	-5.07	108.97	111.50
26	1H	2319	G	N3-C4-N9	5.07	129.04	126.00
26	1H	2641	G	C6-C5-N7	-5.07	127.36	130.40
27	16	41	U	N3-C4-O4	-5.07	115.85	119.40
1	1G	691	G	C5-C6-O6	-5.07	125.56	128.60
26	14	1955	U	C2-N3-C4	-5.07	123.96	127.00
27	1J	84	C	C5-C6-N1	-5.07	118.46	121.00
23	2K	5	G	N3-C4-C5	5.07	131.13	128.60
26	1H	177	G	N3-C4-C5	-5.07	126.06	128.60
26	1H	2393	A	O5'-P-OP1	-5.07	101.14	105.70
26	1H	2512	C	C6-N1-C2	5.07	122.33	120.30
26	14	488	G	O5'-P-OP2	-5.07	101.14	105.70
1	13	306	G	C6-C5-N7	5.07	133.44	130.40
1	13	768	A	N1-C2-N3	5.07	131.83	129.30
26	1H	9	U	C2-N1-C1'	5.07	123.78	117.70
26	1H	1304	C	C2-N1-C1'	-5.07	113.22	118.80
26	1H	1909	C	OP1-P-O3'	5.07	116.35	105.20
1	1G	52	G	N1-C6-O6	-5.07	116.86	119.90
1	1G	1390	U	C5-C4-O4	5.07	128.94	125.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	962	G	N1-C6-O6	5.07	122.94	119.90
26	14	1029	A	N1-C2-N3	-5.07	126.77	129.30
26	14	1314	C	C4-C5-C6	-5.07	114.87	117.40
26	14	2242	G	C8-N9-C4	-5.07	104.37	106.40
26	14	2275	C	N3-C2-O2	-5.07	118.35	121.90
1	13	319	G	N1-C6-O6	-5.07	116.86	119.90
1	13	1096	C	C5-C6-N1	5.07	123.53	121.00
26	1H	649	G	O5'-P-OP2	-5.07	101.14	105.70
26	1H	691	C	N1-C2-O2	-5.07	115.86	118.90
26	1H	912	C	N3-C4-N4	5.07	121.55	118.00
26	1H	1281	G	C5-C6-N1	5.07	114.03	111.50
26	1H	1610	A	C4-C5-N7	5.07	113.23	110.70
26	1H	2712	U	C2-N3-C4	-5.07	123.96	127.00
1	1G	1200	C	N3-C2-O2	-5.07	118.35	121.90
2	12	196	LEU	CA-CB-CG	5.07	126.95	115.30
26	14	1594	G	O5'-P-OP2	5.07	116.78	110.70
26	14	1612	C	C5-C4-N4	-5.07	116.66	120.20
26	14	1617	C	N1-C2-O2	-5.07	115.86	118.90
26	14	2430	A	C4-C5-N7	5.07	113.23	110.70
26	14	2596	U	O5'-P-OP1	-5.07	101.14	105.70
26	1H	606	U	C5-C6-N1	-5.06	120.17	122.70
26	1H	1926	U	C5-C4-O4	5.06	128.94	125.90
26	1H	2491	U	C4-C5-C6	-5.06	116.66	119.70
26	14	781	A	O5'-P-OP1	-5.06	101.14	105.70
26	14	2330	G	C5-C6-O6	-5.06	125.56	128.60
27	1J	78	A	N1-C6-N6	5.06	121.64	118.60
1	13	857	C	C4-C5-C6	5.06	119.93	117.40
24	3K	36	A	P-O3'-C3'	5.06	125.77	119.70
26	1H	1144	G	C5-C6-O6	-5.06	125.56	128.60
26	14	470	A	C4-C5-N7	5.06	113.23	110.70
26	14	1642	G	N1-C6-O6	-5.06	116.86	119.90
26	14	1977	A	N1-C2-N3	5.06	131.83	129.30
26	14	2419	U	OP1-P-O3'	5.06	116.34	105.20
26	1H	655	A	N1-C2-N3	5.06	131.83	129.30
26	1H	860	U	O5'-P-OP2	-5.06	101.14	105.70
26	14	236	C	C6-N1-C2	5.06	122.32	120.30
26	14	921	G	N7-C8-N9	5.06	115.63	113.10
1	13	738	C	C5-C6-N1	5.06	123.53	121.00
26	1H	101	G	C8-N9-C4	5.06	108.42	106.40
26	1H	255	A	C5-C6-N6	-5.06	119.65	123.70
26	1H	1398	C	OP2-P-O3'	5.06	116.33	105.20
26	1H	2256	G	N3-C2-N2	5.06	123.44	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	2595	G	C5-C6-N1	5.06	114.03	111.50
1	1G	221	C	C5-C6-N1	5.06	123.53	121.00
1	1G	481	G	C8-N9-C1'	-5.06	120.42	127.00
26	14	1698	A	C4-C5-C6	5.06	119.53	117.00
26	14	2048	G	N1-C6-O6	-5.06	116.86	119.90
26	14	2417	C	O5'-P-OP2	-5.06	101.15	105.70
26	14	2447	G	C2-N3-C4	-5.06	109.37	111.90
1	13	772	U	C5-C4-O4	5.06	128.93	125.90
1	13	1505	G	O5'-P-OP1	-5.06	101.15	105.70
26	1H	262	A	N9-C4-C5	-5.06	103.78	105.80
26	1H	858	U	OP2-P-O3'	5.06	116.33	105.20
26	1H	907	U	C5-C6-N1	-5.06	120.17	122.70
26	1H	949	C	C6-N1-C2	5.06	122.32	120.30
26	1H	1934	C	N3-C2-O2	5.06	125.44	121.90
1	1G	354	G	C4-C5-N7	5.06	112.82	110.80
26	14	46	C	N3-C4-N4	-5.06	114.46	118.00
26	14	1085	A	P-O3'-C3'	5.06	125.77	119.70
28	19	41	GLY	C-N-CA	-5.06	111.68	122.30
1	13	266	G	C8-N9-C4	-5.06	104.38	106.40
26	1H	696	G	N1-C2-N2	-5.06	111.65	116.20
26	1H	906	G	N9-C4-C5	5.06	107.42	105.40
1	1G	1247	U	C5-C6-N1	5.06	125.23	122.70
26	14	1790	C	N3-C2-O2	5.06	125.44	121.90
26	14	2600	A	N1-C6-N6	-5.06	115.57	118.60
26	14	2723	C	N1-C2-O2	-5.06	115.87	118.90
1	13	975	A	N1-C6-N6	5.05	121.63	118.60
1	13	1414	U	OP2-P-O3'	5.05	116.32	105.20
26	1H	1365	A	C8-N9-C4	-5.05	103.78	105.80
26	1H	1379	A	C8-N9-C4	-5.05	103.78	105.80
26	1H	1389	G	OP1-P-O3'	5.05	116.32	105.20
26	1H	1425	G	C5-C6-O6	-5.05	125.57	128.60
26	1H	1599	C	N3-C4-N4	-5.05	114.46	118.00
26	1H	2455	G	N1-C2-N3	5.05	126.93	123.90
26	1H	2823	A	N7-C8-N9	5.05	116.33	113.80
25	4L	14	A	C8-N9-C4	5.05	107.82	105.80
26	14	2053	G	C2-N3-C4	5.05	114.43	111.90
1	13	571	U	O5'-P-OP2	5.05	116.76	110.70
26	1H	1123	C	N3-C4-C5	5.05	123.92	121.90
26	1H	1971	A	C5-C6-N6	-5.05	119.66	123.70
24	3L	16	U	N1-C2-O2	5.05	126.34	122.80
26	14	462	C	N1-C2-N3	5.05	122.74	119.20
26	14	570	G	C4-C5-C6	5.05	121.83	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	858	G	C8-N9-C4	-5.05	104.38	106.40
26	1H	270(C)	C	C6-N1-C2	5.05	122.32	120.30
26	1H	536	A	C6-N1-C2	-5.05	115.57	118.60
26	1H	2611	U	N3-C4-O4	-5.05	115.86	119.40
42	D8	40	LEU	CA-CB-CG	5.05	126.92	115.30
26	14	621	A	C6-C5-N7	-5.05	128.76	132.30
26	14	984	A	C8-N9-C4	5.05	107.82	105.80
26	14	1247	A	O5'-P-OP1	5.05	116.76	110.70
26	14	1551	C	C2-N3-C4	5.05	122.43	119.90
1	13	1214	C	C2-N1-C1'	-5.05	113.25	118.80
26	1H	586	A	N7-C8-N9	-5.05	111.28	113.80
26	1H	663	G	C4-N9-C1'	5.05	133.06	126.50
26	1H	1569	A	C8-N9-C4	-5.05	103.78	105.80
1	1G	265	G	C8-N9-C4	5.05	108.42	106.40
1	1G	341	C	C6-N1-C2	5.05	122.32	120.30
26	14	2238	G	N3-C2-N2	-5.05	116.36	119.90
26	14	2256	G	C6-C5-N7	-5.05	127.37	130.40
26	1H	1708	C	C5-C4-N4	-5.05	116.67	120.20
26	1H	2039	C	C6-N1-C2	-5.05	118.28	120.30
26	1H	2377	A	N3-C4-C5	5.05	130.33	126.80
26	14	573	G	C2-N3-C4	5.05	114.42	111.90
1	13	558	G	C5-C6-N1	-5.05	108.98	111.50
1	13	690	G	C5-C6-O6	-5.05	125.57	128.60
1	13	1286	A	N7-C8-N9	5.05	116.32	113.80
26	1H	77	C	N3-C4-N4	5.05	121.53	118.00
26	1H	734	A	N7-C8-N9	5.05	116.32	113.80
26	1H	1550	C	N3-C2-O2	5.05	125.43	121.90
26	1H	1839	G	N1-C6-O6	5.05	122.93	119.90
26	1H	2246	G	C6-C5-N7	5.05	133.43	130.40
26	1H	2447	G	P-O3'-C3'	5.05	125.76	119.70
26	1H	2456	C	N3-C4-N4	5.05	121.53	118.00
26	1H	2772	C	C5-C6-N1	-5.05	118.48	121.00
52	N8	40	LYS	C-N-CD	5.05	139.00	128.40
1	1G	422	C	O4'-C1'-N1	5.05	112.24	108.20
9	82	128	ARG	NE-CZ-NH1	5.05	122.82	120.30
26	14	2039	C	C2-N1-C1'	5.05	124.35	118.80
1	13	299	G	N3-C4-N9	-5.04	122.97	126.00
26	1H	138	G	N1-C6-O6	-5.04	116.87	119.90
26	1H	2426	A	O5'-P-OP2	-5.04	101.16	105.70
27	1J	7	G	C4-C5-N7	5.04	112.82	110.80
1	13	20	U	O5'-P-OP2	-5.04	101.16	105.70
26	1H	536	A	O5'-P-OP1	5.04	116.75	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1256	G	N1-C6-O6	5.04	122.93	119.90
26	1H	1317	A	O5'-P-OP2	-5.04	101.16	105.70
26	1H	1605	C	C2-N3-C4	-5.04	117.38	119.90
26	1H	2070	G	N9-C4-C5	-5.04	103.38	105.40
26	14	747	U	C6-N1-C2	5.04	124.03	121.00
26	14	1648	C	C4-C5-C6	5.04	119.92	117.40
1	13	1082	G	N1-C6-O6	5.04	122.92	119.90
26	1H	1554	A	OP1-P-O3'	5.04	116.29	105.20
26	1H	2655	G	C8-N9-C1'	5.04	133.56	127.00
41	C8	74	LEU	CA-CB-CG	5.04	126.90	115.30
26	14	1262	A	N1-C6-N6	5.04	121.62	118.60
26	14	2047	U	N3-C4-C5	5.04	117.62	114.60
26	14	2435	A	C5-C6-N1	5.04	120.22	117.70
1	13	758	G	N3-C4-C5	5.04	131.12	128.60
26	1H	786	C	N3-C4-N4	-5.04	114.47	118.00
26	1H	2566	A	P-O3'-C3'	5.04	125.75	119.70
1	1G	44	G	C4-N9-C1'	5.04	133.05	126.50
26	14	949	C	OP2-P-O3'	5.04	116.29	105.20
26	14	970	C	C2-N1-C1'	-5.04	113.26	118.80
26	14	2445	G	C8-N9-C4	-5.04	104.38	106.40
26	1H	908	C	C4-C5-C6	5.04	119.92	117.40
26	1H	1704	G	C8-N9-C4	5.04	108.42	106.40
26	1H	1947	C	N3-C4-N4	5.04	121.53	118.00
26	1H	2197	U	OP2-P-O3'	5.04	116.29	105.20
27	16	8	U	O4'-C1'-N1	5.04	112.23	108.20
1	1G	1025	U	C2-N1-C1'	5.04	123.75	117.70
26	14	621	A	C6-N1-C2	5.04	121.62	118.60
26	14	1349	A	C2-N3-C4	-5.04	108.08	110.60
26	14	2318	G	N7-C8-N9	5.04	115.62	113.10
27	1J	83	G	C5-C6-N1	-5.04	108.98	111.50
26	1H	946	G	C4-C5-N7	-5.04	108.78	110.80
26	1H	1380	G	N3-C4-N9	5.04	129.02	126.00
26	14	560	C	C6-N1-C2	5.04	122.31	120.30
26	14	1900	A	C6-C5-N7	-5.04	128.77	132.30
26	14	2014	A	C8-N9-C4	5.04	107.81	105.80
1	13	1488	G	C4-C5-C6	5.04	121.82	118.80
26	1H	802	A	C5-N7-C8	-5.04	101.38	103.90
26	1H	2332	U	OP1-P-OP2	-5.04	112.05	119.60
26	1H	2457	U	N3-C2-O2	5.04	125.72	122.20
26	14	270(Y)	G	C5-C6-O6	5.04	131.62	128.60
26	14	2444	G	C2-N3-C4	-5.04	109.38	111.90
1	13	788	U	N3-C2-O2	-5.03	118.68	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	1298	C	C4-C5-C6	-5.03	114.88	117.40
26	1H	2417	C	OP2-P-O3'	5.03	116.27	105.20
26	14	2249	U	C5-C6-N1	5.03	125.22	122.70
1	1G	894	G	C5-C6-O6	-5.03	125.58	128.60
26	14	631	A	N7-C8-N9	-5.03	111.28	113.80
1	13	816	A	N9-C4-C5	5.03	107.81	105.80
26	1H	586	A	C6-C5-N7	5.03	135.82	132.30
26	1H	1895	C	N1-C2-O2	-5.03	115.88	118.90
26	1H	2274	A	OP2-P-O3'	5.03	116.27	105.20
1	1G	505	G	C4-C5-N7	5.03	112.81	110.80
26	14	2321	G	C4-N9-C1'	5.03	133.04	126.50
26	1H	223	A	N1-C6-N6	5.03	121.62	118.60
26	1H	530	G	C8-N9-C1'	5.03	133.54	127.00
26	1H	1373	A	OP1-P-OP2	-5.03	112.06	119.60
26	1H	2491	U	C6-N1-C2	5.03	124.02	121.00
26	14	2373	G	N3-C2-N2	-5.03	116.38	119.90
1	13	191	G	C8-N9-C4	-5.03	104.39	106.40
1	13	660	G	O5'-P-OP2	5.03	116.73	110.70
26	1H	130	C	C6-N1-C1'	-5.03	114.77	120.80
26	1H	500	G	O5'-P-OP1	-5.03	101.18	105.70
26	1H	1487	G	OP1-P-O3'	5.03	116.26	105.20
26	14	453	C	C5-C6-N1	-5.03	118.49	121.00
26	14	675	A	C5-C6-N6	-5.03	119.68	123.70
26	14	1298	C	C5-C6-N1	5.03	123.51	121.00
26	1H	37	C	C4-C5-C6	5.03	119.91	117.40
26	1H	1398	C	OP1-P-OP2	-5.03	112.06	119.60
26	1H	1672	C	N3-C2-O2	5.03	125.42	121.90
26	1H	1894	C	N3-C4-C5	-5.03	119.89	121.90
26	1H	2352	A	O5'-P-OP1	-5.03	101.18	105.70
26	1H	2363	C	C2-N1-C1'	-5.03	113.27	118.80
26	14	629	G	N1-C6-O6	-5.03	116.88	119.90
26	14	980	A	C4-C5-C6	5.03	119.51	117.00
26	14	1380	G	O5'-P-OP2	-5.03	101.18	105.70
26	14	2036	C	O5'-P-OP1	5.03	116.73	110.70
26	14	2420	C	N3-C4-C5	5.03	123.91	121.90
26	1H	800	A	N7-C8-N9	-5.02	111.29	113.80
26	1H	906	G	C5-C6-O6	5.02	131.61	128.60
26	1H	1697	G	OP1-P-O3'	5.02	116.25	105.20
26	14	127	A	C6-N1-C2	-5.02	115.58	118.60
26	14	1914	C	N3-C2-O2	-5.02	118.38	121.90
1	13	781	A	C8-N9-C4	5.02	107.81	105.80
1	13	812	C	C4-C5-C6	5.02	119.91	117.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	13	1522	U	C5-C6-N1	-5.02	120.19	122.70
26	1H	642	G	N3-C4-N9	-5.02	122.99	126.00
26	1H	827	U	N3-C2-O2	5.02	125.72	122.20
26	1H	1278	A	C8-N9-C4	5.02	107.81	105.80
1	1G	175	C	C5-C6-N1	5.02	123.51	121.00
22	1L	34	G	C4-N9-C1'	5.02	133.03	126.50
26	14	213	A	N3-C4-C5	5.02	130.32	126.80
26	14	214	G	N7-C8-N9	5.02	115.61	113.10
26	14	479	A	C5-C6-N6	5.02	127.72	123.70
26	14	778	G	O5'-P-OP2	-5.02	101.18	105.70
26	14	811	U	C5-C6-N1	-5.02	120.19	122.70
26	14	1482	U	C6-N1-C1'	5.02	128.23	121.20
26	14	1598	C	OP1-P-OP2	-5.02	112.07	119.60
26	14	2054	A	N7-C8-N9	5.02	116.31	113.80
26	14	2555	U	N1-C2-O2	-5.02	119.28	122.80
26	14	2642	G	C5-C6-O6	-5.02	125.59	128.60
26	1H	1010	A	OP2-P-O3'	5.02	116.25	105.20
26	14	2426	A	C5-N7-C8	-5.02	101.39	103.90
1	13	540	G	O5'-P-OP2	-5.02	101.18	105.70
1	13	813	U	N3-C4-C5	5.02	117.61	114.60
26	1H	180	G	C4-C5-N7	5.02	112.81	110.80
26	1H	375	C	O5'-P-OP1	5.02	116.72	110.70
26	1H	683	C	C4-C5-C6	-5.02	114.89	117.40
26	1H	827	U	C6-N1-C2	5.02	124.01	121.00
26	1H	2275	C	O4'-C1'-N1	-5.02	104.18	108.20
26	1H	2518	A	N3-C4-C5	5.02	130.31	126.80
26	1H	2763	G	N9-C4-C5	-5.02	103.39	105.40
1	1G	913	A	C8-N9-C4	-5.02	103.79	105.80
26	14	947	G	N3-C4-N9	-5.02	122.99	126.00
26	14	1958	C	N3-C2-O2	5.02	125.41	121.90
26	14	2225	A	N9-C1'-C2'	-5.02	106.48	112.00
26	14	2615	U	N3-C2-O2	-5.02	118.69	122.20
26	1H	866	A	N1-C6-N6	5.02	121.61	118.60
26	1H	935	C	C5-C6-N1	-5.02	118.49	121.00
26	1H	1780	A	OP1-P-OP2	-5.02	112.08	119.60
26	1H	1781	C	N3-C4-C5	5.02	123.91	121.90
27	16	81	G	C2-N3-C4	-5.02	109.39	111.90
1	1G	1237	C	C5-C6-N1	5.02	123.51	121.00
1	1G	1502	A	C8-N9-C4	-5.02	103.79	105.80
26	14	665	C	C6-N1-C2	5.02	122.31	120.30
26	14	2296	U	OP1-P-O3'	5.02	116.24	105.20
1	13	266	G	C6-C5-N7	-5.02	127.39	130.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	1H	435	C	N3-C2-O2	-5.02	118.39	121.90
26	1H	593	G	C5-C6-O6	-5.02	125.59	128.60
26	1H	2751	G	C6-C5-N7	-5.02	127.39	130.40
26	14	188	G	C8-N9-C4	5.02	108.41	106.40
26	14	514	A	C5-N7-C8	5.02	106.41	103.90
26	14	1893	C	C6-N1-C2	-5.02	118.29	120.30
26	14	1939	U	C2-N3-C4	-5.02	123.99	127.00
26	14	2011	U	C2-N1-C1'	-5.02	111.68	117.70
1	13	909	A	C5-C6-N1	-5.01	115.19	117.70
23	2K	68	C	C5-C4-N4	-5.01	116.69	120.20
26	1H	132	G	C5-N7-C8	-5.01	101.79	104.30
26	1H	212	G	C8-N9-C4	5.01	108.41	106.40
26	1H	404	C	N1-C2-O2	-5.01	115.89	118.90
26	1H	919	G	C6-N1-C2	-5.01	122.09	125.10
26	1H	992	C	C5-C6-N1	5.01	123.51	121.00
26	1H	1300	U	OP2-P-O3'	-5.01	94.17	105.20
26	1H	1681	G	C6-C5-N7	-5.01	127.39	130.40
26	1H	2040	C	C6-N1-C2	5.01	122.31	120.30
1	1G	22	G	C8-N9-C4	-5.01	104.39	106.40
1	1G	865	A	C8-N9-C4	-5.01	103.79	105.80
1	1G	1400	C	N1-C2-O2	5.01	121.91	118.90
1	1G	1524	C	C2-N3-C4	-5.01	117.39	119.90
18	9A	31	LEU	CA-CB-CG	5.01	126.83	115.30
26	14	240	G	N3-C4-C5	5.01	131.11	128.60
26	14	468	G	N3-C2-N2	-5.01	116.39	119.90
26	14	2315	G	OP1-P-O3'	5.01	116.23	105.20
26	1H	205	G	OP2-P-O3'	5.01	116.23	105.20
26	1H	1368	G	N3-C4-C5	-5.01	126.09	128.60
26	1H	2219	G	O5'-P-OP1	-5.01	101.19	105.70
1	1G	1300	G	P-O3'-C3'	5.01	125.72	119.70
26	14	762	U	OP1-P-O3'	5.01	116.23	105.20
26	14	1210	A	N7-C8-N9	5.01	116.31	113.80
26	14	2011	U	O5'-P-OP1	-5.01	101.19	105.70
1	13	1251	A	N1-C6-N6	-5.01	115.59	118.60
26	1H	145	G	N1-C6-O6	5.01	122.91	119.90
26	1H	623	G	C8-N9-C4	5.01	108.41	106.40
26	1H	2247	A	C5-C6-N1	-5.01	115.19	117.70
26	1H	2700	C	C5-C6-N1	-5.01	118.49	121.00
28	11	131	LEU	CB-CG-CD2	-5.01	102.48	111.00
1	1G	310	G	C8-N9-C1'	5.01	133.51	127.00
1	1G	497	U	N1-C2-O2	5.01	126.31	122.80
1	1G	769	G	N3-C4-C5	-5.01	126.09	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	660	G	N1-C6-O6	-5.01	116.89	119.90
26	14	2544	G	C4-C5-N7	5.01	112.80	110.80
1	13	496	A	OP2-P-O3'	5.01	116.22	105.20
1	13	503	C	C2-N1-C1'	5.01	124.31	118.80
26	1H	111	A	N1-C6-N6	-5.01	115.59	118.60
26	1H	219	G	C4-C5-N7	-5.01	108.80	110.80
26	1H	665	C	C5-C6-N1	-5.01	118.50	121.00
26	1H	1308	A	N7-C8-N9	5.01	116.31	113.80
26	1H	2071	A	C6-C5-N7	-5.01	128.79	132.30
24	3L	4	C	C6-N1-C2	-5.01	118.30	120.30
26	14	456	C	C5-C6-N1	-5.01	118.50	121.00
26	14	589	C	N3-C4-C5	-5.01	119.90	121.90
26	14	782	A	O5'-P-OP1	-5.01	101.19	105.70
26	14	1602	U	C4-C5-C6	5.01	122.70	119.70
1	13	1306	A	N1-C6-N6	5.01	121.61	118.60
1	13	1388	C	C6-N1-C2	-5.01	118.30	120.30
26	1H	1977	A	OP1-P-O3'	5.01	116.22	105.20
26	1H	2502	G	C4-N9-C1'	5.01	133.01	126.50
38	98	4	LEU	CA-CB-CG	-5.01	103.78	115.30
1	1G	509	A	C5-N7-C8	-5.01	101.40	103.90
26	1H	973	A	C5-C6-N1	-5.01	115.20	117.70
26	1H	1162	G	C5-N7-C8	-5.01	101.80	104.30
26	1H	1245	G	N3-C4-C5	-5.01	126.10	128.60
26	1H	1537	C	C5-C6-N1	5.01	123.50	121.00
1	1G	111	G	C6-C5-N7	-5.01	127.40	130.40
1	1G	702	A	O5'-P-OP2	-5.01	101.19	105.70
1	1G	1056	U	N1-C2-O2	-5.01	119.30	122.80
26	14	138	G	C4-N9-C1'	5.01	133.01	126.50
26	14	1636	C	N3-C4-C5	-5.01	119.90	121.90
26	14	1798	U	O5'-P-OP2	-5.01	101.19	105.70
26	14	1950	G	C6-C5-N7	-5.01	127.40	130.40
26	14	2445	G	N3-C4-C5	-5.01	126.10	128.60
26	14	2820	A	OP1-P-O3'	5.01	116.22	105.20
26	1H	19	C	C6-N1-C2	5.00	122.30	120.30
26	1H	1280	G	N9-C1'-C2'	-5.00	106.49	112.00
26	1H	1319	G	OP1-P-O3'	5.00	116.21	105.20
26	14	186	G	C6-N1-C2	-5.00	122.10	125.10
1	13	312	C	C6-N1-C2	-5.00	118.30	120.30
1	13	1530	G	C4-N9-C1'	-5.00	120.00	126.50
26	1H	579	G	OP2-P-O3'	5.00	116.21	105.20
26	1H	657	U	C2-N3-C4	-5.00	124.00	127.00
26	1H	1355	G	N3-C2-N2	5.00	123.40	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	14	791	C	OP2-P-O3'	5.00	116.21	105.20
26	14	1204	A	N1-C6-N6	5.00	121.60	118.60
26	14	1668	A	C8-N9-C4	5.00	107.80	105.80
26	14	2385	C	N3-C4-N4	5.00	121.50	118.00
26	14	2477	C	N1-C2-O2	5.00	121.90	118.90
26	14	2574	G	C5-C6-O6	-5.00	125.60	128.60
27	1J	94	C	N1-C2-O2	5.00	121.90	118.90
1	13	528	C	C6-N1-C2	5.00	122.30	120.30
1	13	713	G	OP1-P-O3'	5.00	116.20	105.20
1	13	800	G	C6-C5-N7	-5.00	127.40	130.40
1	13	1406	U	OP2-P-O3'	5.00	116.20	105.20
26	1H	734	A	OP1-P-OP2	5.00	127.10	119.60
26	1H	867	C	C6-N1-C2	5.00	122.30	120.30
26	1H	1202	C	OP1-P-OP2	5.00	127.10	119.60
26	1H	1299	G	C6-C5-N7	-5.00	127.40	130.40
26	1H	1357	U	O5'-P-OP2	-5.00	101.20	105.70
24	3L	16	U	C2-N1-C1'	5.00	123.70	117.70
26	14	373	U	C2-N1-C1'	5.00	123.70	117.70
26	14	802	A	N1-C2-N3	5.00	131.80	129.30
26	14	1193	G	N1-C6-O6	5.00	122.90	119.90
26	14	2385	C	N1-C2-O2	-5.00	115.90	118.90
36	35	62	LEU	CB-CA-C	-5.00	100.70	110.20

There are no chirality outliers.

All (142) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
28	11	197	GLY	Peptide
2	12	22	LYS	Peptide
34	15	41	ASP	Peptide
34	15	50	ASP	Peptide
28	19	237	GLU	Peptide
28	19	271	ILE	Peptide
28	19	32	SER	Peptide
28	19	37	LEU	Peptide
10	1A	55	LYS	Peptide
10	1A	87	THR	Peptide
2	1E	15	VAL	Peptide
29	21	56	PRO	Peptide
29	21	57	LYS	Peptide
29	21	82	ARG	Peptide
29	29	139	GLY	Peptide

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Mol	Chain	Res	Type	Group
29	29	186	GLY	Peptide
29	29	201	THR	Peptide
29	29	65	GLY	Peptide
11	2A	126	ARG	Peptide
11	2A	49	GLY	Peptide
30	31	133	ASN	Peptide
4	32	152	SER	Peptide
4	32	28	SER	Peptide
4	32	30	LYS	Peptide
4	32	33	MET	Peptide
36	35	110	TYR	Peptide
36	35	36	LYS	Peptide
36	35	64	LYS	Peptide
36	35	70	GLN	Peptide
30	39	127	GLU	Peptide
30	39	166	ALA	Peptide
30	39	20	LEU	Peptide
30	39	24	LEU	Peptide
30	39	26	ALA	Peptide
30	39	85	GLY	Peptide
12	3A	18	VAL	Peptide
4	3E	31	CYS	Peptide
12	3I	87	GLY	Peptide
31	41	85	GLY	Peptide
31	41	95	ARG	Peptide
37	45	134	ARG	Peptide
37	45	25	ASP	Peptide
37	45	26	TYR	Peptide
37	45	79	LEU	Peptide
37	45	86	GLY	Peptide
37	45	87	LYS	Peptide
31	49	13	GLU	Peptide
38	55	106	GLY	Peptide
34	58	56	ASN	Peptide
34	58	95	PRO	Peptide
32	59	155	SER	Peptide
32	59	89	ILE	Peptide
32	59	90	LYS	Peptide
33	61	11	ASN	Peptide
33	61	114	LEU	Peptide
33	61	134	PRO	Peptide
33	61	143	SER	Peptide

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Mol	Chain	Res	Type	Group
33	61	82	ARG	Peptide
39	65	59	LYS	Peptide
39	65	87	PHE	Peptide
35	68	27	GLY	Peptide
33	69	101	LEU	Peptide
33	69	112	LYS	Peptide
15	61	87	ILE	Peptide
40	75	4	GLY	Peptide
36	78	11	GLY	Peptide
36	78	5	ASP	Peptide
9	82	117	HIS	Peptide
9	82	118	LYS	Peptide
41	85	72	HIS	Peptide
41	85	98	LEU	Peptide
37	88	58	PHE	Peptide
37	88	78	PRO	Peptide
42	95	44	LYS	Peptide
38	98	1	MET	Peptide
38	98	44	LEU	Peptide
38	98	8	ARG	Peptide
43	A5	109	GLU	Peptide
43	A5	43	GLY	Peptide
43	A5	93	ALA	Peptide
39	A8	106	ARG	Peptide
39	A8	110	LEU	Peptide
19	AI	6	LYS	Peptide
19	AI	7	LYS	Peptide
44	B5	24	GLY	Peptide
44	B5	61	GLY	Peptide
40	B8	58	ASN	Peptide
20	BA	101	GLY	Peptide
20	BA	11	SER	Peptide
20	BA	72	LEU	Peptide
45	C5	100	ALA	Peptide
45	C5	102	CYS	Peptide
45	C5	81	LYS	Peptide
41	C8	90	VAL	Peptide
41	C8	92	ARG	Peptide
46	D5	142	SER	Peptide
46	D5	61	LEU	Peptide
48	F5	85	LEU	Peptide
44	F8	24	GLY	Peptide

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Mol	Chain	Res	Type	Group
49	G5	15	LYS	Peptide
49	G5	17	SER	Peptide
49	G5	43	GLN	Peptide
45	G8	5	MET	Peptide
45	G8	53	PRO	Peptide
45	G8	54	LYS	Peptide
46	H8	59	LEU	Peptide
46	H8	63	ASP	Peptide
51	I5	26	SER	Peptide
51	I5	36	CYS	Peptide
47	I8	83	PRO	Peptide
47	I8	9	SER	Peptide
48	J8	75	GLU	Peptide
48	J8	85	LEU	Peptide
53	K5	16	CYS	Peptide
53	K5	43	CYS	Peptide
53	K5	44	ARG	Peptide
49	K8	17	SER	Peptide
49	K8	46	GLN	Peptide
49	K8	5	GLU	Peptide
55	M5	30	ARG	Peptide
55	M5	33	ASN	Peptide
55	M5	40	GLU	Peptide
51	M8	38	LYS	Peptide
51	M8	40	HIS	Peptide
52	N8	41	PRO	Peptide
52	N8	5	PRO	Peptide
53	O8	15	GLU	Peptide
53	O8	27	LYS	Peptide
53	O8	44	ARG	Peptide
55	Q8	18	ALA	Peptide
55	Q8	19	SER	Peptide
55	Q8	27	THR	Peptide
55	Q8	33	ASN	Peptide
55	Q8	37	SER	Peptide
55	Q8	49	VAL	Peptide
55	Q8	51	ALA	Peptide
55	Q8	52	LYS	Peptide
55	Q8	56	GLU	Peptide
55	Q8	57	ARG	Peptide
55	Q8	6	THR	Peptide
55	Q8	7	HIS	Peptide

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Mol	Chain	Res	Type	Group
55	Q8	9	GLY	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	13	32207	0	16256	882	0
1	1G	32182	0	16244	875	0
2	12	1924	0	1975	85	0
2	1E	1924	0	1975	103	0
3	22	1612	0	1677	79	0
3	2E	1605	0	1668	65	0
4	32	1703	0	1763	94	0
4	3E	1703	0	1762	100	0
5	42	1155	0	1213	62	0
5	4E	1155	0	1213	54	0
6	52	843	0	857	44	0
6	5E	843	0	857	39	0
7	62	1257	0	1296	62	0
7	6E	1257	0	1296	49	0
8	72	1116	0	1177	57	0
8	7E	1116	0	1177	64	0
9	82	1010	0	1037	76	0
9	8E	1010	0	1037	61	0
10	1A	801	0	849	43	0
10	1I	801	0	849	62	0
11	2A	885	0	904	40	0
11	2I	885	0	904	44	0
12	3A	975	0	1062	47	0
12	3I	975	0	1062	43	0
13	4A	933	0	992	66	0
13	4I	938	0	997	50	0
14	5A	476	0	511	30	0
14	5I	492	0	529	29	0
15	6A	734	0	771	25	0
15	6I	734	0	771	29	0
16	7A	705	0	725	43	0
16	7I	705	0	725	54	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
17	8A	834	0	904	25	0
17	8I	834	0	904	51	0
18	9A	591	0	662	27	0
18	9I	591	0	662	29	0
19	AA	624	0	636	47	0
19	AI	647	0	665	57	0
20	BA	763	0	861	45	0
20	BI	763	0	861	44	0
21	1B	217	0	234	11	0
21	1F	217	0	234	13	0
22	1K	1626	0	834	28	0
22	1L	1626	0	834	41	0
23	2K	1646	0	845	38	0
23	2L	1646	0	845	39	0
24	3K	1619	0	822	52	0
24	3L	1619	0	822	60	0
25	4K	301	0	153	11	0
25	4L	131	0	66	5	0
26	14	62647	0	31580	1604	1
26	1H	62707	0	31607	1779	1
27	16	2617	0	1328	71	0
27	1J	2617	0	1328	99	0
28	11	2120	0	2197	122	0
28	19	2120	0	2197	117	0
29	21	1568	0	1634	107	0
29	29	1568	0	1634	117	0
30	31	1585	0	1632	104	0
30	39	1627	0	1680	101	0
31	41	1473	0	1535	77	0
31	49	1473	0	1535	70	0
32	51	1336	0	1418	94	0
32	59	1307	0	1382	73	0
33	61	1136	0	1223	63	0
33	69	1136	0	1223	56	0
34	15	1104	0	1180	55	0
34	58	1104	0	1180	70	0
35	25	933	0	996	48	0
35	68	933	0	996	41	0
36	35	1145	0	1228	113	0
36	78	1145	0	1228	119	0
37	45	1122	0	1179	84	0
37	88	1087	0	1129	83	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
38	55	960	0	1021	65	0
38	98	968	0	1033	72	0
39	65	882	0	943	62	0
39	A8	882	0	943	67	0
40	75	1141	0	1202	67	0
40	B8	1141	0	1202	76	0
41	85	964	0	1022	54	0
41	C8	964	0	1022	65	0
42	95	779	0	852	66	0
42	D8	779	0	852	40	0
43	A5	900	0	964	34	0
43	E8	900	0	964	50	0
44	B5	730	0	780	32	0
44	F8	742	0	803	56	0
45	C5	794	0	884	58	0
45	G8	791	0	882	57	0
46	D5	1428	0	1454	88	0
46	H8	1397	0	1430	85	0
47	E5	613	0	633	48	0
47	I8	627	0	642	34	0
48	F5	763	0	848	32	0
48	J8	763	0	848	38	0
49	G5	558	0	610	32	1
49	K8	563	0	612	41	1
50	H5	469	0	518	16	0
50	L8	452	0	503	33	0
51	I5	515	0	514	40	0
51	M8	533	0	526	48	0
52	J5	459	0	480	32	0
52	N8	454	0	475	33	0
53	K5	389	0	404	21	0
53	O8	389	0	404	31	0
54	L5	398	0	441	19	0
54	P8	418	0	467	22	0
55	M5	477	0	540	47	0
55	Q8	480	0	549	132	0
56	11	4	0	0	0	0
56	13	151	0	0	0	0
56	14	389	0	0	0	0
56	16	13	0	0	0	0
56	1G	103	0	0	0	0
56	1H	529	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
56	1J	9	0	0	0	0
56	1K	2	0	0	0	0
56	2I	2	0	0	0	0
56	29	2	0	0	0	0
56	2K	3	0	0	0	0
56	2L	2	0	0	0	0
56	3I	1	0	0	0	0
56	35	1	0	0	0	0
56	3E	2	0	0	0	0
56	3I	1	0	0	0	0
56	4I	2	0	0	0	0
56	49	1	0	0	0	0
56	4A	1	0	0	0	0
56	5E	1	0	0	0	0
56	6A	1	0	0	0	0
56	75	1	0	0	0	0
56	78	2	0	0	0	0
56	85	1	0	0	0	0
56	88	2	0	0	0	0
56	98	1	0	0	0	0
56	AI	1	0	0	0	0
56	C5	1	0	0	0	0
56	C8	1	0	0	0	0
56	E8	1	0	0	0	0
56	F5	1	0	0	0	0
56	I8	2	0	0	0	0
56	J8	2	0	0	0	0
56	L8	1	0	0	0	0
56	P8	1	0	0	0	0
56	Q8	1	0	0	0	0
57	32	1	0	0	0	0
57	3E	1	0	0	0	0
57	5A	1	0	0	0	0
57	5I	1	0	0	0	0
57	C5	1	0	0	0	0
57	G8	1	0	0	0	0
58	11	9	0	0	0	0
58	13	212	0	0	36	0
58	14	730	0	0	169	0
58	16	16	0	0	2	0
58	19	11	0	0	0	0
58	1G	99	0	0	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
58	1H	1097	0	0	264	0
58	1I	2	0	0	0	0
58	1J	12	0	0	3	0
58	1K	1	0	0	0	0
58	2I	5	0	0	3	0
58	29	6	0	0	0	0
58	3I	10	0	0	0	0
58	39	7	0	0	0	0
58	3E	1	0	0	0	0
58	3I	1	0	0	0	0
58	3K	1	0	0	0	0
58	4E	2	0	0	0	0
58	4K	4	0	0	0	0
58	55	1	0	0	0	0
58	58	1	0	0	0	0
58	5I	2	0	0	0	0
58	78	4	0	0	0	0
58	7A	2	0	0	0	0
58	85	1	0	0	0	0
58	8E	2	0	0	0	0
58	98	1	0	0	0	0
58	A5	1	0	0	0	0
58	BA	1	0	0	0	0
58	C8	2	0	0	1	0
58	D8	1	0	0	0	0
58	E8	2	0	0	0	0
58	G8	3	0	0	0	0
58	H5	1	0	0	0	0
58	I8	6	0	0	0	0
58	L5	1	0	0	0	0
58	L8	2	0	0	0	0
58	P8	2	0	0	0	0
58	Q8	2	0	0	0	0
All	All	300537	0	200491	9898	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 21.

All (9898) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:32:9:CYS:SG	4:32:31:CYS:HB2	1.81	1.19
26:1H:2714:G:OP2	58:1H:3670:HOH:O	1.68	1.11
26:14:2593:U:O4	58:14:4084:HOH:O	1.72	1.07
36:78:19:VAL:HG12	36:78:21:ARG:H	1.15	1.07
26:1H:1614:A:OP1	58:1H:3921:HOH:O	1.73	1.06
26:14:2685:G:O6	58:14:4089:HOH:O	1.75	1.04
26:14:450:G:O6	58:14:4019:HOH:O	1.74	1.04
26:1H:1359:A:N1	26:1H:1372:U:N3	2.06	1.03
26:14:274:G:H2'	26:14:275:G:H4'	1.41	1.02
26:1H:730:C:OP2	58:1H:3689:HOH:O	1.76	1.02
55:Q8:46:ARG:HH21	55:Q8:48:PHE:HB3	1.19	1.01
26:1H:585:G:OP2	58:1H:4547:HOH:O	1.78	1.00
26:14:1582:C:HO2'	26:14:1586:A:H8	1.03	1.00
26:1H:810:U:OP1	58:1H:3713:HOH:O	1.77	1.00
46:H8:128:VAL:HA	46:H8:161:VAL:HG11	1.44	0.99
26:14:2505:G:O6	58:14:3717:HOH:O	1.81	0.99
26:14:2017:U:OP1	58:14:3577:HOH:O	1.80	0.99
26:14:1689:A:H62	26:14:1698:A:H2	1.09	0.98
39:65:85:VAL:H	39:65:110:LEU:HB3	1.29	0.97
26:14:676:A:H8	26:14:2069:G:H21	1.07	0.97
26:1H:2593:U:O4	58:1H:3683:HOH:O	1.81	0.97
26:1H:973:A:OP2	58:1H:4553:HOH:O	1.81	0.97
26:14:1614:A:OP1	58:14:3409:HOH:O	1.83	0.97
26:14:1771:C:HO2'	26:14:1786:A:H8	0.98	0.96
26:1H:563:G:OP2	58:1H:3636:HOH:O	1.83	0.95
2:1E:104:ASN:HD21	2:1E:107:THR:HB	1.29	0.95
26:14:568:U:O4	58:14:4004:HOH:O	1.81	0.95
26:1H:135:G:N7	58:1H:4152:HOH:O	1.97	0.94
26:1H:1601:G:N7	58:1H:4021:HOH:O	1.98	0.94
26:1H:2017:U:OP1	58:1H:4641:HOH:O	1.83	0.94
26:1H:2577:A:OP1	58:1H:3800:HOH:O	1.83	0.94
1:13:1348:U:H3	1:13:1374:A:H2	1.09	0.94
26:1H:2582:G:OP2	58:1H:3812:HOH:O	1.84	0.94
26:1H:511:U:OP2	58:1H:4662:HOH:O	1.85	0.94
26:14:517:C:OP1	52:J5:16:ARG:NH2	2.00	0.94
27:1J:80:U:H2'	27:1J:81:G:H21	1.30	0.94
26:1H:2248:C:OP2	58:1H:3722:HOH:O	1.85	0.93
26:1H:586:A:OP2	58:1H:3885:HOH:O	1.84	0.93
26:1H:2705:A:OP2	58:1H:4669:HOH:O	1.86	0.93
26:1H:751:A:OP1	58:1H:3923:HOH:O	1.85	0.92
45:C5:97:ARG:NH1	45:C5:104:GLY:O	2.02	0.92
1:13:1502:A:H2	1:13:1505:G:H1	1.11	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2298:A:H62	26:1H:2318:G:H8	1.14	0.92
26:1H:2308:G:H1	26:1H:2311:A:H2	1.10	0.92
1:13:601:C:H2'	1:13:602:A:H8	1.35	0.92
2:1E:185:ILE:HG22	2:1E:199:TYR:HB2	1.50	0.92
26:1H:1327:C:OP2	58:1H:3645:HOH:O	1.88	0.92
26:1H:450:G:O6	58:1H:3896:HOH:O	1.88	0.91
26:1H:567:A:OP1	58:1H:3606:HOH:O	1.88	0.91
41:C8:8:VAL:HG23	41:C8:11:ARG:HH21	1.32	0.91
32:59:159:GLU:O	32:59:163:TYR:OH	1.86	0.91
36:35:19:VAL:HG13	36:35:21:ARG:H	1.35	0.91
26:14:741:G:OP1	58:14:4071:HOH:O	1.88	0.91
26:1H:1771:C:HO2'	26:1H:1786:A:H8	0.94	0.91
26:1H:741:G:OP1	58:1H:3970:HOH:O	1.87	0.91
26:14:323:G:HO2'	26:14:1205:U:H3	0.97	0.91
26:1H:2271:G:N7	58:1H:4245:HOH:O	2.02	0.91
26:1H:2334:G:O6	47:I8:74:ARG:NH2	2.03	0.90
26:1H:1622:G:OP2	58:1H:4308:HOH:O	1.88	0.90
26:14:1043:C:N3	26:14:1112:G:N2	2.20	0.90
28:11:17:THR:HG22	28:11:205:VAL:H	1.36	0.90
26:14:2720:U:H3	26:14:2873:A:H2	1.19	0.90
26:14:882:G:H22	26:14:894:C:H42	1.20	0.90
13:4I:23:TYR:HB3	13:4I:67:GLU:HB2	1.54	0.90
1:13:1128:C:O2	1:13:1146:A:N6	2.04	0.89
31:41:21:ARG:HG2	31:41:21:ARG:HH11	1.37	0.89
26:14:1616:A:O2'	58:14:3542:HOH:O	1.88	0.89
42:95:85:LYS:HD2	42:95:86:GLY:H	1.36	0.89
26:1H:1021:A:H2'	26:1H:1023:U:H5'	1.55	0.89
37:88:82:ARG:HD2	37:88:82:ARG:N	1.85	0.89
26:1H:1997:G:OP2	58:1H:4005:HOH:O	1.89	0.89
54:P8:9:ARG:HH21	54:P8:47:ARG:HD2	1.35	0.89
40:B8:26:ASP:HB3	40:B8:92:GLY:H	1.38	0.89
26:1H:399:G:OP2	58:1H:4045:HOH:O	1.90	0.89
26:1H:730:C:H3'	58:1H:3687:HOH:O	1.71	0.89
26:1H:2249:U:O4	58:1H:3722:HOH:O	1.91	0.89
1:13:1506:U:O2'	58:13:1803:HOH:O	1.89	0.89
1:1G:1435:G:H2'	1:1G:1436:U:C6	2.07	0.88
26:1H:1639:U:OP1	58:1H:3675:HOH:O	1.91	0.88
26:1H:1187:G:OP2	58:1H:3882:HOH:O	1.90	0.88
3:2E:60:ALA:H	3:2E:63:ASN:HB3	1.38	0.88
46:H8:30:ASN:HD22	46:H8:90:VAL:HB	1.38	0.88
1:1G:1502:A:H2	1:1G:1505:G:H1	1.20	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:235:C:H5'	17:8I:70:ARG:HG2	1.55	0.88
26:1H:330:A:HO2'	26:1H:331:A:H8	0.93	0.88
1:13:737:A:H2'	1:13:738:C:H6	1.35	0.88
26:14:2701:C:H3'	26:14:2702:U:H5''	1.54	0.88
26:1H:1778:U:H2'	26:1H:1784:A:N6	1.88	0.88
26:1H:265:A:H8	26:1H:266:G:H1'	1.39	0.88
26:1H:2407:G:OP1	58:1H:4630:HOH:O	1.92	0.88
26:1H:1327:C:OP2	58:1H:3643:HOH:O	1.91	0.88
1:13:74:C:H42	1:13:96:G:H1	1.19	0.88
26:1H:620:G:H4'	26:1H:621:A:H5''	1.54	0.87
26:1H:846:C:O2'	58:1H:3757:HOH:O	1.90	0.87
26:1H:2656:U:H3	26:1H:2665:A:H2	1.22	0.87
7:62:70:LYS:HG2	7:62:96:GLN:HB3	1.57	0.87
4:3E:107:ARG:HH22	4:3E:194:LEU:HD22	1.38	0.87
1:1G:21:G:OP1	58:1G:1836:HOH:O	1.92	0.87
26:1H:2533:A:OP2	58:1H:4420:HOH:O	1.92	0.87
26:1H:510:C:OP1	58:1H:4662:HOH:O	1.90	0.87
1:13:352:C:OP1	58:13:1899:HOH:O	1.91	0.87
26:14:1997:G:OP2	58:14:3494:HOH:O	1.91	0.87
14:5I:29:ARG:HH21	14:5I:41:ARG:HH12	1.18	0.87
41:85:92:ARG:HD3	41:85:94:ASN:HB3	1.57	0.87
26:14:2035:G:OP1	58:14:3603:HOH:O	1.92	0.87
26:1H:882:G:H22	26:1H:894:C:H42	1.20	0.86
26:14:654(D):G:H22	26:14:654(Q):C:H42	1.23	0.86
26:1H:1453:A:OP2	58:1H:4658:HOH:O	1.92	0.86
26:14:446:G:OP2	58:14:3695:HOH:O	1.93	0.86
1:13:768:A:OP2	58:13:1840:HOH:O	1.92	0.86
26:1H:748:G:OP2	58:1H:4460:HOH:O	1.92	0.86
37:88:12:GLN:HE21	37:88:72:LYS:HG3	1.38	0.86
1:1G:521:G:O6	1:1G:528:C:N4	2.07	0.86
13:4A:81:LEU:HD13	13:4A:88:ARG:HD2	1.56	0.86
26:14:751:A:OP1	58:14:3408:HOH:O	1.94	0.86
26:14:2431:U:OP1	58:14:3415:HOH:O	1.94	0.86
26:1H:945:A:OP1	58:1H:4113:HOH:O	1.94	0.86
29:21:105:THR:OG1	29:21:199:ARG:NH2	2.08	0.86
26:14:1019:U:H3	26:14:1142(A):A:H62	1.24	0.86
26:14:249:C:OP1	58:14:3411:HOH:O	1.94	0.86
26:14:1828:G:OP1	58:14:3449:HOH:O	1.94	0.86
2:12:42:ILE:HD11	2:12:202:PRO:HB2	1.57	0.85
26:14:662:G:H5'	36:35:15:ARG:HA	1.57	0.85
26:14:1330:C:OP1	58:14:3641:HOH:O	1.94	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:51:83:TYR:HB3	32:51:135:GLY:H	1.40	0.85
26:1H:1253:A:N7	58:1H:3713:HOH:O	2.08	0.85
42:95:71:LEU:N	42:95:86:GLY:HA2	1.91	0.85
1:13:1003:G:N2	1:13:1037:C:N3	2.23	0.85
26:1H:2432:A:OP2	58:1H:3905:HOH:O	1.92	0.85
26:14:2588:G:OP1	58:14:3471:HOH:O	1.94	0.85
39:A8:34:HIS:HB2	39:A8:36:TYR:HE1	1.41	0.85
17:8I:76:LEU:HD12	17:8I:77:VAL:H	1.42	0.85
1:13:21:G:OP1	58:13:1836:HOH:O	1.93	0.85
27:1J:15:A:H5'	27:1J:16:G:H8	1.42	0.85
26:14:1327:C:OP2	58:14:3551:HOH:O	1.93	0.85
26:14:1664:A:OP2	58:14:3497:HOH:O	1.91	0.85
26:14:400:G:N7	58:14:4115:HOH:O	2.10	0.85
26:1H:2608:G:N7	58:1H:3814:HOH:O	2.08	0.85
26:1H:860:U:H5	26:1H:917:A:C2	1.94	0.85
26:1H:2035:G:OP1	58:1H:3793:HOH:O	1.94	0.85
26:1H:574:C:OP1	58:1H:4544:HOH:O	1.94	0.85
11:2A:29:ILE:HG22	11:2A:44:SER:HB2	1.58	0.85
26:1H:1689:A:H62	26:1H:1698:A:H2	1.24	0.85
14:5A:21:TYR:HE1	14:5A:23:ARG:HE	1.23	0.85
55:Q8:49:VAL:HB	55:Q8:52:LYS:O	1.77	0.84
26:1H:945:A:OP2	58:1H:4578:HOH:O	1.93	0.84
26:1H:442:G:H1'	30:31:48:THR:HG21	1.58	0.84
55:M5:52:LYS:HZ3	55:M5:53:PRO:HA	1.40	0.84
1:13:1157:A:H61	1:13:1178:G:H21	1.25	0.84
26:1H:1828:G:OP1	58:1H:4688:HOH:O	1.93	0.84
26:1H:2502:G:OP2	58:1H:3629:HOH:O	1.92	0.84
55:Q8:49:VAL:HG21	55:Q8:53:PRO:HD2	1.59	0.84
26:1H:1639:U:OP2	58:1H:4526:HOH:O	1.94	0.84
1:1G:1129:C:H41	1:1G:1141:C:H42	1.25	0.84
26:14:848:G:H2'	26:14:849:A:H8	1.42	0.84
1:1G:396:G:O2'	1:1G:398:C:OP1	1.94	0.84
1:13:812:C:N3	58:13:1809:HOH:O	2.09	0.84
1:1G:975:A:H4'	1:1G:976:G:H5''	1.59	0.84
26:14:1327:C:OP2	58:14:3556:HOH:O	1.95	0.84
32:51:150:ALA:O	32:51:153:LYS:NZ	2.10	0.84
26:1H:1164:G:H2'	26:1H:1165:U:C6	2.12	0.84
26:14:1845:G:N7	58:14:3856:HOH:O	2.10	0.84
29:29:81:ILE:HG22	29:29:82:ARG:H	1.43	0.84
26:14:907:U:O2'	37:45:101:ARG:NH2	2.10	0.84
36:35:39:LYS:HD2	36:35:45:LEU:HD21	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1434:A:H61	26:1H:1558:A:N6	1.76	0.84
30:39:116:ASP:OD2	36:35:1:MET:N	2.09	0.84
26:1H:2068:U:H3	26:1H:2430:A:H2	1.22	0.84
26:14:910:A:H62	37:45:12:GLN:HA	1.41	0.84
8:7E:34:GLU:OE1	8:7E:37:ARG:NH1	2.10	0.84
26:1H:2452:C:OP1	58:1H:4435:HOH:O	1.95	0.84
26:1H:780:G:H21	26:1H:783:A:H62	1.20	0.84
26:1H:2400:G:H2'	26:1H:2401:U:H6	1.43	0.84
12:3I:32:PHE:HB3	12:3I:84:LEU:HD11	1.60	0.83
26:14:1037:G:O6	26:14:1118:C:N4	2.11	0.83
26:1H:577:G:O6	58:1H:4650:HOH:O	1.96	0.83
26:1H:574:C:OP2	58:1H:4652:HOH:O	1.94	0.83
26:1H:1525:G:H2'	26:1H:1526:G:H8	1.43	0.83
26:14:1622:G:OP2	58:14:3866:HOH:O	1.95	0.83
28:11:60:ARG:HD3	28:11:86:PRO:HB2	1.58	0.83
28:11:93:ALA:HB3	28:11:105:ILE:HG22	1.58	0.83
39:A8:7:TYR:HA	39:A8:10:ARG:HH21	1.43	0.83
26:1H:1496:A:H8	26:1H:1577:C:HO2'	1.22	0.83
6:5E:50:TYR:OH	18:9I:74:ARG:O	1.94	0.83
26:1H:226:G:H21	26:1H:228:A:H2	1.26	0.83
26:14:1382:G:N7	58:14:3691:HOH:O	2.11	0.83
26:14:2267:A:OP2	58:14:3824:HOH:O	1.94	0.83
26:1H:1019:U:HO2'	26:1H:1021:A:H2	1.27	0.83
9:82:128:ARG:NH2	23:2L:34:U:OP2	2.11	0.83
26:1H:1332:G:OP1	58:1H:4561:HOH:O	1.96	0.83
49:K8:50:ILE:HD12	49:K8:51:ARG:H	1.43	0.83
26:1H:450:G:OP2	58:1H:3893:HOH:O	1.96	0.83
1:13:664:G:H22	1:13:741:G:H1	1.26	0.83
26:1H:2781:A:H5''	26:1H:2782:G:H5'	1.60	0.83
1:13:975:A:H4'	1:13:976:G:H5''	1.58	0.83
26:1H:1249:U:OP1	58:1H:3886:HOH:O	1.97	0.83
41:85:66:ASN:HB2	41:85:76:TYR:HB2	1.61	0.83
1:13:606:G:N3	1:13:632:A:N6	2.27	0.83
26:1H:1794:U:H2'	26:1H:1795:C:H6	1.43	0.83
26:1H:1658:C:OP1	58:1H:4531:HOH:O	1.95	0.83
26:14:2331:G:H4'	47:E5:43:THR:H	1.44	0.83
30:31:6:VAL:N	30:31:24:LEU:O	2.12	0.83
1:13:144:G:N2	1:13:178:C:O2	2.09	0.83
26:1H:2056:G:OP2	58:1H:3624:HOH:O	1.96	0.83
27:1J:40:U:O2	27:1J:45:A:N6	2.10	0.83
26:1H:2447:G:OP2	58:1H:3853:HOH:O	1.96	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:963:U:OP1	58:1H:3857:HOH:O	1.96	0.83
18:9I:59:SER:HB3	18:9I:62:GLU:HB2	1.61	0.82
36:35:85:LEU:HA	36:35:88:LEU:HB3	1.59	0.82
1:13:1145:C:H4'	1:13:1146:A:H5'	1.61	0.82
26:14:94:G:N3	49:G5:47:ASN:ND2	2.27	0.82
1:13:510:A:OP2	4:3E:49:ARG:NH2	2.11	0.82
40:B8:6:LEU:HA	40:B8:9:LEU:HB2	1.61	0.82
26:14:848:G:H2'	26:14:849:A:C8	2.13	0.82
26:1H:1900:A:H5'	26:1H:1900:A:H8	1.45	0.82
1:13:972:C:OP1	58:13:1830:HOH:O	1.98	0.82
1:1G:827:U:H3	1:1G:872:A:H62	1.27	0.82
46:D5:111:VAL:HB	46:D5:143:GLY:HA3	1.61	0.82
26:1H:676:A:H8	26:1H:2069:G:H21	1.27	0.82
26:1H:411:G:OP1	58:1H:4630:HOH:O	1.98	0.82
41:85:92:ARG:HD2	41:85:95:LEU:HD12	1.60	0.82
1:13:524:G:H2'	1:13:525:C:C6	2.15	0.82
8:7E:120:THR:H	8:7E:123:GLU:HB2	1.43	0.82
24:3K:76:A:H8	26:1H:2394:C:H42	1.25	0.82
54:L5:35:ARG:HG3	54:L5:42:LEU:HD11	1.59	0.82
1:1G:957:U:O2'	1:1G:959:A:N7	2.12	0.82
26:14:2499:C:OP1	58:14:4009:HOH:O	1.96	0.82
49:G5:25:VAL:HG12	49:G5:60:LEU:HD23	1.60	0.82
46:H8:76:LEU:HD22	46:H8:76:LEU:H	1.44	0.82
1:1G:998:G:N2	1:1G:1043:C:N3	2.27	0.82
26:14:1041:C:H42	26:14:1114:G:H22	1.25	0.82
26:1H:910:A:N7	37:88:13:GLN:HG3	1.94	0.82
33:69:75:LEU:HD22	33:69:76:THR:H	1.44	0.82
26:14:2127:G:O6	26:14:2161:C:N4	2.13	0.82
32:51:30:LYS:HE3	32:51:81:GLU:H	1.45	0.82
26:1H:1410:G:O6	26:1H:1592:C:N4	2.09	0.82
24:3L:52:G:H1	24:3L:62:C:H42	1.26	0.81
26:1H:2033:A:OP1	58:1H:4060:HOH:O	1.98	0.81
46:D5:5:LEU:HG	46:D5:47:VAL:HG21	1.61	0.81
1:13:736:C:H2'	1:13:737:A:C8	2.14	0.81
26:1H:2502:G:N7	58:1H:3855:HOH:O	2.13	0.81
37:88:14:ARG:HG2	37:88:41:TRP:HH2	1.45	0.81
1:1G:976:G:N2	1:1G:1362(A):C:OP2	2.09	0.81
26:14:1022:G:O2'	26:14:1024:G:O6	1.96	0.81
26:14:2264:C:N4	47:E5:15:ASP:OD2	2.11	0.81
26:14:631:A:OP2	55:M5:47:LYS:NZ	2.10	0.81
26:1H:624:C:OP1	58:1H:4303:HOH:O	1.97	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:69:102:SER:O	33:69:106:GLY:N	2.13	0.81
26:1H:602:G:HO2'	26:1H:604:G:HO2'	1.24	0.81
2:1E:21:ARG:HB2	2:1E:39:ILE:HA	1.61	0.81
1:1G:1162:C:H42	1:1G:1174:G:H1	1.29	0.81
30:31:29:ASN:H	30:31:112:MET:HE1	1.45	0.81
26:14:1309:G:N7	58:14:3847:HOH:O	2.12	0.81
46:H8:108:PRO:HB2	46:H8:112:ARG:HA	1.62	0.81
26:1H:2712(A):A:OP2	58:1H:3670:HOH:O	1.97	0.81
26:1H:187:G:OP2	58:1H:4410:HOH:O	1.98	0.81
26:14:958:U:OP2	37:45:14:ARG:NH1	2.13	0.81
26:14:1970:A:OP1	58:14:3455:HOH:O	1.98	0.81
26:14:2873:A:H8	38:55:6:SER:H	1.29	0.81
1:13:1110:A:OP2	58:13:1960:HOH:O	1.98	0.81
1:1G:673:G:H2'	1:1G:674:G:C8	2.16	0.81
48:J8:83:GLU:HG2	48:J8:85:LEU:H	1.46	0.81
55:Q8:53:PRO:HA	55:Q8:55:ALA:N	1.96	0.81
1:13:1004:A:O5'	1:13:1025:U:N3	2.12	0.81
1:1G:1452:C:H4'	1:1G:1453:G:H5'	1.63	0.81
26:14:1658:C:OP1	58:14:3490:HOH:O	1.99	0.81
7:62:20:ASP:HB3	7:62:23:VAL:HG23	1.62	0.81
26:1H:1828:G:OP2	58:1H:4634:HOH:O	1.97	0.81
26:1H:732:C:OP2	58:1H:4074:HOH:O	1.98	0.81
30:39:5:ALA:HB1	30:39:125:LEU:HD21	1.63	0.81
36:78:47:ASP:OD2	36:78:50:ARG:NH2	2.14	0.81
1:13:352:C:O2'	1:13:354:G:OP1	1.98	0.81
45:G8:30:VAL:HG22	45:G8:37:VAL:HG12	1.63	0.81
1:13:36:C:OP1	12:3I:123:LYS:NZ	2.13	0.81
1:1G:475:G:OP1	16:7A:81:ARG:NH2	2.13	0.81
38:98:12:ARG:HD3	38:98:16:HIS:CG	2.15	0.81
19:AA:60:VAL:HG21	19:AA:74:PHE:HB3	1.63	0.81
1:13:1240:U:OP2	7:6E:116:ALA:N	2.13	0.81
26:1H:2256:G:N7	58:1H:4119:HOH:O	2.12	0.81
26:1H:1952:A:C5	35:68:22:ILE:HD11	2.15	0.80
38:98:12:ARG:HG2	38:98:12:ARG:HH11	1.44	0.80
19:AA:29:ARG:HH11	19:AA:48:THR:HG23	1.46	0.80
1:13:1027:C:N4	1:13:1032(A):G:O6	2.14	0.80
26:1H:2598:A:OP1	58:1H:3652:HOH:O	2.00	0.80
26:1H:2392:A:H2	26:1H:2424:C:H42	1.30	0.80
11:2A:32:ILE:HD11	11:2A:68:ALA:HB1	1.63	0.80
1:1G:411:A:H62	1:1G:413:G:H21	1.28	0.80
26:1H:1257:C:H4'	30:31:83:PHE:CD1	2.17	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2210:G:H5'	26:1H:2211:G:C8	2.17	0.80
26:14:259:G:H21	26:14:621:A:H8	1.29	0.80
28:11:146:GLU:HB2	28:11:189:CYS:HB3	1.63	0.80
26:1H:2584:U:H2'	26:1H:2585:U:H2'	1.64	0.80
8:7E:27:PRO:O	8:7E:32:LYS:NZ	2.15	0.80
27:1J:5:C:H42	27:1J:115:G:H1	1.30	0.80
26:14:2074:U:OP1	58:14:3403:HOH:O	1.98	0.80
26:1H:2713:A:OP2	58:1H:3673:HOH:O	2.00	0.80
26:1H:265:A:C8	26:1H:266:G:H1'	2.15	0.80
26:1H:1780:A:OP1	58:1H:3626:HOH:O	1.99	0.80
7:6E:111:ARG:NH1	7:6E:113:GLU:OE2	2.14	0.80
26:14:2392:A:H2	26:14:2424:C:H42	1.26	0.80
26:1H:945:A:OP1	58:1H:4115:HOH:O	2.00	0.80
21:1B:8:THR:HG22	21:1B:11:GLY:H	1.47	0.80
26:14:833:U:O2	36:35:55:ARG:NH1	2.14	0.80
26:1H:1484:G:N2	26:1H:1505:C:O2	2.14	0.80
26:14:2210:G:O5'	26:14:2211:G:N2	2.14	0.80
26:14:459:U:H5''	54:L5:40:TRP:CD2	2.17	0.79
26:1H:1152:C:H3'	58:1H:4003:HOH:O	1.80	0.79
1:1G:1298:C:OP2	7:62:114:ARG:NH2	2.15	0.79
1:13:504:C:OP1	58:13:1880:HOH:O	2.00	0.79
26:14:141:A:H8	26:14:1595:G:H21	1.28	0.79
28:19:69:ARG:NH2	28:19:128:GLY:O	2.15	0.79
26:14:1774:C:OP1	58:14:4078:HOH:O	1.98	0.79
36:78:19:VAL:HB	36:78:27:HIS:HB2	1.62	0.79
17:8I:76:LEU:HD11	17:8I:79:SER:HB3	1.64	0.79
1:13:659:U:H2'	1:13:660:G:C8	2.17	0.79
26:14:1678:G:H22	26:14:1989:G:H22	1.28	0.79
3:22:141:VAL:HA	3:22:144:SER:HB3	1.64	0.79
26:1H:1061:U:H4'	26:1H:1070:A:H1'	1.64	0.79
1:13:1500:A:OP1	58:13:1804:HOH:O	2.00	0.79
29:21:105:THR:HG22	29:21:106:GLY:H	1.44	0.79
44:B5:27:THR:HB	44:B5:80:ILE:HG22	1.63	0.79
30:39:78:ILE:HA	30:39:83:PHE:CD2	2.17	0.79
26:1H:1386:C:H2'	26:1H:1387:C:H6	1.47	0.79
35:25:115:VAL:HG13	35:25:121:VAL:HG21	1.63	0.79
26:14:453:C:OP1	58:14:4019:HOH:O	2.00	0.79
51:15:22:ILE:HG12	51:15:23:GLU:H	1.47	0.79
29:21:135:HIS:CD2	58:21:405:HOH:O	2.35	0.79
23:2K:62:C:H2'	23:2K:63:C:H6	1.47	0.79
36:78:29:LYS:HD3	36:78:30:THR:HG22	1.63	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:N8:40:LYS:HZ3	52:N8:46:CYS:HB3	1.47	0.79
30:39:25:PRO:HB2	30:39:27:GLU:H	1.47	0.79
20:BI:71:THR:HG22	20:BI:72:LEU:H	1.45	0.79
1:1G:1141:C:H2'	1:1G:1142:G:H8	1.47	0.79
1:13:538:G:H5''	12:3I:114:LYS:HB2	1.64	0.79
42:95:71:LEU:H	42:95:86:GLY:HA2	1.47	0.78
5:42:31:LEU:HA	5:42:45:PHE:HB2	1.66	0.78
26:14:2610:C:O2'	58:14:3716:HOH:O	1.97	0.78
2:1E:82:ARG:NE	2:1E:92:TYR:OH	2.16	0.78
33:61:39:ALA:HB1	33:61:44:LEU:HD13	1.64	0.78
29:29:66:HIS:NE2	29:29:73:GLU:OE1	2.16	0.78
26:1H:2593:U:H2'	26:1H:2594:C:H6	1.48	0.78
5:4E:33:VAL:HG11	5:4E:109:ILE:HG12	1.65	0.78
10:1I:22:LYS:NZ	10:1I:88:LEU:O	2.17	0.78
26:1H:1658:C:OP1	58:1H:4532:HOH:O	1.99	0.78
26:1H:2057:A:OP2	58:1H:3621:HOH:O	2.00	0.78
12:3A:60:LEU:HB2	12:3A:64:TYR:HB2	1.65	0.78
26:1H:801:G:OP2	58:1H:4182:HOH:O	1.99	0.78
20:BI:57:ARG:HH22	20:BI:102:GLY:HA2	1.49	0.78
34:58:65:LYS:HB3	34:58:69:GLN:HG3	1.64	0.78
30:31:167:ALA:HB1	30:31:173:VAL:HG11	1.65	0.78
26:1H:674:G:H1'	30:31:74:ARG:HD3	1.64	0.78
5:42:81:GLU:HG2	5:42:90:VAL:HG13	1.66	0.78
5:4E:110:LEU:HD13	5:4E:118:ILE:HD13	1.64	0.78
1:1G:961:U:O2	1:1G:1201:A:N6	2.16	0.78
30:31:29:ASN:N	30:31:112:MET:HE1	1.99	0.78
29:29:61:ARG:HA	29:29:63:LEU:HG	1.66	0.78
12:3A:59:ARG:NH2	12:3A:65:GLU:OE1	2.17	0.78
30:39:103:LYS:HA	30:39:106:ARG:HG3	1.66	0.78
26:1H:1643:G:N7	58:1H:3942:HOH:O	2.16	0.78
26:1H:1138:G:H21	34:58:106:MET:HE3	1.45	0.78
26:14:1858:G:O2'	26:14:1884:A:N6	2.16	0.78
26:1H:2431:U:OP2	58:1H:3908:HOH:O	2.01	0.78
1:1G:36:C:OP1	12:3A:123:LYS:NZ	2.16	0.78
1:13:1023:G:H3'	1:13:1024:G:H5''	1.64	0.78
26:1H:860:U:C5	26:1H:917:A:H2	2.02	0.78
31:41:105:LYS:HE2	51:M8:26:SER:HB2	1.65	0.78
5:42:33:VAL:HG21	5:42:109:ILE:HG12	1.63	0.78
26:1H:761:A:N7	58:1H:4073:HOH:O	2.17	0.78
26:14:1225:C:O3'	42:95:85:LYS:HA	1.83	0.78
1:13:736:C:H2'	1:13:737:A:H8	1.48	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2502:G:OP2	58:1H:3631:HOH:O	2.01	0.78
29:21:135:HIS:HE1	58:21:404:HOH:O	1.45	0.78
22:1L:74:C:N4	26:14:2554:U:O2	2.16	0.78
46:D5:27:VAL:HG12	46:D5:87:ASP:HB3	1.65	0.78
42:95:35:LEU:O	42:95:37:VAL:HG22	1.83	0.78
26:1H:1516:U:H2'	26:1H:1517:G:H8	1.48	0.78
26:14:273(C):C:H42	26:14:363(C):G:H1	1.29	0.78
26:1H:878:A:N6	26:1H:899:A:O2'	2.17	0.78
1:13:262:A:H2'	1:13:263:A:C8	2.19	0.78
26:1H:763:G:OP1	58:1H:3691:HOH:O	2.01	0.78
26:1H:510:C:OP1	58:1H:4667:HOH:O	2.02	0.78
9:82:9:ARG:HG2	9:82:14:VAL:HG22	1.66	0.78
41:85:28:ARG:NH1	41:85:38:THR:OG1	2.16	0.78
48:F5:24:ALA:HB3	48:F5:27:GLU:HG3	1.66	0.78
32:51:4:ILE:HD13	32:51:4:ILE:H	1.49	0.78
26:1H:67:U:H3	26:1H:74:A:H2	1.30	0.78
3:22:11:ARG:NH2	3:22:177:THR:O	2.17	0.78
1:13:1367:C:H5'	10:11:60:ARG:HH21	1.49	0.77
26:14:607:U:H3	26:14:621:A:H2	1.29	0.77
26:1H:2636:U:OP1	29:21:79:ARG:HA	1.84	0.77
7:62:79:ARG:HG2	7:62:84:ASN:HB3	1.66	0.77
26:1H:248:G:H5'	26:1H:250:G:N7	1.98	0.77
30:39:101:LEU:O	30:39:106:ARG:NH1	2.17	0.77
13:4A:57:ARG:HH12	51:I5:17:GLY:HA3	1.49	0.77
26:1H:1171:G:N2	26:1H:1178:C:N3	2.32	0.77
24:3L:18:G:H2'	24:3L:57:G:H1	1.46	0.77
39:65:3:ARG:HE	39:65:4:LEU:N	1.82	0.77
1:1G:560:U:H5'	1:1G:566:G:N2	1.99	0.77
44:F8:24:GLY:HA3	44:F8:82:GLN:HE22	1.50	0.77
46:H8:62:PRO:C	46:H8:64:GLY:HA2	2.05	0.77
26:14:2448:A:OP1	58:14:4009:HOH:O	2.03	0.77
26:14:1890:A:OP2	58:14:3914:HOH:O	2.02	0.77
26:14:660:G:H21	36:35:12:ALA:HB2	1.49	0.77
41:85:90:VAL:HA	42:95:39:LEU:HD23	1.64	0.77
26:1H:878:A:N6	26:1H:899:A:HO2'	1.82	0.77
26:1H:1727:U:H3	26:1H:1733:G:H1	1.32	0.77
26:1H:654(D):G:H1	26:1H:654(Q):C:H42	1.31	0.77
26:14:2415:G:H4'	36:35:67:MET:H	1.50	0.77
20:BA:50:GLU:HA	20:BA:100:ILE:HG21	1.67	0.77
5:4E:8:GLU:OE1	5:4E:63:ARG:NH2	2.17	0.77
46:H8:163:LEU:HB3	46:H8:165:VAL:H	1.50	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:324:G:N7	58:1G:1848:HOH:O	2.17	0.77
26:1H:2057:A:OP2	58:1H:3624:HOH:O	2.01	0.77
40:75:1:MET:HB3	40:75:5:ALA:HB3	1.66	0.77
1:1G:1310:G:H5'	13:4A:77:ASN:HD21	1.48	0.77
13:4A:82:MET:HG2	13:4A:83:ASP:H	1.49	0.77
19:AI:15:LEU:HB2	19:AI:31:ILE:HD11	1.66	0.77
17:8I:88:TYR:HD1	17:8I:89:LEU:HD23	1.48	0.77
26:1H:987:G:OP2	58:1H:3995:HOH:O	2.02	0.77
28:19:60:ARG:NH1	28:19:86:PRO:O	2.17	0.77
33:61:73:GLU:HG3	33:61:136:VAL:HG23	1.67	0.77
1:1G:999:U:H2'	1:1G:1000:A:H8	1.50	0.77
1:1G:1149:C:OP2	9:82:9:ARG:NH1	2.18	0.77
9:82:112:LYS:HA	9:82:119:ALA:HB2	1.66	0.77
30:31:66:PRO:O	30:31:67:GLN:HB3	1.84	0.77
26:14:2250:G:C4	37:45:82:ARG:HG3	2.20	0.77
46:H8:7:ALA:HB2	46:H8:59:LEU:HD22	1.67	0.77
1:13:177:C:OP1	20:BI:65:LYS:NZ	2.16	0.77
37:45:26:TYR:O	37:45:28:ALA:N	2.17	0.77
26:14:240:G:O6	58:14:3613:HOH:O	2.03	0.77
29:29:12:THR:O	29:29:23:VAL:HG23	1.85	0.77
43:E8:14:PRO:HB3	43:E8:18:ARG:HH12	1.49	0.77
26:14:1970:A:OP2	58:14:3458:HOH:O	2.03	0.76
1:1G:179:A:OP2	58:1G:1895:HOH:O	2.03	0.76
1:1G:500:G:N2	1:1G:545:C:O2	2.18	0.76
26:14:1676:A:OP2	58:14:3433:HOH:O	2.02	0.76
26:1H:2685:G:N7	58:1H:4089:HOH:O	2.18	0.76
26:14:140:A:H8	26:14:1408:C:HO2'	1.27	0.76
1:1G:353:A:H8	1:1G:353:A:H5'	1.50	0.76
26:14:802:A:OP1	58:14:3908:HOH:O	2.02	0.76
13:4A:97:PRO:HA	13:4A:110:ARG:HD3	1.67	0.76
1:1G:1028(A):C:O2	1:1G:1033:G:N2	2.17	0.76
26:1H:2785:C:OP1	29:21:41:LYS:NZ	2.16	0.76
26:14:1664:A:OP1	58:14:3504:HOH:O	2.04	0.76
26:14:2207:C:H42	26:14:2217:G:H1	1.33	0.76
39:65:106:ARG:NH1	39:65:107:GLU:OE2	2.18	0.76
1:1G:588:G:H1	1:1G:651:C:H42	1.34	0.76
27:16:42:C:O3'	31:41:67:LYS:NZ	2.16	0.76
46:H8:19:ARG:NH1	46:H8:84:GLU:O	2.19	0.76
26:14:152:G:H1	26:14:174:C:H42	1.30	0.76
1:13:601:C:H2'	1:13:602:A:C8	2.19	0.76
46:D5:44:PHE:HE2	46:D5:86:VAL:HG11	1.51	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:69:130:TYR:HB3	33:69:136:VAL:HG13	1.67	0.76
26:1H:2718:G:N7	58:1H:4512:HOH:O	2.17	0.76
1:1G:426:G:OP1	4:32:36:ARG:NH2	2.18	0.76
1:1G:979:C:H3'	1:1G:980:C:H5''	1.67	0.76
26:14:629:G:O6	26:14:634:C:N4	2.17	0.76
26:14:1048:A:N6	26:14:1112:G:O2'	2.18	0.76
26:1H:1557:C:OP2	26:1H:1558:A:O2'	2.04	0.76
26:14:509:C:OP1	58:14:3884:HOH:O	2.04	0.76
26:14:397:G:N7	58:14:3932:HOH:O	2.19	0.76
26:14:2287:A:H62	26:14:2344:U:H3	1.32	0.76
26:1H:761:A:OP1	58:1H:3689:HOH:O	2.03	0.76
26:14:630:G:N2	26:14:633:A:OP2	2.16	0.76
26:1H:2759:G:OP2	58:1H:4474:HOH:O	2.02	0.76
26:14:578:A:OP2	58:14:3576:HOH:O	2.02	0.76
26:14:1060:U:H4'	26:14:1061:U:H5''	1.68	0.76
26:14:2120:G:N2	26:14:2178:C:N3	2.34	0.76
26:1H:1701:A:OP2	58:1H:4389:HOH:O	2.04	0.76
26:1H:2028:U:O4	58:1H:4060:HOH:O	2.03	0.76
28:11:223:GLY:HA3	28:11:231:HIS:ND1	2.01	0.76
1:13:692:U:O2'	1:13:694:A:N7	2.19	0.76
31:49:41:GLN:NE2	31:49:154:GLY:O	2.19	0.76
29:21:111:ARG:HD3	29:21:160:TYR:CD2	2.21	0.76
1:1G:1074:G:O2'	1:1G:1101:A:N1	2.18	0.76
4:3E:9:CYS:HB3	4:3E:32:ALA:HB2	1.67	0.76
55:Q8:30:ARG:HG3	55:Q8:30:ARG:HH11	1.51	0.76
55:Q8:49:VAL:HG12	55:Q8:52:LYS:HB2	1.68	0.76
26:14:2375:G:N7	58:14:3706:HOH:O	2.17	0.76
26:1H:1952:A:C6	35:68:22:ILE:HD11	2.20	0.76
29:21:131:ALA:HB1	58:21:404:HOH:O	1.86	0.76
26:14:2786:U:O2'	29:29:62:PRO:O	2.02	0.76
3:22:152:ILE:HG12	3:22:167:TRP:HB2	1.67	0.76
26:14:2114:A:N6	26:14:2119:A:N7	2.34	0.76
35:25:73:ASP:OD2	40:75:32:TYR:OH	2.04	0.76
26:14:2272:U:O4	58:14:3822:HOH:O	2.00	0.76
1:13:1366:C:O2'	10:1I:60:ARG:NH2	2.20	0.75
26:1H:862:G:OP2	58:1H:3993:HOH:O	2.02	0.75
19:AI:78:ARG:HE	19:AI:78:ARG:H	1.28	0.75
1:1G:1238:A:N3	1:1G:1241:G:O2'	2.19	0.75
1:1G:195:A:OP2	58:1G:1894:HOH:O	2.04	0.75
26:1H:270(E):G:H1	26:1H:270(U):C:H42	1.32	0.75
33:61:110:ASP:OD1	33:61:110:ASP:N	2.19	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:272:C:H2'	1:13:273:A:H8	1.51	0.75
1:1G:156:G:N2	1:1G:165:C:O2	2.16	0.75
26:1H:1478:G:H2'	26:1H:1479:G:H8	1.49	0.75
26:14:2782:G:OP2	58:14:3786:HOH:O	2.03	0.75
1:13:1306:A:H61	1:13:1331:G:H1'	1.52	0.75
26:1H:1676:A:OP2	58:1H:3708:HOH:O	2.04	0.75
11:2A:48:ILE:HD11	11:2A:64:ALA:HA	1.68	0.75
34:15:56:ASN:HA	34:15:125:GLY:H	1.52	0.75
26:1H:1381:G:N7	58:1H:4562:HOH:O	2.18	0.75
11:2I:34:ASP:HB3	11:2I:40:ILE:HD11	1.68	0.75
4:3E:25:ARG:NH1	4:3E:30:LYS:O	2.20	0.75
26:1H:330:A:O2'	26:1H:331:A:H8	1.69	0.75
26:1H:1314:C:OP1	58:1H:4561:HOH:O	2.03	0.75
34:58:96:GLU:HG2	34:58:97:ARG:N	1.99	0.75
26:1H:883:G:O6	26:1H:893:C:N4	2.18	0.75
1:1G:547:A:OP1	58:1G:1804:HOH:O	2.05	0.75
31:49:56:ALA:HA	31:49:59:GLU:HB2	1.69	0.75
1:1G:768:A:OP2	58:1G:1811:HOH:O	2.03	0.75
16:7A:53:VAL:HG13	16:7A:79:VAL:HG22	1.67	0.75
31:49:161:THR:HG22	31:49:163:ALA:H	1.51	0.75
26:1H:620:G:H4'	26:1H:621:A:C5'	2.17	0.75
26:1H:1900:A:C8	26:1H:1900:A:H5'	2.22	0.75
26:1H:1153:C:OP2	58:1H:4000:HOH:O	2.03	0.75
26:1H:2709:G:O2'	58:1H:3679:HOH:O	2.04	0.75
26:1H:999:U:OP2	58:1H:4004:HOH:O	2.04	0.75
40:75:4:GLY:HA2	40:75:8:LYS:HZ3	1.52	0.75
36:35:55:ARG:HG2	36:35:56:SER:H	1.52	0.75
26:1H:270(E):G:N2	26:1H:270(U):C:N3	2.33	0.75
26:14:912:C:OP1	37:45:8:LYS:NZ	2.20	0.75
7:62:113:GLU:O	7:62:119:ARG:HD3	1.86	0.75
26:14:1729:A:H2'	26:14:1731:G:N2	2.02	0.75
30:39:120:GLU:HG3	30:39:122:LYS:HZ2	1.51	0.74
26:14:2705:A:OP2	58:14:3522:HOH:O	2.05	0.74
1:13:737:A:H2'	1:13:738:C:C6	2.22	0.74
26:14:654(B):C:O2	26:14:654(S):G:N2	2.19	0.74
1:13:1305:G:H21	1:13:1331:G:H2'	1.51	0.74
34:58:39:ARG:NH2	34:58:41:ASP:OD2	2.20	0.74
1:13:717:C:H5''	1:13:717:C:H6	1.51	0.74
40:B8:108:ARG:HA	40:B8:111:ARG:NE	2.02	0.74
26:14:1636:C:OP2	58:14:4091:HOH:O	2.04	0.74
40:B8:26:ASP:HB2	40:B8:91:ARG:HA	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:2K:47:7MG:H81	23:2K:48:U:H5	1.53	0.74
46:H8:151:HIS:ND1	46:H8:169:GLU:O	2.19	0.74
9:82:111:ARG:HB3	9:82:113:LYS:HE2	1.68	0.74
55:Q8:34:TRP:CB	55:Q8:35:GLN:HA	2.17	0.74
1:1G:785:G:N7	58:1G:1860:HOH:O	2.19	0.74
1:13:1003:G:H1	1:13:1037:C:H42	1.33	0.74
41:85:90:VAL:HG22	42:95:39:LEU:HB3	1.68	0.74
1:13:926:G:N2	25:4K:15:A:OP2	2.19	0.74
26:14:2581:G:O6	58:14:3718:HOH:O	2.02	0.74
38:98:29:LEU:HB3	38:98:75:LEU:HD11	1.67	0.74
1:13:1286:A:H5''	21:1F:26:LYS:HB2	1.69	0.74
55:Q8:46:ARG:NH2	55:Q8:48:PHE:HB3	1.99	0.74
26:14:973:A:OP2	58:14:4007:HOH:O	2.04	0.74
39:A8:34:HIS:HB2	39:A8:36:TYR:CE1	2.23	0.74
1:13:1454:G:OP1	20:BI:39:LYS:NZ	2.14	0.74
48:J8:92:LYS:HA	48:J8:95:LEU:HB2	1.70	0.74
26:1H:860:U:H5	26:1H:917:A:H2	1.32	0.74
29:29:23:VAL:HG11	29:29:183:LEU:HD23	1.68	0.74
26:14:2498:C:OP2	58:14:4053:HOH:O	2.06	0.74
26:14:1298:C:OP2	58:14:4064:HOH:O	2.04	0.74
36:35:30:THR:HG21	36:35:35:HIS:H	1.53	0.74
27:16:7:G:H4'	39:A8:29:PHE:CD2	2.23	0.74
1:13:1129:C:H4'	1:13:1130:A:H5'	1.70	0.74
26:1H:1516:U:H2'	26:1H:1517:G:C8	2.23	0.74
41:C8:92:ARG:O	41:C8:94:ASN:N	2.20	0.74
46:D5:71:VAL:HA	46:D5:88:PHE:HA	1.69	0.74
17:8I:13:ASP:OD1	17:8I:14:LYS:NZ	2.19	0.74
15:6I:6:GLU:HA	15:6I:9:GLN:HB2	1.69	0.74
26:14:2357:U:OP1	47:E5:20:ARG:NH1	2.21	0.74
1:1G:1125:U:O4	10:1A:5:ARG:NH1	2.21	0.74
26:1H:392:C:OP1	58:1H:3749:HOH:O	2.06	0.74
53:K5:24:GLU:HG3	53:K5:25:LYS:H	1.52	0.74
26:1H:4:C:O2	26:1H:2899:G:N2	2.20	0.74
1:1G:1213:A:N6	1:1G:1215:G:N3	2.35	0.74
11:2I:107:SER:HA	18:9I:87:ARG:HD2	1.70	0.74
26:14:450:G:OP2	58:14:4104:HOH:O	2.05	0.73
26:1H:1359:A:H2	26:1H:1372:U:O4	1.70	0.73
1:13:1500:A:OP1	58:13:1802:HOH:O	2.06	0.73
26:1H:2115:G:N2	26:1H:2172:U:O2	2.21	0.73
26:1H:802:A:OP1	58:1H:4587:HOH:O	2.06	0.73
37:88:135:ASP:HB3	37:88:137:TYR:H	1.51	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:673:G:H2'	1:13:674:G:C8	2.23	0.73
26:1H:2580:U:H4'	29:21:130:GLY:HA3	1.70	0.73
1:13:1224:G:C6	1:13:1322:C:H1'	2.23	0.73
26:14:654:A:O2'	26:14:654(A):A:N7	2.21	0.73
26:14:1642:G:N7	58:14:3951:HOH:O	2.21	0.73
36:35:71:VAL:HG13	36:35:72:PRO:HD3	1.70	0.73
26:1H:1036:G:N2	26:1H:1119:C:O2	2.20	0.73
26:1H:563:G:OP2	58:1H:3635:HOH:O	2.05	0.73
26:1H:1430:C:H2'	26:1H:1431:U:C6	2.22	0.73
4:3E:15:GLU:OE2	4:3E:59:ARG:NH2	2.20	0.73
26:14:399:G:OP2	58:14:4113:HOH:O	2.05	0.73
26:1H:1899:G:H22	26:1H:1902:C:H5	1.37	0.73
26:14:2000:G:OP2	58:14:3947:HOH:O	2.05	0.73
10:1I:28:ARG:HG3	10:1I:34:VAL:HG22	1.69	0.73
26:1H:2794:C:N4	26:1H:2802:G:O6	2.21	0.73
10:1A:3:LYS:N	10:1A:74:ILE:O	2.21	0.73
28:11:142:VAL:HG23	28:11:193:VAL:HA	1.69	0.73
26:1H:2308:G:N1	26:1H:2311:A:H2	1.86	0.73
26:14:619:G:H5''	26:14:620:G:N2	2.04	0.73
29:29:9:VAL:HG12	40:75:8:LYS:HZ1	1.52	0.73
48:J8:91:LYS:O	48:J8:94:LEU:N	2.20	0.73
23:2K:16:C:OP2	23:2K:17:C:N4	2.21	0.73
1:13:150:C:H2'	1:13:151:A:H8	1.52	0.73
26:14:2652:C:H42	26:14:2668:G:H1	1.36	0.73
31:41:64:THR:HG23	31:41:94:LEU:HD13	1.69	0.73
53:K5:28:ARG:HB3	53:K5:32:ASN:HA	1.68	0.73
40:75:4:GLY:O	40:75:7:ILE:N	2.22	0.73
1:1G:639:G:H2'	1:1G:640:A:H8	1.52	0.73
1:1G:1251:A:H2'	1:1G:1252:A:C8	2.23	0.73
20:BI:69:GLY:O	20:BI:73:HIS:NE2	2.22	0.73
9:8E:50:LEU:HD23	9:8E:85:LEU:HD11	1.70	0.73
26:1H:120:U:OP2	58:1H:4135:HOH:O	2.05	0.73
26:1H:1113:U:H5'	32:51:2:SER:HB2	1.70	0.73
46:H8:165:VAL:HB	46:H8:166:SER:HA	1.71	0.73
26:14:1839:G:OP2	58:14:3985:HOH:O	2.06	0.73
26:1H:400:G:O6	58:1H:4042:HOH:O	2.03	0.73
1:1G:1003:G:N2	1:1G:1037:C:O2	2.15	0.73
26:14:1754:C:OP1	40:75:96:ARG:NH1	2.20	0.73
26:1H:2593:U:H2'	26:1H:2594:C:C6	2.23	0.73
1:1G:426:G:OP1	4:32:38:TYR:OH	2.05	0.73
26:1H:409:C:OP1	58:1H:3749:HOH:O	2.06	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:192:U:H2'	1:1G:193:C:H6	1.52	0.73
36:35:14:LYS:O	36:35:16:ARG:N	2.22	0.73
26:14:84:A:N6	26:14:102:G:O2'	2.16	0.73
55:M5:40:GLU:H	55:M5:43:GLN:HG3	1.53	0.73
42:95:57:VAL:HG23	42:95:99:ILE:H	1.53	0.73
4:32:18:LYS:HD3	4:32:31:CYS:SG	2.29	0.73
26:1H:2392:A:H8	36:78:61:ARG:HG2	1.53	0.73
26:14:882:G:H22	26:14:894:C:N4	1.87	0.73
26:1H:2400:G:H2'	26:1H:2401:U:C6	2.23	0.73
26:14:1257:C:H4'	30:39:83:PHE:CE1	2.24	0.73
30:39:25:PRO:HB3	30:39:28:ILE:HG23	1.70	0.73
26:14:2121:G:H1	26:14:2177:C:H42	1.35	0.73
26:1H:49:A:N7	26:1H:120:U:H5	1.85	0.73
26:1H:1077:A:H3'	26:1H:1078:U:H5'	1.70	0.73
1:13:533:A:OP1	58:13:1870:HOH:O	2.06	0.73
39:A8:25:ARG:NH1	39:A8:42:ASP:OD2	2.22	0.73
5:4E:45:PHE:CE2	5:4E:47:LYS:HD2	2.24	0.73
46:D5:139:VAL:HG13	46:D5:156:LYS:HE2	1.70	0.73
38:98:56:LYS:NZ	38:98:90:ARG:O	2.20	0.73
42:95:70:ILE:N	42:95:86:GLY:O	2.18	0.73
1:13:1007:C:H42	1:13:1022:G:H1	1.36	0.73
27:16:80:U:H2'	27:16:81:G:H21	1.53	0.73
10:1I:34:VAL:HG12	10:1I:74:ILE:HG23	1.70	0.73
26:1H:1055:G:O2'	26:1H:1086:A:N6	2.22	0.73
26:1H:176:G:O2'	26:1H:177:G:H5'	1.89	0.73
26:14:2502:G:OP2	58:14:4030:HOH:O	2.06	0.73
36:78:52:GLU:HG3	36:78:57:THR:HA	1.71	0.73
24:3L:11:C:H2'	24:3L:12:U:H6	1.54	0.73
55:Q8:46:ARG:HH21	55:Q8:48:PHE:CB	1.97	0.73
55:M5:52:LYS:NZ	55:M5:53:PRO:HA	2.03	0.73
29:21:111:ARG:HD3	29:21:160:TYR:CE2	2.24	0.73
26:1H:2347:C:OP1	53:O8:39:TYR:OH	2.06	0.73
26:1H:1438:U:H2'	26:1H:1439:A:H8	1.54	0.73
40:75:50:ILE:HD11	40:75:102:ILE:HD11	1.69	0.73
32:59:6:ARG:H	32:59:6:ARG:HH11	1.35	0.73
1:1G:1316:G:H22	1:1G:1319:A:H5''	1.52	0.73
26:14:1899:G:N2	26:14:1902:C:N4	2.36	0.73
45:C5:19:LYS:HG3	45:C5:20:TYR:H	1.53	0.73
19:AA:20:LEU:O	19:AA:23:ASN:ND2	2.21	0.73
55:Q8:57:ARG:HD3	55:Q8:57:ARG:N	2.03	0.72
26:1H:2492:U:H2'	26:1H:2493:U:C6	2.24	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2343:C:O2'	26:14:2373:G:O2'	2.05	0.72
26:1H:1803:A:O2'	28:11:259:THR:HG21	1.89	0.72
55:Q8:21:LYS:HE2	55:Q8:48:PHE:HE1	1.54	0.72
26:1H:2312:U:H5'	31:41:88:ILE:HD12	1.70	0.72
1:13:1358:U:OP2	1:13:1359:C:N4	2.20	0.72
24:3L:8:U:H3	24:3L:14:A:H62	1.33	0.72
2:1E:8:LYS:HG2	2:1E:9:GLU:H	1.55	0.72
26:1H:1495:A:OP2	58:1H:4431:HOH:O	2.06	0.72
44:B5:5:TYR:HB3	49:G5:33:MET:HB2	1.70	0.72
55:Q8:31:HIS:ND1	55:Q8:31:HIS:O	2.22	0.72
32:51:124:GLU:HB3	32:51:132:ARG:HB3	1.70	0.72
27:1J:80:U:H2'	27:1J:81:G:N2	2.05	0.72
26:1H:331:A:N1	58:1H:3844:HOH:O	2.22	0.72
26:1H:2033:A:H8	58:1H:4060:HOH:O	1.71	0.72
44:F8:3:THR:OG1	44:F8:4:ALA:HA	1.88	0.72
1:1G:371:G:H1	1:1G:390:C:H42	1.35	0.72
44:B5:41:ASN:HA	44:B5:44:GLU:HB2	1.71	0.72
39:65:84:GLN:HA	39:65:110:LEU:HD12	1.70	0.72
26:14:323:G:O2'	26:14:1205:U:N3	2.13	0.72
9:8E:96:LEU:HD23	9:8E:102:LEU:HD21	1.70	0.72
46:H8:165:VAL:HB	46:H8:167:PRO:HD3	1.71	0.72
26:1H:1994:C:OP1	58:1H:4006:HOH:O	2.06	0.72
43:E8:96:ILE:HD13	43:E8:96:ILE:H	1.52	0.72
1:1G:947:G:O3'	13:4A:109:THR:OG1	2.07	0.72
1:13:1008:C:N4	1:13:1021:G:O6	2.18	0.72
26:14:1308:A:OP2	58:14:3849:HOH:O	2.08	0.72
45:G8:87:LYS:HB3	45:G8:94:LYS:HG2	1.70	0.72
26:14:2468:G:H3'	26:14:2476:A:N1	2.04	0.72
26:1H:1525:G:H2'	26:1H:1526:G:C8	2.24	0.72
23:2K:47:7MG:H5'	23:2K:47:7MG:H82	1.72	0.72
26:14:731:C:OP1	58:14:4108:HOH:O	2.08	0.72
1:1G:707:C:H2'	1:1G:708:C:H6	1.53	0.72
26:14:2572:A:C8	29:29:144:ARG:HD2	2.23	0.72
26:1H:1828:G:OP1	58:1H:4686:HOH:O	2.07	0.72
1:13:963:G:N2	1:13:972:C:N3	2.31	0.72
1:1G:1248:A:N3	9:82:70:LYS:NZ	2.37	0.72
1:13:1213:A:O2'	1:13:1215:G:N7	2.17	0.72
11:2A:27:ASN:OD1	11:2A:28:THR:N	2.22	0.72
55:Q8:49:VAL:HG23	55:Q8:50:LEU:H	1.55	0.72
55:Q8:6:THR:N	55:Q8:59:LYS:HE3	2.04	0.72
26:14:1061:U:H4'	26:14:1070:A:H1'	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:22:65:ALA:HA	3:22:100:ALA:HB3	1.71	0.72
1:1G:1311:G:N2	1:1G:1326:C:O2	2.20	0.72
1:1G:8:A:C6	4:32:209:ARG:HB2	2.24	0.72
16:7A:37:GLY:HA2	16:7A:50:LYS:HD3	1.70	0.72
27:1J:15:A:H5'	27:1J:16:G:C8	2.25	0.72
28:11:10:THR:OG1	28:11:13:ARG:HB2	1.90	0.72
9:82:28:VAL:HG22	9:82:63:ILE:HB	1.72	0.72
9:8E:112:LYS:HA	9:8E:119:ALA:HB2	1.70	0.72
31:49:40:ASN:HB2	31:49:91:ARG:HG3	1.72	0.72
1:1G:376:G:O3'	16:7A:5:ARG:NH1	2.22	0.72
45:C5:52:SER:HA	45:C5:55:TYR:O	1.89	0.72
24:3K:19:G:OP1	24:3K:60:U:N3	2.19	0.72
26:1H:761:A:OP1	58:1H:3686:HOH:O	2.08	0.72
26:1H:1639:U:O2'	26:1H:1640:C:H5'	1.89	0.72
24:3K:72:C:H3'	24:3K:73:A:H5''	1.72	0.72
1:13:963:G:N3	10:1I:55:LYS:NZ	2.38	0.72
26:14:2393:A:H4'	36:35:62:LEU:H	1.52	0.72
26:14:2134:A:O2'	26:14:2159:G:N2	2.22	0.72
26:14:2340:G:H2'	26:14:2341:G:H8	1.53	0.72
26:14:1264:G:OP1	52:J5:19:ARG:NH2	2.21	0.72
38:98:67:LEU:HD13	38:98:76:VAL:HG21	1.71	0.72
26:14:1899:G:H21	26:14:1902:C:N4	1.87	0.72
26:14:2327:A:H2'	26:14:2328:A:C8	2.25	0.72
26:1H:2693:A:H2'	26:1H:2694:G:H8	1.54	0.72
26:1H:974(A):C:OP1	58:1H:4189:HOH:O	2.07	0.72
30:31:155:LEU:HD11	30:31:176:LEU:HD13	1.72	0.72
1:1G:1432:G:N2	58:1G:1858:HOH:O	2.23	0.72
29:21:116:VAL:HG11	29:21:138:PRO:HB3	1.71	0.71
26:1H:764:A:H2	28:11:219:PRO:HG3	1.53	0.71
1:13:547:A:OP1	4:3E:73:ARG:NH2	2.23	0.71
26:1H:951:C:O2	26:1H:966:G:N2	2.20	0.71
5:42:102:ALA:HB1	5:42:106:PRO:HG2	1.73	0.71
1:1G:1256:A:N6	1:1G:1278:U:OP2	2.22	0.71
26:1H:240:G:O6	58:1H:4452:HOH:O	2.06	0.71
14:5I:29:ARG:HH21	14:5I:41:ARG:NH1	1.88	0.71
26:14:1434:A:H61	26:14:1558:A:H62	1.37	0.71
29:21:201:THR:HG22	29:21:203:LYS:H	1.55	0.71
26:14:674:G:O2'	30:39:74:ARG:HG3	1.89	0.71
40:75:92:GLY:HA2	40:75:116:ALA:HA	1.71	0.71
26:1H:731:C:OP2	58:1H:3690:HOH:O	2.07	0.71
26:1H:1479:G:N7	26:1H:1510:A:N6	2.37	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1298:C:H41	7:62:114:ARG:HB3	1.54	0.71
38:98:100:LEU:HD11	38:98:113:LEU:HD13	1.72	0.71
14:5I:26:ARG:NH1	14:5I:43:CYS:SG	2.63	0.71
26:1H:2871:C:N3	58:1H:4092:HOH:O	2.22	0.71
26:14:34:C:O2'	26:14:35:G:OP1	2.06	0.71
26:1H:910:A:H62	37:88:12:GLN:HA	1.55	0.71
26:1H:252:G:OP2	36:78:50:ARG:NH1	2.24	0.71
20:BI:26:ASN:HB2	20:BI:71:THR:HG23	1.72	0.71
12:3A:70:ILE:HD13	12:3A:77:LEU:HD12	1.73	0.71
1:13:877:C:OP1	8:7E:88:LYS:NZ	2.18	0.71
26:1H:2053:G:OP1	58:1H:3801:HOH:O	2.08	0.71
26:14:2096:U:H3	26:14:2193:G:H1	1.37	0.71
1:13:1194:U:H2'	1:13:1195:C:C6	2.26	0.71
26:1H:993:G:OP1	41:C8:50:ARG:NH2	2.24	0.71
4:3E:15:GLU:OE1	4:3E:66:ARG:NH1	2.22	0.71
1:1G:1157:A:H61	1:1G:1178:G:H21	1.38	0.71
26:1H:256:A:OP2	58:1H:4457:HOH:O	2.07	0.71
39:65:34:HIS:HE2	39:65:54:LEU:HD13	1.56	0.71
1:13:559:A:OP1	5:4E:126:ARG:NH2	2.23	0.71
55:M5:33:ASN:OD1	55:M5:33:ASN:N	2.23	0.71
20:BA:64:ASP:OD2	20:BA:81:LYS:NZ	2.20	0.71
31:41:112:PRO:HB3	51:M8:37:SER:H	1.55	0.71
52:N8:40:LYS:HE2	52:N8:47:PRO:HD2	1.71	0.71
27:1J:42:C:N3	31:49:91:ARG:NH2	2.39	0.71
47:I8:11:ARG:O	47:I8:14:ARG:NH2	2.24	0.71
24:3K:50:U:H3	24:3K:64:A:H61	1.37	0.71
26:1H:2597:G:O3'	58:1H:3651:HOH:O	2.07	0.71
20:BA:51:GLU:HA	20:BA:54:LYS:HE3	1.73	0.71
37:45:22:LYS:HG2	37:45:23:GLY:HA2	1.72	0.71
37:88:32:TYR:CE2	37:88:133:ARG:HG3	2.26	0.71
26:14:1332:G:N2	26:14:1609:A:O2'	2.23	0.71
29:21:117:MET:HE1	29:21:136:ARG:HA	1.72	0.71
26:14:2392:A:H1'	36:35:61:ARG:HH21	1.54	0.71
46:D5:30:ASN:HA	46:D5:89:PHE:HE1	1.54	0.71
26:1H:2327:A:H2'	26:1H:2328:A:C8	2.25	0.71
34:58:130:HIS:C	34:58:134:ARG:HH22	1.93	0.71
6:5E:23:LYS:HD3	6:5E:61:LEU:HD21	1.73	0.71
23:2K:40:C:O2'	24:3K:35:A:O2'	2.08	0.71
26:1H:958:U:OP2	37:88:14:ARG:NH1	2.21	0.71
30:31:6:VAL:HG11	30:31:119:ARG:HA	1.72	0.71
26:1H:2286:A:H2'	53:O8:31:PRO:HD3	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:67:U:H3	26:14:74:A:H2	1.35	0.71
30:39:135:LYS:HB3	30:39:138:GLU:HG3	1.73	0.71
1:13:14:U:OP2	58:13:1860:HOH:O	2.09	0.71
38:98:91:GLN:O	38:98:91:GLN:NE2	2.21	0.71
26:1H:516:C:OP1	52:N8:13:LYS:NZ	2.24	0.71
26:14:2499:C:OP2	58:14:4052:HOH:O	2.08	0.71
37:45:85:LYS:HG2	37:45:86:GLY:H	1.56	0.71
26:1H:2419:U:O4	55:Q8:29:LYS:NZ	2.13	0.71
26:1H:2887:U:H2'	26:1H:2888:C:H6	1.55	0.71
22:1K:27:G:N2	22:1K:43:C:O2	2.14	0.71
15:6A:87:ILE:HG22	15:6A:88:ARG:H	1.55	0.71
10:1I:57:LYS:HD2	10:1I:60:ARG:HH12	1.56	0.70
46:D5:53:ILE:HG22	46:D5:71:VAL:HG13	1.71	0.70
6:52:7:ASN:HD22	18:9A:76:LEU:HD11	1.54	0.70
26:1H:2758:A:OP2	58:1H:4479:HOH:O	2.08	0.70
53:O8:9:LEU:HB3	53:O8:27:LYS:HA	1.73	0.70
10:1I:48:THR:HA	10:1I:62:HIS:HB3	1.72	0.70
46:D5:30:ASN:HD22	46:D5:90:VAL:HB	1.55	0.70
1:1G:1002:G:H2'	1:1G:1003:G:C8	2.26	0.70
1:1G:690:G:H1	11:2A:55:LYS:HZ3	1.38	0.70
26:14:1412:A:H2'	26:14:1413:G:C8	2.26	0.70
31:49:97:ASP:HA	31:49:100:TRP:HD1	1.55	0.70
26:14:2420:C:N4	55:M5:31:HIS:O	2.24	0.70
6:52:23:LYS:HE3	6:52:61:LEU:HD21	1.71	0.70
26:1H:2135:A:N6	26:1H:2156:G:O2'	2.23	0.70
26:1H:1418:G:OP1	26:1H:1588:C:O2'	2.09	0.70
26:1H:2292:C:OP1	39:A8:17:ARG:NH2	2.23	0.70
26:1H:1434:A:H61	26:1H:1558:A:H61	1.39	0.70
1:1G:1348:U:H4'	9:82:120:ARG:HD2	1.73	0.70
36:78:138:LEU:HD12	36:78:144:GLU:HG3	1.73	0.70
39:A8:37:ALA:HB2	39:A8:101:LEU:HD21	1.73	0.70
26:1H:1918:A:O2'	26:1H:1920:C:N4	2.24	0.70
26:1H:534:U:H5'	41:C8:42:ALA:HB1	1.74	0.70
30:39:188:ARG:HA	36:35:3:LEU:HD11	1.73	0.70
1:1G:501:C:H2'	1:1G:502:G:H8	1.57	0.70
26:14:2846:G:N7	58:14:3519:HOH:O	2.23	0.70
26:1H:1330:C:OP1	58:1H:3963:HOH:O	2.10	0.70
17:8I:67:LYS:HA	17:8I:70:ARG:HH12	1.54	0.70
26:1H:910:A:C5	37:88:13:GLN:HG3	2.26	0.70
26:1H:1332:G:N2	26:1H:1609:A:O2'	2.23	0.70
26:1H:1494:A:H2'	26:1H:1495:A:C8	2.27	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:59:10:PRO:HD2	32:59:50:VAL:HG13	1.71	0.70
26:14:2878:U:O4	58:14:3855:HOH:O	2.09	0.70
32:59:106:THR:HG22	32:59:112:PRO:HB3	1.73	0.70
26:1H:848:G:H2'	26:1H:849:A:C8	2.26	0.70
26:14:2331:G:O2'	47:E5:43:THR:HB	1.92	0.70
1:13:659:U:H2'	1:13:660:G:H8	1.55	0.70
23:2K:47:7MG:H81	23:2K:48:U:C5	2.25	0.70
55:Q8:37:SER:HA	55:Q8:39:LYS:O	1.90	0.70
41:C8:92:ARG:NH1	41:C8:94:ASN:OD1	2.25	0.70
26:1H:1899:G:N2	26:1H:1902:C:H5	1.90	0.70
7:6E:95:ARG:HH21	7:6E:99:LEU:HD11	1.55	0.70
1:13:233:C:H2'	1:13:234:C:H6	1.57	0.70
1:1G:330:C:O2	58:1G:1821:HOH:O	2.10	0.70
26:1H:142:G:H1'	44:F8:37:THR:HG21	1.71	0.70
31:49:16:ARG:NH2	31:49:28:VAL:O	2.23	0.70
26:1H:1385:G:HO2'	26:1H:1396:U:H6	1.37	0.70
51:I5:34:GLU:HG2	51:I5:35:VAL:H	1.57	0.70
1:1G:539:A:OP2	12:3A:115:LYS:NZ	2.24	0.70
26:1H:2838:G:N7	58:1H:4251:HOH:O	2.25	0.70
1:13:1391:U:H2'	1:13:1392:G:C8	2.27	0.70
39:A8:48:LEU:HD23	39:A8:82:ILE:HD11	1.74	0.70
27:16:101:A:OP2	58:16:303:HOH:O	2.08	0.70
26:1H:1156:A:C8	41:C8:51:LYS:HD2	2.26	0.70
26:14:796:C:H2'	26:14:797:C:C6	2.26	0.70
53:O8:32:ASN:N	53:O8:32:ASN:OD1	2.23	0.70
26:14:2837:G:N7	58:14:3851:HOH:O	2.24	0.70
26:14:2818:G:OP2	38:55:42:LYS:NZ	2.25	0.70
4:32:31:CYS:C	4:32:33:MET:H	1.93	0.70
26:1H:2711:A:OP2	58:1H:3670:HOH:O	2.09	0.70
26:1H:568:U:O4	58:1H:4553:HOH:O	2.04	0.70
45:C5:99:CYS:SG	45:C5:100:ALA:N	2.65	0.70
26:1H:860:U:C5	26:1H:917:A:C2	2.78	0.70
26:1H:76:C:O2'	49:K8:62:THR:HG21	1.91	0.70
28:19:223:GLY:HA2	28:19:226:MET:HG3	1.71	0.70
1:13:1423:G:OP1	35:68:49:ARG:NH2	2.25	0.70
49:K8:18:PRO:HA	49:K8:21:LEU:HB2	1.73	0.70
1:1G:371:G:O2'	1:1G:373:A:N7	2.24	0.70
1:13:403:C:OP1	4:3E:137:SER:OG	2.08	0.70
26:14:1170:G:O6	26:14:1179:C:N4	2.25	0.70
12:3I:70:ILE:HD13	12:3I:77:LEU:HD12	1.72	0.70
44:F8:2:LYS:HG2	49:K8:26:ARG:HE	1.55	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:154:G:O6	26:14:172:C:N4	2.24	0.70
28:19:72:LYS:NZ	28:19:99:ASP:OD2	2.22	0.70
26:14:1111:A:H4'	32:59:3:ARG:HD3	1.73	0.69
1:1G:1132:C:H42	1:1G:1142:G:H1	1.40	0.69
26:1H:2134:A:OP2	26:1H:2157:G:N2	2.25	0.69
26:14:1891:G:O6	58:14:3913:HOH:O	2.04	0.69
26:1H:1858:G:O6	58:1H:4447:HOH:O	2.10	0.69
38:98:87:TYR:HE1	38:98:117:VAL:HG12	1.56	0.69
1:13:77:C:H1'	1:13:92:G:H1	1.56	0.69
26:1H:578:A:OP2	58:1H:4640:HOH:O	2.08	0.69
11:2I:86:GLY:N	11:2I:112:THR:OG1	2.19	0.69
26:1H:1794:U:H2'	26:1H:1795:C:C6	2.26	0.69
1:1G:957:U:H1'	1:1G:960:U:C5	2.27	0.69
43:E8:14:PRO:HG2	43:E8:78:GLU:HB2	1.73	0.69
26:14:424:G:O6	58:14:3828:HOH:O	2.09	0.69
2:1E:226:ARG:HG3	2:1E:227:GLY:H	1.57	0.69
26:1H:1382:G:O6	58:1H:4567:HOH:O	2.07	0.69
1:1G:64:G:H4'	1:1G:65:U:H5''	1.74	0.69
2:12:91:PRO:HG3	2:12:154:LEU:HB2	1.74	0.69
27:16:44:G:O2'	27:16:47:C:N4	2.25	0.69
27:1J:2:C:H2'	27:1J:3:C:H6	1.58	0.69
1:13:56:U:H4'	33:69:82:ARG:HH21	1.57	0.69
43:E8:73:ALA:HB3	43:E8:106:ILE:HB	1.73	0.69
45:C5:73:ARG:NH2	45:C5:81:LYS:O	2.26	0.69
55:M5:22:VAL:HB	55:M5:53:PRO:HG3	1.74	0.69
1:13:1157:A:N6	1:13:1178:G:H21	1.88	0.69
26:14:1257:C:H4'	30:39:83:PHE:CD1	2.27	0.69
1:13:1368:G:H5''	9:8E:112:LYS:HB3	1.73	0.69
18:9A:32:ARG:HD3	18:9A:65:ILE:HD11	1.73	0.69
26:1H:1204:A:H61	26:1H:1240:U:H2'	1.56	0.69
51:M8:48:ARG:HH11	51:M8:48:ARG:HA	1.57	0.69
36:35:121:LYS:HG3	36:35:122:PRO:HD2	1.73	0.69
1:13:1508:G:OP1	58:13:1802:HOH:O	2.09	0.69
26:14:1828:G:OP1	58:14:3447:HOH:O	2.10	0.69
52:N8:41:PRO:HD2	52:N8:44:THR:HG21	1.74	0.69
1:1G:1286:A:H5'	21:1B:25:LYS:HZ3	1.57	0.69
34:58:96:GLU:C	34:58:98:VAL:H	1.96	0.69
1:1G:519:C:H2'	1:1G:520:A:O4'	1.91	0.69
31:49:106:LEU:HA	31:49:110:ALA:HB3	1.74	0.69
34:15:104:LYS:HA	34:15:107:LEU:HD12	1.73	0.69
26:14:2364:C:H4'	47:E5:56:ASP:OD1	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1435:G:H2'	1:13:1436:U:C6	2.27	0.69
26:1H:2334:G:H5'	39:A8:9:ARG:HG2	1.74	0.69
46:H8:4:ARG:NH1	46:H8:60:GLU:OE1	2.25	0.69
26:14:1021:A:H62	26:14:1141:U:H3	1.38	0.69
26:1H:878:A:H61	26:1H:899:A:HO2'	1.39	0.69
1:13:111:G:H5''	16:7I:27:LYS:HG2	1.74	0.69
26:1H:2213:U:O2	48:J8:52:ARG:NH2	2.25	0.69
40:B8:64:ARG:HB2	40:B8:73:GLU:HG2	1.74	0.69
36:35:105:LEU:O	36:35:106:LEU:HB3	1.92	0.69
32:51:74:ASN:HA	32:51:77:LYS:HD3	1.74	0.69
55:Q8:53:PRO:HB3	55:Q8:56:GLU:H	1.58	0.69
26:14:2210:G:H3'	26:14:2211:G:C2	2.28	0.69
24:3L:33:U:N3	24:3L:36:A:OP2	2.25	0.69
26:14:2777:G:H5''	26:14:2778:A:H5'	1.75	0.69
41:85:49:HIS:HA	41:85:52:ARG:HB2	1.75	0.69
16:7I:74:LEU:HA	16:7I:77:ALA:HB2	1.75	0.69
26:1H:1678:G:H22	26:1H:1989:G:H22	1.38	0.69
26:1H:760:G:OP1	58:1H:3829:HOH:O	2.09	0.69
26:14:802:A:H4'	58:14:3622:HOH:O	1.93	0.69
1:13:1160:G:H22	1:13:1177:G:H22	1.40	0.69
1:1G:1132:C:N3	1:1G:1142:G:N2	2.40	0.69
26:1H:2210:G:H3'	26:1H:2211:G:C8	2.28	0.69
20:BI:26:ASN:O	20:BI:30:LYS:HB2	1.93	0.69
1:1G:979:C:H5	1:1G:980:C:C6	2.11	0.69
27:16:44:G:H1'	27:16:47:C:H42	1.58	0.69
26:14:2689:U:P	26:14:2719:G:H22	2.14	0.69
45:G8:55:TYR:HB2	45:G8:58:GLY:HA3	1.75	0.69
1:13:523:A:H61	12:3I:92:ASP:HB2	1.55	0.69
33:61:8:PRO:HA	33:61:14:ASP:HA	1.73	0.69
32:51:15:VAL:HG12	32:51:28:GLY:HA3	1.73	0.69
26:1H:280:C:N3	26:1H:361:G:N2	2.41	0.69
30:39:79:GLY:HA2	30:39:86:GLY:HA2	1.74	0.69
26:14:1757:U:H3	26:14:1762:A:H2	1.39	0.69
27:16:100:G:OP1	58:16:315:HOH:O	2.09	0.69
1:1G:1235:U:O2'	1:1G:1305:G:O5'	2.09	0.69
2:12:178:ARG:NH1	2:12:196:LEU:O	2.20	0.69
26:1H:323:G:C8	30:31:171:PRO:HG3	2.28	0.69
26:1H:2031:A:O2'	26:1H:2454:G:N2	2.26	0.69
43:A5:62:HIS:HB2	43:A5:64:MET:HG3	1.75	0.69
30:39:28:ILE:HA	30:39:112:MET:HG2	1.73	0.69
39:A8:24:LEU:HB2	39:A8:85:VAL:HG12	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:42:9:LYS:HB2	5:42:112:LEU:HD11	1.74	0.69
1:13:504:C:OP1	58:13:1883:HOH:O	2.11	0.69
26:14:1187:G:OP2	58:14:3586:HOH:O	2.10	0.69
1:13:560:U:O2'	1:13:561:U:OP2	2.11	0.69
37:88:75:THR:HG21	37:88:85:LYS:HD2	1.73	0.69
1:13:1302:U:OP2	13:4I:21:TYR:OH	2.08	0.69
33:61:131:LYS:HB3	33:61:132:PRO:HA	1.74	0.69
2:1E:100:GLY:O	2:1E:104:ASN:N	2.25	0.68
26:14:973:A:OP2	58:14:4004:HOH:O	2.09	0.68
26:1H:2249:U:O4	58:1H:3724:HOH:O	2.10	0.68
46:H8:152:ALA:HB3	46:H8:167:PRO:HA	1.73	0.68
16:7A:11:SER:H	16:7A:14:ASN:HB3	1.57	0.68
16:7I:50:LYS:HD3	16:7I:51:VAL:H	1.58	0.68
43:A5:14:PRO:HG2	43:A5:78:GLU:HG3	1.73	0.68
1:1G:114:U:H2'	1:1G:115:G:C8	2.28	0.68
26:1H:2233:U:H2'	26:1H:2234:G:C8	2.28	0.68
22:1K:76:A:H2	26:1H:2602:A:H2	1.40	0.68
1:1G:345:C:H1'	1:1G:346:G:C2	2.28	0.68
49:G5:47:ASN:O	49:G5:49:LYS:N	2.25	0.68
1:1G:1149:C:O2'	1:1G:1280:A:N1	2.24	0.68
13:4A:82:MET:HG2	13:4A:83:ASP:N	2.09	0.68
26:1H:2469:A:H2	26:1H:2481:G:H21	1.41	0.68
9:8E:49:PRO:HA	9:8E:52:ALA:HB3	1.74	0.68
26:14:2061:G:OP2	58:14:4027:HOH:O	2.10	0.68
30:39:88:VAL:HG23	30:39:89:VAL:O	1.93	0.68
1:1G:452:A:O2'	1:1G:453:A:O4'	2.07	0.68
20:BA:26:ASN:HD22	20:BA:71:THR:HA	1.58	0.68
43:E8:24:ILE:HG12	43:E8:36:LEU:HD21	1.74	0.68
37:88:104:PHE:HE2	37:88:125:LEU:HD11	1.58	0.68
45:C5:76:CYS:SG	45:C5:97:ARG:HG3	2.33	0.68
26:14:96:G:H4'	49:G5:48:HIS:CE1	2.28	0.68
26:1H:2027:G:N7	58:1H:4063:HOH:O	2.25	0.68
40:B8:16:ARG:HE	40:B8:19:LEU:HD21	1.58	0.68
2:12:212:GLN:HE22	2:12:216:SER:HB2	1.58	0.68
31:49:109:VAL:HG11	31:49:142:PRO:HG3	1.75	0.68
26:14:943:U:OP2	36:35:36:LYS:HG3	1.93	0.68
37:45:25:ASP:HB3	37:45:102:VAL:HG23	1.73	0.68
48:J8:18:ILE:HG13	48:J8:37:ILE:HG23	1.75	0.68
1:13:1159:U:O4'	1:13:1182:G:N2	2.26	0.68
17:8I:81:ARG:NH2	17:8I:83:ASP:OD2	2.26	0.68
26:1H:1188:U:H4'	42:D8:79:VAL:HG22	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:AA:38:SER:HB2	19:AA:71:LEU:HG	1.74	0.68
26:14:1678:G:H22	26:14:1989:G:N2	1.92	0.68
26:1H:566:U:OP1	36:78:29:LYS:NZ	2.22	0.68
1:1G:631:G:H3'	1:1G:632:A:H8	1.57	0.68
26:1H:625:G:N7	36:78:107:LYS:NZ	2.42	0.68
26:14:792:G:O6	58:14:3469:HOH:O	2.06	0.68
28:11:30:GLU:HB3	28:11:33:LEU:HB2	1.74	0.68
44:F8:36:LYS:HG2	44:F8:54:VAL:HB	1.75	0.68
13:4A:60:VAL:HG13	13:4A:64:TRP:HE1	1.58	0.68
35:25:47:ILE:HG13	35:25:48:PRO:HD2	1.76	0.68
1:1G:554:C:H2'	1:1G:555:C:C6	2.28	0.68
40:75:56:GLY:O	40:75:59:THR:HG23	1.92	0.68
26:1H:2126:A:N6	26:1H:2163:C:O2	2.26	0.68
34:15:13:TRP:O	34:15:135:PRO:HD2	1.94	0.68
16:7I:47:ASP:OD1	16:7I:47:ASP:N	2.24	0.68
26:14:2392:A:H2	26:14:2424:C:N4	1.91	0.68
37:45:21:THR:HG22	37:45:23:GLY:HA3	1.75	0.68
26:1H:2287:A:N6	26:1H:2344:U:H3	1.92	0.68
52:J5:20:ARG:HG2	52:J5:23:HIS:CD2	2.28	0.68
28:11:136:ILE:O	28:11:168:ARG:NH2	2.27	0.68
26:14:309:G:N3	26:14:329:G:O2'	2.27	0.68
45:G8:76:CYS:O	45:G8:78:ALA:N	2.27	0.68
1:1G:1154:G:H2'	1:1G:1155:G:H8	1.58	0.68
37:45:57:HIS:NE2	37:45:116:GLU:HB3	2.08	0.68
40:B8:26:ASP:CB	40:B8:92:GLY:H	2.06	0.68
32:51:83:TYR:HB3	32:51:135:GLY:N	2.08	0.68
30:39:117:ARG:HH12	36:35:1:MET:H2	1.39	0.68
26:1H:850:C:O3'	50:L8:49:LYS:NZ	2.26	0.68
1:1G:664:G:H22	1:1G:741:G:H1	1.42	0.68
26:1H:1857:G:N7	58:1H:4448:HOH:O	2.26	0.68
26:1H:71:A:H2	44:F8:31:HIS:HE1	1.40	0.68
38:55:20:LEU:HD21	38:55:40:LYS:HD3	1.74	0.68
26:1H:1771:C:OP2	58:1H:3986:HOH:O	2.12	0.68
31:41:67:LYS:HE2	51:M8:6:HIS:CE1	2.29	0.68
26:14:2750:A:OP2	32:59:6:ARG:NH2	2.27	0.68
26:14:674:G:OP2	58:14:3905:HOH:O	2.10	0.68
26:1H:1678:G:N2	26:1H:1989:G:H22	1.92	0.68
49:G5:17:SER:N	49:G5:20:GLU:OE2	2.26	0.68
32:51:113:VAL:HG11	32:51:151:ILE:HD13	1.76	0.68
26:1H:2313:C:H4'	31:41:91:ARG:HG3	1.74	0.68
26:1H:810:U:OP1	58:1H:3711:HOH:O	2.11	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
54:P8:9:ARG:NH2	54:P8:47:ARG:HD2	2.07	0.68
46:D5:69:THR:HG22	46:D5:90:VAL:HG22	1.76	0.68
1:1G:1130:A:H62	1:1G:1144:G:H21	1.40	0.68
26:14:2157:G:HO2'	26:14:2158:A:H8	1.39	0.68
26:14:67:U:H2'	26:14:68:G:H8	1.58	0.68
23:2L:24:C:H2'	23:2L:25:U:C6	2.28	0.68
31:41:37:VAL:H	31:41:99:MET:HE3	1.58	0.68
26:14:71:A:H2	44:B5:31:HIS:HE2	1.42	0.68
28:11:182:LEU:H	28:11:272:ALA:HB3	1.58	0.68
37:88:5:ARG:H	37:88:5:ARG:HD3	1.59	0.68
23:2K:54:G:H2'	23:2K:55:5MU:H6	1.57	0.68
26:1H:879:G:O6	26:1H:898:C:N4	2.26	0.68
26:14:2720:U:N3	26:14:2873:A:H2	1.90	0.68
1:1G:978:A:O2'	1:1G:1322:C:N3	2.27	0.68
9:82:128:ARG:NH1	23:2L:36:A:OP2	2.27	0.68
22:1L:76:A:H62	26:14:2583:G:H21	1.41	0.68
26:14:1890:A:OP2	58:14:3912:HOH:O	2.12	0.68
12:3A:117:ARG:HB3	12:3A:122:THR:HB	1.76	0.68
26:14:2228:G:O6	58:14:4035:HOH:O	2.12	0.68
26:14:1300:U:OP1	58:14:4093:HOH:O	2.12	0.68
29:29:151:TYR:HD2	29:29:154:LYS:HZ2	1.41	0.68
26:14:1778:U:H2'	26:14:1784:A:N6	2.09	0.68
26:1H:376:C:OP1	58:1H:3754:HOH:O	2.12	0.67
26:14:2445:G:N7	58:14:4029:HOH:O	2.26	0.67
20:BI:43:LEU:HD13	20:BI:51:GLU:HB3	1.75	0.67
26:14:1849:G:H2'	26:14:1850:G:H8	1.60	0.67
26:1H:836:G:H5''	26:1H:837:C:OP2	1.94	0.67
1:13:1422:G:H5''	35:68:48:PRO:HB3	1.75	0.67
1:13:287:U:H2'	1:13:288:A:H8	1.57	0.67
26:14:1188:U:O2'	26:14:1189:A:H5'	1.93	0.67
26:1H:1253:A:N7	58:1H:3710:HOH:O	2.27	0.67
26:1H:376:C:OP2	58:1H:3750:HOH:O	2.12	0.67
26:14:1828:G:OP2	58:14:3421:HOH:O	2.12	0.67
36:35:57:THR:HB	36:35:60:MET:HB2	1.76	0.67
29:21:120:TRP:CD2	29:21:155:LYS:HD3	2.29	0.67
26:14:34:C:HO2'	26:14:35:G:P	2.16	0.67
1:1G:538:G:H5''	12:3A:114:LYS:HB2	1.77	0.67
26:1H:299:A:H5'	26:1H:300:A:OP2	1.94	0.67
28:11:181:GLU:HA	28:11:272:ALA:HB3	1.75	0.67
2:12:144:ARG:HH21	2:12:148:TYR:HD2	1.43	0.67
26:14:2709:G:N3	58:14:3434:HOH:O	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:12:19:HIS:NE2	2:12:205:ASP:OD1	2.25	0.67
1:13:1305:G:N2	1:13:1331:G:H2'	2.09	0.67
26:1H:2364:C:H4'	47:I8:56:ASP:OD1	1.95	0.67
30:39:4:VAL:HA	30:39:19:GLU:HB3	1.77	0.67
33:61:110:ASP:HB2	33:61:112:LYS:H	1.59	0.67
1:1G:513:C:H42	1:1G:538:G:H1	1.40	0.67
22:1K:76:A:H8	26:1H:2507:C:H1'	1.58	0.67
26:14:1784:A:H5''	58:14:4067:HOH:O	1.93	0.67
1:13:396:G:O2'	1:13:398:C:OP1	2.09	0.67
26:1H:2074:U:P	58:1H:4530:HOH:O	2.53	0.67
1:13:156:G:H1	1:13:165:C:H42	1.42	0.67
43:E8:60:ASN:HD22	43:E8:60:ASN:N	1.90	0.67
26:1H:751:A:OP1	58:1H:3922:HOH:O	2.11	0.67
26:14:2688:U:H5	26:14:2720:U:OP2	1.77	0.67
26:1H:2262:U:OP1	26:1H:2387:U:O2'	2.10	0.67
26:1H:2287:A:H62	26:1H:2344:U:H3	1.39	0.67
1:1G:1306:A:N6	1:1G:1331:G:O2'	2.27	0.67
4:3E:167:GLY:HA2	28:19:135:PHE:HE1	1.60	0.67
46:D5:17:ALA:HA	46:D5:20:ARG:HD2	1.76	0.67
46:H8:63:ASP:OD2	46:H8:65:GLN:NE2	2.27	0.67
26:1H:155:C:H42	26:1H:171:G:H1	1.40	0.67
1:1G:793:U:OP1	58:1G:1859:HOH:O	2.10	0.67
1:13:1348:U:H4'	9:8E:120:ARG:HD2	1.74	0.67
26:14:1047:G:H2'	26:14:1110:G:H1	1.58	0.67
37:88:72:LYS:HB3	37:88:94:VAL:HG23	1.75	0.67
45:G8:76:CYS:SG	45:G8:97:ARG:HG2	2.35	0.67
1:13:624:C:O3'	16:7I:10:GLY:HA2	1.93	0.67
1:1G:963:G:N3	10:1A:55:LYS:NZ	2.41	0.67
40:B8:56:GLY:O	40:B8:59:THR:HG22	1.94	0.67
1:13:405:U:O4	4:3E:2:GLY:N	2.28	0.67
26:1H:2410:G:N7	58:1H:3729:HOH:O	2.28	0.67
26:14:1406:U:H5''	26:14:1406:U:H6	1.58	0.67
26:1H:587:C:O2	36:78:33:ARG:NH1	2.27	0.67
26:1H:1997:G:H5''	58:1H:4005:HOH:O	1.94	0.67
26:1H:1697:G:OP2	26:1H:1698:A:O2'	2.07	0.67
24:3L:76:A:H8	26:14:2394:C:H42	1.38	0.67
26:1H:1517:G:H5''	26:1H:1518:C:OP2	1.94	0.67
33:61:110:ASP:HB2	33:61:112:LYS:HG2	1.76	0.67
12:3I:90:VAL:HG12	12:3I:92:ASP:H	1.59	0.67
1:1G:186(A):C:H5''	20:BA:86:ARG:HH21	1.59	0.67
1:13:395:C:N4	58:13:1901:HOH:O	2.28	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:353:A:H5'	1:13:353:A:H8	1.58	0.67
26:1H:1646:C:O3'	58:1H:3926:HOH:O	2.13	0.67
26:14:1130:U:O2	29:29:149:ARG:NH2	2.27	0.67
26:14:305:U:H2'	26:14:306:U:C6	2.29	0.67
31:49:124:SER:HB2	31:49:131:TYR:CE2	2.30	0.67
26:14:1971:A:OP1	58:14:3452:HOH:O	2.12	0.67
1:1G:625:G:H2'	1:1G:626:U:H6	1.60	0.67
26:14:1416:G:O2'	26:14:1417:C:O5'	2.12	0.67
26:1H:1510:A:O2'	26:1H:1512:G:N7	2.27	0.67
26:14:2286:A:H2'	53:K5:28:ARG:HH21	1.60	0.67
53:O8:10:LEU:HD23	55:Q8:32:LEU:HD22	1.76	0.67
26:1H:1418:G:OP2	58:1H:4023:HOH:O	2.10	0.67
1:13:1182:G:H4'	1:13:1183:A:H5''	1.74	0.67
22:1K:51:U:H2'	22:1K:52:G:H8	1.59	0.67
1:13:1266:G:N7	58:13:2005:HOH:O	2.27	0.67
26:1H:2690:C:H5''	26:1H:2872:G:H21	1.60	0.67
46:D5:4:ARG:HA	46:D5:58:VAL:HB	1.76	0.67
35:68:63:VAL:HG12	35:68:106:LEU:HD11	1.75	0.67
6:52:81:ILE:HD11	28:19:125:ILE:HG12	1.77	0.67
26:14:1226:G:OP1	42:95:69:LYS:NZ	2.26	0.67
26:14:1024:G:H3'	26:14:1025:G:H5''	1.77	0.67
27:1J:13:A:H5''	27:1J:15:A:C6	2.29	0.67
10:1A:40:LEU:HG	10:1A:41:PRO:HD2	1.75	0.67
1:13:272:C:H2'	1:13:273:A:C8	2.28	0.67
38:55:33:ARG:NH2	38:55:115:GLU:OE2	2.26	0.67
26:14:987:G:O2'	26:14:1000:A:N3	2.28	0.67
38:55:74:LYS:HE3	38:55:77:ARG:HH21	1.59	0.67
26:1H:1633:G:OP2	58:1H:4371:HOH:O	2.12	0.67
38:55:86:ARG:NH2	38:55:87:TYR:OH	2.27	0.67
24:3L:71:G:O2'	26:14:1851:U:O2'	2.01	0.67
3:2E:84:ILE:HG13	3:2E:101:LEU:HD22	1.75	0.67
1:13:1216:G:OP2	58:13:1998:HOH:O	2.11	0.67
1:1G:1503:A:N3	25:4L:13:A:N6	2.42	0.67
26:1H:1657:C:OP2	29:21:136:ARG:HG3	1.94	0.67
1:1G:1126:U:H4'	1:1G:1127:G:N7	2.10	0.67
26:14:399:G:OP2	58:14:4118:HOH:O	2.12	0.67
26:14:1754:C:H2'	26:14:1755:A:C8	2.29	0.67
26:1H:2074:U:OP1	58:1H:4530:HOH:O	2.12	0.67
13:4I:3:ARG:HE	13:4I:9:ILE:HD11	1.60	0.67
30:31:127:GLU:HG2	30:31:196:LEU:HB2	1.75	0.67
26:1H:2788:C:O2'	26:1H:2809:A:N3	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1419:A:H2'	26:14:1421:G:N7	2.09	0.67
1:13:417:C:H2'	1:13:418:C:H6	1.60	0.67
37:88:21:THR:HG22	37:88:99:PRO:O	1.95	0.67
3:22:87:LEU:HA	3:22:90:GLU:HG2	1.76	0.67
29:21:105:THR:HG21	29:21:164:ARG:NH1	2.10	0.66
26:1H:1509:C:H2'	26:1H:1511:A:C8	2.29	0.66
1:1G:1127:G:H22	1:1G:1144:G:H1	1.43	0.66
11:2I:98:LEU:O	11:2I:101:SER:OG	2.11	0.66
26:14:2598:A:OP1	58:14:4081:HOH:O	2.12	0.66
1:13:490:G:OP2	4:3E:132:ARG:NH2	2.26	0.66
1:1G:310:G:OP2	16:7A:27:LYS:NZ	2.21	0.66
5:42:122:GLU:O	5:42:126:ARG:NH1	2.28	0.66
49:K8:32:LEU:HD11	49:K8:54:LYS:HG2	1.75	0.66
26:1H:731:C:OP2	58:1H:3689:HOH:O	2.13	0.66
37:88:12:GLN:HG2	37:88:73:PRO:HD2	1.76	0.66
19:AA:9:VAL:HG13	51:I5:63:TYR:HE1	1.59	0.66
26:1H:567:A:OP1	58:1H:3601:HOH:O	2.13	0.66
26:1H:1298:C:H5''	26:1H:1299:G:OP2	1.96	0.66
26:1H:446:G:OP2	58:1H:3702:HOH:O	2.13	0.66
26:1H:1636:C:OP2	58:1H:3609:HOH:O	2.13	0.66
26:14:1111:A:H4'	32:59:3:ARG:HH11	1.61	0.66
1:1G:957:U:H1'	1:1G:960:U:H5	1.61	0.66
30:39:123:LEU:O	30:39:125:LEU:N	2.27	0.66
1:1G:411:A:C5	1:1G:413:G:H1'	2.31	0.66
26:1H:138:G:N2	44:F8:44:GLU:OE2	2.22	0.66
36:35:125:VAL:O	36:35:144:GLU:HB3	1.96	0.66
1:13:1118:C:H1'	1:13:1179:A:C4	2.31	0.66
26:14:760:G:OP1	58:14:4111:HOH:O	2.12	0.66
26:1H:298:G:P	45:G8:84:ARG:HH12	2.18	0.66
39:A8:88:ASP:O	39:A8:89:ARG:HB3	1.95	0.66
6:5E:38:GLU:OE1	6:5E:64:GLN:NE2	2.27	0.66
46:H8:17:ALA:HA	46:H8:20:ARG:HB2	1.78	0.66
26:14:2873:A:H8	38:55:6:SER:N	1.92	0.66
26:1H:1021:A:H62	26:1H:1141:U:H3	1.43	0.66
26:1H:991:C:H2'	26:1H:992:C:H6	1.61	0.66
26:1H:2492:U:H2'	26:1H:2493:U:H6	1.58	0.66
33:61:7:GLU:HA	33:61:15:VAL:HG22	1.76	0.66
47:E5:49:LYS:HG3	47:E5:80:HIS:HB3	1.77	0.66
26:14:1443:G:N7	58:14:3772:HOH:O	2.28	0.66
33:69:76:THR:HG21	33:69:140:LEU:HD12	1.77	0.66
22:1L:74:C:H1'	22:1L:75:C:H5'	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1438:U:H2'	26:1H:1439:A:C8	2.30	0.66
38:55:38:VAL:HG12	38:55:42:LYS:HD2	1.77	0.66
31:41:131:TYR:O	31:41:159:VAL:HG22	1.96	0.66
26:14:2598:A:OP1	58:14:4079:HOH:O	2.11	0.66
29:21:70:ALA:O	29:21:73:GLU:N	2.29	0.66
35:25:68:GLU:HB3	35:25:78:ARG:HD3	1.77	0.66
26:1H:250:G:H2'	26:1H:251:A:C8	2.31	0.66
26:1H:1165:U:H2'	26:1H:1166:C:C6	2.31	0.66
26:1H:1153:C:OP2	58:1H:4003:HOH:O	2.12	0.66
39:A8:14:VAL:HG21	39:A8:89:ARG:HE	1.59	0.66
33:69:70:GLU:O	33:69:74:ASN:ND2	2.28	0.66
26:14:1949:G:O6	58:14:3816:HOH:O	2.10	0.66
49:K8:15:LYS:H	49:K8:67:LYS:NZ	1.94	0.66
26:14:2080:G:H5'	48:F5:35:THR:OG1	1.95	0.66
43:E8:70:TYR:H	43:E8:70:TYR:HD1	1.44	0.66
1:13:143:A:H2	1:13:220:G:H1	1.43	0.66
26:14:1225:C:O2'	42:95:85:LYS:N	2.28	0.66
26:1H:2701:C:H2'	26:1H:2702:U:H2'	1.77	0.66
26:14:531:C:H4'	26:14:532:A:H5''	1.78	0.66
26:1H:1332:G:N2	26:1H:1610:A:C8	2.64	0.66
1:13:630:G:N2	58:13:1982:HOH:O	2.28	0.66
26:1H:1257:C:H4'	30:31:83:PHE:CE1	2.31	0.66
24:3L:56:C:H5'	26:14:2169:A:C8	2.30	0.66
39:65:34:HIS:NE2	39:65:54:LEU:HD13	2.11	0.66
29:29:65:GLY:O	29:29:67:PHE:N	2.27	0.66
26:1H:2659:G:H4'	32:51:175:LYS:HD2	1.77	0.66
51:I5:48:ARG:HH22	51:I5:51:ASP:HB3	1.61	0.66
29:21:82:ARG:O	29:21:84:PHE:N	2.28	0.66
26:14:900:A:H3'	26:14:901:A:H8	1.61	0.66
48:J8:87:PRO:HA	48:J8:90:ILE:HG13	1.78	0.66
26:1H:2701:C:H3'	26:1H:2702:U:H5''	1.78	0.66
1:13:1160:G:H1	1:13:1177:G:H22	1.44	0.66
26:1H:1331:A:O3'	58:1H:4559:HOH:O	2.13	0.66
32:51:54:ARG:HD3	32:51:65:HIS:ND1	2.11	0.66
26:14:39:C:O2	30:39:46:ARG:NH2	2.29	0.66
26:1H:1693:U:H1'	28:11:14:ARG:NH2	2.11	0.66
43:A5:13:SER:HB3	43:A5:16:LYS:HD2	1.78	0.66
10:1A:48:THR:HA	10:1A:62:HIS:HB3	1.76	0.66
1:13:1429:C:H2'	1:13:1430:C:C6	2.31	0.66
26:14:1599:C:H2'	26:14:1600:C:H6	1.59	0.66
26:1H:1334:G:OP2	58:1H:3955:HOH:O	2.14	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:G8:81:LYS:HB3	45:G8:82:PRO:HA	1.78	0.66
26:14:987:G:OP2	58:14:4040:HOH:O	2.14	0.66
26:14:1464:C:HO2'	26:14:1528:A:H8	1.41	0.66
40:75:77:PRO:HG2	40:75:80:SER:HB2	1.77	0.66
5:42:70:PRO:HB3	5:42:144:THR:HG22	1.77	0.66
26:1H:213:A:OP2	58:1H:4290:HOH:O	2.14	0.66
51:M8:65:ASP:N	51:M8:65:ASP:OD1	2.28	0.66
17:8A:5:VAL:HG22	17:8A:60:ILE:HG12	1.78	0.66
46:D5:152:ALA:HB3	46:D5:167:PRO:HA	1.78	0.66
51:I5:11:PRO:HA	51:I5:25:TYR:HA	1.77	0.66
1:13:376:G:O3'	16:7I:5:ARG:NH1	2.29	0.66
4:32:24:GLU:OE2	4:32:24:GLU:N	2.28	0.65
55:Q8:6:THR:H	55:Q8:59:LYS:HE3	1.60	0.65
1:13:835:U:H3	1:13:851:G:H1	1.44	0.65
39:A8:28:VAL:HG11	39:A8:98:VAL:HG12	1.77	0.65
1:13:627:G:H2'	1:13:628:G:H8	1.61	0.65
22:1K:76:A:H2	26:1H:2602:A:C2	2.13	0.65
26:14:780:G:H21	26:14:783:A:H62	1.44	0.65
19:AI:39:THR:HG22	19:AI:40:ILE:H	1.60	0.65
36:35:113:LYS:HD3	36:35:115:LEU:HD21	1.78	0.65
4:32:8:VAL:HA	4:32:11:LEU:HD12	1.78	0.65
2:12:70:PHE:O	2:12:93:VAL:N	2.20	0.65
26:1H:570:G:OP2	58:1H:3873:HOH:O	2.12	0.65
40:75:20:PRO:HD2	40:75:86:ILE:HG23	1.78	0.65
26:14:1079:C:N4	26:14:1088:A:OP1	2.29	0.65
26:1H:1671:U:OP2	58:1H:3665:HOH:O	2.13	0.65
26:1H:770:G:OP2	58:1H:4400:HOH:O	2.13	0.65
26:1H:2165:G:N7	26:1H:2166:G:N2	2.44	0.65
30:31:67:GLN:HG3	30:31:67:GLN:O	1.96	0.65
1:1G:1145:C:O2'	1:1G:1146:A:N7	2.29	0.65
45:G8:87:LYS:O	45:G8:94:LYS:HB2	1.95	0.65
37:45:75:THR:HB	37:45:86:GLY:HA3	1.78	0.65
26:14:1847:A:OP1	26:14:1847:A:H8	1.79	0.65
26:1H:2789:C:O2	26:1H:2894:G:N2	2.28	0.65
13:4A:66:LEU:HA	13:4A:70:LEU:HB2	1.77	0.65
1:13:1062:U:H2'	1:13:1063:C:C6	2.30	0.65
10:1I:46:ARG:NH2	10:1I:64:GLU:OE1	2.29	0.65
27:1J:6:C:H2'	27:1J:7:G:H5''	1.78	0.65
55:Q8:7:HIS:O	55:Q8:7:HIS:ND1	2.29	0.65
32:51:6:ARG:HB3	32:51:65:HIS:CG	2.32	0.65
26:14:1073:A:OP2	26:14:1094:U:N3	2.29	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1056:G:H21	26:1H:1103:A:H62	1.44	0.65
1:13:292:G:N7	1:13:293:G:H1'	2.12	0.65
26:14:270(Q):C:H5''	33:69:45:LYS:HE3	1.78	0.65
22:1K:14:A:N6	22:1K:22:G:N3	2.44	0.65
26:1H:1586:A:H3'	26:1H:1587:A:H8	1.60	0.65
28:19:16:MET:HE1	28:19:208:LYS:HG2	1.77	0.65
1:13:191(F):U:H2'	1:13:191:G:H8	1.61	0.65
26:1H:453:C:OP1	58:1H:3896:HOH:O	2.15	0.65
1:13:191(F):U:H2'	1:13:191:G:C8	2.31	0.65
51:M8:52:THR:OG1	51:M8:53:GLU:N	2.29	0.65
34:15:47:ALA:HB2	34:15:112:LEU:HD21	1.78	0.65
1:1G:940:C:H2'	1:1G:941:G:C8	2.32	0.65
34:58:56:ASN:N	34:58:125:GLY:O	2.15	0.65
36:35:47:ASP:OD2	36:35:50:ARG:NH1	2.29	0.65
49:K8:5:GLU:OE2	49:K8:5:GLU:HA	1.93	0.65
5:42:143:ARG:NH1	8:72:77:GLU:OE2	2.26	0.65
5:4E:142:LEU:O	5:4E:143:ARG:NH1	2.28	0.65
26:14:910:A:C5	37:45:13:GLN:HG3	2.31	0.65
32:51:4:ILE:HG21	32:51:6:ARG:NH1	2.12	0.65
4:3E:22:LYS:HB2	4:3E:26:CYS:SG	2.37	0.65
1:13:926:G:H5'	1:13:927:G:O5'	1.96	0.65
1:1G:359:U:H2'	1:1G:360:A:C8	2.31	0.65
26:1H:646:A:H2'	26:1H:647:G:O4'	1.96	0.65
26:1H:2849:U:H4'	26:1H:2868:A:C2	2.32	0.65
26:1H:1024:G:H3'	26:1H:1025:G:H5''	1.78	0.65
4:3E:88:VAL:HB	4:3E:91:SER:HB2	1.78	0.65
9:8E:97:LYS:HB2	9:8E:102:LEU:HD12	1.79	0.65
40:B8:3:ARG:O	40:B8:7:ILE:N	2.27	0.65
26:1H:187:G:N7	58:1H:4406:HOH:O	2.29	0.65
26:1H:1386:C:H2'	26:1H:1387:C:C6	2.31	0.65
1:13:1352:C:H2'	1:13:1353:G:C8	2.31	0.65
26:1H:2032:G:H21	29:21:146:THR:HG23	1.61	0.65
8:72:29:SER:HB3	8:72:32:LYS:HG3	1.77	0.65
26:14:2681:C:H5	26:14:2725:A:H62	1.42	0.65
1:13:859:A:H2'	1:13:860:A:H8	1.60	0.65
40:75:98:LYS:HB3	40:75:100:TYR:CE2	2.31	0.65
1:13:407:G:OP1	4:3E:115:ARG:NH1	2.30	0.65
24:3K:22:G:N7	24:3K:46:G:N2	2.45	0.65
26:14:2718:G:N7	58:14:3760:HOH:O	2.29	0.65
26:1H:731:C:H5''	58:1H:3827:HOH:O	1.96	0.65
26:1H:2887:U:H2'	26:1H:2888:C:C6	2.31	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2837:G:N7	58:1H:4253:HOH:O	2.28	0.65
32:51:12:PRO:HG2	32:51:13:LYS:HG2	1.78	0.65
26:14:2150:U:H2'	26:14:2151:G:H8	1.62	0.65
48:F5:14:VAL:HG11	48:F5:39:LYS:HD3	1.79	0.65
4:32:156:GLU:HA	4:32:159:ARG:HD2	1.78	0.65
26:14:1211:U:H2'	26:14:1211:U:OP1	1.97	0.65
33:61:144:VAL:HG22	33:61:145:VAL:HG23	1.79	0.65
5:4E:91:LEU:HD12	5:4E:120:THR:HG22	1.79	0.65
55:Q8:49:VAL:HG12	55:Q8:52:LYS:CB	2.27	0.65
26:1H:1253:A:C8	58:1H:3713:HOH:O	2.48	0.65
26:14:2615:U:C2	52:J5:7:PRO:HA	2.32	0.65
26:1H:945:A:N3	58:1H:3865:HOH:O	2.28	0.65
1:1G:1316:G:N2	1:1G:1319:A:H5''	2.11	0.65
26:1H:1889:A:N1	26:1H:2234:G:H1'	2.12	0.65
46:D5:60:GLU:HA	46:D5:66:SER:HA	1.78	0.65
26:1H:439:G:O6	58:1H:4674:HOH:O	2.10	0.65
23:2K:3:C:H42	23:2K:71:G:H1	1.44	0.65
1:1G:1179:A:H4'	9:82:103:THR:HA	1.78	0.65
28:11:108:PRO:HG3	28:11:143:HIS:CE1	2.31	0.65
26:1H:2263:C:OP2	58:1H:4140:HOH:O	2.14	0.65
29:29:8:LYS:HB3	29:29:193:GLY:H	1.62	0.65
13:4A:37:THR:O	13:4A:55:ARG:NH2	2.29	0.65
1:13:74:C:N4	1:13:96:G:H1	1.93	0.65
1:13:974:A:OP2	14:5I:41:ARG:NH1	2.30	0.65
1:13:505:G:OP1	58:13:1880:HOH:O	2.14	0.65
26:14:2836:U:H2'	26:14:2837:G:C8	2.32	0.65
26:1H:2331:G:O3'	47:I8:43:THR:HG22	1.95	0.65
32:51:149:ARG:HA	32:51:162:ILE:HG21	1.78	0.65
14:5I:9:LYS:HA	14:5I:12:ARG:HG2	1.78	0.65
26:1H:1377:G:OP2	58:1H:4125:HOH:O	2.15	0.65
19:AA:63:THR:OG1	19:AA:65:ASN:O	2.15	0.65
30:31:160:ASN:OD1	30:31:163:VAL:HG23	1.96	0.65
1:13:339:C:OP2	35:68:97:ARG:NH1	2.28	0.65
5:42:82:VAL:HG12	5:42:89:ILE:HG22	1.77	0.65
1:13:649:G:H2'	1:13:650:G:H8	1.62	0.65
42:95:85:LYS:CD	42:95:86:GLY:H	2.08	0.65
26:1H:1164:G:H2'	26:1H:1165:U:H6	1.60	0.65
1:13:963:G:H21	10:1I:55:LYS:HE2	1.62	0.65
38:98:63:ARG:HB2	38:98:80:PHE:HE2	1.62	0.65
26:14:2064:C:H2'	26:14:2065:C:C6	2.32	0.65
27:1J:4:C:H42	27:1J:116:G:H1	1.45	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:555:C:H2'	1:1G:556:C:C6	2.32	0.65
26:14:38:A:H1'	30:39:48:THR:HB	1.77	0.65
50:H5:46:ASN:O	50:H5:50:VAL:HG22	1.97	0.65
26:1H:779:U:OP1	28:11:49:ILE:HG13	1.95	0.65
26:1H:847:U:OP2	58:1H:3755:HOH:O	2.14	0.65
36:78:13:ASN:ND2	36:78:15:ARG:HD3	2.12	0.65
26:14:1945:G:H2'	26:14:1946:U:H6	1.61	0.65
26:1H:270(L):U:H3	33:61:50:ARG:HG2	1.62	0.65
10:1A:51:ARG:HB2	10:1A:60:ARG:HA	1.78	0.65
26:14:2453:A:H61	26:14:2500:U:H3	1.44	0.65
29:29:33:VAL:HG12	29:29:89:ASP:HB3	1.77	0.65
26:14:801:G:OP2	58:14:3736:HOH:O	2.14	0.65
55:Q8:54:GLU:O	55:Q8:56:GLU:N	2.30	0.64
26:14:2415:G:H4'	36:35:67:MET:N	2.11	0.64
40:75:55:ASN:N	40:75:59:THR:HG22	2.12	0.64
29:21:64:LYS:O	29:21:70:ALA:HB2	1.97	0.64
19:AI:30:LEU:HD22	19:AI:30:LEU:H	1.62	0.64
26:1H:1400:G:H2'	26:1H:1401:G:C8	2.33	0.64
26:1H:2855:C:H2'	26:1H:2856:C:H6	1.61	0.64
47:I8:49:LYS:HG3	47:I8:80:HIS:HB3	1.80	0.64
42:95:44:LYS:O	42:95:46:VAL:N	2.24	0.64
11:2A:33:THR:OG1	11:2A:34:ASP:O	2.16	0.64
1:1G:458:C:H42	1:1G:474:G:H1	1.44	0.64
1:1G:54:C:N4	1:1G:353:A:OP2	2.29	0.64
13:4I:3:ARG:HB3	13:4I:9:ILE:HG12	1.79	0.64
1:13:1000:A:H2'	1:13:1001:G:H8	1.61	0.64
29:21:3:GLY:HA3	29:21:81:ILE:HG21	1.79	0.64
26:1H:1932:A:H2'	26:1H:1933:G:O4'	1.96	0.64
1:13:474:G:H2'	1:13:475:G:C8	2.33	0.64
11:2I:99:GLN:HG2	11:2I:105:VAL:HG21	1.78	0.64
3:22:182:ILE:HG22	3:22:203:PHE:HA	1.80	0.64
3:22:81:GLY:HA2	3:22:85:ARG:HD3	1.79	0.64
38:98:51:LEU:HD22	38:98:66:VAL:HG13	1.79	0.64
36:78:101:VAL:HG12	36:78:106:LEU:HD12	1.79	0.64
1:13:1346:A:H5''	9:8E:120:ARG:HH12	1.61	0.64
30:31:6:VAL:HG21	30:31:119:ARG:HB2	1.79	0.64
28:19:148:GLU:HB2	28:19:151:LYS:HD2	1.78	0.64
46:D5:139:VAL:HG22	46:D5:156:LYS:HG2	1.78	0.64
26:14:2261:C:O2'	26:14:2262:U:H5'	1.97	0.64
26:14:1332:G:N2	26:14:1610:A:C8	2.65	0.64
28:11:177:LEU:HD11	28:11:183:ARG:HG2	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:16:90:C:H5'	37:88:18:LYS:HA	1.77	0.64
3:22:91:LEU:O	3:22:95:THR:OG1	2.15	0.64
48:F5:87:PRO:HA	48:F5:90:ILE:HG13	1.78	0.64
1:13:1505:G:OP1	58:13:1804:HOH:O	2.14	0.64
1:13:1002:G:O6	1:13:1038:C:N4	2.30	0.64
26:1H:2061:G:OP2	26:1H:2502:G:H5'	1.98	0.64
26:14:2471:C:N4	26:14:2476:A:O2'	2.31	0.64
1:13:1429:C:H2'	1:13:1430:C:H6	1.61	0.64
44:B5:36:LYS:HG2	44:B5:54:VAL:HB	1.79	0.64
1:13:939:G:H5''	7:6E:102:ARG:NH2	2.12	0.64
26:1H:654(H):G:N7	26:1H:654(N):G:N2	2.45	0.64
1:13:501:C:OP2	12:3I:124:LYS:NZ	2.25	0.64
1:13:1059:C:O3'	14:5I:45:ARG:NH2	2.31	0.64
36:78:1:MET:HE2	36:78:6:LEU:HD13	1.78	0.64
26:1H:1230:C:H2'	26:1H:1231:G:C8	2.32	0.64
26:1H:2161:C:H2'	26:1H:2162:G:H8	1.61	0.64
1:13:1190:G:OP1	3:2E:4:LYS:HA	1.98	0.64
42:95:14:VAL:HB	42:95:96:ILE:HG13	1.79	0.64
26:1H:1702:G:O6	58:1H:4393:HOH:O	2.12	0.64
26:1H:450:G:OP2	58:1H:3889:HOH:O	2.15	0.64
26:14:1113:U:OP1	32:59:3:ARG:N	2.31	0.64
10:1A:8:LEU:HD22	10:1A:20:ALA:HB2	1.80	0.64
30:31:152:GLU:HG3	30:31:191:ARG:HD2	1.80	0.64
26:1H:2306:C:H3'	26:1H:2307:G:H5'	1.79	0.64
1:13:323:U:H2'	1:13:324:G:O4'	1.98	0.64
2:1E:141:GLU:O	2:1E:145:LEU:HB2	1.98	0.64
26:1H:1971:A:C4	28:11:241:PRO:HD3	2.33	0.64
26:14:1040:C:O2	26:14:1115:G:N2	2.20	0.64
26:14:884:C:N4	26:14:892:G:O6	2.30	0.64
18:9I:59:SER:OG	18:9I:60:ALA:N	2.30	0.64
55:Q8:34:TRP:HB3	55:Q8:35:GLN:HA	1.78	0.64
1:1G:1126:U:H4'	1:1G:1127:G:C8	2.33	0.64
23:2L:20:G:C2	23:2L:58:A:C2	2.86	0.64
20:BA:87:LYS:HE3	20:BA:91:LEU:HD11	1.79	0.64
5:42:36:ASP:HB2	5:42:38:GLN:HG3	1.80	0.64
4:32:127:THR:HG21	4:32:149:ALA:HB2	1.79	0.64
1:13:881:G:OP2	12:3I:12:ARG:NH2	2.30	0.64
20:BI:86:ARG:NH1	20:BI:90:GLN:OE1	2.30	0.64
1:13:991:U:O2'	1:13:992:U:O5'	2.15	0.64
26:1H:2080:G:O6	58:1H:4320:HOH:O	2.11	0.64
26:14:996:A:N6	26:14:1160:G:C6	2.66	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:3K:34:G:N7	24:3K:35:A:H8	1.95	0.64
24:3L:36:A:N7	24:3L:37:A:N6	2.46	0.64
26:1H:2352:A:H2'	26:1H:2353:G:O4'	1.98	0.64
26:1H:1533:C:H3'	26:1H:1534:G:H5''	1.80	0.64
9:8E:121:ARG:NH1	9:8E:122:ALA:O	2.31	0.64
26:1H:495:G:O2'	43:E8:62:HIS:HE1	1.80	0.64
55:Q8:49:VAL:CG2	55:Q8:53:PRO:HD2	2.27	0.64
26:1H:250:G:P	36:78:60:MET:HE1	2.38	0.64
26:14:2304:G:O2'	31:49:156:ASP:OD1	2.11	0.64
45:C5:17:SER:HB3	45:C5:71:LYS:HB3	1.79	0.64
29:29:33:VAL:HG11	29:29:36:ARG:HH21	1.62	0.64
26:1H:1534:G:H2'	26:1H:1535:U:H4'	1.80	0.64
2:1E:178:ARG:HG3	8:7E:72:PRO:HA	1.80	0.64
3:22:136:GLN:O	3:22:140:ARG:N	2.29	0.64
7:6E:50:ILE:HB	7:6E:58:PRO:HB3	1.79	0.64
41:C8:8:VAL:HG23	41:C8:11:ARG:NH2	2.11	0.64
26:14:994:C:OP2	41:85:54:LYS:NZ	2.27	0.64
26:14:1022:G:H22	26:14:1142(A):A:H2	1.45	0.64
1:1G:297:G:N2	1:1G:300:A:OP2	2.31	0.64
3:22:152:ILE:HB	3:22:199:LYS:HB2	1.80	0.64
4:32:4:TYR:CE2	4:32:11:LEU:HD11	2.33	0.64
1:13:477:G:N7	58:13:2008:HOH:O	2.29	0.64
26:1H:547:A:H2'	26:1H:548:A:C8	2.33	0.64
15:6A:54:ARG:NH1	15:6A:58:MET:SD	2.71	0.64
26:14:590:A:H2'	26:14:591:C:C6	2.33	0.64
41:85:100:VAL:O	41:85:101:ARG:HG2	1.98	0.64
30:39:129:PHE:HA	30:39:142:TRP:CD1	2.33	0.64
45:C5:87:LYS:H	45:C5:94:LYS:HG2	1.62	0.64
3:22:14:ILE:HG12	3:22:15:THR:H	1.62	0.64
32:51:83:TYR:CG	32:51:134:SER:HA	2.33	0.63
26:14:619:G:H5''	26:14:620:G:H21	1.61	0.63
26:14:2074:U:H2'	26:14:2075:U:C6	2.33	0.63
44:B5:55:ASN:HB2	44:B5:80:ILE:HG13	1.79	0.63
7:62:84:ASN:OD1	7:62:84:ASN:N	2.31	0.63
1:13:750:G:N3	15:6I:23:GLY:HA3	2.13	0.63
26:1H:2002:G:N7	58:1H:4225:HOH:O	2.30	0.63
3:2E:42:LEU:O	3:2E:46:GLU:HG2	1.98	0.63
26:14:2056:G:H1	52:J5:4:HIS:HB3	1.62	0.63
4:32:61:LYS:NZ	4:32:72:GLU:OE2	2.29	0.63
26:14:2646:C:H2'	26:14:2647:U:O4'	1.98	0.63
55:Q8:16:ILE:HD13	55:Q8:56:GLU:OE2	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1028:C:H42	1:1G:1033:G:H1	1.45	0.63
39:A8:29:PHE:HD1	39:A8:30:ARG:N	1.95	0.63
42:D8:59:ALA:HB2	42:D8:96:ILE:HD13	1.79	0.63
4:3E:95:GLY:O	4:3E:99:SER:OG	2.16	0.63
1:1G:1218:C:OP2	14:5A:9:LYS:NZ	2.26	0.63
2:1E:28:PHE:HD1	2:1E:194:PRO:HD3	1.63	0.63
1:1G:532:A:N6	1:1G:1206:G:O2'	2.31	0.63
26:1H:981:A:H8	26:1H:982:C:C5	2.15	0.63
13:4I:23:TYR:HD2	13:4I:67:GLU:HA	1.64	0.63
24:3L:48:C:N4	24:3L:59:U:O2	2.31	0.63
32:59:6:ARG:H	32:59:6:ARG:HD2	1.64	0.63
37:88:51:ARG:HH12	37:88:52:VAL:HG23	1.63	0.63
40:B8:58:ASN:C	40:B8:58:ASN:HD22	2.02	0.63
10:1I:40:LEU:HB2	10:1I:69:ASN:HB3	1.80	0.63
26:1H:2377:A:H2'	26:1H:2378:A:C8	2.33	0.63
17:8I:76:LEU:HD12	17:8I:77:VAL:N	2.13	0.63
13:4A:3:ARG:HB2	51:I5:34:GLU:HG3	1.80	0.63
26:1H:1430:C:H2'	26:1H:1431:U:H6	1.61	0.63
8:7E:87:SER:HB2	8:7E:93:VAL:N	2.13	0.63
27:1J:116:G:H5'	39:65:55:ALA:HB2	1.80	0.63
26:1H:1316:U:H2'	26:1H:1317:A:C8	2.34	0.63
45:G8:99:CYS:SG	45:G8:100:ALA:N	2.72	0.63
45:G8:100:ALA:HB1	45:G8:101:LYS:HB2	1.81	0.63
26:14:1593:G:H2'	26:14:1594:G:C8	2.33	0.63
26:1H:410:G:H3'	58:1H:4628:HOH:O	1.99	0.63
26:1H:1018:C:H2'	26:1H:1019:U:H6	1.64	0.63
36:35:11:GLY:C	36:35:13:ASN:H	2.01	0.63
26:1H:2068:U:N3	26:1H:2430:A:H2	1.94	0.63
55:M5:40:GLU:HA	55:M5:43:GLN:HB2	1.79	0.63
1:13:686:U:O4	1:13:703:G:H1'	1.99	0.63
1:1G:690:G:H1	11:2A:55:LYS:NZ	1.96	0.63
29:29:89:ASP:O	29:29:91:VAL:N	2.29	0.63
1:13:1263:C:H2'	1:13:1264:C:H6	1.64	0.63
1:13:200:G:H1	1:13:217:C:H42	1.45	0.63
26:14:779:U:O4	58:14:3960:HOH:O	2.10	0.63
38:98:104:ARG:NH1	38:98:107:ASP:OD2	2.32	0.63
26:1H:2056:G:C2	26:1H:2057:A:C8	2.87	0.63
42:95:6:LYS:H	42:95:37:VAL:HG12	1.64	0.63
34:58:6:PRO:HG3	34:58:41:ASP:HB2	1.80	0.63
26:14:1926:U:H2'	26:14:1928:A:OP2	1.99	0.63
5:4E:13:ILE:HD12	5:4E:30:ALA:HA	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:859:A:H2'	1:13:860:A:C8	2.34	0.63
2:12:77:ALA:HB2	2:12:211:ILE:HD13	1.79	0.63
26:14:588:U:H2'	26:14:589:C:C6	2.32	0.63
41:C8:28:ARG:NH1	41:C8:38:THR:OG1	2.25	0.63
23:2L:41:C:H2'	23:2L:42:C:H6	1.64	0.63
4:32:173:TRP:CZ3	4:32:193:ASP:HB3	2.34	0.63
26:1H:2360:A:OP1	55:Q8:48:PHE:HB2	1.98	0.63
26:1H:2594:C:N4	58:1H:3683:HOH:O	2.32	0.63
1:13:1505:G:P	58:13:1804:HOH:O	2.57	0.63
26:14:1828:G:OP2	58:14:3417:HOH:O	2.15	0.63
26:1H:2061:G:P	58:1H:3633:HOH:O	2.57	0.63
1:1G:998:G:H2'	1:1G:998(A):C:C6	2.34	0.63
32:59:6:ARG:HH11	32:59:6:ARG:N	1.97	0.63
26:1H:1405:U:H2'	26:1H:1406:U:C6	2.33	0.63
26:1H:141:A:C8	26:1H:1408:C:H1'	2.33	0.63
1:13:77:C:O2'	1:13:92:G:N2	2.32	0.63
46:D5:153:SER:HB3	46:D5:167:PRO:HG3	1.80	0.63
1:13:179:A:H2'	1:13:180:U:H6	1.64	0.63
34:15:95:PRO:O	34:15:98:VAL:HG22	1.98	0.63
38:55:52:ILE:O	38:55:55:ALA:N	2.32	0.63
34:58:18:ALA:HA	34:58:21:LYS:HG3	1.80	0.63
2:12:149:LEU:HD23	2:12:152:PHE:HB3	1.80	0.63
26:14:760:G:OP1	58:14:4109:HOH:O	2.15	0.63
29:29:47:VAL:HG21	29:29:86:PRO:HD2	1.81	0.63
13:4I:4:ILE:HG22	13:4I:5:ALA:H	1.63	0.63
9:8E:42:ARG:HH11	9:8E:71:SER:HB3	1.63	0.63
15:6I:39:LEU:HD13	15:6I:56:LEU:HB2	1.80	0.63
32:59:26:VAL:HG12	32:59:33:LEU:H	1.63	0.63
36:78:19:VAL:HG12	36:78:21:ARG:N	2.01	0.63
1:13:1178:G:N7	9:8E:97:LYS:NZ	2.46	0.63
30:39:117:ARG:NH1	36:35:1:MET:H2	1.96	0.63
23:2K:52:C:O2	23:2K:64:G:N2	2.28	0.63
37:88:66:ILE:HG22	37:88:67:ARG:H	1.64	0.63
23:2K:54:G:H2'	23:2K:55:5MU:C6	2.34	0.63
26:14:1784:A:H4'	26:14:1785:A:O5'	1.99	0.63
26:14:2645:G:H3'	26:14:2646:C:H5'	1.81	0.63
3:2E:40:ARG:O	3:2E:44:GLU:HG2	1.98	0.63
46:H8:73:GLN:HB2	46:H8:87:ASP:HB2	1.81	0.63
1:1G:677:U:H3	1:1G:713:G:H22	1.47	0.63
26:1H:1728:G:H8	26:1H:1732:A:H62	1.47	0.62
1:1G:707:C:H2'	1:1G:708:C:C6	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1406:U:H2'	26:1H:1407:C:C6	2.33	0.62
1:1G:554:C:H2'	1:1G:555:C:H6	1.62	0.62
1:1G:963:G:N2	1:1G:972:C:N3	2.44	0.62
19:AI:40:ILE:HD11	19:AI:62:ILE:HG23	1.79	0.62
51:I5:56:VAL:HG22	51:I5:57:GLU:HG3	1.81	0.62
1:13:1318:A:H1'	19:AI:37:ARG:HH21	1.64	0.62
1:13:131:C:OP1	1:13:190:G:N2	2.32	0.62
26:1H:1252:G:N3	41:C8:33:ARG:HD2	2.14	0.62
29:29:197:ILE:HD11	29:29:199:ARG:HE	1.64	0.62
7:62:92:SER:HB2	7:62:95:ARG:HB2	1.81	0.62
26:14:1329:U:H5''	26:14:1330:C:H5	1.65	0.62
1:1G:1139:G:H4'	1:1G:1140:C:H5'	1.80	0.62
26:14:2415:G:O3'	36:35:66:GLY:HA3	1.99	0.62
26:14:2275:C:O2'	37:45:83:MET:HA	1.99	0.62
26:1H:1798:U:C5'	28:11:259:THR:HG22	2.27	0.62
39:A8:74:ALA:HB1	39:A8:107:GLU:O	1.99	0.62
26:14:579:G:H2'	26:14:580:C:C6	2.34	0.62
26:1H:1101:U:H2'	26:1H:1102:C:C6	2.34	0.62
1:1G:760:G:N2	17:8A:94:ASN:OD1	2.32	0.62
7:62:21:VAL:HG23	7:62:22:LEU:HD12	1.79	0.62
31:49:60:LEU:O	31:49:64:THR:HG22	1.98	0.62
27:1J:62:C:H2'	27:1J:63:G:H8	1.62	0.62
32:59:82:GLY:HA3	32:59:135:GLY:O	1.99	0.62
26:1H:571:A:OP2	58:1H:4551:HOH:O	2.16	0.62
38:98:10:LEU:O	38:98:12:ARG:N	2.32	0.62
39:65:3:ARG:HE	39:65:4:LEU:H	1.47	0.62
1:13:674:G:H2'	1:13:675:A:H8	1.64	0.62
26:1H:49:A:N7	26:1H:120:U:C5	2.67	0.62
36:78:68:GLN:NE2	55:Q8:12:LYS:HB3	2.14	0.62
26:14:330:A:H2	26:14:1210:A:O2'	1.82	0.62
26:1H:1316:U:H2'	26:1H:1317:A:H8	1.64	0.62
45:C5:87:LYS:HB2	45:C5:94:LYS:HA	1.80	0.62
1:13:1316:G:N2	1:13:1318:A:H3'	2.15	0.62
26:1H:2096:U:H3	26:1H:2193:G:H1	1.47	0.62
27:1J:52:A:H62	39:65:33:LYS:HG3	1.63	0.62
23:2L:10:G:N2	23:2L:27:G:H1'	2.14	0.62
26:1H:1412:A:H2'	26:1H:1413:G:C8	2.35	0.62
50:L8:39:ASP:OD2	50:L8:44:ARG:NH1	2.33	0.62
7:6E:115:ARG:O	7:6E:118:VAL:HG12	1.98	0.62
26:1H:1729:A:H2'	26:1H:1731:G:N7	2.13	0.62
1:13:1286:A:C8	1:13:1287:A:H4'	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:67:U:H2'	26:14:68:G:C8	2.34	0.62
7:6E:95:ARG:NH2	7:6E:99:LEU:HD11	2.14	0.62
26:14:819:A:OP2	26:14:1187:G:N2	2.32	0.62
24:3L:5:G:H1	24:3L:68:C:H42	1.46	0.62
26:1H:1264:G:OP1	52:N8:19:ARG:NH2	2.31	0.62
37:45:34:LEU:HB2	37:45:118:LEU:HD13	1.81	0.62
35:68:7:TYR:CZ	35:68:44:LYS:HG3	2.34	0.62
26:14:990:A:H8	26:14:990:A:H5'	1.63	0.62
13:4I:10:PRO:HB2	13:4I:18:ALA:HB1	1.81	0.62
1:1G:1335:C:OP1	1:1G:1337:G:N2	2.30	0.62
50:L8:38:GLU:H	50:L8:38:GLU:CD	2.01	0.62
29:29:55:ASN:O	29:29:57:LYS:NZ	2.32	0.62
24:3L:30:G:H2'	24:3L:31:A:H8	1.65	0.62
26:14:1040:C:H2'	26:14:1041:C:C6	2.35	0.62
1:1G:1220:G:O3'	19:AA:36:ARG:HD3	1.99	0.62
27:1J:70:C:H2'	27:1J:71:C:H6	1.63	0.62
26:1H:862:G:OP2	58:1H:3992:HOH:O	2.16	0.62
26:1H:2057:A:P	58:1H:3621:HOH:O	2.56	0.62
1:1G:1279:A:O2'	1:1G:1282:C:N4	2.32	0.62
1:13:156:G:H1	1:13:165:C:N4	1.97	0.62
38:98:41:ALA:O	38:98:44:LEU:N	2.33	0.62
5:42:151:LEU:HD22	8:72:67:PRO:HD3	1.81	0.62
26:14:1340:U:H4'	26:14:1394:U:O2'	1.99	0.62
2:12:5:ILE:HA	2:12:221:LEU:HD21	1.80	0.62
4:32:98:GLU:OE2	4:32:103:ASN:ND2	2.31	0.62
26:14:491:G:H2'	26:14:492:A:C8	2.35	0.62
8:72:25:ASP:OD1	8:72:25:ASP:N	2.32	0.62
36:78:78:PRO:HB3	36:78:111:ARG:NH2	2.14	0.62
26:1H:998:C:H3'	58:1H:4004:HOH:O	1.99	0.62
45:G8:85:VAL:HG22	45:G8:98:VAL:HB	1.80	0.62
1:1G:940:C:H2'	1:1G:941:G:H8	1.62	0.62
1:13:474:G:H2'	1:13:475:G:H8	1.61	0.62
45:C5:48:ALA:HB3	45:C5:59:GLY:C	2.20	0.62
26:14:2291:U:H5''	26:14:2380:C:O2'	1.98	0.62
29:29:101:ARG:CZ	29:29:171:GLU:HB2	2.30	0.62
2:1E:5:ILE:HG13	2:1E:6:THR:N	2.14	0.62
26:14:2010:G:N7	58:14:3546:HOH:O	2.31	0.62
51:M8:13:ARG:H	51:M8:30:GLU:H	1.47	0.62
55:Q8:6:THR:HA	55:Q8:58:ILE:H	1.64	0.62
1:13:1346:A:H5''	9:8E:120:ARG:NH1	2.14	0.62
2:1E:163:PHE:CD1	2:1E:185:ILE:HG13	2.35	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:273(C):C:N4	26:14:363(C):G:H1	1.96	0.62
26:14:661:C:H1'	36:35:12:ALA:HA	1.81	0.62
34:58:4:TYR:O	41:C8:64:ARG:NH1	2.32	0.62
1:1G:1130:A:N6	1:1G:1144:G:H21	1.97	0.62
45:C5:68:HIS:HB3	45:C5:71:LYS:HG3	1.81	0.62
13:4A:23:TYR:CZ	13:4A:71:ARG:HG3	2.35	0.62
26:1H:270(I):G:H1	26:1H:270(Q):C:H42	1.47	0.62
16:7I:26:ARG:HH21	16:7I:31:LYS:HD3	1.63	0.62
1:13:1412:C:H42	1:13:1488:G:H1	1.48	0.62
10:1I:8:LEU:HD12	10:1I:20:ALA:HB2	1.80	0.62
1:1G:1423:G:H2'	1:1G:1424:C:C6	2.35	0.62
27:1J:8:U:H3	27:1J:112:G:H1	1.47	0.62
46:H8:10:ARG:HG3	46:H8:36:LYS:HB3	1.81	0.62
55:Q8:53:PRO:HB3	55:Q8:56:GLU:N	2.14	0.62
14:5I:29:ARG:NH2	14:5I:41:ARG:HH12	1.94	0.62
1:1G:1224:G:C6	1:1G:1322:C:H1'	2.34	0.62
14:5A:21:TYR:OH	14:5A:23:ARG:NH2	2.31	0.62
39:65:34:HIS:HD2	39:65:54:LEU:HB2	1.64	0.62
40:B8:54:ARG:HA	40:B8:59:THR:HB	1.82	0.62
19:AI:40:ILE:HG23	19:AI:41:VAL:HG13	1.81	0.62
2:12:7:VAL:HG22	2:12:8:LYS:H	1.63	0.62
27:1J:21:G:H2'	27:1J:22:U:O4'	2.00	0.62
20:BA:49:ALA:HA	20:BA:52:ALA:HB3	1.81	0.62
26:1H:2428:G:N3	36:78:61:ARG:NH1	2.48	0.62
42:95:87:HIS:CE1	42:95:89:GLN:HB2	2.35	0.62
26:1H:2698:U:H2'	26:1H:2699:C:C6	2.35	0.62
26:1H:851:U:P	50:L8:49:LYS:HZ2	2.22	0.62
37:88:86:GLY:HA3	37:88:87:LYS:HD3	1.81	0.62
2:12:12:GLU:HB3	2:12:213:LEU:HD13	1.82	0.62
7:62:62:PHE:HA	7:62:124:LEU:HD13	1.81	0.62
26:1H:2787:C:H1'	29:21:62:PRO:HG3	1.81	0.62
40:75:26:ASP:O	40:75:49:VAL:HG22	1.99	0.62
1:13:404:U:OP1	4:3E:118:ARG:NH1	2.33	0.62
39:65:15:ARG:O	39:65:19:LYS:HD3	2.00	0.62
50:L8:31:LEU:O	50:L8:32:GLN:HB2	2.00	0.62
33:69:80:PRO:HA	33:69:143:SER:HA	1.80	0.62
28:11:72:LYS:HD2	28:11:75:ILE:HD12	1.81	0.62
35:68:88:ASN:HD22	35:68:88:ASN:H	1.47	0.62
2:1E:74:LYS:HE3	2:1E:166:ASP:HB2	1.81	0.62
17:8I:65:ILE:HG21	17:8I:69:LYS:HE2	1.80	0.62
1:1G:300:A:H1'	1:1G:565:U:O2	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:980:C:H5'	1:1G:981:U:C5	2.35	0.62
7:62:113:GLU:HB2	7:62:119:ARG:HG2	1.80	0.62
51:M8:18:CYS:HB3	51:M8:39:CYS:SG	2.40	0.62
26:14:571:A:H5'	26:14:2030:A:N7	2.15	0.62
46:D5:23:LYS:HD3	46:D5:40:ASP:HA	1.81	0.62
1:1G:1007:C:O2	1:1G:1023:G:N2	2.33	0.62
23:2K:21:U:H4'	23:2K:22:A:OP1	1.98	0.62
12:3I:52:LEU:O	12:3I:54:LYS:NZ	2.32	0.62
41:85:76:TYR:CZ	41:85:80:ILE:HG13	2.35	0.61
17:8I:66:SER:OG	17:8I:69:LYS:HB2	1.99	0.61
5:4E:102:ALA:HB1	5:4E:106:PRO:HG2	1.81	0.61
24:3L:19:G:O5'	24:3L:57:G:N2	2.33	0.61
26:14:479:A:N3	26:14:481:G:H5''	2.15	0.61
40:B8:111:ARG:H	40:B8:111:ARG:HD3	1.65	0.61
1:13:1349:A:OP2	9:8E:118:LYS:NZ	2.33	0.61
51:M8:39:CYS:SG	51:M8:41:PRO:HD3	2.40	0.61
20:BA:86:ARG:CZ	20:BA:86:ARG:HB2	2.30	0.61
26:1H:2531:A:C8	32:51:175:LYS:HG2	2.34	0.61
1:13:1280:A:C8	10:1I:41:PRO:HD3	2.35	0.61
19:AI:5:LEU:HB3	19:AI:10:PHE:HE1	1.65	0.61
13:4A:49:THR:HG22	13:4A:51:ALA:H	1.65	0.61
10:1I:30:SER:OG	10:1I:84:GLN:NE2	2.33	0.61
26:1H:2793:G:N2	26:1H:2804:C:O2	2.33	0.61
33:61:1:MET:O	33:61:20:ASP:HA	1.98	0.61
29:21:149:ARG:HH11	29:21:149:ARG:HG3	1.64	0.61
34:15:38:HIS:CD2	34:15:39:ARG:HG3	2.35	0.61
26:1H:517:C:OP1	52:N8:16:ARG:NH2	2.33	0.61
26:1H:607:U:H3	26:1H:621:A:H2	1.42	0.61
2:12:47:THR:HA	2:12:202:PRO:HG2	1.80	0.61
24:3K:69:G:H2'	24:3K:70:G:C8	2.35	0.61
32:51:4:ILE:HG13	32:51:6:ARG:NE	2.14	0.61
44:F8:1:MET:O	44:F8:3:THR:HG23	1.99	0.61
1:13:1423:G:P	35:68:49:ARG:HH22	2.22	0.61
4:32:104:VAL:HA	4:32:107:ARG:HB2	1.82	0.61
35:25:25:LEU:HB2	35:25:38:VAL:HG23	1.82	0.61
30:31:103:LYS:HA	30:31:106:ARG:HG3	1.82	0.61
52:N8:58:LEU:HD22	52:N8:60:VAL:HG13	1.82	0.61
26:14:1252:G:N3	41:85:33:ARG:HD2	2.16	0.61
26:14:581:C:H2'	26:14:582:G:H8	1.65	0.61
10:1I:32:ALA:HB3	10:1I:76:ASN:O	2.00	0.61
26:1H:1843:C:H5'	28:11:253:GLN:OE1	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:72:19:VAL:HG23	8:72:21:LYS:HB2	1.82	0.61
1:1G:434:U:H2'	1:1G:435:C:C6	2.35	0.61
18:9A:22:VAL:HG12	18:9A:56:THR:HA	1.82	0.61
1:1G:407:G:O2'	4:32:116:GLN:HG3	2.01	0.61
26:14:676:A:H8	26:14:2069:G:N2	1.90	0.61
26:14:273(D):C:H42	26:14:363(B):G:H1	1.48	0.61
1:1G:1268:A:H2'	1:1G:1269:A:C8	2.35	0.61
4:32:4:TYR:HE2	4:32:11:LEU:HD11	1.64	0.61
26:14:824:A:H1'	26:14:2358:G:N7	2.14	0.61
22:1L:26:A:H61	22:1L:44:G:H22	1.48	0.61
42:D8:60:GLU:HB2	42:D8:97:LYS:HE2	1.82	0.61
29:21:170:LEU:HD21	29:21:187:ALA:HB3	1.83	0.61
32:51:92:ILE:CD1	32:51:93:GLY:H	2.14	0.61
26:1H:1021:A:C8	26:1H:1021:A:H3'	2.35	0.61
1:13:664:G:N2	1:13:741:G:H1	1.97	0.61
40:B8:2:ASN:O	40:B8:3:ARG:HG2	2.00	0.61
55:Q8:35:GLN:O	55:Q8:37:SER:N	2.26	0.61
29:21:201:THR:CG2	29:21:203:LYS:H	2.13	0.61
18:9A:34:TYR:HB3	18:9A:69:THR:HB	1.81	0.61
26:14:2567:G:H2'	26:14:2568:C:C6	2.35	0.61
26:14:739:G:OP1	58:14:3635:HOH:O	2.16	0.61
38:55:97:VAL:HG12	38:55:114:VAL:HG22	1.81	0.61
2:1E:52:GLU:HG2	2:1E:56:ARG:HH21	1.66	0.61
7:62:50:ILE:HD12	7:62:61:VAL:HG11	1.81	0.61
26:14:2537:U:H2'	26:14:2538:C:C6	2.35	0.61
2:1E:67:THR:HG21	2:1E:155:LEU:HG	1.83	0.61
26:1H:607:U:OP1	30:31:102:PRO:HA	2.01	0.61
26:14:1569:A:H5'	28:19:61:LEU:HD21	1.82	0.61
35:68:88:ASN:ND2	35:68:92:GLU:H	1.97	0.61
32:59:42:ARG:NH1	32:59:53:GLU:O	2.32	0.61
4:32:64:LEU:HB2	4:32:198:VAL:HG11	1.83	0.61
10:1I:61:GLU:OE1	14:5I:58:LYS:NZ	2.27	0.61
26:14:19:C:H2'	26:14:20:C:H6	1.64	0.61
55:Q8:4:MET:HB3	55:Q8:59:LYS:HZ2	1.66	0.61
1:13:637:G:H2'	1:13:638:G:C8	2.36	0.61
26:1H:1899:G:H1	26:1H:1902:C:H41	1.48	0.61
26:1H:141:A:OP2	26:1H:141(A):C:N4	2.28	0.61
26:14:71:A:C8	26:14:71:A:H5'	2.36	0.61
26:1H:814:C:N4	26:1H:1193:G:O6	2.17	0.61
26:1H:468:G:N7	54:P8:39:ARG:NH2	2.49	0.61
46:H8:53:ILE:HG22	46:H8:71:VAL:HG22	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:29:112:GLY:O	29:29:159:HIS:HA	2.01	0.61
1:13:243:A:H4'	1:13:244:U:H5''	1.83	0.61
26:1H:2255:G:OP2	58:1H:4121:HOH:O	2.16	0.61
29:29:93:VAL:HG22	29:29:182:LEU:HD13	1.82	0.61
38:98:21:TYR:HB3	38:98:47:PHE:CD2	2.35	0.61
5:4E:144:THR:H	5:4E:147:ASP:HB2	1.66	0.61
36:78:62:LEU:HD13	55:Q8:23:VAL:HG21	1.81	0.61
55:Q8:53:PRO:HA	55:Q8:54:GLU:C	2.21	0.61
34:15:56:ASN:H	34:15:125:GLY:HA3	1.66	0.61
15:6I:6:GLU:H	15:6I:6:GLU:CD	2.04	0.61
24:3L:9:A:H2'	24:3L:11:C:H41	1.65	0.61
26:1H:1798:U:H5'	28:11:259:THR:HG22	1.82	0.61
1:1G:690:G:H2'	1:1G:691:G:O4'	2.00	0.61
13:4I:49:THR:HG22	13:4I:51:ALA:H	1.65	0.61
26:14:2815:C:H5'	52:J5:29:THR:HG21	1.83	0.61
26:1H:1351:C:OP2	58:1H:4124:HOH:O	2.16	0.61
31:41:61:ALA:O	51:M8:7:PRO:HG2	2.01	0.61
1:1G:1376:U:OP1	7:62:98:SER:OG	2.15	0.61
3:2E:52:LEU:HA	3:2E:70:VAL:HG12	1.81	0.61
1:13:1077:G:N2	1:13:1080:A:OP2	2.24	0.61
1:1G:1105:A:H2'	1:1G:1106:G:H8	1.66	0.61
26:1H:2688:U:H5	26:1H:2720:U:OP2	1.83	0.61
26:14:1954:G:C2	26:14:2551:C:H5''	2.35	0.61
35:25:63:VAL:HG12	35:25:106:LEU:HD11	1.81	0.61
1:13:619:U:H3	4:3E:134:ASP:HB2	1.66	0.61
1:1G:547:A:OP1	58:1G:1801:HOH:O	2.16	0.61
37:88:137:TYR:CE1	46:H8:83:PRO:HG3	2.36	0.61
5:42:37:ARG:HG2	5:42:112:LEU:HA	1.83	0.61
1:13:994:A:N7	1:13:1216:G:H4'	2.16	0.61
31:49:60:LEU:HD21	31:49:92:VAL:HG11	1.83	0.61
26:14:2557:G:H2'	26:14:2558:C:C6	2.36	0.61
35:25:14:THR:HG22	35:25:52:VAL:HB	1.81	0.61
4:3E:150:GLU:HG3	4:3E:153:ARG:HE	1.64	0.61
4:3E:153:ARG:HH11	4:3E:181:MET:HB2	1.65	0.61
1:13:1497:G:C2'	1:13:1498:U:H5'	2.30	0.61
40:75:105:LEU:HD23	40:75:109:GLU:HG3	1.81	0.61
40:75:45:PHE:CE2	40:75:74:ARG:HD3	2.36	0.61
32:51:8:PRO:HG2	32:51:69:ARG:NH2	2.16	0.61
26:14:2432:A:H5''	26:14:2433:A:OP2	2.00	0.61
49:K8:47:ASN:O	49:K8:49:LYS:N	2.34	0.61
8:7E:87:SER:HB2	8:7E:93:VAL:H	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:358:U:H2'	1:1G:359:U:H6	1.66	0.61
26:1H:2836:U:H2'	26:1H:2837:G:C8	2.36	0.61
37:45:34:LEU:HD11	37:45:129:THR:HB	1.83	0.61
50:L8:5:LYS:HD2	50:L8:34:GLU:OE1	2.00	0.61
26:1H:637:A:H2'	36:78:117:GLU:OE1	2.00	0.61
1:1G:728:A:H2'	1:1G:729:A:C8	2.34	0.61
13:4I:80:ARG:NH1	19:AI:65:ASN:O	2.34	0.61
28:19:146:GLU:HB2	28:19:189:CYS:HB3	1.82	0.61
1:13:1122:U:O4	1:13:1123:A:N6	2.33	0.61
1:1G:1014:A:H4'	19:AA:14:HIS:CE1	2.36	0.61
26:1H:587:C:OP2	36:78:21:ARG:NH2	2.34	0.61
26:1H:1786:A:H2	26:1H:2606:C:H1'	1.64	0.61
26:14:994:C:O2'	26:14:996:A:OP1	2.17	0.61
26:1H:1332:G:N2	26:1H:1610:A:H8	1.97	0.61
26:1H:1055:G:H1'	26:1H:1085:A:C2	2.36	0.61
26:14:1332:G:N2	26:14:1610:A:H8	1.98	0.61
26:1H:1221:C:H2'	26:1H:1222:C:C6	2.36	0.61
19:AI:5:LEU:HD13	19:AI:10:PHE:CD1	2.35	0.61
26:1H:1252:G:O4'	41:C8:33:ARG:HD3	2.01	0.61
26:14:1688:U:O2	26:14:1700:A:H5'	2.00	0.61
12:3A:27:LEU:HD23	12:3A:33:ARG:HG2	1.82	0.61
26:1H:2683:C:OP1	40:B8:53:ARG:NH2	2.33	0.61
8:7E:10:LEU:HD22	8:7E:83:ILE:HD11	1.81	0.61
26:14:1316:U:O2'	26:14:1317:A:H5'	2.00	0.61
26:1H:1839:G:H8	26:1H:1839:G:H5''	1.64	0.61
26:1H:1143:A:OP1	34:58:25:ARG:NH2	2.15	0.61
4:32:31:CYS:O	4:32:33:MET:N	2.34	0.60
1:13:637:G:H2'	1:13:638:G:H8	1.66	0.60
13:4A:78:ILE:HD13	13:4A:92:HIS:CE1	2.36	0.60
26:14:140:A:C8	26:14:1408:C:O2'	2.51	0.60
1:13:429:U:OP2	4:3E:36:ARG:NH2	2.34	0.60
37:88:66:ILE:HG22	37:88:67:ARG:N	2.15	0.60
1:1G:362:G:H4'	12:3A:33:ARG:HH21	1.66	0.60
1:13:50:A:H1'	1:13:52:G:C8	2.36	0.60
26:14:2126:A:N1	26:14:2163:C:O2'	2.24	0.60
16:7I:3:LYS:HG3	16:7I:24:ALA:HB2	1.83	0.60
47:E5:53:MET:HG3	47:E5:59:LEU:HD23	1.83	0.60
1:1G:680:C:H42	1:1G:710:G:H1	1.48	0.60
40:B8:66:VAL:HA	40:B8:71:GLY:HA2	1.83	0.60
37:88:109:VAL:HG13	37:88:113:GLN:HB3	1.83	0.60
10:1A:79:ARG:HD3	10:1A:79:ARG:H	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1359:A:C2	26:1H:1372:U:O4	2.51	0.60
55:Q8:14:VAL:HG11	55:Q8:21:LYS:HZ1	1.65	0.60
30:31:28:ILE:HG22	30:31:112:MET:HE3	1.83	0.60
1:13:505:G:N7	58:13:1882:HOH:O	2.31	0.60
26:14:2296:U:H4'	26:14:2297:C:OP1	2.01	0.60
34:58:46:VAL:HG11	34:58:48:MET:HG3	1.82	0.60
1:1G:1315:U:H2'	1:1G:1316:G:O4'	2.01	0.60
37:45:57:HIS:CD2	37:45:117:ALA:HB2	2.35	0.60
32:51:92:ILE:HD12	32:51:93:GLY:H	1.66	0.60
26:14:1516:U:H2'	26:14:1517:G:C8	2.36	0.60
26:14:2823:A:OP1	29:29:113:PHE:HB2	2.01	0.60
9:82:42:ARG:HD3	9:82:71:SER:HB3	1.82	0.60
30:39:133:ASN:HA	30:39:162:LEU:HD23	1.81	0.60
26:14:1955:U:OP2	58:14:3818:HOH:O	2.16	0.60
4:32:119:GLN:HG2	4:32:123:HIS:HE1	1.66	0.60
35:68:60:ALA:HB1	35:68:84:ALA:HB1	1.83	0.60
1:13:1004:A:P	1:13:1025:U:H3	2.25	0.60
1:1G:1132:C:H2'	1:1G:1133:G:H8	1.65	0.60
26:1H:270(E):G:H1	26:1H:270(U):C:N4	1.99	0.60
8:72:68:ARG:NH1	8:72:70:GLN:HG2	2.16	0.60
3:2E:84:ILE:HD11	3:2E:88:ARG:HH21	1.67	0.60
26:14:1014:U:H2'	26:14:1015:G:H8	1.65	0.60
46:D5:8:TYR:CD1	46:D5:62:PRO:HG3	2.36	0.60
3:2E:119:ARG:O	3:2E:123:GLN:HG3	2.00	0.60
26:14:2068:U:H3	26:14:2430:A:H2	1.45	0.60
11:2I:73:MET:HG2	11:2I:103:LEU:HD13	1.82	0.60
1:1G:1226:C:H4'	19:AA:80:TYR:CZ	2.36	0.60
26:1H:2413:G:O6	58:1H:3735:HOH:O	2.16	0.60
37:45:35:VAL:HG12	37:45:36:ALA:H	1.65	0.60
27:1J:88:C:H3'	27:1J:89:G:N7	2.16	0.60
26:14:1288:U:C2	26:14:1327:C:O2	2.54	0.60
34:58:46:VAL:CG1	34:58:48:MET:HG3	2.31	0.60
26:14:1434:A:H61	26:14:1558:A:N6	2.00	0.60
2:1E:84:GLU:HB3	2:1E:219:VAL:HG21	1.83	0.60
26:14:1849:G:H2'	26:14:1850:G:C8	2.36	0.60
30:39:18:ARG:HG2	30:39:19:GLU:H	1.67	0.60
33:61:123:LEU:HD23	33:61:143:SER:HA	1.83	0.60
32:51:17:VAL:HG21	32:51:50:VAL:HG11	1.82	0.60
12:3A:27:LEU:HB3	12:3A:33:ARG:HG2	1.83	0.60
14:5A:26:ARG:HG2	14:5A:26:ARG:O	2.01	0.60
27:16:30:C:H2'	27:16:31:C:H5'	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1414:U:H2'	1:13:1415:G:H8	1.66	0.60
5:42:42:GLY:HA3	5:42:65:ASN:O	2.01	0.60
26:1H:1901:A:OP2	28:11:255:LYS:HE2	2.01	0.60
29:21:152:LYS:HD3	34:58:77:GLY:HA3	1.83	0.60
26:14:1337:G:H2'	26:14:1338:G:H8	1.67	0.60
48:F5:76:ARG:HG3	48:F5:94:LEU:HD13	1.81	0.60
49:K8:14:ARG:NH1	49:K8:66:GLU:OE2	2.35	0.60
4:32:31:CYS:C	4:32:33:MET:N	2.54	0.60
45:C5:74:PRO:HG2	45:C5:82:PRO:HG2	1.83	0.60
1:13:1004:A:H1'	1:13:1036:G:N2	2.17	0.60
28:19:93:ALA:HB3	28:19:105:ILE:HG22	1.84	0.60
26:14:1889:A:H3'	58:14:3914:HOH:O	2.01	0.60
4:3E:62:GLN:O	4:3E:66:ARG:HB2	2.02	0.60
1:13:1349:A:H2'	1:13:1350:A:H8	1.66	0.60
3:2E:77:ILE:O	3:2E:84:ILE:HG22	2.01	0.60
26:1H:2330:G:H2'	26:1H:2331:G:O4'	2.01	0.60
1:1G:430:A:OP2	4:32:8:VAL:HG23	2.01	0.60
26:1H:2378:A:H2'	39:A8:21:THR:HG21	1.83	0.60
22:1L:27:G:N2	22:1L:43:C:O2	2.34	0.60
1:13:1239:A:H62	1:13:1299:A:H62	1.49	0.60
26:1H:2562:U:O2'	35:68:23:ARG:NH1	2.34	0.60
26:1H:84:A:OP2	45:G8:8:LYS:NZ	2.30	0.60
39:A8:66:ALA:HA	39:A8:69:VAL:HG12	1.83	0.60
29:21:101:ARG:HG2	29:21:169:ASN:OD1	2.01	0.60
1:1G:1095:U:P	1:1G:1108:G:H1	2.24	0.60
55:M5:59:LYS:O	55:M5:60:LEU:HG	2.01	0.60
36:35:39:LYS:HD2	36:35:45:LEU:CD2	2.30	0.60
44:B5:5:TYR:CZ	49:G5:30:ARG:HG3	2.37	0.60
1:1G:539:A:H2'	1:1G:540:G:C8	2.37	0.60
26:1H:2469:A:H61	26:1H:2481:G:H1'	1.65	0.60
4:32:126:ILE:HG22	4:32:127:THR:H	1.66	0.60
46:H8:93:ASP:HA	46:H8:130:PRO:HG2	1.83	0.60
29:29:76:ARG:HG3	29:29:195:LEU:HD22	1.83	0.60
26:1H:1355:G:O6	58:1H:3717:HOH:O	2.16	0.60
1:13:1395:C:O2'	1:13:1401:G:O2'	2.17	0.60
26:14:839:U:H2'	26:14:840:C:C6	2.37	0.60
1:13:667:G:H4'	15:6I:51:HIS:CE1	2.36	0.60
49:G5:10:LEU:HD13	49:G5:59:ARG:HG2	1.84	0.60
20:BI:89:ARG:HD2	20:BI:104:LEU:HD21	1.83	0.60
1:1G:338:A:H2	1:1G:351:G:H22	1.50	0.60
1:13:368:U:OP1	33:69:91:SER:OG	2.19	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2656:U:H3	26:14:2665:A:H2	1.45	0.60
26:14:1041:C:H42	26:14:1114:G:N2	1.98	0.60
27:1J:13:A:N1	27:1J:69:G:O2'	2.29	0.60
49:K8:48:HIS:N	49:K8:50:ILE:HD11	2.16	0.60
36:35:55:ARG:HG2	36:35:56:SER:N	2.16	0.60
1:1G:520:A:N1	1:1G:536:C:H1'	2.16	0.60
40:B8:50:ILE:HD13	40:B8:64:ARG:HB3	1.82	0.60
29:21:167:VAL:HG21	29:21:187:ALA:HB3	1.84	0.60
9:82:33:PHE:HE1	9:82:37:PHE:HD2	1.49	0.60
26:14:336:C:H5''	45:C5:6:HIS:CD2	2.37	0.60
35:25:4:PRO:O	35:25:5:GLN:HB2	2.01	0.60
8:72:97:VAL:HA	8:72:100:ILE:HD11	1.83	0.60
26:1H:2774:C:H2'	26:1H:2775:A:O4'	2.01	0.60
1:1G:958:A:N3	1:1G:985:C:O2'	2.30	0.60
1:13:1256:A:OP2	3:2E:26:LYS:NZ	2.35	0.60
26:1H:1022:G:H4'	26:1H:1023:U:O5'	2.02	0.60
26:1H:2345:G:H4'	26:1H:2346:A:O5'	2.01	0.60
26:14:2776:A:OP1	26:14:2776:A:H3'	2.01	0.60
24:3L:72:C:H3'	24:3L:73:A:H5''	1.83	0.60
29:29:31:CYS:HB3	29:29:49:LEU:HB3	1.81	0.60
26:14:801:G:OP2	58:14:3734:HOH:O	2.16	0.60
18:9A:53:ARG:HH21	18:9A:60:ALA:N	1.99	0.60
2:12:15:VAL:O	2:12:209:ARG:NH2	2.34	0.60
5:4E:70:PRO:O	5:4E:72:GLN:NE2	2.34	0.60
33:69:14:ASP:OD1	33:69:15:VAL:N	2.34	0.60
5:42:60:TYR:HB3	5:42:64:ARG:HH21	1.64	0.60
9:8E:115:GLY:HA2	10:1I:58:ASP:OD1	2.01	0.60
1:13:240:C:H2'	1:13:241:C:H6	1.67	0.60
41:C8:69:CYS:HB3	41:C8:74:LEU:HD13	1.84	0.60
1:13:345:C:O2'	1:13:346:G:N2	2.33	0.60
20:BI:49:ALA:HB2	20:BI:99:LEU:HD23	1.84	0.60
17:8A:66:SER:OG	17:8A:69:LYS:HB2	2.02	0.60
26:1H:1337:G:H2'	26:1H:1338:G:H8	1.65	0.60
26:1H:2392:A:H2	26:1H:2424:C:N4	1.97	0.60
26:14:1250:G:OP2	36:35:21:ARG:HD3	2.02	0.60
26:14:892:G:N2	26:14:894:C:OP1	2.35	0.60
24:3K:76:A:H5''	48:J8:30:VAL:HG11	1.83	0.60
4:3E:26:CYS:HA	4:3E:31:CYS:HB2	1.82	0.60
1:13:1287:A:H2'	1:13:1288:A:C8	2.37	0.60
26:14:1479:G:O2'	26:14:1558:A:H5'	2.02	0.60
30:39:46:ARG:HG2	30:39:46:ARG:HH11	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1182:G:H5'	1:1G:1183:A:H5'	1.82	0.60
26:14:1653:G:C6	38:55:9:LYS:HB2	2.36	0.60
26:14:2784:C:H2'	26:14:2785:C:C6	2.37	0.60
36:35:59:LEU:O	36:35:59:LEU:HD22	2.02	0.60
1:13:1086:U:H3	1:13:1099:G:H22	1.49	0.60
26:1H:427:U:OP2	58:1H:4623:HOH:O	2.17	0.60
28:19:17:THR:O	28:19:211:ARG:NH2	2.34	0.60
51:15:23:GLU:HG3	51:15:24:THR:H	1.67	0.60
34:15:56:ASN:HB2	34:15:59:LYS:HG3	1.83	0.60
34:58:96:GLU:HG2	34:58:97:ARG:H	1.64	0.60
26:1H:2053:G:OP1	29:21:144:ARG:HD3	2.02	0.60
26:1H:419:C:H2'	26:1H:420:C:O4'	2.02	0.60
6:5E:61:LEU:HD23	6:5E:63:TYR:HE1	1.67	0.60
27:1J:2:C:H2'	27:1J:3:C:C6	2.36	0.60
33:61:143:SER:HB2	33:61:144:VAL:HG12	1.82	0.60
1:13:991:U:C4	1:13:1212:U:H1'	2.36	0.60
5:42:61:TYR:HA	5:42:64:ARG:HB2	1.82	0.60
2:12:75:LYS:HA	2:12:78:GLN:HB2	1.84	0.60
26:1H:1266:G:O2'	26:1H:2012:G:O6	2.19	0.60
1:13:671:G:H2'	1:13:672:U:H6	1.67	0.60
26:1H:492:A:H2'	26:1H:493:G:O4'	2.02	0.60
26:14:1709:U:H2'	26:14:1710:C:C6	2.37	0.60
27:16:12:C:O2'	47:18:74:ARG:HG2	2.02	0.59
26:14:459:U:H2'	26:14:460:A:C8	2.37	0.59
26:14:2317:C:H2'	26:14:2318:G:O4'	2.02	0.59
34:58:96:GLU:O	34:58:98:VAL:HG12	2.02	0.59
37:45:75:THR:HA	37:45:89:ASN:HA	1.84	0.59
1:13:1263:C:H2'	1:13:1264:C:C6	2.37	0.59
37:45:34:LEU:HD12	37:45:130:LYS:O	2.02	0.59
26:1H:2829:C:H2'	26:1H:2830:G:H5''	1.84	0.59
44:F8:89:ILE:HG22	44:F8:92:LEU:HB2	1.83	0.59
26:14:524:U:H2'	26:14:525:U:H6	1.66	0.59
1:13:116:A:H61	1:13:313:A:H1'	1.67	0.59
22:1L:12:U:O2	22:1L:24:G:N2	2.35	0.59
26:1H:943:U:P	58:1H:4508:HOH:O	2.60	0.59
1:1G:692:U:O2'	1:1G:694:A:N7	2.32	0.59
1:13:1503:A:O2'	25:4K:13:A:N6	2.35	0.59
26:14:176:G:O2'	26:14:177:G:H5'	2.02	0.59
14:5I:21:TYR:OH	14:5I:23:ARG:NH2	2.35	0.59
36:35:78:PRO:HA	36:35:110:TYR:CD2	2.36	0.59
1:13:642:A:N3	8:7E:113:SER:OG	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:55:106:GLY:O	38:55:107:ASP:HB3	2.02	0.59
1:13:1157:A:H61	1:13:1178:G:N2	1.96	0.59
26:14:831:G:H5''	26:14:832:G:OP2	2.02	0.59
26:14:459:U:H2'	26:14:460:A:H8	1.67	0.59
32:51:4:ILE:HB	32:51:6:ARG:HG3	1.84	0.59
26:1H:2685:G:OP2	40:B8:51:ARG:NH2	2.34	0.59
53:K5:26:ASN:OD1	53:K5:28:ARG:HG2	2.02	0.59
26:1H:946:G:OP2	58:1H:4576:HOH:O	2.16	0.59
9:8E:42:ARG:NH1	9:8E:71:SER:O	2.35	0.59
36:35:146:VAL:HG13	36:35:147:LEU:HD13	1.85	0.59
26:1H:1093:G:O2'	26:1H:1099:G:N2	2.35	0.59
26:14:1665:A:H4'	35:25:67:LYS:HB2	1.84	0.59
1:13:1410:G:H2'	1:13:1411:C:H6	1.67	0.59
33:69:18:VAL:HG21	33:69:44:LEU:HD21	1.84	0.59
1:1G:1396:A:H4'	1:1G:1397:C:OP2	2.01	0.59
27:16:80:U:H2'	27:16:81:G:N2	2.17	0.59
42:95:35:LEU:HB3	42:95:37:VAL:CG1	2.32	0.59
3:2E:3:ASN:OD1	3:2E:3:ASN:N	2.36	0.59
1:1G:987:G:N2	1:1G:1218:C:N3	2.38	0.59
26:1H:229:A:H4'	26:1H:230:U:H5'	1.83	0.59
24:3K:58:A:O2'	24:3K:59:U:OP1	2.21	0.59
26:14:2037:G:H2'	26:14:2038:G:C8	2.37	0.59
2:12:84:GLU:HB3	2:12:219:VAL:HG11	1.85	0.59
38:55:82:GLU:H	38:55:85:PRO:HG2	1.66	0.59
26:14:1818:U:H2'	28:19:157:ARG:HD3	1.83	0.59
3:22:18:TRP:CD1	14:5A:54:PRO:HA	2.37	0.59
26:14:125:G:H5''	54:L5:19:ARG:HD3	1.83	0.59
10:1I:38:ILE:HG23	10:1I:71:LEU:HB3	1.84	0.59
26:14:1798:U:H5'	28:19:259:THR:OG1	2.02	0.59
26:1H:1290:C:H2'	26:1H:1291:C:C6	2.38	0.59
49:K8:23:LYS:NZ	49:K8:27:GLU:OE2	2.23	0.59
3:2E:15:THR:HG21	3:2E:181:ASN:HA	1.83	0.59
26:1H:1478:G:H2'	26:1H:1479:G:C8	2.35	0.59
26:1H:2210:G:H5'	26:1H:2211:G:N7	2.17	0.59
26:14:1774:C:OP1	58:14:4076:HOH:O	2.16	0.59
23:2K:62:C:H2'	23:2K:63:C:C6	2.34	0.59
31:49:43:LEU:HD12	31:49:45:GLU:OE1	2.02	0.59
10:1A:3:LYS:NZ	10:1A:75:ILE:O	2.36	0.59
26:1H:833:U:O2	36:78:55:ARG:NH1	2.31	0.59
55:Q8:9:GLY:O	55:Q8:13:ARG:HD2	2.03	0.59
53:O8:15:GLU:HB3	53:O8:49:HIS:CD2	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:672:C:H5'	30:39:89:VAL:HG13	1.84	0.59
1:13:406:G:H2'	1:13:407:G:H8	1.66	0.59
1:13:1014:A:H4'	19:AI:14:HIS:CD2	2.38	0.59
4:32:119:GLN:O	4:32:123:HIS:ND1	2.35	0.59
1:13:313:A:H2'	1:13:314:C:C6	2.37	0.59
26:1H:1602:U:O4	58:1H:4018:HOH:O	2.16	0.59
31:49:36:LYS:HG2	31:49:38:VAL:HG23	1.85	0.59
26:14:1292:U:H2'	26:14:1293:C:C6	2.37	0.59
20:BI:50:GLU:HB2	20:BI:100:ILE:HD13	1.85	0.59
1:1G:243:A:H4'	1:1G:244:U:H5''	1.83	0.59
30:31:8:GLN:CD	30:31:8:GLN:H	2.06	0.59
13:4I:39:ILE:HD12	13:4I:56:LEU:HG	1.83	0.59
55:Q8:48:PHE:CG	55:Q8:49:VAL:HG13	2.37	0.59
26:1H:1021:A:H8	26:1H:1021:A:H3'	1.68	0.59
26:1H:2702:U:H6	26:1H:2702:U:OP1	1.85	0.59
24:3K:19:G:C6	26:1H:2112:G:H1'	2.38	0.59
58:1H:4503:HOH:O	27:16:100:G:H5''	2.03	0.59
23:2L:24:C:H2'	23:2L:25:U:H6	1.67	0.59
38:55:33:ARG:HD3	38:55:113:LEU:HD12	1.85	0.59
6:5E:36:ARG:HH21	6:5E:38:GLU:HG2	1.66	0.59
1:13:1000:A:H2'	1:13:1001:G:C8	2.37	0.59
29:21:51:PHE:O	29:21:74:PRO:HB2	2.02	0.59
4:3E:29:PRO:HA	4:3E:34:GLU:HG3	1.84	0.59
26:14:2690:C:OP2	38:55:14:SER:HB3	2.01	0.59
35:25:31:LYS:HB3	35:25:32:TYR:CD1	2.37	0.59
40:75:53:ARG:HH11	40:75:60:THR:HG23	1.67	0.59
42:D8:65:GLY:HA3	42:D8:91:TYR:CE1	2.36	0.59
39:A8:26:LEU:HD23	39:A8:87:PHE:CD1	2.37	0.59
26:1H:1027:A:C2	26:1H:2488:A:H5'	2.37	0.59
26:1H:125:G:C8	26:1H:125:G:H5'	2.38	0.59
1:13:1502:A:H2	1:13:1505:G:N1	1.92	0.59
26:1H:1577:C:OP2	58:1H:4025:HOH:O	2.17	0.59
39:A8:93:LYS:HG2	39:A8:95:HIS:HB2	1.84	0.59
46:D5:58:VAL:HG12	46:D5:60:GLU:HB3	1.83	0.59
26:1H:1252:G:N7	41:C8:36:ARG:NH1	2.51	0.59
1:1G:1004:A:H8	1:1G:1036:G:H22	1.51	0.59
26:1H:125:G:H5'	26:1H:125:G:H8	1.68	0.59
42:D8:41:GLY:O	42:D8:45:THR:HA	2.02	0.59
47:I8:64:ASP:HB2	47:I8:85:ALA:HB1	1.84	0.59
26:1H:2264:C:H5''	26:1H:2265:U:OP2	2.03	0.59
26:1H:512:G:N7	58:1H:4665:HOH:O	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:607:U:OP1	30:39:102:PRO:HA	2.02	0.59
26:1H:1178:C:H4'	26:1H:1179:C:OP1	2.01	0.59
46:H8:126:VAL:HG12	46:H8:163:LEU:HA	1.83	0.59
1:1G:179:A:H2'	1:1G:180:U:H6	1.66	0.59
31:41:67:LYS:HE2	31:41:67:LYS:H	1.67	0.59
31:41:94:LEU:HD23	31:41:94:LEU:H	1.67	0.59
40:75:24:PRO:HA	40:75:49:VAL:HG23	1.84	0.59
26:14:1952:A:C6	35:25:22:ILE:HD11	2.37	0.59
26:1H:2815:C:H2'	26:1H:2816:C:H6	1.67	0.59
52:N8:20:ARG:HG2	52:N8:23:HIS:CE1	2.38	0.59
9:82:99:LEU:HB3	9:82:101:PHE:CD1	2.38	0.59
1:1G:736:C:H2'	1:1G:737:A:C8	2.38	0.59
14:5I:48:ALA:HB2	14:5I:53:LEU:HD12	1.84	0.59
1:1G:280:C:H3'	1:1G:281:G:H5'	1.85	0.59
1:13:947:G:O3'	13:4I:109:THR:OG1	2.20	0.59
16:7I:20:VAL:HG21	16:7I:32:TYR:CD2	2.37	0.59
1:1G:108:G:H5'	1:1G:109:A:H5''	1.83	0.59
11:2I:30:VAL:HG21	11:2I:65:ALA:HA	1.83	0.59
28:11:16:MET:HE3	28:11:211:ARG:HD3	1.84	0.59
3:22:155:GLY:HA3	3:22:196:LEU:HD13	1.84	0.59
1:13:1096:C:H2'	1:13:1097:C:H6	1.66	0.59
33:6I:92:VAL:HG13	33:6I:120:ILE:HG23	1.84	0.59
13:4I:40:ASN:HB3	13:4I:43:THR:HG23	1.85	0.59
26:1H:428:A:OP1	58:1H:4617:HOH:O	2.17	0.59
1:13:973:G:H3'	1:13:974:A:H5''	1.85	0.59
26:1H:2431:U:OP2	58:1H:3905:HOH:O	2.16	0.59
27:1J:40:U:C5	51:15:1:MET:HB3	2.36	0.59
22:1L:76:A:O5'	26:14:2602:A:N6	2.36	0.59
26:1H:65:C:H2'	26:1H:66:C:H6	1.67	0.59
1:1G:1127:G:N3	1:1G:1147:C:N4	2.50	0.59
26:1H:2801:A:H2'	26:1H:2802:G:H4'	1.83	0.59
16:7A:38:TYR:CE2	16:7A:50:LYS:HD2	2.38	0.59
34:58:127:ASP:OD1	34:58:127:ASP:N	2.35	0.59
26:14:329:G:H8	26:14:329:G:OP1	1.84	0.59
47:E5:48:GLY:HA3	47:E5:80:HIS:ND1	2.17	0.59
26:1H:2306:C:H3'	26:1H:2307:G:C5'	2.33	0.59
7:62:16:LEU:HD12	9:82:42:ARG:HA	1.85	0.59
48:F5:92:LYS:O	48:F5:95:LEU:N	2.31	0.59
26:1H:1831:G:H2'	26:1H:1832:C:H6	1.68	0.59
1:1G:764:C:H5''	15:6A:50:HIS:ND1	2.17	0.59
26:1H:535:C:O3'	41:C8:53:ARG:NH1	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:68:93:PRO:HG3	35:68:114:ILE:HG12	1.84	0.59
36:35:5:ASP:O	36:35:7:ARG:NH1	2.36	0.59
26:1H:1216:G:OP2	41:C8:12:ARG:NH2	2.34	0.59
1:1G:617:G:H1	1:1G:623:C:H42	1.50	0.59
16:7I:11:SER:HB2	16:7I:14:ASN:HB3	1.82	0.59
1:13:827:U:C5	1:13:870:U:C4	2.91	0.59
26:1H:764:A:O4'	28:11:213:ARG:HG3	2.03	0.59
1:1G:1423:G:H2'	1:1G:1424:C:H6	1.66	0.59
6:52:100:ASN:OD1	18:9A:23:LYS:NZ	2.23	0.59
21:1B:2:GLY:O	21:1B:5:ASP:N	2.29	0.59
26:14:1197:G:H2'	26:14:1198:U:H6	1.68	0.59
21:1F:9:ARG:HH22	21:1F:23:PRO:HD2	1.68	0.59
26:14:584:C:OP2	41:85:10:ARG:NH2	2.36	0.59
26:1H:2138:C:H42	26:1H:2153:G:H1	1.50	0.59
1:1G:954:G:H21	1:1G:1227:A:H62	1.51	0.59
44:F8:15:GLU:HG3	44:F8:16:LYS:N	2.18	0.59
6:5E:69:GLU:O	6:5E:72:VAL:HG12	2.03	0.59
5:42:145:LYS:O	5:42:149:GLU:N	2.34	0.59
7:6E:5:ARG:NH1	7:6E:8:GLU:OE2	2.36	0.59
46:D5:163:LEU:HD23	46:D5:163:LEU:H	1.67	0.59
42:95:85:LYS:HB3	42:95:87:HIS:H	1.68	0.59
17:8I:88:TYR:CD1	17:8I:89:LEU:HD23	2.35	0.59
5:42:7:GLU:OE1	5:42:37:ARG:NH2	2.35	0.59
31:41:97:ASP:O	31:41:100:TRP:N	2.35	0.59
45:C5:87:LYS:HG2	45:C5:88:LYS:HD2	1.85	0.59
1:1G:1108:G:H5'	3:22:176:HIS:HD2	1.68	0.59
45:G8:15:VAL:HG21	45:G8:42:VAL:HG21	1.84	0.59
36:78:114:ILE:HD11	36:78:130:PHE:HD2	1.68	0.59
27:1J:24:G:H5'	27:1J:25:A:C8	2.38	0.59
26:1H:929:G:O6	58:1H:3758:HOH:O	2.17	0.59
41:C8:106:PHE:O	41:C8:109:LEU:HB2	2.03	0.59
55:Q8:30:ARG:CG	55:Q8:30:ARG:HH11	2.13	0.58
26:14:1025:G:H8	26:14:1025:G:OP1	1.85	0.58
1:1G:861:G:C5	1:1G:862:C:H5	2.20	0.58
55:M5:30:ARG:O	55:M5:32:LEU:N	2.35	0.58
26:1H:1385:G:O2'	26:1H:1396:U:H6	1.86	0.58
26:1H:2684:U:C4	26:1H:2685:G:N7	2.71	0.58
26:14:78:A:OP1	49:G5:14:ARG:NH2	2.36	0.58
29:21:33:VAL:O	29:21:69:LYS:HD2	2.03	0.58
4:3E:83:SER:HA	4:3E:89:THR:HG23	1.84	0.58
6:52:36:ARG:NH2	6:52:38:GLU:OE2	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2131:G:H5''	26:1H:2133:G:H4'	1.85	0.58
6:5E:22:GLU:O	6:5E:26:ILE:HG13	2.03	0.58
46:D5:130:PRO:HA	46:D5:133:ILE:HD11	1.84	0.58
28:19:222:ARG:HH11	28:19:222:ARG:HB2	1.66	0.58
26:1H:2734:A:H5'	26:1H:2735:G:OP2	2.02	0.58
1:13:1505:G:H5'	58:13:1801:HOH:O	2.02	0.58
49:G5:29:LYS:HG3	49:G5:57:ILE:HD13	1.85	0.58
23:2K:64:G:H2'	23:2K:65:G:H8	1.67	0.58
52:N8:30:LEU:HD23	52:N8:41:PRO:HA	1.85	0.58
26:14:1858:G:H2'	26:14:1883:G:H22	1.68	0.58
26:14:635:C:O2'	26:14:639:U:OP1	2.20	0.58
55:Q8:34:TRP:CE3	55:Q8:36:LYS:HD3	2.38	0.58
21:1B:2:GLY:O	21:1B:4:GLY:N	2.36	0.58
1:1G:522:C:OP2	12:3A:69:TYR:OH	2.21	0.58
26:14:2018:G:P	52:J5:9:LYS:HZ1	2.26	0.58
26:1H:2591:C:OP2	28:11:239:ARG:HB2	2.04	0.58
17:8A:22:LEU:HD11	17:8A:39:SER:HB2	1.83	0.58
26:14:1328:G:O6	58:14:3555:HOH:O	2.16	0.58
26:1H:1575:C:H2'	26:1H:1576:U:C6	2.38	0.58
26:1H:1069:A:H4'	26:1H:1070:A:H5''	1.84	0.58
1:13:1369:C:H2'	1:13:1370:G:C8	2.37	0.58
9:82:63:ILE:HD11	9:82:81:ILE:HD11	1.85	0.58
9:82:17:VAL:HA	9:82:63:ILE:HG12	1.85	0.58
38:98:87:TYR:HD1	38:98:90:ARG:HD2	1.67	0.58
44:F8:3:THR:HA	44:F8:6:ASP:OD2	2.03	0.58
45:G8:85:VAL:HG23	45:G8:96:ILE:HB	1.83	0.58
51:M8:38:LYS:NZ	51:M8:44:THR:OG1	2.36	0.58
1:1G:501:C:H2'	1:1G:502:G:C8	2.38	0.58
40:B8:20:PRO:HD2	40:B8:86:ILE:HG23	1.85	0.58
26:1H:155:C:N4	26:1H:171:G:H1	2.00	0.58
26:14:2645:G:N2	26:14:2767:C:OP2	2.35	0.58
26:1H:1831:G:H2'	26:1H:1832:C:C6	2.38	0.58
26:1H:2845:G:OP2	58:1H:4683:HOH:O	2.17	0.58
26:1H:1375:C:H2'	26:1H:1376:C:H6	1.67	0.58
1:13:1226:C:H4'	19:AI:80:TYR:OH	2.02	0.58
35:25:2:ILE:HD12	35:25:6:THR:HG21	1.85	0.58
1:1G:777:A:H2	11:2A:119:CYS:HB3	1.69	0.58
1:13:1074:G:O2'	1:13:1101:A:N1	2.31	0.58
55:Q8:27:THR:O	55:Q8:29:LYS:HA	2.02	0.58
39:A8:35:ILE:HG22	39:A8:97:ARG:HH21	1.66	0.58
8:72:32:LYS:HA	8:72:35:ILE:HD12	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2823:A:OP1	29:29:159:HIS:NE2	2.37	0.58
51:15:49:PHE:HD2	51:15:50:VAL:HG22	1.67	0.58
1:13:536:C:H2'	1:13:537:G:C8	2.38	0.58
26:1H:698:C:O2'	26:1H:734:A:N6	2.36	0.58
39:A8:61:ASN:HB3	39:A8:64:GLU:HG3	1.84	0.58
10:1I:4:ILE:HG12	10:1I:100:THR:HA	1.83	0.58
2:12:166:ASP:CG	2:12:169:LYS:HB2	2.24	0.58
26:1H:270(M):U:OP1	26:1H:270(N):G:N2	2.36	0.58
1:13:727:G:N2	1:13:730:G:OP2	2.21	0.58
10:1I:6:ILE:HG22	10:1I:98:ILE:HG13	1.85	0.58
11:2I:121:PRO:HG2	11:2I:126:ARG:HG3	1.85	0.58
46:H8:128:VAL:HB	46:H8:161:VAL:HG21	1.84	0.58
26:1H:1329:U:H5'	26:1H:1330:C:H5	1.68	0.58
20:BI:71:THR:HG22	20:BI:72:LEU:N	2.18	0.58
1:1G:1286:A:C8	1:1G:1287:A:H4'	2.38	0.58
26:14:2749:A:N1	26:14:2750:A:N6	2.51	0.58
26:14:2157:G:O2'	26:14:2158:A:H8	1.85	0.58
36:35:3:LEU:H	36:35:3:LEU:HD12	1.68	0.58
12:3I:53:ARG:HH12	12:3I:92:ASP:CB	2.17	0.58
26:1H:71:A:C2	44:F8:31:HIS:HE1	2.21	0.58
26:1H:2331:G:O2'	26:1H:2336:A:N1	2.30	0.58
26:14:602:G:O2'	26:14:655:A:N6	2.36	0.58
4:32:103:ASN:OD1	4:32:114:ARG:NH2	2.32	0.58
2:1E:11:LEU:HB3	2:1E:213:LEU:HD11	1.85	0.58
1:13:1497:G:H2'	1:13:1498:U:H5'	1.86	0.58
28:19:67:PHE:HB3	28:19:153:ALA:H	1.68	0.58
39:A8:26:LEU:HD11	39:A8:73:LEU:HD13	1.86	0.58
26:14:287:C:H2'	26:14:288:C:C6	2.38	0.58
26:14:2232:U:P	48:F5:40:ARG:HH12	2.26	0.58
6:5E:24:GLU:HG3	6:5E:28:ARG:NH2	2.18	0.58
28:11:132:PRO:HD3	28:11:190:TYR:CZ	2.38	0.58
39:65:83:LYS:HG3	39:65:84:GLN:HG2	1.85	0.58
26:1H:1785:A:H4'	26:1H:1982:C:O2'	2.04	0.58
35:68:4:PRO:O	35:68:5:GLN:HB2	2.01	0.58
1:13:1325:C:H2'	1:13:1326:C:H6	1.68	0.58
1:13:1453:G:C8	20:BI:39:LYS:HE2	2.38	0.58
4:3E:62:GLN:HB3	4:3E:66:ARG:NH1	2.18	0.58
1:1G:1154:G:H2'	1:1G:1155:G:C8	2.38	0.58
24:3L:5:G:H2'	24:3L:6:G:H8	1.68	0.58
1:13:313:A:H2'	1:13:314:C:H6	1.69	0.58
1:13:1503:A:HO2'	25:4K:13:A:N6	2.00	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1467:C:H2'	26:14:1468:C:H6	1.69	0.58
9:82:18:PHE:HD2	9:82:62:TYR:HD2	1.49	0.58
26:1H:639:U:H2'	26:1H:640:C:C6	2.39	0.58
5:42:78:HIS:HA	8:72:105:ARG:HG3	1.86	0.58
5:4E:35:GLY:HA3	5:4E:112:LEU:HB3	1.86	0.58
26:1H:1567:A:H5''	28:11:58:HIS:ND1	2.19	0.58
26:14:1486:A:H2'	26:14:1487:G:H8	1.68	0.58
26:14:2812:G:N2	26:14:2889:C:O2	2.36	0.58
36:35:98:GLU:HA	36:35:101:VAL:HG12	1.86	0.58
26:14:1510:A:H2'	26:14:1511:A:C8	2.39	0.58
36:78:122:PRO:HA	36:78:142:GLY:HA3	1.84	0.58
33:61:75:LEU:HB3	33:61:105:HIS:CD2	2.39	0.58
6:52:96:PRO:HB3	18:9A:30:ASP:CG	2.23	0.58
28:11:17:THR:HG22	28:11:205:VAL:N	2.13	0.58
1:13:963:G:H21	10:1I:55:LYS:CE	2.16	0.58
28:11:145:VAL:HG12	28:11:146:GLU:O	2.04	0.58
26:1H:1388:G:H2'	26:1H:1389:G:H8	1.67	0.58
1:1G:980:C:H3'	1:1G:981:U:C6	2.39	0.58
26:1H:1800:C:OP2	28:11:183:ARG:NH2	2.32	0.58
2:12:19:HIS:CG	2:12:20:GLU:H	2.22	0.58
48:F5:85:LEU:HA	48:F5:87:PRO:HD2	1.85	0.58
27:1J:88:C:H3'	27:1J:89:G:C8	2.38	0.58
28:19:206:LEU:HD22	28:19:211:ARG:HG2	1.85	0.58
44:B5:8:ILE:O	49:G5:36:ARG:NH2	2.37	0.58
26:1H:2105:C:H2'	26:1H:2106:G:C8	2.38	0.58
26:1H:1270:C:H5''	26:1H:1271:G:H5'	1.83	0.58
47:E5:66:VAL:HG12	47:E5:67:VAL:H	1.69	0.58
11:2I:18:ARG:HB3	11:2I:33:THR:OG1	2.03	0.58
1:1G:481:G:O2'	1:1G:483:C:N4	2.36	0.58
16:7A:22:THR:HA	16:7A:33:ILE:HG12	1.84	0.58
26:14:1759:A:H4'	26:14:2715:C:O4'	2.04	0.58
55:M5:29:LYS:HB3	55:M5:44:LYS:HB2	1.86	0.58
5:42:101:ILE:HD11	5:42:119:LEU:HD23	1.85	0.58
1:1G:964:A:N3	1:1G:969:A:O2'	2.32	0.58
26:1H:588:U:H2'	26:1H:589:C:C6	2.39	0.58
24:3L:18:G:H2'	24:3L:57:G:N1	2.18	0.58
1:13:674:G:H2'	1:13:675:A:C8	2.38	0.58
44:B5:11:PRO:HB3	44:B5:92:LEU:HD21	1.86	0.58
45:G8:55:TYR:CE1	45:G8:61:ILE:HD11	2.39	0.58
39:A8:89:ARG:HG2	39:A8:89:ARG:O	2.03	0.58
31:49:64:THR:HG23	31:49:66:GLN:H	1.67	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:948:C:O2'	1:13:949:A:H5'	2.03	0.58
1:1G:619:U:N3	4:32:134:ASP:OD1	2.35	0.58
26:1H:527:C:H4'	26:1H:528:A:H5'	1.85	0.58
26:14:1520:U:H2'	26:14:1521:G:O4'	2.03	0.58
3:2E:134:ILE:O	3:2E:138:VAL:HG12	2.03	0.58
26:1H:634:C:H2'	26:1H:635:C:C6	2.37	0.58
13:4I:74:VAL:O	13:4I:78:ILE:HD12	2.03	0.58
6:52:11:ASN:ND2	6:52:84:ASN:OD1	2.36	0.58
26:1H:1110:G:O2'	26:1H:1111:A:H8	1.87	0.58
26:14:605:C:O2	26:14:657:U:O2'	2.21	0.58
26:14:718:A:H3'	26:14:719:C:H6	1.68	0.58
12:3A:46:LYS:HE2	12:3A:47:LYS:HB2	1.86	0.58
46:D5:157:LEU:HG	46:D5:161:VAL:HG11	1.86	0.58
36:35:84:ASN:ND2	36:35:117:GLU:HB3	2.19	0.58
1:13:677:U:H3	1:13:713:G:H22	1.52	0.58
26:14:270(J):G:C2	26:14:270(K):C:H1'	2.39	0.58
26:14:1093:G:O2'	26:14:1099:G:N2	2.37	0.58
26:14:453:C:OP1	58:14:4015:HOH:O	2.17	0.58
41:85:47:TYR:HA	41:85:50:ARG:NH2	2.19	0.58
26:1H:2019:A:H2	26:1H:2035:G:H22	1.52	0.58
1:13:413:G:N2	1:13:428:G:H1'	2.19	0.58
38:98:72:ASP:O	38:98:76:VAL:HG23	2.02	0.58
26:14:2340:G:H2'	26:14:2341:G:C8	2.38	0.58
24:3K:50:U:H3	24:3K:64:A:N6	2.00	0.58
26:14:1570:A:H2'	26:14:1571:A:C8	2.39	0.58
26:1H:270(N):G:OP1	33:61:57:ARG:NH2	2.37	0.58
46:H8:72:ARG:NH2	46:H8:97:GLU:O	2.27	0.58
5:42:10:MET:HA	5:42:32:VAL:HG22	1.86	0.58
30:31:64:ILE:HG23	30:31:65:TRP:CD1	2.39	0.58
26:1H:286:C:H2'	26:1H:287:C:H6	1.68	0.58
26:14:854:G:H2'	26:14:855:G:H8	1.69	0.58
32:59:8:PRO:HG2	32:59:69:ARG:HH21	1.67	0.58
50:H5:10:LYS:NZ	50:H5:15:TYR:OH	2.29	0.58
51:I5:18:CYS:SG	51:I5:19:GLY:HA2	2.43	0.58
38:98:103:ARG:HD2	38:98:108:GLY:O	2.03	0.58
2:1E:33:TYR:HB2	2:1E:43:ASP:HB2	1.85	0.58
1:1G:1324:A:H2'	1:1G:1325:C:C6	2.39	0.58
26:14:2298:A:H1'	26:14:2321:G:N2	2.18	0.58
9:82:29:ASN:N	9:82:63:ILE:O	2.36	0.58
24:3K:64:A:C2	24:3K:65:G:H1'	2.39	0.58
6:52:7:ASN:N	6:52:7:ASN:OD1	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:2I:85:ARG:HD3	11:2I:113:PRO:HD3	1.86	0.58
22:1K:76:A:C8	26:1H:2507:C:H1'	2.39	0.58
1:1G:632:A:H1'	1:1G:633:G:OP2	2.04	0.58
5:42:140:ARG:O	5:42:143:ARG:NH2	2.37	0.58
26:1H:1342:A:OP2	58:1H:4018:HOH:O	2.17	0.58
26:1H:1653:G:H3'	38:98:2:ARG:HG2	1.85	0.58
38:98:13:HIS:CE1	38:98:15:SER:HB3	2.38	0.58
47:I8:55:ARG:HH11	47:I8:55:ARG:HG3	1.69	0.58
22:1L:9:A:OP2	22:1L:13:C:N4	2.36	0.58
7:62:67:GLU:OE2	7:62:70:LYS:NZ	2.35	0.57
26:14:751:A:OP1	58:14:3410:HOH:O	2.17	0.57
1:13:1004:A:H1'	1:13:1036:G:H22	1.69	0.57
26:1H:780:G:H21	26:1H:783:A:N6	1.97	0.57
1:13:604:G:H2'	1:13:605:U:O4'	2.04	0.57
1:1G:960:U:H4'	1:1G:961:U:H5''	1.85	0.57
1:1G:1298:C:O2'	1:1G:1299:A:OP2	2.19	0.57
20:BI:61:SER:O	20:BI:65:LYS:HB2	2.04	0.57
47:E5:18:ALA:HB3	47:E5:20:ARG:HH21	1.67	0.57
26:14:1927:A:H2'	26:14:1928:A:C8	2.38	0.57
29:29:8:LYS:HG2	29:29:192:ASN:HA	1.86	0.57
44:B5:36:LYS:HG3	44:B5:56:THR:HG23	1.86	0.57
26:1H:582:G:H2'	26:1H:583:G:C8	2.39	0.57
26:1H:814:C:N3	26:1H:1193:G:N1	2.46	0.57
19:AA:80:TYR:CZ	19:AA:82:GLY:HA2	2.38	0.57
3:22:44:GLU:HA	3:22:52:LEU:HD11	1.86	0.57
9:82:51:ARG:HG2	9:82:56:LEU:HB2	1.84	0.57
1:13:451:A:N6	1:13:480:U:H2'	2.18	0.57
30:39:67:GLN:O	30:39:67:GLN:HG3	2.04	0.57
40:B8:31:SER:OG	40:B8:85:LYS:NZ	2.36	0.57
26:1H:2300:G:H1	26:1H:2316:C:H42	1.52	0.57
33:69:6:LEU:HB2	33:69:36:ALA:HA	1.86	0.57
1:1G:600:C:H2'	1:1G:601:C:C6	2.39	0.57
27:16:15:A:H5'	27:16:16:G:C8	2.38	0.57
29:21:8:LYS:HA	29:21:26:ILE:HG22	1.86	0.57
26:1H:507:A:H5''	26:1H:508:G:H5'	1.86	0.57
3:22:57:ILE:HG12	3:22:66:VAL:HG22	1.86	0.57
4:32:176:LEU:HG	4:32:178:VAL:HG13	1.85	0.57
26:14:996:A:H4'	41:85:92:ARG:CZ	2.34	0.57
41:85:50:ARG:HH22	42:95:72:VAL:HG23	1.68	0.57
1:13:276:G:O3'	17:8I:68:ARG:NH1	2.37	0.57
26:1H:2062:A:N3	26:1H:2062:A:H2'	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:732:C:H3'	58:1H:4070:HOH:O	2.04	0.57
26:14:2233:U:H2'	26:14:2234:G:C8	2.39	0.57
1:13:412:A:OP2	4:3E:35:ARG:NH2	2.37	0.57
26:14:877:U:O4	26:14:899:A:N6	2.37	0.57
9:8E:112:LYS:CA	9:8E:119:ALA:HB2	2.33	0.57
1:1G:1256:A:N6	1:1G:1277:C:H3'	2.18	0.57
26:1H:2283:C:H2'	26:1H:2284:C:O4'	2.04	0.57
12:3I:47:LYS:HA	12:3I:49:ASN:H	1.68	0.57
2:12:145:LEU:O	2:12:149:LEU:HB2	2.04	0.57
6:5E:39:LYS:HB2	6:5E:64:GLN:HB2	1.86	0.57
26:1H:1139:G:O2'	26:1H:1143:A:N1	2.32	0.57
27:1J:87:G:H3'	27:1J:88:C:C5'	2.32	0.57
36:78:114:ILE:HD11	36:78:130:PHE:CD2	2.38	0.57
38:55:67:LEU:HD12	38:55:76:VAL:HG11	1.86	0.57
26:1H:1807:G:N7	58:1H:4385:HOH:O	2.32	0.57
26:1H:1283:G:N2	26:1H:1285:G:H3'	2.20	0.57
15:6I:78:TYR:CZ	15:6I:82:ILE:HD11	2.39	0.57
47:E5:54:GLY:O	47:E5:57:PHE:N	2.36	0.57
26:1H:2023:G:H5'	26:1H:2617:C:H4'	1.86	0.57
26:1H:701:G:N2	26:1H:731:C:O2	2.33	0.57
42:95:85:LYS:HD2	42:95:86:GLY:N	2.14	0.57
1:1G:999:U:H2'	1:1G:1000:A:C8	2.34	0.57
1:1G:457:C:H2'	1:1G:458:C:C6	2.39	0.57
26:14:2392:A:OP2	55:M5:32:LEU:HD23	2.04	0.57
1:1G:1238:A:H62	1:1G:1301:U:H3	1.50	0.57
24:3L:58:A:O2'	24:3L:59:U:O5'	2.21	0.57
33:61:69:LYS:HG3	33:61:136:VAL:HB	1.86	0.57
26:14:1087:G:H2'	26:14:1089:G:H1'	1.86	0.57
27:16:29:A:H2'	27:16:30:C:O4'	2.04	0.57
3:2E:167:TRP:CD1	3:2E:168:ALA:N	2.72	0.57
1:13:1396:A:H4'	1:13:1397:C:H5''	1.86	0.57
26:14:1430:C:H2'	26:14:1431:U:C6	2.40	0.57
9:8E:9:ARG:HD2	9:8E:14:VAL:HG13	1.84	0.57
46:H8:52:SER:O	46:H8:52:SER:OG	2.13	0.57
26:1H:1403:C:C5'	26:1H:1471:A:H1'	2.34	0.57
6:5E:101:ALA:HB2	18:9I:28:GLU:HG2	1.86	0.57
9:8E:21:PRO:HA	9:8E:59:PHE:HA	1.84	0.57
42:95:69:LYS:HB3	42:95:88:ARG:HG3	1.85	0.57
26:1H:607:U:C2	26:1H:621:A:N1	2.71	0.57
26:1H:574:C:OP1	58:1H:4541:HOH:O	2.17	0.57
26:1H:1657:C:H2'	26:1H:1658:C:C6	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1119:C:OP2	9:82:9:ARG:NH2	2.37	0.57
26:1H:1919:A:H5'	26:1H:1920:C:OP2	2.04	0.57
17:8I:83:ASP:OD1	17:8I:83:ASP:N	2.37	0.57
26:14:1600:C:OP1	44:B5:58:HIS:NE2	2.28	0.57
50:L8:8:LEU:HD13	50:L8:31:LEU:HD23	1.85	0.57
1:1G:1190:G:H5'	3:22:176:HIS:CE1	2.40	0.57
44:F8:11:PRO:HB3	44:F8:92:LEU:HD21	1.86	0.57
38:55:72:ASP:O	38:55:76:VAL:HG12	2.04	0.57
9:8E:34:ASN:O	9:8E:38:GLN:HB2	2.04	0.57
44:F8:26:TYR:O	44:F8:81:VAL:HG12	2.05	0.57
19:AA:79:THR:OG1	19:AA:79:THR:O	2.23	0.57
22:1L:51:U:H2'	22:1L:52:G:C8	2.39	0.57
1:1G:1499:A:H1'	1:1G:1520:G:H5'	1.85	0.57
26:1H:1771:C:H1'	26:1H:1786:A:C8	2.39	0.57
14:5A:21:TYR:HE1	14:5A:23:ARG:NE	1.98	0.57
26:1H:1424:G:H2'	26:1H:1425:G:O4'	2.05	0.57
46:D5:39:VAL:HG21	46:D5:44:PHE:HD2	1.69	0.57
24:3L:76:A:H8	26:14:2394:C:N4	2.02	0.57
26:14:2506:U:O5'	26:14:2506:U:H6	1.87	0.57
26:1H:2684:U:H1'	35:68:70:LYS:HD2	1.87	0.57
27:16:42:C:O2'	31:41:67:LYS:HE3	2.04	0.57
45:C5:42:VAL:HG13	45:C5:65:ALA:HB3	1.86	0.57
37:45:22:LYS:N	37:45:23:GLY:HA3	2.19	0.57
37:88:66:ILE:O	37:88:104:PHE:N	2.38	0.57
1:1G:60:A:N6	1:1G:110:C:N3	2.50	0.57
4:3E:79:PHE:O	4:3E:83:SER:HB2	2.04	0.57
26:14:288:C:H2'	26:14:289:A:C8	2.38	0.57
26:1H:1062:G:N1	26:1H:1076:C:O2	2.36	0.57
26:14:918:A:O2'	27:1J:96:G:N2	2.37	0.57
26:1H:443:A:H1'	26:1H:1201:C:O4'	2.04	0.57
26:1H:172:C:H2'	26:1H:173:G:H8	1.70	0.57
33:61:21:VAL:HG21	33:61:25:TYR:HD2	1.68	0.57
26:1H:2232:U:P	48:J8:40:ARG:HH12	2.27	0.57
26:1H:1965:C:H2'	26:1H:1966:A:C8	2.39	0.57
36:78:49:ARG:HE	55:Q8:57:ARG:HG2	1.69	0.57
26:14:1041:C:N4	26:14:1114:G:H22	1.99	0.57
37:88:14:ARG:HG2	37:88:41:TRP:CH2	2.34	0.57
27:1J:15:A:H1'	27:1J:109:G:C4	2.39	0.57
26:1H:2062:A:OP1	58:1H:4649:HOH:O	2.17	0.57
44:F8:49:VAL:HG12	44:F8:50:LYS:N	2.20	0.57
38:98:55:ALA:HA	38:98:80:PHE:CE1	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:950:G:H2'	26:1H:951:C:C6	2.40	0.57
1:1G:345:C:O2'	1:1G:346:G:O5'	2.21	0.57
37:45:25:ASP:OD1	37:45:25:ASP:N	2.38	0.57
23:2L:63:C:H2'	23:2L:64:G:C8	2.39	0.57
26:14:2860:A:H5''	26:14:2861:G:OP2	2.04	0.57
30:39:3:GLU:HG3	30:39:20:LEU:O	2.04	0.57
16:7I:34:GLU:HG2	16:7I:35:LYS:N	2.19	0.57
1:1G:1111:A:H2'	1:1G:1112:C:C6	2.39	0.57
2:1E:24:TRP:CZ3	2:1E:26:PRO:HA	2.39	0.57
26:1H:303:U:H2'	26:1H:304:G:H8	1.69	0.57
32:59:81:GLU:HG2	32:59:83:TYR:H	1.69	0.57
6:52:69:GLU:CD	6:52:69:GLU:H	2.07	0.57
50:L8:28:LEU:HD23	50:L8:33:GLN:HG2	1.86	0.57
23:2L:47:7MG:H3'	23:2L:48:U:C6	2.39	0.57
17:8I:66:SER:O	17:8I:70:ARG:NH1	2.38	0.57
26:14:1677:A:H2'	26:14:1678:G:C8	2.40	0.57
1:13:110:C:H2'	1:13:111:G:O4'	2.05	0.57
30:31:127:GLU:OE1	30:31:128:ALA:N	2.37	0.57
26:1H:298:G:OP2	45:G8:84:ARG:NH1	2.36	0.57
29:29:36:ARG:NH1	29:29:85:ASN:OD1	2.33	0.57
3:2E:40:ARG:HG3	3:2E:40:ARG:HH11	1.69	0.57
27:1J:62:C:H2'	27:1J:63:G:C8	2.38	0.57
5:42:93:PRO:HG2	8:72:105:ARG:NE	2.19	0.57
47:E5:66:VAL:HG12	47:E5:67:VAL:N	2.20	0.57
26:1H:1045:A:H1'	26:1H:1047:G:N3	2.19	0.57
28:11:71:ASP:OD2	28:11:103:ARG:NH2	2.37	0.57
9:82:4:TYR:HB2	9:82:19:LEU:HB2	1.85	0.57
5:4E:11:ILE:HD11	5:4E:31:LEU:HD13	1.85	0.57
26:14:2311:A:N6	31:49:44:GLY:HA3	2.19	0.57
48:F5:29:GLY:O	48:F5:30:VAL:HG22	2.04	0.57
26:1H:7:G:H2'	26:1H:8:A:O4'	2.04	0.57
26:1H:2404:C:OP2	58:1H:4219:HOH:O	2.17	0.57
26:14:2745:C:O2	32:59:139:GLN:NE2	2.29	0.57
13:4A:86:CYS:SG	13:4A:88:ARG:HB3	2.45	0.57
26:1H:1331:A:O2'	26:1H:1332:G:H8	1.88	0.57
29:21:120:TRP:CE3	29:21:155:LYS:HD3	2.40	0.57
4:3E:18:LYS:NZ	4:3E:26:CYS:O	2.34	0.57
40:B8:107:ASP:O	40:B8:110:ILE:HG23	2.05	0.57
45:G8:94:LYS:HG3	45:G8:95:LYS:H	1.69	0.57
26:1H:141:A:H8	26:1H:1408:C:H1'	1.70	0.57
1:1G:359:U:H2'	1:1G:360:A:H8	1.66	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:2L:63:C:H2'	23:2L:64:G:H8	1.68	0.57
3:2E:19:GLU:O	3:2E:40:ARG:NH2	2.38	0.57
31:49:39:ILE:HG23	31:49:157:ILE:HG23	1.85	0.57
26:14:2693:A:H2'	26:14:2694:G:H8	1.70	0.57
26:1H:1448:G:H1'	26:1H:1528:A:H62	1.70	0.57
3:22:129:ALA:O	3:22:133:ALA:N	2.38	0.57
4:32:59:ARG:HA	4:32:62:GLN:HB2	1.87	0.57
26:1H:1654:A:OP2	38:98:1:MET:N	2.27	0.57
40:B8:15:VAL:HG23	40:B8:79:HIS:CE1	2.39	0.57
26:1H:2064:C:H2'	26:1H:2065:C:C6	2.39	0.57
47:E5:27:GLU:HB2	47:E5:69:PHE:HD1	1.69	0.57
13:4I:20:THR:HG23	13:4I:26:GLY:HA3	1.87	0.57
26:1H:1520:U:H2'	26:1H:1521:G:O4'	2.05	0.57
26:1H:1250:G:OP2	36:78:18:ARG:NH1	2.38	0.57
42:95:69:LYS:HD3	42:95:86:GLY:HA3	1.86	0.57
26:1H:2496:C:OP1	37:88:82:ARG:HD3	2.05	0.57
4:3E:107:ARG:NH2	4:3E:194:LEU:HD22	2.15	0.57
26:14:1036:G:OP1	32:59:59:ARG:N	2.35	0.57
1:13:271:C:H2'	1:13:272:C:H6	1.70	0.57
26:14:794:G:H2'	26:14:795:C:C6	2.39	0.57
1:1G:1081:G:OP1	5:42:18:ARG:HB2	2.05	0.57
26:1H:2489:G:O2'	26:1H:2518:A:N6	2.38	0.57
26:14:1485:G:H2'	26:14:1486:A:C8	2.40	0.57
26:1H:172:C:H2'	26:1H:173:G:C8	2.40	0.57
27:16:75:G:H21	46:H8:85:HIS:CE1	2.23	0.57
32:59:23:ARG:HA	32:59:36:PRO:HA	1.87	0.57
26:1H:1049:C:H2'	26:1H:1050:A:H5'	1.87	0.57
3:22:8:ILE:HD11	3:22:184:TYR:HB3	1.85	0.57
3:22:8:ILE:HG23	3:22:16:ARG:HG2	1.87	0.57
30:31:51:THR:HG23	30:31:92:PRO:HG2	1.87	0.57
4:3E:43:HIS:HA	4:3E:46:LYS:HZ3	1.70	0.57
32:59:137:ASP:HB3	32:59:140:LYS:HB3	1.87	0.57
28:19:40:THR:OG1	28:19:41:GLY:N	2.37	0.57
17:8I:48:GLU:O	17:8I:50:LYS:HG2	2.03	0.57
12:3A:24:VAL:O	12:3A:26:ALA:N	2.37	0.57
26:14:443:A:H1'	26:14:1201:C:O4'	2.04	0.57
26:14:1819:A:H4'	26:14:1820:U:O5'	2.04	0.57
26:14:31:C:OP1	58:14:3729:HOH:O	2.18	0.57
10:1I:48:THR:OG1	10:1I:62:HIS:ND1	2.38	0.57
9:82:111:ARG:HD2	14:5A:61:TRP:OXT	2.05	0.57
29:29:25:VAL:HG12	29:29:26:ILE:H	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1037:C:H2'	1:1G:1038:C:C6	2.40	0.57
24:3K:60:U:O2'	24:3K:61:C:OP1	2.18	0.57
55:Q8:27:THR:HG22	55:Q8:29:LYS:HB3	1.86	0.57
1:1G:1206:G:O2'	3:22:192:THR:O	2.12	0.57
36:78:80:TYR:CZ	36:78:111:ARG:HD3	2.40	0.57
27:16:15:A:H1'	27:16:109:G:C8	2.40	0.57
45:G8:35:TYR:CD2	45:G8:69:ALA:HB3	2.39	0.57
3:22:20:SER:HB2	3:22:40:ARG:HH12	1.70	0.57
32:59:20:ALA:O	32:59:22:GLY:N	2.36	0.57
10:11:16:LEU:HD23	10:11:94:VAL:HG13	1.86	0.57
26:14:688:U:C2'	26:14:689:A:H5'	2.35	0.57
26:14:184:C:H2'	26:14:185:U:H6	1.70	0.57
26:14:144:C:H2'	26:14:145:G:H8	1.69	0.57
26:1H:1607:C:H4'	26:1H:1608:A:O5'	2.05	0.57
55:Q8:53:PRO:HB3	55:Q8:56:GLU:HG3	1.86	0.56
22:1L:55:U:H2'	22:1L:57:G:OP2	2.05	0.56
1:13:255:G:P	17:8I:69:LYS:HZ3	2.28	0.56
26:14:95:G:O2'	49:G5:48:HIS:HB3	2.05	0.56
49:G5:53:LEU:O	49:G5:57:ILE:HG13	2.05	0.56
1:1G:1163:C:N3	1:1G:1174:G:N2	2.52	0.56
1:1G:1191:A:OP1	3:22:4:LYS:NZ	2.34	0.56
41:C8:92:ARG:HD3	41:C8:94:ASN:HB3	1.86	0.56
8:7E:86:ILE:HG22	8:7E:87:SER:H	1.70	0.56
1:1G:503:C:OP2	12:3A:116:SER:HB3	2.05	0.56
26:14:424:G:OP1	48:F5:96:LYS:NZ	2.38	0.56
30:31:139:PHE:HB2	30:31:166:ALA:HB1	1.86	0.56
50:L8:7:LYS:HB2	50:L8:34:GLU:HG2	1.87	0.56
26:1H:286:C:H2'	26:1H:287:C:C6	2.40	0.56
28:11:79:VAL:HG12	28:11:113:VAL:HA	1.87	0.56
26:14:2849:U:O4	40:75:23:ARG:NH2	2.38	0.56
38:55:8:ARG:HH11	38:55:39:PRO:HB3	1.68	0.56
26:14:1796:U:H2'	26:14:1797:C:C6	2.40	0.56
27:1J:101:A:OP2	58:1J:311:HOH:O	2.17	0.56
27:1J:104:A:H2'	27:1J:105:G:O4'	2.05	0.56
26:1H:2272:U:H5''	26:1H:2273:A:OP1	2.05	0.56
46:H8:76:LEU:CD2	46:H8:76:LEU:H	2.15	0.56
46:D5:44:PHE:CE2	46:D5:86:VAL:HG11	2.37	0.56
46:D5:5:LEU:HD11	46:D5:44:PHE:HA	1.87	0.56
7:62:78:ARG:CZ	7:62:80:VAL:HB	2.36	0.56
26:14:660:G:H21	36:35:12:ALA:CB	2.17	0.56
40:75:4:GLY:HA2	40:75:8:LYS:HB2	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:61:47:LEU:HA	33:61:50:ARG:HB2	1.87	0.56
39:A8:51:ALA:HB3	39:A8:73:LEU:HG	1.86	0.56
1:1G:954:G:H2'	1:1G:955:U:C6	2.40	0.56
26:1H:1568:G:H4'	28:11:59:LYS:HB3	1.87	0.56
32:59:19:VAL:HG12	32:59:20:ALA:H	1.69	0.56
11:2I:95:ILE:HD13	11:2I:108:ILE:HG21	1.86	0.56
8:72:120:THR:HG23	8:72:123:GLU:H	1.70	0.56
26:14:2410:G:C2	26:14:2411:A:H1'	2.40	0.56
9:82:43:ALA:HA	9:82:74:ILE:HD13	1.86	0.56
1:13:586:C:O2'	1:13:878:G:H4'	2.05	0.56
26:1H:1105:U:H2'	26:1H:1106:G:C8	2.40	0.56
1:1G:142:G:H2'	1:1G:143:A:H8	1.69	0.56
34:15:15:LEU:HB2	34:15:134:ARG:HG2	1.85	0.56
34:58:22:THR:HG22	34:58:23:LEU:N	2.20	0.56
55:Q8:49:VAL:HG12	55:Q8:52:LYS:CG	2.36	0.56
26:14:1021:A:H8	26:14:1021:A:H3'	1.70	0.56
1:13:1026:G:H1	1:13:1035:A:H61	1.54	0.56
1:13:1026:G:H1	1:13:1035:A:N6	2.03	0.56
8:72:12:ARG:HE	8:72:26:VAL:HA	1.70	0.56
32:51:46:GLU:OE1	32:51:51:ARG:NH2	2.38	0.56
24:3L:18:G:H1'	24:3L:58:A:H2	1.70	0.56
43:E8:14:PRO:HB3	43:E8:18:ARG:NH1	2.20	0.56
26:14:2286:A:H4'	26:14:2287:A:O4'	2.04	0.56
1:13:1453:G:H2'	20:BI:39:LYS:HE2	1.86	0.56
26:14:389:G:N1	36:35:71:VAL:HG12	2.20	0.56
26:1H:1042:G:H1	26:1H:1113:U:H3	1.53	0.56
44:F8:1:MET:C	44:F8:3:THR:H	2.08	0.56
1:1G:1329:A:H5'	13:4A:29:ARG:HE	1.70	0.56
33:69:81:VAL:HG23	33:69:143:SER:HB2	1.87	0.56
26:14:582:G:H2'	26:14:583:G:C8	2.41	0.56
26:1H:443:A:N7	30:31:45:ARG:HG2	2.21	0.56
28:11:71:ASP:CB	28:11:103:ARG:HH22	2.18	0.56
26:14:1864:U:OP1	26:14:2410:G:O2'	2.14	0.56
1:13:652:U:O4	1:13:752:G:O2'	2.14	0.56
26:14:920:G:H2'	26:14:921:G:C8	2.40	0.56
30:39:64:ILE:HG13	30:39:65:TRP:CD1	2.40	0.56
4:32:76:ARG:NH2	4:32:80:GLU:OE1	2.37	0.56
40:75:99:LEU:HD22	40:75:101:PHE:HE1	1.71	0.56
44:F8:12:VAL:HG13	44:F8:27:THR:O	2.06	0.56
4:3E:155:LEU:HD12	4:3E:158:ILE:HD11	1.86	0.56
31:41:33:ARG:O	31:41:162:THR:HG23	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:41:161:THR:HG22	31:41:163:ALA:H	1.68	0.56
26:1H:2566:A:H4'	26:1H:2567:G:O5'	2.05	0.56
1:13:390:C:O3'	16:7I:28:ARG:NH2	2.38	0.56
26:14:667:U:O2	55:M5:2:PRO:HD2	2.05	0.56
26:14:575:A:OP2	26:14:2055:C:N4	2.33	0.56
4:32:53:ASP:O	4:32:57:ARG:NH1	2.39	0.56
26:1H:1016:G:O6	58:1H:4298:HOH:O	2.17	0.56
1:1G:222:U:H2'	1:1G:223:U:C6	2.40	0.56
30:39:183:VAL:O	30:39:187:VAL:HG23	2.05	0.56
26:1H:2257:U:O2'	26:1H:2258:C:H5'	2.05	0.56
32:59:152:ARG:HD2	32:59:153:LYS:HG3	1.86	0.56
7:62:115:ARG:O	7:62:118:VAL:HG22	2.04	0.56
55:Q8:48:PHE:O	55:Q8:49:VAL:HG22	2.05	0.56
36:78:49:ARG:HA	55:Q8:54:GLU:HG2	1.88	0.56
1:1G:1321:C:C2	1:1G:1322:C:H5	2.24	0.56
1:13:36:C:O2'	12:3I:117:ARG:NH2	2.38	0.56
51:I5:12:ALA:O	51:I5:24:THR:OG1	2.22	0.56
34:58:43:THR:O	34:58:46:VAL:HG12	2.05	0.56
55:Q8:34:TRP:HB3	55:Q8:36:LYS:H	1.69	0.56
53:O8:44:ARG:O	53:O8:45:LYS:HG2	2.05	0.56
45:G8:82:PRO:HG3	45:G8:97:ARG:HG3	1.88	0.56
40:B8:57:PHE:O	40:B8:58:ASN:ND2	2.38	0.56
26:14:823:G:H2'	26:14:824:A:C8	2.41	0.56
36:78:130:PHE:HE1	36:78:146:VAL:HG23	1.69	0.56
26:14:2311:A:H62	31:49:44:GLY:HA3	1.71	0.56
26:1H:1608:A:H4'	58:1H:3840:HOH:O	2.04	0.56
34:15:128:HIS:CE1	34:15:130:HIS:HA	2.41	0.56
26:14:527:C:OP2	26:14:2779:U:H5	1.89	0.56
1:13:1149:C:H2'	1:13:1150:U:H6	1.71	0.56
26:1H:322:A:H5'	26:1H:340:A:H1'	1.87	0.56
3:22:138:VAL:HG23	3:22:151:VAL:HG23	1.86	0.56
26:14:61:G:OP2	49:G5:54:LYS:NZ	2.39	0.56
30:39:181:LEU:HD23	30:39:186:ILE:HD11	1.87	0.56
12:3I:62:SER:HB2	12:3I:64:TYR:CD1	2.40	0.56
26:14:1167:U:C2	26:14:1183:G:N2	2.73	0.56
24:3K:30:G:N2	24:3K:40:C:O2	2.36	0.56
55:Q8:48:PHE:CD2	55:Q8:49:VAL:HG13	2.41	0.56
55:Q8:5:LYS:H	55:Q8:59:LYS:HZ2	1.53	0.56
26:1H:1190:G:N7	58:1H:3874:HOH:O	2.33	0.56
42:95:69:LYS:HB2	42:95:86:GLY:HA3	1.87	0.56
46:H8:67:LEU:HD22	46:H8:90:VAL:HG11	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:277:C:P	17:8I:68:ARG:HH12	2.29	0.56
26:1H:763:G:O2'	26:1H:764:A:H3'	2.05	0.56
32:51:4:ILE:HD11	32:51:7:LEU:HD21	1.87	0.56
38:98:63:ARG:HB2	38:98:80:PHE:CE2	2.40	0.56
24:3L:11:C:H2'	24:3L:12:U:C6	2.37	0.56
26:14:2134:A:H62	26:14:2157:G:H1'	1.69	0.56
39:A8:88:ASP:OD1	39:A8:90:GLY:N	2.33	0.56
7:6E:5:ARG:HG2	7:6E:7:ALA:H	1.70	0.56
26:14:854:G:H2'	26:14:855:G:C8	2.39	0.56
32:59:152:ARG:HG3	32:59:153:LYS:HB2	1.88	0.56
26:14:736:C:OP1	58:14:3867:HOH:O	2.18	0.56
26:14:746:A:H2'	26:14:2612:C:H5''	1.87	0.56
53:K5:9:LEU:HD13	53:K5:11:LEU:HD21	1.85	0.56
6:52:82:ARG:HB2	6:52:85:VAL:HG23	1.86	0.56
6:5E:16:GLN:HG2	6:5E:17:SER:N	2.19	0.56
53:K5:36:LEU:HD23	53:K5:50:ARG:HD3	1.87	0.56
26:14:2277:G:H2'	26:14:2278:A:H5''	1.87	0.56
33:69:112:LYS:HA	33:69:114:LEU:H	1.69	0.56
7:6E:150:ALA:HB2	11:2I:50:TYR:OH	2.05	0.56
30:31:164:ARG:HH11	30:31:164:ARG:HG2	1.71	0.56
11:2A:100:ALA:O	11:2A:102:GLY:N	2.38	0.56
1:13:828:A:H2'	1:13:829:G:O4'	2.06	0.56
1:1G:1443:G:O2'	40:75:122:ASP:OD2	2.23	0.56
36:78:58:THR:HG21	55:Q8:52:LYS:HE2	1.86	0.56
1:13:963:G:H21	10:1I:55:LYS:NZ	2.03	0.56
5:4E:110:LEU:CD1	5:4E:118:ILE:HD13	2.35	0.56
16:7A:57:ARG:HG3	16:7A:79:VAL:HG13	1.86	0.56
26:1H:993:G:C4	26:1H:994:C:H5	2.23	0.56
46:D5:138:GLU:O	46:D5:156:LYS:HG3	2.05	0.56
26:14:328:U:H4'	45:C5:68:HIS:CD2	2.40	0.56
26:1H:1221:C:H2'	26:1H:1222:C:H6	1.70	0.56
4:32:8:VAL:O	4:32:11:LEU:N	2.32	0.56
1:1G:713:G:H2'	1:1G:714:G:C8	2.39	0.56
13:4I:105:THR:OG1	13:4I:106:ASN:N	2.39	0.56
1:1G:750:G:N3	15:6A:23:GLY:HA3	2.20	0.56
26:1H:1748:G:H2'	26:1H:1749:A:C8	2.40	0.56
26:14:959:A:N6	26:14:960:A:N1	2.54	0.56
37:45:32:TYR:O	37:45:105:GLU:HB2	2.05	0.56
26:14:2273:A:H2'	26:14:2274:A:C8	2.41	0.56
55:Q8:48:PHE:N	55:Q8:48:PHE:CD1	2.74	0.56
36:35:85:LEU:HB3	36:35:114:ILE:HD11	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:41:67:LYS:HE3	31:41:67:LYS:O	2.06	0.56
41:C8:92:ARG:HD2	42:D8:11:GLN:HB2	1.87	0.56
36:78:52:GLU:OE1	36:78:55:ARG:NE	2.28	0.56
26:14:966:G:H2'	26:14:967:C:H6	1.70	0.56
40:75:55:ASN:ND2	40:75:55:ASN:O	2.31	0.56
1:1G:680:C:N4	1:1G:710:G:H1	2.02	0.56
1:13:1278:U:H5''	1:13:1279:A:O4'	2.05	0.56
21:1F:9:ARG:O	21:1F:13:ILE:HG13	2.06	0.56
28:19:52:ARG:HB2	28:19:53:PHE:CD2	2.41	0.56
4:3E:173:TRP:CZ3	4:3E:193:ASP:HB3	2.41	0.56
32:59:130:ARG:HH12	32:59:132:ARG:HH22	1.51	0.56
45:G8:9:LYS:HA	45:G8:27:VAL:HG22	1.88	0.56
52:J5:3:LYS:HE3	52:J5:3:LYS:HA	1.88	0.56
8:72:110:ALA:HB3	8:72:121:ASP:HB3	1.86	0.56
42:95:71:LEU:H	42:95:86:GLY:CA	2.16	0.56
1:13:1022:G:H2'	1:13:1023:G:H8	1.71	0.56
26:1H:442:G:C4	26:1H:444:C:C5	2.94	0.56
9:82:112:LYS:HA	9:82:119:ALA:CB	2.35	0.56
33:69:73:GLU:HG3	33:69:136:VAL:HG23	1.87	0.56
26:14:2120:G:H2'	26:14:2121:G:H8	1.70	0.56
26:14:2119:A:N6	26:14:2170:A:N7	2.53	0.56
55:M5:34:TRP:CG	55:M5:35:GLN:N	2.61	0.56
46:D5:152:ALA:HB2	46:D5:169:GLU:O	2.06	0.56
2:12:9:GLU:HA	2:12:12:GLU:HG3	1.88	0.56
2:1E:189:ASP:HB2	2:1E:205:ASP:CG	2.27	0.56
49:K8:63:VAL:HA	49:K8:66:GLU:HG3	1.88	0.56
14:5I:23:ARG:HD2	14:5I:28:GLY:O	2.05	0.56
1:1G:737:A:H2'	1:1G:738:C:H6	1.69	0.56
26:14:2880:C:H1'	38:55:92:GLY:HA3	1.87	0.56
26:1H:2498:C:O2'	26:1H:2499:C:H5'	2.06	0.56
1:1G:1513:A:H2'	1:1G:1514:C:C6	2.39	0.56
4:3E:61:LYS:HG3	4:3E:203:VAL:HG13	1.87	0.56
26:14:1514:U:O2'	26:14:1515:C:H5'	2.06	0.56
42:95:60:GLU:HG3	42:95:97:LYS:HE2	1.86	0.56
26:14:1268:A:H2'	26:14:1269:A:O4'	2.05	0.56
24:3K:37:A:O2'	24:3K:38:A:O5'	2.23	0.56
1:1G:167:G:H2'	1:1G:168:G:H8	1.70	0.56
12:3I:58:VAL:O	12:3I:65:GLU:HA	2.06	0.56
43:E8:58:ALA:O	43:E8:64:MET:HB2	2.06	0.56
26:1H:2594:C:H2'	26:1H:2595:G:C8	2.41	0.56
1:13:1508:G:P	58:13:1803:HOH:O	2.64	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:B8:26:ASP:O	40:B8:49:VAL:HG12	2.06	0.56
26:1H:399:G:OP2	58:1H:4043:HOH:O	2.18	0.56
1:13:767:A:H3'	58:13:1840:HOH:O	2.05	0.56
1:1G:1239:A:H62	1:1G:1299:A:H62	1.54	0.56
42:95:35:LEU:HB3	42:95:37:VAL:HG11	1.88	0.56
29:21:79:ARG:HG3	29:21:80:GLU:O	2.06	0.56
29:21:119:ARG:HH11	29:21:120:TRP:HE1	1.54	0.56
24:3L:13:C:H2'	24:3L:14:A:H8	1.71	0.56
1:1G:1305:G:H22	1:1G:1331:G:C2'	2.19	0.56
26:14:309:G:O3'	45:C5:18:GLY:HA3	2.05	0.56
24:3L:71:G:H2'	24:3L:72:C:O4'	2.05	0.56
1:13:919:A:O2'	1:13:920:U:H5'	2.06	0.56
26:1H:1349:A:OP1	58:1H:4126:HOH:O	2.18	0.56
2:12:109:SER:HA	2:12:112:VAL:HG23	1.86	0.56
2:1E:59:GLU:HG3	2:1E:225:ALA:HB2	1.87	0.56
26:14:1966:A:H4'	26:14:1967:C:OP1	2.06	0.56
26:14:1693:U:O2'	28:19:14:ARG:NH2	2.39	0.56
6:52:6:VAL:HG13	6:52:90:VAL:HG22	1.87	0.56
26:1H:185:U:H4'	26:1H:218:A:H4'	1.88	0.56
26:14:1427:A:H4'	26:14:1428:C:O4'	2.05	0.56
28:11:52:ARG:HB2	28:11:53:PHE:CD2	2.41	0.56
26:14:2294:C:OP1	39:65:89:ARG:NH2	2.39	0.56
37:45:66:ILE:HG13	37:45:67:ARG:H	1.69	0.56
1:13:804:U:H5''	1:13:805:C:OP2	2.05	0.56
26:1H:2431:U:H3'	58:1H:3905:HOH:O	2.05	0.56
36:35:64:LYS:HB2	55:M5:30:ARG:HH22	1.70	0.56
51:15:22:ILE:HG12	51:15:23:GLU:N	2.20	0.56
26:1H:674:G:C1'	30:31:74:ARG:HD3	2.35	0.56
26:14:649:G:H2'	26:14:650:C:C6	2.41	0.56
26:1H:2371:G:H4'	53:O8:45:LYS:HG3	1.88	0.56
48:F5:86:SER:N	48:F5:87:PRO:HD2	2.20	0.56
33:69:123:LEU:HD22	33:69:143:SER:HB3	1.88	0.56
23:2K:20:G:C2	23:2K:58:A:N3	2.74	0.56
28:19:153:ALA:O	28:19:157:ARG:NH1	2.37	0.56
36:78:122:PRO:HA	36:78:142:GLY:CA	2.36	0.56
9:82:24:GLY:HA2	9:82:59:PHE:O	2.06	0.56
37:45:87:LYS:HB3	37:45:90:VAL:HG23	1.88	0.56
31:41:47:LYS:HD3	31:41:81:LYS:CB	2.36	0.56
26:14:1496:A:H8	26:14:1577:C:HO2'	1.51	0.56
26:1H:2629:A:OP1	26:1H:2629:A:H4'	2.05	0.56
26:14:214:G:OP1	26:14:214:G:H4'	2.03	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:58:12:ARG:HH21	34:58:14:VAL:HG22	1.71	0.56
26:1H:1514:U:H2'	26:1H:1515:C:H6	1.71	0.56
1:1G:1075:C:OP1	2:12:179:LYS:NZ	2.30	0.56
26:1H:241:A:H5''	58:1H:4304:HOH:O	2.06	0.56
28:19:71:ASP:OD2	28:19:103:ARG:NH1	2.32	0.56
43:E8:88:ARG:NH1	43:E8:94:ASP:OD2	2.31	0.55
1:1G:1220:G:H5'	19:AA:35:SER:HA	1.87	0.55
26:1H:1163:G:C2	26:1H:1164:G:C8	2.95	0.55
26:14:2324:C:H5''	26:14:2325:G:H5'	1.88	0.55
1:1G:827:U:H3	1:1G:872:A:N6	2.00	0.55
31:49:42:GLY:O	31:49:43:LEU:HD13	2.04	0.55
45:G8:87:LYS:HD2	45:G8:96:ILE:HD11	1.87	0.55
26:1H:2695:C:H2'	26:1H:2696:U:H6	1.71	0.55
53:O8:15:GLU:OE2	53:O8:44:ARG:NH2	2.39	0.55
1:1G:538:G:H3'	12:3A:115:LYS:HZ3	1.71	0.55
2:12:115:LEU:HB2	2:12:145:LEU:HD12	1.87	0.55
19:AI:42:PRO:HD3	51:M8:63:TYR:OH	2.06	0.55
2:1E:189:ASP:HB3	2:1E:191:ASP:HB2	1.89	0.55
6:52:100:ASN:ND2	18:9A:26:LEU:O	2.39	0.55
26:1H:2883:A:H5'	26:1H:2884:U:H5'	1.87	0.55
26:14:2875:C:OP1	40:75:3:ARG:NH1	2.39	0.55
47:E5:51:VAL:HG23	47:E5:81:VAL:HG23	1.87	0.55
40:B8:77:PRO:HG2	40:B8:80:SER:HB2	1.87	0.55
55:Q8:48:PHE:CD1	55:Q8:49:VAL:HG13	2.42	0.55
26:14:2378:A:O5'	26:14:2378:A:H8	1.89	0.55
26:1H:2019:A:C6	26:1H:2020:A:N7	2.74	0.55
43:E8:12:ILE:HG13	43:E8:42:ARG:NH1	2.22	0.55
26:14:2029:G:H2'	26:14:2031:A:OP1	2.07	0.55
45:G8:94:LYS:HZ2	45:G8:95:LYS:H	1.53	0.55
1:13:113:G:H2'	1:13:114:U:H6	1.71	0.55
1:1G:649:G:H2'	1:1G:650:G:H8	1.71	0.55
51:I5:20:ASN:HD21	51:I5:36:CYS:HB2	1.70	0.55
29:21:97:LYS:N	29:21:100:GLU:OE1	2.39	0.55
1:13:1013:G:N2	1:13:1016:A:OP2	2.40	0.55
31:41:35:GLU:OE1	31:41:36:LYS:N	2.39	0.55
26:14:573:G:O2'	26:14:574:C:H3'	2.06	0.55
3:2E:58:GLU:HB2	3:2E:65:ALA:HB3	1.88	0.55
26:1H:50:U:H3'	26:1H:51:G:H5'	1.88	0.55
30:31:129:PHE:HA	30:31:142:TRP:NE1	2.21	0.55
30:31:40:GLN:NE2	30:31:184:TYR:HB3	2.21	0.55
48:J8:3:LYS:O	48:J8:12:PRO:HD3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1128:C:H6	1:13:1139:G:C6	2.24	0.55
36:35:13:ASN:C	36:35:15:ARG:H	2.08	0.55
36:35:15:ARG:CZ	36:35:15:ARG:HB2	2.36	0.55
36:35:85:LEU:HB2	36:35:88:LEU:HD23	1.88	0.55
24:3K:76:A:H8	26:1H:2394:C:N4	2.01	0.55
24:3L:52:G:N2	24:3L:62:C:N3	2.39	0.55
26:1H:2572:A:N7	29:21:145:LYS:HB2	2.22	0.55
5:4E:126:ARG:HH11	5:4E:126:ARG:HG3	1.71	0.55
37:45:25:ASP:CB	37:45:102:VAL:H	2.20	0.55
26:14:2345:G:N3	26:14:2381:C:H2'	2.21	0.55
1:13:724:G:C2	1:13:725:G:C8	2.94	0.55
1:13:654:G:C2'	1:13:655:A:H5'	2.36	0.55
1:1G:853:G:H2'	1:1G:854:G:H8	1.71	0.55
17:8A:99:SER:O	17:8A:100:LYS:NZ	2.36	0.55
1:1G:196:A:OP1	20:BA:68:LYS:NZ	2.40	0.55
1:1G:1128:C:H4'	9:82:16:ARG:HH22	1.72	0.55
27:1J:3:C:H2'	27:1J:4:C:C6	2.42	0.55
26:14:2052:G:O4'	29:29:142:GLY:HA3	2.06	0.55
26:14:1419:A:N6	26:14:1494:A:N1	2.45	0.55
13:4I:65:LYS:HZ2	51:M8:52:THR:HB	1.70	0.55
26:14:1945:G:H2'	26:14:1946:U:C6	2.39	0.55
26:14:2564:A:OP1	26:14:2648:C:H4'	2.06	0.55
2:12:12:GLU:HB2	2:12:16:HIS:CG	2.42	0.55
14:5I:24:CYS:HB3	14:5I:28:GLY:H	1.71	0.55
1:1G:1255:G:O2'	1:1G:1258:G:O2'	2.25	0.55
1:13:669:U:C2	1:13:670:G:C8	2.95	0.55
26:1H:1279:G:H4'	38:98:31:HIS:CD2	2.40	0.55
26:1H:686:G:OP1	54:P8:11:LYS:NZ	2.40	0.55
26:1H:831:G:N2	36:78:53:GLY:O	2.39	0.55
26:1H:1433:U:O2	26:1H:1561:G:C2	2.60	0.55
27:1J:38:C:H42	27:1J:44:G:H1	1.52	0.55
1:13:419:C:H5'	1:13:420:U:OP2	2.07	0.55
15:6I:17:ARG:HD3	15:6I:26:GLU:HG3	1.87	0.55
26:14:1233:C:H2'	26:14:1234:U:H6	1.71	0.55
6:52:70:ASP:OD1	6:52:70:ASP:N	2.39	0.55
1:13:187:C:O2	1:13:191(A):G:N1	2.39	0.55
26:1H:2645:G:H3'	26:1H:2646:C:H5'	1.88	0.55
55:Q8:48:PHE:N	55:Q8:48:PHE:HD1	2.04	0.55
22:1L:18:G:O2'	22:1L:19:G:OP1	2.24	0.55
1:13:1157:A:H8	1:13:1158:C:N4	2.03	0.55
1:1G:1298:C:H1'	1:1G:1299:A:C6	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
31:41:62:LEU:HD12	51:M8:27:THR:HG21	1.89	0.55
5:42:107:ARG:O	5:42:110:LEU:N	2.39	0.55
26:1H:1668:A:H4'	26:1H:1669:A:O5'	2.05	0.55
46:D5:70:LEU:O	46:D5:89:PHE:N	2.39	0.55
1:1G:1326:C:H2'	1:1G:1327:C:C6	2.41	0.55
8:7E:7:ALA:HB2	8:7E:85:ARG:HH11	1.72	0.55
8:7E:6:ILE:HB	8:7E:85:ARG:HH12	1.71	0.55
1:1G:513:C:N4	1:1G:538:G:H1	2.03	0.55
27:16:78:A:C2	27:16:99:A:C4	2.94	0.55
1:13:406:G:H2'	1:13:407:G:C8	2.40	0.55
3:2E:3:ASN:O	3:2E:4:LYS:HG2	2.07	0.55
26:14:315:G:H2'	26:14:316:C:C6	2.40	0.55
1:1G:532:A:H62	1:1G:1207:G:H5'	1.71	0.55
55:M5:60:LEU:HD12	55:M5:61:LEU:H	1.71	0.55
22:1L:23:A:H2'	22:1L:24:G:H8	1.72	0.55
26:14:2256:G:H2'	26:14:2257:U:H6	1.71	0.55
26:1H:638:G:C5	26:1H:651:G:C2	2.95	0.55
27:16:15:A:H3'	27:16:16:G:H5'	1.88	0.55
26:14:2015:A:N3	52:J5:2:ALA:N	2.54	0.55
26:1H:1441:G:H2'	26:1H:1442:G:H8	1.71	0.55
26:14:755:C:H2'	26:14:756:C:C6	2.41	0.55
2:12:127:ILE:HA	2:12:130:ARG:HG3	1.88	0.55
1:1G:1478:C:H2'	1:1G:1479:C:H6	1.71	0.55
1:1G:994:A:C5	1:1G:1216:G:H4'	2.40	0.55
2:1E:15:VAL:HG21	2:1E:209:ARG:HB3	1.89	0.55
2:12:162:ILE:HD11	2:12:184:VAL:HG13	1.88	0.55
32:59:149:ARG:HH21	32:59:154:PRO:HB2	1.72	0.55
18:9I:66:LEU:O	18:9I:70:ILE:HG13	2.07	0.55
11:2I:19:ALA:O	11:2I:82:VAL:HA	2.05	0.55
26:1H:2392:A:C8	36:78:61:ARG:HG2	2.40	0.55
55:Q8:48:PHE:HD1	55:Q8:48:PHE:H	1.54	0.55
1:13:1160:G:N2	1:13:1177:G:H22	2.03	0.55
1:13:1160:G:H1	1:13:1177:G:N2	2.04	0.55
26:14:832:G:H5'	36:35:45:LEU:HD11	1.89	0.55
36:35:63:PRO:O	55:M5:30:ARG:NH2	2.40	0.55
32:51:77:LYS:NZ	32:51:82:GLY:O	2.37	0.55
26:14:1360:A:H5''	26:14:1361:G:OP2	2.07	0.55
45:G8:76:CYS:HB2	45:G8:82:PRO:HD3	1.88	0.55
3:22:6:HIS:HB3	14:5A:49:HIS:CD2	2.42	0.55
45:C5:87:LYS:HG2	45:C5:88:LYS:H	1.71	0.55
36:35:127:ALA:O	36:35:147:LEU:N	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1410:G:H2'	1:13:1411:C:C6	2.41	0.55
6:5E:100:ASN:HB2	18:9I:27:GLY:O	2.05	0.55
42:D8:3:ALA:HB1	42:D8:38:LEU:HD11	1.89	0.55
26:1H:1788:C:H2'	26:1H:1789:A:H8	1.71	0.55
26:14:247:G:H4'	26:14:386:G:C5	2.41	0.55
26:1H:1694:C:H4'	26:1H:1695:G:O5'	2.06	0.55
7:62:146:GLU:O	7:62:149:ARG:HB2	2.07	0.55
26:1H:484:C:H2'	26:1H:485:C:C6	2.41	0.55
26:14:2439:A:H5'	26:14:2439:A:C8	2.40	0.55
20:BA:73:HIS:HB3	20:BA:74:LYS:HG2	1.89	0.55
34:15:136:GLU:O	34:15:137:LYS:NZ	2.40	0.55
26:1H:969:U:OP1	50:L8:17:LYS:HG2	2.06	0.55
1:13:343:U:N3	1:13:347:G:C6	2.74	0.55
33:61:102:SER:HA	33:61:107:VAL:O	2.07	0.55
1:1G:111:G:H8	1:1G:111:G:O5'	1.89	0.55
1:1G:1435:G:H2'	1:1G:1436:U:H6	1.63	0.55
7:62:23:VAL:O	7:62:27:ILE:HG12	2.07	0.55
19:AI:78:ARG:N	19:AI:78:ARG:HE	2.03	0.55
26:1H:1797:C:O2'	28:11:259:THR:HB	2.07	0.55
26:1H:225:A:N6	26:1H:419:C:H4'	2.22	0.55
1:13:1014:A:C2	1:13:1219:U:H1'	2.42	0.55
40:75:24:PRO:HD3	40:75:52:ILE:HG13	1.89	0.55
33:61:120:ILE:HD12	33:61:121:LYS:H	1.71	0.55
26:14:289:A:H3'	26:14:290:G:H8	1.71	0.55
1:13:452:A:H2'	1:13:453:A:C8	2.42	0.55
53:K5:9:LEU:N	53:K5:27:LYS:HG3	2.21	0.55
20:BA:16:HIS:O	20:BA:19:SER:N	2.39	0.55
26:14:646:A:H2'	26:14:647:G:O4'	2.07	0.55
46:H8:9:TYR:CE1	46:H8:35:ARG:HD3	2.41	0.55
29:29:92:THR:O	29:29:95:ILE:HG13	2.06	0.55
26:14:208:C:H2'	26:14:209:C:H6	1.71	0.55
26:14:2528:U:O2'	26:14:2530:A:OP1	2.15	0.55
32:59:144:VAL:HA	32:59:147:ASN:HB2	1.88	0.55
42:95:76:LYS:HB2	42:95:79:VAL:HG23	1.88	0.55
26:14:1666:G:OP1	35:25:66:LYS:HD3	2.07	0.55
55:Q8:23:VAL:HG13	55:Q8:46:ARG:HG3	1.89	0.55
55:Q8:52:LYS:N	55:Q8:52:LYS:HD2	2.21	0.55
26:14:1689:A:N6	26:14:1698:A:H2	1.92	0.55
24:3K:71:G:H2'	24:3K:72:C:O4'	2.07	0.55
24:3L:18:G:H2'	24:3L:57:G:H22	1.72	0.55
44:F8:24:GLY:O	44:F8:83:VAL:HG22	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2171:A:O2'	26:1H:2172:U:O5'	2.24	0.55
1:1G:1281:U:H3'	1:1G:1282:C:C5	2.42	0.55
26:14:2697:G:H2'	26:14:2698:U:O4'	2.07	0.55
8:72:29:SER:H	8:72:32:LYS:HB2	1.71	0.55
1:13:130:A:O2'	1:13:131:C:O5'	2.24	0.55
8:7E:83:ILE:HB	8:7E:137:VAL:HG13	1.88	0.55
46:D5:8:TYR:HD1	46:D5:62:PRO:HG3	1.70	0.55
38:55:81:ASP:O	38:55:82:GLU:HB3	2.07	0.55
39:A8:26:LEU:HD23	39:A8:87:PHE:HD1	1.70	0.55
26:14:606:U:H4'	26:14:658:C:H4'	1.89	0.55
26:14:2494:G:H2'	26:14:2495:G:H8	1.71	0.55
19:AI:50:ALA:HB1	19:AI:57:HIS:HB3	1.88	0.55
29:21:2:LYS:HD2	29:21:200:GLU:OE2	2.07	0.55
41:85:8:VAL:HG11	41:85:12:ARG:HH21	1.72	0.55
28:19:20:ASP:OD2	28:19:22:SER:OG	2.24	0.55
29:29:111:ARG:HA	38:55:2:ARG:HH12	1.70	0.55
37:88:68:ILE:HG23	37:88:103:MET:HB3	1.88	0.55
28:19:43:ARG:HH11	28:19:43:ARG:CG	2.19	0.55
26:14:337:C:H2'	26:14:338:G:O4'	2.06	0.55
28:19:118:VAL:HG22	28:19:119:ALA:H	1.72	0.55
3:22:34:LEU:HD11	3:22:38:ARG:HH21	1.71	0.55
19:AA:18:LYS:O	19:AA:22:LEU:HB2	2.07	0.55
12:3A:100:ILE:HG22	12:3A:101:VAL:N	2.21	0.55
26:14:2432:A:C8	48:F5:33:LYS:HD2	2.41	0.55
29:29:81:ILE:HG22	29:29:82:ARG:N	2.18	0.55
24:3K:70:G:C6	24:3K:71:G:C5	2.95	0.55
33:61:4:ILE:HG12	33:61:18:VAL:HG22	1.89	0.55
1:1G:588:G:H1	1:1G:651:C:N4	2.02	0.55
26:1H:1803:A:H4'	28:11:259:THR:CG2	2.36	0.55
26:1H:2032:G:H21	29:21:146:THR:CG2	2.20	0.55
26:14:2420:C:N4	55:M5:31:HIS:HB3	2.22	0.55
26:1H:2157:G:O2'	26:1H:2158:A:O5'	2.21	0.55
1:1G:1305:G:H22	1:1G:1331:G:H2'	1.72	0.55
26:14:972:G:O2'	58:14:3594:HOH:O	2.18	0.55
40:75:54:ARG:HA	40:75:59:THR:HB	1.89	0.55
26:14:986:C:O2'	26:14:1001:A:O2'	2.25	0.55
32:51:154:PRO:HD3	32:51:162:ILE:O	2.06	0.55
26:1H:582:G:H2'	26:1H:583:G:H8	1.72	0.55
29:29:101:ARG:O	29:29:201:THR:OG1	2.24	0.55
26:1H:2376:A:N1	39:A8:87:PHE:HD2	2.05	0.55
15:6A:39:LEU:HD12	15:6A:56:LEU:HD13	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1485:G:H2'	26:14:1486:A:H8	1.70	0.55
26:14:2294:C:P	39:65:89:ARG:HH22	2.30	0.55
13:4I:108:ARG:N	13:4I:108:ARG:HD2	2.20	0.55
3:2E:191:THR:HG22	3:2E:192:THR:H	1.72	0.55
14:5I:4:LYS:O	14:5I:7:ILE:HG13	2.06	0.55
44:B5:84:ALA:O	44:B5:87:GLN:HG3	2.06	0.55
26:14:2363:C:O2	47:E5:39:ARG:NH2	2.39	0.55
2:1E:174:VAL:HG13	2:1E:184:VAL:HG11	1.88	0.55
26:1H:673:C:H5''	30:31:81:PRO:HD2	1.88	0.55
26:1H:937:U:H2'	26:1H:938:G:O4'	2.07	0.55
26:14:871:U:OP1	37:45:5:ARG:HG3	2.07	0.55
1:1G:1411:C:H2'	1:1G:1412:C:C6	2.42	0.55
1:13:814:A:N7	1:13:816:A:C4	2.75	0.55
10:1A:32:ALA:HA	10:1A:76:ASN:HB2	1.88	0.55
26:14:2540:C:O2'	26:14:2740:A:N3	2.37	0.55
26:14:1537:C:H2'	26:14:1538:G:C8	2.41	0.55
26:14:2103:C:H2'	26:14:2104:G:C8	2.42	0.55
40:75:21:GLU:O	40:75:91:ARG:NH2	2.37	0.55
1:1G:1320:C:H1'	19:AA:73:GLU:HG2	1.89	0.55
26:1H:1478:G:O2'	26:1H:1558:A:H2	1.90	0.55
51:I5:34:GLU:HG2	51:I5:35:VAL:N	2.22	0.55
39:65:78:LEU:HD12	39:65:107:GLU:HB3	1.89	0.55
34:58:58:ASP:HB2	34:58:95:PRO:HB2	1.89	0.55
1:1G:1125:U:OP2	1:1G:1145:C:N4	2.39	0.55
9:82:77:ILE:O	9:82:81:ILE:HG12	2.07	0.55
26:1H:2341:G:H2'	26:1H:2342:C:C6	2.42	0.55
40:B8:58:ASN:ND2	40:B8:58:ASN:O	2.36	0.55
26:14:975:G:C2	26:14:990:A:C8	2.95	0.55
26:14:739:G:P	58:14:3635:HOH:O	2.65	0.55
2:12:219:VAL:O	2:12:222:ILE:HG13	2.07	0.55
26:1H:1062:G:H1'	26:1H:1088:A:C5	2.42	0.55
1:1G:1365:G:H2'	1:1G:1366:C:H6	1.71	0.55
1:13:486:U:H2'	1:13:487:A:H8	1.71	0.55
1:13:438:G:H4'	4:3E:123:HIS:CD2	2.41	0.55
53:O8:36:LEU:HD22	53:O8:50:ARG:HG2	1.88	0.55
35:68:75:SER:OG	40:B8:74:ARG:NH2	2.40	0.55
46:D5:175:VAL:HA	46:D5:177:PRO:HD3	1.88	0.55
26:1H:796:C:H2'	26:1H:797:C:C6	2.42	0.55
12:3A:41:ARG:HB3	12:3A:41:ARG:HH11	1.72	0.55
26:1H:881:G:O6	26:1H:895:U:N3	2.38	0.54
26:1H:2773:C:H5''	29:21:164:ARG:HG2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:674:G:H21	11:2A:116:HIS:HB2	1.71	0.54
1:13:826:C:H2'	1:13:827:U:O2	2.07	0.54
1:13:626:U:C2	1:13:627:G:C8	2.95	0.54
1:13:417:C:H2'	1:13:418:C:C6	2.42	0.54
15:6A:55:GLY:HA2	15:6A:58:MET:HE2	1.88	0.54
18:9I:44:LEU:HD11	18:9I:70:ILE:HG21	1.88	0.54
8:7E:21:LYS:O	8:7E:65:TYR:OH	2.22	0.54
1:1G:527:G:O6	12:3A:49:ASN:ND2	2.37	0.54
4:32:18:LYS:HG3	4:32:20:TYR:H	1.71	0.54
26:14:1034:G:H2'	26:14:1035:U:O4'	2.07	0.54
52:N8:50:GLY:H	52:N8:56:LYS:HG3	1.72	0.54
29:21:48:GLN:OE1	29:21:77:ILE:HD13	2.07	0.54
26:1H:1055:G:H1'	26:1H:1085:A:H2	1.71	0.54
55:Q8:9:GLY:N	55:Q8:12:LYS:HG3	2.23	0.54
6:5E:61:LEU:HD23	6:5E:63:TYR:CE1	2.41	0.54
16:7I:77:ALA:HB3	16:7I:79:VAL:H	1.72	0.54
1:13:376:G:H5''	16:7I:5:ARG:HD2	1.89	0.54
27:1J:52:A:O2'	27:1J:53:A:N7	2.37	0.54
26:1H:1265:A:H3'	52:N8:19:ARG:NH1	2.23	0.54
49:G5:10:LEU:O	49:G5:14:ARG:HB2	2.08	0.54
1:13:392:G:H5'	16:7I:12:LYS:HE3	1.89	0.54
46:D5:158:PRO:HD2	46:D5:161:VAL:HG13	1.88	0.54
38:55:70:LEU:HD13	38:55:75:LEU:HD13	1.89	0.54
28:19:3:VAL:H	28:19:20:ASP:HB2	1.72	0.54
29:29:68:ALA:O	29:29:69:LYS:HB2	2.07	0.54
1:1G:888:G:O2'	1:1G:1488:G:O2'	2.23	0.54
33:69:101:LEU:HB2	33:69:105:HIS:HB2	1.88	0.54
31:49:37:VAL:HG23	31:49:99:MET:HE3	1.89	0.54
4:3E:117:ALA:O	4:3E:120:LEU:HB2	2.08	0.54
1:1G:881:G:P	12:3A:12:ARG:HH12	2.29	0.54
26:14:2295:C:H5	39:65:13:ARG:HH22	1.55	0.54
26:1H:1639:U:H4'	26:1H:2699:C:H4'	1.89	0.54
1:1G:1248:A:H2'	9:82:70:LYS:NZ	2.22	0.54
1:1G:1279:A:HO2'	1:1G:1282:C:N4	2.05	0.54
23:2K:2:G:H2'	23:2K:3:C:C6	2.42	0.54
26:14:571:A:OP2	58:14:4005:HOH:O	2.18	0.54
46:D5:40:ASP:OD2	46:D5:43:GLU:HG2	2.07	0.54
26:14:1337:G:H2'	26:14:1338:G:C8	2.42	0.54
26:14:839:U:H2'	26:14:840:C:H6	1.72	0.54
1:13:240:C:H2'	1:13:241:C:C6	2.41	0.54
15:6A:50:HIS:O	15:6A:53:HIS:HB3	2.06	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2849:U:H4'	26:14:2868:A:C2	2.42	0.54
26:14:920:G:H2'	26:14:921:G:H8	1.72	0.54
17:8I:22:LEU:HD11	17:8I:39:SER:HB3	1.89	0.54
29:21:143:ASN:HD22	29:21:147:PRO:HD2	1.72	0.54
26:14:1686:C:H2'	26:14:1687:G:O4'	2.07	0.54
1:1G:141:A:H1'	1:1G:182:U:O2	2.08	0.54
17:8I:31:LEU:HD23	17:8I:32:TYR:CZ	2.42	0.54
1:1G:607:A:H2'	1:1G:608:A:O4'	2.07	0.54
3:22:150:LYS:HG3	3:22:169:ALA:HB2	1.90	0.54
27:16:89:G:H2'	27:16:89(A):A:C8	2.43	0.54
40:B8:93:ARG:HH11	40:B8:93:ARG:HG3	1.70	0.54
39:65:85:VAL:HG22	39:65:110:LEU:HG	1.88	0.54
26:1H:1641:A:H2'	26:1H:1642:G:O4'	2.08	0.54
26:1H:428:A:P	58:1H:4617:HOH:O	2.66	0.54
26:1H:2035:G:H4'	26:1H:2036:C:OP2	2.07	0.54
49:K8:58:ALA:O	49:K8:62:THR:HG22	2.07	0.54
49:K8:59:ARG:O	49:K8:62:THR:HG23	2.07	0.54
29:29:63:LEU:O	29:29:66:HIS:CG	2.60	0.54
1:1G:1369:C:OP2	9:82:112:LYS:N	2.40	0.54
26:14:2250:G:C6	37:45:82:ARG:HD2	2.42	0.54
1:1G:547:A:OP2	4:32:2:GLY:N	2.40	0.54
47:E5:18:ALA:HB3	47:E5:20:ARG:HE	1.73	0.54
37:88:30:GLY:HA2	37:88:107:ALA:HB2	1.90	0.54
32:59:41:MET:SD	32:59:41:MET:N	2.80	0.54
29:21:51:PHE:CD2	29:21:52:LEU:HG	2.42	0.54
26:14:1486:A:H2'	26:14:1487:G:C8	2.42	0.54
26:1H:1472:A:H2'	26:1H:1473:G:C8	2.42	0.54
26:14:2745:C:H4'	32:59:142:GLY:O	2.07	0.54
11:2I:46:GLY:HA2	11:2I:50:TYR:O	2.08	0.54
28:19:75:ILE:O	28:19:118:VAL:HG23	2.07	0.54
1:13:688:G:H2'	1:13:689:C:H6	1.72	0.54
19:AA:66:MET:HB3	19:AA:69:HIS:CD2	2.42	0.54
1:13:1510:U:H2'	1:13:1511:G:C8	2.43	0.54
4:3E:192:GLU:OE1	4:3E:192:GLU:N	2.31	0.54
2:1E:238:LEU:HD12	2:1E:238:LEU:H	1.71	0.54
4:32:15:GLU:OE1	4:32:66:ARG:NH1	2.41	0.54
37:45:110:THR:HG23	37:45:113:GLN:NE2	2.22	0.54
1:13:820:U:H4'	1:13:821:G:OP2	2.07	0.54
25:4K:23:A:H3'	25:4K:24:A:C8	2.43	0.54
26:1H:2592:G:C5	26:1H:2593:U:C4	2.95	0.54
26:1H:2250:G:C8	26:1H:2496:C:H5"	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:39:112:MET:O	30:39:115:ALA:HB3	2.07	0.54
1:13:1288:A:N3	1:13:1352:C:O2'	2.37	0.54
44:B5:5:TYR:HD1	49:G5:33:MET:SD	2.31	0.54
43:E8:82:LEU:HD13	43:E8:84:ARG:HH21	1.72	0.54
26:14:17:G:H2'	26:14:18:C:H6	1.72	0.54
26:14:172:C:H2'	26:14:173:G:H8	1.72	0.54
26:1H:1204:A:H2	26:1H:1241:A:N1	2.04	0.54
1:1G:114:U:O2'	1:1G:115:G:H5'	2.07	0.54
29:29:154:LYS:HA	29:29:154:LYS:HE3	1.88	0.54
26:1H:821:A:H2'	26:1H:946:G:H5''	1.88	0.54
40:B8:55:ASN:H	40:B8:59:THR:HB	1.71	0.54
19:AA:9:VAL:HG13	51:I5:63:TYR:CE1	2.41	0.54
26:1H:296:C:H2'	26:1H:297:C:H6	1.72	0.54
26:1H:1401:G:H2'	26:1H:1402:C:C6	2.42	0.54
20:BA:87:LYS:O	20:BA:91:LEU:HG	2.07	0.54
1:1G:1217:C:H5''	14:5A:9:LYS:HE2	1.90	0.54
26:14:19:C:H2'	26:14:20:C:C6	2.42	0.54
23:2L:22:A:N6	23:2L:47:7MG:H2'	2.23	0.54
12:3A:41:ARG:HD2	12:3A:42:THR:H	1.71	0.54
28:11:76:PRO:HB2	28:11:116:GLN:NE2	2.23	0.54
1:1G:1096:C:H2'	1:1G:1097:C:H6	1.71	0.54
26:1H:2505:G:H2'	26:1H:2576:G:O6	2.07	0.54
10:1I:24:VAL:HG12	10:1I:25:GLU:HG3	1.88	0.54
8:72:86:ILE:HD12	8:72:133:LEU:HD11	1.88	0.54
52:J5:31:VAL:HG13	52:J5:42:PRO:HG3	1.89	0.54
40:75:85:LYS:NZ	40:75:87:ASP:OD2	2.24	0.54
26:14:2446:G:C3'	26:14:2447:G:H5''	2.38	0.54
4:3E:209:ARG:HA	4:3E:209:ARG:HH11	1.73	0.54
43:E8:88:ARG:HB3	43:E8:92:ARG:HB3	1.88	0.54
26:14:568:U:O2'	58:14:4012:HOH:O	2.18	0.54
26:14:1021:A:H3'	26:14:1021:A:C8	2.43	0.54
26:14:2275:C:H6	26:14:2275:C:H5'	1.72	0.54
26:14:638:G:H2'	26:14:639:U:O4'	2.07	0.54
26:14:2287:A:N6	26:14:2344:U:H3	2.03	0.54
26:14:1636:C:H2'	26:14:1637:A:C8	2.42	0.54
26:1H:1265:A:OP1	26:1H:1265:A:H8	1.89	0.54
1:13:1412:C:H2'	1:13:1413:A:C8	2.43	0.54
1:13:1486:G:H2'	1:13:1487:G:O4'	2.07	0.54
38:55:72:ASP:OD2	38:55:75:LEU:HB2	2.08	0.54
44:F8:12:VAL:HG13	44:F8:27:THR:HG23	1.88	0.54
3:22:45:LYS:O	3:22:48:TYR:HB3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:F8:67:GLY:O	44:F8:69:TYR:N	2.41	0.54
1:1G:564:C:O2'	8:72:91:ARG:NH2	2.39	0.54
1:13:201:C:H42	1:13:216:G:H1	1.54	0.54
1:1G:298:A:H5''	1:1G:299:G:OP2	2.06	0.54
28:19:255:LYS:H	28:19:255:LYS:HE3	1.72	0.54
26:14:1856:G:H1	26:14:1886:C:N4	2.06	0.54
40:B8:81:PRO:HG2	40:B8:82:LEU:HD12	1.89	0.54
26:14:2542:A:N3	26:14:2542:A:H5''	2.23	0.54
1:1G:1342:C:H4'	9:82:125:TYR:HB3	1.90	0.54
30:39:153:SER:OG	30:39:190:GLU:HB2	2.07	0.54
26:1H:761:A:H5''	58:1H:3686:HOH:O	2.07	0.54
26:14:1113:U:H2'	26:14:1114:G:O4'	2.08	0.54
26:14:1141:U:H5''	26:14:1142(A):A:O4'	2.08	0.54
52:N8:40:LYS:NZ	52:N8:46:CYS:HB3	2.19	0.54
13:4A:2:ALA:HB1	51:15:34:GLU:HB3	1.89	0.54
43:E8:18:ARG:HD3	43:E8:76:VAL:HG13	1.90	0.54
26:14:152:G:H1	26:14:174:C:N4	2.04	0.54
26:14:1298:C:H5''	26:14:1299:G:OP2	2.08	0.54
1:13:407:G:H1	1:13:435:C:H42	1.55	0.54
1:1G:1272:G:C2	1:1G:1273:G:H1'	2.42	0.54
20:BA:49:ALA:HB2	20:BA:92:LEU:HD22	1.88	0.54
46:D5:7:ALA:O	46:D5:8:TYR:CG	2.60	0.54
26:14:2054:A:H5''	26:14:2055:C:O5'	2.06	0.54
1:13:1016:A:H2'	1:13:1017:G:O4'	2.08	0.54
26:1H:618:G:H2'	26:1H:618(A):C:H6	1.71	0.54
1:1G:1411:C:H2'	1:1G:1412:C:H6	1.71	0.54
31:41:110:ALA:HA	31:41:140:ILE:O	2.07	0.54
35:25:97:ARG:CZ	35:25:99:PHE:HE1	2.20	0.54
30:31:7:TYR:HA	30:31:22:ALA:O	2.08	0.54
46:H8:103:ARG:HG3	46:H8:136:PHE:CD2	2.43	0.54
1:13:1309:G:N2	1:13:1329:A:H1'	2.22	0.54
1:1G:1360:A:OP1	1:1G:1360:A:H8	1.91	0.54
36:35:8:PRO:HG2	36:35:13:ASN:HB3	1.90	0.54
26:14:910:A:N3	26:14:2264:C:O2'	2.39	0.54
29:21:181:LEU:HD21	40:B8:7:ILE:HG23	1.90	0.54
43:E8:17:VAL:HG13	43:E8:76:VAL:HG11	1.89	0.54
26:14:898:C:H3'	26:14:899:A:H5''	1.90	0.54
40:75:18:ASP:N	40:75:18:ASP:OD1	2.38	0.54
1:13:179:A:H2'	1:13:180:U:C6	2.43	0.54
27:1J:9:G:H5'	39:65:25:ARG:HH12	1.72	0.54
1:13:44:G:C2	1:13:45:U:H1'	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:78:130:PHE:CE1	36:78:146:VAL:HG23	2.43	0.54
1:1G:994:A:C2	14:5A:5:ALA:HB2	2.42	0.54
18:9I:50:ILE:HG12	18:9I:70:ILE:HD12	1.90	0.54
1:13:438:G:O2'	1:13:494:U:O4	2.21	0.54
1:13:444:C:H2'	1:13:445:G:H8	1.73	0.54
19:AI:6:LYS:O	19:AI:7:LYS:HB3	2.06	0.54
31:41:109:VAL:O	31:41:113:ARG:HG3	2.08	0.54
33:61:3:VAL:HG12	33:61:38:LEU:HA	1.89	0.54
1:1G:187:C:H2'	1:1G:188:U:O4'	2.08	0.54
1:13:544:G:H2'	1:13:545:C:H6	1.72	0.54
2:12:185:ILE:HG22	2:12:199:TYR:HB2	1.90	0.54
26:1H:589:C:H2'	26:1H:590:A:C8	2.43	0.54
55:Q8:46:ARG:NH2	55:Q8:47:LYS:O	2.41	0.54
55:Q8:49:VAL:HB	55:Q8:52:LYS:C	2.27	0.54
26:1H:1590:U:H2'	26:1H:1591:G:C8	2.42	0.54
26:1H:1593:G:H2'	26:1H:1594:G:C8	2.42	0.54
29:21:77:ILE:O	29:21:79:ARG:N	2.39	0.54
26:14:2304:G:N2	26:14:2312:U:O4	2.36	0.54
55:Q8:34:TRP:HB3	55:Q8:36:LYS:N	2.22	0.54
26:14:731:C:H5''	58:14:4110:HOH:O	2.07	0.54
55:M5:33:ASN:O	55:M5:36:LYS:HE3	2.08	0.54
34:58:133:GLN:OE1	34:58:133:GLN:N	2.37	0.54
1:1G:590:C:H2'	1:1G:591:U:H6	1.72	0.54
1:13:1318:A:H5''	19:AI:10:PHE:CD2	2.43	0.54
36:78:111:ARG:HG2	36:78:128:HIS:CG	2.43	0.54
46:D5:6:LYS:HE2	46:D5:43:GLU:HG3	1.89	0.54
29:21:167:VAL:HG13	29:21:189:PRO:HD3	1.90	0.54
24:3K:48:C:C5	24:3K:59:U:H1'	2.43	0.54
26:1H:270(K):C:O2'	26:1H:270(N):G:N2	2.39	0.54
26:14:566:U:H2'	26:14:567:A:O4'	2.08	0.54
26:1H:1514:U:H2'	26:1H:1515:C:C6	2.42	0.54
4:32:13:ARG:C	4:32:15:GLU:H	2.11	0.54
1:13:924:C:H2'	1:13:925:G:C8	2.43	0.54
33:61:99:GLU:HG2	33:61:103:ARG:HH21	1.73	0.54
1:1G:198:G:H2'	1:1G:199:G:H8	1.73	0.54
26:14:303:U:H2'	26:14:304:G:C8	2.43	0.54
39:A8:67:ARG:HB2	39:A8:67:ARG:NH1	2.23	0.54
1:1G:1027:C:O2'	1:1G:1035:A:N6	2.41	0.54
29:29:11:MET:HA	29:29:24:THR:HA	1.88	0.54
9:8E:89:ASN:O	9:8E:91:ASP:N	2.39	0.54
1:13:186(E):C:H42	1:13:191(B):G:H1	1.55	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:7E:39:LEU:HB3	8:7E:45:ILE:HD11	1.90	0.54
1:13:630:G:H2'	1:13:631:G:O4'	2.07	0.54
26:1H:729:G:C8	28:11:208:LYS:HD3	2.43	0.54
34:58:4:TYR:CE2	41:C8:100:VAL:HG11	2.42	0.54
1:1G:1177:G:O2'	1:1G:1178:G:O4'	2.26	0.54
45:C5:17:SER:HB2	45:C5:71:LYS:HD2	1.90	0.54
38:55:98:LEU:HB2	38:55:113:LEU:CD2	2.38	0.54
26:1H:2808:U:H5'	26:1H:2891:G:O6	2.08	0.54
10:1I:46:ARG:HB2	10:1I:46:ARG:HH11	1.73	0.54
15:6A:43:LEU:HD11	15:6A:53:HIS:HA	1.89	0.54
26:1H:1047:G:O2'	26:1H:1111:A:N6	2.40	0.54
23:2L:47:7MG:H3'	23:2L:48:U:H6	1.73	0.54
32:51:23:ARG:HH12	32:51:25:LYS:HD2	1.72	0.54
20:BA:97:ALA:O	20:BA:99:LEU:HD12	2.08	0.54
27:1J:31:C:H2'	27:1J:32:C:H5'	1.90	0.54
30:39:141:ALA:O	30:39:144:LYS:HB3	2.07	0.54
26:14:2165:G:H2'	26:14:2166:G:H5'	1.89	0.54
5:42:41:VAL:O	5:42:67:VAL:N	2.41	0.54
26:1H:988:A:P	50:L8:11:SER:HB2	2.48	0.54
1:13:1342:C:O2'	9:8E:124:GLN:HG2	2.07	0.54
47:E5:72:ARG:HB3	47:E5:75:LEU:HB2	1.90	0.54
28:19:200:ASP:OD1	28:19:203:ASN:ND2	2.41	0.54
1:1G:11:G:C5	1:1G:12:U:C5	2.96	0.54
1:1G:861:G:C4	1:1G:862:C:H5	2.27	0.53
36:35:30:THR:HG21	36:35:35:HIS:N	2.22	0.53
44:F8:3:THR:CB	44:F8:6:ASP:HB2	2.38	0.53
20:BA:85:MET:HB2	20:BA:104:LEU:HD21	1.90	0.53
26:14:2567:G:H2'	26:14:2568:C:H6	1.72	0.53
4:32:107:ARG:HE	4:32:173:TRP:HZ2	1.56	0.53
26:14:1654:A:H1'	26:14:2823:A:H5'	1.90	0.53
26:1H:2105:C:H2'	26:1H:2106:G:H8	1.72	0.53
5:42:101:ILE:HD13	5:42:101:ILE:H	1.73	0.53
3:22:36:ASP:HA	3:22:39:ILE:HD12	1.89	0.53
26:14:1167:U:O2	26:14:1183:G:N2	2.41	0.53
26:14:2103:C:H2'	26:14:2104:G:H8	1.73	0.53
28:19:253:GLN:HB3	28:19:255:LYS:NZ	2.23	0.53
26:1H:2475:C:H4'	26:1H:2476:A:OP1	2.08	0.53
26:14:1007:C:OP1	34:15:37:LYS:NZ	2.40	0.53
26:14:2094:G:H5'	33:69:25:TYR:CD1	2.43	0.53
1:1G:382:A:H2'	1:1G:383:A:C8	2.43	0.53
40:B8:29:ARG:NH1	40:B8:46:GLU:OE2	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:22:70:VAL:HG11	3:22:76:VAL:HG11	1.90	0.53
19:AI:53:ASN:O	19:AI:77:THR:HG22	2.07	0.53
26:14:1347:G:H5''	26:14:1348:G:OP2	2.08	0.53
1:13:237:C:H5''	17:8I:25:ARG:CZ	2.37	0.53
47:I8:44:ARG:HB2	47:I8:44:ARG:HH11	1.73	0.53
6:5E:97:PHE:HD2	18:9I:31:LEU:HD21	1.71	0.53
33:69:63:ALA:HA	33:69:66:GLU:HG2	1.91	0.53
26:14:2001:A:H2'	26:14:2002:G:C8	2.43	0.53
26:1H:1642:G:N7	58:1H:3947:HOH:O	2.34	0.53
51:I5:12:ALA:HB1	51:I5:29:PRO:HA	1.90	0.53
2:1E:8:LYS:HE2	2:1E:8:LYS:H	1.73	0.53
35:25:15:GLY:O	35:25:47:ILE:HB	2.08	0.53
26:14:1229:G:H2'	26:14:1229(A):G:H5''	1.90	0.53
42:D8:59:ALA:CB	42:D8:96:ILE:HD13	2.38	0.53
26:14:218:A:H2	26:14:235:U:H4'	1.72	0.53
28:11:46:GLN:HB2	28:11:48:ARG:HG2	1.90	0.53
32:59:9:ILE:HG22	32:59:51:ARG:HA	1.89	0.53
9:8E:25:LYS:NZ	9:8E:26:VAL:O	2.39	0.53
26:14:678:C:H2'	26:14:679:C:C6	2.43	0.53
26:1H:381:G:C4	26:1H:394:A:C2	2.96	0.53
53:K5:51:GLU:HG2	53:K5:52:VAL:H	1.73	0.53
36:78:2:LYS:HE3	36:78:4:SER:HB3	1.91	0.53
31:41:166:ASP:N	31:41:166:ASP:OD1	2.41	0.53
26:1H:2784:C:O2'	29:21:37:ARG:NH1	2.40	0.53
1:13:554:C:H2'	1:13:555:C:H6	1.73	0.53
44:B5:40:LYS:HA	44:B5:51:VAL:HG11	1.90	0.53
17:8I:43:LEU:HD12	17:8I:68:ARG:HG2	1.90	0.53
26:1H:2408:U:H2'	26:1H:2409:G:C8	2.43	0.53
55:M5:22:VAL:HG12	55:M5:50:LEU:HB2	1.90	0.53
1:1G:1131:G:H2'	1:1G:1132:C:C6	2.43	0.53
26:14:1774:C:P	58:14:4078:HOH:O	2.65	0.53
1:13:412:A:H4'	1:13:413:G:O5'	2.08	0.53
26:14:1056:G:H1'	26:14:1103:A:H61	1.73	0.53
8:7E:87:SER:CB	8:7E:93:VAL:H	2.22	0.53
53:O8:16:CYS:O	53:O8:17:LYS:HB2	2.08	0.53
26:14:2844:G:H3'	26:14:2845:G:H8	1.72	0.53
26:1H:2747:G:OP1	32:51:138:LYS:NZ	2.40	0.53
1:1G:428:G:C5	1:1G:430:A:C6	2.96	0.53
13:4I:65:LYS:NZ	51:M8:52:THR:HB	2.22	0.53
26:1H:1401:G:H2'	26:1H:1402:C:H6	1.73	0.53
26:14:588:U:O4	26:14:670:A:H1'	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:59:17:VAL:HG23	32:59:26:VAL:HA	1.89	0.53
26:14:288:C:O2	26:14:353:G:N2	2.35	0.53
26:1H:2022:U:O2'	26:1H:2617:C:H5'	2.09	0.53
37:45:110:THR:OG1	37:45:112:GLU:OE1	2.23	0.53
2:12:163:PHE:HD1	2:12:185:ILE:HG13	1.73	0.53
31:49:55:LYS:O	31:49:58:GLN:HG3	2.07	0.53
3:22:112:SER:O	3:22:116:VAL:HG23	2.08	0.53
1:13:863:U:O2'	1:13:865:A:N7	2.34	0.53
44:F8:80:ILE:HG13	44:F8:80:ILE:O	2.08	0.53
3:2E:175:LEU:HD21	3:2E:201:TYR:CE2	2.44	0.53
54:P8:24:THR:HG23	54:P8:27:GLY:H	1.73	0.53
26:1H:606:U:H4'	26:1H:658:C:H4'	1.89	0.53
11:2I:48:ILE:HG12	11:2I:63:LEU:HB2	1.90	0.53
42:D8:46:VAL:HG12	42:D8:47:VAL:H	1.73	0.53
26:1H:2335:A:C8	26:1H:2337:G:C5	2.96	0.53
26:14:1542:G:O6	26:14:1543:A:N6	2.42	0.53
26:14:2378:A:H4'	39:65:23:ARG:HH11	1.73	0.53
26:1H:2611:U:H6	26:1H:2611:U:H5'	1.72	0.53
27:1J:13:A:H2'	27:1J:70:C:O2'	2.09	0.53
26:1H:631:A:OP1	36:78:65:ARG:NH2	2.41	0.53
7:62:27:ILE:HA	7:62:30:ILE:HD12	1.90	0.53
26:1H:699:A:H2'	26:1H:700:G:O4'	2.08	0.53
19:AA:28:LYS:HG2	19:AA:29:ARG:H	1.71	0.53
36:35:62:LEU:HG	55:M5:25:MET:HB2	1.90	0.53
1:1G:1062:U:H2'	1:1G:1063:C:C6	2.43	0.53
33:69:109:ILE:HB	33:69:130:TYR:CZ	2.43	0.53
28:11:223:GLY:HA3	28:11:231:HIS:CE1	2.44	0.53
26:1H:2347:C:P	55:Q8:36:LYS:HE3	2.48	0.53
38:98:63:ARG:NH2	38:98:80:PHE:HD2	2.07	0.53
26:1H:1796:U:H2'	26:1H:1797:C:C6	2.43	0.53
7:6E:92:SER:O	7:6E:96:GLN:HG3	2.08	0.53
37:88:51:ARG:NH1	37:88:52:VAL:HG23	2.23	0.53
1:13:491:G:H2'	1:13:492:G:O4'	2.08	0.53
6:5E:36:ARG:NH2	6:5E:38:GLU:HG2	2.24	0.53
29:29:105:THR:OG1	29:29:199:ARG:NH2	2.41	0.53
22:1L:10:G:H2'	22:1L:11:C:C6	2.44	0.53
26:14:1665:A:OP2	58:14:3506:HOH:O	2.18	0.53
16:7I:34:GLU:HG2	16:7I:35:LYS:H	1.72	0.53
40:75:99:LEU:HD22	40:75:101:PHE:CE1	2.43	0.53
20:BA:72:LEU:O	20:BA:73:HIS:HB2	2.07	0.53
39:65:67:ARG:NH1	39:65:67:ARG:HB2	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1935:G:H1'	26:14:1964:G:N2	2.23	0.53
26:14:842:G:H8	26:14:842:G:O5'	1.91	0.53
11:2I:44:SER:OG	11:2I:47:VAL:HG23	2.07	0.53
3:2E:18:TRP:H	3:2E:18:TRP:HE3	1.56	0.53
26:14:25:U:H5'	43:A5:79:GLY:HA2	1.90	0.53
26:1H:1339:G:H21	26:1H:1603:A:H1'	1.74	0.53
1:13:280:C:C2	17:8I:38:ARG:HG3	2.44	0.53
26:1H:2461:C:H2'	26:1H:2462:U:C6	2.43	0.53
33:69:77:LEU:HD13	33:69:141:LYS:HD2	1.91	0.53
8:7E:12:ARG:NH1	8:7E:27:PRO:HD2	2.24	0.53
43:E8:38:TYR:HE2	52:N8:41:PRO:HD3	1.73	0.53
26:1H:1138:G:H21	34:58:106:MET:CE	2.18	0.53
17:8I:14:LYS:HD2	17:8I:14:LYS:H	1.72	0.53
26:1H:2579:C:H2'	26:1H:2580:U:O4'	2.08	0.53
24:3L:46:G:H2'	24:3L:47:U:H4'	1.91	0.53
26:14:780:G:OP1	28:19:218:ARG:NH2	2.41	0.53
50:L8:9:VAL:HG22	50:L8:54:VAL:HA	1.91	0.53
30:31:32:LEU:CD2	30:31:105:VAL:HG13	2.39	0.53
22:1L:54:U:H5'	37:45:51:ARG:HH22	1.73	0.53
14:5A:25:VAL:O	14:5A:26:ARG:HB3	2.09	0.53
26:1H:1337:G:H2'	26:1H:1338:G:C8	2.43	0.53
36:35:80:TYR:CD2	36:35:111:ARG:HB3	2.43	0.53
11:2I:21:ILE:HD13	11:2I:84:VAL:HG12	1.90	0.53
2:12:74:LYS:HE3	2:12:166:ASP:HB2	1.89	0.53
54:P8:16:HIS:HB2	54:P8:44:PRO:HG2	1.91	0.53
1:13:486:U:H2'	1:13:487:A:C8	2.44	0.53
27:16:87:G:N2	27:16:89(A):A:OP2	2.40	0.53
26:1H:2291:U:O2'	26:1H:2374:C:O2	2.25	0.53
1:1G:1275:A:H2'	1:1G:1276:G:O4'	2.09	0.53
1:13:448:A:P	1:13:485:G:H22	2.30	0.53
3:2E:79:ARG:NH1	18:9A:87:ARG:HH12	2.06	0.53
26:14:826:U:H2'	26:14:828:U:O4'	2.08	0.53
11:2A:14:VAL:HG11	11:2A:35:PRO:HD3	1.90	0.53
26:14:1776:G:OP2	58:14:3427:HOH:O	2.19	0.53
2:1E:94:ASN:OD1	2:1E:95:GLN:N	2.40	0.53
26:1H:1766:U:H2'	26:1H:1767:C:H6	1.73	0.53
7:62:15:ASP:OD1	7:62:44:TYR:OH	2.25	0.53
1:13:1504:G:OP1	1:13:1507:A:H4'	2.09	0.53
26:14:2702:U:O2'	26:14:2703:C:C5	2.60	0.53
29:21:105:THR:HG22	29:21:106:GLY:N	2.21	0.53
27:1J:16:G:H2'	27:1J:17:C:C6	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:K8:47:ASN:O	49:K8:49:LYS:HG3	2.08	0.53
12:3I:117:ARG:HB3	12:3I:122:THR:HB	1.89	0.53
26:1H:2347:C:P	53:O8:39:TYR:HH	2.27	0.53
1:1G:589:C:N3	1:1G:650:G:N2	2.54	0.53
30:39:89:VAL:HG12	30:39:90:PHE:H	1.73	0.53
9:82:103:THR:HG22	9:82:105:ASP:H	1.73	0.53
29:29:89:ASP:C	29:29:91:VAL:H	2.11	0.53
41:C8:6:THR:N	58:C8:302:HOH:O	2.40	0.53
7:62:65:ALA:HB1	7:62:127:ALA:HB3	1.91	0.53
30:39:132:VAL:O	30:39:134:GLY:N	2.42	0.53
1:1G:1226:C:H4'	19:AA:80:TYR:OH	2.07	0.53
1:1G:617:G:N7	58:1G:1839:HOH:O	2.33	0.53
26:14:31:C:OP1	58:14:3732:HOH:O	2.19	0.53
7:62:15:ASP:HB3	7:62:19:GLY:H	1.73	0.53
28:11:68:LYS:HB3	28:11:70:TRP:CH2	2.43	0.53
36:78:39:LYS:HG3	36:78:45:LEU:HD23	1.91	0.53
33:61:67:ARG:O	33:61:71:ILE:HG22	2.08	0.53
1:1G:821:G:H5''	58:1G:1865:HOH:O	2.09	0.53
26:14:576:U:H5	58:14:3567:HOH:O	1.91	0.53
34:15:67:LEU:O	34:15:88:GLU:HG3	2.08	0.53
10:1I:42:THR:HG23	10:1I:67:THR:O	2.08	0.53
26:14:2712:U:H2'	26:14:2714:G:H5''	1.90	0.53
34:58:38:HIS:O	41:C8:67:ALA:HB1	2.08	0.53
29:29:56:PRO:HD2	29:29:58:ARG:HD3	1.90	0.53
43:A5:50:VAL:HG22	43:A5:105:VAL:HG23	1.89	0.53
55:Q8:5:LYS:H	55:Q8:59:LYS:NZ	2.06	0.53
19:AA:35:SER:O	19:AA:71:LEU:HD12	2.09	0.53
26:14:1019:U:H3	26:14:1142(A):A:N6	1.99	0.53
1:13:1160:G:H22	1:13:1177:G:N2	2.07	0.53
8:7E:41:ARG:NH1	8:7E:123:GLU:OE1	2.42	0.53
29:29:51:PHE:O	29:29:52:LEU:HB2	2.07	0.53
9:82:112:LYS:HE3	9:82:118:LYS:H	1.72	0.53
29:21:119:ARG:HD2	29:21:120:TRP:CE2	2.44	0.53
1:1G:157:G:H1	1:1G:164:U:H3	1.57	0.53
1:13:1349:A:H2'	1:13:1350:A:C8	2.44	0.53
26:14:2689:U:OP2	26:14:2719:G:N2	2.42	0.53
26:14:1638:C:H1'	26:14:2698:U:O2'	2.09	0.53
28:11:108:PRO:HG3	28:11:143:HIS:HE1	1.72	0.53
50:L8:7:LYS:O	50:L8:54:VAL:HG23	2.08	0.53
20:BI:97:ALA:O	20:BI:99:LEU:N	2.42	0.53
1:1G:778:G:O2'	11:2A:119:CYS:HB2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:765:G:H5''	1:13:766:A:OP1	2.08	0.53
26:14:353:G:H2'	26:14:354:G:C8	2.43	0.53
26:1H:85:G:OP2	45:G8:9:LYS:HB2	2.08	0.53
8:7E:45:ILE:HB	8:7E:47:GLY:H	1.73	0.53
26:14:194:G:H2'	26:14:195:A:O4'	2.09	0.53
21:1B:9:ARG:NE	21:1B:10:ARG:HG2	2.23	0.53
27:16:2:C:H2'	27:16:3:C:C6	2.43	0.53
26:1H:396:G:O4'	48:J8:13:ILE:HD11	2.09	0.53
24:3K:28:G:H2'	24:3K:29:G:H8	1.74	0.53
46:H8:117:LEU:H	46:H8:117:LEU:HD13	1.74	0.53
44:B5:35:THR:HB	44:B5:38:GLU:H	1.73	0.53
52:J5:16:ARG:HG3	52:J5:17:ASP:N	2.24	0.53
26:14:2299:G:H2'	26:14:2300:G:C8	2.43	0.53
26:14:2502:G:H8	58:14:3687:HOH:O	1.92	0.53
53:O8:26:ASN:ND2	53:O8:35:GLU:OE2	2.42	0.53
26:1H:1417:C:H2'	26:1H:1418:G:O4'	2.08	0.53
27:16:73:A:C4	27:16:104:A:C2	2.97	0.53
4:3E:112:VAL:HG12	4:3E:116:GLN:OE1	2.08	0.53
26:1H:654(G):C:O2	26:1H:654(N):G:N2	2.41	0.53
18:9A:53:ARG:HA	18:9A:56:THR:OG1	2.09	0.53
49:G5:59:ARG:O	49:G5:63:VAL:HG23	2.09	0.53
26:14:1358:G:O2'	26:14:1359:A:H5'	2.08	0.53
26:1H:1653:G:H3'	38:98:2:ARG:CG	2.39	0.53
26:14:688:U:H2'	26:14:689:A:H5'	1.91	0.53
1:13:652:U:O2'	1:13:653:A:O5'	2.26	0.53
15:6I:74:ASP:HB3	15:6I:77:ARG:HB3	1.91	0.53
45:C5:39:VAL:HG23	45:C5:41:GLY:H	1.73	0.53
1:13:1418:A:C2	1:13:1483:A:C2	2.97	0.53
39:65:64:GLU:O	39:65:68:GLN:HG3	2.09	0.53
28:19:166:GLN:HB3	28:19:174:ILE:HG22	1.91	0.53
26:14:251:A:C5	26:14:252:G:H1'	2.44	0.53
26:1H:389:G:H22	36:78:72:PRO:HD3	1.73	0.53
1:1G:688:G:H2'	1:1G:689:C:H6	1.74	0.53
6:5E:44:GLY:HA2	6:5E:59:TYR:CE2	2.44	0.53
48:F5:8:SER:HB3	48:F5:66:HIS:CD2	2.44	0.53
26:1H:1013:C:O2'	26:1H:1014:U:H5'	2.08	0.53
31:49:172:LEU:O	31:49:176:LEU:HB2	2.08	0.53
41:C8:95:LEU:HG	42:D8:4:ILE:HD13	1.91	0.53
26:1H:963:U:H5''	58:1H:3858:HOH:O	2.09	0.53
26:1H:415:A:H2'	26:1H:416:C:O4'	2.08	0.53
7:62:91:VAL:HG12	7:62:95:ARG:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:9I:58:LEU:HD23	18:9I:62:GLU:HB3	1.91	0.53
55:M5:40:GLU:N	55:M5:43:GLN:HG3	2.22	0.53
11:2A:86:GLY:N	11:2A:112:THR:OG1	2.38	0.53
24:3K:24:G:H2'	24:3K:25:C:C6	2.44	0.53
26:1H:1500:G:N2	28:11:99:ASP:O	2.36	0.53
13:4I:49:THR:O	13:4I:52:GLU:HG3	2.09	0.53
26:14:2371:G:O2'	53:K5:46:HIS:ND1	2.32	0.53
26:1H:528:A:C2	26:1H:2043:C:H4'	2.44	0.53
1:13:452:A:H62	1:13:480:U:H3	1.56	0.53
26:1H:2617:C:H2'	26:1H:2618:G:H8	1.73	0.53
26:1H:1441:G:H2'	26:1H:1442:G:C8	2.43	0.53
26:1H:2461:C:H2'	26:1H:2462:U:H6	1.74	0.53
1:13:719:C:H1'	18:9I:49:LYS:HB3	1.91	0.53
30:31:52:LYS:O	30:31:88:VAL:HG23	2.09	0.53
1:13:1492:A:H1'	25:4K:20:U:O2'	2.08	0.53
29:29:96:PHE:O	29:29:175:VAL:HG11	2.08	0.53
16:7A:19:ILE:HB	16:7A:36:ILE:O	2.08	0.53
26:14:2390:U:O2'	26:14:2391:G:H5'	2.09	0.53
26:14:1278:A:OP1	38:55:36:THR:HG22	2.09	0.53
1:1G:407:G:OP1	4:32:115:ARG:NE	2.41	0.53
55:Q8:57:ARG:HA	55:Q8:58:ILE:C	2.29	0.53
26:14:2016:U:OP1	58:14:3792:HOH:O	2.18	0.53
26:1H:512:G:C8	58:1H:4665:HOH:O	2.53	0.53
36:35:39:LYS:HB2	36:35:45:LEU:HD21	1.90	0.53
26:14:1678:G:N2	26:14:1989:G:H22	2.01	0.53
55:Q8:38:GLY:HA2	55:Q8:39:LYS:O	2.09	0.53
26:1H:2693:A:H2'	26:1H:2694:G:C8	2.40	0.53
1:1G:513:C:N3	1:1G:538:G:N2	2.50	0.53
26:14:1169:G:N2	26:14:1180:C:O2	2.41	0.53
2:12:193:ASP:OD1	2:12:196:LEU:HG	2.09	0.53
28:11:182:LEU:N	28:11:272:ALA:HB3	2.23	0.53
26:1H:821:A:H5''	26:1H:822:U:H6	1.74	0.53
26:1H:527:C:OP2	26:1H:2779:U:H5	1.91	0.53
26:1H:302:C:H2'	26:1H:303:U:C6	2.44	0.53
1:1G:142:G:H2'	1:1G:143:A:C8	2.44	0.53
26:14:1449(A):G:H2'	26:14:1450:C:H6	1.73	0.53
1:13:7:G:H5'	1:13:298:A:O4'	2.09	0.53
41:C8:14:HIS:O	41:C8:18:LEU:HD12	2.09	0.53
26:1H:270(V):G:H2'	26:1H:270(W):G:H8	1.74	0.53
45:C5:49:VAL:O	45:C5:51:VAL:HG12	2.09	0.53
46:H8:45:ASP:OD2	46:H8:49:ARG:NH1	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:B8:99:LEU:HB3	40:B8:101:PHE:CE1	2.44	0.53
26:14:1999:C:H4'	26:14:2723:C:O2	2.08	0.53
5:42:5:ASP:HA	5:42:63:ARG:HH12	1.73	0.53
1:13:984:C:H2'	1:13:985:C:C6	2.44	0.53
13:4I:11:ARG:HG2	13:4I:12:ASN:H	1.73	0.53
1:13:615:C:C2	1:13:616:G:C8	2.97	0.53
1:13:616:G:C2	1:13:617:G:N7	2.77	0.53
36:78:113:LYS:HG2	36:78:115:LEU:HD23	1.89	0.53
12:3I:66:VAL:HG21	12:3I:98:TYR:CE1	2.43	0.53
36:78:19:VAL:CB	36:78:27:HIS:HB2	2.37	0.52
42:95:7:THR:HG23	42:95:22:VAL:HG21	1.89	0.52
26:1H:1784:A:OP1	58:1H:3967:HOH:O	2.19	0.52
1:13:93:U:H2'	1:13:95:G:O4'	2.09	0.52
1:1G:179:A:H2'	1:1G:180:U:C6	2.43	0.52
1:13:427:U:H3'	1:13:428:G:H2'	1.90	0.52
26:1H:999:U:P	58:1H:3999:HOH:O	2.67	0.52
1:13:558:G:H5''	1:13:559:A:OP2	2.08	0.52
26:1H:2155:G:H2'	26:1H:2156:G:H5'	1.90	0.52
26:1H:569:U:C4	26:1H:570:G:C6	2.97	0.52
26:1H:2807:G:H3'	26:1H:2808:U:H5''	1.92	0.52
26:1H:661:C:HO2'	36:78:14:LYS:H	1.55	0.52
26:1H:1839:G:C8	26:1H:1839:G:H5''	2.43	0.52
1:1G:1226:C:N4	13:4A:104:ARG:HD2	2.24	0.52
26:1H:1093:G:H1'	26:1H:1099:G:H22	1.74	0.52
26:1H:1290:C:H2'	26:1H:1291:C:H6	1.73	0.52
26:1H:1062:G:H2'	26:1H:1063:G:C8	2.44	0.52
26:1H:988:A:O5'	50:L8:11:SER:HB2	2.09	0.52
26:1H:394:A:C6	26:1H:395:U:N3	2.76	0.52
31:49:118:ARG:HB3	31:49:181:ARG:HG3	1.91	0.52
30:39:140:LEU:HD21	30:39:170:LEU:HD21	1.90	0.52
29:21:32:PRO:HD2	29:21:50:GLY:O	2.09	0.52
1:1G:1490:C:H2'	1:1G:1491:G:O4'	2.09	0.52
26:14:275:G:H2'	26:14:276:A:C8	2.44	0.52
10:1I:54:PHE:CD2	10:1I:55:LYS:HG2	2.43	0.52
20:BI:30:LYS:HE2	20:BI:72:LEU:HD22	1.92	0.52
5:4E:100:VAL:O	5:4E:107:ARG:NH2	2.42	0.52
37:88:135:ASP:HB3	37:88:137:TYR:N	2.23	0.52
26:14:71:A:H3'	26:14:71:A:OP2	2.08	0.52
27:1J:9:G:P	39:65:25:ARG:HH22	2.32	0.52
26:1H:2138:C:N4	26:1H:2153:G:H1	2.07	0.52
26:1H:507:A:H5''	26:1H:508:G:H3'	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:3L:35:A:H2	25:4L:14:A:C6	2.27	0.52
1:13:4:U:O4	8:7E:105:ARG:HD2	2.10	0.52
26:1H:2724:C:OP1	29:21:118:LYS:HE3	2.09	0.52
44:B5:67:GLY:O	44:B5:69:TYR:N	2.40	0.52
26:14:748:G:C8	43:A5:89:ALA:HB1	2.45	0.52
2:1E:111:ARG:HG2	2:1E:111:ARG:HH11	1.72	0.52
34:15:62:VAL:HG22	34:15:66:LYS:HD2	1.91	0.52
26:14:696:G:H2'	26:14:697:C:H6	1.75	0.52
31:49:95:ARG:HG2	31:49:96:ARG:HG2	1.91	0.52
6:52:19:LEU:HD11	6:52:59:TYR:CD2	2.44	0.52
30:39:164:ARG:O	30:39:167:ALA:HB3	2.08	0.52
48:F5:52:ARG:HH11	48:F5:57:GLU:HB2	1.74	0.52
26:14:2448:A:N1	58:14:3682:HOH:O	2.34	0.52
26:14:2315:G:H2'	26:14:2316:C:C6	2.45	0.52
5:4E:8:GLU:HG2	5:4E:34:VAL:HG22	1.91	0.52
26:1H:518:G:H5'	43:E8:18:ARG:NH2	2.25	0.52
29:29:179:GLU:HB3	29:29:181:LEU:HD22	1.90	0.52
55:Q8:39:LYS:O	55:Q8:40:GLU:HB3	2.09	0.52
41:C8:92:ARG:NE	42:D8:11:GLN:H	2.06	0.52
26:14:2572:A:H62	29:29:145:LYS:HD2	1.74	0.52
26:1H:280:C:N3	26:1H:361:G:C2	2.77	0.52
26:1H:300:A:H2'	26:1H:334:C:H1'	1.92	0.52
26:1H:654(A):A:H2	26:1H:654(T):A:N1	2.07	0.52
41:85:98:LEU:HA	41:85:100:VAL:O	2.08	0.52
26:1H:1649:G:O2'	38:98:107:ASP:OD1	2.13	0.52
18:9A:22:VAL:C	18:9A:24:ALA:H	2.13	0.52
14:5I:6:LEU:HB3	14:5I:23:ARG:NH2	2.23	0.52
5:42:145:LYS:O	5:42:149:GLU:HG2	2.09	0.52
3:22:39:ILE:O	3:22:43:LEU:HB2	2.10	0.52
23:2L:46:G:H4'	23:2L:47:7MG:OP1	2.09	0.52
13:4I:106:ASN:O	13:4I:106:ASN:ND2	2.35	0.52
26:14:2440:C:H5'	58:14:3470:HOH:O	2.09	0.52
1:1G:1096:C:H2'	1:1G:1097:C:C6	2.44	0.52
46:D5:15:PRO:HB2	46:D5:19:ARG:NH2	2.24	0.52
1:1G:103:C:OP2	20:BA:14:LYS:HE2	2.08	0.52
37:45:39:PRO:HD3	37:45:99:PRO:HD3	1.90	0.52
28:11:121:PRO:HB3	28:11:135:PHE:CE2	2.44	0.52
26:1H:35:G:H2'	26:1H:36:G:O4'	2.08	0.52
5:42:12:LEU:O	5:42:30:ALA:HA	2.09	0.52
26:1H:857:C:H4'	47:I8:23:VAL:HG21	1.90	0.52
47:E5:37:LEU:HG	47:E5:60:PHE:HA	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:8I:29:HIS:N	17:8I:34:LYS:O	2.39	0.52
48:J8:65:SER:OG	48:J8:66:HIS:ND1	2.31	0.52
32:59:121:ILE:HG23	32:59:133:VAL:HG13	1.91	0.52
2:12:50:GLU:HB3	2:12:200:ILE:O	2.10	0.52
2:1E:100:GLY:N	2:1E:176:GLU:OE2	2.42	0.52
26:1H:621:A:O4'	26:1H:621:A:N3	2.41	0.52
13:4A:81:LEU:HB3	13:4A:88:ARG:HG3	1.92	0.52
26:14:1327:C:O3'	38:55:105:ARG:NH2	2.43	0.52
9:82:16:ARG:O	9:82:63:ILE:HG23	2.10	0.52
26:1H:4:C:H2'	26:1H:5:A:C8	2.44	0.52
1:1G:1328:C:OP1	21:1B:21:TYR:OH	2.24	0.52
26:14:2131:G:C5'	26:14:2158:A:H61	2.22	0.52
1:1G:518:C:H5''	1:1G:519:C:C6	2.45	0.52
1:1G:1080:A:H5'	5:42:14:ARG:NH2	2.24	0.52
23:2L:54:G:H3'	23:2L:55:5MU:H71	1.90	0.52
10:1I:40:LEU:HB2	10:1I:69:ASN:CB	2.39	0.52
5:4E:68:GLU:O	5:4E:70:PRO:HD3	2.10	0.52
1:1G:763:G:H2'	1:1G:764:C:C6	2.44	0.52
26:1H:1566:A:O2'	26:1H:1567:A:H5'	2.09	0.52
26:1H:1403:C:H5'	26:1H:1471:A:H1'	1.92	0.52
1:1G:854:G:O6	58:1G:1881:HOH:O	2.19	0.52
19:AI:51:VAL:O	19:AI:57:HIS:HA	2.09	0.52
26:14:1536:A:H3'	26:14:1537:C:O4'	2.09	0.52
26:14:1538:G:H2'	26:14:1539:G:H8	1.73	0.52
26:1H:2150:U:H2'	26:1H:2151:G:C8	2.45	0.52
37:45:39:PRO:HA	37:45:97:VAL:O	2.09	0.52
43:A5:69:LEU:HD22	43:A5:107:LEU:HB3	1.92	0.52
28:11:69:ARG:HH21	28:11:130:ALA:H	1.57	0.52
1:13:757:U:OP1	1:13:822:C:O2'	2.27	0.52
9:8E:79:LEU:HD22	9:8E:83:ARG:HG3	1.91	0.52
33:61:12:LEU:HG	33:61:19:VAL:HG21	1.92	0.52
1:13:958:A:O4'	19:AI:55:LYS:NZ	2.43	0.52
1:1G:56:U:H2'	1:1G:57:G:C8	2.45	0.52
15:6I:27:VAL:O	15:6I:31:LEU:HB2	2.08	0.52
2:1E:97:TRP:CZ3	2:1E:172:ILE:HB	2.45	0.52
26:14:2023:G:OP2	26:14:2617:C:H4'	2.08	0.52
26:14:1963:U:H5''	26:14:1963:U:O2	2.10	0.52
47:I8:60:PHE:CD1	47:I8:60:PHE:N	2.77	0.52
1:13:971:G:N2	1:13:1363:A:OP2	2.34	0.52
1:1G:804:U:H5''	1:1G:805:C:OP2	2.08	0.52
1:13:1125:U:H5'	1:13:1126:U:C5	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1187:G:P	58:1H:3882:HOH:O	2.66	0.52
26:14:2638:G:OP2	29:29:82:ARG:NH2	2.42	0.52
26:1H:1510:A:H4'	26:1H:1511:A:N7	2.25	0.52
24:3K:70:G:H2'	24:3K:71:G:O4'	2.10	0.52
8:72:12:ARG:NH2	8:72:27:PRO:HD3	2.23	0.52
26:14:2506:U:H4'	26:14:2507:C:OP1	2.07	0.52
9:82:117:HIS:O	9:82:118:LYS:HB2	2.07	0.52
26:1H:2785:C:H2'	26:1H:2786:U:O4'	2.10	0.52
1:13:693:G:H2'	1:13:694:A:C8	2.45	0.52
1:1G:195:A:C6	1:1G:196:A:N1	2.78	0.52
26:1H:1899:G:H22	26:1H:1902:C:H41	1.58	0.52
26:14:138:G:N2	44:B5:44:GLU:OE2	2.43	0.52
13:4A:29:ARG:HB3	13:4A:64:TRP:CZ2	2.45	0.52
13:4A:60:VAL:HG13	13:4A:64:TRP:NE1	2.24	0.52
1:13:10:A:OP2	5:4E:126:ARG:HD3	2.10	0.52
39:A8:35:ILE:HD11	39:A8:101:LEU:HD23	1.90	0.52
4:3E:100:ARG:NH1	4:3E:137:SER:HA	2.25	0.52
40:75:55:ASN:H	40:75:59:THR:HG22	1.74	0.52
43:E8:60:ASN:N	43:E8:60:ASN:ND2	2.57	0.52
1:13:1063:C:H3'	1:13:1064:G:H2'	1.90	0.52
32:51:154:PRO:HB3	32:51:163:TYR:CE2	2.45	0.52
38:98:44:LEU:HD22	38:98:48:VAL:HG22	1.92	0.52
2:12:77:ALA:O	2:12:81:VAL:HG23	2.08	0.52
50:L8:4:LEU:HG	50:L8:39:ASP:HB2	1.92	0.52
12:3A:28:LYS:HZ1	12:3A:33:ARG:HH22	1.58	0.52
26:14:2860:A:N7	26:14:2861:G:H1'	2.25	0.52
1:1G:993:G:O2'	1:1G:994:A:N7	2.42	0.52
19:AI:51:VAL:HG12	19:AI:52:TYR:H	1.74	0.52
43:A5:51:LEU:HD23	43:A5:105:VAL:HG11	1.92	0.52
3:22:47:LEU:HG	3:22:68:VAL:HG11	1.91	0.52
35:25:59:LYS:HB3	35:25:87:ILE:HG22	1.91	0.52
47:I8:72:ARG:HB3	47:I8:75:LEU:HB2	1.90	0.52
43:E8:74:ALA:HA	43:E8:104:THR:O	2.09	0.52
1:1G:1372:U:OP1	9:82:72:GLY:N	2.43	0.52
28:11:26:LYS:HE2	28:11:94:LEU:HD12	1.90	0.52
39:65:102:ALA:HA	39:65:105:ALA:HB3	1.91	0.52
1:1G:604:G:H2'	1:1G:605:U:O4'	2.10	0.52
37:88:23:GLY:O	37:88:25:ASP:HB3	2.10	0.52
7:6E:15:ASP:OD2	7:6E:44:TYR:OH	2.28	0.52
46:D5:117:LEU:HD12	46:D5:119:GLU:HG2	1.91	0.52
5:4E:51:VAL:O	5:4E:55:VAL:HG23	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1683:C:H2'	26:1H:1684:C:C6	2.44	0.52
52:J5:11:THR:HG23	52:J5:15:ARG:HB3	1.92	0.52
1:13:31:G:O2'	1:13:48:C:N4	2.43	0.52
1:13:767:A:H2'	1:13:768:A:O4'	2.10	0.52
27:1J:14:U:H4'	27:1J:70:C:O2	2.10	0.52
24:3K:71:G:H2'	24:3K:72:C:H5''	1.91	0.52
1:13:1028(A):C:H42	1:13:1032(A):G:H1	1.58	0.52
26:14:1492:G:OP1	26:14:2210:G:N1	2.31	0.52
26:1H:2233:U:H2'	26:1H:2234:G:H8	1.72	0.52
28:19:237:GLU:HG3	28:19:239:ARG:H	1.75	0.52
39:A8:11:LYS:HD3	39:A8:91:PRO:HD3	1.91	0.52
26:14:40:C:H2'	26:14:41:C:H6	1.75	0.52
1:13:375:U:OP1	16:71:69:THR:HG21	2.10	0.52
30:31:37:VAL:HG21	36:78:6:LEU:HD21	1.91	0.52
1:1G:1207:G:H2'	1:1G:1208:C:C6	2.44	0.52
31:49:39:ILE:HD11	31:49:94:LEU:HD11	1.91	0.52
27:1J:52:A:N6	39:65:33:LYS:HG3	2.25	0.52
46:H8:11:GLU:HA	46:H8:36:LYS:HE3	1.92	0.52
26:1H:1127:A:O2'	26:1H:2518:A:OP1	2.20	0.52
26:1H:1049:C:C2'	26:1H:1050:A:H5'	2.40	0.52
1:13:1148:U:H2'	1:13:1149:C:O4'	2.09	0.52
17:8I:28:PRO:HA	17:8I:34:LYS:O	2.09	0.52
26:1H:1931:U:H5	26:1H:1969:A:N7	2.08	0.52
1:1G:32:A:H2'	1:1G:33:A:C8	2.44	0.52
26:1H:1826:G:OP2	28:11:222:ARG:HB3	2.10	0.52
34:15:42:TRP:O	41:85:64:ARG:NH2	2.33	0.52
26:1H:1547:C:H2'	26:1H:1548:C:C6	2.43	0.52
26:1H:859:G:O2'	26:1H:916:G:O6	2.27	0.52
2:12:95:GLN:OE1	2:12:147:LYS:HG2	2.09	0.52
32:51:97:ARG:NH2	32:51:104:GLU:OE2	2.40	0.52
26:14:2353:G:H2'	26:14:2354:G:O4'	2.10	0.52
1:13:360:A:H2'	1:13:361:G:C8	2.44	0.52
26:1H:2611:U:H2'	52:N8:3:LYS:HG3	1.91	0.52
26:14:363(B):G:H2'	26:14:363(C):G:C8	2.45	0.52
39:A8:30:ARG:HG3	39:A8:30:ARG:O	2.09	0.52
43:E8:84:ARG:HB3	43:E8:96:ILE:HD11	1.91	0.52
39:A8:48:LEU:CD2	39:A8:82:ILE:HD11	2.39	0.52
1:1G:590:C:H2'	1:1G:591:U:C6	2.45	0.52
26:14:782:A:H5'	26:14:783:A:C2	2.45	0.52
1:1G:358:U:H2'	1:1G:359:U:C6	2.45	0.52
1:13:407:G:O2'	4:3E:116:GLN:HG3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:654(M):C:H5'	26:1H:654(N):G:C5	2.45	0.52
47:E5:25:ARG:HD2	47:E5:29:GLN:NE2	2.25	0.52
26:14:2418:A:OP1	55:M5:29:LYS:NZ	2.38	0.52
26:14:2447:G:H3'	58:14:4011:HOH:O	2.08	0.52
36:78:2:LYS:HG2	36:78:4:SER:HB3	1.92	0.52
19:AI:11:VAL:HG13	19:AI:16:LEU:HD22	1.92	0.52
3:2E:7:PRO:O	3:2E:11:ARG:NH1	2.43	0.52
54:L5:10:ARG:O	54:L5:14:LYS:HB2	2.10	0.52
26:14:1936:A:C8	26:14:1940:U:O2	2.63	0.52
21:1B:6:ARG:HD3	21:1B:15:ARG:CZ	2.39	0.52
11:2I:20:TYR:HB2	11:2I:31:THR:HG22	1.92	0.52
32:59:105:LEU:HG	32:59:113:VAL:HG13	1.92	0.52
26:1H:1855:G:N7	58:1H:4282:HOH:O	2.33	0.52
10:1A:4:ILE:HA	10:1A:100:THR:HG22	1.91	0.52
43:E8:86:LEU:HD12	43:E8:86:LEU:C	2.29	0.52
32:51:105:LEU:H	32:51:105:LEU:HD23	1.73	0.52
17:8I:59:ILE:HG22	17:8I:73:VAL:HA	1.91	0.52
43:E8:57:ASN:O	43:E8:61:ASN:HB2	2.10	0.52
28:11:17:THR:CG2	28:11:204:ILE:HA	2.39	0.52
26:1H:607:U:O2	26:1H:621:A:N1	2.43	0.52
1:13:974:A:OP2	14:5I:29:ARG:NH2	2.43	0.52
1:13:976:G:H5'	1:13:1358:U:O2'	2.10	0.52
26:1H:1591:G:H2'	26:1H:1592:C:C6	2.44	0.52
24:3L:52:G:H1	24:3L:62:C:N4	2.03	0.52
26:1H:566:U:P	36:78:29:LYS:HZ2	2.32	0.52
10:1A:5:ARG:HG3	10:1A:73:ASP:OD1	2.10	0.52
38:98:52:ILE:O	38:98:55:ALA:N	2.42	0.52
1:1G:538:G:H2'	1:1G:539:A:C8	2.45	0.52
1:1G:538:G:H2'	1:1G:539:A:H8	1.74	0.52
26:1H:1858:G:H2'	26:1H:1883:G:N2	2.25	0.52
1:13:77:C:H2'	1:13:78:G:H5''	1.91	0.52
38:98:21:TYR:OH	38:98:43:GLU:HG2	2.10	0.52
26:1H:1012:U:O4	34:58:25:ARG:HA	2.09	0.52
1:1G:616:G:N3	1:1G:617:G:C8	2.78	0.52
34:15:128:HIS:NE2	34:15:134:ARG:HD2	2.25	0.52
1:1G:909:A:H2'	1:1G:910:C:O4'	2.10	0.52
24:3L:64:A:C2	24:3L:65:G:H1'	2.45	0.52
1:13:1081:G:H2'	1:13:1082:G:H8	1.75	0.52
1:13:222:U:H2'	1:13:223:U:H6	1.74	0.52
26:1H:2314:C:H2'	26:1H:2315:G:H8	1.75	0.52
26:1H:1950:G:N2	58:1H:4011:HOH:O	1.99	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:4E:122:GLU:HG2	5:4E:131:ILE:HD12	1.92	0.52
26:1H:2367:G:H2'	26:1H:2368:C:H6	1.75	0.52
26:14:1270:C:H5''	26:14:1271:G:H5'	1.92	0.52
11:2A:17:GLY:HA3	11:2A:77:MET:SD	2.50	0.52
42:95:85:LYS:HG3	42:95:87:HIS:N	2.25	0.52
26:1H:839:U:H2'	26:1H:840:C:C6	2.44	0.52
26:1H:194:G:H2'	26:1H:195:A:O4'	2.09	0.52
26:1H:987:G:O2'	26:1H:1000:A:N3	2.39	0.52
1:1G:1191:A:OP1	3:22:3:ASN:ND2	2.34	0.52
26:1H:674:G:P	58:1H:4586:HOH:O	2.67	0.52
29:21:46:ALA:HB1	29:21:80:GLU:HB3	1.91	0.52
4:3E:25:ARG:NH1	4:3E:30:LYS:HG3	2.24	0.52
34:58:128:HIS:HB2	34:58:129:PRO:HD2	1.90	0.52
26:14:796:C:H2'	26:14:797:C:H6	1.73	0.52
26:1H:580:C:H2'	26:1H:581:C:C6	2.45	0.52
37:45:37:LEU:HD11	37:45:130:LYS:HB2	1.91	0.52
2:12:9:GLU:HB2	2:12:217:ARG:NH2	2.24	0.52
26:1H:2845:G:O2'	26:1H:2846:G:H5'	2.09	0.52
2:1E:59:GLU:HB2	2:1E:221:LEU:HD11	1.91	0.52
1:13:654:G:H2'	1:13:655:A:H5'	1.92	0.52
32:59:44:VAL:HG13	32:59:51:ARG:HB2	1.92	0.52
26:1H:2343:C:O2'	26:1H:2373:G:O2'	2.23	0.52
28:19:80:ALA:HB3	28:19:94:LEU:HB3	1.92	0.52
26:1H:11:G:H2'	26:1H:12:U:H5'	1.92	0.52
43:A5:18:ARG:HG3	43:A5:76:VAL:HG13	1.92	0.52
26:14:1028:A:N3	26:14:2486:G:O2'	2.36	0.52
1:1G:587:G:N2	1:1G:754:C:OP2	2.43	0.52
13:4I:82:MET:O	13:4I:84:ILE:N	2.41	0.52
28:19:267:SER:O	28:19:268:ARG:HG2	2.10	0.52
1:1G:983:A:O2'	1:1G:1050:G:OP2	2.20	0.52
26:1H:1853:A:H2'	26:1H:1854:A:C8	2.45	0.52
37:88:37:LEU:HD21	37:88:130:LYS:HE2	1.92	0.52
2:1E:17:PHE:H	2:1E:17:PHE:HD1	1.56	0.52
1:1G:340:U:H3	1:1G:349:A:H61	1.56	0.52
29:21:1:MET:N	29:21:83:ASP:O	2.42	0.52
26:1H:588:U:C2	30:31:90:PHE:CE1	2.97	0.52
26:14:1204:A:O2'	26:14:1205:U:OP2	2.26	0.52
42:95:10:LYS:NZ	42:95:23:GLU:OE1	2.43	0.52
26:14:654(D):G:N2	26:14:654(Q):C:H42	2.01	0.52
1:1G:975:A:C4'	1:1G:976:G:H5''	2.38	0.52
1:1G:1132:C:H2'	1:1G:1133:G:C8	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1729:A:C2	26:14:1730:U:H5	2.27	0.52
24:3K:18:G:O2'	24:3K:19:G:OP1	2.26	0.52
1:1G:1281:U:H3'	1:1G:1282:C:H5	1.75	0.52
30:39:4:VAL:HG11	30:39:17:ARG:CZ	2.40	0.52
26:1H:2001:A:H2'	26:1H:2002:G:C8	2.45	0.52
7:62:46:ALA:O	7:62:50:ILE:HG12	2.09	0.52
26:14:2557:G:H2'	26:14:2558:C:H6	1.75	0.52
23:2L:48:U:H4'	23:2L:49:C:H5'	1.92	0.52
38:55:21:TYR:OH	38:55:43:GLU:HG2	2.09	0.52
26:1H:1107:G:H2'	26:1H:1108:U:C6	2.45	0.52
26:14:566:U:OP1	36:35:29:LYS:HD2	2.09	0.52
26:1H:1788:C:H2'	26:1H:1789:A:C8	2.45	0.52
27:16:37:C:C2'	27:16:38:C:H5'	2.40	0.52
2:1E:124:SER:HB2	2:1E:125:PRO:HD2	1.91	0.52
31:41:5:VAL:H	51:M8:25:TYR:HE2	1.58	0.52
1:1G:838:G:N2	1:1G:849:C:N3	2.57	0.52
46:H8:105:VAL:O	46:H8:140:ASP:HA	2.10	0.52
11:2A:21:ILE:HD12	11:2A:95:ILE:HD13	1.92	0.52
28:19:37:LEU:HA	28:19:38:LYS:CG	2.40	0.52
26:1H:102:G:OP1	49:K8:7:ARG:NH2	2.42	0.52
35:68:122:LEU:HD23	40:B8:43:GLN:HE22	1.75	0.52
26:1H:600:G:N2	26:1H:605:C:O3'	2.42	0.52
33:61:129:THR:HA	33:61:137:PRO:HA	1.91	0.52
55:Q8:48:PHE:CE1	55:Q8:49:VAL:HG13	2.46	0.51
26:1H:882:G:N2	26:1H:894:C:H42	2.00	0.51
26:14:1678:G:N2	26:14:1989:G:N2	2.56	0.51
23:2K:64:G:H2'	23:2K:65:G:C8	2.44	0.51
52:N8:31:VAL:HB	52:N8:41:PRO:O	2.10	0.51
26:1H:65:C:H2'	26:1H:66:C:C6	2.45	0.51
1:1G:1286:A:H3'	1:1G:1286:A:H8	1.75	0.51
26:14:2207:C:N3	26:14:2217:G:N2	2.47	0.51
38:98:55:ALA:HB2	38:98:79:LEU:HD13	1.93	0.51
26:1H:1086:A:H1'	26:1H:1103:A:H61	1.76	0.51
1:1G:1157:A:H61	1:1G:1178:G:N2	2.07	0.51
1:1G:1268:A:N3	1:1G:1326:C:O2'	2.43	0.51
5:4E:126:ARG:NH1	5:4E:126:ARG:HG3	2.25	0.51
26:14:1180:C:H2'	26:14:1181:C:C6	2.44	0.51
2:12:178:ARG:NH2	8:72:68:ARG:HH22	2.08	0.51
26:14:1188:U:C2'	26:14:1189:A:H5'	2.40	0.51
39:A8:15:ARG:HD2	39:A8:88:ASP:OD2	2.09	0.51
26:1H:661:C:O3'	36:78:15:ARG:NE	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:68:88:ASN:HD21	35:68:92:GLU:H	1.56	0.51
35:25:16:ALA:HB2	35:25:52:VAL:HG11	1.92	0.51
26:14:718:A:H3'	26:14:719:C:C6	2.44	0.51
32:59:137:ASP:HB2	32:59:140:LYS:HE2	1.90	0.51
39:65:27:SER:HA	39:65:88:ASP:HB2	1.91	0.51
47:E5:51:VAL:N	47:E5:62:LEU:HD12	2.26	0.51
1:13:444:C:H2'	1:13:445:G:C8	2.45	0.51
26:14:2636:U:OP1	29:29:80:GLU:HB2	2.10	0.51
1:1G:830:G:H2'	1:1G:831:U:O4'	2.10	0.51
26:14:198:C:H5'	26:14:2244:U:OP1	2.09	0.51
26:14:2139:C:N4	26:14:2152:G:O6	2.43	0.51
41:85:102:GLU:HG3	42:95:2:PHE:CE2	2.45	0.51
1:1G:1122:U:C4	1:1G:1123:A:N7	2.78	0.51
26:1H:1028:A:N6	26:1H:1125:G:H2'	2.25	0.51
27:1J:50:G:OP1	39:65:63:THR:HG23	2.09	0.51
26:14:1901:A:H2'	26:14:1901:A:N3	2.25	0.51
13:4I:27:LYS:HD3	13:4I:31:LYS:HZ3	1.74	0.51
7:62:69:VAL:HG12	7:62:103:TRP:HE3	1.75	0.51
26:1H:1242:A:H5''	26:1H:1243:G:OP2	2.10	0.51
1:13:806:C:H2'	1:13:807:A:H8	1.74	0.51
24:3K:7:A:O2'	24:3K:49:C:O4'	2.21	0.51
26:1H:1636:C:H2'	26:1H:1637:A:C8	2.45	0.51
26:1H:962:G:H2'	26:1H:963:U:C6	2.45	0.51
32:51:83:TYR:CB	32:51:134:SER:HA	2.40	0.51
26:1H:2210:G:H5'	26:1H:2211:G:H8	1.73	0.51
26:14:621:A:H3'	26:14:622:G:H8	1.74	0.51
26:1H:1388:G:O2'	26:1H:1389:G:H5'	2.10	0.51
16:7A:74:LEU:HD13	16:7A:79:VAL:HG21	1.93	0.51
1:13:1453:G:H8	20:BI:39:LYS:HE2	1.75	0.51
6:52:23:LYS:NZ	6:52:42:GLU:OE2	2.34	0.51
26:14:2068:U:N3	26:14:2430:A:H2	2.07	0.51
44:F8:11:PRO:HD3	49:K8:37:PHE:CD2	2.45	0.51
1:1G:1111:A:H2'	1:1G:1112:C:H6	1.75	0.51
28:19:68:LYS:HD3	28:19:70:TRP:CZ2	2.45	0.51
28:19:255:LYS:CE	28:19:255:LYS:H	2.23	0.51
1:13:237:C:H5''	17:8I:25:ARG:NH2	2.25	0.51
26:14:2245:U:H5''	26:14:2246:G:H5'	1.92	0.51
8:7E:58:TYR:O	8:7E:59:LEU:HD23	2.09	0.51
26:14:2461:C:H2'	26:14:2462:U:C6	2.45	0.51
26:1H:719:C:H2'	26:1H:720:C:C6	2.45	0.51
1:1G:284:G:H2'	1:1G:285:G:H8	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:567:G:O2'	58:1G:1835:HOH:O	2.19	0.51
27:1J:29:A:H5''	27:1J:30:C:OP2	2.10	0.51
55:Q8:53:PRO:HG3	55:Q8:56:GLU:HG3	1.92	0.51
46:H8:58:VAL:O	46:H8:60:GLU:N	2.42	0.51
13:4A:81:LEU:HD22	13:4A:88:ARG:HG2	1.91	0.51
26:1H:1509:C:H3'	26:1H:1510:A:H5''	1.93	0.51
26:1H:764:A:N3	28:11:213:ARG:HD3	2.25	0.51
1:1G:1369:C:H2'	1:1G:1370:G:O4'	2.10	0.51
41:C8:92:ARG:CZ	42:D8:11:GLN:H	2.24	0.51
1:1G:1316:G:H4'	14:5A:18:VAL:HG11	1.93	0.51
26:14:2261:C:H1'	26:14:2388:A:N3	2.26	0.51
26:1H:141(A):C:H2'	26:1H:142:G:O4'	2.11	0.51
2:1E:215:LEU:O	2:1E:219:VAL:HG23	2.10	0.51
26:14:943:U:P	36:35:36:LYS:HG3	2.50	0.51
4:32:108:LEU:HD11	4:32:174:LEU:HB3	1.92	0.51
24:3L:30:G:H2'	24:3L:31:A:C8	2.44	0.51
26:14:2690:C:OP1	38:55:17:ARG:NH1	2.42	0.51
38:55:21:TYR:HB3	38:55:47:PHE:CD2	2.45	0.51
28:11:134:ARG:HD3	28:11:135:PHE:CZ	2.46	0.51
3:2E:6:HIS:ND1	3:2E:7:PRO:HD2	2.25	0.51
33:61:77:LEU:HD11	33:61:140:LEU:HD23	1.91	0.51
1:1G:31:G:O2'	1:1G:48:C:N4	2.43	0.51
32:51:101:ARG:NH2	32:51:121:ILE:O	2.43	0.51
9:8E:43:ALA:HA	9:8E:74:ILE:HD13	1.91	0.51
1:13:1443:G:O2'	40:B8:122:ASP:OD2	2.23	0.51
42:D8:43:GLU:OE2	42:D8:44:LYS:NZ	2.38	0.51
23:2K:24:C:H2'	23:2K:25:U:C6	2.46	0.51
23:2L:21:U:O2	23:2L:21:U:H2'	2.10	0.51
40:75:107:ASP:N	40:75:107:ASP:OD1	2.44	0.51
33:69:92:VAL:HB	33:69:120:ILE:HG13	1.91	0.51
26:1H:1891:G:N7	58:1H:4325:HOH:O	2.34	0.51
26:14:1113:U:P	32:59:3:ARG:H	2.33	0.51
10:1I:5:ARG:HB2	10:1I:73:ASP:HA	1.92	0.51
1:13:21:G:OP1	58:13:1838:HOH:O	2.19	0.51
27:1J:15:A:H3'	27:1J:16:G:H5'	1.92	0.51
26:14:1900:A:OP2	58:14:3461:HOH:O	2.19	0.51
1:1G:411:A:H62	1:1G:413:G:N2	2.01	0.51
29:29:51:PHE:CE2	29:29:52:LEU:HG	2.45	0.51
51:M8:26:SER:OG	51:M8:27:THR:O	2.26	0.51
4:3E:31:CYS:SG	4:3E:32:ALA:N	2.83	0.51
26:1H:2345:G:N3	26:1H:2381:C:H2'	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2453:A:H2'	26:1H:2454:G:O4'	2.11	0.51
37:45:25:ASP:HB3	37:45:102:VAL:CG2	2.39	0.51
1:1G:630:G:H3'	1:1G:631:G:C8	2.45	0.51
2:12:92:TYR:CD1	2:12:151:GLY:HA3	2.46	0.51
26:1H:1047:G:H2'	26:1H:1110:G:N2	2.25	0.51
26:1H:2065:C:H2'	26:1H:2066:C:C6	2.46	0.51
1:1G:1446:A:OP1	1:1G:1446:A:H4'	2.11	0.51
1:1G:765:G:N2	1:1G:813:U:OP2	2.35	0.51
7:6E:79:ARG:NE	7:6E:84:ASN:HB2	2.25	0.51
26:1H:181:A:H1'	26:1H:435:C:H5'	1.91	0.51
34:15:111:PRO:HA	34:15:114:ARG:NH1	2.25	0.51
46:H8:111:VAL:HG11	46:H8:146:ILE:HD11	1.92	0.51
1:1G:581:G:OP1	15:6A:61:GLY:HA3	2.10	0.51
26:14:142:G:H2'	26:14:143:C:H6	1.75	0.51
1:1G:1338:G:N7	1:1G:1339:A:C6	2.78	0.51
26:14:903:C:H2'	26:14:904:C:C6	2.46	0.51
1:1G:442:C:H2'	1:1G:443:C:C6	2.46	0.51
55:Q8:6:THR:HG22	55:Q8:59:LYS:HG3	1.91	0.51
26:14:2016:U:O2	52:J5:7:PRO:HG2	2.11	0.51
9:8E:97:LYS:HD3	9:8E:102:LEU:HB2	1.92	0.51
28:19:69:ARG:NE	28:19:105:ILE:HD11	2.25	0.51
16:7A:57:ARG:HA	16:7A:60:LEU:HD12	1.92	0.51
44:F8:3:THR:HB	44:F8:6:ASP:HB2	1.93	0.51
27:1J:42:C:C4	31:49:91:ARG:NH2	2.79	0.51
26:14:2280:G:O2'	26:14:2388:A:N1	2.32	0.51
18:9A:32:ARG:HA	18:9A:69:THR:HG21	1.93	0.51
16:7I:53:VAL:O	16:7I:57:ARG:HG3	2.11	0.51
26:14:2575:C:H2'	26:14:2578:G:O6	2.11	0.51
22:1K:76:A:C2	26:1H:2602:A:H2	2.25	0.51
20:BA:26:ASN:HA	20:BA:29:LYS:HG2	1.92	0.51
36:78:101:VAL:HA	36:78:105:LEU:O	2.11	0.51
36:78:98:GLU:O	36:78:101:VAL:HG13	2.11	0.51
45:C5:86:ARG:HG3	45:C5:87:LYS:N	2.24	0.51
2:1E:6:THR:OG1	2:1E:7:VAL:N	2.44	0.51
26:1H:2815:C:H5'	52:N8:29:THR:HG21	1.93	0.51
28:11:71:ASP:HB2	28:11:103:ARG:HH22	1.76	0.51
26:14:117:G:OP1	26:14:124:G:N1	2.38	0.51
26:14:2542:A:H4'	26:14:2542:A:OP1	2.11	0.51
17:8I:29:HIS:CD2	17:8I:30:PRO:HD2	2.45	0.51
26:1H:1346:G:H2'	26:1H:1347:G:H8	1.74	0.51
24:3K:51:U:H2'	24:3K:52:G:C8	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2443:C:OP1	30:39:68:LYS:HG3	2.11	0.51
27:16:43:C:H5''	51:M8:1:MET:HG2	1.93	0.51
1:1G:287:U:H2'	1:1G:288:A:C8	2.46	0.51
48:F5:18:ILE:HG12	48:F5:37:ILE:HG13	1.92	0.51
32:59:55:PRO:HG2	32:59:61:HIS:CD2	2.46	0.51
22:1K:36:A:H62	22:1K:37:MIA:H122	1.75	0.51
1:1G:313:A:H2'	1:1G:314:C:C6	2.46	0.51
4:3E:128:VAL:HG23	4:3E:146:ILE:HG23	1.92	0.51
18:9A:41:LYS:O	18:9A:41:LYS:HD3	2.11	0.51
26:1H:2320:A:N3	26:1H:2320:A:H2'	2.25	0.51
26:1H:1359:A:N3	26:1H:1359:A:O4'	2.42	0.51
55:Q8:45:GLY:N	55:Q8:46:ARG:O	2.43	0.51
1:1G:859:A:OP2	1:1G:869:G:N2	2.37	0.51
30:39:5:ALA:CB	30:39:125:LEU:HD21	2.39	0.51
26:1H:566:U:P	36:78:29:LYS:NZ	2.84	0.51
20:BI:57:ARG:HH12	20:BI:102:GLY:C	2.14	0.51
26:14:2602:A:H4'	26:14:2603:G:O5'	2.10	0.51
24:3L:58:A:O2'	24:3L:60:U:OP2	2.27	0.51
1:13:1325:C:H2'	1:13:1326:C:C6	2.46	0.51
26:1H:1899:G:H22	26:1H:1902:C:N4	2.08	0.51
26:1H:1899:G:H1	26:1H:1902:C:N4	2.07	0.51
32:59:54:ARG:HD3	32:59:65:HIS:ND1	2.26	0.51
26:14:2130:U:H2'	26:14:2158:A:N1	2.26	0.51
8:72:30:ARG:O	8:72:34:GLU:HG2	2.10	0.51
26:1H:654:A:H2'	26:1H:654(A):A:H5''	1.93	0.51
1:13:511:C:H4'	4:3E:43:HIS:CD2	2.45	0.51
26:1H:1140:C:OP1	34:58:23:LEU:HB3	2.10	0.51
1:13:191(C):G:H2'	1:13:191(D):U:O4'	2.11	0.51
26:1H:618:G:H2'	26:1H:618(A):C:C6	2.46	0.51
1:13:484:G:O2'	1:13:485:G:OP2	2.25	0.51
26:14:1753:G:N1	26:14:1756:G:C2	2.79	0.51
1:1G:445:G:H2'	1:1G:446:G:C8	2.45	0.51
1:13:1385:G:N7	58:13:1950:HOH:O	2.34	0.51
1:13:1208:C:H2'	1:13:1209:C:O4'	2.10	0.51
33:69:29:TYR:O	33:69:32:PRO:HD2	2.10	0.51
26:14:686:G:OP1	54:L5:11:LYS:HE2	2.11	0.51
30:39:146:ALA:HB3	30:39:148:LEU:HG	1.92	0.51
28:19:182:LEU:H	28:19:272:ALA:HB2	1.76	0.51
9:8E:10:ARG:NH1	9:8E:75:ASP:OD2	2.44	0.51
55:M5:8:LYS:HB3	55:M5:12:LYS:HE3	1.92	0.51
30:39:40:GLN:HE22	30:39:182:ASN:HB2	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:C8:97:ASP:O	41:C8:101:ARG:N	2.41	0.51
26:14:1366:A:H2'	26:14:1367:A:O4'	2.10	0.51
26:14:2016:U:H2'	26:14:2017:U:C6	2.45	0.51
26:14:1161:C:H2'	26:14:1162:G:H8	1.76	0.51
29:21:116:VAL:O	29:21:117:MET:HB3	2.10	0.51
26:14:2331:G:H4'	47:E5:43:THR:N	2.21	0.51
1:13:963:G:H1	1:13:972:C:H42	1.59	0.51
26:1H:1388:G:H2'	26:1H:1389:G:C8	2.45	0.51
32:51:4:ILE:HG21	32:51:6:ARG:HH11	1.76	0.51
1:1G:1286:A:H3'	1:1G:1286:A:C8	2.46	0.51
1:13:1305:G:O2'	1:13:1331:G:N2	2.44	0.51
26:14:1557:C:H5''	26:14:1558:A:OP2	2.11	0.51
22:1L:14:A:C2	22:1L:22:G:H1'	2.46	0.51
26:1H:300:A:N3	26:1H:319:C:H1'	2.26	0.51
26:1H:71:A:H2	44:F8:31:HIS:CE1	2.26	0.51
49:K8:15:LYS:H	49:K8:67:LYS:HZ3	1.58	0.51
13:4A:70:LEU:O	13:4A:74:VAL:HG23	2.10	0.51
26:1H:779:U:O4	58:1H:4569:HOH:O	2.19	0.51
26:14:817:C:H2'	26:14:818:G:O4'	2.10	0.51
26:14:2655:G:N2	26:14:2665:A:OP2	2.44	0.51
26:14:126:A:O5'	54:L5:19:ARG:HG3	2.11	0.51
26:14:287:C:H2'	26:14:288:C:H6	1.76	0.51
26:14:218:A:C2	26:14:235:U:H4'	2.45	0.51
1:13:922:G:C6	1:13:923:A:C6	2.98	0.51
26:1H:1260:G:H2'	26:1H:1261:C:C6	2.45	0.51
26:1H:784:A:H5''	28:11:227:ASN:ND2	2.25	0.51
38:55:78:LYS:O	38:55:83:ILE:HG13	2.11	0.51
26:1H:2841:C:H2'	26:1H:2842:G:H8	1.74	0.51
26:14:2788:C:O2'	26:14:2809:A:N3	2.44	0.51
26:14:1386:C:H2'	26:14:1387:C:C6	2.46	0.51
27:16:24:G:N7	27:16:56:G:H2'	2.25	0.51
26:14:863:A:H2	26:14:914:C:N4	2.08	0.51
6:52:41:GLU:HG3	6:52:62:TRP:CE3	2.46	0.51
26:14:1266:G:O5'	43:A5:15:ARG:NH2	2.44	0.51
12:3I:27:LEU:HD23	12:3I:33:ARG:HB2	1.92	0.51
1:13:998:G:H2'	1:13:998(A):C:C6	2.46	0.51
26:14:2383:G:O2'	26:14:2384:G:H5'	2.10	0.51
26:14:1639:U:P	58:14:4066:HOH:O	2.67	0.51
2:12:98:LEU:O	2:12:101:MET:HG2	2.10	0.51
36:78:49:ARG:NH2	55:Q8:59:LYS:HZ3	2.09	0.51
26:1H:751:A:P	58:1H:3922:HOH:O	2.69	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:H8:3:TYR:O	46:H8:58:VAL:HG23	2.11	0.51
26:14:137(A):G:H2'	26:14:139:G:C8	2.46	0.51
30:39:83:PHE:C	30:39:85:GLY:H	2.13	0.51
13:4A:3:ARG:HG2	13:4A:9:ILE:HG13	1.93	0.51
37:45:26:TYR:O	37:45:26:TYR:CD1	2.64	0.51
1:1G:353:A:H5'	1:1G:353:A:C8	2.40	0.51
1:1G:403:C:N4	58:1G:1804:HOH:O	2.44	0.51
26:1H:993:G:C4	26:1H:994:C:C5	2.99	0.51
23:2K:17:C:H2'	23:2K:18:C:H2'	1.92	0.51
26:1H:2053:G:H5'	29:21:144:ARG:O	2.10	0.51
12:3I:53:ARG:HH12	12:3I:92:ASP:HB2	1.76	0.51
1:1G:625:G:H2'	1:1G:626:U:C6	2.44	0.51
19:AI:41:VAL:H	19:AI:44:MET:HB2	1.75	0.51
32:51:149:ARG:HA	32:51:162:ILE:CG2	2.40	0.51
1:1G:1055:A:C6	1:1G:1056:U:O2	2.64	0.51
26:1H:580:C:H2'	26:1H:581:C:H6	1.76	0.51
1:1G:1105:A:H2'	1:1G:1106:G:C8	2.45	0.51
52:J5:3:LYS:CE	52:J5:3:LYS:HA	2.38	0.51
26:14:1449(A):G:H2'	26:14:1450:C:C6	2.46	0.51
1:13:295:C:H2'	1:13:296:U:O4'	2.10	0.51
1:1G:1194:U:H2'	1:1G:1195:C:C6	2.46	0.51
46:H8:5:LEU:HD23	46:H8:47:VAL:HG21	1.93	0.51
26:1H:192:C:P	58:1H:4537:HOH:O	2.68	0.51
3:22:84:ILE:HG12	3:22:88:ARG:NH2	2.26	0.51
16:7I:45:THR:HG22	16:7I:46:PRO:HD2	1.93	0.51
34:58:90:MET:O	34:58:94:HIS:N	2.30	0.51
10:1A:38:ILE:HB	10:1A:71:LEU:HB3	1.93	0.51
26:14:1171:G:O2'	26:14:1173:G:O4'	2.29	0.51
42:D8:76:LYS:HB2	42:D8:81:TYR:HB3	1.93	0.51
26:14:2757:A:N1	32:59:67:LEU:HD22	2.26	0.51
18:9I:53:ARG:NE	18:9I:59:SER:O	2.43	0.51
1:13:524:G:O5'	1:13:524:G:H8	1.94	0.51
26:1H:2033:A:OP1	58:1H:4062:HOH:O	2.18	0.51
32:51:51:ARG:HG2	32:51:52:VAL:H	1.76	0.51
1:1G:1028(A):C:H5	1:1G:1029:G:C2	2.29	0.51
40:75:8:LYS:HZ3	40:75:8:LYS:HB2	1.76	0.51
44:F8:3:THR:HA	44:F8:6:ASP:CG	2.31	0.51
26:14:2061:G:P	58:14:4027:HOH:O	2.68	0.51
3:22:6:HIS:CG	14:5A:49:HIS:HB3	2.46	0.51
26:14:1378:A:O2'	26:14:1380:G:N7	2.40	0.51
45:C5:87:LYS:CB	45:C5:94:LYS:HA	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:3E:92:VAL:O	4:3E:96:LEU:HD22	2.11	0.51
27:16:15:A:O2'	27:16:109:G:C8	2.60	0.51
1:1G:1478:C:H2'	1:1G:1479:C:C6	2.45	0.51
29:29:120:TRP:CE3	29:29:155:LYS:HD3	2.45	0.51
31:49:32:PRO:HB2	31:49:172:LEU:HD22	1.92	0.51
1:1G:448:A:P	1:1G:485:G:H22	2.34	0.51
26:14:1726:G:N2	26:14:1734:C:O2	2.40	0.51
14:5A:43:CYS:O	14:5A:46:GLU:HG3	2.11	0.51
3:2E:64:VAL:HG23	3:2E:98:ASN:O	2.11	0.51
38:98:18:LEU:HD11	38:98:22:ARG:CZ	2.41	0.51
16:7A:75:ARG:O	16:7A:78:GLY:N	2.40	0.51
22:1K:44:G:H2'	22:1K:45:U:C6	2.46	0.51
9:82:26:VAL:HG13	9:82:61:ALA:HB3	1.93	0.51
26:1H:325:G:H2'	26:1H:326:G:H8	1.76	0.51
19:AA:11:VAL:HG12	19:AA:39:THR:HB	1.93	0.51
5:4E:81:GLU:HG2	5:4E:90:VAL:HG23	1.93	0.51
29:29:52:LEU:C	29:29:74:PRO:HB3	2.31	0.51
26:1H:654(D):G:N2	26:1H:654(R):C:N3	2.59	0.51
4:3E:30:LYS:CB	4:3E:35:ARG:HE	2.24	0.51
45:G8:85:VAL:O	45:G8:86:ARG:HD3	2.11	0.51
26:14:2387:U:P	47:E5:55:ARG:HH22	2.34	0.51
1:13:1194:U:H2'	1:13:1195:C:H6	1.75	0.51
34:58:132:ALA:HB1	34:58:133:GLN:OE1	2.11	0.51
26:1H:1407:C:H2'	26:1H:1408:C:H6	1.75	0.51
26:14:71:A:H2	44:B5:31:HIS:NE2	2.07	0.51
19:AA:10:PHE:O	51:I5:63:TYR:OH	2.28	0.51
2:1E:178:ARG:HB2	2:1E:178:ARG:HH11	1.76	0.51
50:L8:7:LYS:HE3	50:L8:32:GLN:HA	1.93	0.51
27:16:31:C:H2'	27:16:32:C:H6	1.76	0.51
4:3E:173:TRP:CD1	4:3E:189:PRO:HG3	2.46	0.51
37:45:110:THR:HG23	37:45:113:GLN:HE21	1.76	0.51
38:98:32:GLY:HA2	38:98:116:LEU:HD12	1.92	0.51
1:13:1446:A:H4'	1:13:1446:A:OP1	2.11	0.51
26:1H:1805:U:O2	28:11:50:THR:HB	2.11	0.51
30:31:53:THR:O	30:31:56:GLU:N	2.39	0.51
26:14:706:A:H2'	26:14:707:G:O4'	2.11	0.51
32:51:33:LEU:HD21	32:51:136:ILE:HB	1.93	0.51
4:3E:52:SER:O	4:3E:55:ALA:HB3	2.11	0.51
32:59:60:ARG:O	32:59:63:SER:OG	2.29	0.51
1:13:793:U:H5'	1:13:794:A:H5''	1.92	0.51
26:1H:1200:C:H5'	58:1H:3697:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:671:C:OP1	36:78:42:SER:O	2.29	0.51
41:85:74:LEU:HD13	41:85:79:PHE:HB2	1.92	0.51
48:F5:51:VAL:HG23	48:F5:58:ILE:HB	1.93	0.51
1:13:1447:G:O5'	1:13:1447:G:H8	1.94	0.51
55:Q8:48:PHE:CE2	55:Q8:49:VAL:HG13	2.46	0.50
39:65:109:GLY:O	39:65:110:LEU:HD22	2.11	0.50
26:14:751:A:H2'	26:14:789:A:C2	2.47	0.50
32:51:83:TYR:CG	32:51:84:SER:N	2.78	0.50
26:1H:674:G:OP2	58:1H:4586:HOH:O	2.18	0.50
13:4A:3:ARG:HG2	13:4A:9:ILE:CG1	2.41	0.50
41:85:90:VAL:HG22	42:95:39:LEU:CB	2.38	0.50
26:14:2115:G:H2'	26:14:2116:G:C8	2.46	0.50
27:16:7:G:O5'	39:A8:29:PHE:CE2	2.64	0.50
34:58:133:GLN:HG2	34:58:134:ARG:H	1.76	0.50
24:3K:33:U:H3'	24:3K:34:G:H5''	1.93	0.50
53:O8:26:ASN:OD1	53:O8:28:ARG:HB2	2.11	0.50
1:13:114:U:O2'	1:13:115:G:H5'	2.10	0.50
26:14:330:A:H2	26:14:1210:A:HO2'	1.56	0.50
30:39:132:VAL:C	30:39:134:GLY:H	2.14	0.50
1:13:1226:C:H2'	13:4I:103:THR:HB	1.93	0.50
26:14:1357:U:H2'	26:14:1358:G:O4'	2.11	0.50
26:14:959:A:C6	26:14:960:A:C2	2.99	0.50
28:19:70:TRP:C	28:19:70:TRP:CD1	2.85	0.50
19:AI:36:ARG:NH1	19:AI:52:TYR:O	2.41	0.50
26:14:1856:G:H1	26:14:1886:C:H42	1.59	0.50
1:13:1221:G:O3'	19:AI:77:THR:HG21	2.11	0.50
1:13:757:U:H2'	1:13:758:G:O4'	2.11	0.50
1:1G:1449:C:H3'	1:1G:1450:U:H4'	1.93	0.50
1:13:1530:G:O2'	1:13:1531:A:OP2	2.24	0.50
26:1H:627:A:H4'	26:1H:628:G:OP1	2.11	0.50
1:1G:1151:A:O2'	1:1G:1152:A:O5'	2.28	0.50
44:B5:24:GLY:HA3	44:B5:82:GLN:HG3	1.93	0.50
47:E5:23:VAL:HG13	47:E5:38:VAL:HG23	1.93	0.50
26:14:1994:C:OP1	58:14:3493:HOH:O	2.19	0.50
49:K8:8:LYS:HZ3	49:K8:12:GLU:HG2	1.76	0.50
26:1H:757:U:H2'	26:1H:758:C:O4'	2.11	0.50
7:62:32:ARG:O	7:62:34:GLY:N	2.44	0.50
1:13:813:U:OP2	1:13:813:U:H6	1.94	0.50
26:1H:2577:A:H4'	52:N8:3:LYS:HB3	1.93	0.50
10:1I:35:SER:OG	10:1I:73:ASP:HB2	2.11	0.50
41:85:92:ARG:HH22	42:95:10:LYS:HA	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1002:G:C4	1:13:1003:G:C8	3.00	0.50
30:31:108:LYS:O	30:31:112:MET:HG3	2.10	0.50
30:39:25:PRO:C	30:39:27:GLU:N	2.65	0.50
24:3L:48:C:C5	24:3L:59:U:H1'	2.46	0.50
44:F8:24:GLY:CA	44:F8:82:GLN:HE22	2.19	0.50
1:13:1243:C:OP1	21:1F:10:ARG:HG3	2.11	0.50
48:J8:91:LYS:CG	48:J8:92:LYS:H	2.23	0.50
20:BI:73:HIS:HB3	20:BI:74:LYS:HZ3	1.76	0.50
26:14:2134:A:N6	26:14:2157:G:O2'	2.43	0.50
26:1H:1800:C:OP1	28:11:266:SER:OG	2.24	0.50
1:1G:589:C:C2	1:1G:650:G:N2	2.77	0.50
40:B8:20:PRO:HG2	40:B8:86:ILE:O	2.12	0.50
26:14:71:A:H8	26:14:71:A:H5'	1.74	0.50
1:1G:265:G:O3'	17:8A:66:SER:HA	2.11	0.50
34:15:128:HIS:HB2	34:15:129:PRO:HD2	1.94	0.50
26:14:1530:G:O6	26:14:1542:G:N2	2.44	0.50
13:4I:84:ILE:HG13	19:AI:74:PHE:HE2	1.76	0.50
51:M8:54:GLY:O	51:M8:58:ARG:N	2.43	0.50
5:4E:15:ARG:NH1	25:4K:25:A:H3'	2.25	0.50
1:1G:1293:G:H2'	1:1G:1294:G:C8	2.46	0.50
26:1H:556:G:H2'	26:1H:557:U:C6	2.45	0.50
26:14:2808:U:H5''	26:14:2891:G:O6	2.11	0.50
30:31:123:LEU:HD12	30:31:124:LEU:H	1.75	0.50
1:13:690:G:H2'	1:13:691:G:O4'	2.12	0.50
1:13:1152:A:H4'	10:1I:13:HIS:CD2	2.46	0.50
1:1G:1040:U:H2'	1:1G:1041:A:H8	1.75	0.50
1:13:1106:G:H5''	3:2E:172:ARG:HG2	1.92	0.50
26:1H:1142(A):A:C4	26:1H:1144:G:C8	2.99	0.50
30:39:117:ARG:NH1	30:39:120:GLU:OE1	2.44	0.50
29:21:117:MET:HE1	29:21:136:ARG:CA	2.38	0.50
24:3L:51:U:H2'	24:3L:52:G:H8	1.76	0.50
36:78:29:LYS:CD	36:78:30:THR:HG22	2.37	0.50
43:E8:38:TYR:OH	52:N8:47:PRO:HG3	2.12	0.50
12:3A:60:LEU:HB2	12:3A:64:TYR:CB	2.38	0.50
43:E8:12:ILE:HD13	43:E8:17:VAL:HB	1.93	0.50
1:1G:164:U:H2'	1:1G:165:C:H6	1.77	0.50
1:13:684:A:C6	1:13:685:G:C6	3.00	0.50
26:14:1333:C:H2'	26:14:1334:G:H8	1.77	0.50
27:16:48:A:H4'	39:A8:95:HIS:HD2	1.76	0.50
36:35:89:ALA:HB1	36:35:121:LYS:HD2	1.93	0.50
1:13:560:U:H5'	1:13:566:G:N2	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:88:75:THR:HA	37:88:87:LYS:O	2.12	0.50
1:13:628:G:H2'	1:13:629:G:C8	2.46	0.50
26:1H:2469:A:O2'	37:88:56:ARG:NE	2.41	0.50
45:G8:82:PRO:HG3	45:G8:97:ARG:HA	1.93	0.50
45:G8:97:ARG:N	45:G8:97:ARG:HD2	2.26	0.50
30:31:34:TRP:HB2	36:78:6:LEU:HG	1.94	0.50
4:32:172:PRO:HD2	4:32:173:TRP:CZ3	2.46	0.50
1:13:878:G:H5'	8:7E:89:PRO:HG2	1.93	0.50
1:1G:749:C:H2'	1:1G:750:G:H8	1.77	0.50
3:22:70:VAL:HG21	3:22:76:VAL:HG11	1.94	0.50
26:14:1348:G:C6	26:14:1349:A:N1	2.79	0.50
47:I8:44:ARG:HB2	47:I8:44:ARG:NH1	2.26	0.50
27:1J:113:C:O2'	39:65:46:VAL:HG13	2.12	0.50
8:7E:49:GLU:O	8:7E:51:VAL:HG13	2.10	0.50
1:13:587:G:N2	1:13:755:G:C5	2.79	0.50
26:14:57:C:H2'	26:14:58:G:O4'	2.12	0.50
30:39:114:VAL:HG21	30:39:202:PHE:CE1	2.46	0.50
26:1H:276:A:C8	26:1H:278:A:H2	2.29	0.50
26:14:433:C:C4	26:14:434:U:O4	2.65	0.50
1:13:266:G:H5''	1:13:267:C:C5	2.46	0.50
26:14:2239:G:P	58:14:3404:HOH:O	2.69	0.50
1:1G:685:G:C2	1:1G:686:U:C4	2.99	0.50
11:2I:57:THR:HG22	11:2I:60:ALA:H	1.76	0.50
1:1G:986:A:H1'	19:AA:54:GLY:O	2.11	0.50
38:55:103:ARG:HH21	38:55:110:PRO:HB3	1.76	0.50
1:1G:1260:C:H6	1:1G:1260:C:H3'	1.76	0.50
54:P8:9:ARG:HE	54:P8:47:ARG:HB2	1.76	0.50
10:1I:57:LYS:O	10:1I:60:ARG:NH1	2.45	0.50
26:1H:910:A:C4	37:88:13:GLN:NE2	2.80	0.50
1:1G:1224:G:N1	1:1G:1322:C:H1'	2.27	0.50
28:11:85:ASP:HB2	28:11:92:ILE:HG12	1.93	0.50
33:69:76:THR:HG23	33:69:77:LEU:N	2.26	0.50
22:1L:75:C:H2'	22:1L:76:A:H5''	1.93	0.50
40:75:5:ALA:O	40:75:9:LEU:HB2	2.11	0.50
1:1G:1153:C:H2'	1:1G:1154:G:C8	2.46	0.50
23:2K:70:C:H2'	23:2K:71:G:O4'	2.11	0.50
28:11:108:PRO:HD2	28:11:111:LEU:HG	1.94	0.50
26:14:602:G:N2	26:14:655:A:C8	2.77	0.50
3:22:91:LEU:HB2	3:22:99:VAL:HG11	1.94	0.50
1:1G:1095:U:OP1	1:1G:1108:G:N2	2.44	0.50
1:1G:777:A:C2	11:2A:119:CYS:HB3	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:656:G:H2'	26:14:657:U:O4'	2.11	0.50
54:P8:15:THR:HG22	54:P8:16:HIS:CE1	2.46	0.50
29:29:120:TRP:CD2	29:29:155:LYS:HD3	2.47	0.50
13:4I:108:ARG:NH1	13:4I:112:GLY:O	2.44	0.50
33:61:68:LEU:HA	33:61:71:ILE:CG2	2.41	0.50
34:58:36:GLY:HA2	34:58:38:HIS:CE1	2.46	0.50
1:1G:688:G:O2'	1:1G:704:A:N1	2.40	0.50
9:8E:7:THR:O	9:8E:83:ARG:HD2	2.11	0.50
17:8I:45:HIS:O	17:8I:73:VAL:HG23	2.12	0.50
26:14:686:G:N7	54:L5:5:TRP:CH2	2.80	0.50
26:1H:2294:C:C4	26:1H:2295:C:C5	3.00	0.50
1:1G:1231:G:O3'	9:82:126:SER:OG	2.24	0.50
1:1G:662:G:O2'	1:1G:836:G:OP1	2.28	0.50
31:41:104:GLU:OE1	51:M8:23:GLU:HG3	2.11	0.50
7:6E:18:TYR:HD2	7:6E:59:LEU:HD22	1.76	0.50
3:2E:124:ILE:HG12	3:2E:130:VAL:HG22	1.93	0.50
26:14:1794:U:O2'	26:14:1795:C:H5'	2.12	0.50
1:13:153:C:N4	1:13:168:G:H1	2.09	0.50
26:14:1152:C:H4'	41:85:77:SER:HA	1.94	0.50
26:14:547:A:H3'	26:14:548:A:C8	2.47	0.50
2:1E:163:PHE:HA	2:1E:185:ILE:O	2.12	0.50
26:1H:2700:C:C2'	26:1H:2701:C:H5'	2.40	0.50
36:35:15:ARG:NH1	36:35:15:ARG:HB2	2.27	0.50
30:31:29:ASN:H	30:31:112:MET:CE	2.22	0.50
11:2I:106:LYS:O	11:2I:107:SER:OG	2.24	0.50
26:14:2472:G:H8	26:14:2472:G:O5'	1.94	0.50
16:7A:18:ARG:HA	16:7A:38:TYR:HA	1.92	0.50
12:3A:55:VAL:HA	12:3A:70:ILE:HG13	1.92	0.50
6:5E:10:LEU:HD13	6:5E:61:LEU:HD13	1.93	0.50
26:1H:2344:U:O2'	53:O8:37:ARG:HG2	2.11	0.50
12:3I:93:LEU:O	12:3I:96:VAL:HG13	2.12	0.50
26:1H:817:C:H4'	26:1H:932:G:C5	2.47	0.50
1:13:1414:U:H2'	1:13:1415:G:C8	2.45	0.50
47:E5:26:TYR:O	47:E5:29:GLN:HB2	2.12	0.50
26:1H:475:U:C4	26:1H:481:G:O6	2.65	0.50
26:1H:1465:G:H5'	26:1H:1528:A:H1'	1.94	0.50
26:14:1871:A:H2'	26:14:1872:A:C8	2.46	0.50
21:1B:9:ARG:CZ	21:1B:10:ARG:HG2	2.42	0.50
15:6I:12:ILE:HG12	15:6I:31:LEU:HD11	1.94	0.50
39:65:87:PHE:CZ	39:65:102:ALA:HB2	2.46	0.50
16:7I:17:TYR:HE2	16:7I:41:PRO:HG3	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:392:C:H5''	26:14:409:C:H5''	1.93	0.50
28:11:147:LEU:HD13	28:11:155:LEU:HD21	1.93	0.50
46:H8:6:LYS:HE3	46:H8:8:TYR:HE1	1.75	0.50
16:7A:68:ASP:O	16:7A:71:ARG:HB3	2.12	0.50
6:5E:75:LEU:HD22	6:5E:79:LEU:HG	1.94	0.50
1:13:79:G:H2'	1:13:79:G:N3	2.25	0.50
26:14:2660:A:OP1	26:14:2660:A:H8	1.94	0.50
26:1H:1636:C:P	58:1H:3609:HOH:O	2.69	0.50
55:Q8:25:MET:HG2	55:Q8:46:ARG:HG2	1.94	0.50
55:Q8:7:HIS:CD2	55:Q8:57:ARG:NH2	2.79	0.50
26:14:323:G:H5'	30:39:169:ASN:HD21	1.77	0.50
2:12:43:ASP:O	2:12:47:THR:OG1	2.29	0.50
26:14:1729:A:H2	26:14:1730:U:H5	1.58	0.50
46:D5:30:ASN:HA	46:D5:89:PHE:CE1	2.41	0.50
1:13:625:G:H2'	1:13:626:U:H6	1.75	0.50
2:12:141:GLU:O	2:12:145:LEU:HB2	2.12	0.50
26:14:39:C:O2'	26:14:40:C:H5'	2.12	0.50
1:13:407:G:H2'	1:13:408:A:C8	2.46	0.50
19:AI:30:LEU:HA	19:AI:48:THR:O	2.11	0.50
4:32:61:LYS:HA	4:32:203:VAL:HG22	1.93	0.50
26:1H:581:C:H2'	26:1H:582:G:H8	1.77	0.50
11:2I:78:GLN:O	11:2I:103:LEU:HA	2.11	0.50
26:1H:1420:U:HO2'	26:1H:1421:G:P	2.35	0.50
36:35:101:VAL:HG23	36:35:107:LYS:N	2.26	0.50
26:1H:527:C:OP2	26:1H:2779:U:C5	2.64	0.50
26:14:184:C:H2'	26:14:185:U:C6	2.46	0.50
35:25:13:ASN:HD21	35:25:97:ARG:H	1.60	0.50
19:AI:7:LYS:HB3	19:AI:7:LYS:NZ	2.27	0.50
26:1H:1346:G:H2'	26:1H:1347:G:C8	2.46	0.50
7:6E:69:VAL:HG12	7:6E:100:ALA:HA	1.93	0.50
30:39:51:THR:HG23	30:39:92:PRO:HG2	1.93	0.50
26:14:2867:G:OP2	40:75:119:LYS:NZ	2.31	0.50
27:16:91:C:H5''	46:H8:79:ARG:NH2	2.26	0.50
32:59:76:VAL:HA	32:59:79:VAL:HG22	1.94	0.50
31:49:103:LEU:HD22	31:49:178:PHE:HZ	1.76	0.50
1:13:1157:A:H1'	1:13:1158:C:C4	2.46	0.50
26:1H:631:A:H61	26:1H:2402:C:N4	2.09	0.50
19:AA:29:ARG:HD3	19:AA:48:THR:H	1.76	0.50
30:39:83:PHE:O	30:39:85:GLY:N	2.45	0.50
26:1H:2124:G:H2'	26:1H:2125:G:H5'	1.92	0.50
11:2A:48:ILE:HG23	11:2A:63:LEU:HD22	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2472:G:H21	26:14:2478:A:H62	1.58	0.50
26:1H:2327:A:H2'	26:1H:2328:A:H8	1.77	0.50
49:K8:22:GLU:OE2	49:K8:68:ARG:NH2	2.37	0.50
16:7I:53:VAL:HG13	16:7I:79:VAL:HG22	1.93	0.50
1:1G:963:G:H21	10:1A:55:LYS:NZ	2.09	0.50
2:12:70:PHE:HB2	2:12:92:TYR:HB2	1.94	0.50
1:1G:987:G:H1	1:1G:1218:C:H42	1.59	0.50
1:1G:1207:G:H2'	1:1G:1208:C:H6	1.77	0.50
12:3I:42:THR:HG22	12:3I:54:LYS:HG3	1.92	0.50
26:14:185:U:H4'	26:14:218:A:H4'	1.93	0.50
26:14:528:A:C2	26:14:2043:C:H4'	2.47	0.50
42:95:76:LYS:HD2	42:95:80:GLN:O	2.11	0.50
28:19:267:SER:HA	28:19:270:ILE:HD12	1.94	0.50
37:88:35:VAL:HG13	37:88:130:LYS:HB3	1.94	0.50
26:14:2677:G:H2'	26:14:2678:C:C6	2.46	0.50
13:4I:107:ALA:HB3	13:4I:111:LYS:HD2	1.93	0.50
26:1H:2109:U:N3	26:1H:2110:G:O6	2.45	0.50
26:14:107:C:H2'	26:14:108:U:H6	1.76	0.50
26:1H:855:G:O2'	47:I8:27:GLU:HG2	2.12	0.50
52:J5:36:CYS:HG	52:J5:49:CYS:HB3	1.77	0.50
40:B8:47:GLY:HA3	40:B8:65:LYS:HB3	1.93	0.50
26:1H:565:C:H4'	58:1H:3712:HOH:O	2.11	0.50
42:D8:20:LEU:HD21	42:D8:22:VAL:HG23	1.94	0.50
26:1H:59:U:O2'	26:1H:73:A:H2'	2.11	0.50
37:45:17:LEU:HD21	37:45:41:TRP:HE1	1.75	0.50
26:14:2688:U:H1'	26:14:2721:A:N6	2.27	0.50
41:85:92:ARG:O	41:85:94:ASN:N	2.44	0.50
17:8I:65:ILE:HB	17:8I:69:LYS:HB3	1.94	0.50
26:14:654(C):G:H2'	26:14:654(D):G:O4'	2.11	0.50
26:14:1025:G:C4	26:14:1135:C:H1'	2.46	0.50
49:G5:25:VAL:O	49:G5:29:LYS:HB2	2.10	0.50
13:4A:57:ARG:NH1	51:I5:34:GLU:O	2.45	0.50
1:1G:979:C:H3'	1:1G:980:C:C5'	2.39	0.50
28:11:231:HIS:CD2	28:11:249:PRO:HA	2.47	0.50
29:29:62:PRO:C	29:29:64:LYS:H	2.15	0.50
32:59:6:ARG:HD3	32:59:54:ARG:HH12	1.77	0.50
1:1G:1316:G:H5''	14:5A:17:LYS:NZ	2.27	0.50
44:F8:1:MET:C	44:F8:3:THR:N	2.65	0.50
26:1H:2261:C:O2'	26:1H:2262:U:H5'	2.11	0.50
26:14:1352:U:O2	26:14:1570:A:H2	1.95	0.50
41:C8:6:THR:HG21	41:C8:10:ARG:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:31:101:LEU:CD1	30:31:105:VAL:HB	2.42	0.50
32:51:10:PRO:HG2	32:51:50:VAL:O	2.12	0.50
3:22:18:TRP:HE1	14:5A:55:GLY:H	1.60	0.50
26:14:1430:C:H2'	26:14:1431:U:H6	1.75	0.50
30:39:36:VAL:HG11	30:39:183:VAL:HG11	1.93	0.50
1:1G:485:G:O2'	1:1G:486:U:O5'	2.30	0.50
26:1H:1913:A:H4'	26:1H:1914:C:H5'	1.93	0.50
26:1H:459:U:H5''	54:P8:40:TRP:CD2	2.46	0.50
2:12:97:TRP:CZ3	2:12:99:GLY:HA2	2.47	0.50
15:6I:63:ARG:O	15:6I:67:LEU:HD12	2.12	0.50
30:39:93:LYS:HB3	30:39:94:PRO:HD2	1.93	0.50
26:1H:2243:U:O2'	26:1H:2244:U:H5'	2.12	0.50
26:1H:1335:U:H2'	26:1H:1336:A:O4'	2.12	0.50
22:1L:19:G:N2	22:1L:56:C:H42	2.09	0.50
1:1G:1502:A:H2	1:1G:1505:G:N1	2.01	0.50
1:1G:1131:G:H2'	1:1G:1132:C:H6	1.77	0.50
28:11:60:ARG:HD3	28:11:86:PRO:CB	2.38	0.50
26:1H:225:A:H61	26:1H:419:C:H4'	1.77	0.50
31:41:112:PRO:HB3	51:M8:36:CYS:HA	1.93	0.50
31:41:112:PRO:HG3	51:M8:38:LYS:HD2	1.94	0.50
53:O8:17:LYS:O	53:O8:19:ARG:N	2.42	0.50
1:13:55:A:H2	33:69:82:ARG:HG2	1.76	0.50
27:16:73:A:N3	27:16:73:A:H2'	2.27	0.50
26:1H:212:G:H2'	26:1H:213:A:O4'	2.12	0.50
26:1H:981:A:C8	26:1H:982:C:C5	2.98	0.50
2:12:12:GLU:HB3	2:12:213:LEU:HD22	1.94	0.50
50:L8:6:VAL:HG12	50:L8:56:VAL:HG22	1.94	0.50
1:1G:1228:C:P	13:4A:115:LYS:HZ3	2.35	0.50
26:1H:2591:C:OP1	28:11:239:ARG:HG3	2.12	0.50
30:31:65:TRP:CZ3	30:31:72:ARG:HB3	2.47	0.50
26:14:747:U:C6	52:J5:2:ALA:HB3	2.47	0.50
26:1H:184:C:H2'	26:1H:185:U:C6	2.47	0.50
1:1G:994:A:N3	14:5A:5:ALA:HB2	2.26	0.50
26:14:2528:U:O3'	26:14:2529:G:N2	2.29	0.50
9:8E:26:VAL:HG13	9:8E:61:ALA:HB3	1.94	0.50
1:13:985:C:H2'	1:13:986:A:C8	2.47	0.50
8:7E:53:VAL:HB	8:7E:58:TYR:CD2	2.46	0.50
1:1G:313:A:H2'	1:1G:314:C:H6	1.77	0.50
26:1H:564:C:H2'	26:1H:565:C:O4'	2.11	0.50
38:98:60:LEU:O	38:98:64:ARG:HG3	2.12	0.50
26:14:1646:C:O3'	58:14:3541:HOH:O	2.20	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:3I:126:LYS:HG3	12:3I:128:ALA:H	1.76	0.50
1:1G:971:G:N2	1:1G:1363:A:OP2	2.34	0.50
26:14:2306:C:H3'	26:14:2307:G:H5''	1.94	0.50
27:16:25:A:H5''	27:16:26:A:OP2	2.12	0.50
26:1H:33:U:H4'	26:1H:34:C:OP1	2.11	0.50
39:A8:5:THR:HG23	39:A8:8:GLU:OE2	2.11	0.50
26:1H:2766:G:N3	26:1H:2766:G:H2'	2.27	0.50
4:32:138:TYR:HD1	4:32:139:ARG:N	2.10	0.50
22:1K:16:U:O2'	22:1K:18:G:OP2	2.28	0.50
28:19:172:TYR:CD1	28:19:186:HIS:HA	2.47	0.50
48:J8:64:ALA:HA	48:J8:67:ILE:HG13	1.93	0.50
40:B8:62:THR:HG22	40:B8:75:ILE:HG12	1.94	0.50
4:3E:90:GLY:HA3	4:3E:204:ILE:HD11	1.92	0.50
27:16:116:G:C5'	39:A8:55:ALA:HB2	2.42	0.50
26:14:270(L):U:O2	33:69:50:ARG:HD2	2.12	0.50
7:6E:27:ILE:HA	7:6E:30:ILE:HD12	1.94	0.50
1:1G:1418:A:H2	26:14:1948:G:N3	2.10	0.50
26:1H:2169:A:H3'	26:1H:2170:A:C8	2.47	0.50
36:78:18:ARG:O	36:78:19:VAL:HG22	2.11	0.49
7:62:95:ARG:HH21	7:62:99:LEU:HD11	1.77	0.49
26:14:531:C:OP1	26:14:561:G:N2	2.45	0.49
26:1H:573:G:O2'	26:1H:574:C:H3'	2.11	0.49
26:14:2113:U:N3	26:14:2114:A:O2'	2.42	0.49
26:14:2262:U:H4'	26:14:2328:A:C2	2.46	0.49
51:M8:38:LYS:H	51:M8:38:LYS:HD2	1.77	0.49
2:1E:226:ARG:HG3	2:1E:227:GLY:N	2.25	0.49
26:14:2639:A:O3'	34:15:97:ARG:NH2	2.36	0.49
26:14:987:G:OP2	58:14:4038:HOH:O	2.20	0.49
10:1A:48:THR:OG1	10:1A:62:HIS:HB3	2.12	0.49
2:12:70:PHE:HB2	2:12:92:TYR:CB	2.42	0.49
40:75:16:ARG:HH21	40:75:19:LEU:HD21	1.77	0.49
40:75:16:ARG:HD3	40:75:79:HIS:HA	1.94	0.49
1:1G:1118:C:OP1	9:82:104:ARG:NH1	2.45	0.49
26:1H:1971:A:N3	28:11:241:PRO:HD3	2.27	0.49
26:14:1071:G:N2	26:14:1087:G:H22	2.10	0.49
28:11:238:GLY:O	28:11:240:ALA:N	2.44	0.49
26:1H:1528:A:C6	26:1H:1545:A:N1	2.80	0.49
4:3E:148:VAL:HG21	4:3E:158:ILE:HG21	1.93	0.49
10:1A:32:ALA:HB2	10:1A:81:THR:HG21	1.94	0.49
27:16:37:C:H2'	27:16:38:C:H5'	1.94	0.49
26:1H:193:U:H5	58:1H:4537:HOH:O	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1152:A:O3'	10:1A:13:HIS:NE2	2.45	0.49
26:1H:557:U:H2'	26:1H:558:G:H8	1.75	0.49
8:7E:103:VAL:HG12	8:7E:104:ARG:HG3	1.92	0.49
37:45:136:ALA:HB1	37:45:138:ASP:OD1	2.11	0.49
26:14:476:G:H4'	26:14:502:A:N1	2.27	0.49
11:2I:13:GLN:HG3	11:2I:75:TYR:O	2.12	0.49
35:25:17:ARG:HB2	35:25:45:GLU:HG2	1.94	0.49
26:1H:1836:C:H2'	26:1H:1837:C:H6	1.78	0.49
42:D8:24:LYS:HA	42:D8:92:THR:HG23	1.92	0.49
1:13:598:U:H4'	8:7E:94:TYR:CD2	2.47	0.49
13:4I:32:GLU:O	13:4I:35:GLU:HG2	2.12	0.49
38:55:59:ASP:OD2	38:55:61:HIS:HB3	2.12	0.49
4:32:31:CYS:O	4:32:31:CYS:SG	2.70	0.49
36:78:18:ARG:HG2	36:78:18:ARG:HH21	1.77	0.49
26:1H:2392:A:P	55:Q8:30:ARG:HH22	2.34	0.49
26:1H:511:U:H5''	26:1H:512:G:OP2	2.12	0.49
1:13:1130:A:O2'	9:8E:3:GLN:NE2	2.45	0.49
26:1H:2699:C:H2'	26:1H:2700:C:O4'	2.12	0.49
26:14:832:G:H5'	36:35:45:LEU:CD1	2.42	0.49
1:13:975:A:H8	1:13:975:A:H5'	1.78	0.49
30:39:102:PRO:HB2	30:39:105:VAL:HG23	1.93	0.49
22:1L:76:A:H62	26:14:2583:G:N2	2.10	0.49
26:14:239:U:H2'	26:14:240:G:O4'	2.12	0.49
29:21:119:ARG:HG2	29:21:160:TYR:HB2	1.94	0.49
1:13:1321:C:H3'	1:13:1322:C:H5''	1.93	0.49
22:1L:14:A:H3'	22:1L:15:G:H8	1.77	0.49
26:1H:1858:G:O2'	26:1H:1884:A:N6	2.45	0.49
26:14:2061:G:H5''	26:14:2503:A:C2	2.46	0.49
26:14:672:C:H5'	30:39:89:VAL:CG1	2.41	0.49
49:K8:15:LYS:HD3	49:K8:67:LYS:HZ1	1.77	0.49
19:AI:63:THR:OG1	19:AI:64:GLU:N	2.45	0.49
26:1H:932:G:H4'	26:1H:933:A:O5'	2.12	0.49
26:1H:2127:G:H2'	26:1H:2128:C:O4'	2.11	0.49
4:32:126:ILE:HG22	4:32:127:THR:N	2.26	0.49
26:14:779:U:OP1	28:19:49:ILE:HG22	2.12	0.49
5:42:147:ASP:O	5:42:151:LEU:HG	2.12	0.49
1:1G:15:G:H4'	5:42:24:ARG:NH1	2.28	0.49
21:1F:9:ARG:NH2	21:1F:23:PRO:HD2	2.26	0.49
40:B8:79:HIS:N	40:B8:79:HIS:CD2	2.79	0.49
26:1H:2273:A:H2'	26:1H:2274:A:C8	2.46	0.49
7:6E:40:ALA:O	7:6E:44:TYR:HB2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:4I:27:LYS:HA	13:4I:31:LYS:NZ	2.27	0.49
1:1G:1151:A:O2'	1:1G:1152:A:O4'	2.30	0.49
26:14:370:G:H4'	26:14:371:A:OP2	2.11	0.49
1:1G:818:G:N2	1:1G:873:A:OP1	2.44	0.49
53:K5:14:THR:O	53:K5:49:HIS:HA	2.11	0.49
26:14:618:G:H2'	26:14:618(A):C:O4'	2.12	0.49
46:H8:44:PHE:CE2	46:H8:86:VAL:HG11	2.47	0.49
26:1H:2864:G:H2'	26:1H:2865:U:C6	2.47	0.49
45:G8:29:GLU:HB3	45:G8:38:ILE:CG2	2.43	0.49
26:14:363(E):U:H5'	26:14:363(F):A:OP2	2.11	0.49
46:D5:99:TYR:CZ	46:D5:125:LEU:HD12	2.46	0.49
26:1H:2577:A:P	58:1H:3800:HOH:O	2.64	0.49
27:16:11:C:O5'	27:16:12:C:H5	1.95	0.49
1:13:1177:G:O2'	1:13:1178:G:O5'	2.28	0.49
1:1G:1298:C:H4'	1:1G:1299:A:C8	2.48	0.49
26:14:1858:G:H2'	26:14:1883:G:N2	2.27	0.49
26:14:639:U:H3	26:14:649:G:H1	1.60	0.49
34:58:95:PRO:O	34:58:96:GLU:CD	2.50	0.49
34:58:46:VAL:HG13	34:58:47:ALA:N	2.27	0.49
1:1G:371:G:H1	1:1G:390:C:N4	2.07	0.49
26:1H:2345:G:H1'	26:1H:2382:G:H5'	1.93	0.49
12:3A:52:LEU:O	12:3A:54:LYS:NZ	2.45	0.49
4:3E:100:ARG:HH12	4:3E:137:SER:HA	1.76	0.49
2:1E:88:ALA:O	2:1E:226:ARG:NH1	2.45	0.49
29:29:8:LYS:HB3	29:29:193:GLY:N	2.27	0.49
26:1H:1265:A:OP2	58:1H:3618:HOH:O	2.18	0.49
1:1G:1014:A:H2'	1:1G:1015:A:C8	2.47	0.49
8:72:99:GLU:OE2	8:72:100:ILE:N	2.33	0.49
36:35:78:PRO:HA	36:35:110:TYR:CE2	2.47	0.49
26:1H:1469:A:H2'	26:1H:1470:G:H8	1.78	0.49
12:3I:60:LEU:HB2	12:3I:64:TYR:HB2	1.93	0.49
26:1H:1348:G:H5''	26:1H:1349:A:OP2	2.12	0.49
48:F5:52:ARG:NH1	48:F5:57:GLU:HB2	2.27	0.49
40:B8:45:PHE:HE2	40:B8:63:VAL:HB	1.77	0.49
1:1G:890:G:O2'	1:1G:906:G:O6	2.20	0.49
42:D8:19:LYS:HG3	42:D8:95:LEU:HD23	1.94	0.49
1:1G:323:U:H5'	20:BA:23:ARG:HB2	1.94	0.49
1:1G:922:G:N3	1:1G:1398:A:H2	2.09	0.49
26:14:297:C:H2'	26:14:298:G:O4'	2.12	0.49
2:12:142:LEU:O	2:12:142:LEU:HD23	2.12	0.49
1:13:1137:C:O2	1:13:1138:G:N2	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:766:A:H2'	1:1G:767:A:O4'	2.13	0.49
26:1H:660:G:H21	36:78:12:ALA:HA	1.77	0.49
3:2E:121:ALA:O	3:2E:125:GLU:HG3	2.12	0.49
1:13:883:C:C2'	1:13:884:U:H5'	2.41	0.49
7:6E:43:PHE:O	7:6E:46:ALA:HB3	2.12	0.49
31:49:145:THR:O	31:49:146:TYR:HB3	2.12	0.49
5:42:16:THR:OG1	5:42:16:THR:O	2.28	0.49
5:4E:26:PHE:CD1	5:4E:26:PHE:N	2.80	0.49
30:31:149:ASP:OD1	30:31:149:ASP:N	2.33	0.49
26:1H:327:G:H2'	26:1H:328:U:C6	2.47	0.49
11:2A:109:VAL:HG13	18:9A:86:VAL:HG13	1.94	0.49
28:19:108:PRO:HG2	28:19:111:LEU:HB2	1.93	0.49
1:1G:676:A:H1'	11:2A:115:PRO:HB3	1.93	0.49
26:1H:2689:U:OP2	26:1H:2719:G:N2	2.33	0.49
26:1H:510:C:H3'	58:1H:4662:HOH:O	2.12	0.49
1:13:1132:C:H2'	1:13:1133:G:H8	1.77	0.49
1:1G:1321:C:H3'	1:1G:1322:C:H5''	1.94	0.49
51:M8:26:SER:OG	51:M8:27:THR:N	2.45	0.49
39:65:106:ARG:O	39:65:106:ARG:HD2	2.13	0.49
26:14:2116:G:N3	26:14:2117:A:N6	2.60	0.49
26:1H:1899:G:N2	26:1H:1902:C:C5	2.65	0.49
24:3K:64:A:H8	24:3K:64:A:O5'	1.95	0.49
26:1H:140:A:H8	26:1H:1408:C:HO2'	1.60	0.49
1:13:1301:U:O2'	1:13:1302:U:H5'	2.12	0.49
26:14:1945:G:C4	26:14:1946:U:C5	3.01	0.49
1:1G:1367:C:H5'	10:1A:60:ARG:NH2	2.28	0.49
26:14:1952:A:C5	35:25:22:ILE:HD11	2.47	0.49
1:13:1154:G:H2'	1:13:1155:G:H8	1.78	0.49
26:14:289:A:H5'	26:14:290:G:OP2	2.12	0.49
26:1H:1568:G:H5''	28:11:61:LEU:HD22	1.94	0.49
26:1H:1965:C:H3'	26:1H:1966:A:H2'	1.95	0.49
26:1H:2405:G:OP1	36:78:77:ARG:NH2	2.45	0.49
26:14:2852:G:H2'	26:14:2853:C:C6	2.46	0.49
4:3E:173:TRP:CD1	4:3E:174:LEU:HG	2.47	0.49
4:3E:190:ASP:HB2	4:3E:193:ASP:OD1	2.13	0.49
26:14:118:A:N3	26:14:178:G:H1'	2.28	0.49
8:7E:21:LYS:O	8:7E:63:LEU:HD23	2.13	0.49
26:14:303:U:H2'	26:14:304:G:H8	1.77	0.49
29:29:11:MET:SD	29:29:24:THR:HG22	2.52	0.49
3:2E:11:ARG:HE	3:2E:180:ALA:HB3	1.77	0.49
26:14:2674:G:H4'	35:25:30:ALA:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:722:A:C8	1:1G:724:G:H1'	2.47	0.49
1:13:188:U:O2'	1:13:189:U:H5'	2.12	0.49
3:2E:156:ARG:HG2	3:2E:163:ALA:HB2	1.93	0.49
1:13:739:C:O2'	15:6I:42:HIS:ND1	2.29	0.49
26:1H:1007:C:OP1	34:58:35:ARG:NH1	2.42	0.49
26:1H:2058:A:N6	58:1H:3619:HOH:O	2.42	0.49
34:58:68:GLU:HG3	34:58:88:GLU:OE2	2.12	0.49
28:19:124:PRO:HG2	28:19:129:ASN:ND2	2.28	0.49
26:14:279:C:N4	26:14:361:G:H1	2.10	0.49
26:14:1204:A:H2	26:14:1241:A:N1	2.11	0.49
17:8I:18:THR:HG21	17:8I:69:LYS:HD2	1.95	0.49
1:1G:827:U:H5'	1:1G:828:A:OP2	2.13	0.49
1:13:524:G:H2'	1:13:525:C:C5	2.47	0.49
1:1G:1300:G:O2'	1:1G:1301:U:O5'	2.25	0.49
26:14:1678:G:N3	26:14:1678:G:H2'	2.28	0.49
23:2K:63:C:O2	23:2K:64:G:C8	2.65	0.49
43:E8:37:ARG:HH22	52:N8:47:PRO:CB	2.25	0.49
41:85:28:ARG:HH11	41:85:38:THR:HG1	1.53	0.49
26:14:2301:C:H6	26:14:2301:C:H3'	1.77	0.49
1:1G:1368:G:OP2	9:82:112:LYS:HD2	2.11	0.49
26:1H:518:G:H5'	43:E8:18:ARG:HH21	1.77	0.49
26:14:2271:G:OP1	47:E5:18:ALA:HB1	2.12	0.49
31:41:37:VAL:HG22	31:41:159:VAL:HG12	1.93	0.49
37:88:21:THR:HG21	37:88:101:ARG:N	2.27	0.49
4:32:150:GLU:C	4:32:152:SER:H	2.16	0.49
1:13:575:G:C5	1:13:881:G:C2	3.00	0.49
26:1H:579:G:H2'	26:1H:580:C:C6	2.47	0.49
2:12:221:LEU:HA	2:12:224:GLN:HB2	1.94	0.49
52:N8:49:CYS:SG	52:N8:60:VAL:HB	2.53	0.49
1:13:244:U:H4'	1:13:245:C:O5'	2.12	0.49
1:1G:1108:G:OP1	3:22:175:LEU:HD12	2.13	0.49
26:14:1669:A:H5''	26:14:1670:C:OP2	2.11	0.49
1:1G:617:G:OP2	58:1G:1842:HOH:O	2.18	0.49
46:H8:28:MET:HG3	46:H8:37:VAL:HG11	1.94	0.49
41:85:8:VAL:O	41:85:12:ARG:HG3	2.13	0.49
26:1H:805:G:OP2	36:78:41:ARG:HG2	2.13	0.49
26:1H:806:C:OP2	36:78:41:ARG:HD3	2.12	0.49
3:2E:8:ILE:HD13	3:2E:184:TYR:HB3	1.94	0.49
32:51:118:PRO:HD2	32:51:121:ILE:HB	1.93	0.49
26:1H:2243:U:H2'	26:1H:2244:U:C6	2.47	0.49
7:6E:45:ASP:O	7:6E:49:ILE:HG12	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:278:A:H8	26:14:278:A:OP2	1.96	0.49
26:14:278:A:HO2'	26:14:279:C:H5	1.60	0.49
26:14:192:C:P	58:14:4102:HOH:O	2.70	0.49
41:C8:75:ASN:HB3	41:C8:77:SER:H	1.77	0.49
51:I5:43:TYR:CG	51:I5:43:TYR:O	2.66	0.49
39:65:5:THR:O	39:65:8:GLU:N	2.44	0.49
26:1H:29:U:H2'	26:1H:30:G:C8	2.47	0.49
26:14:1993:U:H4'	29:29:128:SER:HB3	1.95	0.49
28:19:29:PRO:HA	28:19:83:GLU:OE1	2.12	0.49
1:1G:419:C:H5'	1:1G:420:U:OP2	2.13	0.49
43:A5:20:VAL:CG2	43:A5:47:VAL:HG21	2.43	0.49
26:1H:309:G:N3	26:1H:329:G:O2'	2.41	0.49
31:49:125:PHE:HB3	31:49:166:ASP:HB2	1.94	0.49
9:8E:18:PHE:HB2	9:8E:62:TYR:HB3	1.93	0.49
2:1E:47:THR:HG22	2:1E:51:LEU:HD12	1.93	0.49
55:Q8:48:PHE:CZ	55:Q8:49:VAL:HG13	2.47	0.49
26:14:1019:U:HO2'	26:14:1021:A:H2	1.59	0.49
26:14:662:G:H5'	36:35:15:ARG:CA	2.36	0.49
24:3L:53:G:H2'	24:3L:54:U:C6	2.48	0.49
1:1G:565:U:OP2	1:1G:566:G:O2'	2.23	0.49
26:1H:1668:A:N6	26:1H:1676:A:H61	2.11	0.49
40:75:4:GLY:O	40:75:6:LEU:N	2.45	0.49
40:B8:108:ARG:HA	40:B8:111:ARG:CZ	2.42	0.49
27:16:7:G:H4'	39:A8:29:PHE:HD2	1.73	0.49
1:1G:1127:G:C2'	1:1G:1147:C:H42	2.26	0.49
26:14:2716:U:O2'	26:14:2717:G:H5'	2.12	0.49
38:98:77:ARG:O	38:98:80:PHE:N	2.46	0.49
26:14:17:G:H2'	26:14:18:C:C6	2.48	0.49
40:75:64:ARG:CB	40:75:73:GLU:HG2	2.42	0.49
19:AI:41:VAL:HG11	19:AI:67:VAL:HA	1.93	0.49
19:AI:64:GLU:O	19:AI:67:VAL:HG13	2.12	0.49
20:BA:45:GLN:HB2	20:BA:91:LEU:HD13	1.94	0.49
4:32:149:ALA:O	4:32:153:ARG:HG2	2.12	0.49
4:32:173:TRP:HB3	4:32:187:ARG:HH11	1.76	0.49
9:82:33:PHE:CE2	9:82:47:LEU:HD23	2.47	0.49
26:14:2019:A:N7	52:J5:9:LYS:HD2	2.27	0.49
50:H5:8:LEU:HD22	50:H5:10:LYS:O	2.13	0.49
26:1H:303:U:H2'	26:1H:304:G:C8	2.46	0.49
26:14:30:G:O6	58:14:3893:HOH:O	2.18	0.49
34:15:15:LEU:HD13	34:15:16:ILE:N	2.28	0.49
37:45:29:PHE:HB3	37:45:65:PHE:CE2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2439:A:H5'	26:14:2439:A:H8	1.77	0.49
28:19:43:ARG:HG2	28:19:43:ARG:HH11	1.77	0.49
26:1H:784:A:OP1	26:1H:2588:G:H5''	2.13	0.49
10:1A:13:HIS:O	10:1A:17:ASP:HB2	2.12	0.49
26:1H:459:U:H2'	26:1H:460:A:C8	2.46	0.49
26:14:2027:G:H2'	26:14:2028:U:O4'	2.13	0.49
50:H5:9:VAL:HG11	50:H5:55:ARG:HB2	1.94	0.49
31:41:139:LEU:HD22	31:41:146:TYR:HD1	1.76	0.49
46:H8:31:ARG:HD2	46:H8:32:HIS:CE1	2.47	0.49
26:14:5:A:H2'	26:14:6:A:C8	2.48	0.49
1:13:953:G:H2'	1:13:954:G:O4'	2.13	0.49
26:1H:1754:C:H5	40:B8:96:ARG:CZ	2.25	0.49
3:22:73:PRO:HG3	3:22:105:GLU:HG3	1.93	0.49
26:14:1217:C:OP1	41:85:15:LYS:NZ	2.37	0.49
1:13:1147:C:O2	9:8E:16:ARG:NH1	2.45	0.49
37:88:17:LEU:HD21	37:88:96:VAL:HG13	1.95	0.49
1:1G:229:U:H2'	1:1G:230:G:H8	1.78	0.49
15:6A:77:ARG:HA	15:6A:80:ALA:HB3	1.94	0.49
26:14:1441:G:H2'	26:14:1442:G:H8	1.78	0.49
38:55:51:LEU:HD22	38:55:66:VAL:HG13	1.93	0.49
1:1G:408:A:H2'	1:1G:409:G:O4'	2.12	0.49
26:14:2593:U:H2'	26:14:2594:C:C6	2.48	0.49
26:14:1771:C:H1'	26:14:1786:A:C8	2.47	0.49
26:1H:1327:C:H2'	26:1H:1328:G:O4'	2.13	0.49
26:14:1462:C:H4'	26:14:2703:C:H5'	1.94	0.49
1:13:1366:C:H2'	1:13:1367:C:C6	2.47	0.49
1:13:1034:G:N2	1:13:1035:A:N7	2.60	0.49
26:1H:250:G:OP1	36:78:60:MET:HE1	2.13	0.49
1:1G:1129:C:H41	1:1G:1141:C:N4	2.03	0.49
26:14:831:G:N7	58:14:3673:HOH:O	2.34	0.49
1:1G:652:U:H2'	1:1G:653:A:H5''	1.95	0.49
26:14:389:G:H1	36:35:71:VAL:HG12	1.75	0.49
38:98:53:HIS:ND1	38:98:94:TYR:OH	2.34	0.49
19:AA:23:ASN:HA	19:AA:27:GLU:HG3	1.95	0.49
26:1H:1798:U:H5''	28:11:259:THR:HG22	1.94	0.49
1:13:232:G:H2'	1:13:233:C:H6	1.78	0.49
26:1H:1239:G:H2'	26:1H:1240:U:O4'	2.12	0.49
26:1H:57:C:H2'	26:1H:58:G:O4'	2.13	0.49
23:2L:24:C:C2	23:2L:25:U:C5	3.00	0.49
1:13:142:G:C2	1:13:143:A:C5	3.00	0.49
1:13:457:C:H2'	1:13:458:C:H6	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:581:C:H2'	26:1H:582:G:C8	2.47	0.49
34:15:38:HIS:CD2	34:15:39:ARG:H	2.30	0.49
1:13:619:U:O2	4:3E:135:LEU:HD22	2.12	0.49
26:14:1015:G:O2'	26:14:1016:G:H5'	2.12	0.49
5:42:60:TYR:CB	5:42:64:ARG:HH21	2.26	0.49
38:55:107:ASP:C	38:55:107:ASP:OD1	2.50	0.49
45:G8:39:VAL:HB	45:G8:42:VAL:HG13	1.93	0.49
26:1H:1403:C:H5''	26:1H:1471:A:H1'	1.95	0.49
26:14:208:C:H2'	26:14:209:C:C6	2.47	0.49
29:29:120:TRP:CD1	29:29:155:LYS:HB3	2.48	0.49
27:16:3:C:H2'	27:16:4:C:C6	2.48	0.49
20:BA:11:SER:HA	20:BA:13:LEU:H	1.77	0.49
33:69:120:ILE:HG22	33:69:122:GLU:H	1.78	0.49
1:13:792:A:H4'	1:13:793:U:O5'	2.12	0.49
32:59:30:LYS:HE2	32:59:136:ILE:HG22	1.95	0.49
26:1H:459:U:H2'	26:1H:460:A:H8	1.77	0.49
26:1H:311:A:C6	26:1H:328:U:C4	3.01	0.49
26:1H:30:G:H2'	26:1H:31:C:C6	2.48	0.49
26:14:1826:G:H4'	28:19:242:ARG:NH1	2.28	0.49
32:51:20:ALA:HB1	32:51:21:PRO:HD2	1.94	0.49
35:25:24:VAL:HB	35:25:33:ALA:HB2	1.93	0.49
53:O8:47:THR:HG22	53:O8:48:VAL:H	1.77	0.49
40:B8:100:TYR:HB3	40:B8:103:ARG:NH1	2.26	0.49
26:1H:1680:U:O2	26:1H:1763:G:H3'	2.12	0.49
26:14:146:G:H2'	26:14:147:U:O4'	2.12	0.49
46:D5:128:VAL:HG22	46:D5:129:SER:H	1.77	0.49
26:14:2629:A:N3	26:14:2629:A:H2'	2.28	0.49
46:D5:11:GLU:CG	46:D5:12:GLY:H	2.25	0.49
26:1H:731:C:P	58:1H:3689:HOH:O	2.70	0.49
55:Q8:6:THR:OG1	55:Q8:6:THR:O	2.23	0.49
1:13:1129:C:N4	1:13:1139:G:H1	2.11	0.49
31:41:21:ARG:NH1	31:41:21:ARG:HG2	2.16	0.49
26:14:1161:C:H2'	26:14:1162:G:C8	2.48	0.49
7:62:92:SER:HB3	7:62:94:ARG:HG3	1.93	0.49
49:G5:25:VAL:HA	49:G5:28:LYS:HB2	1.94	0.49
26:14:1889:A:O2'	26:14:2087:G:H5'	2.13	0.49
1:13:413:G:H22	1:13:428:G:H1'	1.77	0.49
26:14:1331:A:O2'	26:14:1332:G:H8	1.95	0.49
45:G8:49:VAL:HG21	45:G8:55:TYR:HE2	1.78	0.49
26:1H:654(A):A:C2	26:1H:654(T):A:N1	2.81	0.49
32:51:154:PRO:HB3	32:51:163:TYR:CZ	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:324:G:N1	1:13:327:A:OP2	2.44	0.49
1:1G:1376:U:H2'	1:1G:1377:A:C8	2.47	0.49
39:A8:63:THR:O	39:A8:66:ALA:HB3	2.12	0.49
1:1G:735:C:H2'	1:1G:736:C:H6	1.78	0.49
1:13:947:G:H2'	1:13:948:C:O4'	2.13	0.49
1:1G:952:U:H4'	1:1G:964:A:N1	2.28	0.49
1:1G:599:C:H2'	1:1G:600:C:C6	2.47	0.49
26:1H:530:G:C5	26:1H:2022:U:H5''	2.48	0.49
26:14:528:A:C2	26:14:2042:A:H2'	2.47	0.49
36:78:113:LYS:HA	36:78:129:ALA:O	2.12	0.49
37:45:41:TRP:HZ3	37:45:74:TYR:HE1	1.59	0.49
26:14:1439:A:C2	26:14:1553:A:C5	3.00	0.49
26:1H:1903:G:C2'	26:1H:1904:G:H5'	2.43	0.49
8:72:44:PHE:HD1	8:72:80:ILE:HG12	1.77	0.49
7:6E:94:ARG:O	7:6E:97:GLN:HB3	2.13	0.49
28:11:27:THR:HG23	28:11:28:GLU:HG2	1.93	0.49
26:1H:886:C:H5'	26:1H:887:A:OP2	2.13	0.49
28:19:79:VAL:HG12	28:19:113:VAL:HA	1.95	0.49
26:14:27:G:O2'	26:14:28:A:OP2	2.27	0.49
13:4A:27:LYS:HD2	13:4A:31:LYS:HE3	1.93	0.49
1:1G:409:G:OP1	4:32:25:ARG:HB2	2.13	0.49
26:1H:1189:A:OP2	58:1H:3874:HOH:O	2.19	0.49
26:1H:2577:A:H5''	26:1H:2578:G:H5'	1.95	0.49
45:C5:97:ARG:HG2	45:C5:102:CYS:O	2.13	0.49
26:1H:1328:G:H2'	26:1H:1330:C:C5	2.48	0.49
26:14:1204:A:C2	26:14:1241:A:N1	2.80	0.49
1:13:1128:C:C6	1:13:1139:G:C6	3.00	0.49
46:H8:4:ARG:HD3	46:H8:60:GLU:OE2	2.12	0.49
26:14:1024:G:C3'	26:14:1025:G:H5''	2.41	0.49
26:1H:445:C:O2'	26:1H:446:G:H5'	2.13	0.49
31:49:104:GLU:HG2	51:I5:23:GLU:HG3	1.93	0.49
2:1E:69:LEU:HD22	2:1E:91:PRO:HB2	1.95	0.49
5:4E:100:VAL:HG22	5:4E:115:VAL:HG12	1.95	0.49
26:14:273(D):C:N4	26:14:363(B):G:H1	2.10	0.49
4:3E:9:CYS:O	4:3E:13:ARG:HG2	2.13	0.49
29:29:176:ILE:HB	29:29:181:LEU:HB2	1.95	0.49
1:13:1285:A:H4'	1:13:1286:A:C5'	2.43	0.49
26:14:2128:C:H42	26:14:2160:G:H1	1.60	0.49
26:14:1322:A:N1	26:14:1333:C:O2'	2.43	0.49
26:1H:1678:G:N2	26:1H:1989:G:N2	2.60	0.49
26:14:2575:C:H5'	29:29:143:ASN:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:115:G:H1'	1:1G:116:A:N7	2.27	0.49
26:14:2080:G:O2'	26:14:2081:C:H5'	2.12	0.49
22:1K:14:A:C6	22:1K:22:G:N3	2.81	0.49
37:45:118:LEU:HD12	37:45:131:ILE:HG23	1.93	0.49
37:45:103:MET:O	37:45:104:PHE:HB2	2.12	0.49
40:75:53:ARG:NH1	40:75:60:THR:HG23	2.27	0.49
1:13:725:G:H2'	1:13:726:C:H6	1.78	0.49
26:1H:1568:G:P	28:11:63:ARG:HH12	2.36	0.49
26:14:600:G:N2	26:14:605:C:O3'	2.46	0.49
30:39:66:PRO:O	30:39:67:GLN:HB3	2.13	0.49
4:3E:72:GLU:OE1	4:3E:207:TYR:OH	2.27	0.49
27:1J:50:G:OP1	39:65:62:LYS:HB2	2.13	0.49
30:39:146:ALA:CB	30:39:148:LEU:HG	2.42	0.49
38:55:57:ARG:NE	38:55:59:ASP:OD1	2.38	0.49
46:D5:99:TYR:HA	46:D5:124:ILE:O	2.12	0.49
26:1H:2324:C:H5''	26:1H:2325:G:H5''	1.95	0.49
13:4I:81:LEU:HD22	13:4I:88:ARG:HB3	1.94	0.49
20:BI:53:LEU:HA	20:BI:56:MET:HB3	1.94	0.49
22:1K:10:G:H2'	22:1K:11:C:C6	2.48	0.49
1:1G:490:G:OP2	4:32:132:ARG:NH2	2.23	0.49
16:7I:71:ARG:O	16:7I:75:ARG:N	2.46	0.49
48:J8:81:LYS:N	48:J8:81:LYS:HD2	2.28	0.49
28:11:6:PHE:HE1	28:11:18:VAL:HG23	1.78	0.49
46:H8:127:LYS:NZ	46:H8:162:GLU:OE2	2.42	0.49
26:14:2378:A:O2'	39:65:21:THR:HG21	2.13	0.49
41:85:92:ARG:C	41:85:94:ASN:H	2.16	0.49
46:H8:7:ALA:HB3	46:H8:61:LEU:HB3	1.93	0.49
26:14:2701:C:H3'	26:14:2702:U:C5'	2.37	0.49
27:1J:13:A:H5''	27:1J:15:A:N6	2.28	0.49
11:2A:22:HIS:HB3	11:2A:29:ILE:HG12	1.95	0.49
26:14:639:U:H2'	26:14:640:C:C6	2.47	0.49
33:61:110:ASP:OD1	33:61:111:PRO:HA	2.12	0.49
11:2I:34:ASP:HB2	11:2I:35:PRO:HD2	1.94	0.49
1:1G:1211:U:H1'	1:1G:1213:A:C2	2.48	0.49
19:AA:20:LEU:HA	19:AA:23:ASN:HD22	1.77	0.49
1:1G:373:A:C2	1:1G:374:A:C8	3.01	0.49
49:K8:18:PRO:O	49:K8:22:GLU:HG3	2.12	0.49
12:3I:43:VAL:HG23	12:3I:93:LEU:HD22	1.95	0.49
1:1G:302:G:O2'	1:1G:556:C:H5''	2.13	0.49
26:1H:70:G:H21	26:1H:71:A:N6	2.11	0.49
47:E5:53:MET:HG3	47:E5:59:LEU:CD2	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:16:13:A:N6	27:16:70:C:H5'	2.27	0.49
1:13:511:C:C2	1:13:512:U:C5	3.00	0.49
26:1H:184:C:H2'	26:1H:185:U:H6	1.78	0.49
26:1H:2460:U:H2'	26:1H:2461:C:H6	1.78	0.49
1:13:296:U:H2'	1:13:297:G:C8	2.48	0.49
37:88:20:ALA:HB2	46:H8:79:ARG:HG2	1.93	0.49
31:49:103:LEU:HD22	31:49:178:PHE:CZ	2.48	0.49
37:45:38:GLU:HB2	37:45:127:ILE:HG22	1.94	0.49
8:72:20:TYR:CE2	8:72:75:ARG:HB3	2.48	0.49
26:1H:1443:G:C2	26:1H:1549:C:N3	2.81	0.49
22:1K:5:G:H1'	22:1K:69:G:N2	2.26	0.49
1:1G:1375:A:H4'	7:62:29:LYS:HE3	1.95	0.49
2:1E:73:THR:OG1	2:1E:170:GLU:OE2	2.11	0.49
26:1H:2780:G:OP1	34:58:118:LYS:NZ	2.30	0.49
35:68:112:MET:HA	35:68:115:VAL:HG13	1.93	0.49
26:14:2185:C:H2'	26:14:2186:G:C8	2.47	0.49
17:8A:63:ARG:HG2	17:8A:64:PRO:HD2	1.95	0.49
26:1H:2705:A:O2'	26:1H:2852:G:OP1	2.23	0.48
26:1H:2319:G:N1	26:1H:2334:G:OP2	2.46	0.48
1:13:1126:U:H3	1:13:1127:G:N2	2.11	0.48
30:39:120:GLU:HG3	30:39:122:LYS:NZ	2.23	0.48
27:1J:40:U:H1'	27:1J:46:A:N1	2.28	0.48
1:13:745:C:OP1	1:13:851:G:O2'	2.30	0.48
1:13:963:G:C2	10:1I:55:LYS:NZ	2.81	0.48
26:1H:1170:G:N2	26:1H:1180:C:C2	2.81	0.48
40:75:4:GLY:HA2	40:75:8:LYS:NZ	2.26	0.48
49:G5:37:PHE:O	49:G5:41:ILE:HG12	2.13	0.48
1:13:232:G:H2'	1:13:233:C:C6	2.47	0.48
26:14:666:G:H5''	36:35:47:ASP:O	2.13	0.48
9:82:102:LEU:O	9:82:103:THR:OG1	2.28	0.48
29:29:33:VAL:HG11	29:29:36:ARG:NH2	2.27	0.48
36:78:111:ARG:HG2	36:78:128:HIS:CD2	2.48	0.48
29:21:60:ASN:OD1	29:21:62:PRO:HD2	2.13	0.48
50:L8:6:VAL:HB	50:L8:54:VAL:HG21	1.94	0.48
28:11:12:SER:O	28:11:16:MET:HB2	2.12	0.48
4:3E:148:VAL:HG12	4:3E:149:ALA:O	2.13	0.48
4:3E:155:LEU:O	4:3E:157:LEU:N	2.46	0.48
26:1H:321:G:O3'	30:31:168:ARG:NH2	2.44	0.48
26:1H:322:A:P	30:31:168:ARG:HH21	2.35	0.48
4:3E:108:LEU:HD11	4:3E:174:LEU:HD22	1.94	0.48
38:55:24:GLN:OE1	38:55:36:THR:HG21	2.11	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:11:G:C2'	26:1H:12:U:H5'	2.44	0.48
4:32:138:TYR:CD1	4:32:139:ARG:N	2.81	0.48
38:55:32:GLY:HA2	38:55:116:LEU:CD1	2.43	0.48
8:72:17:THR:O	8:72:78:GLN:NE2	2.46	0.48
27:1J:10:C:C4	27:1J:11:C:C5	3.01	0.48
26:14:2412:A:H2'	26:14:2413:G:O4'	2.13	0.48
26:1H:2406:U:H3'	58:1H:3728:HOH:O	2.12	0.48
35:25:105:GLU:HA	35:25:108:GLU:OE1	2.13	0.48
26:1H:1213:A:H1'	26:1H:1238:G:N3	2.27	0.48
26:14:861:A:N3	27:1J:79:C:O2'	2.46	0.48
26:14:2569:G:C2	26:14:2570:G:C8	3.01	0.48
1:1G:1346:A:OP2	1:1G:1346:A:H3'	2.13	0.48
36:78:91:PHE:CD1	36:78:91:PHE:N	2.81	0.48
26:1H:449:A:OP2	58:1H:3897:HOH:O	2.20	0.48
26:1H:1771:C:OP1	58:1H:3983:HOH:O	2.20	0.48
26:14:320:A:H4'	26:14:322:A:C8	2.48	0.48
26:14:1043:C:H42	26:14:1112:G:H1	1.60	0.48
26:14:2873:A:C8	38:55:5:LYS:HA	2.48	0.48
26:1H:2401:U:H2'	26:1H:2402:C:C6	2.49	0.48
26:14:2394:C:OP1	36:35:63:PRO:HG2	2.13	0.48
26:14:1423:G:OP1	26:14:1492:G:O2'	2.31	0.48
51:I5:22:ILE:HD13	51:I5:22:ILE:H	1.76	0.48
23:2K:52:C:H2'	23:2K:53:G:O4'	2.13	0.48
1:1G:1249:C:O2'	9:82:73:GLN:OE1	2.31	0.48
1:1G:1028(A):C:H42	1:1G:1032(B):G:H22	1.61	0.48
11:2I:34:ASP:OD1	11:2I:36:ASP:N	2.40	0.48
11:2I:79:SER:HB2	11:2I:106:LYS:HD2	1.94	0.48
49:G5:33:MET:O	49:G5:37:PHE:HD1	1.95	0.48
26:14:2131:G:OP1	26:14:2132:U:H3'	2.13	0.48
26:1H:2694:G:C4	26:1H:2695:C:C5	3.01	0.48
6:5E:23:LYS:HB2	6:5E:23:LYS:NZ	2.28	0.48
38:98:91:GLN:CD	38:98:91:GLN:H	2.15	0.48
26:1H:851:U:OP1	50:L8:49:LYS:NZ	2.45	0.48
1:13:55:A:C5	1:13:56:U:C5	3.01	0.48
26:14:1210:A:H5'	26:14:1212:G:O4'	2.13	0.48
26:14:39:C:H2'	26:14:40:C:C6	2.49	0.48
1:13:1203:C:H2'	1:13:1204:A:O4'	2.13	0.48
1:13:648:A:H2'	1:13:649:G:C8	2.48	0.48
29:29:31:CYS:O	29:29:91:VAL:HG22	2.14	0.48
1:13:244:U:H4'	1:13:245:C:C5'	2.43	0.48
12:3A:27:LEU:HD21	12:3A:61:THR:OG1	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:524:U:H2'	26:14:525:U:C6	2.47	0.48
35:68:85:VAL:HG11	35:68:114:ILE:HD13	1.94	0.48
26:1H:2627:G:N2	26:1H:2777:G:OP2	2.46	0.48
38:98:13:HIS:ND1	38:98:15:SER:HB3	2.28	0.48
3:2E:58:GLU:HB2	3:2E:65:ALA:CB	2.43	0.48
32:59:149:ARG:HG2	32:59:162:ILE:HG22	1.95	0.48
26:14:1536:A:C8	26:14:1537:C:H1'	2.48	0.48
26:1H:658:C:H2'	26:1H:659:C:C6	2.48	0.48
26:1H:324:A:C2'	26:1H:325:G:H5'	2.42	0.48
30:31:123:LEU:HD12	30:31:124:LEU:N	2.28	0.48
26:14:5:A:H2'	26:14:6:A:N7	2.27	0.48
7:62:111:ARG:NH2	7:62:122:HIS:HB3	2.27	0.48
1:13:1478:C:H2'	1:13:1479:C:C6	2.48	0.48
26:1H:1268:A:H2'	26:1H:1269:A:O4'	2.13	0.48
24:3K:66:U:H2'	24:3K:67:C:C6	2.48	0.48
2:1E:76:GLN:HA	2:1E:208:ILE:HG12	1.95	0.48
45:C5:63:LYS:HA	45:C5:63:LYS:NZ	2.28	0.48
4:3E:53:ASP:HB3	4:3E:57:ARG:NH1	2.28	0.48
30:31:138:GLU:O	30:31:141:ALA:HB3	2.13	0.48
34:15:96:GLU:CD	34:15:96:GLU:H	2.17	0.48
26:14:1425:G:H2'	26:14:1426:G:C8	2.47	0.48
30:31:179:GLU:H	30:31:179:GLU:CD	2.16	0.48
40:B8:78:LEU:O	40:B8:78:LEU:HD13	2.12	0.48
1:13:38:G:C2	1:13:397:A:C2	3.01	0.48
26:1H:2689:U:H5''	26:1H:2713:A:C2	2.49	0.48
55:Q8:47:LYS:NZ	55:Q8:47:LYS:HB2	2.27	0.48
26:14:568:U:H5''	26:14:568:U:H6	1.79	0.48
1:13:600:C:H2'	1:13:601:C:C6	2.48	0.48
2:1E:163:PHE:HD1	2:1E:185:ILE:HG13	1.78	0.48
32:59:3:ARG:HD2	32:59:4:ILE:HG13	1.95	0.48
42:95:85:LYS:CB	42:95:87:HIS:H	2.27	0.48
1:13:255:G:H1'	17:8I:16:GLN:OE1	2.13	0.48
26:14:751:A:P	58:14:3410:HOH:O	2.71	0.48
26:14:248:G:H5'	26:14:250:G:N7	2.27	0.48
1:1G:1141:C:H2'	1:1G:1142:G:C8	2.38	0.48
27:1J:46:A:H2'	27:1J:47:C:C6	2.48	0.48
26:14:94:G:OP1	45:C5:54:LYS:NZ	2.32	0.48
26:1H:1593:G:H2'	26:1H:1594:G:H8	1.79	0.48
26:1H:6:A:H4'	34:58:129:PRO:HB3	1.94	0.48
26:1H:2371:G:C4'	53:O8:45:LYS:HG3	2.43	0.48
39:A8:38:GLN:HG3	39:A8:47:THR:HG21	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:E8:72:LYS:N	43:E8:106:ILE:O	2.46	0.48
1:13:113:G:H2'	1:13:114:U:C6	2.47	0.48
37:88:66:ILE:O	37:88:67:ARG:HB2	2.14	0.48
26:1H:69:C:H2'	26:1H:70:G:H8	1.78	0.48
13:4I:3:ARG:CZ	13:4I:7:VAL:HG13	2.43	0.48
32:51:12:PRO:HD3	32:51:48:GLY:O	2.12	0.48
1:1G:1006:C:N3	1:1G:1007:C:N4	2.61	0.48
14:5I:6:LEU:HB3	14:5I:23:ARG:HH22	1.79	0.48
26:1H:1432:C:H2'	26:1H:1433:U:O4'	2.11	0.48
26:14:1533:C:H42	26:14:1538:G:H1	1.60	0.48
29:29:11:MET:HE3	29:29:187:ALA:H	1.78	0.48
46:D5:19:ARG:NH1	46:D5:84:GLU:O	2.45	0.48
26:1H:1028:A:N3	26:1H:2486:G:O2'	2.37	0.48
10:1A:38:ILE:N	10:1A:71:LEU:O	2.37	0.48
6:5E:33:TYR:CD2	6:5E:75:LEU:HD23	2.48	0.48
4:3E:19:LEU:HD23	4:3E:21:LEU:HD21	1.95	0.48
49:K8:42:GLY:O	49:K8:44:LEU:N	2.46	0.48
27:1J:12:C:O2'	47:E5:74:ARG:HG2	2.12	0.48
40:75:51:ARG:HB3	40:75:62:THR:HG23	1.94	0.48
1:13:590:C:H2'	1:13:591:U:H6	1.78	0.48
26:14:2335:A:C8	26:14:2337:G:C5	3.01	0.48
26:14:1127:A:N3	26:14:2518:A:H5'	2.28	0.48
41:C8:58:ARG:HA	41:C8:61:TRP:CE3	2.48	0.48
46:D5:103:ARG:HE	46:D5:136:PHE:HB3	1.78	0.48
44:B5:57:LEU:N	44:B5:57:LEU:HD23	2.28	0.48
26:1H:2721:A:H2'	26:1H:2722:G:O4'	2.13	0.48
26:1H:1427:A:C2	28:11:31:LYS:HE2	2.48	0.48
45:G8:44:ILE:HA	45:G8:64:GLU:HA	1.93	0.48
18:9I:47:THR:O	18:9I:83:GLU:N	2.37	0.48
26:1H:1705:G:C6	26:1H:1706:U:C4	3.01	0.48
1:1G:1362:C:H2'	1:1G:1362(A):C:H5''	1.96	0.48
26:14:2318:G:H5'	26:14:2319:G:OP2	2.13	0.48
26:14:2275:C:C5'	26:14:2275:C:H6	2.25	0.48
26:14:2286:A:H8	26:14:2287:A:N6	2.12	0.48
26:14:1641:A:H5''	26:14:1642:G:OP2	2.12	0.48
47:E5:21:LEU:HD11	47:E5:41:ARG:NH1	2.28	0.48
1:1G:1277:C:O2'	1:1G:1279:A:H8	1.95	0.48
26:1H:141:A:H1'	26:1H:1408:C:O4'	2.14	0.48
26:1H:2213:U:O4'	48:J8:52:ARG:NH2	2.46	0.48
45:G8:53:PRO:O	45:G8:54:LYS:NZ	2.40	0.48
26:14:821:A:O2'	26:14:946:G:OP2	2.21	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2789:C:H1'	26:1H:2892:A:H2	1.78	0.48
32:51:149:ARG:HG3	32:51:162:ILE:HG22	1.95	0.48
26:1H:654(H):G:H2'	26:1H:654(H):G:N3	2.28	0.48
3:22:14:ILE:HG12	3:22:15:THR:N	2.29	0.48
2:12:208:ILE:HA	2:12:211:ILE:HD12	1.94	0.48
2:1E:19:HIS:HB3	2:1E:189:ASP:OD2	2.14	0.48
1:1G:433:C:H2'	1:1G:434:U:C6	2.49	0.48
1:13:1485:U:O2'	1:13:1486:G:H5'	2.13	0.48
1:1G:316:G:OP2	1:1G:351:G:O2'	2.31	0.48
6:52:25:ILE:HD13	6:52:82:ARG:CD	2.43	0.48
29:29:111:ARG:HB2	29:29:160:TYR:O	2.13	0.48
1:1G:1489:G:H2'	1:1G:1490:C:O4'	2.13	0.48
1:13:806:C:H2'	1:13:807:A:C8	2.48	0.48
26:14:1344:G:O2'	26:14:1385:G:H2'	2.13	0.48
6:5E:82:ARG:HB2	6:5E:85:VAL:HG23	1.95	0.48
26:1H:1301:A:H2'	26:1H:1301:A:N3	2.29	0.48
1:13:1103:C:H2'	1:13:1104:G:O4'	2.13	0.48
48:J8:25:LYS:HB3	48:J8:25:LYS:NZ	2.28	0.48
7:62:18:TYR:HB3	7:62:59:LEU:HD12	1.94	0.48
26:1H:2860:A:C8	26:1H:2861:G:H1'	2.48	0.48
26:14:587:C:O2	36:35:33:ARG:NH1	2.47	0.48
9:82:53:VAL:HG22	9:82:95:LYS:NZ	2.29	0.48
34:58:73:THR:HB	34:58:82:LEU:HD11	1.95	0.48
1:13:1425:U:H2'	1:13:1426:C:C6	2.49	0.48
26:14:1416:G:H21	26:14:1586:A:N6	2.12	0.48
26:1H:1021:A:H8	26:1H:1022:G:H5''	1.79	0.48
26:1H:1142(A):A:C4	26:1H:1144:G:N7	2.82	0.48
26:1H:956:G:OP2	37:88:14:ARG:NH2	2.46	0.48
19:AA:51:VAL:O	19:AA:57:HIS:HA	2.13	0.48
1:13:1178:G:H5''	9:8E:93:ARG:NH2	2.28	0.48
38:98:29:LEU:HD23	38:98:70:LEU:HD11	1.94	0.48
51:M8:38:LYS:O	51:M8:39:CYS:HB3	2.12	0.48
46:H8:63:ASP:HB2	46:H8:65:GLN:HG3	1.95	0.48
26:14:38:A:H2'	26:14:39:C:C6	2.49	0.48
1:13:337:C:H2'	1:13:338:A:C8	2.48	0.48
30:31:34:TRP:CZ3	30:31:35:GLU:HG2	2.49	0.48
40:75:105:LEU:CD2	40:75:109:GLU:HG3	2.44	0.48
1:1G:1189:C:H5''	1:1G:1190:G:OP2	2.12	0.48
26:1H:2846:G:N7	58:1H:4680:HOH:O	2.34	0.48
17:8A:99:SER:C	17:8A:100:LYS:HZ3	2.16	0.48
26:14:2495:G:O3'	37:45:81:VAL:HG12	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:11:118:VAL:HG22	28:11:119:ALA:H	1.79	0.48
8:72:20:TYR:HA	8:72:65:TYR:CZ	2.47	0.48
13:4I:50:GLU:O	13:4I:54:VAL:HG23	2.13	0.48
32:51:86:GLU:CD	32:51:86:GLU:H	2.16	0.48
26:1H:1087:G:C5	26:1H:1089:G:H1'	2.48	0.48
26:14:1006:C:H5'	34:15:28:THR:HG23	1.95	0.48
26:14:953:A:H2'	26:14:954:G:H8	1.77	0.48
27:1J:41:U:C4	31:49:70:VAL:HG23	2.48	0.48
5:42:39:GLY:HA2	5:42:71:LEU:HD12	1.95	0.48
41:C8:52:ARG:HA	41:C8:55:ARG:HG3	1.96	0.48
26:1H:217:G:P	58:1H:4616:HOH:O	2.70	0.48
26:14:511:U:C5	26:14:512:G:C5	3.02	0.48
30:31:174:VAL:HG11	30:31:189:THR:HG21	1.96	0.48
1:1G:1203:C:H2'	1:1G:1204:A:O4'	2.14	0.48
26:1H:243:U:OP2	55:Q8:8:LYS:NZ	2.38	0.48
26:1H:668:G:H2'	26:1H:670:A:H62	1.78	0.48
55:Q8:4:MET:HB2	55:Q8:59:LYS:HD2	1.95	0.48
26:14:996:A:C2	26:14:997:G:C8	3.02	0.48
24:3K:5:G:H1'	24:3K:69:G:N2	2.29	0.48
26:1H:1591:G:H2'	26:1H:1592:C:H6	1.79	0.48
31:41:67:LYS:CE	51:M8:6:HIS:CE1	2.95	0.48
26:14:1069:A:H2'	26:14:1073:A:N7	2.29	0.48
1:13:428:G:O4'	1:13:430:A:C8	2.67	0.48
5:4E:13:ILE:HD12	5:4E:13:ILE:HA	1.74	0.48
26:14:2134:A:OP2	26:14:2157:G:N2	2.45	0.48
26:14:955:C:OP1	37:45:85:LYS:NZ	2.47	0.48
1:13:291:C:N4	1:13:309:G:H1	2.12	0.48
1:1G:451:A:N6	1:1G:480:U:H2'	2.28	0.48
26:14:943:U:OP2	36:35:36:LYS:HE3	2.13	0.48
26:14:307:G:N2	26:14:310:A:O5'	2.38	0.48
26:14:329:G:P	45:C5:71:LYS:HD3	2.54	0.48
1:13:992:U:O2'	1:13:993:G:OP2	2.29	0.48
26:14:2647:U:H2'	26:14:2648:C:C6	2.49	0.48
2:1E:168:THR:OG1	2:1E:192:SER:HB2	2.13	0.48
26:14:2537:U:H2'	26:14:2538:C:H6	1.75	0.48
1:13:1123:A:O2'	10:1I:37:PRO:O	2.24	0.48
12:3A:27:LEU:HB3	12:3A:33:ARG:CG	2.44	0.48
26:14:1027:A:C2	26:14:2488:A:H5'	2.49	0.48
5:4E:68:GLU:HG3	5:4E:68:GLU:O	2.13	0.48
1:1G:737:A:H2'	1:1G:738:C:C6	2.47	0.48
1:1G:617:G:C2	1:1G:618:C:C5	3.01	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:622:A:C8	1:1G:623:C:C6	3.02	0.48
15:6I:82:ILE:O	15:6I:86:GLY:N	2.44	0.48
1:13:1150:U:O2	10:1I:39:PRO:HG2	2.14	0.48
28:19:253:GLN:HB2	28:19:257:LEU:HD12	1.94	0.48
27:1J:31:C:C2'	27:1J:32:C:H5'	2.44	0.48
33:61:68:LEU:HA	33:61:71:ILE:HG22	1.96	0.48
26:1H:2109:U:O2'	26:1H:2181:G:N2	2.47	0.48
26:1H:2864:G:H2'	26:1H:2865:U:H6	1.79	0.48
45:G8:40:GLU:O	45:G8:40:GLU:HG3	2.13	0.48
1:13:1417:G:N2	1:13:1482:G:H2'	2.29	0.48
3:2E:43:LEU:O	3:2E:47:LEU:HB2	2.13	0.48
26:1H:2206:C:H2'	26:1H:2207:C:H6	1.77	0.48
12:3A:89:ARG:O	12:3A:99:HIS:HE1	1.96	0.48
1:13:135:C:H2'	1:13:136:C:H5'	1.95	0.48
1:1G:1114:C:H2'	1:1G:1115:C:H6	1.78	0.48
26:1H:86:C:H4'	26:1H:104:U:H1'	1.95	0.48
51:M8:14:ILE:HA	51:M8:31:ILE:O	2.14	0.48
26:1H:1598:C:O2'	26:1H:1599:C:H5'	2.13	0.48
37:45:42:ILE:O	37:45:95:ALA:N	2.44	0.48
29:21:6:GLY:O	29:21:195:LEU:HD23	2.13	0.48
1:1G:1072:G:C6	1:1G:1073:U:C4	3.01	0.48
41:C8:93:LYS:O	41:C8:96:ALA:HB2	2.13	0.48
1:1G:675:A:N1	1:1G:716:A:H2	2.12	0.48
26:1H:1142(A):A:N3	26:1H:1144:G:C8	2.82	0.48
1:1G:468:A:O2'	16:7A:82:GLN:HG2	2.13	0.48
1:13:872:A:C4	1:13:874:G:N7	2.82	0.48
26:1H:2347:C:H4'	53:O8:39:TYR:HE2	1.79	0.48
38:98:26:LYS:HE2	38:98:70:LEU:O	2.13	0.48
1:13:1452:C:H4'	1:13:1453:G:O5'	2.13	0.48
26:1H:2340:G:O2'	26:1H:2341:G:H5'	2.14	0.48
26:1H:142:G:C1'	44:F8:37:THR:HG21	2.42	0.48
26:14:1179:C:H2'	26:14:1180:C:C6	2.49	0.48
1:1G:1305:G:H1	1:1G:1331:G:H2'	1.79	0.48
40:B8:16:ARG:HH21	40:B8:19:LEU:HD21	1.77	0.48
1:1G:625:G:C4	1:1G:626:U:C5	3.02	0.48
26:1H:2791:C:N3	26:1H:2807:G:N2	2.61	0.48
10:1I:46:ARG:HB2	10:1I:46:ARG:NH1	2.29	0.48
19:AA:61:TYR:CE2	19:AA:63:THR:HA	2.48	0.48
10:1A:49:VAL:O	10:1A:60:ARG:HB2	2.13	0.48
27:16:90:C:P	37:88:16:ARG:HH21	2.36	0.48
1:1G:742:G:H5'	15:6A:58:MET:HE1	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1052:U:H5''	1:1G:1053:G:OP2	2.14	0.48
34:15:39:ARG:NH1	34:15:41:ASP:HB2	2.29	0.48
23:2L:9:G:N3	23:2L:46:G:H2'	2.29	0.48
45:C5:39:VAL:HG23	45:C5:41:GLY:N	2.29	0.48
1:1G:485:G:H1'	1:1G:486:U:H5	1.78	0.48
26:14:863:A:H2'	26:14:864:G:H8	1.78	0.48
28:19:124:PRO:HG2	28:19:129:ASN:HD21	1.78	0.48
26:1H:1680:U:H2'	26:1H:1681:G:O4'	2.13	0.48
7:62:87:VAL:HG22	7:62:88:PRO:O	2.14	0.48
2:1E:112:VAL:O	2:1E:115:LEU:N	2.47	0.48
26:1H:1564:C:H2'	26:1H:1565:C:C6	2.49	0.48
45:G8:68:HIS:O	45:G8:71:LYS:HB2	2.14	0.48
36:35:52:GLU:O	36:35:54:GLY:N	2.36	0.48
23:2K:26:C:H2'	23:2K:27:G:O4'	2.14	0.48
3:22:21:ARG:O	3:22:58:GLU:HA	2.13	0.48
1:13:450:G:OP1	16:7I:43:LYS:NZ	2.47	0.48
26:14:2563:U:H4'	35:25:28:SER:HA	1.96	0.48
42:95:85:LYS:HG3	42:95:87:HIS:HA	1.95	0.48
1:13:1036:G:N7	1:13:1037:C:N4	2.61	0.48
26:1H:917:A:N6	27:16:80:U:H4'	2.29	0.48
1:1G:1129:C:N4	1:1G:1141:C:H42	2.04	0.48
26:1H:2402:C:OP1	26:1H:2402:C:H4'	2.12	0.48
1:13:1027:C:H1'	1:13:1028:C:C5	2.49	0.48
24:3L:18:G:C2'	24:3L:57:G:H22	2.26	0.48
26:1H:1728:G:C2	26:1H:1730:U:OP2	2.67	0.48
31:41:67:LYS:HE2	51:M8:6:HIS:HE1	1.75	0.48
3:22:199:LYS:HB3	3:22:201:TYR:HE1	1.78	0.48
1:1G:1248:A:H2'	9:82:70:LYS:HZ1	1.78	0.48
1:13:558:G:C4	1:13:559:A:C2	3.01	0.48
26:1H:2126:A:H2'	26:1H:2126:A:N3	2.28	0.48
49:K8:15:LYS:HD3	49:K8:67:LYS:NZ	2.29	0.48
1:13:501:C:H2'	1:13:502:G:H8	1.79	0.48
33:61:2:LYS:NZ	33:61:20:ASP:HB3	2.29	0.48
3:22:18:TRP:HD1	14:5A:54:PRO:HA	1.77	0.48
27:16:15:A:H1'	27:16:109:G:N9	2.29	0.48
26:14:2869:G:H2'	26:14:2870:C:O4'	2.12	0.48
30:39:181:LEU:CD2	30:39:186:ILE:HD11	2.43	0.48
1:1G:509:A:C8	1:1G:509:A:H3'	2.49	0.48
26:1H:484:C:H2'	26:1H:485:C:H6	1.79	0.48
46:H8:9:TYR:HE1	46:H8:28:MET:HE1	1.78	0.48
26:14:1534:G:O2'	26:14:1538:G:N2	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:32:A:C2	1:1G:33:A:C4	3.02	0.48
1:1G:563:A:N6	58:1G:1833:HOH:O	2.45	0.48
26:1H:324:A:H2'	26:1H:325:G:H5'	1.95	0.48
1:1G:1151:A:HO2'	1:1G:1152:A:C5'	2.25	0.48
32:59:30:LYS:HB3	32:59:79:VAL:O	2.14	0.48
26:1H:1709:U:H2'	26:1H:1710:C:C6	2.48	0.48
26:1H:356:G:H2'	26:1H:357:A:H8	1.78	0.48
26:14:2238:G:N7	58:14:3607:HOH:O	2.35	0.48
7:6E:146:GLU:O	7:6E:149:ARG:HB2	2.13	0.48
42:D8:15:GLU:HG3	42:D8:16:PRO:HD2	1.96	0.48
26:14:2608:G:H8	26:14:2608:G:O5'	1.96	0.48
26:1H:747:U:O2	26:1H:2014:A:H1'	2.13	0.48
26:1H:2309:A:C5	26:1H:2310:A:H8	2.32	0.48
26:1H:934:G:H2'	26:1H:935:C:H6	1.78	0.48
48:F5:67:ILE:N	48:F5:68:PRO:HD2	2.28	0.48
7:62:45:ASP:HB3	7:62:117:ALA:HB1	1.95	0.48
42:95:22:VAL:HG22	42:95:23:GLU:H	1.78	0.48
26:14:1142:U:O2	26:14:1142:U:H2'	2.13	0.48
24:3K:1:G:H3'	24:3K:2:C:H5''	1.96	0.48
26:14:2211:G:O2'	26:14:2212:A:OP1	2.32	0.48
38:98:117:VAL:O	38:98:118:GLU:HB2	2.12	0.48
12:3A:110:VAL:HG23	12:3A:120:TYR:HB3	1.95	0.48
26:1H:141:A:H8	26:1H:1595:G:H21	1.62	0.48
26:1H:1533:C:H3'	26:1H:1534:G:C5'	2.44	0.48
26:1H:2232:U:OP1	48:J8:40:ARG:NH1	2.45	0.48
4:3E:108:LEU:HD23	4:3E:110:PHE:HE1	1.78	0.48
27:1J:94:C:H2'	27:1J:95:U:C6	2.49	0.48
28:11:44:ASN:OD1	28:11:46:GLN:HG3	2.14	0.48
26:14:841:A:H2'	26:14:842:G:C8	2.49	0.48
22:1K:44:G:H2'	22:1K:45:U:C5	2.49	0.48
8:7E:104:ARG:HB2	8:7E:107:LEU:HB3	1.95	0.48
46:H8:44:PHE:HE2	46:H8:86:VAL:HG11	1.78	0.48
30:31:135:LYS:HB3	30:31:138:GLU:HG3	1.96	0.48
26:14:940:G:N3	26:14:1191:G:H4'	2.29	0.48
26:14:925:C:H2'	26:14:926:A:H8	1.78	0.48
26:1H:1168:G:C2	26:1H:1182:A:C2	3.02	0.48
1:1G:583:A:H2'	1:1G:584:G:O4'	2.13	0.48
3:2E:83:ARG:O	3:2E:86:VAL:HG22	2.13	0.48
6:52:30:LEU:HD23	6:52:75:LEU:HD11	1.95	0.48
1:13:157:G:H1	1:13:164:U:H3	1.62	0.48
31:41:16:ARG:O	31:41:20:ILE:HG12	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1637:A:H4'	26:1H:2711:A:O2'	2.14	0.48
55:Q8:46:ARG:HB2	55:Q8:46:ARG:NH1	2.29	0.48
55:Q8:6:THR:H	55:Q8:59:LYS:CE	2.24	0.48
45:C5:81:LYS:HB2	45:C5:99:CYS:SG	2.54	0.48
45:C5:82:PRO:HB3	45:C5:97:ARG:HB3	1.96	0.48
1:1G:1312:G:N2	1:1G:1325:C:O2	2.40	0.48
1:1G:868:C:H2'	1:1G:869:G:O4'	2.13	0.48
31:49:7:LEU:HB2	31:49:104:GLU:OE1	2.14	0.48
52:N8:50:GLY:N	52:N8:56:LYS:HG3	2.28	0.48
24:3L:56:C:H3'	24:3L:57:G:H8	1.78	0.48
21:1B:25:LYS:HE3	21:1B:26:LYS:HG2	1.95	0.48
26:14:2114:A:N6	26:14:2119:A:H62	2.12	0.48
1:1G:1316:G:N2	1:1G:1318:A:H3'	2.29	0.48
31:41:96:ARG:O	31:41:97:ASP:HB2	2.14	0.48
28:19:239:ARG:HB2	28:19:240:ALA:H	1.44	0.48
1:13:1216:G:H5''	14:5I:5:ALA:HB2	1.96	0.48
1:13:1117:G:O3'	9:8E:104:ARG:HD3	2.13	0.48
1:13:376:G:OP1	16:7I:5:ARG:HB2	2.13	0.48
10:1A:96:ILE:HD11	10:1A:98:ILE:HG12	1.96	0.48
26:14:2291:U:O2'	26:14:2374:C:O2	2.29	0.48
1:1G:1065:U:H6	1:1G:1190:G:H21	1.62	0.48
22:1L:10:G:O2'	22:1L:11:C:OP1	2.29	0.48
26:14:175:G:C2'	26:14:176:G:H5'	2.43	0.48
47:I8:63:VAL:HG23	47:I8:64:ASP:O	2.14	0.48
26:1H:1215:G:C5	26:1H:1216:G:C8	3.02	0.48
31:41:33:ARG:HB2	31:41:162:THR:HG21	1.96	0.48
15:6I:26:GLU:OE2	15:6I:77:ARG:NH1	2.45	0.48
26:1H:396:G:O3'	48:J8:44:PRO:HA	2.14	0.48
13:4I:27:LYS:HA	13:4I:31:LYS:HZ2	1.79	0.48
19:AA:11:VAL:HB	19:AA:12:ASP:OD1	2.14	0.48
1:1G:1448:C:H2'	1:1G:1449:C:O4'	2.13	0.48
51:M8:23:GLU:OE1	51:M8:24:THR:N	2.47	0.48
31:49:173:LEU:HD22	31:49:178:PHE:CE1	2.49	0.48
4:3E:57:ARG:HE	4:3E:205:GLU:HB3	1.78	0.48
1:13:57:G:H2'	1:13:58:C:C6	2.49	0.48
50:H5:23:LEU:HD23	50:H5:28:LEU:HB2	1.96	0.48
20:BI:35:THR:O	20:BI:38:LYS:HB2	2.13	0.48
22:1K:3:C:O5'	22:1K:3:C:H6	1.97	0.48
44:F8:34:ALA:O	44:F8:77:LYS:NZ	2.47	0.48
1:1G:1295:G:H21	1:1G:1302:U:H3	1.61	0.48
38:98:45:ARG:HB3	38:98:46:GLY:H	1.57	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:2E:71:ALA:HB2	3:2E:109:PRO:HB3	1.96	0.48
29:29:4:ILE:HD12	29:29:198:VAL:HB	1.96	0.48
29:21:25:VAL:HB	29:21:183:LEU:HD12	1.96	0.48
7:6E:31:MET:SD	7:6E:36:LYS:HG2	2.54	0.48
23:2L:44:A:H2'	23:2L:45:A:C8	2.49	0.48
26:1H:2392:A:N1	26:1H:2424:C:N3	2.61	0.47
26:1H:567:A:N6	58:1H:4551:HOH:O	2.46	0.47
26:1H:1021:A:C3'	26:1H:1021:A:C8	2.97	0.47
7:62:72:ARG:NH2	7:62:96:GLN:OE1	2.46	0.47
1:13:22:G:H2'	1:13:23:C:C6	2.49	0.47
26:1H:1900:A:N1	26:1H:1970:A:C6	2.82	0.47
1:1G:464:G:C6	1:1G:466:C:H5'	2.48	0.47
26:1H:1006:C:O2	34:58:106:MET:HG2	2.14	0.47
5:42:109:ILE:HG22	5:42:110:LEU:HD23	1.96	0.47
22:1L:73:A:O2'	22:1L:75:C:OP2	2.32	0.47
32:51:43:VAL:HB	32:51:52:VAL:HG22	1.96	0.47
37:45:26:TYR:O	37:45:26:TYR:HD1	1.96	0.47
1:1G:1256:A:H62	1:1G:1277:C:H3'	1.78	0.47
20:BA:81:LYS:O	20:BA:85:MET:HG2	2.14	0.47
26:14:775:G:C4	26:14:794:G:C8	3.02	0.47
1:13:112:G:H2'	1:13:113:G:H5'	1.96	0.47
26:1H:1219:G:OP2	41:C8:19:LYS:NZ	2.39	0.47
26:1H:1230:C:H2'	26:1H:1231:G:H8	1.77	0.47
5:42:6:PHE:HD1	5:42:36:ASP:HB3	1.79	0.47
26:1H:1533:C:C3'	26:1H:1534:G:H5''	2.44	0.47
26:1H:2378:A:H4'	39:A8:23:ARG:NH1	2.29	0.47
1:13:1291:G:OP1	7:6E:37:ASN:ND2	2.47	0.47
22:1L:51:U:H2'	22:1L:52:G:H8	1.77	0.47
26:14:2850:A:C2	26:14:2851:A:C4	3.01	0.47
26:14:2520:C:H41	26:14:2542:A:H62	1.61	0.47
26:14:2165:G:C2'	26:14:2166:G:H5'	2.44	0.47
7:6E:40:ALA:HB1	7:6E:44:TYR:CD2	2.49	0.47
39:65:62:LYS:O	39:65:66:ALA:N	2.42	0.47
31:49:107:LEU:HD21	31:49:178:PHE:CE2	2.49	0.47
37:88:17:LEU:HA	37:88:17:LEU:HD23	1.53	0.47
26:14:1138:G:H21	34:15:106:MET:HE3	1.79	0.47
26:14:244:A:C2	26:14:255:A:C4	3.02	0.47
38:98:45:ARG:HD2	38:98:97:VAL:HG21	1.95	0.47
45:C5:33:LYS:HE2	45:C5:34:LYS:HG3	1.96	0.47
17:8A:86:GLU:O	17:8A:90:ILE:HG12	2.13	0.47
26:1H:900:A:H5'	26:1H:901:A:OP2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:455:C:N3	26:1H:473:G:H5'	2.29	0.47
3:22:113:ALA:HB2	3:22:202:ILE:HG13	1.95	0.47
26:1H:2114:A:H5'	26:1H:2117:A:OP2	2.14	0.47
1:13:977:A:O2'	1:13:979:C:OP2	2.28	0.47
26:1H:1300:U:H3'	58:1H:3608:HOH:O	2.14	0.47
15:6A:75:PRO:HB2	15:6A:79:ARG:HH12	1.79	0.47
28:11:38:LYS:HE2	28:11:39:LYS:O	2.14	0.47
2:1E:157:ARG:NH1	2:1E:157:ARG:HB3	2.28	0.47
2:12:23:ARG:C	2:12:23:ARG:HD3	2.34	0.47
6:5E:7:ASN:ND2	6:5E:62:TRP:HD1	2.12	0.47
26:14:2187:G:C5	26:14:2188:C:C4	3.02	0.47
26:1H:706:A:OP1	28:11:7:LYS:HE3	2.14	0.47
4:3E:104:VAL:O	4:3E:107:ARG:N	2.47	0.47
26:1H:1479:G:O2'	26:1H:1558:A:H5'	2.14	0.47
27:1J:39:A:N1	51:I5:1:MET:N	2.56	0.47
24:3L:51:U:H2'	24:3L:52:G:C8	2.49	0.47
32:51:46:GLU:CD	32:51:51:ARG:HH22	2.18	0.47
31:41:94:LEU:N	31:41:94:LEU:HD23	2.28	0.47
26:1H:1077:A:H5'	26:1H:1078:U:OP2	2.13	0.47
26:1H:1103:A:H3'	26:1H:1104:C:C6	2.49	0.47
26:14:2134:A:C2	26:14:2159:G:H1'	2.49	0.47
26:14:2134:A:H62	26:14:2157:G:C1'	2.26	0.47
26:14:1607:C:H4'	26:14:1608:A:O5'	2.13	0.47
26:1H:2346:A:O2'	53:O8:24:GLU:OE2	2.31	0.47
26:14:2776:A:H4'	26:14:2777:G:O5'	2.14	0.47
26:14:1778:U:P	58:14:3630:HOH:O	2.71	0.47
26:14:1665:A:C4'	35:25:67:LYS:HB2	2.44	0.47
16:7I:18:ARG:NH1	16:7I:32:TYR:OH	2.47	0.47
26:1H:1568:G:N3	28:11:58:HIS:HE1	2.12	0.47
1:13:166:G:H2'	1:13:167:G:C8	2.48	0.47
7:62:69:VAL:HG13	7:62:134:ALA:O	2.13	0.47
1:1G:446:G:H2'	1:1G:447:G:O4'	2.14	0.47
26:1H:557:U:H2'	26:1H:558:G:C8	2.49	0.47
1:1G:922:G:H4'	5:42:20:GLN:HA	1.96	0.47
26:14:2854:G:C2	26:14:2864:G:C2	3.01	0.47
26:14:478:A:N1	26:14:500:G:H4'	2.28	0.47
35:68:104:ARG:HD3	40:B8:36:GLU:HG2	1.94	0.47
1:13:1052:U:O2'	1:13:1055:A:OP2	2.31	0.47
15:6I:87:ILE:HG22	15:6I:88:ARG:N	2.29	0.47
33:69:61:ARG:HA	33:69:64:GLU:HB2	1.96	0.47
13:4I:57:ARG:HB2	13:4I:57:ARG:HH11	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:355:G:H2'	26:14:356:G:C8	2.49	0.47
55:Q8:56:GLU:HA	55:Q8:58:ILE:HG22	1.95	0.47
26:1H:1453:A:O2'	26:1H:1454:U:H2'	2.13	0.47
19:AA:51:VAL:HG12	19:AA:52:TYR:H	1.79	0.47
26:1H:197:A:N6	26:1H:2430:A:H2'	2.27	0.47
1:13:1323:G:H4'	1:13:1362(A):C:N3	2.29	0.47
30:31:114:VAL:HG21	30:31:202:PHE:CZ	2.50	0.47
32:51:30:LYS:HE3	32:51:81:GLU:N	2.22	0.47
1:1G:673:G:H5''	6:52:87:ARG:NH1	2.29	0.47
26:14:1658:C:H2'	26:14:1659:U:C6	2.49	0.47
9:82:111:ARG:HG2	9:82:112:LYS:H	1.79	0.47
42:D8:10:LYS:NZ	42:D8:23:GLU:OE1	2.47	0.47
26:1H:2032:G:N2	29:21:146:THR:HG23	2.29	0.47
1:13:10:A:O2'	1:13:11:G:H5'	2.14	0.47
20:BA:51:GLU:HA	20:BA:54:LYS:HB3	1.96	0.47
1:13:626:U:H5''	16:7I:38:TYR:CD2	2.49	0.47
26:14:2503:A:OP1	58:14:3571:HOH:O	2.19	0.47
26:14:2012:G:OP1	43:A5:11:ARG:NH2	2.44	0.47
29:29:89:ASP:CG	29:29:90:THR:H	2.18	0.47
26:14:20:C:OP1	41:85:22:LYS:NZ	2.33	0.47
26:1H:528:A:N1	26:1H:2042:A:H2'	2.29	0.47
26:1H:1111:A:H5'	32:51:3:ARG:HD2	1.96	0.47
27:16:16:G:N2	27:16:69:G:H1'	2.29	0.47
26:14:2712:U:OP1	26:14:2714:G:H4'	2.14	0.47
1:1G:447:G:O6	1:1G:485:G:H2'	2.14	0.47
1:13:1442:G:C6	1:13:1446:A:N6	2.82	0.47
1:1G:229:U:H2'	1:1G:230:G:C8	2.49	0.47
27:1J:100:G:O5'	58:1J:309:HOH:O	2.20	0.47
1:1G:1198:G:H2'	1:1G:1199:U:H6	1.80	0.47
33:61:29:TYR:O	33:61:32:PRO:HD2	2.13	0.47
5:42:72:GLN:O	5:42:75:THR:HG22	2.14	0.47
27:1J:56:G:H4'	27:1J:57:A:C8	2.49	0.47
1:13:1339:A:H2'	1:13:1340:A:O4'	2.15	0.47
28:19:273:ARG:O	28:19:273:ARG:HG2	2.15	0.47
28:11:112:GLN:H	28:11:115:GLN:HG3	1.79	0.47
55:Q8:24:ALA:O	55:Q8:46:ARG:HG2	2.14	0.47
26:1H:1188:U:C4'	42:D8:79:VAL:HG22	2.43	0.47
26:1H:2311:A:H1'	31:41:88:ILE:HD13	1.95	0.47
32:51:83:TYR:CD2	32:51:134:SER:HA	2.50	0.47
26:1H:2452:C:OP1	58:1H:4436:HOH:O	2.20	0.47
26:14:460:A:H5''	26:14:461:C:OP2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1774:C:O5'	26:14:1774:C:H6	1.97	0.47
13:4A:57:ARG:O	13:4A:61:GLU:HB2	2.14	0.47
26:1H:1726:G:H2'	26:1H:1727:U:O4'	2.14	0.47
29:29:25:VAL:O	29:29:26:ILE:HG12	2.14	0.47
34:58:41:ASP:O	41:C8:64:ARG:NH2	2.47	0.47
55:Q8:34:TRP:HB2	55:Q8:35:GLN:HA	1.95	0.47
1:13:377:G:H5'	16:7I:5:ARG:HH12	1.78	0.47
29:29:33:VAL:HG23	29:29:47:VAL:HG13	1.95	0.47
2:12:6:THR:OG1	2:12:7:VAL:N	2.47	0.47
26:14:1252:G:O4'	41:85:33:ARG:HD3	2.15	0.47
26:14:1016:G:H2'	26:14:1017:G:O4'	2.15	0.47
32:51:3:ARG:HA	32:51:3:ARG:CZ	2.44	0.47
31:41:77:ILE:O	31:41:81:LYS:O	2.32	0.47
26:1H:1203:G:H5'	36:78:3:LEU:HD13	1.96	0.47
26:14:2002:G:O6	58:14:3797:HOH:O	2.20	0.47
26:1H:2783:G:H2'	26:1H:2784:C:C6	2.49	0.47
37:45:20:ALA:HA	37:45:99:PRO:HG2	1.96	0.47
29:21:45:THR:O	29:21:83:ASP:N	2.48	0.47
35:68:120:GLU:HG2	35:68:122:LEU:HG	1.95	0.47
1:13:791:G:C6	1:13:792:A:C2	3.02	0.47
26:14:1439:A:C2	26:14:1553:A:C4	3.02	0.47
1:13:380:G:N2	1:13:384:G:C5	2.83	0.47
1:1G:1454:G:OP1	20:BA:39:LYS:NZ	2.38	0.47
27:16:40:U:H5	51:M8:2:LYS:HE2	1.79	0.47
34:58:99:LEU:O	34:58:103:VAL:HG23	2.14	0.47
2:1E:87:ARG:NH2	2:1E:233:SER:HB2	2.30	0.47
54:L5:29:LYS:HA	54:L5:32:LYS:HB2	1.95	0.47
26:14:2259:G:C2	26:14:2282:G:N1	2.83	0.47
32:51:37:VAL:HG22	32:51:68:THR:HG22	1.96	0.47
26:14:603:A:H8	26:14:604:G:H1'	1.79	0.47
6:52:35:ALA:HA	6:52:67:MET:HB3	1.95	0.47
7:62:101:LEU:O	7:62:105:VAL:HG23	2.14	0.47
49:G5:5:GLU:H	49:G5:5:GLU:CD	2.18	0.47
11:2A:87:THR:O	11:2A:87:THR:OG1	2.26	0.47
41:C8:47:TYR:C	41:C8:47:TYR:CD1	2.88	0.47
54:P8:37:LYS:O	54:P8:37:LYS:HG3	2.14	0.47
31:41:84:LYS:O	31:41:84:LYS:HG3	2.14	0.47
1:1G:251:G:H4'	1:1G:252:U:O5'	2.14	0.47
9:8E:29:ASN:OD1	9:8E:65:VAL:N	2.46	0.47
42:95:62:LEU:HD21	42:95:95:LEU:HB2	1.97	0.47
26:1H:589:C:H2'	26:1H:590:A:H8	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:78:18:ARG:C	36:78:19:VAL:HG22	2.34	0.47
1:13:1504:G:H3'	1:13:1504:G:P	2.54	0.47
26:1H:1324:G:C4	26:1H:1328:G:O6	2.67	0.47
26:1H:861:A:H5''	58:1H:3993:HOH:O	2.13	0.47
1:13:1156:G:H2'	1:13:1157:A:H5''	1.96	0.47
26:1H:1510:A:OP1	26:1H:1511:A:H5'	2.13	0.47
11:2I:34:ASP:OD2	11:2I:38:ASN:HB2	2.15	0.47
44:F8:3:THR:CB	44:F8:4:ALA:HA	2.43	0.47
1:1G:1327:C:H2'	1:1G:1328:C:C6	2.50	0.47
26:1H:2284:C:C2'	26:1H:2285:C:H5'	2.45	0.47
26:1H:2383:G:O2'	26:1H:2384:G:H5'	2.15	0.47
1:1G:537:G:H2'	1:1G:538:G:C8	2.50	0.47
26:1H:1404:C:O2'	26:1H:1405:U:H5'	2.14	0.47
45:G8:54:LYS:O	45:G8:55:TYR:CG	2.67	0.47
12:3I:47:LYS:HA	12:3I:49:ASN:N	2.29	0.47
1:1G:1099:G:OP1	2:12:148:TYR:OH	2.32	0.47
26:14:1971:A:H5''	58:14:3452:HOH:O	2.12	0.47
26:1H:296:C:H2'	26:1H:297:C:C6	2.50	0.47
5:42:18:ARG:HH21	5:42:25:ARG:HB3	1.79	0.47
19:AI:41:VAL:HB	19:AI:42:PRO:HA	1.95	0.47
7:6E:16:LEU:HD12	9:8E:42:ARG:HA	1.96	0.47
26:14:579:G:H2'	26:14:580:C:H6	1.76	0.47
26:14:2342:C:O2'	26:14:2374:C:H5''	2.15	0.47
26:14:581:C:H2'	26:14:582:G:C8	2.46	0.47
1:1G:438:G:C4'	4:32:123:HIS:HD2	2.27	0.47
26:1H:1215:G:C4	26:1H:1216:G:C8	3.02	0.47
1:13:15:G:OP1	1:13:1396:A:O2'	2.24	0.47
26:14:2311:A:C8	31:49:88:ILE:HG12	2.49	0.47
26:1H:1106:G:H2'	26:1H:1107:G:O4'	2.13	0.47
3:2E:191:THR:HG21	3:2E:193:TYR:CZ	2.50	0.47
2:12:54:THR:HG23	2:12:199:TYR:HB3	1.97	0.47
9:8E:25:LYS:HZ1	9:8E:32:ASP:HA	1.80	0.47
26:1H:805:G:P	36:78:41:ARG:HG2	2.54	0.47
26:14:2154:G:H2'	26:14:2155:G:O4'	2.13	0.47
26:1H:2367:G:H2'	26:1H:2368:C:C6	2.49	0.47
1:13:807:A:H2'	1:13:808:C:C6	2.50	0.47
26:14:297:C:O2	26:14:341:G:N2	2.42	0.47
27:1J:99:A:C4	27:1J:100:G:C8	3.03	0.47
26:1H:1811:G:H2'	26:1H:1812:A:O4'	2.13	0.47
26:1H:270(G):C:H2'	26:1H:270(H):C:O4'	2.14	0.47
28:11:140:THR:HG22	28:11:141:VAL:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1188:A:O3'	14:5A:58:LYS:HE3	2.15	0.47
26:14:2883:A:H5'	26:14:2884:U:H5'	1.94	0.47
1:1G:1013:G:N2	1:1G:1017:G:O6	2.46	0.47
27:16:54:G:H2'	27:16:55:U:H6	1.78	0.47
35:68:8:LEU:HB2	35:68:19:ILE:HG13	1.96	0.47
26:1H:1665:A:N7	58:1H:4084:HOH:O	2.35	0.47
26:14:698:C:O2'	26:14:734:A:N6	2.47	0.47
26:1H:207:A:H2'	26:1H:208:C:O4'	2.14	0.47
50:H5:35:ARG:HB3	50:H5:37:LEU:HD21	1.96	0.47
4:32:112:VAL:HG12	4:32:116:GLN:OE1	2.14	0.47
26:14:996:A:N3	26:14:997:G:C8	2.82	0.47
26:14:2702:U:HO2'	26:14:2703:C:H5	1.55	0.47
26:14:93:C:H5'	26:14:94:G:OP2	2.15	0.47
43:E8:38:TYR:CE2	52:N8:41:PRO:HD3	2.48	0.47
26:1H:1171:G:N2	26:1H:1179:C:N3	2.62	0.47
26:14:2208:U:H2'	26:14:2209:C:C6	2.49	0.47
41:C8:88:ILE:O	41:C8:90:VAL:N	2.47	0.47
1:1G:192:U:H2'	1:1G:193:C:C6	2.41	0.47
26:1H:2695:C:H2'	26:1H:2696:U:C6	2.49	0.47
1:1G:1279:A:H61	3:22:26:LYS:HZ1	1.61	0.47
2:12:144:ARG:HG3	2:12:145:LEU:N	2.28	0.47
47:E5:48:GLY:HA3	47:E5:80:HIS:CE1	2.48	0.47
1:13:196:A:O2'	1:13:221:C:O2'	2.29	0.47
13:4A:23:TYR:HE1	13:4A:70:LEU:HD12	1.79	0.47
33:61:46:ALA:O	33:61:50:ARG:N	2.42	0.47
6:52:100:ASN:ND2	18:9A:23:LYS:O	2.45	0.47
26:1H:813:U:H2'	26:1H:814:C:C6	2.49	0.47
1:13:618:C:H5''	1:13:619:U:H5''	1.96	0.47
27:16:29:A:OP2	39:A8:31:SER:HB2	2.15	0.47
26:1H:2815:C:H2'	26:1H:2816:C:C6	2.49	0.47
32:59:22:GLY:O	32:59:37:VAL:HG12	2.14	0.47
26:14:2447:G:O2'	58:14:4050:HOH:O	2.20	0.47
26:14:1450:C:H2'	26:14:1451:C:C6	2.50	0.47
11:2A:120:ARG:HA	11:2A:121:PRO:HD3	1.67	0.47
26:14:2025:C:N4	58:14:3606:HOH:O	2.47	0.47
1:1G:918:A:H2'	1:1G:919:A:O4'	2.15	0.47
5:4E:76:ILE:HG13	5:4E:93:PRO:HB3	1.96	0.47
1:13:1084:G:C5	1:13:1085:U:C4	3.02	0.47
22:1L:46:G:H2'	22:1L:48:C:H1'	1.96	0.47
1:13:127:G:O2'	17:8I:2:PRO:O	2.33	0.47
33:69:135:GLU:CD	33:69:135:GLU:H	2.17	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1647:G:OP2	26:1H:1647:G:H3'	2.13	0.47
26:1H:2081:C:H2'	26:1H:2082:A:C8	2.50	0.47
32:51:94:TYR:HA	32:51:106:THR:O	2.14	0.47
17:8A:79:SER:OG	17:8A:80:GLY:N	2.47	0.47
26:1H:1250:G:OP2	36:78:21:ARG:NH1	2.46	0.47
1:13:1347:G:OP2	9:8E:107:ARG:HG2	2.14	0.47
28:11:3:VAL:HG12	28:11:17:THR:HG23	1.97	0.47
1:13:254:G:OP1	17:8I:68:ARG:HB3	2.15	0.47
26:14:654(A):A:H2'	26:14:654(B):C:C6	2.50	0.47
1:1G:1320:C:OP1	19:AA:70:LYS:HE3	2.14	0.47
26:1H:2432:A:C4	48:J8:33:LYS:HG2	2.49	0.47
26:1H:1163:G:C2	26:1H:1164:G:N7	2.83	0.47
26:14:2266:A:H5'	26:14:2267:A:C5	2.49	0.47
40:B8:3:ARG:O	40:B8:6:LEU:N	2.48	0.47
1:1G:957:U:H2'	1:1G:959:A:OP2	2.15	0.47
26:1H:2394:C:H2'	26:1H:2395:C:H6	1.78	0.47
26:14:609(A):G:H2'	26:14:610:C:C6	2.50	0.47
20:BI:30:LYS:NZ	20:BI:80:ARG:HH12	2.12	0.47
26:1H:66:C:H2'	26:1H:67:U:H6	1.79	0.47
29:29:62:PRO:O	29:29:64:LYS:N	2.46	0.47
34:58:96:GLU:C	34:58:98:VAL:N	2.66	0.47
46:D5:30:ASN:ND2	46:D5:90:VAL:HB	2.28	0.47
1:1G:1127:G:N2	1:1G:1144:G:H22	2.13	0.47
1:1G:390:C:O3'	16:7A:28:ARG:NH2	2.45	0.47
53:O8:30:THR:HA	53:O8:31:PRO:C	2.35	0.47
55:Q8:28:GLY:N	55:Q8:29:LYS:HB3	2.29	0.47
32:51:77:LYS:HE2	32:51:138:LYS:HD2	1.96	0.47
40:B8:16:ARG:HH21	40:B8:19:LEU:CD2	2.28	0.47
1:1G:630:G:H3'	1:1G:631:G:H8	1.80	0.47
3:22:9:GLY:HA3	14:5A:49:HIS:HA	1.95	0.47
14:5A:45:ARG:O	14:5A:49:HIS:ND1	2.47	0.47
26:14:1000:A:C6	26:14:1001:A:N1	2.83	0.47
24:3L:70:G:H2'	24:3L:71:G:O4'	2.14	0.47
36:35:144:GLU:N	36:35:144:GLU:CD	2.68	0.47
19:AI:40:ILE:HA	19:AI:44:MET:SD	2.54	0.47
28:19:34:VAL:HG11	28:19:61:LEU:HG	1.95	0.47
28:19:16:MET:CE	28:19:208:LYS:HG2	2.44	0.47
1:13:648:A:H2'	1:13:649:G:H8	1.80	0.47
1:13:1218:C:H2'	1:13:1219:U:C6	2.50	0.47
1:13:339:C:H2'	1:13:340:U:H6	1.80	0.47
4:32:173:TRP:O	4:32:174:LEU:HD23	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1360:A:H2'	1:13:1361:G:C8	2.50	0.47
1:13:1360:A:O2'	1:13:1361:G:H5'	2.14	0.47
2:1E:7:VAL:HB	2:1E:217:ARG:HH11	1.79	0.47
2:1E:74:LYS:NZ	2:1E:205:ASP:O	2.47	0.47
26:1H:2680:C:H5'	29:21:189:PRO:HA	1.97	0.47
26:14:2555:U:C5	26:14:2556:C:C2	3.03	0.47
30:39:132:VAL:HG13	30:39:133:ASN:ND2	2.29	0.47
27:1J:88:C:P	27:1J:88:C:H6	2.38	0.47
26:14:839:U:H3	26:14:939:G:H1	1.63	0.47
38:55:104:ARG:HB2	38:55:107:ASP:OD1	2.15	0.47
1:1G:954:G:H21	1:1G:1227:A:N6	2.10	0.47
1:1G:4:U:O4	8:72:105:ARG:HD3	2.15	0.47
11:2I:18:ARG:O	11:2I:33:THR:N	2.45	0.47
3:2E:134:ILE:HG22	3:2E:168:ALA:HB3	1.97	0.47
26:14:235:U:H2'	26:14:236:C:C6	2.49	0.47
30:39:65:TRP:CZ3	30:39:72:ARG:HB3	2.50	0.47
26:14:2611:U:H5'	26:14:2611:U:H6	1.79	0.47
26:14:116:C:H2'	26:14:117:G:O4'	2.13	0.47
26:14:2495:G:H2'	26:14:2496:C:H6	1.80	0.47
19:AA:66:MET:SD	19:AA:66:MET:N	2.88	0.47
44:F8:55:ASN:HB2	44:F8:80:ILE:HG13	1.96	0.47
34:15:36:GLY:H	34:15:42:TRP:HZ3	1.62	0.47
1:1G:1123:A:O2'	10:1A:37:PRO:O	2.15	0.47
1:1G:685:G:N1	1:1G:686:U:O4	2.47	0.47
8:7E:49:GLU:HG2	8:7E:62:TYR:HE2	1.79	0.47
31:41:104:GLU:CD	51:M8:23:GLU:HG3	2.34	0.47
26:1H:460:A:H5''	26:1H:461:C:OP2	2.14	0.47
4:32:105:VAL:HG12	4:32:117:ALA:HB1	1.95	0.47
35:68:112:MET:HA	35:68:115:VAL:HG22	1.97	0.47
20:BI:34:LYS:O	20:BI:38:LYS:HG3	2.14	0.47
26:1H:1523:U:C2	26:1H:1524:G:C8	3.02	0.47
1:1G:186(C):G:C6	1:1G:186(D):C:C4	3.03	0.47
1:13:1389:C:H2'	1:13:1390:U:O4'	2.15	0.47
19:AI:68:GLY:HA3	51:M8:59:PHE:CZ	2.49	0.47
4:3E:147:ALA:HB2	4:3E:182:LYS:HB3	1.97	0.47
1:13:945:G:C2	1:13:946:A:C8	3.02	0.47
26:1H:1296:G:O2'	26:1H:1297:C:H5'	2.14	0.47
47:I8:54:GLY:O	47:I8:57:PHE:N	2.44	0.47
9:8E:86:VAL:O	9:8E:90:PRO:HB3	2.15	0.47
2:12:21:ARG:HA	2:12:39:ILE:HA	1.95	0.47
26:14:2586:C:C2'	26:14:2587:A:H5'	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2224:G:H4'	26:1H:2226:C:C2	2.49	0.47
11:2A:103:LEU:HD12	11:2A:103:LEU:HA	1.69	0.47
36:35:93:GLY:H	36:35:123:LEU:HD12	1.80	0.47
30:39:29:ASN:HA	30:39:30:PRO:HD3	1.67	0.47
28:11:67:PHE:HB3	28:11:153:ALA:H	1.80	0.47
55:Q8:14:VAL:O	55:Q8:15:LYS:HD3	2.14	0.47
55:Q8:49:VAL:HA	55:Q8:52:LYS:HG2	1.97	0.47
26:1H:2592:G:C6	26:1H:2593:U:C4	3.03	0.47
26:1H:1144:G:C4	26:1H:1145:C:C5	3.02	0.47
26:14:2702:U:O2'	26:14:2703:C:H5	1.97	0.47
1:13:631:G:C8	1:13:632:A:C2	3.03	0.47
26:14:1970:A:OP2	58:14:3459:HOH:O	2.20	0.47
1:13:872:A:C5	1:13:874:G:C8	3.03	0.47
36:35:63:PRO:CA	55:M5:13:ARG:HG2	2.45	0.47
32:51:4:ILE:HG13	32:51:6:ARG:CZ	2.45	0.47
26:14:2300:G:C2	26:14:2301:C:N3	2.82	0.47
1:1G:1287:A:N3	1:1G:1353:G:O2'	2.38	0.47
29:29:60:ASN:C	29:29:62:PRO:HD3	2.36	0.47
26:14:2114:A:H61	26:14:2170:A:N6	2.13	0.47
26:14:912:C:C2	26:14:913:U:C5	3.03	0.47
38:98:94:TYR:O	38:98:117:VAL:HB	2.14	0.47
26:14:2328:A:H2'	26:14:2329:G:C8	2.50	0.47
12:3I:49:ASN:ND2	12:3I:92:ASP:OD2	2.34	0.47
33:61:131:LYS:HB3	33:61:132:PRO:CA	2.43	0.47
37:45:117:ALA:HA	37:45:120:ILE:HB	1.96	0.47
1:1G:1503:A:H1'	25:4L:13:A:H61	1.79	0.47
1:13:490:G:H2'	1:13:491:G:H8	1.79	0.47
19:AI:44:MET:O	19:AI:47:HIS:HB2	2.15	0.47
26:1H:1024:G:C3'	26:1H:1025:G:H5''	2.43	0.47
1:13:1057:G:H2'	1:13:1058:G:O4'	2.15	0.47
45:C5:87:LYS:HB2	45:C5:96:ILE:HD11	1.96	0.47
2:1E:11:LEU:HD12	2:1E:217:ARG:HH12	1.80	0.47
18:9A:22:VAL:HG22	18:9A:23:LYS:H	1.80	0.47
26:1H:2264:C:N4	47:I8:15:ASP:OD2	2.45	0.47
1:1G:763:G:H2'	1:1G:764:C:H6	1.80	0.47
26:1H:2152:G:H2'	26:1H:2153:G:O4'	2.14	0.47
31:41:47:LYS:HD3	31:41:81:LYS:HB3	1.97	0.47
51:I5:20:ASN:HB3	51:I5:39:CYS:SG	2.55	0.47
8:7E:45:ILE:HD12	8:7E:47:GLY:HA2	1.95	0.47
32:59:46:GLU:OE2	32:59:51:ARG:NH1	2.47	0.47
3:2E:6:HIS:HB2	14:5I:49:HIS:ND1	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:446:G:H1	1:1G:488:C:H42	1.63	0.47
4:32:117:ALA:O	4:32:121:VAL:HG23	2.14	0.47
9:8E:18:PHE:HD2	9:8E:62:TYR:HD2	1.63	0.47
26:14:1006:C:H1'	34:15:106:MET:HE3	1.96	0.47
1:13:834:C:C2	1:13:853:G:C2	3.02	0.47
44:F8:5:TYR:CE1	49:K8:30:ARG:HD2	2.50	0.47
43:A5:73:ALA:H	43:A5:106:ILE:HG12	1.78	0.47
26:14:2774:C:H2'	26:14:2775:A:O4'	2.14	0.47
26:1H:108:U:C2	26:1H:109:G:C8	3.03	0.47
26:1H:2612:C:C5	26:1H:2613:U:H5	2.33	0.47
7:62:60:LYS:HD2	7:62:60:LYS:HA	1.65	0.47
5:4E:10:MET:HG3	5:4E:10:MET:O	2.15	0.47
1:13:1276:G:N3	1:13:1282:C:O2'	2.42	0.47
33:69:129:THR:HA	33:69:137:PRO:HA	1.97	0.47
36:35:65:ARG:HE	55:M5:15:LYS:HB3	1.79	0.47
2:1E:195:ASP:O	8:7E:74:PRO:HG3	2.15	0.47
11:2I:87:THR:HG22	11:2I:88:GLY:H	1.79	0.47
55:Q8:23:VAL:O	55:Q8:46:ARG:HG3	2.15	0.47
26:14:1114:G:H2'	26:14:1115:G:C8	2.50	0.47
42:95:71:LEU:O	42:95:72:VAL:HG12	2.15	0.47
26:1H:1144:G:H2'	26:1H:1145:C:H6	1.79	0.47
1:1G:1392:G:H21	1:1G:1502:A:H8	1.62	0.47
26:14:531:C:OP1	26:14:561:G:C2	2.67	0.47
27:1J:14:U:H5'	27:1J:71:C:O4'	2.15	0.47
26:14:1664:A:OP1	58:14:3507:HOH:O	2.20	0.47
26:1H:562:U:O4	26:1H:2036:C:H1'	2.15	0.47
26:14:932:G:OP1	50:H5:29:ARG:NH1	2.48	0.47
49:K8:47:ASN:HB2	49:K8:50:ILE:HD11	1.97	0.47
1:13:975:A:H5'	1:13:975:A:C8	2.49	0.47
26:1H:1386:C:OP2	26:1H:1396:U:H5	1.98	0.47
41:85:90:VAL:CG2	42:95:39:LEU:HB3	2.43	0.47
1:1G:1177:G:OP2	9:82:97:LYS:NZ	2.48	0.47
26:14:1558:A:H4'	26:14:1559:G:H5'	1.97	0.47
37:88:32:TYR:CZ	37:88:111:GLU:HG3	2.50	0.47
26:14:69:C:H2'	26:14:70:G:H8	1.80	0.47
1:1G:691:G:N7	11:2A:26:ASN:HB3	2.30	0.47
26:1H:2157:G:HO2'	26:1H:2158:A:P	2.38	0.47
39:A8:47:THR:HG22	39:A8:48:LEU:H	1.78	0.47
1:1G:1154:G:N3	1:1G:1155:G:C8	2.82	0.47
26:14:2228:G:OP1	28:19:261:LYS:HE3	2.15	0.47
37:45:118:LEU:HD22	37:45:118:LEU:HA	1.68	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:D8:21:ARG:HG2	42:D8:91:TYR:CE2	2.50	0.47
15:6A:39:LEU:HD12	15:6A:56:LEU:HB2	1.96	0.47
26:1H:2846:G:H2'	26:1H:2847:U:O4'	2.14	0.47
10:1I:6:ILE:HG22	10:1I:98:ILE:CG1	2.44	0.47
16:7A:21:VAL:HG22	16:7A:33:ILE:HB	1.97	0.47
26:14:270(J):G:N2	26:14:270(K):C:H1'	2.30	0.47
32:59:85:LYS:HG2	32:59:141:VAL:HG12	1.96	0.47
45:G8:35:TYR:CE2	45:G8:69:ALA:HB3	2.49	0.47
1:13:390:C:H2'	1:13:391:G:C8	2.50	0.47
40:B8:60:THR:HG22	40:B8:77:PRO:HA	1.96	0.47
3:2E:58:GLU:H	3:2E:65:ALA:HB3	1.80	0.47
10:1A:45:ARG:HB3	10:1A:65:LEU:HB3	1.97	0.47
26:14:2542:A:H1'	26:14:2543:G:C8	2.50	0.47
2:1E:17:PHE:HB3	2:1E:44:LEU:HD11	1.97	0.47
46:H8:116:VAL:H	46:H8:174:VAL:HG13	1.80	0.47
26:14:142:G:H2'	26:14:143:C:C6	2.50	0.47
54:L5:12:ARG:NH2	54:L5:44:PRO:HB3	2.30	0.47
27:16:116:G:H5''	39:A8:55:ALA:HB2	1.96	0.47
26:1H:2309:A:C6	26:1H:2310:A:C8	3.03	0.47
2:1E:87:ARG:NH1	2:1E:220:ASP:OD1	2.47	0.47
26:1H:2081:C:H2'	26:1H:2082:A:H8	1.79	0.47
1:1G:577:G:C8	1:1G:816:A:C6	3.02	0.47
52:J5:48:GLU:HG2	52:J5:48:GLU:H	1.35	0.47
26:14:110:G:C2	26:14:111:A:C8	3.02	0.47
46:H8:133:ILE:HA	46:H8:134:PRO:HD2	1.76	0.47
34:15:121:LYS:HB3	34:15:123:TYR:CE1	2.50	0.47
5:42:79:GLU:OE1	8:72:104:ARG:HA	2.15	0.47
26:14:2173:A:H8	26:14:2173:A:P	2.38	0.47
8:72:40:ALA:HA	8:72:45:ILE:HG13	1.96	0.47
26:1H:1973:G:H2'	26:1H:1974:C:H6	1.80	0.47
18:9A:45:SER:OG	18:9A:46:GLU:N	2.48	0.47
26:14:1354:A:H2'	26:14:1355:G:O4'	2.15	0.47
1:13:1376:U:H2'	1:13:1377:A:C8	2.49	0.47
55:Q8:57:ARG:HB2	55:Q8:59:LYS:HE2	1.97	0.47
26:14:1416:G:H21	26:14:1586:A:H62	1.62	0.47
26:14:676:A:H2	26:14:802:A:H61	1.61	0.47
26:1H:1288:U:H4'	26:1H:1289:C:OP2	2.15	0.47
46:H8:59:LEU:O	46:H8:60:GLU:HB2	2.14	0.47
26:14:654(B):C:H2'	26:14:654(C):G:C8	2.49	0.47
26:14:249:C:H4'	26:14:250:G:O5'	2.15	0.47
26:1H:444:C:H4'	30:31:49:ALA:HB2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:459:U:H5''	54:L5:40:TRP:CE2	2.49	0.47
30:39:25:PRO:HB2	30:39:27:GLU:O	2.15	0.47
20:BA:50:GLU:HA	20:BA:100:ILE:HG12	1.97	0.47
1:13:427:U:H5'	4:3E:41:GLY:HA2	1.97	0.47
48:J8:92:LYS:HA	48:J8:95:LEU:CB	2.43	0.47
26:14:2502:G:OP2	58:14:4028:HOH:O	2.20	0.47
26:14:2750:A:H8	26:14:2752:C:H41	1.63	0.47
24:3L:4:C:H2'	24:3L:5:G:O4'	2.15	0.47
26:1H:1317:A:H2'	26:1H:1318:C:C6	2.50	0.47
26:1H:182:A:H2'	26:1H:183:C:C6	2.50	0.47
22:1K:14:A:N1	22:1K:22:G:H1'	2.30	0.47
50:L8:31:LEU:HB3	50:L8:32:GLN:OE1	2.15	0.47
1:1G:1004:A:H2	1:1G:1024:G:C5	2.32	0.47
1:1G:1092:A:C2	1:1G:1183:A:H2	2.32	0.47
1:13:44:G:OP1	16:7I:12:LYS:HB3	2.15	0.47
47:E5:66:VAL:CG1	47:E5:67:VAL:H	2.27	0.47
26:14:2693:A:H2'	26:14:2694:G:C8	2.49	0.47
26:14:55:G:C2	26:14:116:C:C2	3.03	0.47
43:E8:3:ALA:HB3	43:E8:58:ALA:HB2	1.95	0.47
28:11:69:ARG:NH2	28:11:128:GLY:O	2.29	0.47
26:14:864:G:C6	26:14:865:C:N4	2.83	0.47
11:2I:12:ARG:HG3	11:2I:13:GLN:N	2.31	0.47
26:1H:2396:G:H5''	48:J8:25:LYS:HE2	1.96	0.47
26:1H:2557:G:O2'	26:1H:2558:C:H5'	2.15	0.47
26:1H:2468:G:OP1	37:88:119:ARG:NH2	2.44	0.47
45:G8:34:LYS:HD3	45:G8:36:ALA:HB2	1.97	0.47
1:13:1386:G:O2'	1:13:1387:G:H5'	2.14	0.47
26:1H:1456:G:OP2	58:1H:4659:HOH:O	2.19	0.47
26:14:2467:C:H4'	37:45:123:HIS:CE1	2.50	0.47
26:14:1251:C:H5	58:14:3962:HOH:O	1.98	0.47
1:1G:843:U:H3'	1:1G:848:C:O4'	2.15	0.47
26:14:2197:U:H1'	26:14:2198:A:C8	2.50	0.47
1:1G:397:A:N3	1:1G:397:A:H3'	2.30	0.47
46:H8:113:ALA:N	46:H8:114:GLY:HA2	2.29	0.47
26:14:196:A:C8	36:35:46:LYS:HD2	2.50	0.47
1:13:1332:A:H2'	1:13:1333:A:C8	2.50	0.47
50:H5:19:GLN:OE1	50:H5:52:HIS:NE2	2.47	0.47
33:69:117:GLU:CD	33:69:117:GLU:H	2.18	0.47
26:1H:2542:A:H4'	26:1H:2543:G:C8	2.50	0.47
26:1H:2182:G:H2'	26:1H:2183:C:C6	2.50	0.47
1:1G:791:G:C6	1:1G:792:A:N7	2.82	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:4A:35:GLU:O	13:4A:38:GLY:N	2.44	0.47
55:Q8:45:GLY:CA	55:Q8:46:ARG:C	2.83	0.46
55:Q8:53:PRO:CB	55:Q8:56:GLU:HG3	2.45	0.46
26:1H:265:A:H1'	26:1H:266:G:O4'	2.15	0.46
1:1G:826:C:H5'	8:72:12:ARG:NH1	2.30	0.46
32:51:51:ARG:HG2	32:51:52:VAL:N	2.30	0.46
26:14:2299:G:N1	26:14:2318:G:C8	2.82	0.46
29:29:23:VAL:HA	29:29:184:VAL:O	2.15	0.46
36:78:59:LEU:HD12	36:78:59:LEU:O	2.15	0.46
26:14:2472:G:C4	26:14:2475:C:N4	2.83	0.46
26:14:2279:G:N2	26:14:2280:G:H1'	2.31	0.46
8:7E:87:SER:HB2	8:7E:93:VAL:HG23	1.97	0.46
8:72:68:ARG:HH11	8:72:70:GLN:HG2	1.79	0.46
26:14:307:G:H2'	26:14:309:G:OP2	2.15	0.46
5:4E:91:LEU:HD12	5:4E:120:THR:CG2	2.45	0.46
15:6A:55:GLY:O	15:6A:59:MET:HG3	2.14	0.46
1:1G:1263:C:N3	1:1G:1273:G:N2	2.64	0.46
2:12:8:LYS:HG2	2:12:11:LEU:HD22	1.97	0.46
35:68:88:ASN:ND2	35:68:92:GLU:O	2.47	0.46
7:62:50:ILE:O	7:62:54:THR:HG23	2.15	0.46
9:82:99:LEU:HB3	9:82:101:PHE:CE1	2.49	0.46
1:13:1073:U:H2'	1:13:1074:G:C8	2.50	0.46
47:E5:36:ILE:HD11	47:E5:39:ARG:HG2	1.97	0.46
1:1G:881:G:C6	1:1G:882:C:C4	3.03	0.46
1:13:544:G:H2'	1:13:545:C:C6	2.50	0.46
36:78:38:GLN:O	36:78:41:ARG:HB2	2.15	0.46
26:1H:1260:G:H2'	26:1H:1261:C:H6	1.80	0.46
2:1E:115:LEU:HD21	2:1E:153:ARG:HD3	1.98	0.46
26:1H:2309:A:C4	26:1H:2310:A:H8	2.33	0.46
14:5I:39:LEU:HD11	14:5I:47:LEU:HD12	1.96	0.46
29:29:117:MET:HB2	29:29:122:PHE:O	2.15	0.46
26:1H:2756:U:H1'	26:1H:2757:A:H5''	1.97	0.46
12:3A:109:GLY:HA3	12:3A:121:GLY:O	2.15	0.46
26:14:1012:U:O4	34:15:25:ARG:HA	2.15	0.46
46:D5:106:GLY:HA3	46:D5:140:ASP:OD2	2.15	0.46
26:14:1590:U:H2'	26:14:1591:G:H8	1.80	0.46
23:2L:13:C:O2'	26:14:1924:C:H4'	2.16	0.46
26:1H:775:G:O5'	26:1H:777:A:H1'	2.15	0.46
26:1H:1835:G:N3	26:1H:1835:G:H2'	2.29	0.46
48:F5:20:ARG:HG2	48:F5:20:ARG:HH11	1.80	0.46
1:13:181:G:HO2'	1:13:182:U:H6	1.61	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:32:A:C2	1:13:33:A:C4	3.04	0.46
55:Q8:48:PHE:O	55:Q8:49:VAL:CG2	2.63	0.46
26:14:1786:A:C2	26:14:2606:C:H1'	2.50	0.46
26:1H:2577:A:H2'	26:1H:2614:A:N6	2.30	0.46
26:14:1049:C:O2	26:14:1113:U:H4'	2.15	0.46
22:1L:19:G:N7	26:14:882:G:H4'	2.30	0.46
22:1L:56:C:N4	22:1L:57:G:O6	2.49	0.46
26:1H:2250:G:C4	37:88:82:ARG:HG2	2.50	0.46
26:14:1019:U:OP1	26:14:1035:U:O2'	2.24	0.46
20:BI:30:LYS:CE	20:BI:80:ARG:HH12	2.29	0.46
51:M8:9:LEU:HD12	51:M8:27:THR:N	2.29	0.46
1:13:1305:G:H8	1:13:1305:G:OP2	1.97	0.46
34:58:96:GLU:O	34:58:98:VAL:N	2.44	0.46
29:29:26:ILE:HG22	29:29:27:LEU:C	2.35	0.46
34:58:42:TRP:N	41:C8:64:ARG:HD2	2.31	0.46
42:95:57:VAL:HG23	42:95:99:ILE:N	2.28	0.46
1:1G:1158:C:N4	1:1G:1181:G:H22	2.13	0.46
26:14:2052:G:H4'	29:29:143:ASN:N	2.30	0.46
1:13:627:G:H2'	1:13:628:G:C8	2.45	0.46
6:52:81:ILE:HD11	28:19:125:ILE:CG1	2.43	0.46
1:13:490:G:H2'	1:13:491:G:C8	2.50	0.46
4:32:173:TRP:CD1	4:32:174:LEU:HG	2.51	0.46
23:2L:26:C:H2'	23:2L:27:G:O4'	2.16	0.46
5:42:57:LYS:O	5:42:60:TYR:HB2	2.15	0.46
1:1G:616:G:C2	1:1G:617:G:C8	3.04	0.46
1:1G:599:C:H2'	1:1G:600:C:H6	1.81	0.46
1:1G:600:C:H2'	1:1G:601:C:H6	1.80	0.46
15:6I:71:GLN:HG2	15:6I:71:GLN:O	2.14	0.46
1:1G:1442:G:C6	1:1G:1446:A:N6	2.83	0.46
53:K5:52:VAL:HG22	53:K5:53:LYS:H	1.80	0.46
36:78:38:GLN:HG2	36:78:45:LEU:HD13	1.97	0.46
47:E5:60:PHE:N	47:E5:60:PHE:CD1	2.83	0.46
2:12:118:LEU:HB3	2:12:142:LEU:HD12	1.96	0.46
28:19:83:GLU:OE2	28:19:104:TYR:HE1	1.98	0.46
36:78:91:PHE:O	36:78:123:LEU:HD11	2.15	0.46
1:1G:811:C:H4'	1:1G:900:A:N6	2.29	0.46
26:1H:2564:A:C2	26:1H:2647:U:H4'	2.50	0.46
46:D5:10:ARG:NH2	46:D5:26:GLY:O	2.48	0.46
1:13:103:C:C2	1:13:104:G:C8	3.04	0.46
2:1E:98:LEU:HB2	2:1E:101:MET:HG3	1.96	0.46
26:14:1656:C:H2'	26:14:1657:C:H6	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:194:C:H5''	20:BA:65:LYS:HZ2	1.80	0.46
8:72:11:THR:HG23	8:72:14:ARG:HH12	1.81	0.46
54:P8:35:ARG:NH1	54:P8:42:LEU:HD11	2.30	0.46
41:85:86:ALA:HB2	41:85:116:ALA:HB2	1.97	0.46
26:14:1131:G:H4'	34:15:82:LEU:HB2	1.97	0.46
1:13:250:A:H4'	1:13:251:G:C5'	2.45	0.46
10:1A:27:ALA:HB2	10:1A:85:LEU:HD11	1.96	0.46
31:49:120:LEU:HG	31:49:179:PRO:O	2.16	0.46
26:14:2666:C:H3'	26:14:2667:C:H6	1.80	0.46
26:1H:1962:C:H4'	26:1H:1963:U:H5	1.80	0.46
50:H5:18:ASP:HA	50:H5:21:ALA:HB3	1.97	0.46
1:13:1124:G:N7	1:13:1145:C:H2'	2.30	0.46
26:14:1161:C:O2'	42:95:8:GLY:HA2	2.15	0.46
42:95:67:GLY:O	42:95:88:ARG:HG2	2.15	0.46
1:1G:976:G:H5'	1:1G:1358:U:O2'	2.15	0.46
26:14:662:G:OP1	36:35:15:ARG:NH2	2.47	0.46
26:14:933:A:C5	26:14:934:G:C8	3.03	0.46
30:39:122:LYS:HD3	30:39:191:ARG:HH21	1.80	0.46
1:13:963:G:N2	10:1I:55:LYS:NZ	2.63	0.46
1:1G:862:C:H2'	1:1G:862:C:O2	2.16	0.46
26:1H:655:A:H8	26:1H:656:G:O4'	1.98	0.46
1:1G:1300:G:HO2'	1:1G:1301:U:P	2.38	0.46
1:13:538:G:OP2	12:3I:115:LYS:HG3	2.14	0.46
5:4E:106:PRO:O	5:4E:110:LEU:HG	2.15	0.46
1:13:271:C:H2'	1:13:272:C:C6	2.50	0.46
1:1G:1127:G:N3	1:1G:1127:G:H2'	2.30	0.46
53:O8:43:CYS:HB3	53:O8:44:ARG:HH11	1.80	0.46
37:45:86:GLY:O	37:45:88:GLY:N	2.47	0.46
1:1G:536:C:H2'	1:1G:537:G:C8	2.50	0.46
45:G8:49:VAL:HG21	45:G8:55:TYR:CE2	2.50	0.46
26:1H:626:U:O4	36:78:107:LYS:HD3	2.15	0.46
26:1H:821:A:C2'	26:1H:946:G:H5''	2.45	0.46
40:75:16:ARG:HB3	40:75:18:ASP:OD1	2.15	0.46
32:51:11:VAL:HG12	32:51:12:PRO:HD2	1.97	0.46
1:13:97:U:H2'	1:13:99:C:C6	2.50	0.46
1:1G:362:G:O2'	12:3A:33:ARG:NH2	2.48	0.46
27:16:31:C:H2'	27:16:32:C:C6	2.51	0.46
46:D5:126:VAL:HA	46:D5:163:LEU:HA	1.97	0.46
26:14:2019:A:OP2	52:J5:9:LYS:NZ	2.44	0.46
26:14:2231:C:OP1	48:F5:42:GLN:HA	2.16	0.46
1:13:711:G:O2'	1:13:712:A:H5'	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:527:C:H4'	26:14:528:A:O5'	2.15	0.46
6:52:25:ILE:HD13	6:52:82:ARG:HD2	1.97	0.46
1:1G:197:A:H8	1:1G:198:G:N9	2.13	0.46
26:14:1278:A:O2'	38:55:34:ILE:HD11	2.16	0.46
34:15:35:ARG:HB2	34:15:42:TRP:CZ3	2.49	0.46
26:1H:2302:G:C6	26:1H:2315:G:C6	3.03	0.46
29:29:37:ARG:NE	29:29:42:ASP:OD2	2.48	0.46
26:14:1441:G:H2'	26:14:1442:G:C8	2.51	0.46
26:1H:2288:A:C2	26:1H:2325:G:C8	3.03	0.46
26:1H:900:A:H3'	26:1H:901:A:H8	1.81	0.46
6:52:67:MET:HB2	6:52:68:PRO:HD2	1.98	0.46
26:1H:16:G:H2'	26:1H:17:G:H8	1.80	0.46
44:F8:57:LEU:HG	44:F8:78:LYS:HG3	1.97	0.46
45:G8:20:TYR:CE1	45:G8:43:ASN:HA	2.51	0.46
2:1E:187:LEU:HA	2:1E:201:ILE:O	2.16	0.46
26:14:1342:A:C2	26:14:1397:U:C2	3.03	0.46
30:31:132:VAL:HG23	30:31:133:ASN:N	2.29	0.46
26:14:2360:A:H2'	26:14:2361:A:O4'	2.15	0.46
30:39:205:ARG:HB2	30:39:205:ARG:NH1	2.30	0.46
23:2L:32:G:C5	23:2L:33:OMC:C5	3.04	0.46
26:1H:2450:A:C2	26:1H:2451:A:C4	3.04	0.46
1:13:192:U:O4'	20:BI:103:GLY:HA2	2.15	0.46
40:B8:26:ASP:CB	40:B8:91:ARG:HA	2.42	0.46
26:1H:958:U:OP2	37:88:14:ARG:HD3	2.15	0.46
36:35:13:ASN:O	36:35:15:ARG:N	2.48	0.46
1:13:1007:C:N4	1:13:1022:G:H1	2.08	0.46
14:5A:23:ARG:HD2	14:5A:29:ARG:O	2.16	0.46
26:14:830:G:H4'	26:14:831:G:OP2	2.16	0.46
26:14:95:G:H4'	49:G5:46:GLN:HB2	1.97	0.46
26:14:631:A:H2'	26:14:632:A:O4'	2.15	0.46
26:1H:729:G:OP2	28:11:13:ARG:NH1	2.46	0.46
26:14:506:G:H5''	26:14:509:C:O2'	2.15	0.46
26:1H:1162:G:H1'	42:D8:23:GLU:OE2	2.16	0.46
5:4E:30:ALA:O	5:4E:45:PHE:HA	2.15	0.46
1:1G:1328:C:O2'	13:4A:29:ARG:NE	2.48	0.46
1:13:77:C:H1'	1:13:92:G:H22	1.79	0.46
28:19:121:PRO:HB3	28:19:135:PHE:CD2	2.50	0.46
2:1E:192:SER:OG	2:1E:193:ASP:N	2.46	0.46
3:22:195:VAL:O	3:22:196:LEU:HD22	2.16	0.46
48:F5:92:LYS:O	48:F5:93:GLU:C	2.54	0.46
7:6E:5:ARG:NH1	7:6E:7:ALA:HA	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:59:8:PRO:HG2	32:59:69:ARG:HE	1.80	0.46
26:1H:412:A:H5''	26:1H:413:C:OP2	2.15	0.46
26:14:2779:U:OP1	58:14:3844:HOH:O	2.20	0.46
1:1G:852:G:C6	1:1G:853:G:N7	2.83	0.46
26:14:873:G:N2	26:14:905:U:C2	2.84	0.46
1:13:924:C:H2'	1:13:925:G:H8	1.78	0.46
1:1G:10:A:H2'	1:1G:11:G:H8	1.79	0.46
10:1A:61:GLU:OE1	14:5A:58:LYS:HE2	2.16	0.46
44:F8:57:LEU:HD23	44:F8:57:LEU:N	2.31	0.46
11:2A:124:LYS:HG2	11:2A:124:LYS:H	1.28	0.46
26:1H:709:U:H2'	26:1H:710:G:C8	2.50	0.46
26:14:2376:A:H2'	26:14:2377:A:C8	2.50	0.46
39:65:77:ALA:O	39:65:80:LEU:HB2	2.15	0.46
28:19:18:VAL:HG12	28:19:19:ALA:N	2.31	0.46
27:16:94:C:H2'	27:16:95:U:C6	2.50	0.46
26:1H:576:U:H5	58:1H:4654:HOH:O	1.99	0.46
26:1H:2516:G:O2'	26:1H:2517:C:H5'	2.15	0.46
26:1H:866:A:H5''	26:1H:867:C:OP2	2.15	0.46
48:J8:78:LYS:H	48:J8:78:LYS:HE3	1.80	0.46
6:52:54:LYS:HE2	6:52:54:LYS:HB2	1.66	0.46
26:1H:975:G:H1'	26:1H:990:A:C2	2.50	0.46
19:AI:25:LYS:HG2	19:AI:27:GLU:OE1	2.15	0.46
36:78:135:LEU:HD13	36:78:139:LYS:NZ	2.31	0.46
55:Q8:46:ARG:HA	55:Q8:46:ARG:CZ	2.44	0.46
55:Q8:49:VAL:HG12	55:Q8:52:LYS:HG2	1.96	0.46
26:1H:2705:A:H3'	26:1H:2706:G:H8	1.80	0.46
26:1H:1019:U:O2'	26:1H:1021:A:H2	1.92	0.46
1:1G:926:G:C6	1:1G:1505:G:C6	3.03	0.46
1:1G:1324:A:H4'	1:1G:1362:C:H4'	1.98	0.46
13:4A:81:LEU:O	13:4A:89:GLY:HA3	2.16	0.46
19:AA:36:ARG:HD2	19:AA:52:TYR:O	2.16	0.46
26:1H:445:C:OP1	41:C8:2:PRO:HA	2.15	0.46
1:13:976:G:N2	1:13:1362(A):C:OP2	2.49	0.46
26:1H:1387:C:C2	26:1H:1388:G:C8	3.04	0.46
3:22:11:ARG:HH21	3:22:180:ALA:HB3	1.81	0.46
26:1H:2635:C:H5''	29:21:78:LEU:HA	1.98	0.46
26:14:2420:C:H41	55:M5:31:HIS:HB3	1.81	0.46
26:1H:2156:G:H2'	26:1H:2157:G:N3	2.30	0.46
26:14:2845:G:H2'	26:14:2846:G:C8	2.51	0.46
40:75:64:ARG:HB2	40:75:73:GLU:HG2	1.96	0.46
23:2L:56:PSU:N3	23:2L:59:A:OP2	2.36	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:2L:62:C:H2'	23:2L:63:C:H6	1.81	0.46
1:13:457:C:H2'	1:13:458:C:C6	2.49	0.46
1:13:474:G:H5''	16:7I:81:ARG:CZ	2.45	0.46
26:1H:1534:G:C2'	26:1H:1535:U:H4'	2.45	0.46
35:68:44:LYS:HD3	35:68:44:LYS:HA	1.72	0.46
5:4E:147:ASP:HA	5:4E:150:ARG:NH2	2.30	0.46
3:2E:167:TRP:C	3:2E:167:TRP:CD1	2.88	0.46
40:B8:31:SER:HB2	40:B8:84:GLN:HB3	1.96	0.46
6:52:79:LEU:HB3	6:52:88:VAL:HG21	1.98	0.46
1:1G:298:A:O5'	1:1G:298:A:H8	1.98	0.46
29:29:58:ARG:HD2	29:29:58:ARG:H	1.80	0.46
20:BA:13:LEU:O	20:BA:17:ARG:HG3	2.15	0.46
46:D5:124:ILE:HD11	46:D5:165:VAL:HG11	1.97	0.46
38:55:32:GLY:HA2	38:55:116:LEU:HD12	1.96	0.46
45:G8:40:GLU:HB2	45:G8:64:GLU:OE1	2.14	0.46
7:62:45:ASP:HB3	7:62:117:ALA:CB	2.46	0.46
2:1E:54:THR:HG21	2:1E:201:ILE:HD11	1.96	0.46
18:9A:59:SER:HB2	18:9A:62:GLU:H	1.79	0.46
26:1H:803:U:C4	26:1H:804:A:N7	2.84	0.46
19:AA:41:VAL:HG12	19:AA:42:PRO:HD2	1.97	0.46
1:1G:1401:G:C2	1:1G:1402:C:H1'	2.50	0.46
1:13:258:G:H2'	1:13:259:G:H8	1.79	0.46
26:1H:2358:G:C5	26:1H:2359:C:C5	3.04	0.46
26:1H:1232:G:C6	26:1H:1233:C:C4	3.03	0.46
37:88:59:ARG:C	37:88:61:GLY:H	2.18	0.46
26:14:1260:G:C6	26:14:1261:C:C4	3.03	0.46
1:13:1346:A:C6	7:6E:10:ARG:HD3	2.50	0.46
26:14:997:G:H2'	26:14:998:C:H6	1.81	0.46
26:14:1328:G:H2'	26:14:1330:C:C5	2.50	0.46
26:1H:1575:C:H2'	26:1H:1576:U:H6	1.77	0.46
26:1H:602:G:O2'	26:1H:655:A:N6	2.48	0.46
27:1J:5:C:N4	27:1J:115:G:H1	2.06	0.46
26:1H:1070:A:N7	26:1H:1096:A:H2'	2.31	0.46
5:4E:74:GLY:O	5:4E:115:VAL:HA	2.16	0.46
44:F8:24:GLY:HA3	44:F8:82:GLN:NE2	2.24	0.46
34:58:96:GLU:HB2	34:58:122:VAL:HG12	1.96	0.46
40:75:8:LYS:O	40:75:11:GLU:HB2	2.15	0.46
55:Q8:39:LYS:HE3	55:Q8:40:GLU:H	1.80	0.46
26:1H:2388:A:C2'	26:1H:2389:G:H5'	2.46	0.46
26:1H:2286:A:OP1	53:O8:28:ARG:NH2	2.48	0.46
1:1G:539:A:H2'	1:1G:540:G:H8	1.78	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:112:G:C2'	1:13:113:G:H5'	2.46	0.46
16:7I:74:LEU:HD22	16:7I:79:VAL:HG21	1.96	0.46
37:88:104:PHE:CE2	37:88:125:LEU:HD11	2.45	0.46
4:3E:167:GLY:HA2	28:19:135:PHE:CE1	2.47	0.46
49:K8:15:LYS:H	49:K8:67:LYS:HZ1	1.63	0.46
1:1G:360:A:H2'	1:1G:361:G:C8	2.50	0.46
39:A8:18:ILE:O	39:A8:21:THR:OG1	2.32	0.46
26:14:1951:U:N3	26:14:1954:G:OP2	2.39	0.46
26:14:1014:U:H2'	26:14:1015:G:C8	2.48	0.46
30:31:8:GLN:N	30:31:8:GLN:CD	2.68	0.46
26:14:1197:G:H2'	26:14:1198:U:C6	2.49	0.46
26:14:1431:U:H2'	26:14:1432:C:C6	2.51	0.46
27:1J:101:A:OP2	27:1J:101:A:H8	1.99	0.46
1:13:342:C:N3	1:13:348:G:C2	2.83	0.46
26:14:2142:C:H2'	26:14:2143:C:C6	2.50	0.46
26:1H:2052:G:H4'	29:21:143:ASN:O	2.16	0.46
1:1G:197:A:OP2	1:1G:197:A:H3'	2.15	0.46
26:1H:2335:A:C8	26:1H:2337:G:C6	3.04	0.46
26:1H:2841:C:H2'	26:1H:2842:G:C8	2.49	0.46
26:1H:556:G:H2'	26:1H:557:U:H6	1.80	0.46
1:13:598:U:H4'	8:7E:94:TYR:CG	2.51	0.46
8:7E:114:THR:HG22	8:7E:130:GLY:O	2.16	0.46
1:1G:659:U:N3	1:1G:660:G:N7	2.63	0.46
26:1H:370:G:H4'	26:1H:371:A:OP2	2.15	0.46
3:22:135:LYS:NZ	5:42:52:PRO:HG2	2.30	0.46
1:1G:216:G:H2'	1:1G:217:C:C6	2.50	0.46
26:1H:2235:G:H2'	26:1H:2236:C:C6	2.50	0.46
4:32:155:LEU:O	4:32:158:ILE:HD12	2.16	0.46
26:1H:1581:G:C6	26:1H:1582:C:C4	3.04	0.46
1:13:894:G:H2'	1:13:895:G:O4'	2.16	0.46
43:A5:29:LEU:HD11	43:A5:33:ARG:HE	1.81	0.46
26:1H:500:G:N2	26:1H:502:A:H3'	2.30	0.46
1:1G:790:A:H8	1:1G:790:A:O5'	1.98	0.46
39:A8:80:LEU:HA	39:A8:80:LEU:HD13	1.65	0.46
46:H8:119:GLU:CD	46:H8:119:GLU:H	2.19	0.46
26:14:442:G:C6	26:14:444:C:N4	2.84	0.46
26:14:448:U:H4'	58:14:4020:HOH:O	2.15	0.46
33:69:97:ILE:O	33:69:100:ALA:HB3	2.16	0.46
8:72:42:GLU:HG3	8:72:109:ILE:HD12	1.97	0.46
26:1H:748:G:OP2	43:E8:88:ARG:HG3	2.15	0.46
26:1H:1359:A:C2	26:1H:1372:U:C4	3.04	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
55:Q8:25:MET:HB2	55:Q8:42:ARG:HB2	1.98	0.46
2:1E:185:ILE:HA	2:1E:199:TYR:O	2.16	0.46
22:1L:18:G:O2'	22:1L:60:U:O4	2.31	0.46
26:14:1160:G:C6	26:14:1161:C:C4	3.04	0.46
1:13:1365:G:H2'	1:13:1366:C:C6	2.51	0.46
30:31:114:VAL:HG21	30:31:202:PHE:CE1	2.51	0.46
26:14:1970:A:H4'	26:14:1970:A:OP1	2.16	0.46
46:D5:29:TYR:CE2	46:D5:87:ASP:HB2	2.51	0.46
41:85:88:ILE:HG22	41:85:90:VAL:HG23	1.98	0.46
45:C5:19:LYS:CG	45:C5:20:TYR:H	2.17	0.46
26:1H:2261:C:H1'	26:1H:2388:A:N3	2.31	0.46
26:14:2419:U:O4	55:M5:31:HIS:CD2	2.69	0.46
26:14:2845:G:N2	26:14:2871:C:O2	2.43	0.46
45:G8:52:SER:OG	45:G8:53:PRO:HD2	2.15	0.46
16:7I:49:LEU:HD12	16:7I:50:LYS:H	1.81	0.46
26:1H:2602:A:H4'	26:1H:2603:G:O5'	2.15	0.46
26:14:1599:C:H2'	26:14:1600:C:C6	2.47	0.46
26:14:1599:C:C5	26:14:1600:C:H5	2.34	0.46
26:1H:182:A:H2'	26:1H:183:C:H6	1.81	0.46
33:69:41:GLU:O	33:69:45:LYS:HG2	2.16	0.46
31:49:60:LEU:HD22	31:49:68:PRO:HB3	1.98	0.46
32:51:8:PRO:O	32:51:10:PRO:HD3	2.16	0.46
2:12:209:ARG:HD3	2:12:240:GLN:HE22	1.79	0.46
26:14:717:G:H2'	26:14:718:A:O4'	2.16	0.46
26:1H:1749:A:C4	26:1H:1750:G:C8	3.03	0.46
40:75:85:LYS:HD2	40:75:87:ASP:OD1	2.16	0.46
26:14:1011:G:H2'	26:14:1013:C:O4'	2.15	0.46
26:14:1857:G:O2'	26:14:1885:A:N6	2.48	0.46
26:1H:389:G:H8	26:1H:389:G:O5'	1.99	0.46
26:1H:1259:G:H2'	26:1H:1260:G:H8	1.80	0.46
26:14:548:A:H2'	26:14:549:G:O4'	2.16	0.46
7:6E:26:PHE:CE2	7:6E:30:ILE:HD11	2.50	0.46
26:14:26:G:OP1	43:A5:80:PRO:HB3	2.16	0.46
26:1H:1972:A:H2'	26:1H:1973:G:H8	1.81	0.46
26:14:196:A:O4'	36:35:46:LYS:HE2	2.15	0.46
26:1H:2844:G:O6	58:1H:4096:HOH:O	2.20	0.46
26:14:1104:C:H2'	26:14:1105:U:C6	2.49	0.46
47:E5:34:GLY:HA2	47:E5:61:ALA:O	2.16	0.46
26:1H:749:C:H5''	58:1H:4461:HOH:O	2.14	0.46
16:7A:43:LYS:HB2	16:7A:43:LYS:HE2	1.62	0.46
1:1G:236:G:H1'	17:8A:4:LYS:NZ	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:17:U:H2'	1:13:18:C:C6	2.51	0.46
32:51:85:LYS:HA	32:51:85:LYS:HD3	1.44	0.46
1:1G:1530:G:H5''	1:1G:1530:G:H8	1.81	0.46
45:G8:88:LYS:HA	45:G8:88:LYS:HD3	1.69	0.46
33:69:2:LYS:H	33:69:2:LYS:HG2	1.39	0.46
1:13:1186:G:H21	14:5I:61:TRP:C	2.19	0.46
26:14:1115:G:H2'	26:14:1116:C:C6	2.51	0.46
33:69:77:LEU:HA	33:69:141:LYS:HB3	1.98	0.46
24:3L:76:A:O2'	26:14:2394:C:N3	2.45	0.46
31:41:67:LYS:NZ	51:M8:6:HIS:CE1	2.84	0.46
29:29:181:LEU:HA	29:29:181:LEU:HD12	1.77	0.46
26:1H:903:C:O2'	46:H8:169:GLU:OE1	2.34	0.46
42:D8:9:GLY:O	42:D8:10:LYS:HG3	2.16	0.46
26:1H:239:U:H2'	26:1H:240:G:O4'	2.16	0.46
7:6E:91:VAL:HB	7:6E:96:GLN:HG2	1.98	0.46
1:13:403:C:H4'	4:3E:122:ARG:NH1	2.30	0.46
1:1G:345:C:HO2'	1:1G:346:G:P	2.39	0.46
1:1G:973:G:O3'	14:5A:41:ARG:NH2	2.45	0.46
26:14:1421:G:C2	26:14:1422:G:C8	3.04	0.46
26:14:1569:A:H2'	26:14:1570:A:O4'	2.15	0.46
4:32:159:ARG:HG3	4:32:159:ARG:H	1.55	0.46
1:13:42:G:H2'	1:13:43:C:O4'	2.16	0.46
3:22:136:GLN:O	3:22:139:GLN:N	2.49	0.46
26:14:492:A:H2'	26:14:493:G:O4'	2.15	0.46
39:65:11:LYS:HE3	39:65:15:ARG:HH22	1.80	0.46
26:1H:289:A:H2'	26:1H:290:G:O4'	2.15	0.46
9:82:42:ARG:NH1	9:82:71:SER:O	2.48	0.46
1:1G:15:G:H2'	1:1G:16:A:C8	2.51	0.46
1:13:730:G:C5	1:13:731:G:H1'	2.51	0.46
1:13:713:G:H2'	1:13:714:G:C8	2.51	0.46
38:98:13:HIS:CD2	38:98:13:HIS:H	2.33	0.46
1:1G:1512:U:C2	1:1G:1513:A:C8	3.04	0.46
2:1E:209:ARG:HG3	2:1E:240:GLN:NE2	2.30	0.46
26:14:1598:C:O3'	44:B5:35:THR:HG22	2.15	0.46
11:2A:72:ALA:HB1	11:2A:77:MET:HE3	1.98	0.46
1:1G:448:A:H2'	1:1G:449:C:O2	2.16	0.46
40:B8:45:PHE:CZ	40:B8:65:LYS:HG2	2.50	0.46
1:1G:922:G:C2	1:1G:923:A:C4	3.04	0.46
51:I5:42:PHE:O	51:I5:43:TYR:HB3	2.15	0.46
32:51:86:GLU:HG3	32:51:165:ALA:HB3	1.98	0.46
26:1H:356:G:H2'	26:1H:357:A:C8	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:901:A:H5'	26:1H:902:C:OP2	2.16	0.46
26:1H:2168:G:OP1	26:1H:2168:G:H4'	2.16	0.46
1:1G:25:C:H2'	1:1G:26:A:H8	1.81	0.46
5:4E:153:LYS:HG3	5:4E:154:GLY:H	1.80	0.46
26:1H:2102:U:H2'	26:1H:2103:C:C6	2.51	0.46
26:1H:2104:G:N1	26:1H:2186:G:C2	2.84	0.46
16:7I:21:VAL:O	16:7I:33:ILE:N	2.44	0.46
29:21:103:ASP:OD2	29:21:168:MET:HG2	2.15	0.46
33:61:55:ALA:HA	33:61:58:LEU:HB3	1.98	0.46
26:1H:2549:G:H5''	26:1H:2549:G:H8	1.81	0.46
11:2I:125:PHE:CD1	11:2I:125:PHE:N	2.82	0.46
52:N8:9:LYS:HA	52:N8:9:LYS:HD3	1.58	0.46
26:1H:236:C:H2'	26:1H:237:C:H6	1.81	0.46
38:55:58:GLY:HA2	38:55:80:PHE:CE2	2.50	0.46
26:1H:2712:U:H1'	26:1H:2712(A):A:C8	2.51	0.46
13:4I:67:GLU:OE2	13:4I:68:GLY:N	2.48	0.46
1:13:1139:G:H4'	1:13:1140:C:C5'	2.46	0.46
1:1G:1320:C:O2	19:AA:36:ARG:NH2	2.49	0.46
27:1J:109:G:C5	27:1J:110:G:N7	2.84	0.46
1:1G:998:G:H22	1:1G:1043:C:H42	1.64	0.46
2:1E:69:LEU:HD13	2:1E:92:TYR:HA	1.98	0.46
33:61:40:THR:O	33:61:44:LEU:HB2	2.16	0.46
26:14:1889:A:N1	26:14:2234:G:H1'	2.31	0.46
46:H8:75:ASN:O	46:H8:84:GLU:N	2.41	0.46
1:1G:37:U:O2'	1:1G:547:A:N1	2.42	0.46
24:3K:57:G:O5'	24:3K:57:G:H8	1.99	0.46
34:58:132:ALA:O	34:58:134:ARG:NH2	2.49	0.46
26:14:2419:U:H5'	53:K5:23:THR:HG21	1.97	0.46
39:A8:95:HIS:O	39:A8:98:VAL:HG23	2.16	0.46
37:88:87:LYS:HA	37:88:87:LYS:HD3	1.70	0.46
27:16:71:C:C2	27:16:72:G:C8	3.03	0.46
26:1H:907:U:O2'	37:88:101:ARG:NH2	2.46	0.46
1:13:1218:C:OP2	14:5I:9:LYS:NZ	2.39	0.46
3:22:91:LEU:CB	3:22:99:VAL:HG11	2.45	0.46
1:1G:438:G:H5'	4:32:123:HIS:HD2	1.81	0.46
26:1H:2590:A:H2'	26:1H:2591:C:H6	1.79	0.46
26:1H:270(J):G:H3'	26:1H:270(K):C:H6	1.81	0.46
1:13:712:A:H2'	1:13:713:G:C8	2.50	0.46
4:3E:142:PRO:HB3	4:3E:186:LEU:O	2.16	0.46
30:31:40:GLN:HE22	30:31:184:TYR:HB3	1.81	0.46
20:BA:67:ALA:HA	20:BA:73:HIS:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:4I:84:ILE:HG23	13:4I:86:CYS:H	1.79	0.46
22:1K:26:A:H61	22:1K:44:G:N2	2.13	0.46
32:51:86:GLU:HB2	32:51:165:ALA:HB3	1.97	0.46
1:1G:259:G:H2'	1:1G:260:G:O4'	2.15	0.46
1:13:1298:C:H2'	7:6E:114:ARG:HH21	1.81	0.46
31:49:27:ASN:HB3	31:49:30:GLU:HG3	1.98	0.46
26:1H:2436:G:C5	26:1H:2437:U:C5	3.04	0.46
26:1H:2436:G:C6	26:1H:2437:U:C4	3.03	0.46
8:72:64:LYS:O	8:72:79:VAL:HB	2.15	0.46
15:6A:17:ARG:HD3	15:6A:26:GLU:HG3	1.98	0.46
26:14:2048:G:N7	58:14:3562:HOH:O	2.36	0.46
1:1G:624:C:O3'	16:7A:10:GLY:HA2	2.16	0.46
16:7A:8:ARG:HD3	16:7A:17:TYR:CE1	2.51	0.46
1:13:358:U:H2'	1:13:359:U:O4'	2.16	0.46
1:1G:134:A:H61	16:7A:25:ARG:NH1	2.13	0.46
29:21:27:LEU:HD22	40:B8:1:MET:HE3	1.98	0.46
26:14:300:A:N6	58:14:3654:HOH:O	2.38	0.46
26:14:212:G:H2'	26:14:213:A:O4'	2.16	0.46
1:13:891:U:H2'	1:13:892:A:H8	1.80	0.46
26:14:1339:G:H21	26:14:1603:A:H1'	1.81	0.46
32:51:42:ARG:NH1	32:51:42:ARG:HB2	2.30	0.46
23:2K:35:C:H5''	23:2K:36:A:OP2	2.16	0.46
4:32:12:CYS:SG	4:32:18:LYS:HG3	2.56	0.46
26:14:2016:U:H2'	26:14:2017:U:H6	1.79	0.46
2:1E:104:ASN:ND2	2:1E:107:THR:HB	2.13	0.46
26:1H:2317:C:H2'	26:1H:2318:G:H5'	1.98	0.46
1:13:1133:G:N2	1:13:1141:C:N3	2.56	0.46
26:14:1225:C:H4'	42:95:85:LYS:HG2	1.98	0.46
1:1G:1323:G:H2'	1:1G:1324:A:C8	2.50	0.46
1:13:1023:G:H3'	1:13:1024:G:C5'	2.42	0.46
26:1H:631:A:H61	26:1H:2402:C:H42	1.64	0.46
26:1H:1591:G:O2'	26:1H:1592:C:H5'	2.15	0.46
5:42:31:LEU:HD22	5:42:43:LEU:HD11	1.98	0.46
31:41:62:LEU:HD12	31:41:62:LEU:HA	1.68	0.46
26:1H:66:C:H2'	26:1H:67:U:C6	2.51	0.46
26:1H:654(C):G:H2'	26:1H:654(D):G:O4'	2.15	0.46
26:14:2781:A:H5''	26:14:2782:G:H5'	1.98	0.46
26:1H:2801:A:H2'	26:1H:2802:G:C4'	2.46	0.46
38:98:117:VAL:HG13	38:98:118:GLU:N	2.31	0.46
45:C5:20:TYR:CZ	45:C5:42:VAL:HA	2.51	0.46
26:14:952:G:C6	26:14:966:G:C6	3.04	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2053:G:P	58:1H:3801:HOH:O	2.74	0.46
37:45:22:LYS:HG2	37:45:23:GLY:CA	2.42	0.46
24:3K:24:G:C6	24:3K:25:C:N4	2.84	0.46
24:3K:25:C:N4	24:3K:26:A:C8	2.84	0.46
18:9A:61:LYS:O	18:9A:65:ILE:HG22	2.15	0.46
1:13:300:A:H1'	1:13:565:U:O2	2.16	0.46
26:1H:155:C:H5'	26:1H:161:U:OP2	2.16	0.46
5:42:68:GLU:OE2	5:42:70:PRO:HG3	2.16	0.46
1:13:435:C:H2'	1:13:436:C:H6	1.81	0.46
26:14:2068:U:N3	26:14:2430:A:C2	2.76	0.46
1:13:536:C:H2'	1:13:537:G:H8	1.77	0.46
26:14:2231:C:H2'	26:14:2232:U:O4'	2.16	0.46
9:8E:9:ARG:HB3	9:8E:14:VAL:HG22	1.97	0.46
1:13:819:A:H4'	1:13:820:U:OP2	2.16	0.46
3:22:70:VAL:HG11	3:22:76:VAL:CG1	2.46	0.46
32:51:105:LEU:HD23	32:51:105:LEU:N	2.31	0.46
13:4I:82:MET:HG2	13:4I:93:ARG:HG3	1.96	0.46
46:H8:146:ILE:HA	46:H8:174:VAL:O	2.15	0.46
26:14:863:A:H2'	26:14:864:G:C8	2.51	0.46
6:5E:33:TYR:HB2	6:5E:75:LEU:HG	1.98	0.46
27:16:116:G:H2'	27:16:117:G:O4'	2.15	0.46
45:G8:28:LYS:HD2	45:G8:40:GLU:HB3	1.97	0.46
26:1H:1709:U:H1'	26:1H:2860:A:N3	2.31	0.46
27:1J:28:C:H42	27:1J:56:G:H1	1.62	0.46
1:1G:1016:A:H3'	1:1G:1017:G:C8	2.52	0.46
1:13:1298:C:P	7:6E:114:ARG:HH22	2.39	0.46
26:1H:2280:G:C2'	26:1H:2281:C:H5'	2.46	0.46
1:13:382:A:H2'	1:13:383:A:C8	2.51	0.46
26:14:384:U:H2'	26:14:385:C:H6	1.80	0.46
8:72:89:PRO:HA	8:72:92:ARG:HE	1.81	0.46
28:11:232:PRO:HB3	28:11:244:ARG:CZ	2.46	0.46
1:1G:570:G:H1'	1:1G:820:U:C4	2.51	0.46
1:1G:578:C:OP1	58:1G:1866:HOH:O	2.20	0.46
26:1H:1161:C:H1'	42:D8:8:GLY:O	2.16	0.46
26:14:1153:C:OP1	41:85:93:LYS:NZ	2.47	0.46
8:72:39:LEU:HD12	8:72:39:LEU:HA	1.78	0.46
7:6E:104:LEU:HD13	7:6E:104:LEU:HA	1.81	0.46
30:39:196:LEU:HA	30:39:196:LEU:HD22	1.82	0.46
30:39:41:LEU:HD23	30:39:41:LEU:HA	1.67	0.46
46:H8:80:ARG:H	46:H8:80:ARG:HG2	1.60	0.46
33:61:81:VAL:HG21	33:61:88:ILE:HD13	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:19:132:PRO:HG3	28:19:190:TYR:CE1	2.51	0.46
26:14:1786:A:H2	26:14:2606:C:H1'	1.81	0.45
26:14:1049:C:H42	26:14:2751:G:H1	1.64	0.45
46:H8:60:GLU:HB3	46:H8:61:LEU:H	1.46	0.45
26:14:847:U:H5'	26:14:848:G:OP2	2.16	0.45
1:1G:826:C:H2'	1:1G:827:U:C6	2.51	0.45
1:1G:857:C:H2'	1:1G:858:G:O4'	2.16	0.45
49:G5:31:GLU:HB2	49:G5:53:LEU:HD11	1.98	0.45
26:14:2393:A:C2'	26:14:2394:C:H5'	2.45	0.45
36:35:63:PRO:HA	55:M5:13:ARG:HA	1.98	0.45
26:14:2212:A:H1'	26:14:2215:G:C5	2.51	0.45
55:Q8:35:GLN:C	55:Q8:37:SER:H	2.15	0.45
31:41:63:ILE:HD12	31:41:141:PHE:CE2	2.51	0.45
26:1H:1437:C:C2	26:1H:1438:U:C5	3.04	0.45
32:59:54:ARG:NH2	32:59:57:ASP:OD1	2.36	0.45
1:1G:1157:A:H4'	1:1G:1158:C:O5'	2.17	0.45
11:2A:85:ARG:HA	11:2A:112:THR:OG1	2.16	0.45
26:14:17:G:H4'	41:85:25:TRP:CH2	2.52	0.45
43:E8:24:ILE:HD12	43:E8:24:ILE:O	2.16	0.45
31:41:99:MET:HG3	31:41:100:TRP:N	2.31	0.45
1:13:142:G:H2'	1:13:143:A:C8	2.50	0.45
4:32:153:ARG:NH1	4:32:181:MET:HB2	2.31	0.45
1:1G:90:C:H2'	1:1G:91:C:C6	2.51	0.45
25:4K:13:A:O2'	25:4K:14:A:OP1	2.33	0.45
1:1G:618:C:H5'	1:1G:619:U:H5''	1.98	0.45
26:1H:2590:A:H2'	26:1H:2591:C:C6	2.52	0.45
1:13:1073:U:H2'	1:13:1074:G:H8	1.81	0.45
5:4E:11:ILE:HD11	5:4E:31:LEU:HD22	1.98	0.45
26:14:443:A:H2	26:14:1245:G:N3	2.14	0.45
26:14:1011:G:C2	26:14:1151:G:C2	3.04	0.45
1:13:1329:A:N7	21:1F:7:ARG:NH2	2.64	0.45
1:13:544:G:C5	1:13:545:C:C5	3.04	0.45
1:13:186(E):C:N4	1:13:191(B):G:H1	2.15	0.45
26:14:2002:G:C6	58:14:3797:HOH:O	2.69	0.45
47:I8:23:VAL:HG13	47:I8:38:VAL:HG23	1.97	0.45
28:11:119:ALA:CB	28:11:130:ALA:HB3	2.46	0.45
8:7E:25:ASP:OD1	8:7E:60:ARG:HG3	2.16	0.45
26:14:2346:A:C2	26:14:2383:G:C2	3.04	0.45
41:C8:75:ASN:HB2	41:C8:78:THR:OG1	2.16	0.45
26:14:1005:C:O2'	34:15:28:THR:HG21	2.16	0.45
26:1H:2114:A:H2'	26:1H:2168:G:H5''	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:1J:99:A:H3'	58:1J:309:HOH:O	2.16	0.45
8:72:114:THR:OG1	8:72:115:SER:N	2.49	0.45
1:13:105:G:H2'	1:13:106:C:C6	2.50	0.45
41:85:21:ALA:HA	41:85:24:TYR:CE2	2.51	0.45
26:1H:483:A:O2'	45:G8:59:GLY:HA2	2.16	0.45
1:13:952:U:O4	13:4I:104:ARG:HD3	2.16	0.45
31:41:98:ARG:HA	31:41:101:ILE:HG23	1.97	0.45
47:18:53:MET:HG3	47:18:59:LEU:CD2	2.46	0.45
38:98:28:LEU:HA	38:98:28:LEU:HD23	1.77	0.45
3:22:27:LYS:HB2	3:22:28:GLN:NE2	2.31	0.45
36:78:86:LYS:HB3	36:78:118:GLY:CA	2.45	0.45
26:1H:989:G:N7	50:L8:13:ILE:HD11	2.31	0.45
1:1G:801:U:H2'	1:1G:802:A:C8	2.51	0.45
13:4A:13:LYS:NZ	13:4A:21:TYR:OH	2.49	0.45
26:1H:1358:G:N2	26:1H:1372:U:C5	2.84	0.45
42:95:85:LYS:HG3	42:95:87:HIS:CA	2.46	0.45
13:4A:78:ILE:HA	13:4A:81:LEU:HD12	1.98	0.45
1:1G:1129:C:N4	1:1G:1142:G:O6	2.35	0.45
1:13:741:G:H2'	1:13:742:G:O4'	2.16	0.45
30:31:23:ASP:CG	30:31:24:LEU:H	2.19	0.45
1:13:825:G:C6	1:13:826:C:C4	3.05	0.45
30:39:27:GLU:O	30:39:28:ILE:HG12	2.16	0.45
26:1H:674:G:H1'	30:31:74:ARG:HH11	1.81	0.45
26:1H:1170:G:N2	26:1H:1180:C:O2	2.50	0.45
26:1H:654(D):G:H1	26:1H:654(Q):C:N4	2.08	0.45
26:14:1060:U:H5''	26:14:1061:U:C6	2.50	0.45
26:1H:2418:A:OP2	55:Q8:39:LYS:NZ	2.41	0.45
16:7A:49:LEU:HD12	16:7A:50:LYS:N	2.31	0.45
9:8E:118:LYS:O	9:8E:119:ALA:HB3	2.17	0.45
24:3K:19:G:O2'	24:3K:57:G:N3	2.47	0.45
39:A8:47:THR:HG22	39:A8:48:LEU:N	2.31	0.45
26:14:972:G:OP2	26:14:974:G:H5''	2.16	0.45
1:1G:591:U:OP2	8:72:30:ARG:HD3	2.17	0.45
26:14:784:A:C8	26:14:792:G:C5	3.05	0.45
1:1G:1086:U:H3	1:1G:1099:G:H22	1.64	0.45
26:1H:1864:U:OP1	26:1H:2410:G:O2'	2.19	0.45
26:1H:2531:A:H5'	32:51:157:TYR:CZ	2.52	0.45
1:13:1057:G:C5	1:13:1204:A:C2	3.04	0.45
1:13:41:G:H2'	1:13:42:G:C8	2.51	0.45
26:14:2805:G:H2'	26:14:2807:G:H8	1.81	0.45
26:1H:1412:A:H2'	26:1H:1413:G:H8	1.77	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:433:C:H2'	1:1G:434:U:H6	1.78	0.45
27:1J:87:G:N2	27:1J:89:G:H3'	2.31	0.45
35:25:3:GLN:HB2	35:25:4:PRO:HD2	1.97	0.45
20:BI:49:ALA:HB1	20:BI:99:LEU:HB2	1.98	0.45
26:14:150:C:H42	26:14:176:G:H1	1.63	0.45
42:D8:21:ARG:HG2	42:D8:91:TYR:HE2	1.80	0.45
26:1H:2590:A:C2	26:1H:2605:U:C2	3.05	0.45
26:14:1093:G:N1	26:14:1097:U:OP2	2.48	0.45
26:1H:1748:G:H2'	26:1H:1749:A:H8	1.81	0.45
26:14:747:U:O2	26:14:2014:A:H1'	2.16	0.45
26:14:2882:A:H5'	38:55:96:ARG:HG3	1.97	0.45
26:14:1164:G:H1	26:14:1185:C:H42	1.63	0.45
28:11:70:TRP:CD1	28:11:70:TRP:C	2.90	0.45
7:6E:15:ASP:O	7:6E:19:GLY:HA2	2.15	0.45
28:19:94:LEU:HA	28:19:94:LEU:HD23	1.77	0.45
33:69:29:TYR:C	33:69:32:PRO:HD2	2.37	0.45
38:55:79:LEU:HA	38:55:83:ILE:HB	1.98	0.45
9:82:95:LYS:HE2	9:82:95:LYS:HB2	1.77	0.45
9:82:53:VAL:HG22	9:82:95:LYS:HZ2	1.81	0.45
30:31:174:VAL:CG1	30:31:189:THR:HG21	2.47	0.45
26:14:2666:C:H3'	26:14:2667:C:C6	2.51	0.45
23:2K:67:C:H2'	23:2K:68:C:O4'	2.15	0.45
30:39:200:GLU:O	30:39:203:GLN:HB2	2.16	0.45
1:13:1113:C:H2'	1:13:1114:C:H6	1.81	0.45
1:1G:661:G:H1	1:1G:744:C:H42	1.64	0.45
29:29:100:GLU:O	29:29:172:VAL:HG23	2.16	0.45
15:6A:4:THR:OG1	15:6A:7:GLU:HB2	2.17	0.45
1:1G:1517:G:C6	1:1G:1518:A:C5	3.04	0.45
16:7A:72:ARG:HD2	16:7A:73:LEU:HD23	1.98	0.45
51:M8:15:ILE:HG22	51:M8:16:CYS:H	1.81	0.45
50:L8:10:LYS:HB3	50:L8:53:LEU:HD23	1.98	0.45
26:1H:1776:G:OP2	58:1H:3656:HOH:O	2.21	0.45
5:42:127:ASN:OD1	5:42:130:ASN:ND2	2.47	0.45
29:21:49:LEU:HD12	29:21:49:LEU:HA	1.72	0.45
34:58:10:GLU:HG3	34:58:11:PRO:HD2	1.99	0.45
8:7E:42:GLU:HG3	8:7E:109:ILE:HD12	1.98	0.45
26:14:1813:G:H1'	28:19:50:THR:OG1	2.16	0.45
55:Q8:21:LYS:HA	55:Q8:21:LYS:HD2	1.70	0.45
55:Q8:22:VAL:HG22	55:Q8:47:LYS:NZ	2.31	0.45
26:14:1225:C:H4'	42:95:85:LYS:CG	2.45	0.45
1:13:973:G:OP1	10:1I:57:LYS:HD3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
30:31:119:ARG:HB3	30:31:119:ARG:HE	1.58	0.45
30:31:6:VAL:N	30:31:24:LEU:HB3	2.31	0.45
29:29:5:LEU:HD23	29:29:51:PHE:HB2	1.98	0.45
22:1L:72:C:H2'	22:1L:73:A:H5''	1.99	0.45
19:AI:78:ARG:NE	19:AI:78:ARG:H	2.04	0.45
55:Q8:34:TRP:CB	55:Q8:35:GLN:CA	2.92	0.45
38:98:72:ASP:OD2	38:98:75:LEU:HB2	2.16	0.45
15:6I:6:GLU:CD	15:6I:6:GLU:N	2.69	0.45
10:1A:35:SER:HB3	10:1A:73:ASP:HB2	1.97	0.45
1:13:685:G:O2'	1:13:686:U:H5'	2.16	0.45
1:13:1020:U:H2'	1:13:1021:G:C8	2.52	0.45
8:7E:85:ARG:HD3	8:7E:86:ILE:O	2.16	0.45
39:65:31:SER:OG	39:65:34:HIS:N	2.48	0.45
27:1J:116:G:C5'	39:65:55:ALA:HB2	2.46	0.45
26:14:307:G:H21	26:14:330:A:H62	1.63	0.45
26:14:311:A:C6	26:14:328:U:C4	3.05	0.45
26:1H:2690:C:H5''	26:1H:2872:G:N2	2.29	0.45
39:A8:11:LYS:O	39:A8:15:ARG:HB2	2.16	0.45
19:AI:40:ILE:O	19:AI:41:VAL:HG22	2.16	0.45
40:75:19:LEU:HB3	40:75:86:ILE:CG2	2.46	0.45
30:31:32:LEU:HD22	30:31:105:VAL:HG13	1.98	0.45
1:13:515:G:N7	58:13:1869:HOH:O	2.36	0.45
1:13:724:G:O2'	1:13:725:G:H5'	2.16	0.45
30:39:63:LYS:HE2	30:39:67:GLN:HB3	1.98	0.45
3:22:37:GLN:O	3:22:40:ARG:HG3	2.17	0.45
51:I5:37:SER:OG	51:I5:38:LYS:N	2.49	0.45
26:14:2762:G:H5'	26:14:2763:G:OP2	2.15	0.45
26:14:906:G:OP1	37:45:140:ALA:HB3	2.16	0.45
26:14:690:G:O2'	28:19:43:ARG:NH2	2.49	0.45
1:13:1312:G:H5'	19:AI:6:LYS:CD	2.46	0.45
1:13:448:A:OP2	1:13:485:G:N2	2.33	0.45
37:45:19:GLY:O	37:45:99:PRO:HD2	2.16	0.45
26:14:1751:C:O2'	26:14:1752:C:H5'	2.15	0.45
2:12:105:PHE:HA	2:12:108:ILE:HB	1.97	0.45
26:14:342:G:H2'	26:14:343:C:H6	1.81	0.45
1:1G:1072:G:H2'	1:1G:1073:U:C6	2.52	0.45
26:1H:2111:C:O2'	26:1H:2119:A:OP1	2.34	0.45
6:5E:62:TRP:CE2	18:9I:35:ARG:NH2	2.85	0.45
26:14:1131:G:C2	26:14:1132:A:C4	3.05	0.45
26:14:1648:C:H42	26:14:2009:G:H1	1.64	0.45
17:8A:87:LYS:O	17:8A:91:ARG:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:25:10:VAL:CG1	35:25:19:ILE:HG23	2.46	0.45
46:D5:24:LEU:HD12	46:D5:25:PRO:O	2.16	0.45
10:1I:17:ASP:O	10:1I:21:GLN:HG2	2.16	0.45
8:72:82:HIS:HB3	8:72:138:TRP:CZ3	2.51	0.45
26:14:484:C:H2'	26:14:485:C:H6	1.82	0.45
17:8A:3:LYS:HB3	17:8A:61:GLU:HB3	1.98	0.45
29:29:13:ARG:HA	29:29:21:VAL:O	2.16	0.45
26:14:1235:G:C6	26:14:1236:G:N1	2.84	0.45
26:14:1364:G:OP2	48:F5:2:SER:N	2.48	0.45
29:21:134:ILE:C	29:21:134:ILE:HD12	2.36	0.45
2:1E:36:ARG:HD2	2:1E:36:ARG:HA	1.81	0.45
48:F5:82:LEU:HB2	48:F5:83:GLU:H	1.59	0.45
45:C5:75:ILE:HG23	45:C5:76:CYS:N	2.32	0.45
26:14:2751:G:N1	32:59:2:SER:HB3	2.30	0.45
1:1G:926:G:C6	1:1G:1505:G:C5	3.04	0.45
26:1H:881:G:H5'	26:1H:882:G:O5'	2.15	0.45
1:13:22:G:H2'	1:13:23:C:H6	1.80	0.45
7:62:26:PHE:CD2	7:62:30:ILE:HD11	2.52	0.45
1:13:1028:C:H42	1:13:1033:G:H1	1.64	0.45
26:14:2074:U:P	58:14:3403:HOH:O	2.70	0.45
22:1L:73:A:O3'	22:1L:74:C:H3'	2.17	0.45
1:13:263:A:OP2	20:BI:79:ARG:NH1	2.49	0.45
32:51:4:ILE:C	32:51:6:ARG:H	2.19	0.45
29:21:47:VAL:O	29:21:80:GLU:HA	2.16	0.45
1:13:468:A:O2'	16:7I:82:GLN:HG2	2.17	0.45
24:3L:9:A:OP1	24:3L:46:G:O2'	2.29	0.45
1:1G:8:A:N6	4:32:209:ARG:HB2	2.30	0.45
26:1H:1204:A:N1	26:1H:1241:A:H2	2.14	0.45
26:1H:1888:G:H5'	26:1H:1889:A:OP1	2.16	0.45
23:2K:77:A:O2'	26:1H:2602:A:N7	2.39	0.45
28:11:136:ILE:HG12	28:11:136:ILE:H	1.61	0.45
1:13:540:G:H2'	1:13:541:G:O4'	2.16	0.45
26:14:2453:A:H2'	26:14:2454:G:O4'	2.16	0.45
7:62:21:VAL:O	7:62:24:THR:HB	2.17	0.45
26:1H:1500:G:O2'	28:11:100:GLY:O	2.30	0.45
2:1E:189:ASP:HB3	2:1E:191:ASP:H	1.82	0.45
30:39:132:VAL:HG22	30:39:133:ASN:OD1	2.16	0.45
29:29:195:LEU:HD12	29:29:195:LEU:HA	1.76	0.45
1:1G:1228:C:H5'	13:4A:114:ARG:HB3	1.99	0.45
1:13:1074:G:C2	1:13:1075:C:C2	3.05	0.45
26:1H:270(K):C:O2'	26:1H:270(M):U:OP1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:55:29:LEU:HD23	38:55:70:LEU:HD11	1.98	0.45
26:1H:1471:A:C2	26:1H:1472:A:C8	3.04	0.45
6:5E:100:ASN:ND2	18:9I:26:LEU:O	2.49	0.45
26:1H:1108:U:O4	26:1H:1109:C:N4	2.49	0.45
2:1E:210:SER:O	2:1E:214:ILE:HG12	2.16	0.45
19:AI:51:VAL:HG12	19:AI:52:TYR:N	2.31	0.45
1:13:555:C:OP2	12:3I:20:LYS:NZ	2.48	0.45
3:2E:149:ALA:HA	3:2E:201:TYR:O	2.15	0.45
39:65:102:ALA:O	39:65:105:ALA:N	2.49	0.45
1:13:222:U:C2	1:13:223:U:C5	3.05	0.45
26:1H:721:C:H2'	26:1H:722:A:C8	2.51	0.45
19:AA:12:ASP:OD1	19:AA:12:ASP:N	2.47	0.45
1:13:153:C:H42	1:13:168:G:H1	1.63	0.45
26:1H:29:U:H2'	26:1H:30:G:H8	1.82	0.45
29:29:128:SER:OG	29:29:129:HIS:N	2.49	0.45
26:1H:2859:G:C6	26:1H:2860:A:N6	2.85	0.45
1:1G:1072:G:C2	1:1G:1073:U:C2	3.05	0.45
1:1G:1198:G:H2'	1:1G:1199:U:C6	2.51	0.45
26:1H:2542:A:H4'	26:1H:2543:G:H8	1.80	0.45
26:14:1196:C:O2'	26:14:1228:G:O2'	2.27	0.45
13:4A:7:VAL:HG11	31:49:115:ARG:HE	1.81	0.45
1:1G:646:U:H2'	1:1G:647:C:C6	2.50	0.45
26:14:2:G:N2	26:14:2901:C:H42	2.14	0.45
8:7E:38:ILE:HD11	8:7E:118:VAL:O	2.17	0.45
26:1H:2208:U:H4'	28:11:151:LYS:HG2	1.97	0.45
26:14:1341:U:O4	44:B5:16:LYS:NZ	2.40	0.45
31:49:81:LYS:HB3	31:49:82:LEU:H	1.55	0.45
1:13:1513:A:H2'	1:13:1514:C:C6	2.51	0.45
1:13:1164:G:C6	1:13:1165:C:C4	3.04	0.45
33:61:79:ILE:HD13	33:61:79:ILE:HA	1.73	0.45
26:1H:1312:U:H4'	26:1H:1313:U:O5'	2.17	0.45
36:78:94:GLU:OE1	36:78:124:LYS:HD3	2.17	0.45
22:1K:74:C:H1'	22:1K:75:C:H5'	1.96	0.45
1:1G:406:G:H5'	4:32:5:ILE:HG22	1.97	0.45
1:13:1399:C:O2	58:13:1909:HOH:O	2.18	0.45
12:3I:84:LEU:HB2	12:3I:105:TYR:CE2	2.52	0.45
27:1J:45:A:N3	27:1J:45:A:H2'	2.30	0.45
1:1G:411:A:N7	1:1G:413:G:N3	2.64	0.45
52:N8:40:LYS:HD3	52:N8:46:CYS:SG	2.56	0.45
32:51:4:ILE:HG12	32:51:4:ILE:O	2.16	0.45
44:F8:82:GLN:HE21	44:F8:83:VAL:N	2.14	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:69:69:LYS:O	33:69:73:GLU:HB2	2.16	0.45
31:49:41:GLN:HB3	31:49:43:LEU:HD21	1.98	0.45
20:BI:39:LYS:HB2	20:BI:55:ILE:HG21	1.97	0.45
20:BI:73:HIS:HB3	20:BI:74:LYS:HG3	1.99	0.45
26:14:1754:C:H5	40:75:96:ARG:NH2	2.15	0.45
38:98:55:ALA:HA	38:98:80:PHE:HE1	1.78	0.45
26:14:2748:A:H2'	26:14:2749:A:H8	1.82	0.45
53:O8:11:LEU:O	53:O8:23:THR:HB	2.17	0.45
26:14:16:G:H2'	26:14:17:G:H8	1.82	0.45
26:14:797:C:H2'	26:14:798:G:O4'	2.15	0.45
26:14:783:A:H8	26:14:784:A:H4'	1.80	0.45
44:B5:32:PRO:HA	44:B5:77:LYS:HB2	1.98	0.45
26:14:1971:A:H1'	28:19:240:ALA:O	2.16	0.45
3:22:90:GLU:OE2	3:22:93:LYS:HD2	2.16	0.45
1:1G:1117:G:O3'	9:82:104:ARG:HD2	2.16	0.45
26:1H:2161:C:H2'	26:1H:2162:G:C8	2.48	0.45
1:13:881:G:P	12:3I:12:ARG:HH22	2.38	0.45
1:1G:1272:G:N3	1:1G:1273:G:H1'	2.31	0.45
27:1J:61:G:C6	27:1J:62:C:C4	3.04	0.45
39:65:11:LYS:O	39:65:15:ARG:HG2	2.16	0.45
3:22:173:VAL:HG12	3:22:175:LEU:HG	1.99	0.45
11:2I:21:ILE:HG13	11:2I:30:VAL:HG12	1.97	0.45
26:14:601:C:O2	26:14:605:C:H4'	2.17	0.45
28:19:71:ASP:CG	28:19:103:ARG:HH12	2.19	0.45
26:14:1885:A:H5'	26:14:1886:C:OP2	2.16	0.45
34:15:67:LEU:HD23	34:15:88:GLU:HG2	1.98	0.45
46:D5:19:ARG:HB2	46:D5:19:ARG:HE	1.34	0.45
1:13:1206:G:C6	1:13:1207:G:C5	3.04	0.45
8:7E:104:ARG:HD2	8:7E:138:TRP:CG	2.52	0.45
26:1H:2208:U:O2'	26:1H:2209:C:H5'	2.16	0.45
36:35:97:PRO:HG3	36:35:112:LEU:HD12	1.98	0.45
30:31:198:ALA:O	30:31:201:VAL:N	2.50	0.45
7:6E:120:ILE:O	7:6E:124:LEU:HB2	2.17	0.45
40:B8:39:ARG:HD2	40:B8:40:THR:H	1.82	0.45
8:7E:122:ARG:O	8:7E:126:LYS:HG3	2.16	0.45
5:4E:54:ALA:HA	5:4E:57:LYS:HG3	1.98	0.45
1:13:687:A:C2'	1:13:701:C:H41	2.30	0.45
50:H5:12:PRO:HB2	50:H5:20:LYS:HG2	1.98	0.45
37:45:43:THR:O	37:45:46:GLN:HB2	2.17	0.45
1:13:1274:G:H2'	1:13:1275:A:H8	1.80	0.45
3:2E:24:ALA:HB2	3:2E:32:LEU:HD12	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:19:213:ARG:HD2	28:19:213:ARG:HA	1.73	0.45
26:1H:2820:A:O2'	26:1H:2821:A:OP1	2.34	0.45
26:1H:2427:C:H5''	26:1H:2428:G:OP1	2.16	0.45
55:Q8:47:LYS:HZ3	55:Q8:47:LYS:HB2	1.82	0.45
26:1H:751:A:C6	26:1H:789:A:C5	3.05	0.45
1:13:22:G:H4'	1:13:885:G:C8	2.52	0.45
26:1H:562:U:C4	26:1H:2036:C:O4'	2.69	0.45
36:35:88:LEU:HD22	36:35:114:ILE:HD13	1.97	0.45
1:13:963:G:N2	1:13:972:C:C2	2.84	0.45
1:1G:860:A:H2'	1:1G:861:G:O4'	2.16	0.45
26:14:2414:G:H21	36:35:67:MET:CE	2.30	0.45
29:29:51:PHE:O	29:29:74:PRO:HB2	2.16	0.45
26:1H:2636:U:P	29:21:79:ARG:HA	2.56	0.45
1:1G:1310:G:H5'	13:4A:77:ASN:ND2	2.23	0.45
26:14:1069:A:H4'	26:14:1070:A:H5''	1.98	0.45
29:29:9:VAL:HG12	40:75:8:LYS:NZ	2.29	0.45
26:1H:2347:C:OP2	55:Q8:36:LYS:HE3	2.16	0.45
39:65:54:LEU:HD12	39:65:54:LEU:HA	1.77	0.45
26:14:1321:A:H2'	26:14:1322:A:O4'	2.16	0.45
26:1H:1858:G:C6	26:1H:1883:G:C6	3.04	0.45
1:13:565:U:H3'	1:13:566:G:H2'	1.98	0.45
1:1G:591:U:H2'	1:1G:592:G:C8	2.51	0.45
26:14:1784:A:H4'	26:14:1785:A:C5'	2.46	0.45
26:14:1420:U:O2'	26:14:1421:G:OP1	2.29	0.45
1:13:939:G:C6	1:13:940:C:C4	3.04	0.45
45:C5:87:LYS:CG	45:C5:88:LYS:H	2.30	0.45
9:8E:42:ARG:HB2	9:8E:42:ARG:HE	1.37	0.45
29:29:103:ASP:OD1	29:29:201:THR:HG23	2.16	0.45
1:1G:438:G:H4'	4:32:123:HIS:HD2	1.82	0.45
46:D5:61:LEU:HB3	46:D5:62:PRO:O	2.16	0.45
15:6A:53:HIS:O	15:6A:56:LEU:HB3	2.15	0.45
26:1H:2153:G:H2'	26:1H:2154:G:O4'	2.16	0.45
12:3A:69:TYR:HB2	12:3A:96:VAL:HG11	1.99	0.45
1:13:724:G:N3	1:13:725:G:C8	2.85	0.45
26:1H:527:C:N4	26:1H:2777:G:O2'	2.50	0.45
3:22:36:ASP:OD1	3:22:57:ILE:HG21	2.16	0.45
26:1H:530:G:O4'	26:1H:530:G:N3	2.49	0.45
1:13:1015:A:H2'	1:13:1016:A:C8	2.52	0.45
27:1J:44:G:C2	27:1J:48:A:C2	3.04	0.45
26:1H:2645:G:C3'	26:1H:2646:C:H5'	2.46	0.45
26:14:2529:G:P	26:14:2529:G:H21	2.38	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:309:G:O2'	1:1G:607:A:N1	2.50	0.45
26:14:1011:G:N3	26:14:1151:G:C2	2.85	0.45
1:1G:219:C:H2'	1:1G:220:G:O4'	2.17	0.45
8:7E:39:LEU:HD13	8:7E:111:ILE:HD11	1.99	0.45
6:5E:97:PHE:O	18:9I:31:LEU:HD23	2.16	0.45
26:1H:1766:U:H2'	26:1H:1767:C:C6	2.51	0.45
26:1H:396:G:C8	48:J8:13:ILE:HD11	2.52	0.45
17:8I:29:HIS:O	17:8I:33:GLY:N	2.36	0.45
1:13:1081:G:H2'	1:13:1082:G:C8	2.51	0.45
1:13:223:U:H2'	1:13:224:C:O4'	2.17	0.45
26:1H:719:C:H2'	26:1H:720:C:H6	1.81	0.45
26:14:1386:C:OP2	26:14:1396:U:H5	1.99	0.45
1:13:189:U:O2	17:8I:63:ARG:NH2	2.49	0.45
1:13:591:U:H2'	1:13:592:G:C8	2.52	0.45
27:1J:99:A:C5	27:1J:100:G:C8	3.04	0.45
26:14:1590:U:H2'	26:14:1591:G:C8	2.52	0.45
23:2L:32:G:C4	23:2L:33:OMC:C5	3.05	0.45
47:I8:53:MET:HG3	47:I8:59:LEU:HD23	1.97	0.45
29:21:57:LYS:HE2	29:21:59:VAL:HG12	1.98	0.45
1:1G:1386:G:C2	1:1G:1387:G:C8	3.05	0.45
26:1H:1001:A:H2'	26:1H:1002:G:O4'	2.16	0.45
20:BA:56:MET:HG2	20:BA:84:LEU:HD13	1.99	0.45
26:1H:244:A:H4'	36:78:74:GLU:HB3	1.97	0.45
40:B8:112:ARG:HA	40:B8:115:ARG:NE	2.32	0.45
4:32:81:GLU:O	4:32:85:LYS:HB2	2.16	0.45
1:1G:967:C:H2'	1:1G:968:A:C8	2.51	0.45
17:8A:45:HIS:NE2	17:8A:47:PRO:HB3	2.31	0.45
32:51:170:ARG:HD3	32:51:170:ARG:HA	1.54	0.45
1:13:19:C:OP1	5:4E:125:SER:OG	2.21	0.45
38:98:9:LYS:HA	38:98:17:ARG:NE	2.32	0.45
26:1H:2712:U:O2'	26:1H:2713:A:H5'	2.17	0.45
45:C5:73:ARG:NH2	45:C5:82:PRO:O	2.49	0.45
40:B8:26:ASP:HB3	40:B8:92:GLY:N	2.18	0.45
1:13:1365:G:H2'	1:13:1366:C:H6	1.82	0.45
26:1H:1511:A:H2'	26:1H:1512:G:O4'	2.17	0.45
54:L5:34:ARG:HB3	54:L5:42:LEU:HD22	1.99	0.45
1:13:826:C:H4'	8:7E:12:ARG:HG2	1.99	0.45
31:41:67:LYS:HE2	31:41:67:LYS:N	2.32	0.45
26:14:635:C:H2'	26:14:636:G:O4'	2.16	0.45
55:Q8:33:ASN:O	55:Q8:33:ASN:ND2	2.50	0.45
1:1G:1127:G:C6	1:1G:1145:C:N3	2.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2262:U:O2'	26:14:2263:C:H5'	2.16	0.45
30:31:176:LEU:HB3	30:31:177:ALA:O	2.17	0.45
26:1H:1405:U:H2'	26:1H:1406:U:H6	1.78	0.45
26:14:775:G:C5	26:14:794:G:C8	3.04	0.45
49:K8:16:LEU:HG	49:K8:17:SER:N	2.31	0.45
24:3L:36:A:H2'	24:3L:37:A:N7	2.32	0.45
26:1H:323:G:C2	26:1H:333:G:H1'	2.51	0.45
1:13:994:A:C8	1:13:1216:G:H4'	2.51	0.45
26:14:223:A:N1	26:14:407:G:O2'	2.40	0.45
52:J5:4:HIS:O	52:J5:5:PRO:C	2.55	0.45
26:14:1338:G:N3	26:14:1393:A:H2	2.14	0.45
26:1H:1392:A:C6	26:1H:1393:A:N1	2.85	0.45
1:1G:619:U:O2	4:32:135:LEU:HD22	2.17	0.45
26:1H:287:C:H2'	26:1H:288:C:H6	1.82	0.45
27:16:69:G:C5	27:16:70:C:C5	3.04	0.45
26:1H:302:C:H2'	26:1H:303:U:H6	1.82	0.45
34:15:15:LEU:HD22	34:15:53:VAL:HB	1.99	0.45
34:15:16:ILE:HB	34:15:54:VAL:HG22	1.99	0.45
39:65:88:ASP:HB3	39:65:89:ARG:H	1.44	0.45
1:13:555:C:P	12:3I:20:LYS:NZ	2.90	0.45
26:1H:2373:G:H1	26:1H:2380:C:H42	1.63	0.45
1:1G:485:G:HO2'	1:1G:486:U:H6	1.59	0.45
26:14:1171:G:O2'	26:14:1173:G:OP2	2.35	0.45
1:13:1425:U:H2'	1:13:1426:C:H6	1.82	0.45
37:45:1:MET:H3	37:45:69:PHE:HE1	1.60	0.45
26:1H:26:G:OP1	43:E8:80:PRO:HB3	2.16	0.45
21:1F:2:GLY:O	21:1F:4:GLY:N	2.50	0.45
12:3A:35:GLY:HA3	12:3A:58:VAL:CG1	2.47	0.45
1:13:861:G:H4'	8:7E:18:ARG:HH21	1.82	0.45
41:85:59:ARG:O	41:85:63:VAL:HG23	2.16	0.45
9:8E:125:TYR:HD1	9:8E:126:SER:H	1.63	0.45
12:3I:8:ASN:HA	12:3I:11:VAL:HG23	1.98	0.45
6:5E:35:ALA:HB1	6:5E:65:VAL:HG21	1.98	0.45
28:11:20:ASP:OD2	28:11:22:SER:HB2	2.17	0.45
30:39:165:ARG:HA	30:39:168:ARG:HH11	1.81	0.45
1:1G:415:A:C5	1:1G:416:G:C5	3.05	0.45
26:1H:1537:C:O5'	26:1H:1537:C:H6	1.99	0.45
26:1H:2672:G:H8	26:1H:2672:G:H5''	1.81	0.45
26:14:1627:G:OP2	58:14:3836:HOH:O	2.21	0.45
37:88:63:LYS:HE2	37:88:65:PHE:CZ	2.52	0.45
23:2K:37:U:H2'	23:2K:38:A:O4'	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:E8:33:ARG:NE	43:E8:52:GLU:OE1	2.50	0.45
43:A5:59:VAL:HG12	43:A5:60:ASN:ND2	2.31	0.45
26:14:1614:A:H61	43:A5:88:ARG:H	1.63	0.45
26:1H:1018:C:O3'	26:1H:1120:G:N2	2.50	0.45
1:13:737:A:C4	1:13:738:C:C5	3.05	0.45
4:3E:107:ARG:HH22	4:3E:194:LEU:CD2	2.19	0.45
26:1H:880:G:O2'	26:1H:881:G:O5'	2.27	0.45
29:21:105:THR:HG21	29:21:164:ARG:CZ	2.46	0.45
26:1H:2062:A:N6	26:1H:2503:A:H62	2.15	0.45
52:N8:33:CYS:SG	52:N8:40:LYS:HD3	2.56	0.45
1:1G:1149:C:H2'	1:1G:1150:U:C6	2.51	0.45
44:F8:82:GLN:HA	44:F8:82:GLN:NE2	2.31	0.45
4:3E:9:CYS:HB3	4:3E:32:ALA:CB	2.44	0.45
40:B8:107:ASP:O	40:B8:111:ARG:NH1	2.50	0.45
1:1G:8:A:C5	4:32:209:ARG:HB2	2.51	0.45
26:1H:2287:A:C2	26:1H:2346:A:C2	3.05	0.45
6:52:7:ASN:ND2	18:9A:76:LEU:HD11	2.28	0.45
39:A8:70:GLY:HA2	39:A8:101:LEU:HD13	1.98	0.45
40:B8:50:ILE:HD11	40:B8:102:ILE:HD11	1.97	0.45
1:13:625:G:H4'	16:7I:16:HIS:CD2	2.51	0.45
37:88:52:VAL:HA	37:88:55:VAL:HG13	1.98	0.45
28:19:237:GLU:HB3	28:19:238:GLY:H	1.59	0.45
1:1G:310:G:P	16:7A:27:LYS:HD3	2.56	0.45
1:1G:427:U:H3'	1:1G:428:G:H2'	1.98	0.45
48:F5:73:LEU:HB3	48:F5:90:ILE:HG22	1.99	0.45
26:1H:1538:G:H2'	26:1H:1539:G:H8	1.82	0.45
26:14:315:G:H2'	26:14:316:C:H6	1.81	0.45
23:2L:5:G:H2'	23:2L:6:G:O4'	2.17	0.45
19:AI:37:ARG:HG3	19:AI:37:ARG:H	1.36	0.45
1:13:1412:C:N4	1:13:1488:G:H1	2.13	0.45
4:3E:153:ARG:O	4:3E:181:MET:HE1	2.17	0.45
27:1J:48:A:H4'	39:65:95:HIS:CD2	2.52	0.45
1:1G:129(A):G:C6	1:1G:188:U:H4'	2.51	0.45
26:14:1348:G:N3	26:14:1348:G:H2'	2.32	0.45
53:K5:19:ARG:NH2	53:K5:52:VAL:HG21	2.32	0.45
1:13:922:G:H1'	5:4E:19:MET:HB2	1.98	0.45
26:1H:1259:G:O2'	26:1H:1260:G:H5'	2.17	0.45
26:1H:325:G:H2'	26:1H:326:G:C8	2.52	0.45
26:14:193:U:H5	58:14:4102:HOH:O	1.98	0.45
41:C8:75:ASN:HB3	41:C8:77:SER:N	2.32	0.45
43:A5:72:LYS:HB3	43:A5:106:ILE:HG13	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:177:C:H2'	1:1G:178:C:H6	1.82	0.45
31:41:178:PHE:O	31:41:180:PHE:HD1	2.00	0.45
30:31:197:ASP:O	30:31:199:TRP:N	2.49	0.45
46:H8:40:ASP:OD1	46:H8:43:GLU:HG3	2.17	0.45
14:5I:42:ILE:O	14:5I:46:GLU:HG3	2.16	0.45
24:3L:44:G:H2'	24:3L:45:U:O4'	2.16	0.45
26:14:2400:G:H3'	26:14:2401:U:H6	1.80	0.45
1:13:760:G:H2'	1:13:761:G:H5'	1.99	0.45
17:8A:10:VAL:HG21	17:8A:52:LYS:O	2.17	0.45
26:1H:1207:C:H2'	26:1H:1208:C:H6	1.81	0.45
33:61:85:GLU:HA	33:61:85:GLU:OE2	2.17	0.45
14:5I:50:LYS:HB3	14:5I:50:LYS:HE3	1.66	0.45
20:BI:48:LYS:HB3	20:BI:48:LYS:HE3	1.73	0.45
10:1A:16:LEU:HA	10:1A:16:LEU:HD23	1.84	0.45
33:61:30:LEU:HA	33:61:30:LEU:HD23	1.81	0.45
26:1H:889:C:H3'	26:1H:890:A:H4'	1.99	0.45
26:14:439:G:H2'	26:14:440:G:H8	1.82	0.45
1:13:909:A:H2'	1:13:910:C:O4'	2.16	0.45
26:14:2469:A:H2	26:14:2481:G:N2	2.15	0.45
26:1H:1329:U:H5''	26:1H:1330:C:C5	2.51	0.45
41:85:92:ARG:CZ	42:95:11:GLN:H	2.30	0.45
1:13:1367:C:H5'	10:1I:60:ARG:NH2	2.26	0.45
1:13:1022:G:H2'	1:13:1023:G:C8	2.50	0.45
9:82:128:ARG:NH2	23:2L:34:U:P	2.90	0.45
12:3A:59:ARG:CZ	12:3A:65:GLU:HB2	2.46	0.45
1:13:542:G:H5'	4:3E:41:GLY:HA3	1.98	0.45
27:16:7:G:H5''	27:16:7:G:H8	1.82	0.45
26:1H:1042:G:C6	26:1H:1043:C:C4	3.05	0.45
19:AA:23:ASN:HA	19:AA:27:GLU:CG	2.47	0.45
44:F8:3:THR:HG22	44:F8:6:ASP:HB2	1.99	0.45
1:13:1368:G:H5''	9:8E:112:LYS:CB	2.43	0.45
24:3K:19:G:N2	26:1H:2112:G:H21	2.15	0.45
37:88:134:ARG:NH2	46:H8:122:ARG:HD2	2.31	0.45
1:13:628:G:H2'	1:13:629:G:H8	1.81	0.45
26:1H:69:C:H2'	26:1H:70:G:C8	2.51	0.45
1:1G:973:G:O4'	10:1A:55:LYS:HB3	2.17	0.45
37:88:21:THR:OG1	37:88:22:LYS:O	2.31	0.45
4:32:19:LEU:HB2	4:32:21:LEU:HD21	1.98	0.45
26:14:2847:U:OP1	40:75:98:LYS:HE2	2.16	0.45
35:68:13:ASN:ND2	35:68:97:ARG:HB3	2.31	0.45
23:2L:16:C:O2'	23:2L:62:C:OP1	2.28	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:602:G:OP2	26:14:602:G:H8	1.99	0.45
11:2I:99:GLN:HA	11:2I:105:VAL:HG21	1.99	0.45
34:58:18:ALA:H	34:58:21:LYS:NZ	2.15	0.45
36:78:80:TYR:CE2	36:78:111:ARG:HD3	2.51	0.45
1:13:66:G:C2	1:13:67:C:C6	3.04	0.45
31:49:38:VAL:HG22	31:49:93:THR:HG23	1.98	0.45
14:5I:37:PHE:CE2	14:5I:53:LEU:HD13	2.51	0.45
26:1H:527:C:H4'	26:1H:528:A:C5'	2.46	0.45
26:1H:1364:G:OP1	48:J8:3:LYS:HG2	2.17	0.45
29:29:119:ARG:HD2	29:29:120:TRP:CE2	2.52	0.45
12:3A:100:ILE:HG22	12:3A:101:VAL:H	1.80	0.45
39:65:86:ALA:O	39:65:87:PHE:HB2	2.17	0.45
27:1J:33:G:N3	27:1J:50:G:C2	2.85	0.45
1:1G:567:G:O6	12:3A:5:PRO:HD3	2.17	0.45
26:14:1396:U:H2'	26:14:1396:U:O2	2.17	0.45
26:1H:277:C:H3'	26:1H:278:A:O4'	2.17	0.45
28:11:155:LEU:HD13	28:11:155:LEU:N	2.32	0.45
46:H8:6:LYS:HE3	46:H8:8:TYR:CE1	2.51	0.45
26:1H:1810:A:H2'	26:1H:1811:G:O4'	2.17	0.45
26:14:1551:C:H2'	26:14:1552:G:O4'	2.17	0.45
26:1H:270(C):C:O2'	26:1H:273(B):C:H5'	2.17	0.45
26:1H:37:C:H2'	26:1H:38:A:C8	2.52	0.45
30:31:68:LYS:O	30:31:69:HIS:HB2	2.17	0.45
26:1H:46:C:H2'	26:1H:47:C:C6	2.52	0.45
30:39:81:PRO:HB3	30:39:87:GLY:O	2.17	0.45
4:32:191:ARG:NH1	4:32:200:GLU:OE1	2.45	0.45
26:14:458:G:O2'	54:L5:39:ARG:HD3	2.17	0.45
16:7I:58:TYR:O	16:7I:62:VAL:HG22	2.17	0.45
34:15:29:LYS:HG2	34:15:29:LYS:H	1.42	0.45
31:41:128:ARG:HB2	31:41:128:ARG:HH21	1.81	0.45
1:13:665:A:N3	1:13:732:C:H2'	2.32	0.45
46:H8:128:VAL:CB	46:H8:161:VAL:HG21	2.47	0.45
41:85:58:ARG:HA	41:85:61:TRP:CE3	2.52	0.45
26:1H:1639:U:C2'	26:1H:1640:C:H5'	2.47	0.45
26:1H:442:G:N3	30:31:48:THR:CG2	2.80	0.45
26:1H:1512:G:H2'	26:1H:1513:C:C6	2.52	0.45
24:3K:4:C:H2'	24:3K:5:G:O4'	2.17	0.45
30:31:24:LEU:HD21	30:31:114:VAL:HG12	1.98	0.45
1:1G:1162:C:C2	1:1G:1175:G:C2	3.05	0.45
6:52:86:ARG:O	6:52:87:ARG:HG2	2.17	0.45
26:14:2301:C:C5	26:14:2302:G:N7	2.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:270(E):G:C6	26:1H:270(F):U:C4	3.05	0.45
20:BI:55:ILE:HD13	20:BI:55:ILE:HA	1.81	0.45
26:14:2716:U:H2'	26:14:2717:G:C8	2.53	0.45
45:G8:93:GLY:O	45:G8:94:LYS:HB2	2.17	0.45
1:1G:1256:A:OP2	3:22:26:LYS:NZ	2.42	0.45
1:13:558:G:H2'	1:13:559:A:H2	1.81	0.45
12:3A:110:VAL:HB	12:3A:113:ARG:HB2	1.99	0.45
24:3L:36:A:H2'	24:3L:37:A:C8	2.52	0.45
13:4A:23:TYR:CE1	13:4A:70:LEU:HD12	2.52	0.45
1:1G:532:A:N7	1:1G:1207:G:H4'	2.32	0.45
31:49:135:LEU:HD11	31:49:157:ILE:HD11	1.99	0.45
7:62:50:ILE:HB	7:62:58:PRO:HG3	1.99	0.45
11:2I:121:PRO:HG2	11:2I:126:ARG:CG	2.47	0.45
26:14:959:A:C6	26:14:960:A:N1	2.85	0.45
26:14:1495:A:O2'	26:14:1496:A:H5'	2.17	0.45
51:I5:37:SER:C	51:I5:39:CYS:H	2.21	0.45
51:I5:37:SER:HB3	51:I5:39:CYS:HB2	1.99	0.45
1:13:438:G:OP1	4:3E:125:HIS:HE1	2.00	0.45
26:14:807:U:H4'	26:14:2446:G:OP1	2.16	0.45
1:1G:57:G:H2'	1:1G:58:C:C6	2.52	0.45
33:61:129:THR:HG22	33:61:137:PRO:HB3	1.99	0.45
12:3I:33:ARG:O	12:3I:85:ILE:HG12	2.17	0.45
26:14:1439:A:C8	26:14:1440:G:C8	3.05	0.45
1:1G:1057:G:H2'	1:1G:1058:G:C8	2.52	0.45
26:14:942:G:H4'	26:14:1190:G:H5'	1.98	0.45
1:13:833:U:H3	1:13:853:G:H1	1.65	0.45
29:29:117:MET:HA	29:29:122:PHE:N	2.32	0.45
41:85:62:ILE:HD11	41:85:93:LYS:HD3	1.98	0.45
26:14:484:C:H2'	26:14:485:C:C6	2.52	0.45
37:45:43:THR:HB	37:45:45:GLN:HG2	1.98	0.45
23:2L:3:C:H5'	26:14:2255:G:O2'	2.17	0.45
29:29:50:GLY:HA2	29:29:78:LEU:HB3	1.99	0.45
26:1H:536:A:H2'	26:1H:537:C:C6	2.52	0.45
39:65:6:ALA:O	39:65:10:ARG:HG3	2.17	0.45
1:1G:569:C:H1'	1:1G:574:A:C4	2.51	0.45
26:1H:2016:U:H1'	52:N8:6:VAL:HG13	1.98	0.45
40:B8:5:ALA:O	40:B8:8:LYS:HG2	2.17	0.45
26:1H:1378:A:O2'	26:1H:1380:G:N7	2.43	0.45
11:2A:56:GLY:O	11:2A:89:ALA:HB3	2.17	0.45
40:B8:120:ARG:HA	40:B8:123:GLN:HG2	1.99	0.45
5:42:129:ILE:O	5:42:132:ALA:HB3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:J8:50:ARG:HD2	48:J8:57:GLU:OE2	2.17	0.45
7:62:131:LYS:HB3	7:62:131:LYS:HE3	1.58	0.45
14:5A:47:LEU:HA	14:5A:47:LEU:HD23	1.68	0.45
27:16:57:A:H8	27:16:57:A:O5'	2.00	0.45
1:1G:1250:A:H4'	9:82:68:GLY:N	2.32	0.45
36:78:95:VAL:O	36:78:126:VAL:HG23	2.17	0.45
4:32:25:ARG:NH1	4:32:31:CYS:HA	2.32	0.44
41:C8:11:ARG:O	41:C8:15:LYS:HG3	2.17	0.44
26:14:1112:G:H2'	26:14:1113:U:C6	2.52	0.44
49:K8:55:ARG:O	49:K8:58:ALA:HB3	2.16	0.44
26:14:2324:C:H42	26:14:2331:G:H1	1.65	0.44
26:1H:1590:U:H2'	26:1H:1591:G:H8	1.79	0.44
26:1H:66:C:C2	26:1H:67:U:C6	3.06	0.44
53:K5:28:ARG:NH1	53:K5:30:THR:O	2.50	0.44
48:J8:76:ARG:HH11	48:J8:94:LEU:HD22	1.81	0.44
38:98:53:HIS:HB2	38:98:94:TYR:HE2	1.82	0.44
29:21:144:ARG:HG3	29:21:144:ARG:HH11	1.82	0.44
26:1H:2388:A:H2'	26:1H:2389:G:H5'	1.99	0.44
24:3K:8:U:O2'	24:3K:9:A:OP1	2.28	0.44
46:D5:60:GLU:HB2	46:D5:66:SER:OG	2.18	0.44
26:14:668:G:H2'	26:14:670:A:H62	1.82	0.44
32:59:82:GLY:O	32:59:134:SER:HA	2.17	0.44
2:1E:11:LEU:HD13	2:1E:217:ARG:HH22	1.81	0.44
28:19:65:ILE:HD11	28:19:67:PHE:CE1	2.52	0.44
12:3A:93:LEU:HB3	12:3A:96:VAL:HG21	1.99	0.44
32:59:83:TYR:O	32:59:84:SER:OG	2.34	0.44
26:14:117:G:C6	26:14:119:A:C6	3.05	0.44
37:45:29:PHE:HD2	37:45:65:PHE:CZ	2.35	0.44
2:12:185:ILE:CG2	2:12:199:TYR:HB2	2.46	0.44
1:13:984:C:H2'	1:13:985:C:H6	1.82	0.44
43:A5:71:VAL:HA	43:A5:107:LEU:HD12	1.99	0.44
1:1G:286:G:C6	1:1G:287:U:C4	3.05	0.44
26:14:1385:G:C4	26:14:1386:C:C5	3.05	0.44
42:D8:25:LEU:H	42:D8:92:THR:CG2	2.30	0.44
26:14:279:C:H42	26:14:361:G:H1	1.64	0.44
43:A5:36:LEU:HD11	43:A5:47:VAL:HG12	1.98	0.44
31:41:16:ARG:N	31:41:17:PRO:HD2	2.31	0.44
2:1E:87:ARG:HH21	2:1E:233:SER:HB2	1.82	0.44
26:1H:609:A:H8	26:1H:609:A:O5'	2.00	0.44
26:1H:2723:C:OP1	29:21:109:LYS:HD3	2.17	0.44
17:8A:95:TYR:HA	17:8A:98:LEU:HG	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:32:43:HIS:O	4:32:46:LYS:HB2	2.17	0.44
1:1G:1343:G:H4'	9:82:122:ALA:HB3	1.98	0.44
26:14:396:G:O4'	48:F5:13:ILE:HD11	2.17	0.44
26:14:1165:U:H2'	26:14:1166:C:C6	2.52	0.44
31:49:130:ASN:HB3	31:49:159:VAL:O	2.16	0.44
26:1H:1945:G:C4	26:1H:1946:U:C5	3.06	0.44
11:2A:38:ASN:HA	11:2A:39:PRO:HD3	1.84	0.44
38:98:24:GLN:HE22	38:98:36:THR:CG2	2.29	0.44
18:9I:37:VAL:O	18:9I:41:LYS:N	2.36	0.44
26:1H:2444:G:O6	58:1H:4648:HOH:O	2.20	0.44
26:1H:2661:G:H8	26:1H:2661:G:OP2	2.00	0.44
33:61:10:GLU:O	33:61:10:GLU:HG3	2.16	0.44
10:1I:90:LEU:N	10:1I:91:PRO:HD3	2.32	0.44
55:Q8:14:VAL:HB	55:Q8:21:LYS:HE3	1.99	0.44
2:1E:185:ILE:CG2	2:1E:199:TYR:HB2	2.35	0.44
1:1G:1322:C:O2'	1:1G:1323:G:H5'	2.18	0.44
1:1G:1355:G:H2'	1:1G:1356:G:C8	2.52	0.44
27:1J:15:A:H1'	27:1J:109:G:N9	2.32	0.44
36:35:39:LYS:CD	36:35:45:LEU:HD21	2.40	0.44
26:1H:1556:C:H2'	26:1H:1557:C:C6	2.51	0.44
26:14:2331:G:H5'	47:E5:44:ARG:HG2	1.99	0.44
46:D5:39:VAL:HG21	46:D5:44:PHE:HB2	1.99	0.44
26:14:2395:C:H2'	26:14:2396:G:O4'	2.17	0.44
26:14:1677:A:H2'	26:14:1678:G:H8	1.81	0.44
26:1H:1899:G:N2	26:1H:1902:C:H41	2.15	0.44
26:1H:1439:A:H3'	26:1H:1440:G:H8	1.82	0.44
26:14:2472:G:N2	26:14:2478:A:H62	2.14	0.44
26:1H:2262:U:H4'	26:1H:2328:A:C2	2.52	0.44
24:3K:35:A:C6	24:3K:36:A:N6	2.86	0.44
1:1G:1054:C:N4	22:1L:34:G:C8	2.86	0.44
1:1G:1272:G:H2'	1:1G:1273:G:O4'	2.17	0.44
26:14:491:G:H2'	26:14:492:A:H8	1.78	0.44
1:1G:1007:C:H2'	1:1G:1008:C:O4'	2.16	0.44
27:16:30:C:H2'	27:16:31:C:C5'	2.46	0.44
46:D5:158:PRO:HB2	46:D5:159:PRO:HD2	1.99	0.44
32:59:66:GLY:O	32:59:69:ARG:HB3	2.17	0.44
26:1H:1469:A:H2'	26:1H:1470:G:C8	2.52	0.44
27:1J:31:C:N4	39:65:32:LEU:HD13	2.33	0.44
5:42:34:VAL:O	5:42:41:VAL:HG12	2.16	0.44
55:M5:8:LYS:HD3	55:M5:8:LYS:N	2.31	0.44
26:1H:1259:G:H2'	26:1H:1260:G:C8	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:32:101:LEU:HA	4:32:101:LEU:HD12	1.69	0.44
26:14:360:G:H2'	26:14:361:G:O4'	2.16	0.44
26:14:2629:A:H4'	26:14:2630:G:H5'	1.99	0.44
26:14:2412:A:C2	26:14:2413:G:H1'	2.53	0.44
26:1H:1625:C:H2'	26:1H:1626:G:O4'	2.18	0.44
26:14:111:A:H4'	49:G5:69:ARG:NH2	2.32	0.44
46:D5:10:ARG:HH21	46:D5:26:GLY:H	1.66	0.44
26:1H:1380:G:N2	26:1H:1570:A:C2	2.85	0.44
1:1G:34:C:H2'	1:1G:35:G:C8	2.52	0.44
17:8I:60:ILE:O	17:8I:62:SER:OG	2.33	0.44
2:12:214:ILE:HG22	2:12:215:LEU:HD22	2.00	0.44
26:14:753:C:H2'	26:14:754:C:H6	1.81	0.44
30:39:124:LEU:HB3	30:39:126:VAL:CG1	2.47	0.44
11:2I:41:THR:HG21	11:2I:71:LYS:HB3	1.99	0.44
42:95:55:ALA:HA	42:95:101:GLY:HA2	1.98	0.44
1:13:161:A:H2'	1:13:162:A:C8	2.52	0.44
48:J8:5:CYS:CB	48:J8:8:SER:HG	2.30	0.44
1:1G:1509:C:H2'	1:1G:1510:U:O4'	2.17	0.44
49:K8:61:LEU:HD23	49:K8:61:LEU:HA	1.59	0.44
1:1G:720:C:O5'	1:1G:720:C:H6	2.00	0.44
6:5E:76:ALA:O	6:5E:80:ARG:HG3	2.18	0.44
1:1G:417:C:H2'	1:1G:418:C:C6	2.53	0.44
26:1H:2097:C:C2'	26:1H:2098:U:H5'	2.47	0.44
29:29:134:ILE:O	29:29:134:ILE:HD12	2.16	0.44
30:31:183:VAL:O	30:31:187:VAL:HG23	2.17	0.44
38:55:35:THR:HG21	38:55:100:LEU:HD11	1.99	0.44
26:1H:973:A:O4'	26:1H:1188:U:C6	2.71	0.44
1:13:1347:G:H22	1:13:1374:A:P	2.39	0.44
26:1H:2253:G:OP1	58:1H:3724:HOH:O	2.21	0.44
26:14:322:A:H3'	30:39:169:ASN:OD1	2.17	0.44
46:H8:60:GLU:O	46:H8:61:LEU:HD23	2.16	0.44
26:1H:1454:U:O2'	26:1H:1455:G:N7	2.50	0.44
26:1H:444:C:C4'	30:31:49:ALA:HB2	2.47	0.44
30:31:57:VAL:HG13	30:31:59:TYR:CD2	2.52	0.44
1:1G:1237:C:O2'	1:1G:1300:G:N2	2.49	0.44
1:1G:1298:C:H5''	7:62:114:ARG:HH12	1.82	0.44
1:1G:651:C:H2'	1:1G:652:U:C6	2.53	0.44
8:7E:87:SER:HB2	8:7E:93:VAL:CB	2.47	0.44
5:4E:126:ARG:CG	5:4E:126:ARG:HH11	2.30	0.44
26:1H:2286:A:H8	53:O8:37:ARG:HH11	1.66	0.44
18:9A:69:THR:HA	18:9A:72:ARG:HB2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:3L:36:A:H5'	24:3L:37:A:OP2	2.18	0.44
26:14:900:A:H2'	26:14:900:A:N3	2.32	0.44
40:75:80:SER:HB3	40:75:83:ILE:HG13	1.99	0.44
19:AI:5:LEU:CB	19:AI:10:PHE:HE1	2.31	0.44
19:AI:4:SER:C	19:AI:5:LEU:HG	2.35	0.44
1:1G:79:G:H2'	1:1G:79:G:N3	2.32	0.44
29:21:149:ARG:CG	29:21:149:ARG:HH11	2.30	0.44
1:1G:1106:G:C2	1:1G:1107:C:C6	3.06	0.44
1:1G:15:G:N7	1:1G:1396:A:C2	2.84	0.44
39:A8:49:VAL:HG12	39:A8:73:LEU:HD23	2.00	0.44
1:13:948:C:C2'	1:13:949:A:H5'	2.47	0.44
27:16:88:C:H2'	27:16:89:G:O4'	2.17	0.44
9:8E:89:ASN:C	9:8E:91:ASP:H	2.21	0.44
11:2I:59:TYR:CE2	11:2I:63:LEU:HD11	2.53	0.44
31:49:95:ARG:HG2	31:49:96:ARG:H	1.82	0.44
1:13:958:A:C6	1:13:959:A:N1	2.85	0.44
28:11:26:LYS:CE	28:11:94:LEU:HD12	2.47	0.44
21:1B:6:ARG:O	21:1B:12:LYS:HE2	2.17	0.44
30:39:68:LYS:HB3	30:39:69:HIS:CD2	2.53	0.44
32:59:75:ALA:O	32:59:79:VAL:HG13	2.18	0.44
26:1H:357:A:H2'	26:1H:358:U:H6	1.82	0.44
26:14:2076:U:H5''	26:14:2238:G:H22	1.82	0.44
13:4A:13:LYS:HA	13:4A:44:ARG:HH11	1.83	0.44
1:1G:1387:G:H2'	1:1G:1388:C:C6	2.53	0.44
26:1H:537:C:H2'	26:1H:539:G:C8	2.52	0.44
33:61:93:THR:O	33:61:97:ILE:HG13	2.18	0.44
13:4A:50:GLU:O	13:4A:54:VAL:HG23	2.16	0.44
26:1H:785:G:C5	26:1H:786:C:C5	3.04	0.44
1:1G:1459:C:H2'	1:1G:1460:A:O4'	2.17	0.44
19:AA:53:ASN:HA	19:AA:77:THR:HG22	1.99	0.44
2:1E:136:VAL:O	2:1E:140:HIS:N	2.50	0.44
26:1H:1756:G:H1'	26:1H:1758:G:C2	2.52	0.44
26:1H:64:A:O3'	44:F8:71:GLY:HA3	2.17	0.44
54:P8:46:VAL:HG22	54:P8:48:LYS:HE3	2.00	0.44
35:68:116:SER:OG	35:68:117:LEU:N	2.50	0.44
26:14:2729:G:H2'	26:14:2730:C:H6	1.81	0.44
26:1H:2303:G:O2'	31:41:132:ASN:HB2	2.17	0.44
26:14:2092:U:H4'	26:14:2093:G:O5'	2.17	0.44
47:I8:19:LYS:HD3	47:I8:19:LYS:HA	1.59	0.44
3:2E:115:LEU:HD23	3:2E:115:LEU:HA	1.66	0.44
41:C8:104:GLN:CD	41:C8:104:GLN:H	2.20	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2275:C:O2'	37:88:83:MET:HA	2.18	0.44
26:1H:2689:U:P	26:1H:2719:G:H22	2.38	0.44
26:1H:587:C:H4'	26:1H:588:U:O5'	2.17	0.44
55:Q8:47:LYS:HD3	55:Q8:47:LYS:HA	1.51	0.44
26:1H:449:A:H3'	58:1H:3893:HOH:O	2.17	0.44
26:1H:374:A:C2	26:1H:401:A:C4	3.06	0.44
1:1G:1321:C:C2	1:1G:1322:C:C5	3.03	0.44
26:1H:1657:C:H2'	26:1H:1658:C:H6	1.80	0.44
1:1G:1068:G:N7	1:1G:1094:G:C8	2.86	0.44
1:1G:1028(A):C:N4	1:1G:1032(B):G:H22	2.15	0.44
26:1H:1675:C:N3	29:21:128:SER:OG	2.49	0.44
35:68:4:PRO:O	35:68:5:GLN:CB	2.66	0.44
1:13:1285:A:H4'	1:13:1286:A:H5'	1.98	0.44
4:3E:59:ARG:HH22	4:3E:66:ARG:HH22	1.66	0.44
1:1G:373:A:N3	1:1G:374:A:C8	2.85	0.44
16:7A:16:HIS:N	16:7A:16:HIS:CD2	2.84	0.44
39:A8:62:LYS:HB3	39:A8:97:ARG:HD2	1.99	0.44
1:13:644:G:H2'	1:13:645:C:O4'	2.18	0.44
16:7I:38:TYR:CE1	16:7I:50:LYS:HB2	2.52	0.44
26:14:2228:G:C5	26:14:2229:C:C4	3.05	0.44
1:13:286:G:C6	1:13:287:U:C4	3.04	0.44
13:4I:7:VAL:HB	31:4I:115:ARG:NH1	2.32	0.44
1:1G:17:U:H2'	1:1G:18:C:C6	2.51	0.44
36:78:15:ARG:CB	36:78:16:ARG:HB2	2.47	0.44
1:13:475:G:H2'	1:13:476:G:H8	1.83	0.44
26:1H:1534:G:H2'	26:1H:1534:G:N3	2.33	0.44
2:1E:166:ASP:C	2:1E:168:THR:H	2.20	0.44
30:31:101:LEU:O	30:31:106:ARG:NH1	2.51	0.44
40:75:45:PHE:CD2	40:75:74:ARG:HD3	2.52	0.44
1:1G:438:G:H5'	4:32:123:HIS:CD2	2.52	0.44
55:M5:6:THR:HG22	55:M5:59:LYS:O	2.17	0.44
22:1L:11:C:C4	22:1L:12:U:C4	3.05	0.44
36:35:128:HIS:HA	36:35:147:LEU:HA	1.99	0.44
7:6E:5:ARG:CZ	7:6E:7:ALA:HA	2.48	0.44
36:35:101:VAL:HG23	36:35:107:LYS:H	1.83	0.44
5:42:92:LYS:HB3	5:42:119:LEU:HB2	1.99	0.44
26:1H:1111:A:H5'	32:51:3:ARG:CD	2.48	0.44
26:1H:1053:C:H42	26:1H:1106:G:H1	1.64	0.44
24:3K:37:A:HO2'	24:3K:38:A:C5'	2.29	0.44
6:52:8:ILE:HG13	6:52:88:VAL:HG22	1.99	0.44
17:8I:31:LEU:HD23	17:8I:32:TYR:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:25:11:ALA:HB1	35:25:99:PHE:O	2.17	0.44
2:12:185:ILE:HG22	2:12:199:TYR:CD1	2.53	0.44
39:65:67:ARG:O	39:65:71:ARG:HB2	2.18	0.44
36:78:39:LYS:HG3	36:78:45:LEU:CD2	2.46	0.44
47:I8:23:VAL:HA	47:I8:38:VAL:HG22	2.00	0.44
32:59:58:GLU:HB2	32:59:61:HIS:CE1	2.52	0.44
37:45:41:TRP:CZ3	37:45:74:TYR:HE1	2.35	0.44
7:6E:46:ALA:O	7:6E:49:ILE:N	2.50	0.44
26:14:1993:U:H4'	29:29:128:SER:CB	2.47	0.44
26:1H:886:C:C6	26:1H:887:A:H1'	2.52	0.44
47:E5:32:ARG:O	47:E5:34:GLY:N	2.43	0.44
16:7A:40:ASP:HB3	16:7A:48:TRP:HB2	1.98	0.44
26:14:1910:G:O2'	26:14:1911:U:H5'	2.18	0.44
26:1H:2650:U:H2'	26:1H:2651:C:C6	2.53	0.44
31:41:26:GLN:HG3	31:41:30:GLU:OE1	2.18	0.44
26:14:1844:C:O3'	28:19:258:LYS:NZ	2.38	0.44
26:1H:2123:G:H1	26:1H:2175:C:H42	1.65	0.44
26:14:120:U:OP2	58:14:3995:HOH:O	2.21	0.44
39:A8:72:ALA:O	39:A8:76:LYS:HG3	2.17	0.44
34:58:121:LYS:HB3	34:58:123:TYR:HE1	1.81	0.44
34:58:74:ARG:CZ	34:58:85:ILE:HD11	2.48	0.44
45:C5:92:ASN:HB3	45:C5:93:GLY:H	1.65	0.44
4:3E:86:LYS:HD2	4:3E:86:LYS:H	1.83	0.44
26:14:1894:C:C2'	26:14:1895:C:H5'	2.47	0.44
11:2A:70:LYS:HB2	11:2A:70:LYS:HE3	1.65	0.44
41:85:39:LEU:HD23	41:85:39:LEU:HA	1.67	0.44
41:C8:39:LEU:HA	41:C8:39:LEU:HD23	1.76	0.44
26:1H:242:G:H5'	55:Q8:60:LEU:CD1	2.48	0.44
26:14:270(R):G:H2'	26:14:270(S):G:C8	2.51	0.44
6:52:2:ARG:HD3	6:52:92:LYS:HE3	2.00	0.44
1:13:636:U:H2'	1:13:637:G:H8	1.83	0.44
39:A8:9:ARG:O	39:A8:12:PHE:N	2.51	0.44
1:13:1124:G:N3	1:13:1127:G:N2	2.65	0.44
26:14:1225:C:H5''	42:95:85:LYS:NZ	2.33	0.44
39:A8:36:TYR:N	39:A8:36:TYR:CD1	2.85	0.44
26:1H:860:U:H2'	26:1H:861:A:H8	1.83	0.44
1:1G:1280:A:P	10:1A:40:LEU:HD21	2.58	0.44
29:21:128:SER:OG	29:21:129:HIS:N	2.50	0.44
48:J8:72:GLU:O	48:J8:76:ARG:HG2	2.18	0.44
26:1H:2287:A:H2	26:1H:2346:A:C2	2.35	0.44
26:1H:1204:A:N6	26:1H:1240:U:H2'	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:115:G:C2	1:13:289:G:N7	2.86	0.44
26:1H:879:G:H8	26:1H:879:G:O5'	2.00	0.44
1:1G:1330:U:H4'	13:4A:23:TYR:CE2	2.53	0.44
1:1G:1118:C:H1'	1:1G:1179:A:C4	2.53	0.44
1:13:649:G:N3	1:13:650:G:C8	2.85	0.44
19:AI:5:LEU:HB3	19:AI:10:PHE:CE1	2.49	0.44
2:12:8:LYS:HE2	2:12:213:LEU:HD21	1.98	0.44
50:L8:50:VAL:HG23	50:L8:54:VAL:HG12	1.99	0.44
50:L8:54:VAL:HG13	50:L8:54:VAL:O	2.17	0.44
28:11:75:ILE:HD13	28:11:99:ASP:OD2	2.18	0.44
10:1I:47:PHE:CE2	14:5I:37:PHE:HE1	2.36	0.44
1:13:720:C:C4	1:13:721:G:C5	3.05	0.44
26:1H:1542:G:OP2	26:1H:1543:A:O2'	2.35	0.44
26:14:747:U:OP1	52:J5:3:LYS:HD3	2.16	0.44
4:3E:61:LYS:HD2	4:3E:207:TYR:OH	2.17	0.44
26:14:756:C:C2'	26:14:757:U:H5'	2.48	0.44
26:14:1011:G:C6	26:14:1013:C:C4	3.06	0.44
30:31:7:TYR:O	30:31:22:ALA:N	2.48	0.44
53:K5:51:GLU:HG2	53:K5:52:VAL:N	2.32	0.44
26:1H:2373:G:H1	26:1H:2380:C:N4	2.16	0.44
26:14:2425:A:H5'	26:14:2427:C:O4'	2.18	0.44
43:A5:27:LYS:O	43:A5:71:VAL:HG23	2.18	0.44
42:D8:44:LYS:NZ	42:D8:44:LYS:HA	2.33	0.44
1:1G:448:A:OP2	1:1G:485:G:N1	2.45	0.44
43:A5:15:ARG:O	43:A5:19:LEU:HD13	2.16	0.44
1:13:1442:G:C6	1:13:1446:A:C6	3.05	0.44
26:14:1005:C:O2	26:14:1143:A:C6	2.71	0.44
26:1H:2111:C:HO2'	26:1H:2119:A:P	2.40	0.44
35:68:10:VAL:HG23	35:68:17:ARG:O	2.18	0.44
47:E5:12:ASN:HA	47:E5:14:ARG:HH21	1.82	0.44
6:52:15:ASP:OD1	6:52:17:SER:N	2.49	0.44
8:7E:77:GLU:HG2	8:7E:78:GLN:H	1.82	0.44
33:61:5:LEU:HD12	33:61:13:GLY:O	2.17	0.44
1:1G:476:G:O2'	1:1G:477:G:H5'	2.17	0.44
26:1H:739:G:OP2	26:1H:739:G:H8	1.99	0.44
26:1H:754:C:H2'	26:1H:755:C:C6	2.53	0.44
26:1H:755:C:H2'	26:1H:756:C:C6	2.53	0.44
1:13:414:A:H2'	1:13:415:A:C8	2.52	0.44
26:1H:478:A:C6	26:1H:480:A:C6	3.05	0.44
26:1H:90:U:H4'	26:1H:91:A:H5'	1.99	0.44
30:39:158:THR:O	30:39:177:ALA:HA	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:513:A:C2	26:14:514:A:C5	3.05	0.44
11:2A:54:ARG:HG3	11:2A:54:ARG:H	1.58	0.44
1:1G:1082:G:OP2	1:1G:1082:G:H8	2.01	0.44
2:12:204:ASN:N	2:12:204:ASN:OD1	2.51	0.44
32:51:125:VAL:HG12	32:51:127:GLU:O	2.18	0.44
26:1H:1799:G:H5'	26:1H:1819:A:H61	1.83	0.44
26:1H:2743:C:H2'	26:1H:2744:G:O4'	2.18	0.44
1:1G:807:A:H2'	1:1G:808:C:C6	2.53	0.44
26:14:1022:G:C6	26:14:1140:C:C4	3.06	0.44
26:1H:1479:G:C4	26:1H:1480:G:C8	3.06	0.44
26:14:2392:A:H1'	36:35:61:ARG:NH2	2.27	0.44
1:1G:1237:C:H5''	1:1G:1238:A:O4'	2.17	0.44
44:F8:47:PHE:O	44:F8:49:VAL:HG23	2.17	0.44
40:B8:51:ARG:HB2	40:B8:98:LYS:HD3	1.99	0.44
46:H8:15:PRO:O	46:H8:19:ARG:HB2	2.18	0.44
1:13:412:A:C8	4:3E:35:ARG:NH1	2.85	0.44
55:Q8:32:LEU:H	55:Q8:32:LEU:HD23	1.83	0.44
1:1G:1147:C:O2	9:82:16:ARG:NE	2.51	0.44
1:1G:1157:A:O5'	1:1G:1158:C:H5	2.01	0.44
1:1G:1157:A:N6	1:1G:1178:G:H21	2.11	0.44
1:13:626:U:H2'	1:13:627:G:H8	1.82	0.44
1:1G:629:G:H2'	1:1G:630:G:O4'	2.18	0.44
1:1G:1153:C:H2'	1:1G:1154:G:H8	1.83	0.44
27:1J:18:G:H2'	27:1J:19:G:H5''	1.99	0.44
1:13:109:A:N7	1:13:326:G:H2'	2.33	0.44
1:13:992:U:H2'	1:13:1043:C:H41	1.83	0.44
26:1H:2000:G:C6	26:1H:2001:A:N7	2.86	0.44
4:32:61:LYS:HE2	4:32:206:PHE:CE2	2.52	0.44
26:14:301:G:C4	26:14:302:C:C5	3.06	0.44
2:1E:28:PHE:O	2:1E:32:ILE:HG22	2.18	0.44
29:29:57:LYS:HD3	29:29:57:LYS:HA	1.66	0.44
26:1H:1500:G:C5	26:1H:1501:C:C4	3.06	0.44
9:82:33:PHE:HE2	9:82:47:LEU:HD23	1.82	0.44
26:14:1709:U:O4'	26:14:2860:A:H1'	2.18	0.44
26:14:49:A:C5	26:14:177:G:C6	3.06	0.44
26:14:829:A:N7	26:14:2248:C:H5'	2.33	0.44
27:16:15:A:H1'	27:16:109:G:C4	2.52	0.44
26:14:1496:A:C8	26:14:1577:C:O2'	2.71	0.44
1:1G:1255:G:C2	1:1G:1283:G:C2	3.05	0.44
26:1H:686:G:H22	54:P8:15:THR:HG21	1.83	0.44
28:19:2:ALA:HB3	28:19:20:ASP:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:B8:93:ARG:NH1	40:B8:93:ARG:HG3	2.33	0.44
54:P8:24:THR:HG23	54:P8:27:GLY:N	2.33	0.44
27:16:25:A:H2'	27:16:26:A:O4'	2.18	0.44
3:22:113:ALA:HA	3:22:202:ILE:HD11	2.00	0.44
26:14:923:C:H2'	26:14:924:C:C6	2.53	0.44
33:69:128:LEU:O	33:69:137:PRO:HA	2.18	0.44
26:1H:500:G:N1	26:1H:503:A:OP2	2.48	0.44
1:1G:744:C:O2'	1:1G:851:G:N2	2.50	0.44
26:1H:1753:G:N1	26:1H:1756:G:C2	2.86	0.44
28:19:85:ASP:OD2	28:19:88:ARG:HD2	2.18	0.44
43:E8:19:LEU:HD23	52:N8:25:LEU:HD21	2.00	0.44
8:7E:1:MET:HB3	8:7E:2:LEU:H	1.61	0.44
26:14:562:U:O4	26:14:2036:C:H1'	2.17	0.44
26:14:2599:G:OP2	28:19:236:GLY:N	2.51	0.44
26:14:81:G:O2'	26:14:295:G:O2'	2.35	0.44
26:1H:2850:A:C2	26:1H:2851:A:C4	3.06	0.44
41:C8:21:ALA:HA	41:C8:24:TYR:CE2	2.53	0.44
54:P8:10:ARG:O	54:P8:10:ARG:HG2	2.18	0.44
52:J5:52:TYR:CD1	52:J5:53:ALA:N	2.86	0.44
43:E8:45:TYR:CE1	43:E8:49:LYS:HD2	2.53	0.44
1:13:322:C:H5	1:13:328:C:C5	2.36	0.44
31:49:144:ILE:HG22	31:49:148:MET:HB3	2.00	0.44
26:14:2473:U:H2'	26:14:2473:U:O2	2.17	0.44
35:25:93:PRO:HD3	35:25:113:LYS:HG3	1.99	0.44
45:G8:17:SER:HA	45:G8:21:LYS:HB2	1.99	0.44
26:14:815:C:H2'	26:14:816:C:C6	2.52	0.44
26:14:723:G:H2'	26:14:724:U:O4'	2.18	0.44
36:78:49:ARG:HG3	55:Q8:57:ARG:HD2	2.00	0.44
55:Q8:6:THR:HG22	55:Q8:59:LYS:CG	2.47	0.44
1:13:1504:G:O3'	58:13:1804:HOH:O	2.21	0.44
26:14:1047:G:N2	26:14:1111:A:H62	2.16	0.44
26:14:886:C:H1'	26:14:890:A:C2	2.53	0.44
13:4A:78:ILE:HD13	13:4A:92:HIS:HE1	1.82	0.44
26:1H:1576:U:N3	26:1H:1577:C:C5	2.85	0.44
1:1G:1300:G:C5	1:1G:1334:G:C6	3.06	0.44
12:3A:60:LEU:HD23	12:3A:64:TYR:HB3	1.99	0.44
26:14:1674:G:H1'	26:14:1676:A:N6	2.32	0.44
1:13:465:A:N7	1:13:467:G:N7	2.65	0.44
26:14:2286:A:C8	26:14:2287:A:N6	2.86	0.44
38:98:92:GLY:H	38:98:94:TYR:HE1	1.66	0.44
26:1H:1086:A:H1'	26:1H:1103:A:N6	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
43:E8:82:LEU:HD13	43:E8:84:ARG:NH2	2.32	0.44
1:1G:683:G:C6	1:1G:684:A:C6	3.06	0.44
12:3A:40:VAL:HG11	12:3A:77:LEU:HB3	2.00	0.44
26:14:1179:C:H2'	26:14:1180:C:H6	1.82	0.44
37:45:25:ASP:HB3	37:45:102:VAL:H	1.81	0.44
26:14:2012:G:OP2	43:A5:16:LYS:NZ	2.50	0.44
26:1H:2818:G:H4'	26:1H:2837:G:O4'	2.18	0.44
4:32:108:LEU:CD2	4:32:183:GLY:HA3	2.47	0.44
38:55:54:LEU:HD12	38:55:54:LEU:HA	1.85	0.44
37:88:109:VAL:HG12	37:88:114:ALA:HB2	2.00	0.44
26:1H:2012:G:OP2	43:E8:16:LYS:NZ	2.51	0.44
1:13:1489:G:H2'	1:13:1490:C:O4'	2.18	0.44
1:1G:109:A:H5'	1:1G:110:C:C5	2.53	0.44
6:52:95:GLU:HA	6:52:96:PRO:HD3	1.83	0.44
5:42:118:ILE:HG12	5:42:119:LEU:N	2.32	0.44
5:42:118:ILE:HG12	5:42:119:LEU:H	1.83	0.44
16:7I:28:ARG:HD2	16:7I:29:ASP:OD1	2.17	0.44
26:14:2611:U:H1'	52:J5:3:LYS:HE2	2.00	0.44
28:19:70:TRP:CD1	28:19:71:ASP:N	2.86	0.44
1:13:942:G:C2	1:13:1342:C:C2	3.06	0.44
26:14:749:C:H4'	26:14:1271:G:N3	2.33	0.44
28:19:37:LEU:HD23	28:19:62:TYR:HB2	1.99	0.44
26:14:2864:G:C6	26:14:2865:U:N3	2.86	0.44
1:13:1084:G:H2'	1:13:1085:U:C5	2.52	0.44
26:14:1131:G:O6	26:14:2040:C:H1'	2.18	0.44
1:1G:476:G:H2'	1:1G:477:G:H8	1.83	0.44
27:1J:43:C:OP1	51:I5:6:HIS:HE1	2.01	0.44
28:19:244:ARG:HB2	28:19:245:PRO:HD2	2.00	0.44
8:7E:64:LYS:HB3	8:7E:79:VAL:HG21	2.00	0.44
26:14:1986:A:H2'	26:14:1987:G:C8	2.53	0.44
1:13:901:A:C5	1:13:902:G:H1'	2.53	0.44
28:19:263:ARG:HB2	28:19:263:ARG:HE	1.30	0.44
27:16:79:C:H6	27:16:79:C:O5'	2.00	0.44
54:L5:8:ASN:C	54:L5:8:ASN:OD1	2.57	0.44
11:2I:114:VAL:HA	11:2I:115:PRO:HD2	1.90	0.44
2:12:190:THR:O	2:12:191:ASP:HB3	2.18	0.44
31:49:15:VAL:O	31:49:19:LEU:HD23	2.17	0.44
26:14:447:A:C8	26:14:473:G:C6	3.06	0.44
26:14:1461:G:H2'	26:14:1462:C:C6	2.52	0.44
13:4A:86:CYS:HB2	19:AA:73:GLU:HB3	1.99	0.44
26:14:1142(A):A:C4	26:14:1144:G:C8	3.06	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:250:G:H5'	36:78:60:MET:CE	2.48	0.44
26:1H:623:G:H2'	26:1H:624:C:C6	2.52	0.44
26:14:607:U:H5	26:14:619:G:C5	2.36	0.44
26:1H:1000:A:C8	26:1H:1154:G:N2	2.86	0.44
1:1G:1299:A:N1	1:1G:1301:U:N3	2.65	0.44
28:19:92:ILE:HG21	28:19:92:ILE:HD13	1.72	0.44
1:1G:51:A:C6	1:1G:353:A:C2	3.06	0.44
26:14:2115:G:H8	26:14:2117:A:OP2	2.00	0.44
26:14:2168:G:N2	26:14:2170:A:OP2	2.51	0.44
41:C8:92:ARG:HD3	41:C8:94:ASN:CB	2.48	0.44
26:1H:2691:C:O3'	26:1H:2871:C:H4'	2.17	0.44
51:M8:40:HIS:CG	51:M8:45:GLY:HA3	2.53	0.44
26:1H:143:C:H2'	26:1H:144:C:H6	1.83	0.44
1:1G:1331:G:OP1	1:1G:1331:G:H4'	2.17	0.44
26:1H:946:G:P	58:1H:4573:HOH:O	2.76	0.44
29:29:89:ASP:OD1	29:29:90:THR:N	2.50	0.44
38:98:44:LEU:HA	38:98:44:LEU:HD23	1.82	0.44
1:13:1280:A:H3'	1:13:1281:U:H5'	1.99	0.44
50:L8:38:GLU:OE2	50:L8:38:GLU:N	2.37	0.44
10:1I:63:PHE:HA	14:5I:58:LYS:HA	1.98	0.44
26:1H:1338:G:O2'	26:1H:1393:A:N1	2.42	0.44
26:14:1665:A:H1'	35:25:1:MET:HG3	2.00	0.44
26:1H:1038:C:H2'	26:1H:1039:G:O4'	2.18	0.44
47:E5:26:TYR:HB2	47:E5:29:GLN:OE1	2.17	0.44
46:D5:157:LEU:HA	46:D5:158:PRO:HD2	1.72	0.44
38:55:29:LEU:HB3	38:55:75:LEU:HD21	2.00	0.44
26:1H:1471:A:N3	26:1H:1472:A:C8	2.85	0.44
26:1H:1528:A:C2	26:1H:1543:A:N1	2.86	0.44
26:1H:1543:A:H3'	26:1H:1543:A:OP2	2.17	0.44
26:14:2850:A:N7	26:14:2868:A:O2'	2.45	0.44
34:15:128:HIS:CE1	34:15:134:ARG:HD2	2.52	0.44
53:K5:36:LEU:HA	53:K5:50:ARG:HG2	1.99	0.44
38:55:92:GLY:O	38:55:94:TYR:CE1	2.71	0.44
1:1G:1255:G:OP1	10:1A:45:ARG:NH1	2.48	0.44
26:1H:686:G:N2	54:P8:15:THR:HG21	2.33	0.44
26:14:2438:U:O3'	26:14:2439:A:H3'	2.16	0.44
26:14:2527:C:C4	26:14:2528:U:C5	3.06	0.44
40:B8:74:ARG:HD3	40:B8:76:PHE:CZ	2.52	0.44
47:I8:44:ARG:HG2	47:I8:45:PHE:CE1	2.53	0.44
42:D8:47:VAL:HG22	42:D8:48:GLY:N	2.33	0.44
42:D8:46:VAL:HG11	42:D8:52:VAL:HG21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:1261:A:H61	1:1G:1274:G:H1'	1.83	0.44
3:2E:7:PRO:O	3:2E:11:ARG:HG2	2.17	0.44
3:2E:8:ILE:O	3:2E:11:ARG:N	2.45	0.44
26:1H:2058:A:H5''	26:1H:2059:A:OP2	2.18	0.44
28:19:242:ARG:N	28:19:242:ARG:HD3	2.32	0.44
8:72:69:ARG:HD3	8:72:76:PRO:HA	1.99	0.44
33:61:29:TYR:C	33:61:32:PRO:HD2	2.38	0.44
26:14:2400:G:H3'	26:14:2401:U:C6	2.52	0.44
1:13:159:G:N2	1:13:161:A:H5''	2.33	0.44
4:3E:85:LYS:HG2	4:3E:86:LYS:N	2.32	0.44
51:15:40:HIS:CG	51:15:45:GLY:HA3	2.53	0.44
26:14:2123:G:C2	26:14:2176:A:C6	3.06	0.44
15:6I:33:THR:HG21	15:6I:85:LEU:HD22	1.99	0.44
26:1H:270(R):G:H2'	26:1H:270(S):G:C8	2.53	0.44
13:4A:76:ALA:HA	13:4A:79:LYS:HB2	2.00	0.44
26:14:1805:U:H5''	28:19:250:TRP:CD2	2.53	0.44
17:8A:21:VAL:HG11	17:8A:59:ILE:HG13	2.00	0.44
1:13:1521:G:H2'	1:13:1522:U:C6	2.53	0.44
5:4E:69:VAL:O	5:4E:71:LEU:N	2.51	0.44
10:1I:78:ASN:N	10:1I:78:ASN:OD1	2.51	0.44
26:14:2422:A:H2'	26:14:2422:A:H8	1.64	0.44
26:1H:320:A:H2'	30:31:136:THR:HG21	2.00	0.44
1:13:1187:G:O5'	9:8E:113:LYS:NZ	2.41	0.44
29:29:106:GLY:HA3	29:29:189:PRO:HB2	1.99	0.44
55:Q8:7:HIS:CD2	55:Q8:57:ARG:HH22	2.36	0.44
27:1J:109:G:C6	27:1J:110:G:C5	3.05	0.44
11:2A:29:ILE:HG22	11:2A:44:SER:CB	2.38	0.44
26:14:932:G:H4'	26:14:933:A:O5'	2.18	0.44
24:3K:4:C:H2'	24:3K:5:G:C8	2.53	0.44
1:1G:828:A:H2'	1:1G:829:G:O4'	2.18	0.44
46:D5:44:PHE:CE1	46:D5:48:PHE:HB2	2.53	0.44
1:13:1028(B):C:OP2	1:13:1029:G:H5''	2.18	0.44
30:39:102:PRO:O	30:39:105:VAL:N	2.48	0.44
26:14:469:G:OP2	58:14:3711:HOH:O	2.21	0.44
26:1H:2636:U:H2'	26:1H:2637:U:C6	2.52	0.44
28:19:148:GLU:CB	28:19:151:LYS:HD2	2.47	0.44
34:15:56:ASN:ND2	34:15:56:ASN:O	2.43	0.44
32:51:126:PRO:HG2	32:51:130:ARG:HH12	1.83	0.44
26:14:1331:A:H2'	26:14:1333:C:C5	2.53	0.44
34:58:127:ASP:O	34:58:128:HIS:HB3	2.18	0.44
11:2A:24:SER:OG	11:2A:26:ASN:N	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:78:144:GLU:HA	36:78:145:PRO:HD3	1.91	0.44
26:14:307:G:N2	26:14:309:G:H3'	2.33	0.44
36:35:126:VAL:HG13	36:35:145:PRO:HB2	2.00	0.44
1:13:1118:C:OP1	9:8E:104:ARG:NH1	2.49	0.44
26:1H:297:C:H2'	26:1H:298:G:O4'	2.18	0.44
26:1H:643:A:O2'	26:1H:644:A:H5'	2.18	0.44
1:13:407:G:O4'	4:3E:119:GLN:NE2	2.51	0.44
29:29:31:CYS:HB2	29:29:91:VAL:CG2	2.48	0.44
1:13:501:C:H2'	1:13:502:G:C8	2.52	0.44
4:32:108:LEU:HB3	4:32:110:PHE:CE1	2.53	0.44
35:68:88:ASN:OD1	35:68:90:GLN:HB2	2.18	0.44
26:1H:128:C:H2'	26:1H:129:C:H6	1.83	0.44
1:13:1226:C:H4'	19:AI:80:TYR:CZ	2.52	0.44
36:35:107:LYS:O	36:35:109:GLY:N	2.42	0.44
1:1G:965:A:C2	1:1G:969:A:C2	3.06	0.44
26:1H:1750:G:C2	26:1H:1751:C:C5	3.06	0.44
32:59:89:ILE:CG2	32:59:130:ARG:HA	2.48	0.44
45:G8:9:LYS:HA	45:G8:27:VAL:CG2	2.46	0.44
1:13:1015:A:H2'	1:13:1016:A:H8	1.83	0.44
26:1H:50:U:H3'	26:1H:51:G:C5'	2.48	0.44
1:1G:854:G:N1	1:1G:855:G:N7	2.66	0.44
1:1G:942:G:C2	1:1G:1342:C:C2	3.06	0.44
3:22:70:VAL:HG12	3:22:72:LYS:H	1.83	0.44
10:1A:99:LYS:CE	10:1A:100:THR:H	2.31	0.44
1:1G:1338:G:N7	1:1G:1339:A:C5	2.86	0.44
2:12:98:LEU:HD23	2:12:98:LEU:HA	1.83	0.44
36:78:42:SER:O	36:78:44:GLY:N	2.51	0.44
20:BI:53:LEU:H	20:BI:53:LEU:HD22	1.82	0.44
6:5E:41:GLU:HB2	6:5E:62:TRP:CE3	2.53	0.44
46:D5:10:ARG:HH21	46:D5:26:GLY:N	2.15	0.44
1:1G:1343:G:H2'	1:1G:1344:C:C6	2.53	0.44
30:39:53:THR:HG22	30:39:56:GLU:CD	2.38	0.44
1:13:61:G:H2'	1:13:62:U:O4'	2.18	0.44
7:62:137:LYS:O	7:62:141:VAL:HG23	2.18	0.44
1:13:1356:G:H2'	1:13:1357:A:C8	2.53	0.44
3:22:109:PRO:HB2	3:22:115:LEU:HD12	1.99	0.44
45:C5:36:ALA:HA	45:C5:67:LEU:O	2.17	0.44
26:14:2780:G:OP1	34:15:118:LYS:HE2	2.18	0.44
26:1H:2048:G:H1'	26:1H:2823:A:N6	2.32	0.44
20:BI:63:ILE:HG22	20:BI:77:ALA:HB1	1.99	0.44
13:4I:13:LYS:O	13:4I:44:ARG:NE	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:19:13:ARG:HD2	28:19:13:ARG:HA	1.59	0.44
1:13:1191:A:H2'	1:13:1192:C:C6	2.53	0.44
36:78:49:ARG:HG3	55:Q8:57:ARG:NE	2.33	0.43
26:1H:2308:G:N3	26:1H:2308:G:H2'	2.33	0.43
26:1H:1022:G:N2	26:1H:1142(A):A:N1	2.65	0.43
1:13:253:U:OP2	17:8I:67:LYS:NZ	2.44	0.43
13:4A:92:HIS:CD2	13:4A:98:VAL:HG11	2.53	0.43
29:21:105:THR:HG1	29:21:199:ARG:NH2	2.12	0.43
26:1H:919:G:H4'	27:16:81:G:H4'	2.00	0.43
26:1H:2400:G:C4	26:1H:2401:U:C5	3.06	0.43
29:21:181:LEU:HA	29:21:181:LEU:HD12	1.59	0.43
48:J8:85:LEU:HD22	48:J8:85:LEU:N	2.33	0.43
36:78:29:LYS:HB3	36:78:30:THR:H	1.50	0.43
29:29:73:GLU:HA	29:29:74:PRO:HD2	1.61	0.43
26:1H:1137:G:H2'	26:1H:1138:G:H8	1.82	0.43
42:95:5:VAL:HB	42:95:37:VAL:HG11	2.00	0.43
53:K5:28:ARG:CD	53:K5:32:ASN:H	2.31	0.43
34:58:96:GLU:HB2	34:58:122:VAL:CG1	2.48	0.43
26:1H:883:G:H2'	26:1H:884:C:H4'	1.99	0.43
1:13:1310:G:O2'	1:13:1311:G:H5'	2.17	0.43
1:13:1284:C:H3'	1:13:1285:A:H8	1.83	0.43
1:13:1453:G:HO2'	1:13:1454:G:P	2.41	0.43
24:3L:8:U:H4'	24:3L:9:A:OP1	2.17	0.43
1:1G:1158:C:N3	1:1G:1160:G:C8	2.86	0.43
1:13:1009:G:C2	1:13:1010:G:C8	3.05	0.43
26:14:2572:A:N7	29:29:144:ARG:HD2	2.32	0.43
26:14:2156:G:O6	26:14:2157:G:N2	2.51	0.43
8:7E:86:ILE:HG22	8:7E:87:SER:N	2.32	0.43
26:14:74:A:H4'	26:14:75:G:O5'	2.17	0.43
30:39:185:ASP:HA	30:39:188:ARG:HE	1.83	0.43
1:13:232:G:C5	1:13:233:C:C5	3.06	0.43
1:13:112:G:OP1	16:7I:27:LYS:HD2	2.18	0.43
1:13:51:A:C6	1:13:353:A:C2	3.06	0.43
1:13:458:C:H42	1:13:474:G:H1	1.66	0.43
26:14:2791:C:C4	26:14:2893:G:C6	3.06	0.43
10:1I:26:ALA:HB1	10:1I:84:GLN:HG2	2.00	0.43
1:13:765:G:H2'	58:13:1807:HOH:O	2.17	0.43
2:1E:25:ASN:HA	2:1E:26:PRO:HD3	1.73	0.43
26:14:1693:U:H4'	26:14:1694:C:OP2	2.19	0.43
31:41:47:LYS:NZ	31:41:81:LYS:HB2	2.33	0.43
26:14:1496:A:H8	26:14:1577:C:O2'	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:21:37:ARG:HA	29:21:42:ASP:OD2	2.18	0.43
2:1E:121:LEU:HA	2:1E:124:SER:OG	2.18	0.43
1:1G:818:G:O2'	1:1G:819:A:H5'	2.17	0.43
26:1H:2863:C:O2'	26:1H:2864:G:H5'	2.18	0.43
37:88:39:PRO:HA	37:88:97:VAL:O	2.17	0.43
1:1G:1346:A:C6	7:62:10:ARG:NH1	2.86	0.43
26:1H:247:G:O6	55:Q8:8:LYS:HB3	2.17	0.43
16:7A:23:ASP:OD1	16:7A:25:ARG:HG3	2.17	0.43
26:1H:483:A:O4'	45:G8:48:ALA:HB1	2.18	0.43
36:78:86:LYS:HB3	36:78:118:GLY:HA3	2.00	0.43
33:61:79:ILE:HD11	33:61:100:ALA:CB	2.48	0.43
7:6E:61:VAL:HG12	7:6E:124:LEU:HD23	1.99	0.43
26:1H:1945:G:H2'	26:1H:1946:U:C6	2.54	0.43
6:52:2:ARG:HB3	6:52:4:TYR:CE1	2.53	0.43
1:13:60:A:H4'	1:13:61:G:H5'	2.00	0.43
26:14:1647:G:OP2	58:14:3540:HOH:O	2.21	0.43
8:72:73:ASP:HA	8:72:74:PRO:HD2	1.77	0.43
26:1H:912:C:OP1	37:88:8:LYS:NZ	2.42	0.43
30:31:111:ALA:HB2	30:31:206:ILE:HG21	2.00	0.43
52:J5:45:VAL:HG22	52:J5:51:TYR:CE2	2.53	0.43
38:55:10:LEU:O	38:55:12:ARG:HG3	2.18	0.43
42:D8:2:PHE:O	42:D8:42:GLY:N	2.51	0.43
36:78:100:LEU:HA	36:78:100:LEU:HD12	1.70	0.43
33:69:38:LEU:HD12	33:69:38:LEU:H	1.83	0.43
46:H8:160:GLY:O	46:H8:161:VAL:HG13	2.18	0.43
26:1H:1288:U:C2	26:1H:1327:C:O2	2.71	0.43
26:14:992:C:H2'	26:14:993:G:H8	1.83	0.43
26:14:997:G:P	41:85:58:ARG:HH21	2.41	0.43
1:13:1367:C:H4'	10:11:48:THR:HG21	1.98	0.43
26:1H:780:G:N2	26:1H:783:A:H62	2.00	0.43
26:1H:1589:C:H2'	26:1H:1590:U:C6	2.53	0.43
30:31:59:TYR:CD1	30:31:78:ILE:HD11	2.54	0.43
26:14:2211:G:O2'	26:14:2212:A:P	2.76	0.43
26:1H:986:C:C2'	26:1H:987:G:H5'	2.48	0.43
30:39:83:PHE:O	30:39:84:VAL:HB	2.18	0.43
51:15:23:GLU:HG3	51:15:24:THR:N	2.33	0.43
52:N8:46:CYS:SG	52:N8:50:GLY:HA2	2.59	0.43
1:13:465:A:H2'	1:13:466:C:H5''	2.00	0.43
34:58:42:TRP:HA	34:58:48:MET:SD	2.58	0.43
17:8I:14:LYS:HD2	17:8I:14:LYS:N	2.33	0.43
26:14:2748:A:H2'	26:14:2749:A:C8	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:B5:5:TYR:CE1	49:G5:30:ARG:HG3	2.53	0.43
16:7A:14:ASN:O	16:7A:16:HIS:N	2.50	0.43
26:14:68:G:H2'	26:14:69:C:C6	2.53	0.43
24:3K:9:A:H3'	24:3K:10:G:C8	2.53	0.43
27:16:44:G:C2	27:16:48:A:C2	3.06	0.43
26:1H:2893:G:H4'	26:1H:2894:G:O4'	2.19	0.43
23:2K:3:C:N4	23:2K:71:G:H1	2.13	0.43
26:1H:1071:G:P	26:1H:1071:G:H8	2.41	0.43
26:14:1954:G:H1'	26:14:1956:U:O4	2.17	0.43
1:13:312:C:H2'	1:13:313:A:H8	1.83	0.43
26:1H:2518:A:H5'	26:1H:2518:A:C8	2.53	0.43
16:7I:12:LYS:C	16:7I:14:ASN:H	2.20	0.43
1:13:724:G:O6	1:13:733:A:N6	2.51	0.43
26:1H:639:U:H3	26:1H:649:G:H1	1.66	0.43
19:AA:66:MET:HA	19:AA:67:VAL:O	2.18	0.43
1:1G:198:G:H2'	1:1G:199:G:C8	2.51	0.43
26:1H:1766:U:O2'	26:1H:1767:C:H5'	2.18	0.43
2:1E:93:VAL:HG21	2:1E:97:TRP:HD1	1.83	0.43
26:14:1756:G:OP2	58:14:3804:HOH:O	2.21	0.43
26:1H:460:A:H3'	26:1H:461:C:H6	1.82	0.43
4:32:101:LEU:HD23	4:32:121:VAL:HG13	1.99	0.43
26:1H:2862:G:H2'	26:1H:2863:C:H6	1.83	0.43
1:13:590:C:H2'	1:13:591:U:C6	2.53	0.43
1:1G:814:A:N7	1:1G:816:A:C4	2.86	0.43
8:7E:122:ARG:HB3	8:7E:126:LYS:HE2	1.99	0.43
3:2E:23:TYR:CD1	3:2E:24:ALA:N	2.87	0.43
31:41:122:PRO:HB3	31:41:180:PHE:HD2	1.81	0.43
11:2I:41:THR:HG21	11:2I:71:LYS:CB	2.49	0.43
36:78:97:PRO:HA	36:78:100:LEU:HB2	1.99	0.43
12:3I:110:VAL:CG2	12:3I:120:TYR:HB3	2.48	0.43
51:M8:43:TYR:O	51:M8:46:GLN:HA	2.18	0.43
1:1G:636:U:H2'	1:1G:637:G:C8	2.54	0.43
40:75:29:ARG:HD3	40:75:44:ASP:OD2	2.18	0.43
29:21:21:VAL:HA	29:21:22:PRO:HD3	1.53	0.43
3:22:61:ALA:C	3:22:63:ASN:H	2.22	0.43
1:1G:552:U:O2'	12:3A:86:ARG:O	2.36	0.43
26:1H:463:G:N2	26:1H:466:A:OP2	2.39	0.43
26:1H:2038:G:H2'	26:1H:2039:C:H6	1.83	0.43
3:2E:108:ASN:ND2	3:2E:110:ASN:HB2	2.33	0.43
16:7A:34:GLU:OE1	16:7A:55:ARG:NH1	2.52	0.43
32:59:99:VAL:HG13	32:59:100:GLY:H	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:135:C:C2	16:7A:1:MET:HB3	2.53	0.43
6:52:39:LYS:HE3	6:52:64:GLN:HE21	1.83	0.43
26:14:1438:U:O2	26:14:1438:U:H2'	2.16	0.43
51:15:8:LYS:HD3	51:15:8:LYS:HA	1.82	0.43
26:14:2743:C:H2'	26:14:2744:G:O4'	2.17	0.43
31:41:10:LYS:O	31:41:15:VAL:HG23	2.18	0.43
55:Q8:23:VAL:HG22	55:Q8:24:ALA:N	2.32	0.43
55:Q8:45:GLY:HA2	55:Q8:46:ARG:C	2.38	0.43
26:1H:375:C:H2'	26:1H:376:C:C6	2.54	0.43
26:1H:895:U:OP1	26:1H:895:U:H4'	2.16	0.43
55:M5:14:VAL:CG1	55:M5:22:VAL:HG13	2.48	0.43
30:39:117:ARG:HG3	30:39:192:LEU:HD12	2.00	0.43
1:13:525:C:O5'	1:13:525:C:H6	2.01	0.43
26:1H:1389:G:C2	26:1H:1399:C:O2	2.71	0.43
26:14:2301:C:N4	26:14:2315:G:O6	2.51	0.43
13:4A:91:ARG:NE	13:4A:97:PRO:O	2.51	0.43
41:C8:88:ILE:O	41:C8:88:ILE:HG22	2.19	0.43
26:14:84:A:H5''	45:C5:8:LYS:HG3	1.99	0.43
26:14:898:C:N4	26:14:899:A:N3	2.66	0.43
26:14:2468:G:H3'	26:14:2476:A:C2	2.53	0.43
26:14:2129:C:C4	26:14:2130:U:C4	3.06	0.43
20:BA:64:ASP:OD1	20:BA:81:LYS:HD2	2.18	0.43
37:88:30:GLY:CA	37:88:107:ALA:HB2	2.47	0.43
26:14:2065:C:H5''	26:14:2252:G:H1'	2.00	0.43
26:1H:140:A:C8	26:1H:1408:C:O2'	2.69	0.43
1:13:611:A:H61	1:13:629:G:H1	1.65	0.43
44:F8:31:HIS:CD2	44:F8:33:LYS:H	2.37	0.43
44:B5:31:HIS:HA	44:B5:32:PRO:HD3	1.77	0.43
46:D5:59:LEU:HA	46:D5:59:LEU:HD22	1.65	0.43
50:H5:50:VAL:O	50:H5:54:VAL:HG23	2.18	0.43
1:13:939:G:C6	1:13:940:C:N4	2.86	0.43
1:13:993:G:N3	1:13:993:G:H2'	2.33	0.43
26:1H:981:A:H1'	26:1H:2037:G:O4'	2.18	0.43
20:BA:88:VAL:O	20:BA:92:LEU:HG	2.19	0.43
26:1H:1491:G:O4'	28:11:99:ASP:HB3	2.18	0.43
26:14:2562:U:H4'	35:25:25:LEU:HD21	2.00	0.43
26:14:1953:A:H5''	26:14:1954:G:OP2	2.18	0.43
1:1G:1225:A:H5''	13:4A:103:THR:OG1	2.17	0.43
26:1H:1337:G:C4	26:1H:1338:G:C8	3.06	0.43
1:13:1098:C:C2	1:13:1099:G:C8	3.06	0.43
26:1H:2137:C:N3	26:1H:2138:C:N4	2.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:32:59:ARG:O	4:32:63:LYS:N	2.52	0.43
46:H8:26:GLY:HA2	46:H8:85:HIS:CD2	2.53	0.43
1:13:1320:C:N4	19:AI:36:ARG:HG3	2.34	0.43
46:D5:148:ASP:OD2	46:D5:175:VAL:HG11	2.18	0.43
26:14:2295:C:H41	39:65:13:ARG:NH2	2.15	0.43
1:13:818:G:C2	1:13:820:U:O2'	2.69	0.43
33:61:11:ASN:O	33:61:12:LEU:HB2	2.17	0.43
26:1H:2862:G:C4	26:1H:2863:C:C5	3.07	0.43
26:14:192:C:OP1	58:14:4099:HOH:O	2.21	0.43
26:1H:1705:G:C6	26:1H:1706:U:N3	2.86	0.43
7:62:105:VAL:O	7:62:108:ALA:HB3	2.18	0.43
26:1H:2358:G:H2'	26:1H:2359:C:H6	1.83	0.43
26:14:2576:G:O2'	26:14:2579:C:OP2	2.30	0.43
26:1H:433:C:H2'	26:1H:434:U:C6	2.54	0.43
1:1G:895:G:H1	1:1G:904:C:H42	1.64	0.43
45:G8:45:VAL:N	45:G8:63:LYS:O	2.43	0.43
1:13:955:U:H1'	1:13:1227:A:N6	2.33	0.43
1:1G:59:A:N1	58:1G:1824:HOH:O	2.36	0.43
26:14:948:G:C2	26:14:970:C:O2	2.72	0.43
1:1G:1103:C:H2'	1:1G:1104:G:O4'	2.18	0.43
46:D5:52:SER:C	46:D5:54:HIS:H	2.21	0.43
26:1H:476:G:N1	26:1H:479:A:OP2	2.52	0.43
49:G5:18:PRO:O	49:G5:21:LEU:HB2	2.19	0.43
1:1G:137:C:C2	1:1G:227:G:C2	3.06	0.43
35:68:91:LEU:HA	35:68:91:LEU:HD12	1.72	0.43
19:AI:81:ARG:HB2	19:AI:81:ARG:HE	1.61	0.43
29:21:14:ILE:HA	29:21:14:ILE:HD13	1.63	0.43
26:1H:1885:A:H2'	26:1H:1886:C:O4'	2.17	0.43
13:4I:22:ILE:HB	13:4I:25:ILE:HD12	2.01	0.43
27:1J:66:A:C6	27:1J:108:C:C6	3.06	0.43
26:1H:2592:G:C6	26:1H:2593:U:N3	2.87	0.43
26:14:881:G:N7	26:14:882:G:C2	2.87	0.43
41:85:61:TRP:O	41:85:65:ILE:HD13	2.18	0.43
26:14:1997:G:OP2	58:14:3495:HOH:O	2.21	0.43
26:1H:2027:G:O2'	26:1H:2028:U:H5'	2.18	0.43
26:1H:1257:C:H4'	30:31:83:PHE:HD1	1.77	0.43
26:1H:74:A:H8	26:1H:74:A:H5''	1.83	0.43
26:14:2287:A:C2	26:14:2289:G:C8	3.07	0.43
26:14:2115:G:H4'	26:14:2168:G:N7	2.33	0.43
34:58:96:GLU:CG	34:58:97:ARG:H	2.31	0.43
26:1H:2692:C:O2'	26:1H:2693:A:H5'	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:966:G:H2'	26:14:967:C:C6	2.52	0.43
26:1H:2572:A:C8	29:21:144:ARG:HB3	2.53	0.43
26:1H:2572:A:C4	29:21:144:ARG:NH1	2.87	0.43
1:1G:537:G:H5''	12:3A:113:ARG:NH1	2.34	0.43
1:13:626:U:H2'	1:13:627:G:C8	2.53	0.43
1:1G:589:C:H42	1:1G:650:G:H1	1.66	0.43
26:14:1638:C:H5''	26:14:2710:C:O2	2.18	0.43
30:39:46:ARG:HG2	30:39:46:ARG:NH1	2.32	0.43
26:1H:1532:C:N3	26:1H:1539:G:N2	2.62	0.43
45:C5:90:LEU:HA	45:C5:91:GLU:HA	1.86	0.43
41:C8:28:ARG:O	41:C8:35:ALA:HA	2.19	0.43
26:1H:515:A:H1'	26:1H:581:C:H1'	2.01	0.43
23:2K:21:U:H2'	23:2K:21:U:H6	1.61	0.43
7:62:50:ILE:HG21	7:62:58:PRO:HA	1.99	0.43
1:1G:1065:U:C5	1:1G:1190:G:N3	2.87	0.43
36:35:147:LEU:HD23	36:35:148:LEU:H	1.83	0.43
1:1G:15:G:H4'	5:42:24:ARG:HH12	1.82	0.43
1:1G:953:G:H2'	1:1G:954:G:O4'	2.18	0.43
55:M5:29:LYS:HB3	55:M5:44:LYS:HD2	2.01	0.43
26:14:637:A:H2'	36:35:117:GLU:OE2	2.19	0.43
26:1H:507:A:C5'	26:1H:508:G:H5'	2.47	0.43
3:22:35:GLU:HG2	3:22:39:ILE:HD11	2.00	0.43
31:41:81:LYS:HB3	31:41:82:LEU:H	1.57	0.43
30:31:184:TYR:CE1	36:78:3:LEU:HD11	2.53	0.43
26:14:1534:G:H3'	26:14:1535:U:C5'	2.49	0.43
19:AA:66:MET:N	19:AA:67:VAL:HB	2.34	0.43
1:1G:198:G:C6	1:1G:220:G:C2	3.06	0.43
3:22:50:ALA:HB3	3:22:76:VAL:HG12	2.00	0.43
26:1H:2337:G:H2'	26:1H:2338:G:H8	1.83	0.43
13:4I:84:ILE:HD11	19:AI:66:MET:CG	2.48	0.43
8:72:69:ARG:NH1	8:72:75:ARG:O	2.51	0.43
26:14:1138:G:H21	34:15:106:MET:CE	2.32	0.43
55:M5:15:LYS:HE3	55:M5:16:ILE:O	2.18	0.43
26:14:1259:G:H2'	26:14:1260:G:C8	2.54	0.43
37:45:1:MET:HB2	37:45:69:PHE:HE1	1.83	0.43
46:H8:23:LYS:HD3	46:H8:40:ASP:HA	1.99	0.43
26:14:120:U:P	58:14:3994:HOH:O	2.76	0.43
26:14:816:C:P	58:14:3587:HOH:O	2.77	0.43
13:4A:15:VAL:HG12	13:4A:45:VAL:HG22	1.99	0.43
41:85:16:LYS:O	41:85:20:LEU:HD13	2.19	0.43
26:14:1833:U:H2'	26:14:1834:U:O4'	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1615:C:C5	26:14:1617:C:C4	3.06	0.43
6:52:14:LEU:HD12	6:52:18:GLN:HB3	2.00	0.43
26:14:664:C:H2'	26:14:665:C:H6	1.82	0.43
36:78:31:ALA:O	36:78:32:THR:HB	2.19	0.43
46:D5:102:LEU:HD22	46:D5:104:PHE:HE1	1.84	0.43
46:D5:67:LEU:HA	46:D5:68:PRO:HD3	1.68	0.43
36:35:51:PHE:CE2	36:35:53:GLY:HA2	2.53	0.43
40:B8:21:GLU:HG3	40:B8:21:GLU:H	1.56	0.43
6:5E:14:LEU:HD23	6:5E:14:LEU:HA	1.53	0.43
26:14:1214:A:O5'	26:14:1214:A:H8	2.01	0.43
28:11:269:PHE:CD1	28:11:269:PHE:N	2.86	0.43
1:13:799:G:C6	1:13:800:G:C4	3.06	0.43
55:Q8:49:VAL:CG2	55:Q8:50:LEU:H	2.22	0.43
26:1H:1329:U:H3'	26:1H:1330:C:H6	1.82	0.43
26:1H:449:A:C6	26:1H:450:G:C5	3.06	0.43
1:13:1126:U:N3	1:13:1127:G:C2	2.86	0.43
1:13:1126:U:O2'	1:13:1127:G:OP1	2.34	0.43
26:1H:444:C:H2'	26:1H:445:C:H6	1.83	0.43
48:J8:85:LEU:HD13	48:J8:85:LEU:HA	1.48	0.43
36:35:62:LEU:CD2	55:M5:27:THR:HA	2.48	0.43
46:D5:87:ASP:N	46:D5:87:ASP:OD1	2.51	0.43
26:14:2300:G:N2	26:14:2317:C:C2	2.87	0.43
9:82:119:ALA:O	9:82:120:ARG:HB2	2.18	0.43
31:41:67:LYS:HZ3	51:M8:6:HIS:CE1	2.35	0.43
26:1H:2125:G:N2	26:1H:2172:U:OP1	2.52	0.43
1:13:926:G:H5'	1:13:927:G:C5'	2.49	0.43
1:1G:1252:A:H61	1:1G:1285:A:H61	1.67	0.43
1:1G:376:G:H5''	16:7A:5:ARG:HD3	2.01	0.43
26:14:2262:U:H4'	26:14:2328:A:H2	1.83	0.43
26:14:35:G:H2'	26:14:36:G:O4'	2.17	0.43
39:65:34:HIS:CD2	39:65:54:LEU:HD13	2.54	0.43
26:1H:142:G:H1'	44:F8:37:THR:CG2	2.44	0.43
26:14:40:C:H2'	26:14:41:C:C6	2.51	0.43
26:1H:1586:A:H3'	26:1H:1587:A:C8	2.46	0.43
26:1H:2849:U:OP2	40:B8:95:ARG:NH1	2.52	0.43
29:29:91:VAL:HG23	29:29:91:VAL:O	2.18	0.43
10:1I:40:LEU:HB3	10:1I:41:PRO:HD2	1.99	0.43
26:1H:583:G:H5''	41:C8:10:ARG:NH1	2.33	0.43
32:59:41:MET:HB3	32:59:42:ARG:H	1.64	0.43
26:14:1953:A:H2	26:14:2549:G:N3	2.17	0.43
26:14:2038:G:H2'	26:14:2039:C:C6	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:2256:G:C5	26:14:2257:U:C5	3.06	0.43
26:14:1485:G:H1	26:14:1504:C:H42	1.67	0.43
32:59:130:ARG:O	32:59:131:VAL:HB	2.18	0.43
26:1H:51:G:N3	26:1H:119:A:C2	2.86	0.43
37:88:68:ILE:CG2	37:88:103:MET:HB3	2.47	0.43
26:14:2391:G:O6	26:14:2425:A:H8	2.01	0.43
6:52:10:LEU:N	6:52:10:LEU:HD12	2.33	0.43
27:1J:113:C:OP2	27:1J:113:C:H6	2.02	0.43
2:12:118:LEU:O	2:12:122:PHE:N	2.51	0.43
26:1H:2206:C:H2'	26:1H:2207:C:C6	2.52	0.43
26:14:2187:G:C6	26:14:2188:C:C4	3.07	0.43
3:2E:24:ALA:HB1	3:2E:28:GLN:HB2	2.00	0.43
9:82:121:ARG:NH1	9:82:122:ALA:O	2.52	0.43
26:1H:738:G:C6	26:1H:739:G:C2	3.06	0.43
3:22:59:ARG:HG2	3:22:64:VAL:HG23	2.01	0.43
1:1G:1288:A:N1	1:1G:1371:G:H1'	2.34	0.43
1:13:595:G:H22	1:13:643:C:H41	1.66	0.43
44:B5:25:LYS:HA	44:B5:81:VAL:O	2.19	0.43
31:41:41:GLN:HG3	31:41:60:LEU:HD11	1.99	0.43
27:16:61:G:C6	27:16:62:C:C4	3.06	0.43
8:72:83:ILE:HB	8:72:137:VAL:HG13	1.99	0.43
1:1G:596:C:H2'	1:1G:597:G:C8	2.53	0.43
21:1F:12:LYS:O	21:1F:16:GLY:N	2.51	0.43
1:13:256:U:H2'	1:13:257:G:C8	2.53	0.43
24:3K:13:C:O2'	24:3K:14:A:H5'	2.18	0.43
29:21:66:HIS:O	29:21:66:HIS:ND1	2.51	0.43
44:F8:66:LEU:O	44:F8:66:LEU:HG	2.19	0.43
10:1I:23:ILE:H	10:1I:23:ILE:HG13	1.67	0.43
1:1G:389:A:N3	1:1G:389:A:H2'	2.33	0.43
6:52:21:LEU:HD22	6:52:21:LEU:HA	1.76	0.43
26:1H:2643:G:H2'	26:1H:2644:G:O4'	2.19	0.43
55:Q8:21:LYS:HE2	55:Q8:21:LYS:C	2.38	0.43
55:Q8:14:VAL:CG1	55:Q8:21:LYS:HZ1	2.31	0.43
26:1H:2334:G:C2	39:A8:12:PHE:CD1	3.07	0.43
26:1H:445:C:OP1	58:1H:3702:HOH:O	2.20	0.43
26:1H:1793:C:O2	26:1H:1900:A:H2	2.02	0.43
33:69:76:THR:HG22	33:69:139:GLN:O	2.18	0.43
38:98:10:LEU:O	38:98:12:ARG:HG2	2.18	0.43
26:1H:1179:C:H2'	26:1H:1180:C:C6	2.54	0.43
33:61:69:LYS:O	33:61:73:GLU:HB2	2.18	0.43
1:13:717:C:H5''	1:13:717:C:C6	2.42	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1284:C:H3'	1:13:1285:A:C8	2.54	0.43
26:14:2031:A:C6	26:14:2498:C:H1'	2.53	0.43
41:C8:92:ARG:NH2	42:D8:10:LYS:HB3	2.33	0.43
38:98:92:GLY:N	38:98:94:TYR:HE1	2.15	0.43
2:1E:88:ALA:HB3	2:1E:90:MET:HG2	2.01	0.43
26:1H:547:A:C6	26:1H:548:A:C6	3.07	0.43
26:14:2056:G:N3	26:14:2056:G:H2'	2.34	0.43
29:29:199:ARG:HB3	29:29:200:GLU:OE1	2.19	0.43
32:51:92:ILE:HG13	32:51:92:ILE:H	1.59	0.43
55:M5:4:MET:HB3	55:M5:60:LEU:HD13	2.00	0.43
35:25:31:LYS:HB3	35:25:32:TYR:CE1	2.54	0.43
1:13:1096:C:H2'	1:13:1097:C:C6	2.51	0.43
15:6I:78:TYR:CE1	15:6I:82:ILE:HD11	2.54	0.43
26:1H:1607:C:N4	26:1H:1621:U:H2'	2.32	0.43
4:3E:173:TRP:CG	4:3E:189:PRO:HG3	2.53	0.43
29:21:97:LYS:O	29:21:100:GLU:HG3	2.18	0.43
26:14:2002:G:N7	58:14:3797:HOH:O	2.36	0.43
24:3K:28:G:H2'	24:3K:29:G:C8	2.53	0.43
13:4I:84:ILE:HG23	13:4I:86:CYS:HB3	2.00	0.43
26:14:2140:C:H1'	26:14:2152:G:N2	2.34	0.43
42:D8:64:HIS:ND1	42:D8:92:THR:HG22	2.33	0.43
2:12:119:GLU:HA	2:12:122:PHE:HD2	1.83	0.43
2:1E:170:GLU:N	2:1E:170:GLU:CD	2.71	0.43
7:62:111:ARG:NH1	7:62:126:ASP:OD2	2.31	0.43
26:14:1138:G:C4	26:14:1139:G:H1'	2.52	0.43
27:1J:28:C:P	39:65:36:TYR:HH	2.42	0.43
1:13:380:G:C2	1:13:384:G:C6	3.06	0.43
8:72:89:PRO:HA	8:72:92:ARG:HH21	1.83	0.43
26:14:513:A:C2	26:14:514:A:C4	3.06	0.43
1:13:595:G:N2	1:13:643:C:N4	2.66	0.43
1:1G:1171:G:H2'	1:1G:1172:C:C6	2.54	0.43
30:31:192:LEU:HD23	30:31:193:VAL:N	2.33	0.43
1:1G:823:G:H21	8:72:1:MET:HE2	1.82	0.43
26:14:2547:U:O2	35:25:23:ARG:NH2	2.50	0.43
26:14:844:C:C5	26:14:845:G:C6	3.07	0.43
1:1G:1313:U:OP1	19:AA:7:LYS:N	2.51	0.43
26:14:480:A:H1'	45:C5:44:ILE:HG12	2.01	0.43
40:75:106:SER:HA	40:75:110:ILE:HD12	1.99	0.43
1:1G:404:U:OP1	4:32:118:ARG:NH1	2.48	0.43
15:6I:69:TYR:CE1	15:6I:73:GLU:HG3	2.54	0.43
41:85:31:SER:OG	41:85:31:SER:O	2.37	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:12:22:LYS:HB3	2:12:40:HIS:CE1	2.54	0.43
37:45:55:VAL:HG23	37:45:64:ILE:HD12	1.99	0.43
7:62:74:GLU:O	7:62:89:MET:N	2.51	0.43
1:1G:407:G:H2'	1:1G:408:A:C8	2.53	0.43
26:14:335:C:H4'	45:C5:73:ARG:CZ	2.48	0.43
26:14:1204:A:HO2'	26:14:1205:U:P	2.41	0.43
1:13:1128:C:O2'	1:13:1139:G:O6	2.37	0.43
41:85:66:ASN:CB	41:85:76:TYR:HB2	2.41	0.43
26:14:1161:C:H1'	42:95:8:GLY:O	2.18	0.43
37:88:77:LYS:HE3	37:88:82:ARG:HA	2.00	0.43
3:2E:59:ARG:HA	3:2E:63:ASN:O	2.19	0.43
19:AA:71:LEU:HA	19:AA:71:LEU:HD23	1.75	0.43
26:1H:2061:G:H5'	58:1H:3633:HOH:O	2.18	0.43
26:1H:1510:A:N3	26:1H:1510:A:H2'	2.33	0.43
26:14:1900:A:N1	26:14:1970:A:C6	2.86	0.43
26:14:607:U:N3	26:14:621:A:C2	2.78	0.43
1:1G:1299:A:C6	1:1G:1301:U:C2	3.06	0.43
26:14:2508:G:HO2'	26:14:2554:U:HO2'	1.66	0.43
26:14:2298:A:N6	26:14:2318:G:C8	2.86	0.43
34:15:17:ASP:O	34:15:18:ALA:HB3	2.18	0.43
55:Q8:38:GLY:HA2	55:Q8:39:LYS:C	2.39	0.43
39:A8:29:PHE:CD1	39:A8:30:ARG:N	2.82	0.43
36:78:57:THR:HB	36:78:59:LEU:H	1.83	0.43
1:1G:1317:C:H5''	1:1G:1318:A:OP2	2.19	0.43
1:13:683:G:C6	1:13:684:A:C6	3.07	0.43
1:1G:1292:U:H5'	9:82:38:GLN:NE2	2.34	0.43
26:14:1608:A:H1'	26:14:1610:A:OP2	2.19	0.43
37:88:34:LEU:HD23	37:88:104:PHE:HD2	1.82	0.43
26:1H:836:G:C5	26:1H:837:C:C4	3.07	0.43
46:D5:59:LEU:HB3	46:D5:60:GLU:H	1.48	0.43
9:82:10:ARG:HD2	9:82:105:ASP:HB3	2.00	0.43
29:29:33:VAL:HG12	29:29:89:ASP:CB	2.44	0.43
48:F5:85:LEU:HB3	48:F5:87:PRO:HG2	1.99	0.43
31:49:64:THR:OG1	31:49:94:LEU:HD21	2.19	0.43
27:1J:9:G:OP1	39:65:25:ARG:NH2	2.51	0.43
26:14:1954:G:N2	26:14:1956:U:C2	2.87	0.43
37:88:110:THR:HG23	37:88:113:GLN:OE1	2.18	0.43
26:1H:1419:A:C8	26:1H:1421:G:C6	3.07	0.43
1:1G:316:G:H2'	1:1G:317:G:C8	2.53	0.43
36:35:110:TYR:HB3	36:35:111:ARG:H	1.47	0.43
3:22:18:TRP:HE1	14:5A:55:GLY:N	2.16	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1198:U:C2	26:14:1199:U:C5	3.07	0.43
43:E8:64:MET:O	43:E8:65:LEU:HB2	2.18	0.43
1:1G:1274:G:N2	1:1G:1275:A:N7	2.57	0.43
7:62:15:ASP:HB3	7:62:19:GLY:N	2.34	0.43
26:1H:1826:G:H4'	28:11:242:ARG:CZ	2.48	0.43
26:14:686:G:N2	26:14:788:A:H61	2.16	0.43
41:C8:97:ASP:HB2	41:C8:101:ARG:HH11	1.83	0.43
26:14:341:G:C6	26:14:342:G:C5	3.06	0.43
2:1E:149:LEU:HD23	2:1E:149:LEU:HA	1.79	0.43
2:1E:149:LEU:O	2:1E:153:ARG:N	2.52	0.43
26:1H:1181:C:O2'	26:1H:1182:A:H5'	2.19	0.43
26:1H:2821:A:H2'	26:1H:2822:G:O4'	2.19	0.43
26:14:2480:C:H5'	26:14:2481:G:OP2	2.18	0.43
18:9I:38:GLU:HA	18:9I:41:LYS:NZ	2.33	0.43
26:14:1375:C:H2'	26:14:1376:C:H6	1.82	0.43
41:C8:66:ASN:CB	41:C8:76:TYR:HB2	2.49	0.43
1:13:75:C:H2'	1:13:76:G:O4'	2.18	0.43
1:13:585:G:O2'	1:13:879:C:H5''	2.19	0.43
47:I8:51:VAL:N	47:I8:62:LEU:HD12	2.33	0.43
30:39:160:ASN:HB3	30:39:163:VAL:HB	2.00	0.43
1:13:349:A:O2'	1:13:350:G:H5'	2.18	0.43
32:51:95:ARG:HB3	32:51:95:ARG:NH1	2.34	0.43
26:1H:1321:A:H2'	26:1H:1322:A:O4'	2.18	0.43
2:1E:223:ILE:HG12	2:1E:223:ILE:H	1.55	0.43
26:14:238:C:C2	26:14:260:G:C2	3.06	0.43
26:14:1812:A:O2'	28:19:45:ASN:N	2.51	0.43
46:H8:18:LEU:O	46:H8:21:ALA:HB3	2.18	0.43
55:Q8:30:ARG:CZ	55:Q8:30:ARG:HB2	2.49	0.43
26:1H:2593:U:C4	58:1H:3683:HOH:O	2.57	0.43
26:14:993:G:N3	42:95:89:GLN:NE2	2.59	0.43
26:1H:265:A:C2	26:1H:428:A:C2	3.07	0.43
26:14:1021:A:C8	26:14:1021:A:C3'	3.02	0.43
26:14:619:G:H5'	26:14:620:G:OP2	2.19	0.43
7:6E:113:GLU:HB2	7:6E:118:VAL:HG13	2.00	0.43
32:51:52:VAL:HG12	32:51:65:HIS:CD2	2.53	0.43
26:14:2299:G:N1	26:14:2318:G:H8	2.17	0.43
26:1H:518:G:H2'	26:1H:519:U:C6	2.54	0.43
26:14:2207:C:N4	26:14:2217:G:H1	2.08	0.43
33:69:109:ILE:HB	33:69:130:TYR:OH	2.18	0.43
26:14:909:A:C8	26:14:912:C:N4	2.87	0.43
26:14:1754:C:N3	26:14:2716:U:O2'	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:375:U:O3'	16:7A:6:LEU:HB2	2.17	0.43
37:88:133:ARG:O	37:88:134:ARG:HB2	2.17	0.43
26:1H:58:G:N2	26:1H:70:G:C4	2.87	0.43
26:14:2228:G:C6	26:14:2229:C:C4	3.07	0.43
26:14:1848:A:C4	26:14:1849:G:C8	3.06	0.43
26:14:900:A:H5''	26:14:901:A:OP2	2.19	0.43
26:1H:183:C:H42	26:1H:213:A:H61	1.67	0.43
1:1G:532:A:H2	3:22:156:ARG:HH12	1.64	0.43
1:13:1316:G:H5''	14:5I:17:LYS:NZ	2.34	0.43
26:1H:581:C:OP1	41:C8:33:ARG:HG3	2.18	0.43
1:13:68:G:N2	1:13:69:G:H1'	2.33	0.43
26:1H:1843:C:H6	26:1H:1843:C:O5'	2.02	0.43
44:F8:21:PHE:HE1	44:F8:92:LEU:HD13	1.83	0.43
36:35:2:LYS:O	36:35:5:ASP:HB2	2.19	0.43
26:1H:481:G:H1'	26:1H:507:A:N1	2.34	0.43
26:1H:8:A:H2'	26:1H:9:U:C6	2.53	0.43
26:14:2015:A:H1'	52:J5:2:ALA:HA	2.01	0.43
18:9I:31:LEU:H	18:9I:31:LEU:HD23	1.84	0.43
26:1H:270(V):G:H2'	26:1H:270(W):G:C8	2.53	0.43
37:45:19:GLY:H	37:45:98:LYS:HZ3	1.66	0.43
8:7E:94:TYR:HE1	8:7E:132:GLU:HB2	1.83	0.43
46:D5:11:GLU:HG3	46:D5:12:GLY:H	1.83	0.43
20:BI:56:MET:HG3	20:BI:88:VAL:HG21	2.00	0.43
13:4I:54:VAL:O	13:4I:58:GLU:HG3	2.19	0.43
26:1H:1665:A:H1'	35:68:1:MET:HG3	2.01	0.43
26:1H:775:G:C4	26:1H:794:G:C8	3.06	0.43
26:14:2040:C:H2'	26:14:2041:U:O4'	2.19	0.43
26:14:2376:A:H2	39:65:112:PHE:HB3	1.84	0.43
26:14:1259:G:H2'	26:14:1260:G:H8	1.84	0.43
35:25:19:ILE:HD12	35:25:41:ALA:CB	2.49	0.43
37:45:1:MET:HB3	37:45:2:LEU:H	1.61	0.43
30:31:125:LEU:HD11	30:31:199:TRP:CD2	2.54	0.43
17:8A:10:VAL:HG23	17:8A:54:GLY:N	2.32	0.43
26:14:2091:U:OP2	26:14:2092:U:O2'	2.24	0.43
31:41:26:GLN:OE1	31:41:27:ASN:HB2	2.19	0.43
1:1G:1172:C:H2'	1:1G:1173:G:H8	1.83	0.43
1:1G:266:G:H5''	1:1G:268:C:H41	1.83	0.43
1:1G:756:C:H2'	1:1G:757:U:O4'	2.19	0.43
40:75:90:GLN:OE1	40:75:121:ILE:HD11	2.18	0.43
26:14:2072:G:H2'	26:14:2073:C:O4'	2.19	0.43
42:D8:35:LEU:O	42:D8:37:VAL:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:59:68:THR:O	32:59:72:ILE:HG13	2.19	0.43
33:69:65:ALA:O	33:69:68:LEU:N	2.51	0.43
1:13:101:A:H8	1:13:101:A:OP2	2.02	0.43
1:1G:620:C:H3'	1:1G:621:A:H8	1.84	0.43
26:14:2592:G:C6	26:14:2593:U:C4	3.07	0.43
39:65:23:ARG:NH2	39:65:84:GLN:HG3	2.34	0.43
26:1H:2017:U:P	58:1H:4643:HOH:O	2.76	0.43
26:1H:2298:A:H2'	26:1H:2299:G:O4'	2.19	0.43
7:62:92:SER:O	7:62:96:GLN:HG3	2.18	0.43
26:14:1141:U:P	34:15:63:THR:HG21	2.59	0.43
1:1G:674:G:N2	11:2A:116:HIS:HB2	2.33	0.43
7:6E:111:ARG:HB3	7:6E:113:GLU:OE2	2.18	0.43
26:1H:1484:G:C2	26:1H:1485:G:C8	3.06	0.43
26:1H:1153:C:C4	26:1H:1154:G:C5	3.07	0.43
26:14:139:G:N2	26:14:141:A:N1	2.67	0.43
5:4E:110:LEU:O	5:4E:115:VAL:HB	2.19	0.43
26:1H:1727:U:H2'	26:1H:1728:G:O4'	2.18	0.43
1:1G:979:C:C5	1:1G:980:C:C6	3.00	0.43
26:14:2305:A:H8	31:49:156:ASP:OD2	2.02	0.43
26:14:2113:U:O4	26:14:2168:G:O2'	2.32	0.43
1:13:412:A:H1'	1:13:413:G:OP2	2.19	0.43
45:G8:94:LYS:HG3	45:G8:95:LYS:N	2.33	0.43
26:1H:950:G:H2'	26:1H:951:C:H6	1.82	0.43
8:7E:88:LYS:O	8:7E:92:ARG:HD3	2.18	0.43
55:M5:35:GLN:O	55:M5:35:GLN:HG3	2.19	0.43
1:13:232:G:C5	1:13:233:C:C4	3.07	0.43
26:14:2640:G:OP1	34:15:97:ARG:NH1	2.45	0.43
16:7I:53:VAL:HG13	16:7I:79:VAL:HA	2.01	0.43
39:A8:41:ASP:OD2	39:A8:44:LYS:HB2	2.19	0.43
1:13:749:C:H2'	1:13:750:G:H8	1.83	0.43
36:78:80:TYR:CD2	36:78:111:ARG:HB3	2.53	0.43
26:14:1956:U:C2	26:14:1957:C:C6	3.06	0.43
26:14:1027:A:H5'	27:1J:88:C:H41	1.84	0.43
26:1H:2830:G:N3	26:1H:2883:A:H2	2.16	0.43
26:1H:1830:C:O2'	26:1H:1831:G:H5'	2.17	0.43
26:1H:1975:G:H5''	58:1H:3984:HOH:O	2.17	0.43
46:D5:94:GLU:O	46:D5:130:PRO:HD3	2.18	0.43
38:55:94:TYR:C	38:55:117:VAL:HG23	2.39	0.43
37:88:103:MET:H	37:88:103:MET:HG2	1.40	0.43
37:45:110:THR:H	37:45:113:GLN:HE21	1.67	0.43
1:13:1342:C:H2'	1:13:1343:G:C8	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1E:111:ARG:NH1	2:1E:111:ARG:HG2	2.33	0.43
2:1E:71:VAL:HG12	2:1E:93:VAL:HB	2.00	0.43
1:1G:1050:G:C6	1:1G:1051:C:N4	2.87	0.43
29:29:35:GLN:HG2	29:29:37:ARG:HG2	2.01	0.43
9:8E:70:LYS:O	9:8E:74:ILE:HG13	2.19	0.43
9:8E:10:ARG:HD2	9:8E:11:LYS:HG3	1.99	0.43
5:4E:15:ARG:HB2	5:4E:28:PHE:CE2	2.54	0.43
26:1H:2843:G:O2'	26:1H:2844:G:H5'	2.18	0.43
31:49:81:LYS:HA	31:49:81:LYS:HD3	1.69	0.43
26:14:947:G:H2'	26:14:948:G:C8	2.54	0.43
26:14:664:C:H2'	26:14:665:C:C6	2.53	0.43
28:19:228:PRO:HG3	28:19:234:GLY:O	2.18	0.43
42:D8:98:GLU:OE2	42:D8:100:ARG:NH1	2.49	0.43
16:7I:13:HIS:C	16:7I:15:PRO:HD3	2.39	0.43
26:1H:2679:A:H61	26:1H:2728:U:H3	1.65	0.43
26:1H:828:U:H2'	26:1H:829:A:C8	2.54	0.43
1:1G:937:A:C5	1:1G:938:A:N7	2.87	0.43
1:1G:1009:G:C2	1:1G:1010:G:C8	3.06	0.43
35:68:2:ILE:HD12	35:68:6:THR:HG21	2.01	0.43
55:Q8:18:ALA:O	55:Q8:19:SER:OG	2.34	0.43
26:14:2352:A:N1	47:E5:33:ALA:O	2.52	0.43
26:14:1507:A:C2	26:14:1508:A:H1'	2.54	0.43
8:72:36:LEU:HD23	8:72:36:LEU:HA	1.86	0.43
6:5E:94:GLN:HG3	6:5E:94:GLN:H	1.55	0.43
26:14:931:G:H3'	26:14:931:G:C8	2.54	0.43
53:O8:25:LYS:HB3	53:O8:25:LYS:HE3	1.80	0.43
26:14:2820:A:C6	38:55:4:LEU:HD11	2.54	0.43
26:1H:1359:A:N1	26:1H:1372:U:C4	2.84	0.43
26:1H:2360:A:O2'	36:78:61:ARG:NH2	2.52	0.43
52:N8:3:LYS:HB2	52:N8:4:HIS:H	1.61	0.43
26:1H:818:G:H5'	26:1H:839:U:OP1	2.19	0.43
49:K8:48:HIS:H	49:K8:50:ILE:HD11	1.83	0.43
1:1G:861:G:C4	1:1G:862:C:C5	3.06	0.43
26:1H:1592:C:H2'	26:1H:1593:G:H8	1.84	0.43
46:D5:44:PHE:HE1	46:D5:48:PHE:CD2	2.37	0.43
1:1G:673:G:H2'	1:1G:674:G:H8	1.78	0.43
26:1H:2256:G:C5	58:1H:4119:HOH:O	2.69	0.43
1:13:658:G:H2'	1:13:659:U:C6	2.54	0.43
26:14:2318:G:N2	39:65:3:ARG:HB2	2.34	0.43
5:4E:63:ARG:O	5:4E:66:MET:HG3	2.19	0.43
4:3E:13:ARG:NH1	4:3E:38:TYR:O	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:75:8:LYS:NZ	40:75:8:LYS:HB2	2.34	0.43
26:14:878:A:N6	26:14:899:A:O2'	2.52	0.43
30:31:177:ALA:HB1	30:31:178:PRO:HD2	2.01	0.43
34:58:133:GLN:O	34:58:134:ARG:HB2	2.18	0.43
26:14:2065:C:H2'	26:14:2066:C:C6	2.54	0.43
18:9A:32:ARG:HB2	18:9A:65:ILE:HD11	2.01	0.43
26:1H:2031:A:HO2'	26:1H:2454:G:N2	2.17	0.43
26:1H:71:A:OP1	26:1H:72:U:H2'	2.19	0.43
44:F8:31:HIS:CD2	44:F8:33:LYS:HB2	2.54	0.43
26:1H:2352:A:N6	26:1H:2365:G:O2'	2.51	0.43
4:3E:167:GLY:CA	28:19:135:PHE:HE1	2.29	0.43
1:1G:963:G:HO2'	10:1A:54:PHE:HZ	1.66	0.43
1:1G:974:A:P	14:5A:41:ARG:HH12	2.42	0.43
1:1G:17:U:H1'	1:1G:1080:A:H1'	2.01	0.43
26:14:2150:U:H2'	26:14:2151:G:C8	2.49	0.43
26:14:2766:G:N3	26:14:2766:G:H2'	2.34	0.43
26:14:2789:C:H3'	26:14:2790:A:H5''	2.00	0.43
1:1G:582:U:H1'	1:1G:760:G:C6	2.54	0.43
2:1E:5:ILE:HG13	2:1E:6:THR:H	1.83	0.43
1:13:671:G:C4	1:13:672:U:C6	3.07	0.43
26:14:2861:G:O2'	26:14:2862:G:H5'	2.19	0.43
47:E5:24:LYS:O	47:E5:25:ARG:HD3	2.18	0.43
1:13:711:G:H2'	1:13:712:A:H8	1.84	0.43
24:3K:38:A:C2	24:3K:39:U:C4	3.07	0.43
26:14:1967:C:H2'	26:14:1968:G:O4'	2.18	0.43
1:13:668:G:O2'	15:6I:46:HIS:HB3	2.19	0.43
29:29:119:ARG:HA	29:29:160:TYR:CD2	2.54	0.43
1:13:983:A:H5'	1:13:984:C:OP2	2.19	0.43
1:13:4:U:C4	8:7E:105:ARG:HD2	2.54	0.43
1:13:790:A:C6	1:13:791:G:C6	3.07	0.43
4:32:101:LEU:HB2	4:32:138:TYR:HB3	2.00	0.43
26:1H:660:G:O3'	30:31:38:ARG:NH2	2.52	0.43
26:1H:1705:G:C2'	26:1H:1706:U:H5'	2.49	0.43
26:1H:1666:G:C2'	26:1H:1667:G:H5'	2.49	0.43
26:1H:2557:G:H2'	26:1H:2558:C:C6	2.54	0.43
11:2A:124:LYS:HB2	11:2A:125:PHE:CD1	2.53	0.43
43:A5:29:LEU:HD21	43:A5:33:ARG:NH2	2.34	0.43
13:4A:13:LYS:HA	13:4A:44:ARG:NH1	2.34	0.43
13:4A:7:VAL:HG11	31:49:115:ARG:NE	2.34	0.43
33:61:79:ILE:HA	33:61:80:PRO:HD2	1.93	0.43
1:1G:636:U:H2'	1:1G:637:G:H8	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:757:U:O2'	1:1G:879:C:O2	2.37	0.43
38:98:8:ARG:NH1	38:98:39:PRO:HB3	2.34	0.43
1:13:8:A:H62	4:3E:208:SER:HB2	1.84	0.43
2:12:186:ALA:O	2:12:201:ILE:HG12	2.19	0.43
38:98:78:LYS:O	38:98:83:ILE:HG13	2.19	0.43
26:14:654(E):C:H42	26:14:654(P):G:H1	1.64	0.43
26:1H:1545(A):A:H2'	26:1H:1546:C:O4'	2.19	0.43
29:21:23:VAL:HA	29:21:185:LYS:HA	2.00	0.43
26:1H:467:G:OP1	54:P8:33:ARG:NH1	2.48	0.43
28:11:248:SER:HB3	28:11:252:TRP:CZ3	2.54	0.43
4:3E:98:GLU:O	4:3E:103:ASN:ND2	2.52	0.43
32:51:116:GLU:HG3	32:51:117:PRO:HD2	2.01	0.43
44:B5:66:LEU:HA	44:B5:66:LEU:HD23	1.88	0.43
1:13:1225:A:N3	1:13:1225:A:H2'	2.34	0.43
1:1G:892:A:O2'	1:1G:1415:G:H4'	2.19	0.43
26:1H:2725:A:C4	26:1H:2727:G:C8	3.07	0.43
26:14:2107:C:H2'	26:14:2108:C:O4'	2.19	0.43
4:32:22:LYS:HG3	4:32:26:CYS:SG	2.59	0.43
26:1H:1869:G:N2	26:1H:1872:A:OP2	2.48	0.43
26:14:2600:A:N6	58:14:4084:HOH:O	2.43	0.42
26:14:751:A:C6	26:14:789:A:C5	3.07	0.42
28:11:105:ILE:HA	28:11:105:ILE:HD12	1.55	0.42
1:13:869:G:H4'	1:13:872:A:O4'	2.18	0.42
42:95:35:LEU:HB3	42:95:37:VAL:HG13	2.00	0.42
34:58:91:LEU:HA	34:58:95:PRO:HA	2.01	0.42
26:14:1925:C:C2'	26:14:1926:U:H5'	2.49	0.42
9:82:34:ASN:O	9:82:38:GLN:HB2	2.19	0.42
34:58:129:PRO:O	34:58:134:ARG:NH1	2.52	0.42
26:1H:1677:A:H2'	26:1H:1678:G:C8	2.54	0.42
1:1G:1305:G:O2'	1:1G:1306:A:H8	2.01	0.42
26:14:820:A:H2'	26:14:821:A:C8	2.54	0.42
35:25:47:ILE:HA	35:25:47:ILE:HD12	1.79	0.42
40:75:54:ARG:HB3	40:75:54:ARG:HH11	1.84	0.42
26:14:1777:U:O2'	26:14:1778:U:H5'	2.19	0.42
30:31:163:VAL:O	30:31:166:ALA:HB3	2.18	0.42
34:58:18:ALA:CA	34:58:21:LYS:HG3	2.49	0.42
2:12:6:THR:H	2:12:221:LEU:HD11	1.84	0.42
49:K8:14:ARG:HG2	49:K8:63:VAL:HG11	2.00	0.42
1:13:1258:G:H2'	1:13:1259:C:C6	2.54	0.42
40:75:53:ARG:O	40:75:53:ARG:HG3	2.19	0.42
26:1H:1282:U:H2'	26:1H:1283:G:C8	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:2617:C:O2'	26:1H:2618:G:H5'	2.19	0.42
26:14:710:G:C2	26:14:722:A:C2	3.06	0.42
53:K5:48:VAL:HG21	53:K5:50:ARG:NH2	2.34	0.42
33:69:110:ASP:HA	33:69:111:PRO:HD3	1.82	0.42
28:11:44:ASN:O	28:11:46:GLN:O	2.36	0.42
1:13:1492:A:H1'	25:4K:20:U:HO2'	1.84	0.42
32:51:97:ARG:O	32:51:103:LEU:HD12	2.18	0.42
30:39:110:LEU:O	30:39:114:VAL:HG23	2.19	0.42
51:M8:22:ILE:C	51:M8:24:THR:HG23	2.40	0.42
6:5E:7:ASN:HD21	6:5E:62:TRP:HD1	1.67	0.42
26:1H:188:G:H1	26:1H:208:C:H42	1.67	0.42
26:1H:1972:A:OP2	58:1H:4636:HOH:O	2.21	0.42
5:42:51:VAL:HG23	5:42:52:PRO:HD3	2.01	0.42
1:13:63:C:N4	1:13:105:G:O6	2.52	0.42
43:E8:29:LEU:HD21	43:E8:33:ARG:NH2	2.34	0.42
26:14:2579:C:H4'	29:29:134:ILE:HG12	2.01	0.42
29:29:134:ILE:HA	29:29:137:HIS:CD2	2.53	0.42
54:P8:10:ARG:HG2	54:P8:14:LYS:HD3	2.01	0.42
26:1H:341:G:H2'	26:1H:342:G:O4'	2.19	0.42
34:58:75:TYR:HA	34:58:81:GLY:O	2.19	0.42
41:85:106:PHE:O	41:85:109:LEU:HB2	2.19	0.42
46:D5:141:VAL:HG11	46:D5:150:LEU:HD22	2.01	0.42
1:13:1060:C:O2'	10:1I:56:HIS:ND1	2.52	0.42
33:69:7:GLU:HG3	33:69:8:PRO:HD2	2.01	0.42
37:88:118:LEU:HD23	37:88:118:LEU:HA	1.89	0.42
20:BI:29:LYS:HB2	20:BI:29:LYS:HE3	1.80	0.42
45:C5:40:GLU:CD	45:C5:40:GLU:N	2.72	0.42
31:49:111:LEU:HD23	31:49:111:LEU:HA	1.83	0.42
48:J8:20:ARG:HH11	48:J8:20:ARG:HG2	1.84	0.42
26:1H:2266:A:H4'	26:1H:2267:A:N3	2.34	0.42
1:1G:1521:G:H2'	1:1G:1522:U:H6	1.84	0.42
32:51:144:VAL:O	32:51:148:ILE:HG12	2.19	0.42
30:39:157:VAL:HB	30:39:194:MET:HB3	1.99	0.42
26:14:273(F):C:H3'	26:14:274:G:H5''	2.02	0.42
26:1H:1188:U:O2'	26:1H:1189:A:H5'	2.19	0.42
26:1H:2017:U:O2	52:N8:10:LYS:HB2	2.19	0.42
26:14:890:A:H2'	26:14:892:G:C8	2.54	0.42
26:1H:1777:U:O2'	26:1H:1778:U:H5'	2.19	0.42
9:8E:114:TYR:HB2	10:1I:60:ARG:HG3	2.02	0.42
1:1G:1357:A:C5	1:1G:1358:U:C4	3.07	0.42
26:14:1022:G:C5	26:14:1140:C:C4	3.07	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:250:G:H5'	36:78:60:MET:HE1	2.01	0.42
26:1H:1163:G:N3	26:1H:1164:G:C8	2.88	0.42
26:1H:1423:G:C6	26:1H:1424:G:N7	2.87	0.42
29:21:117:MET:HE3	29:21:117:MET:HB3	1.81	0.42
1:13:827:U:C5	1:13:872:A:N1	2.87	0.42
1:1G:1148:U:H2'	1:1G:1149:C:O4'	2.19	0.42
43:E8:12:ILE:HG13	43:E8:42:ARG:HH11	1.83	0.42
1:13:694:A:H5''	11:2I:53:SER:OG	2.20	0.42
38:98:29:LEU:HA	38:98:29:LEU:HD12	1.69	0.42
41:C8:90:VAL:HG12	41:C8:91:ASP:HA	2.01	0.42
26:14:1839:G:N3	26:14:1839:G:H2'	2.34	0.42
1:1G:1326:C:H2'	1:1G:1327:C:H6	1.83	0.42
1:13:1350:A:C6	1:13:1351:U:N3	2.87	0.42
20:BA:64:ASP:OD1	20:BA:64:ASP:N	2.52	0.42
26:1H:2287:A:C2	26:1H:2346:A:H2	2.37	0.42
1:1G:502:G:C2	1:1G:503:C:C2	3.07	0.42
20:BA:26:ASN:HB2	20:BA:71:THR:HG23	2.00	0.42
1:1G:1080:A:H5''	1:1G:1081:G:OP2	2.19	0.42
28:11:14:ARG:HD3	28:11:15:PHE:CZ	2.53	0.42
27:1J:7:G:N2	39:65:38:GLN:OE1	2.34	0.42
26:14:1668:A:OP1	35:25:5:GLN:HG2	2.20	0.42
26:1H:2138:C:N3	26:1H:2153:G:N2	2.55	0.42
47:E5:66:VAL:CG1	47:E5:67:VAL:N	2.82	0.42
26:1H:1473:G:H2'	26:1H:1474:C:O4'	2.20	0.42
34:58:22:THR:HG22	34:58:23:LEU:H	1.84	0.42
26:14:55:G:H2'	26:14:56:A:C8	2.55	0.42
26:14:55:G:H2'	26:14:56:A:H8	1.83	0.42
1:1G:247:G:OP2	17:8A:100:LYS:HE2	2.19	0.42
1:13:668:G:C6	1:13:669:U:C5	3.07	0.42
15:6I:70:LEU:HD11	15:6I:77:ARG:HG2	2.01	0.42
1:13:186(F):C:H2'	1:13:187:C:O4'	2.19	0.42
26:1H:1550:C:H2'	26:1H:1551:C:C6	2.54	0.42
27:1J:83:G:H1	27:1J:93:C:H42	1.67	0.42
1:1G:1097:C:H2'	1:1G:1098:C:H6	1.83	0.42
44:B5:67:GLY:C	44:B5:69:TYR:H	2.20	0.42
20:BA:14:LYS:HA	20:BA:17:ARG:CZ	2.49	0.42
17:8I:34:LYS:HD3	17:8I:36:ILE:HG22	2.00	0.42
32:51:101:ARG:NH1	32:51:122:THR:OG1	2.48	0.42
1:1G:812:C:H1'	1:1G:813:U:OP2	2.19	0.42
26:1H:784:A:H5''	28:11:227:ASN:HD21	1.84	0.42
26:14:1794:U:H2'	26:14:1795:C:H6	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:26:G:C6	26:14:27:G:N1	2.87	0.42
1:13:591:U:H2'	1:13:592:G:H8	1.84	0.42
1:1G:1295:G:O2'	13:4A:14:ARG:NH1	2.52	0.42
44:F8:5:TYR:HB3	49:K8:33:MET:HB2	2.01	0.42
26:1H:109:G:H2'	26:1H:110:G:O4'	2.18	0.42
37:45:43:THR:O	37:45:46:GLN:N	2.52	0.42
26:1H:1945:G:H2'	26:1H:1946:U:H6	1.83	0.42
18:9I:38:GLU:HA	18:9I:41:LYS:HE3	2.01	0.42
1:1G:1459:C:OP1	20:BA:31:SER:OG	2.35	0.42
26:14:80:G:C6	26:14:81:G:C5	3.07	0.42
31:49:19:LEU:HD13	31:49:23:PHE:HE2	1.84	0.42
50:H5:7:LYS:HG3	50:H5:34:GLU:CG	2.49	0.42
1:13:931:C:O2'	1:13:932:C:H5'	2.19	0.42
18:9A:74:ARG:HB3	18:9A:81:PHE:CE1	2.54	0.42
9:8E:45:ALA:O	9:8E:48:GLU:HG2	2.19	0.42
52:N8:42:PRO:HB2	52:N8:43:HIS:ND1	2.34	0.42
2:1E:75:LYS:HA	2:1E:78:GLN:HB2	2.00	0.42
20:BA:37:SER:O	20:BA:41:ILE:HG12	2.20	0.42
1:1G:6:G:O2'	1:1G:7:G:O5'	2.37	0.42
1:1G:668:G:O2'	15:6A:46:HIS:HB3	2.18	0.42
40:75:61:PHE:CE1	40:75:76:PHE:HB2	2.54	0.42
31:49:76:SER:OG	31:49:84:LYS:N	2.51	0.42
26:1H:1486:A:H2'	26:1H:1487:G:C8	2.55	0.42
26:1H:1486:A:H2'	26:1H:1487:G:H8	1.84	0.42
1:1G:304:U:H2'	1:1G:305:G:C8	2.54	0.42
5:4E:53:LEU:O	5:4E:56:GLN:HB2	2.18	0.42
3:2E:95:THR:HB	3:2E:97:LYS:H	1.84	0.42
2:12:24:TRP:NE1	2:12:26:PRO:HG3	2.34	0.42
41:C8:110:VAL:O	41:C8:114:LYS:N	2.38	0.42
36:78:121:LYS:HE2	36:78:121:LYS:HB3	1.82	0.42
28:19:64:ILE:O	28:19:64:ILE:HG12	2.18	0.42
26:14:1280:G:H5''	26:14:1280:G:C8	2.53	0.42
26:14:221:A:C4	26:14:266:G:N7	2.87	0.42
4:32:25:ARG:HG2	4:32:30:LYS:O	2.19	0.42
55:Q8:52:LYS:HA	55:Q8:54:GLU:HB2	2.02	0.42
26:14:1043:C:H2'	26:14:1044:G:H5'	2.01	0.42
1:13:1145:C:H5''	1:13:1146:A:OP1	2.20	0.42
42:95:72:VAL:HG13	42:95:72:VAL:O	2.19	0.42
42:95:85:LYS:CG	42:95:86:GLY:H	2.31	0.42
1:13:21:G:H2'	1:13:22:G:C8	2.54	0.42
49:K8:50:ILE:HD12	49:K8:51:ARG:N	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1110:A:OP2	58:13:1958:HOH:O	2.22	0.42
26:14:259:G:N2	26:14:621:A:H8	2.05	0.42
26:1H:674:G:O2'	30:31:74:ARG:HD3	2.19	0.42
22:1L:76:A:N6	26:14:2583:G:H21	2.11	0.42
24:3L:58:A:H4'	24:3L:59:U:OP1	2.20	0.42
1:1G:1187:G:H4'	9:82:111:ARG:HH11	1.83	0.42
26:14:2706:G:C2	26:14:2707:G:H1'	2.54	0.42
48:J8:73:LEU:HD11	48:J8:95:LEU:HD21	2.01	0.42
1:1G:540:G:H2'	1:1G:541:G:O4'	2.19	0.42
26:1H:139:G:N3	26:1H:141:A:N1	2.67	0.42
49:K8:64:LEU:HD22	49:K8:68:ARG:HD2	1.99	0.42
1:13:560:U:H6	1:13:560:U:O5'	2.01	0.42
37:88:56:ARG:HD2	37:88:56:ARG:HA	1.90	0.42
26:1H:835:A:N6	26:1H:836:G:C6	2.88	0.42
36:35:138:LEU:HD22	36:35:138:LEU:HA	1.90	0.42
35:25:78:ARG:NH2	40:75:73:GLU:OE2	2.53	0.42
26:1H:847:U:C5	26:1H:933:A:N1	2.87	0.42
23:2L:5:G:N1	23:2L:69:C:O2	2.52	0.42
26:1H:583:G:H5''	41:C8:10:ARG:HH12	1.85	0.42
23:2K:8:4SU:O2	23:2K:22:A:H2	2.01	0.42
23:2K:56:PSU:O4	23:2K:58:A:C8	2.72	0.42
54:L5:19:ARG:HG2	54:L5:19:ARG:HH11	1.84	0.42
16:7I:4:ILE:HD12	16:7I:66:PRO:HB3	2.02	0.42
26:14:1359:A:N7	26:14:1372:U:O4	2.52	0.42
16:7A:22:THR:CA	16:7A:33:ILE:HG12	2.48	0.42
30:31:63:LYS:HE3	30:31:65:TRP:O	2.18	0.42
30:39:63:LYS:HE3	30:39:75:HIS:O	2.18	0.42
26:1H:1472:A:H2'	26:1H:1473:G:H8	1.84	0.42
26:1H:301:G:C4	26:1H:302:C:C5	3.08	0.42
31:49:108:ASN:O	51:I5:37:SER:HA	2.19	0.42
1:13:347:G:OP2	1:13:347:G:H8	2.03	0.42
4:3E:120:LEU:HD23	4:3E:120:LEU:HA	1.86	0.42
1:13:1309:G:C6	1:13:1329:A:C2	3.07	0.42
53:K5:35:GLU:HB2	53:K5:51:GLU:HB2	2.01	0.42
11:2I:29:ILE:HG13	11:2I:44:SER:HB3	2.01	0.42
47:I8:75:LEU:HA	47:I8:75:LEU:HD23	1.56	0.42
26:1H:1931:U:C5	26:1H:1969:A:N7	2.87	0.42
1:1G:838:G:N2	1:1G:842:C:H1'	2.34	0.42
10:1A:24:VAL:HG21	10:1A:37:PRO:HG3	2.00	0.42
26:1H:671:C:H42	26:1H:809:G:H1	1.65	0.42
16:7I:72:ARG:O	16:7I:75:ARG:HB3	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1426:G:H2'	26:1H:1427:A:C8	2.54	0.42
26:14:2348:U:OP2	55:M5:42:ARG:NH2	2.52	0.42
38:55:58:GLY:HA2	38:55:80:PHE:HE2	1.84	0.42
8:72:92:ARG:HB3	8:72:94:TYR:CE2	2.53	0.42
22:1K:74:C:O2'	22:1K:75:C:O5'	2.37	0.42
33:61:93:THR:OG1	33:61:96:ASP:OD1	2.21	0.42
26:14:1832:C:N4	26:14:1833:U:C4	2.87	0.42
44:B5:26:TYR:O	44:B5:81:VAL:HG22	2.18	0.42
8:72:10:LEU:HD22	8:72:83:ILE:HD11	2.00	0.42
1:13:746:A:C2'	1:13:747:C:H5'	2.49	0.42
26:14:1743:G:C2	26:14:1746:G:C8	3.07	0.42
2:1E:60:ASP:HB3	2:1E:64:ARG:NH1	2.35	0.42
31:49:53:LEU:HD12	31:49:54:GLU:N	2.34	0.42
5:4E:41:VAL:HG13	5:4E:113:ALA:HB2	2.00	0.42
26:1H:665:C:H2'	26:1H:666:G:C8	2.54	0.42
45:G8:79:CYS:HB2	45:G8:80:GLY:H	1.51	0.42
42:95:48:GLY:HA3	42:95:51:VAL:C	2.40	0.42
26:14:1592:C:O5'	26:14:1592:C:H6	2.02	0.42
4:3E:70:ILE:HA	4:3E:70:ILE:HD12	1.82	0.42
1:13:1090:U:H2'	1:13:1091:U:C6	2.54	0.42
26:1H:219:G:H2'	26:1H:220:G:C8	2.54	0.42
26:1H:2072:G:H2'	26:1H:2073:C:O4'	2.19	0.42
1:13:1145:C:H4'	1:13:1146:A:H8	1.84	0.42
26:1H:1142(A):A:C2	26:1H:1144:G:C8	3.08	0.42
37:88:41:TRP:HB3	37:88:94:VAL:HG11	2.02	0.42
26:14:2431:U:O2	26:14:2433:A:C8	2.72	0.42
5:42:43:LEU:HB3	5:42:136:MET:SD	2.60	0.42
3:22:3:ASN:H	3:22:3:ASN:HD22	1.67	0.42
26:14:2275:C:C5'	26:14:2275:C:C6	3.01	0.42
26:1H:1798:U:H5'	28:11:259:THR:CG2	2.48	0.42
26:1H:2040:C:H2'	26:1H:2041:U:O4'	2.18	0.42
44:F8:2:LYS:HG2	49:K8:26:ARG:NE	2.30	0.42
36:35:122:PRO:HB3	36:35:141:ALA:HB1	2.02	0.42
1:13:1434:A:H2'	1:13:1435:G:O4'	2.19	0.42
16:7I:57:ARG:NH2	16:7I:79:VAL:O	2.48	0.42
30:39:89:VAL:O	30:39:90:PHE:C	2.58	0.42
31:41:130:ASN:HB3	31:41:159:VAL:O	2.18	0.42
28:11:182:LEU:H	28:11:272:ALA:CB	2.30	0.42
26:14:2141:G:C6	26:14:2151:G:C6	3.07	0.42
1:13:648:A:C6	1:13:649:G:C6	3.07	0.42
38:98:48:VAL:O	38:98:51:LEU:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:315:G:C6	26:14:316:C:C4	3.07	0.42
26:1H:410:G:H2'	58:1H:4631:HOH:O	2.19	0.42
19:AI:5:LEU:HD13	19:AI:10:PHE:HD1	1.84	0.42
23:2K:20:G:HO2'	23:2K:21:U:H3	1.67	0.42
1:13:1409:C:H2'	1:13:1410:G:H8	1.83	0.42
1:1G:746:A:H2'	1:1G:747:C:C6	2.55	0.42
15:6A:39:LEU:CD1	15:6A:56:LEU:HB2	2.49	0.42
26:1H:1543:A:C8	26:1H:1545:A:H5''	2.54	0.42
1:1G:1512:U:N3	1:1G:1513:A:N7	2.66	0.42
12:3I:59:ARG:HA	12:3I:65:GLU:HA	2.00	0.42
26:14:1163:G:H2'	26:14:1164:G:H8	1.83	0.42
47:E5:72:ARG:CB	47:E5:75:LEU:HB2	2.49	0.42
11:2I:59:TYR:CZ	11:2I:63:LEU:HD11	2.54	0.42
26:14:1217:C:OP2	41:85:15:LYS:HE3	2.20	0.42
34:58:73:THR:HG22	34:58:84:LYS:HG2	2.02	0.42
28:19:120:GLY:HA2	28:19:190:TYR:OH	2.19	0.42
1:1G:476:G:C4	1:1G:477:G:C8	3.07	0.42
26:1H:2746:U:O4	26:1H:2755:C:H4'	2.19	0.42
17:8A:59:ILE:HG22	17:8A:71:PHE:CD2	2.54	0.42
1:1G:135:C:O2	16:7A:1:MET:HB3	2.18	0.42
1:13:8:A:N7	4:3E:208:SER:HB3	2.34	0.42
1:1G:29:G:H5'	1:1G:296:U:OP1	2.20	0.42
23:2L:7:G:H3'	23:2L:8:4SU:C5'	2.48	0.42
8:7E:100:ILE:HA	8:7E:101:PRO:HD3	1.85	0.42
8:7E:121:ASP:HB2	8:7E:125:ARG:NH2	2.34	0.42
26:14:2057:A:H2'	26:14:2058:A:C8	2.54	0.42
26:14:866:A:H5''	26:14:867:C:OP2	2.19	0.42
26:14:999:U:C5	26:14:1154:G:C5	3.07	0.42
32:59:70:THR:O	32:59:74:ASN:N	2.52	0.42
22:1L:38:A:H2'	22:1L:39:PSU:H6	1.84	0.42
8:72:136:GLU:O	8:72:136:GLU:HG3	2.18	0.42
48:F5:11:ARG:HB3	48:F5:11:ARG:CZ	2.49	0.42
1:13:49:U:O2	1:13:362:G:H1'	2.19	0.42
1:13:551:U:H2'	1:13:552:U:C6	2.54	0.42
7:6E:108:ALA:HB2	7:6E:123:GLU:HG2	2.01	0.42
4:32:30:LYS:CB	4:32:35:ARG:HD2	2.48	0.42
39:65:24:LEU:HB2	39:65:85:VAL:HG12	2.01	0.42
26:14:675:A:C6	26:14:676:A:C6	3.07	0.42
45:C5:73:ARG:HB3	45:C5:73:ARG:HE	1.48	0.42
1:13:1500:A:P	58:13:1804:HOH:O	2.73	0.42
10:1I:62:HIS:N	10:1I:62:HIS:CD2	2.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:1689:A:C6	26:1H:1700:A:C2	3.07	0.42
1:13:605:U:C2'	1:13:606:G:H5'	2.49	0.42
1:1G:474:G:C2	1:1G:475:G:C5	3.08	0.42
30:31:78:ILE:HA	30:31:83:PHE:CD2	2.54	0.42
8:7E:9:MET:O	8:7E:12:ARG:N	2.50	0.42
51:I5:21:VAL:HG22	51:I5:22:ILE:H	1.84	0.42
26:14:2344:U:OP1	53:K5:38:LYS:HD3	2.19	0.42
26:14:2305:A:C6	31:49:154:GLY:N	2.88	0.42
29:29:181:LEU:HD11	40:75:7:ILE:HD11	2.01	0.42
55:Q8:34:TRP:HB3	55:Q8:35:GLN:CA	2.48	0.42
46:D5:69:THR:HB	46:D5:88:PHE:HB2	2.01	0.42
26:1H:825:C:O2	36:78:55:ARG:NH2	2.51	0.42
16:7A:9:PHE:HB2	16:7A:16:HIS:O	2.20	0.42
1:13:1368:G:OP2	9:8E:112:LYS:HE2	2.19	0.42
26:14:2388:A:O2'	26:14:2389:G:H5'	2.20	0.42
37:88:30:GLY:O	37:88:134:ARG:HD2	2.19	0.42
26:14:69:C:H2'	26:14:70:G:C8	2.54	0.42
26:1H:138:G:H2'	26:1H:139:G:C8	2.54	0.42
26:1H:1204:A:C2	26:1H:1241:A:N1	2.86	0.42
12:3I:53:ARG:HG3	12:3I:53:ARG:HH11	1.84	0.42
1:13:1238:A:N3	1:13:1241:G:O2'	2.49	0.42
20:BA:26:ASN:CB	20:BA:71:THR:HG23	2.50	0.42
1:13:286:G:C6	1:13:287:U:N3	2.87	0.42
26:1H:2074:U:H2'	26:1H:2075:U:C6	2.54	0.42
1:13:142:G:O2'	1:13:196:A:N1	2.48	0.42
26:1H:644:A:C2	26:1H:646:A:C4	3.08	0.42
26:1H:646:A:C8	26:1H:647:G:H1'	2.55	0.42
23:2L:20:G:H1	23:2L:57:C:N4	2.17	0.42
1:1G:1053:G:O2'	1:1G:1054:C:OP2	2.35	0.42
1:13:1316:G:N1	1:13:1319:A:OP2	2.50	0.42
19:AI:10:PHE:N	19:AI:10:PHE:CD1	2.87	0.42
26:1H:1265:A:OP2	58:1H:3616:HOH:O	2.20	0.42
13:4A:49:THR:N	13:4A:52:GLU:OE1	2.25	0.42
4:3E:135:LEU:HA	4:3E:136:PRO:HD2	1.80	0.42
37:45:36:ALA:HB2	37:45:103:MET:SD	2.59	0.42
27:1J:89:G:C8	27:1J:89:G:OP2	2.73	0.42
55:M5:11:LYS:HA	55:M5:59:LYS:HZ2	1.84	0.42
1:1G:735:C:O2'	1:1G:736:C:H5'	2.20	0.42
29:21:67:PHE:O	29:21:69:LYS:HE2	2.19	0.42
1:1G:778:G:H2'	1:1G:779:C:O4'	2.19	0.42
5:4E:35:GLY:HA3	5:4E:112:LEU:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:6I:71:GLN:HG3	15:6I:78:TYR:CD2	2.55	0.42
26:1H:301:G:C6	26:1H:317:G:C5	3.08	0.42
26:14:30:G:C5	26:14:31:C:C4	3.08	0.42
31:41:81:LYS:HA	31:41:81:LYS:HD3	1.84	0.42
32:51:101:ARG:HE	32:51:101:ARG:HB3	1.60	0.42
1:1G:1151:A:OP1	10:1A:42:THR:N	2.44	0.42
1:1G:1418:A:H5''	1:1G:1419:G:OP2	2.19	0.42
7:6E:48:LYS:HE2	7:6E:48:LYS:HB3	1.80	0.42
26:14:2185:C:H2'	26:14:2186:G:H8	1.84	0.42
5:42:69:VAL:O	5:42:71:LEU:N	2.50	0.42
29:21:182:LEU:HD12	29:21:183:LEU:H	1.84	0.42
35:68:104:ARG:NH1	40:B8:36:GLU:OE2	2.53	0.42
26:1H:236:C:H2'	26:1H:237:C:C6	2.55	0.42
26:1H:1161:C:O2'	42:D8:8:GLY:HA2	2.19	0.42
29:29:97:LYS:O	29:29:100:GLU:HG3	2.19	0.42
48:F5:80:LEU:HD23	48:F5:82:LEU:HD21	2.01	0.42
7:6E:62:PHE:HA	7:6E:124:LEU:HD21	2.00	0.42
37:45:2:LEU:O	37:45:70:PRO:HG2	2.20	0.42
26:1H:26:G:C6	26:1H:27:G:N1	2.87	0.42
31:41:128:ARG:NH2	31:41:128:ARG:HB2	2.35	0.42
26:14:513:A:N3	26:14:514:A:C8	2.88	0.42
26:14:2352:A:C2	47:E5:33:ALA:HB1	2.54	0.42
26:1H:220:G:O2'	26:1H:233:A:N3	2.41	0.42
8:72:51:VAL:HG11	8:72:60:ARG:HH21	1.84	0.42
26:1H:1711:C:H2'	26:1H:1712:C:C6	2.54	0.42
26:1H:773:U:C4'	28:11:47:GLY:HA3	2.49	0.42
34:15:34:LEU:O	34:15:49:GLY:HA3	2.19	0.42
37:88:1:MET:O	37:88:2:LEU:HB2	2.20	0.42
20:BI:59:ALA:HA	20:BI:62:LEU:HD12	2.00	0.42
26:14:1854:A:H2'	26:14:1855:G:O4'	2.20	0.42
26:14:1023:U:H4'	26:14:1123:C:OP1	2.19	0.42
31:41:111:LEU:HD23	31:41:114:ILE:HD12	2.01	0.42
1:13:1234:C:H2'	1:13:1235:U:C6	2.54	0.42
26:14:1788:C:H2'	26:14:1789:A:H8	1.83	0.42
26:14:325:G:H2'	26:14:326:G:O4'	2.19	0.42
41:85:91:ASP:CG	41:85:96:ALA:HB2	2.40	0.42
26:1H:1136:G:N3	26:1H:1136:G:H2'	2.34	0.42
1:1G:149:A:H2'	1:1G:150:C:C6	2.54	0.42
26:14:2654:A:OP1	26:14:2654:A:H8	2.03	0.42
37:45:31:ASP:H	37:45:107:ALA:HB2	1.84	0.42
26:1H:2259:G:C2	26:1H:2282:G:N1	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:69:88:ILE:HG13	33:69:121:LYS:C	2.39	0.42
26:1H:1986:A:C2	26:1H:1987:G:C5	3.08	0.42
15:6I:18:PHE:CZ	15:6I:21:ASP:HB2	2.54	0.42
45:C5:76:CYS:HB2	45:C5:97:ARG:HE	1.84	0.42
1:13:635:G:C6	1:13:636:U:C4	3.08	0.42
1:13:1124:G:C2	1:13:1127:G:N2	2.87	0.42
26:1H:962:G:H4'	26:1H:2496:C:O2'	2.20	0.42
26:1H:960:A:H61	37:88:82:ARG:HH21	1.67	0.42
26:14:531:C:OP1	26:14:561:G:N1	2.52	0.42
1:1G:1320:C:C1'	19:AA:73:GLU:HG2	2.48	0.42
26:14:1035:U:H2'	26:14:1036:G:C8	2.54	0.42
46:D5:111:VAL:HG22	46:D5:112:ARG:HH11	1.85	0.42
30:31:108:LYS:HE2	30:31:108:LYS:HB3	1.76	0.42
38:98:10:LEU:O	38:98:11:ASN:C	2.58	0.42
36:35:63:PRO:HB3	55:M5:13:ARG:HG2	2.02	0.42
26:14:634:C:H2'	26:14:635:C:C6	2.54	0.42
1:1G:683:G:N2	1:1G:707:C:O2	2.44	0.42
26:14:2420:C:OP1	55:M5:34:TRP:HB3	2.18	0.42
1:13:310:G:OP2	16:7I:27:LYS:NZ	2.28	0.42
36:35:138:LEU:HD12	36:35:144:GLU:HG3	2.00	0.42
1:13:375:U:C4	1:13:376:G:N7	2.88	0.42
1:1G:430:A:C4	1:1G:431:A:C8	3.07	0.42
36:78:106:LEU:O	36:78:106:LEU:HD22	2.19	0.42
3:2E:19:GLU:HB3	3:2E:40:ARG:NH2	2.35	0.42
2:1E:11:LEU:O	2:1E:16:HIS:NE2	2.53	0.42
2:1E:7:VAL:HG21	2:1E:11:LEU:HD12	2.02	0.42
7:62:65:ALA:HB3	7:62:124:LEU:HD22	2.01	0.42
1:1G:434:U:H2'	1:1G:435:C:H6	1.80	0.42
1:1G:728:A:C2	1:1G:729:A:C5	3.08	0.42
26:14:2784:C:H2'	26:14:2785:C:H6	1.83	0.42
27:1J:23:G:C2	27:1J:24:G:O6	2.72	0.42
26:14:289:A:H3'	26:14:290:G:C8	2.52	0.42
26:1H:1471:A:C4	26:1H:1472:A:C8	3.08	0.42
38:55:94:TYR:O	38:55:117:VAL:HG23	2.20	0.42
26:14:118:A:OP2	26:14:119:A:H5''	2.20	0.42
26:14:51:G:N3	26:14:119:A:C2	2.88	0.42
26:14:873:G:H1'	37:45:29:PHE:HE2	1.85	0.42
28:11:81:ALA:HB3	28:11:94:LEU:HB3	2.02	0.42
1:1G:754:C:H3'	1:1G:755:G:H5'	2.01	0.42
26:14:902:C:H2'	26:14:903:C:C6	2.55	0.42
22:1K:37:MIA:H112	22:1K:38:A:N3	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:864:G:O5'	26:14:864:G:H8	2.03	0.42
26:1H:2766:G:H5''	26:1H:2767:C:OP2	2.19	0.42
7:62:18:TYR:CD2	7:62:59:LEU:HD12	2.54	0.42
32:51:86:GLU:HG2	32:51:87:LEU:H	1.84	0.42
31:41:12:TYR:HD1	31:41:16:ARG:HD3	1.83	0.42
26:14:2854:G:N2	26:14:2864:G:C4	2.87	0.42
32:51:91:GLY:O	32:51:94:TYR:HB2	2.19	0.42
26:1H:1297:C:OP1	26:1H:2710:C:H4'	2.19	0.42
26:14:300:A:H1'	26:14:319:C:H1'	2.01	0.42
29:29:16:ARG:HB2	29:29:21:VAL:HG21	2.01	0.42
26:1H:1651:G:H5'	38:98:39:PRO:HG2	2.00	0.42
40:75:33:LYS:HD3	40:75:82:LEU:O	2.19	0.42
3:22:111:LEU:HD21	3:22:146:ALA:HB2	2.00	0.42
34:58:57:ALA:C	34:58:59:LYS:H	2.23	0.42
6:52:22:GLU:O	6:52:26:ILE:HG13	2.19	0.42
47:I8:37:LEU:HD11	47:I8:61:ALA:N	2.35	0.42
26:14:729:G:H2'	26:14:1775:U:O2	2.18	0.42
26:1H:53:A:H2'	26:1H:54:G:O4'	2.20	0.42
26:1H:2878:U:O4	58:1H:4250:HOH:O	2.21	0.42
1:13:134:A:H1'	1:13:325:A:C5	2.55	0.42
26:14:271(B):G:N7	26:14:421:U:H2'	2.35	0.42
25:4L:15:A:O5'	25:4L:15:A:H8	2.02	0.42
26:1H:2641:G:H5''	26:1H:2641:G:H8	1.85	0.42
26:1H:1194:A:OP2	26:1H:1194:A:H8	2.03	0.42
20:BI:68:LYS:HB2	20:BI:68:LYS:HE3	1.67	0.42
1:13:46:G:H2'	1:13:366:C:C5	2.54	0.42
26:1H:425:G:C2	26:1H:426:C:C6	3.07	0.42
36:35:131:SER:HB3	36:35:134:ALA:CB	2.50	0.42
1:13:1144:G:H21	1:13:1146:A:H62	1.66	0.42
1:13:254:G:O3'	17:8I:69:LYS:NZ	2.43	0.42
1:1G:21:G:OP2	58:1G:1834:HOH:O	2.22	0.42
28:11:85:ASP:OD1	28:11:86:PRO:HD2	2.20	0.42
1:13:963:G:N2	10:1I:55:LYS:HZ1	2.17	0.42
26:14:945:A:C4	26:14:2448:A:C2	3.07	0.42
24:3L:51:U:C2	24:3L:52:G:C8	3.08	0.42
1:13:1240:U:C5	7:6E:32:ARG:HD2	2.53	0.42
1:13:872:A:C8	1:13:874:G:C8	3.08	0.42
29:21:78:LEU:O	29:21:79:ARG:HB2	2.19	0.42
13:4A:73:GLU:O	13:4A:77:ASN:HB2	2.18	0.42
53:K5:28:ARG:HH11	53:K5:28:ARG:HA	1.84	0.42
26:14:2782:G:N7	58:14:3788:HOH:O	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:3E:30:LYS:HB2	4:3E:35:ARG:HE	1.83	0.42
55:Q8:39:LYS:CD	55:Q8:40:GLU:H	2.32	0.42
1:13:1288:A:O3'	21:1F:10:ARG:NH2	2.53	0.42
46:D5:71:VAL:HB	46:D5:88:PHE:CD2	2.55	0.42
26:1H:2693:A:C4	26:1H:2694:G:C8	3.07	0.42
31:41:112:PRO:HB3	51:M8:37:SER:N	2.28	0.42
26:14:16:G:O2'	26:14:17:G:H5'	2.18	0.42
1:1G:518:C:H5''	1:1G:519:C:H6	1.85	0.42
27:1J:117:G:H8	27:1J:117:G:O5'	2.01	0.42
26:1H:2364:C:H2'	26:1H:2365:G:O4'	2.19	0.42
26:1H:2238:G:H2'	26:1H:2238:G:N3	2.34	0.42
1:13:353:A:C8	1:13:353:A:H5'	2.47	0.42
45:G8:84:ARG:O	45:G8:84:ARG:NE	2.53	0.42
26:1H:2336:A:H61	47:I8:43:THR:HB	1.84	0.42
30:31:34:TRP:CH2	36:78:8:PRO:HB3	2.55	0.42
26:1H:129:C:H2'	26:1H:130:C:C6	2.55	0.42
14:5I:6:LEU:HA	14:5I:6:LEU:HD23	1.91	0.42
26:14:1483:G:C2	26:14:1484:G:C8	3.08	0.42
1:13:515:G:N2	1:13:537:G:C4	2.88	0.42
26:1H:1470:G:H5''	26:1H:1471:A:OP1	2.20	0.42
31:49:88:ILE:HD12	31:49:88:ILE:HA	1.89	0.42
26:14:1268:A:C2	26:14:2013:A:C4	3.08	0.42
2:1E:15:VAL:HB	2:1E:210:SER:HB2	2.01	0.42
28:19:118:VAL:HG22	28:19:119:ALA:N	2.33	0.42
1:1G:608:A:H2'	1:1G:609:A:O4'	2.20	0.42
27:1J:93:C:H2'	27:1J:94:C:H6	1.84	0.42
26:14:1491:G:O2'	28:19:101:GLU:HB2	2.19	0.42
26:14:1011:G:N3	26:14:1151:G:N2	2.67	0.42
19:AI:13:ASP:HA	19:AI:16:LEU:HB3	2.02	0.42
12:3I:24:VAL:HB	12:3I:27:LEU:HD12	2.02	0.42
34:15:96:GLU:HB2	34:15:122:VAL:HG12	2.00	0.42
1:13:57:G:C5	1:13:58:C:C4	3.07	0.42
38:98:45:ARG:HD2	38:98:97:VAL:CG2	2.48	0.42
1:1G:758:G:H4'	1:1G:880:C:H4'	2.02	0.42
27:16:95:U:H2'	27:16:96:G:C8	2.55	0.42
46:D5:52:SER:O	46:D5:54:HIS:N	2.52	0.42
42:95:48:GLY:HA3	42:95:52:VAL:N	2.35	0.42
1:1G:130:A:H1'	1:1G:263:A:O2'	2.19	0.42
43:E8:110:LYS:HG3	43:E8:111:HIS:H	1.84	0.42
26:1H:2197:U:H1'	26:1H:2198:A:C8	2.55	0.42
26:14:1465:G:N2	26:14:1466:G:H1'	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:19:136:ILE:HA	28:19:137:PRO:HD3	1.88	0.42
26:14:1769:G:C5	26:14:1984:G:C6	3.08	0.42
15:6I:61:GLY:O	15:6I:64:ARG:HB3	2.19	0.42
26:14:225:A:N6	26:14:226:G:C2	2.88	0.42
26:14:2632:A:O2'	26:14:2811:G:O2'	2.12	0.42
26:1H:1051:G:OP2	26:1H:1051:G:H8	2.02	0.42
44:F8:41:ASN:OD1	44:F8:41:ASN:N	2.52	0.42
1:13:282:A:H2'	1:13:282:A:N3	2.34	0.42
26:1H:364:C:H6	26:1H:364:C:H5'	1.84	0.42
18:9A:29:PHE:CD1	18:9A:29:PHE:N	2.87	0.42
26:14:1067:A:H2'	26:14:1067:A:N3	2.34	0.42
24:3L:25:C:C4	24:3L:26:A:C8	3.07	0.42
10:1A:78:ASN:OD1	10:1A:80:LYS:HB3	2.20	0.42
1:13:549:C:C2	1:13:550:G:C8	3.08	0.42
1:13:636:U:H2'	1:13:637:G:C8	2.55	0.42
46:H8:61:LEU:CD2	46:H8:67:LEU:HD12	2.49	0.42
26:14:654(C):G:C2	26:14:654(S):G:C2	3.08	0.42
1:13:664:G:P	18:9I:64:ARG:HH21	2.42	0.42
38:98:12:ARG:HD3	38:98:16:HIS:CD2	2.53	0.42
1:1G:1063:C:H3'	1:1G:1064:G:H2'	2.02	0.42
22:1L:73:A:H5'	22:1L:74:C:OP2	2.19	0.42
9:82:111:ARG:HG2	9:82:112:LYS:N	2.34	0.42
26:1H:1825:A:O4'	28:11:254:THR:HG21	2.20	0.42
1:1G:165:C:H2'	1:1G:166:G:C8	2.55	0.42
26:1H:1113:U:H2'	26:1H:1114:G:C8	2.54	0.42
26:14:1902:C:N4	26:14:1903:G:C4	2.87	0.42
32:51:130:ARG:HB3	32:51:130:ARG:NH1	2.35	0.42
26:1H:950:G:C5	26:1H:951:C:C4	3.08	0.42
26:14:1479:G:C2'	26:14:1480:G:H5'	2.49	0.42
26:1H:300:A:H1'	26:1H:319:C:O4'	2.19	0.42
26:14:819:A:H2'	26:14:820:A:H5'	2.01	0.42
24:3L:66:U:H2'	24:3L:67:C:C6	2.55	0.42
3:2E:84:ILE:HD11	3:2E:88:ARG:NH2	2.32	0.42
26:1H:2818:G:O2'	26:1H:2837:G:H5'	2.20	0.42
3:2E:4:LYS:HE3	3:2E:4:LYS:HB3	1.57	0.42
4:32:152:SER:O	4:32:153:ARG:C	2.58	0.42
29:29:197:ILE:HD11	29:29:199:ARG:HH21	1.85	0.42
50:L8:7:LYS:C	50:L8:54:VAL:HG23	2.40	0.42
35:25:22:ILE:HA	35:25:22:ILE:HD12	1.43	0.42
26:14:1357:U:OP2	58:14:3624:HOH:O	2.22	0.42
26:1H:2404:C:O3'	36:78:77:ARG:NH2	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:652:U:C4	1:13:752:G:N3	2.88	0.42
26:14:2273:A:H2'	26:14:2274:A:H8	1.81	0.42
4:3E:173:TRP:CE3	4:3E:193:ASP:HB3	2.54	0.42
40:B8:61:PHE:CE1	40:B8:76:PHE:HB2	2.55	0.42
29:21:15:PHE:HA	29:21:19:ARG:O	2.19	0.42
16:7A:36:ILE:HG13	16:7A:36:ILE:O	2.19	0.42
31:49:95:ARG:CG	31:49:96:ARG:H	2.33	0.42
26:14:1794:U:H2'	26:14:1795:C:C6	2.55	0.42
17:8I:63:ARG:HG3	17:8I:64:PRO:N	2.35	0.42
18:9A:45:SER:N	18:9A:49:LYS:O	2.51	0.42
26:1H:2235:G:H2'	26:1H:2236:C:H6	1.85	0.42
26:1H:989:G:C5	50:L8:13:ILE:HG13	2.54	0.42
26:1H:463:G:N2	26:1H:465:G:H3'	2.35	0.42
1:13:1327:C:OP2	21:1F:12:LYS:NZ	2.44	0.42
2:12:22:LYS:H	2:12:22:LYS:HG3	1.70	0.42
32:59:68:THR:HG22	32:59:72:ILE:HD11	2.02	0.42
45:C5:35:TYR:CE2	45:C5:69:ALA:HB3	2.54	0.42
52:J5:41:PRO:HG2	52:J5:44:THR:OG1	2.19	0.42
44:F8:84:ALA:HB3	44:F8:87:GLN:OE1	2.19	0.42
2:1E:130:ARG:HA	2:1E:131:PRO:HD3	1.88	0.42
26:14:1063:G:C4	26:14:1076:C:N4	2.88	0.42
3:22:62:ASP:O	3:22:97:LYS:HB2	2.20	0.42
2:12:71:VAL:HG23	2:12:164:VAL:HG22	2.02	0.42
26:1H:498:G:C6	26:1H:499:U:C4	3.08	0.42
20:BA:76:ALA:O	20:BA:80:ARG:HG2	2.18	0.42
21:1F:5:ASP:O	21:1F:11:GLY:HA3	2.20	0.42
26:1H:681:G:H2'	26:1H:682:G:O4'	2.20	0.42
36:78:112:LEU:HD23	36:78:112:LEU:HA	1.87	0.42
28:19:33:LEU:HA	28:19:33:LEU:HD12	1.70	0.42
33:69:113:ARG:HA	33:69:113:ARG:HD3	1.76	0.42
5:42:99:GLY:O	5:42:117:ASP:HA	2.19	0.42
5:4E:98:THR:HG22	5:4E:99:GLY:O	2.20	0.42
7:6E:65:ALA:HB1	7:6E:127:ALA:HB3	2.01	0.42
4:32:5:ILE:H	4:32:5:ILE:HG13	1.67	0.42
37:88:81:VAL:C	37:88:82:ARG:HD2	2.38	0.42
1:1G:20:U:H2'	1:1G:21:G:O4'	2.20	0.42
1:1G:1356:G:H2'	1:1G:1357:A:C8	2.55	0.42
26:14:934:G:H2'	26:14:935:C:C6	2.55	0.42
26:14:94:G:O2'	49:G5:46:GLN:HB3	2.20	0.42
10:1I:54:PHE:CZ	10:1I:55:LYS:NZ	2.82	0.42
29:21:77:ILE:HG22	29:21:78:LEU:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
44:F8:52:VAL:N	44:F8:82:GLN:O	2.38	0.42
1:1G:639:G:O2'	1:1G:640:A:H5'	2.20	0.42
26:1H:1437:C:H2'	26:1H:1438:U:C6	2.54	0.42
1:1G:1269:A:H5''	1:1G:1270:C:OP2	2.20	0.42
24:3K:18:G:H2'	24:3K:57:G:N2	2.34	0.42
15:6A:87:ILE:HG22	15:6A:88:ARG:N	2.30	0.42
26:1H:142:G:O3'	44:F8:35:THR:HG21	2.19	0.42
26:14:795:C:O2'	26:14:796:C:H5'	2.20	0.42
12:3I:70:ILE:CD1	12:3I:77:LEU:HD12	2.47	0.42
40:B8:50:ILE:CD1	40:B8:102:ILE:HD11	2.50	0.42
26:1H:361:G:H8	26:1H:361:G:O5'	2.03	0.42
27:16:99:A:C4	27:16:100:G:C8	3.07	0.42
39:A8:24:LEU:CD1	39:A8:41:ASP:HB2	2.50	0.42
26:14:1778:U:OP2	58:14:3630:HOH:O	2.22	0.42
2:12:19:HIS:CG	2:12:20:GLU:N	2.86	0.42
39:A8:14:VAL:HG21	39:A8:89:ARG:NE	2.29	0.42
8:72:31:PHE:CE2	8:72:35:ILE:HD11	2.55	0.42
1:13:323:U:H5'	20:BI:23:ARG:HB2	2.01	0.42
45:C5:88:LYS:O	45:C5:89:PHE:HB3	2.19	0.42
26:1H:448:U:O4	26:1H:583:G:H1'	2.19	0.42
1:1G:1006:C:H2'	1:1G:1007:C:C6	2.55	0.42
1:1G:78:G:C2	1:1G:79:G:C8	3.08	0.42
34:15:38:HIS:H	34:15:38:HIS:CD2	2.36	0.42
22:1L:26:A:H5''	22:1L:27:G:N7	2.34	0.42
1:13:1121:U:H2'	1:13:1122:U:C6	2.55	0.42
11:2I:69:ALA:O	11:2I:73:MET:HG3	2.19	0.42
8:72:97:VAL:HG22	8:72:129:VAL:C	2.40	0.42
1:1G:522:C:H2'	1:1G:523:A:O4'	2.20	0.42
38:55:67:LEU:O	38:55:71:GLN:N	2.48	0.42
26:14:235:U:H2'	26:14:236:C:H6	1.85	0.42
1:1G:1442:G:C5	1:1G:1446:A:C6	3.08	0.42
35:25:66:LYS:HD2	35:25:81:ASP:HA	2.01	0.42
33:61:95:LYS:HE3	33:61:99:GLU:OE2	2.20	0.42
26:14:2831:G:P	29:29:58:ARG:HH11	2.43	0.42
17:8I:59:ILE:HB	17:8I:71:PHE:HD2	1.85	0.42
39:65:66:ALA:HA	39:65:69:VAL:HG12	2.00	0.42
30:39:143:ALA:HB1	30:39:148:LEU:HB2	2.02	0.42
26:14:1426:G:H8	26:14:1426:G:O5'	2.02	0.42
26:1H:483:A:H5''	45:G8:50:ARG:HE	1.84	0.42
46:D5:54:HIS:HB3	46:D5:101:PRO:HG3	2.02	0.42
26:1H:1183:G:H8	26:1H:1183:G:OP2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:996:A:C2	26:1H:997:G:C8	3.08	0.42
1:13:386:C:H2'	1:13:387:U:O4'	2.20	0.42
33:69:78:THR:HG21	33:69:104:GLN:HG3	2.02	0.42
12:3I:34:ARG:HH11	12:3I:82:VAL:HG21	1.84	0.42
26:1H:274:G:H2'	26:1H:275:G:H4'	2.02	0.42
1:1G:45:U:H2'	1:1G:46:G:C8	2.55	0.42
30:39:70:THR:OG1	30:39:71:GLY:N	2.52	0.42
28:19:159:ALA:HB1	28:19:198:ASN:O	2.20	0.42
37:88:58:PHE:O	37:88:60:ARG:N	2.53	0.42
26:1H:2215:G:H2'	26:1H:2216:G:H8	1.85	0.42
15:6I:4:THR:O	15:6I:7:GLU:HB3	2.20	0.42
26:14:869:G:C2	26:14:870:A:C8	3.08	0.42
22:1K:23:A:H2'	22:1K:24:G:O4'	2.19	0.42
42:95:83:ARG:HD2	42:95:83:ARG:HA	1.78	0.42
26:1H:1058:U:H2'	26:1H:1059:G:H8	1.84	0.42
36:35:86:LYS:HB3	36:35:118:GLY:HA3	2.02	0.42
26:14:2100:G:N2	26:14:2190:G:H1'	2.34	0.42
38:98:57:ARG:HB3	38:98:59:ASP:OD1	2.19	0.42
4:32:24:GLU:HG2	4:32:25:ARG:H	1.84	0.42
31:41:43:LEU:N	31:41:88:ILE:O	2.46	0.42
26:1H:250:G:H5'	36:78:60:MET:SD	2.59	0.42
26:1H:1689:A:C2	26:1H:1690:A:C5	3.07	0.42
29:21:116:VAL:HG11	29:21:138:PRO:CB	2.44	0.42
26:1H:732:C:H2'	26:1H:733:G:O4'	2.20	0.42
26:14:2394:C:H2'	26:14:2395:C:C6	2.54	0.42
20:BI:30:LYS:HD2	20:BI:30:LYS:HA	1.78	0.42
26:1H:1728:G:H3'	26:1H:1729:A:H5''	2.02	0.42
16:7A:74:LEU:HB3	16:7A:80:PHE:HE1	1.85	0.42
36:78:59:LEU:O	55:Q8:13:ARG:HD3	2.19	0.42
26:14:2750:A:H5'	32:59:6:ARG:HH22	1.85	0.42
1:1G:1278:U:H5'	1:1G:1279:A:O4'	2.20	0.42
26:1H:2032:G:H1'	29:21:145:LYS:HE3	2.02	0.42
24:3K:32:U:O2	24:3K:33:U:H5''	2.20	0.42
26:1H:849:A:H5''	26:1H:850:C:OP2	2.20	0.42
26:1H:1858:G:H2'	26:1H:1883:G:H22	1.84	0.42
26:1H:299:A:N6	26:1H:300:A:N6	2.68	0.42
31:49:109:VAL:O	31:49:113:ARG:HG3	2.20	0.42
1:1G:186(A):C:H2'	1:1G:186(B):C:C6	2.55	0.42
1:1G:186(A):C:H2'	1:1G:186(B):C:H6	1.85	0.42
26:14:1000:A:C6	26:14:1001:A:C6	3.08	0.42
26:1H:2790:A:H4'	26:1H:2791:C:OP2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:4A:23:TYR:HB3	13:4A:67:GLU:HA	2.02	0.42
26:1H:2849:U:P	40:B8:95:ARG:HH12	2.42	0.42
26:14:2789:C:H1'	26:14:2892:A:H2	1.85	0.42
26:14:975:G:H1'	26:14:990:A:C2	2.55	0.42
24:3L:29:G:H2'	24:3L:30:G:C8	2.55	0.42
50:L8:8:LEU:HD13	50:L8:31:LEU:HA	2.01	0.42
1:1G:1227:A:O2'	13:4A:115:LYS:HD2	2.20	0.42
26:14:1359:A:N7	26:14:1372:U:C4	2.88	0.42
1:13:730:G:N2	1:13:766:A:OP1	2.43	0.42
26:14:30:G:H2'	26:14:31:C:C6	2.55	0.42
32:59:89:ILE:HG23	32:59:130:ARG:HA	2.02	0.42
26:14:2014:A:H2'	26:14:2015:A:C8	2.55	0.42
26:1H:1442:G:C2	26:1H:1550:C:O2	2.73	0.42
26:14:756:C:H2'	26:14:757:U:H5'	2.01	0.42
26:14:905:U:H5''	26:14:906:G:OP2	2.20	0.42
1:1G:887:G:H21	1:1G:1489:G:H4'	1.85	0.42
43:A5:89:ALA:O	43:A5:90:ARG:HB2	2.20	0.42
39:65:62:LYS:HA	39:65:65:VAL:HG12	2.02	0.42
26:14:1750:G:O2'	26:14:1751:C:H5'	2.20	0.42
26:14:914:C:N3	26:14:915:C:H1'	2.35	0.42
26:14:107:C:H2'	26:14:108:U:C6	2.55	0.42
1:1G:1113:C:H2'	1:1G:1114:C:C6	2.55	0.42
6:52:68:PRO:HG2	6:52:71:ARG:HG3	2.01	0.42
28:11:67:PHE:CD1	28:11:153:ALA:HB3	2.55	0.42
1:1G:1402:C:H2'	1:1G:1403:C:O4'	2.19	0.42
33:61:79:ILE:O	33:61:142:VAL:HA	2.20	0.42
1:1G:934:C:O2'	1:1G:1344:C:OP2	2.27	0.42
26:14:1986:A:OP1	58:14:3747:HOH:O	2.21	0.42
26:14:1374:G:H2'	26:14:1375:C:C6	2.55	0.42
1:1G:28:G:C6	1:1G:29:G:C5	3.07	0.42
9:82:48:GLU:N	9:82:49:PRO:HD2	2.35	0.42
26:1H:270:A:OP1	58:1H:4580:HOH:O	2.21	0.42
45:C5:7:VAL:HG21	45:C5:37:VAL:HG11	2.02	0.42
26:14:769:G:H5'	26:14:1379:A:N6	2.34	0.42
26:1H:2076:U:H5	26:1H:2596:U:O2	2.02	0.42
26:14:773:U:H5'	28:19:47:GLY:HA3	2.02	0.42
26:14:691:C:H2'	26:14:692:C:C6	2.55	0.42
28:11:172:TYR:CD1	28:11:186:HIS:HA	2.55	0.42
2:1E:72:GLY:HA3	2:1E:165:VAL:HG22	2.01	0.42
19:AA:15:LEU:HD22	19:AA:15:LEU:HA	1.96	0.42
22:1K:70:G:H4'	22:1K:70:G:OP1	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:3E:188:LEU:HD22	4:3E:188:LEU:HA	1.74	0.42
54:L5:36:GLN:HG2	54:L5:36:GLN:O	2.20	0.42
36:35:6:LEU:HA	36:35:6:LEU:HD13	1.77	0.42
19:AI:20:LEU:O	19:AI:23:ASN:HB2	2.19	0.42
1:1G:146:G:C2	1:1G:147:G:C8	3.08	0.42
42:95:85:LYS:HG3	42:95:87:HIS:H	1.84	0.41
26:1H:818:G:N7	26:1H:1187:G:C6	2.88	0.41
1:13:247:G:OP2	17:8I:100:LYS:HB2	2.20	0.41
26:1H:2408:U:H2'	26:1H:2409:G:H8	1.85	0.41
8:72:12:ARG:HH21	8:72:27:PRO:HD3	1.84	0.41
35:68:22:ILE:HD13	35:68:22:ILE:HA	1.45	0.41
34:58:28:THR:HA	34:58:106:MET:HE2	2.02	0.41
5:42:102:ALA:HB2	5:42:120:THR:HG21	2.02	0.41
26:14:273(C):C:H5'	26:14:273(D):C:OP2	2.20	0.41
26:14:638:G:C6	26:14:639:U:C4	3.07	0.41
29:29:26:ILE:HG22	29:29:28:ALA:N	2.35	0.41
1:13:1268:A:N3	1:13:1326:C:O2'	2.52	0.41
26:1H:2347:C:P	53:O8:39:TYR:OH	2.76	0.41
53:O8:10:LEU:CD2	55:Q8:32:LEU:HD22	2.48	0.41
1:13:1286:A:H5''	21:1F:26:LYS:HD3	2.02	0.41
26:1H:993:G:C6	26:1H:1162:G:C6	3.08	0.41
26:1H:1113:U:OP1	32:51:2:SER:N	2.53	0.41
1:1G:191:G:C2	1:1G:192:U:C2	3.07	0.41
11:2A:85:ARG:HD3	11:2A:113:PRO:HD3	2.02	0.41
38:98:91:GLN:H	38:98:91:GLN:NE2	2.18	0.41
26:14:2844:G:N2	26:14:2874:C:C2	2.88	0.41
26:14:792:G:H3'	26:14:793:A:H5'	2.02	0.41
26:1H:71:A:C2	44:F8:31:HIS:CE1	3.04	0.41
1:13:119:A:N6	1:13:288:A:H1'	2.35	0.41
43:E8:59:VAL:HG12	43:E8:60:ASN:HD22	1.85	0.41
40:75:19:LEU:HB3	40:75:86:ILE:HG21	2.01	0.41
23:2L:20:G:H1	23:2L:57:C:H42	1.68	0.41
1:1G:1095:U:OP1	1:1G:1108:G:N1	2.48	0.41
26:14:817:C:H6	26:14:817:C:O5'	2.03	0.41
1:1G:736:C:OP2	18:9A:68:LYS:HE2	2.20	0.41
1:1G:1227:A:C8	1:1G:1227:A:H3'	2.55	0.41
1:13:1072:G:H2'	1:13:1073:U:O4'	2.20	0.41
51:I5:49:PHE:CD2	51:I5:50:VAL:HG22	2.52	0.41
30:39:66:PRO:O	30:39:67:GLN:CB	2.68	0.41
38:55:8:ARG:NE	38:55:43:GLU:OE2	2.53	0.41
4:3E:142:PRO:HG3	4:3E:187:ARG:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:617:G:H2'	26:1H:618:G:O4'	2.19	0.41
44:B5:35:THR:O	44:B5:39:ILE:HG13	2.19	0.41
1:13:985:C:H2'	1:13:986:A:H8	1.85	0.41
24:3L:35:A:H2	25:4L:14:A:N6	2.18	0.41
24:3L:65:G:N3	24:3L:65:G:H2'	2.35	0.41
28:19:37:LEU:CB	28:19:38:LYS:HG2	2.50	0.41
8:7E:25:ASP:HA	8:7E:59:LEU:O	2.20	0.41
26:1H:721:C:H2'	26:1H:722:A:H8	1.85	0.41
1:1G:567:G:N2	58:1G:1833:HOH:O	2.50	0.41
9:8E:46:ALA:HB2	9:8E:74:ILE:HG23	2.01	0.41
1:1G:580:U:H2'	1:1G:581:G:O4'	2.20	0.41
11:2I:57:THR:HA	11:2I:58:PRO:HD2	1.76	0.41
4:3E:57:ARG:HB3	4:3E:206:PHE:HB2	2.02	0.41
45:G8:44:ILE:H	45:G8:44:ILE:HG13	1.67	0.41
39:65:36:TYR:HA	39:65:52:SER:HB3	2.02	0.41
26:1H:1666:G:O2'	26:1H:1667:G:H5'	2.19	0.41
16:7A:43:LYS:HG3	16:7A:48:TRP:CE3	2.55	0.41
26:1H:1206:G:C6	26:1H:1207:C:C4	3.07	0.41
18:9I:38:GLU:HG2	18:9I:38:GLU:H	1.61	0.41
26:14:2093:G:OP1	33:69:23:PRO:HG2	2.20	0.41
26:1H:2850:A:C8	26:1H:2869:G:O4'	2.72	0.41
34:15:115:ARG:O	34:15:118:LYS:HB2	2.20	0.41
26:14:947:G:N2	26:14:971:C:C2	2.88	0.41
31:49:116:ASP:O	31:49:117:PHE:HB3	2.19	0.41
30:31:116:ASP:O	30:31:120:GLU:HG3	2.20	0.41
28:19:260:ARG:NH2	28:19:264:LYS:HD3	2.34	0.41
22:1L:29:G:N2	22:1L:42:C:O2	2.52	0.41
26:1H:88:G:C2'	26:1H:89:G:H5'	2.49	0.41
28:11:4:LYS:HD3	28:11:4:LYS:HA	1.90	0.41
26:1H:718:A:O5'	26:1H:718:A:H8	2.01	0.41
26:1H:234:C:N3	26:1H:235:U:C5	2.88	0.41
36:35:124:LYS:HE2	36:35:124:LYS:HA	2.02	0.41
26:1H:2546:U:O4	58:1H:4469:HOH:O	2.21	0.41
1:1G:406:G:C2	1:1G:407:G:C8	3.09	0.41
1:13:1399:C:C2	1:13:1502:A:N6	2.88	0.41
27:16:12:C:C2	47:18:74:ARG:NH1	2.88	0.41
41:85:65:ILE:HG22	41:85:66:ASN:N	2.34	0.41
1:1G:1320:C:C4	1:1G:1321:C:C4	3.08	0.41
26:1H:1690:A:H2'	26:1H:1691:C:O4'	2.20	0.41
26:1H:442:G:C6	26:1H:444:C:N4	2.88	0.41
26:1H:2061:G:H5''	26:1H:2503:A:C2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:835:U:OP2	18:9I:60:ALA:HB3	2.19	0.41
7:62:26:PHE:O	7:62:30:ILE:HG13	2.20	0.41
28:19:92:ILE:HA	28:19:105:ILE:O	2.20	0.41
7:62:78:ARG:O	7:62:84:ASN:HA	2.19	0.41
41:85:88:ILE:HG22	41:85:90:VAL:CG2	2.50	0.41
26:14:2250:G:OP1	26:14:2275:C:H2'	2.19	0.41
13:4A:96:LEU:HD22	13:4A:97:PRO:HD2	2.01	0.41
29:29:34:VAL:HG11	29:29:64:LYS:HD3	2.01	0.41
23:2K:9:G:N3	23:2K:46:G:H2'	2.35	0.41
44:B5:11:PRO:HD3	49:G5:37:PHE:CD2	2.55	0.41
1:1G:707:C:OP1	11:2A:85:ARG:NH1	2.54	0.41
47:E5:21:LEU:HA	47:E5:21:LEU:HD23	1.88	0.41
49:K8:17:SER:HB3	49:K8:18:PRO:HD2	2.02	0.41
37:88:11:LYS:HE2	37:88:86:GLY:HA2	2.00	0.41
28:11:182:LEU:HA	28:11:182:LEU:HD23	1.89	0.41
50:H5:50:VAL:HG23	50:H5:54:VAL:HG21	2.02	0.41
1:1G:1055:A:C6	1:1G:1206:G:C4	3.08	0.41
4:32:173:TRP:HA	4:32:187:ARG:HD3	2.02	0.41
37:45:54:MET:HE2	37:45:118:LEU:HD22	2.01	0.41
45:C5:48:ALA:HB1	45:C5:50:ARG:NE	2.36	0.41
23:2K:8:4SU:H6	23:2K:8:4SU:O5'	2.19	0.41
29:21:188:VAL:HA	29:21:189:PRO:HD3	1.72	0.41
36:78:116:GLY:H	36:78:134:ALA:HB2	1.84	0.41
27:16:28:C:OP1	39:A8:31:SER:OG	2.28	0.41
26:14:2857:G:C6	26:14:2861:G:O6	2.73	0.41
26:1H:1215:G:C5	26:1H:1216:G:N7	2.88	0.41
26:14:2256:G:H2'	26:14:2257:U:C6	2.54	0.41
6:5E:72:VAL:CG2	6:5E:90:VAL:HG11	2.51	0.41
33:61:75:LEU:HB3	33:61:105:HIS:HD2	1.84	0.41
3:2E:152:ILE:HG12	3:2E:167:TRP:HB2	2.02	0.41
26:1H:1110:G:O2'	26:1H:1111:A:O5'	2.32	0.41
1:13:712:A:C6	1:13:713:G:C6	3.07	0.41
1:1G:598:U:H2'	1:1G:599:C:C6	2.55	0.41
1:1G:509:A:C6	1:1G:510:A:N1	2.87	0.41
31:49:96:ARG:HH11	31:49:96:ARG:HG2	1.85	0.41
26:1H:1547:C:H2'	26:1H:1548:C:H6	1.84	0.41
13:4I:82:MET:C	13:4I:84:ILE:H	2.22	0.41
2:1E:17:PHE:HB3	2:1E:44:LEU:HD21	2.02	0.41
1:1G:487:A:H2'	1:1G:488:C:O4'	2.20	0.41
26:14:1639:U:OP2	58:14:4066:HOH:O	2.22	0.41
8:7E:82:HIS:CE1	8:7E:138:TRP:NE1	2.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:D5:99:TYR:CE2	46:D5:125:LEU:HD12	2.56	0.41
2:1E:47:THR:O	2:1E:51:LEU:HB2	2.20	0.41
1:1G:1374:A:H2'	1:1G:1375:A:H5'	2.02	0.41
1:13:397:A:N7	1:13:548:G:C8	2.89	0.41
31:41:20:ILE:O	31:41:24:GLY:N	2.52	0.41
1:13:977:A:C8	1:13:1223:C:N3	2.88	0.41
22:1L:21:A:N6	22:1L:46:G:N3	2.68	0.41
26:1H:1647:G:C8	26:1H:1647:G:H5'	2.54	0.41
28:11:65:ILE:HD12	28:11:66:ASP:N	2.34	0.41
23:2L:32:G:H5''	23:2L:33:OMC:OP2	2.20	0.41
1:1G:260:G:H2'	1:1G:261:U:C6	2.55	0.41
1:13:358:U:H5''	33:69:87:LYS:HG2	2.01	0.41
50:L8:10:LYS:HD3	50:L8:53:LEU:HD23	2.00	0.41
37:45:4:PRO:HD3	37:45:70:PRO:O	2.20	0.41
26:14:439:G:H2'	26:14:440:G:C8	2.55	0.41
26:14:844:C:C2'	26:14:845:G:H5'	2.50	0.41
26:1H:2679:A:H4'	29:21:165:VAL:HG11	2.02	0.41
33:69:90:GLY:O	33:69:121:LYS:HD2	2.20	0.41
26:1H:1167:U:C2	26:1H:1183:G:N2	2.88	0.41
26:14:868:U:C4	26:14:869:G:N7	2.88	0.41
26:1H:612:G:H2'	26:1H:613:U:O2	2.20	0.41
6:5E:4:TYR:CD1	6:5E:92:LYS:HA	2.55	0.41
28:19:95:LEU:O	28:19:102:LYS:HA	2.20	0.41
26:1H:774:A:H2	26:1H:787:U:HO2'	1.61	0.41
1:13:260:G:H2'	1:13:261:U:C6	2.55	0.41
1:13:9:G:C2	1:13:26:A:N1	2.88	0.41
1:13:1054:C:N3	22:1K:34:G:H1'	2.34	0.41
1:13:1405:G:H1'	1:13:1519:A:O4'	2.20	0.41
36:78:79:ARG:HB2	36:78:110:TYR:HD1	1.85	0.41
12:3I:21:LYS:HD2	12:3I:21:LYS:N	2.34	0.41
26:1H:146:G:H2'	26:1H:147:U:O4'	2.20	0.41
26:1H:776:G:C8	26:1H:793:A:C2	3.07	0.41
1:1G:944:G:C4	1:1G:1340:A:N1	2.88	0.41
1:1G:991:U:O4	1:1G:1212:U:O2'	2.28	0.41
1:13:646:U:H2'	1:13:647:C:C6	2.54	0.41
4:32:30:LYS:HB2	4:32:35:ARG:HD2	2.02	0.41
26:1H:1614:A:N6	43:E8:88:ARG:H	2.17	0.41
1:13:1129:C:N4	1:13:1142:G:O6	2.52	0.41
54:P8:47:ARG:HH11	54:P8:47:ARG:HB2	1.84	0.41
46:H8:30:ASN:ND2	46:H8:90:VAL:HB	2.19	0.41
26:1H:2408:U:H6	26:1H:2408:U:O5'	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:654:A:OP1	26:14:654:A:H8	2.03	0.41
1:1G:1320:C:N3	19:AA:36:ARG:NH2	2.68	0.41
13:4A:84:ILE:C	13:4A:86:CYS:H	2.23	0.41
26:1H:991:C:H42	26:1H:1163:G:H1	1.68	0.41
26:1H:1332:G:P	58:1H:4559:HOH:O	2.78	0.41
26:1H:1952:A:C6	35:68:22:ILE:CD1	2.99	0.41
33:61:40:THR:HG22	33:61:41:GLU:OE1	2.20	0.41
9:82:14:VAL:O	9:82:65:VAL:HA	2.20	0.41
46:H8:126:VAL:HA	46:H8:164:ALA:H	1.84	0.41
41:C8:88:ILE:C	41:C8:90:VAL:H	2.23	0.41
1:1G:1002:G:H22	1:1G:1038:C:H42	1.67	0.41
1:13:682:G:H2'	1:13:683:G:O4'	2.21	0.41
1:1G:376:G:H5''	16:7A:5:ARG:CD	2.50	0.41
20:BA:89:ARG:HH11	20:BA:104:LEU:HB3	1.86	0.41
26:14:2065:C:H1'	26:14:2449:U:H3	1.86	0.41
36:78:144:GLU:N	36:78:144:GLU:CD	2.74	0.41
26:14:522:G:H2'	26:14:523:C:C6	2.55	0.41
12:3A:110:VAL:CG2	12:3A:120:TYR:HB3	2.51	0.41
26:14:2639:A:C2	26:14:2778:A:C8	3.08	0.41
40:B8:19:LEU:HA	40:B8:20:PRO:HD3	1.81	0.41
13:4A:37:THR:HG21	13:4A:56:LEU:HD23	2.01	0.41
30:31:33:LEU:HD23	36:78:1:MET:HG3	2.01	0.41
1:1G:1424:C:H2'	1:1G:1425:U:O4'	2.20	0.41
50:L8:7:LYS:HD2	50:L8:34:GLU:HG2	2.01	0.41
1:1G:1023:G:H2'	1:1G:1024:G:H5''	2.02	0.41
34:15:12:ARG:HH21	34:15:38:HIS:CE1	2.38	0.41
1:13:1080:A:H5''	5:4E:16:THR:HG21	2.01	0.41
35:25:22:ILE:HG22	35:25:40:VAL:HB	2.01	0.41
1:13:1120:G:C2	1:13:1154:G:C2	3.09	0.41
26:14:1708:C:O2'	26:14:1709:U:H5'	2.19	0.41
1:1G:949:A:H2'	1:1G:950:U:C6	2.55	0.41
1:13:453:A:C6	1:13:454:C:C4	3.08	0.41
26:1H:1448:G:N2	26:1H:1449:A:N6	2.68	0.41
49:G5:50:ILE:HG22	49:G5:54:LYS:HE3	2.03	0.41
54:P8:12:ARG:NH2	54:P8:44:PRO:HB3	2.35	0.41
33:61:107:VAL:HG12	33:61:108:THR:N	2.35	0.41
1:1G:309:G:H1'	1:1G:608:A:C2	2.55	0.41
3:22:72:LYS:HG3	3:22:75:VAL:HG23	2.02	0.41
39:65:67:ARG:CZ	39:65:67:ARG:HB2	2.50	0.41
26:1H:99:U:C6	26:1H:102:G:C2	3.08	0.41
1:1G:284:G:H2'	1:1G:285:G:C8	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:H8:146:ILE:HD13	46:H8:174:VAL:HG12	2.02	0.41
1:13:266:G:H8	1:13:266:G:H2'	1.75	0.41
7:6E:26:PHE:CD2	7:6E:30:ILE:HD11	2.55	0.41
28:11:6:PHE:CE1	28:11:18:VAL:HG23	2.55	0.41
26:1H:1665:A:C2'	26:1H:1666:G:H5'	2.50	0.41
26:1H:2543:G:H2'	26:1H:2544:G:O4'	2.20	0.41
26:1H:18:C:O3'	41:C8:23:GLY:HA2	2.20	0.41
1:13:381:C:H2'	1:13:382:A:O4'	2.20	0.41
26:14:2729:G:H2'	26:14:2730:C:C6	2.56	0.41
13:4A:15:VAL:HA	13:4A:18:ALA:HB3	2.02	0.41
47:I8:47:PRO:HB3	47:I8:51:VAL:HG12	2.01	0.41
38:98:37:THR:OG1	38:98:39:PRO:HD2	2.21	0.41
1:1G:781:A:H4'	1:1G:1522:U:O2'	2.21	0.41
47:I8:31:VAL:HG21	47:I8:61:ALA:HB2	2.02	0.41
30:31:117:ARG:O	30:31:120:GLU:HB2	2.21	0.41
26:1H:255:A:H1'	26:1H:384:U:C6	2.55	0.41
26:1H:2749:A:H1'	32:51:63:SER:OG	2.19	0.41
26:1H:1173:G:N2	26:1H:1175:U:H5	2.18	0.41
53:O8:21:TYR:HB2	53:O8:22:ALA:H	1.50	0.41
26:1H:745:G:OP2	29:21:133:LYS:HE2	2.20	0.41
26:14:962:G:C2	26:14:963:U:C2	3.08	0.41
9:8E:77:ILE:O	9:8E:81:ILE:HG12	2.20	0.41
26:1H:2239:G:P	58:1H:4527:HOH:O	2.78	0.41
26:14:860:U:C2	26:14:2268:A:C8	3.08	0.41
26:14:2366:A:H2'	26:14:2367:G:O4'	2.20	0.41
26:14:725:G:H8	26:14:725:G:O5'	2.03	0.41
26:14:2407:G:C8	26:14:2407:G:O5'	2.74	0.41
51:I5:9:LEU:HD22	51:I5:9:LEU:H	1.85	0.41
26:1H:1572:A:O5'	26:1H:1572:A:H8	2.03	0.41
35:25:90:GLN:O	35:25:91:LEU:HB2	2.19	0.41
26:1H:550:G:O2'	26:1H:1220:A:N3	2.40	0.41
10:1A:68:HIS:HB3	10:1A:70:ARG:HH21	1.85	0.41
1:13:95:G:H3'	1:13:96:G:H8	1.85	0.41
26:1H:94:G:H2'	26:1H:95:G:O4'	2.20	0.41
1:13:976:G:C8	1:13:1362:C:N4	2.87	0.41
46:D5:39:VAL:HG21	46:D5:44:PHE:CD2	2.52	0.41
26:14:2428:G:H21	36:35:61:ARG:HH11	1.69	0.41
1:13:658:G:H2'	1:13:659:U:H6	1.85	0.41
26:1H:1389:G:C2	26:1H:1390:U:C2	3.08	0.41
26:14:660:G:N2	36:35:12:ALA:HB2	2.26	0.41
1:1G:1349:A:P	9:82:118:LYS:HZ1	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:650:C:H5''	26:14:651:G:OP2	2.20	0.41
1:13:428:G:C8	1:13:430:A:C5	3.08	0.41
55:Q8:32:LEU:HG	55:Q8:33:ASN:N	2.35	0.41
26:1H:2348:U:OP1	55:Q8:36:LYS:HD2	2.20	0.41
48:J8:91:LYS:O	48:J8:93:GLU:N	2.53	0.41
26:14:2271:G:H5''	47:E5:20:ARG:NE	2.34	0.41
26:14:2472:G:N1	26:14:2476:A:O2'	2.53	0.41
16:7A:9:PHE:CE1	16:7A:18:ARG:HG3	2.56	0.41
1:13:1350:A:C2	1:13:1351:U:C2	3.08	0.41
26:1H:142:G:H2'	26:1H:143:C:C6	2.55	0.41
26:14:1762:A:H4'	26:14:1763:G:OP2	2.19	0.41
37:45:102:VAL:O	37:45:102:VAL:HG12	2.20	0.41
26:14:270(P):C:H2'	26:14:270(Q):C:C6	2.55	0.41
22:1K:21:A:N7	22:1K:46:G:C6	2.89	0.41
34:58:17:ASP:O	34:58:56:ASN:HB2	2.21	0.41
2:1E:27:LYS:HB2	2:1E:194:PRO:HD2	2.02	0.41
38:55:52:ILE:O	38:55:54:LEU:N	2.53	0.41
26:1H:583:G:OP2	41:C8:10:ARG:HD2	2.20	0.41
2:12:7:VAL:HG13	2:12:8:LYS:HG3	2.03	0.41
1:13:607:A:C2	16:7I:31:LYS:HG3	2.56	0.41
7:62:65:ALA:CB	7:62:124:LEU:HD22	2.51	0.41
26:1H:2811:G:OP1	29:21:60:ASN:HB2	2.21	0.41
26:14:2536:G:C6	26:14:2537:U:C4	3.09	0.41
26:14:1952:A:C2	35:25:22:ILE:HG13	2.55	0.41
4:3E:150:GLU:N	4:3E:150:GLU:OE1	2.53	0.41
13:4I:80:ARG:HH11	19:AI:65:ASN:HB2	1.85	0.41
49:K8:14:ARG:HD3	49:K8:66:GLU:OE1	2.19	0.41
1:1G:337:C:H2'	1:1G:338:A:C8	2.55	0.41
9:82:33:PHE:HE1	9:82:37:PHE:CD2	2.35	0.41
1:13:1277:C:O2'	1:13:1279:A:H1'	2.21	0.41
26:1H:2732:G:H3'	26:1H:2733:A:C4'	2.50	0.41
1:13:1193:G:P	3:2E:167:TRP:HZ3	2.43	0.41
26:1H:633:A:H2'	26:1H:634:C:H5'	2.02	0.41
6:5E:100:ASN:O	18:9I:28:GLU:HG3	2.19	0.41
32:59:85:LYS:HZ2	32:59:142:GLY:HA2	1.85	0.41
4:3E:105:VAL:HG13	4:3E:110:PHE:HB2	2.01	0.41
37:45:66:ILE:HG13	37:45:67:ARG:N	2.35	0.41
1:1G:1254:C:OP1	10:1A:45:ARG:HA	2.19	0.41
26:1H:2475:C:O2'	26:1H:2476:A:H5'	2.20	0.41
1:13:615:C:N3	1:13:616:G:N7	2.68	0.41
43:A5:90:ARG:HG3	43:A5:90:ARG:HH11	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:720:C:H2'	26:1H:721:C:C6	2.56	0.41
1:13:1207:G:C6	1:13:1208:C:C4	3.09	0.41
26:1H:59:U:HO2'	26:1H:73:A:H2'	1.84	0.41
1:1G:1057:G:O2'	1:1G:1058:G:H5'	2.20	0.41
1:13:16:A:O2'	1:13:17:U:H5'	2.20	0.41
37:45:43:THR:HG22	37:45:94:VAL:HG12	2.01	0.41
26:1H:25:U:H2'	26:1H:26:G:C8	2.54	0.41
4:32:42:GLN:HG2	4:32:43:HIS:ND1	2.36	0.41
26:14:663:G:H2'	26:14:664:C:O4'	2.20	0.41
26:14:699:A:H2'	26:14:700:G:O4'	2.20	0.41
28:19:231:HIS:ND1	28:19:232:PRO:HD2	2.35	0.41
51:M8:49:PHE:CD2	51:M8:50:VAL:HB	2.55	0.41
41:C8:17:ILE:O	41:C8:20:LEU:HB2	2.20	0.41
34:15:5:VAL:HG12	34:15:6:PRO:HD2	2.02	0.41
26:1H:200:U:O2	26:1H:386:G:N2	2.53	0.41
26:14:911:A:H2'	37:45:9:TYR:OH	2.20	0.41
1:13:236:G:H5''	17:8I:42:TYR:OH	2.20	0.41
10:1I:3:LYS:N	10:1I:75:ILE:HA	2.34	0.41
17:8I:58:GLU:HB2	17:8I:74:LEU:HB3	2.02	0.41
50:L8:30:ARG:HG3	50:L8:30:ARG:H	1.76	0.41
24:3L:39:U:H2'	24:3L:39:U:O2	2.20	0.41
48:F5:34:THR:O	48:F5:34:THR:HG22	2.20	0.41
28:19:127:VAL:HA	28:19:193:VAL:HG22	2.02	0.41
4:32:113:SER:OG	4:32:116:GLN:HB2	2.20	0.41
45:C5:75:ILE:CG2	45:C5:76:CYS:N	2.84	0.41
46:H8:59:LEU:HD23	46:H8:59:LEU:HA	1.64	0.41
14:5A:21:TYR:HD1	14:5A:22:THR:O	2.03	0.41
36:35:85:LEU:CA	36:35:88:LEU:HB3	2.41	0.41
1:1G:464:G:N2	1:1G:467:G:C8	2.88	0.41
1:13:827:U:C5	1:13:870:U:C5	3.09	0.41
5:42:136:MET:HB3	5:42:136:MET:HE2	1.99	0.41
5:42:106:PRO:O	5:42:110:LEU:HG	2.21	0.41
26:1H:729:G:O2'	26:1H:763:G:H4'	2.21	0.41
26:14:633:A:H2'	26:14:634:C:H5'	2.02	0.41
26:1H:1825:A:OP1	28:11:249:PRO:HD3	2.21	0.41
26:1H:2171:A:H8	26:1H:2171:A:P	2.43	0.41
26:1H:1668:A:OP1	35:68:5:GLN:HG3	2.20	0.41
1:13:429:U:H1'	1:13:430:A:H5''	2.02	0.41
27:16:7:G:O5'	39:A8:29:PHE:HE2	2.03	0.41
1:1G:1316:G:HO2'	1:1G:1318:A:H62	1.69	0.41
32:51:124:GLU:HG2	32:51:126:PRO:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:374:A:C6	1:1G:375:U:C4	3.09	0.41
26:1H:2694:G:H2'	26:1H:2695:C:H6	1.86	0.41
31:41:112:PRO:HB3	51:M8:36:CYS:CA	2.51	0.41
26:1H:2339:G:N2	26:1H:2340:G:C4	2.89	0.41
1:1G:520:A:OP1	12:3A:52:LEU:HB2	2.20	0.41
27:16:44:G:H1'	27:16:47:C:N4	2.32	0.41
37:45:116:GLU:O	37:45:117:ALA:HB3	2.19	0.41
1:1G:1099:G:C6	1:1G:1100:C:N3	2.88	0.41
22:1K:51:U:H2'	22:1K:52:G:C8	2.47	0.41
9:82:10:ARG:NH1	9:82:105:ASP:OD2	2.53	0.41
23:2L:54:G:H2'	23:2L:55:5MU:C6	2.56	0.41
26:1H:2128:C:N4	26:1H:2160:G:H1	2.17	0.41
2:12:217:ARG:HB2	2:12:217:ARG:NH1	2.36	0.41
26:14:2815:C:C5'	52:J5:29:THR:HG21	2.49	0.41
5:4E:16:THR:OG1	5:4E:17:ALA:N	2.54	0.41
1:13:1154:G:H2'	1:13:1155:G:C8	2.56	0.41
26:14:2656:U:C4	26:14:2664:G:N2	2.89	0.41
22:1L:11:C:H2'	22:1L:12:U:C6	2.56	0.41
26:1H:1198:U:H2'	26:1H:1199:U:C6	2.56	0.41
47:E5:29:GLN:O	47:E5:67:VAL:HG23	2.20	0.41
30:31:64:ILE:HA	30:31:64:ILE:HD13	1.70	0.41
28:19:8:PRO:HB3	28:19:14:ARG:HB2	2.02	0.41
6:52:6:VAL:HG12	6:52:8:ILE:HD11	2.03	0.41
1:1G:247:G:OP2	17:8A:100:LYS:HG2	2.20	0.41
29:21:2:LYS:HD3	29:21:200:GLU:HB2	2.03	0.41
1:1G:1097:C:H2'	1:1G:1098:C:C6	2.56	0.41
26:14:2543:G:H2'	26:14:2544:G:C8	2.55	0.41
26:1H:832:G:H4'	36:78:45:LEU:HD11	2.01	0.41
43:A5:90:ARG:HG3	43:A5:90:ARG:NH1	2.35	0.41
43:A5:70:TYR:O	43:A5:107:LEU:HD12	2.19	0.41
37:88:25:ASP:OD2	46:H8:78:LYS:NZ	2.40	0.41
10:1A:4:ILE:HG12	10:1A:100:THR:HG22	2.02	0.41
13:4I:84:ILE:HD11	19:AI:66:MET:HG2	2.03	0.41
54:L5:12:ARG:HH21	54:L5:44:PRO:HB3	1.86	0.41
30:39:40:GLN:NE2	30:39:182:ASN:HB2	2.34	0.41
1:13:998(A):C:O2'	1:13:999:U:H5'	2.21	0.41
1:1G:1260:C:P	1:1G:1284:C:H4'	2.61	0.41
51:M8:22:ILE:HA	51:M8:24:THR:HG23	2.02	0.41
3:22:73:PRO:O	3:22:77:ILE:HG13	2.20	0.41
38:55:51:LEU:HA	38:55:51:LEU:HD23	1.83	0.41
30:31:179:GLU:OE1	30:31:179:GLU:N	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:52:30:LEU:HD23	6:52:30:LEU:HA	1.82	0.41
30:39:29:ASN:OD1	30:39:32:LEU:HB2	2.21	0.41
52:J5:38:ALA:CB	52:J5:48:GLU:HG3	2.51	0.41
26:1H:824:A:H1'	26:1H:2358:G:N7	2.36	0.41
28:19:210:GLY:O	28:19:213:ARG:HB2	2.19	0.41
17:8A:59:ILE:HG22	17:8A:71:PHE:HD2	1.84	0.41
1:1G:781:A:H2'	1:1G:782:A:H5'	2.03	0.41
17:8I:7:THR:OG1	17:8I:58:GLU:HG2	2.19	0.41
1:13:622:A:C8	1:13:623:C:C6	3.09	0.41
26:1H:253:C:H2'	26:1H:254:G:O4'	2.19	0.41
1:1G:1185:G:H2'	1:1G:1186:G:O4'	2.20	0.41
1:1G:1379:G:OP1	7:62:6:ARG:NH1	2.54	0.41
1:13:1171:G:O2'	1:13:1172:C:H5'	2.20	0.41
26:14:767:U:O2'	26:14:768:G:H5'	2.21	0.41
1:1G:1496:C:H2'	1:1G:1497:G:C8	2.55	0.41
26:14:1319:G:C6	26:14:1320:C:N4	2.89	0.41
32:51:24:VAL:HG11	32:51:72:ILE:HD11	2.03	0.41
26:1H:2619:C:H4'	29:21:151:TYR:O	2.20	0.41
26:1H:339:U:H6	26:1H:339:U:O5'	2.03	0.41
4:3E:160:GLN:NE2	4:3E:160:GLN:O	2.53	0.41
26:14:982:C:O5'	26:14:982:C:H6	2.02	0.41
26:14:1716:U:H2'	26:14:1717:G:H8	1.84	0.41
48:J8:75:GLU:O	48:J8:77:ALA:N	2.54	0.41
41:C8:44:ASN:ND2	42:D8:75:PHE:O	2.39	0.41
26:1H:2614:A:H3'	58:1H:3613:HOH:O	2.20	0.41
26:14:322:A:H5'	26:14:340:A:H1'	2.03	0.41
26:1H:1144:G:H2'	26:1H:1145:C:C6	2.55	0.41
54:P8:9:ARG:HH11	54:P8:9:ARG:HD2	1.71	0.41
26:1H:1299:G:H3'	26:1H:1639:U:O4	2.20	0.41
1:1G:1220:G:H1'	19:AA:52:TYR:HD2	1.86	0.41
36:35:11:GLY:C	36:35:13:ASN:N	2.71	0.41
1:1G:1129:C:C4	1:1G:1139:G:N1	2.89	0.41
26:1H:1498:C:O4'	26:1H:1577:C:H4'	2.20	0.41
26:1H:2393:A:H2'	26:1H:2394:C:H6	1.85	0.41
26:14:609(A):G:N2	26:14:619:G:H1'	2.36	0.41
1:1G:1239:A:H4'	1:1G:1240:U:C5'	2.50	0.41
46:D5:29:TYR:HB3	46:D5:34:ASN:OD1	2.20	0.41
26:14:363(B):G:H2'	26:14:363(C):G:H8	1.84	0.41
41:C8:90:VAL:HA	42:D8:39:LEU:HD23	2.02	0.41
26:14:1925:C:O2'	26:14:1926:U:H5'	2.21	0.41
1:1G:1002:G:H1	1:1G:1038:C:H42	1.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:684:A:N6	1:13:685:G:C6	2.88	0.41
45:G8:94:LYS:HZ2	45:G8:95:LYS:N	2.16	0.41
8:7E:87:SER:OG	8:7E:92:ARG:HA	2.21	0.41
24:3K:32:U:H2'	24:3K:33:U:H5''	2.03	0.41
27:16:99:A:C6	27:16:100:G:C5	3.08	0.41
26:14:71:A:H4'	26:14:72:U:H5''	2.03	0.41
2:12:19:HIS:CD2	2:12:20:GLU:HG2	2.56	0.41
26:14:1420:U:O2'	26:14:1421:G:P	2.79	0.41
1:13:1057:G:C4	1:13:1204:A:C2	3.09	0.41
37:88:16:ARG:HD3	37:88:16:ARG:HH11	1.76	0.41
26:14:2516:G:C6	26:14:2517:C:N4	2.88	0.41
26:1H:2127:G:N1	26:1H:2162:G:N3	2.68	0.41
1:1G:1218:C:H2'	1:1G:1219:U:C6	2.56	0.41
26:14:822:U:O2'	26:14:823:G:H5'	2.21	0.41
26:14:1477:A:H2'	26:14:1478:G:O4'	2.21	0.41
1:13:243:A:C2	1:13:246:A:C8	3.08	0.41
26:14:2556:C:H2'	26:14:2557:G:O4'	2.21	0.41
1:13:1120:G:H2'	1:13:1121:U:C6	2.55	0.41
26:14:1027:A:N6	26:14:1126:A:C4	2.88	0.41
26:1H:1354:A:H2'	26:1H:1355:G:O4'	2.21	0.41
26:14:2859:G:H3'	26:14:2859:G:C8	2.55	0.41
1:1G:109:A:H5'	1:1G:110:C:H5	1.85	0.41
1:1G:1228:C:OP1	13:4A:115:LYS:NZ	2.54	0.41
26:1H:2733:A:C2	26:1H:2734:A:H1'	2.56	0.41
33:69:6:LEU:HD13	33:69:37:VAL:CG2	2.51	0.41
38:55:29:LEU:HD12	38:55:29:LEU:HA	1.75	0.41
26:1H:1474:C:H2'	26:1H:1475:G:C8	2.55	0.41
22:1L:52:G:C2	22:1L:63:G:C2	3.09	0.41
32:59:85:LYS:NZ	32:59:142:GLY:HA2	2.35	0.41
28:11:25:THR:HG21	28:11:113:VAL:HG21	2.03	0.41
4:3E:187:ARG:O	4:3E:189:PRO:HD3	2.21	0.41
8:72:103:VAL:HG21	8:72:110:ALA:HB2	2.02	0.41
1:13:654:G:C6	1:13:655:A:C5	3.08	0.41
26:1H:686:G:H4'	26:1H:687:C:OP2	2.21	0.41
26:1H:658:C:H2'	26:1H:659:C:H6	1.84	0.41
48:F5:65:SER:OG	48:F5:66:HIS:ND1	2.30	0.41
2:1E:108:ILE:O	2:1E:111:ARG:HB2	2.20	0.41
15:6I:8:LYS:O	15:6I:12:ILE:HG13	2.21	0.41
1:13:1152:A:C6	1:13:1153:C:C4	3.09	0.41
26:1H:1181:C:C2'	26:1H:1182:A:H5'	2.51	0.41
26:1H:2119:A:H62	26:1H:2168:G:N2	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:16:94:C:H2'	27:16:95:U:H6	1.86	0.41
1:1G:236:G:H2'	1:1G:237:C:O4'	2.21	0.41
1:1G:570:G:H2'	1:1G:571:U:C6	2.55	0.41
26:1H:2820:A:N3	26:1H:2820:A:H2'	2.36	0.41
17:8A:59:ILE:HD13	17:8A:73:VAL:HA	2.02	0.41
27:1J:97:G:H2'	27:1J:98:G:O4'	2.20	0.41
28:19:133:LEU:HB2	28:19:187:GLY:HA2	2.02	0.41
46:D5:178:GLU:HG3	46:D5:179:ASP:H	1.86	0.41
26:1H:2666:C:H5''	26:1H:2667:C:OP2	2.20	0.41
26:14:2408:U:O2'	26:14:2409:G:H5'	2.21	0.41
35:25:101:PRO:HG3	40:75:67:SER:OG	2.20	0.41
26:1H:526:A:N3	26:1H:2044:C:H1'	2.36	0.41
32:51:56:SER:OG	32:51:58:GLU:HG2	2.21	0.41
26:14:2533:A:N6	26:14:2534:A:C6	2.89	0.41
26:14:242:G:C8	55:M5:5:LYS:HG3	2.55	0.41
1:13:811:C:H4'	1:13:900:A:N6	2.35	0.41
26:14:1830:C:O5'	26:14:1830:C:H6	2.04	0.41
48:F5:49:VAL:HG11	48:F5:70:VAL:HG11	2.03	0.41
51:M8:12:ALA:HA	51:M8:29:PRO:HB3	2.02	0.41
26:1H:2298:A:H5''	26:1H:2299:G:OP2	2.21	0.41
17:8I:67:LYS:O	17:8I:68:ARG:HB3	2.21	0.41
26:1H:911:A:H2'	37:88:9:TYR:OH	2.20	0.41
1:1G:1323:G:H4'	1:1G:1362(A):C:C2	2.56	0.41
1:1G:1324:A:H5'	1:1G:1362:C:O2'	2.21	0.41
26:1H:1331:A:H2'	26:1H:1333:C:C5	2.55	0.41
1:1G:828:A:C4	1:1G:859:A:C8	3.08	0.41
44:F8:49:VAL:CG1	44:F8:50:LYS:N	2.82	0.41
44:F8:50:LYS:O	44:F8:83:VAL:HA	2.21	0.41
43:E8:78:GLU:OE1	43:E8:99:ARG:HD3	2.20	0.41
46:H8:19:ARG:NH1	46:H8:84:GLU:HB2	2.36	0.41
1:1G:1001:G:O6	1:1G:1002:G:N2	2.53	0.41
55:M5:39:LYS:HG2	55:M5:43:GLN:HG3	2.02	0.41
38:98:94:TYR:CD1	38:98:94:TYR:N	2.88	0.41
26:1H:175:G:O2'	26:1H:176:G:H5'	2.21	0.41
26:14:2387:U:H1'	47:E5:41:ARG:CD	2.49	0.41
12:3A:113:ARG:HH21	12:3A:116:SER:HG	1.68	0.41
11:2I:112:THR:HA	11:2I:113:PRO:HD3	1.63	0.41
26:1H:319:C:C2	26:1H:333:G:N2	2.89	0.41
26:1H:1025:G:C4	26:1H:1135:C:H1'	2.56	0.41
26:1H:1025:G:H8	26:1H:1025:G:OP1	2.03	0.41
30:31:34:TRP:CZ2	36:78:8:PRO:HB3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
45:C5:89:PHE:O	45:C5:90:LEU:HB2	2.21	0.41
39:A8:66:ALA:HA	39:A8:69:VAL:CG1	2.49	0.41
26:14:817:C:O2'	26:14:839:U:H5''	2.20	0.41
26:1H:129:C:H2'	26:1H:130:C:H6	1.86	0.41
1:1G:693:G:H2'	1:1G:694:A:O4'	2.21	0.41
3:22:18:TRP:HE3	3:22:18:TRP:H	1.68	0.41
16:7I:4:ILE:HA	16:7I:20:VAL:O	2.21	0.41
1:1G:109:A:C6	1:1G:326:G:C6	3.08	0.41
26:14:1356:G:C5	26:14:1357:U:C5	3.09	0.41
26:1H:528:A:N3	26:1H:528:A:O4'	2.50	0.41
30:31:64:ILE:HG23	30:31:65:TRP:NE1	2.36	0.41
26:14:2839:G:H21	38:55:92:GLY:HA2	1.86	0.41
26:14:872:A:C6	26:14:906:G:C2	3.09	0.41
9:8E:23:ASN:OD1	9:8E:25:LYS:HB2	2.20	0.41
23:2K:24:C:H2'	23:2K:25:U:H6	1.86	0.41
52:J5:36:CYS:HG	52:J5:49:CYS:CB	2.33	0.41
26:1H:1301:A:H2	26:1H:1626:G:H21	1.66	0.41
1:1G:1057:G:C5	1:1G:1204:A:C2	3.08	0.41
28:11:65:ILE:HG13	28:11:67:PHE:CE2	2.56	0.41
26:1H:1973:G:H2'	26:1H:1974:C:C6	2.56	0.41
1:1G:791:G:C5	1:1G:792:A:N7	2.89	0.41
26:14:1235:G:C5	26:14:1236:G:C6	3.08	0.41
37:45:69:PHE:CD1	37:45:70:PRO:HD2	2.56	0.41
29:29:48:GLN:OE1	29:29:78:LEU:HD13	2.21	0.41
26:14:1894:C:O2'	26:14:1895:C:H5'	2.21	0.41
26:1H:242:G:H5'	55:Q8:60:LEU:HD11	2.03	0.41
26:14:663:G:C5	26:14:664:C:C5	3.08	0.41
8:7E:100:ILE:HD13	8:7E:112:LEU:HD21	2.02	0.41
1:1G:145:G:H2'	1:1G:146:G:O4'	2.21	0.41
28:19:133:LEU:HA	28:19:133:LEU:HD23	1.94	0.41
26:14:2733:A:H61	29:29:202:LYS:HB3	1.86	0.41
36:78:147:LEU:O	36:78:148:LEU:HD23	2.20	0.41
3:2E:186:PHE:CE2	3:2E:188:LEU:HD23	2.56	0.41
26:14:553:U:C4	26:14:554:U:C4	3.09	0.41
26:1H:2881:C:C2	26:1H:2882:A:C8	3.08	0.41
43:E8:22:ASP:HA	43:E8:25:ARG:NH1	2.35	0.41
26:1H:2246:G:H2'	26:1H:2247:A:C8	2.56	0.41
23:2L:14:A:C2	23:2L:23:G:C4	3.09	0.41
1:13:1251:A:H2'	1:13:1252:A:C8	2.56	0.41
26:14:2459:A:C5	26:14:2460:U:C5	3.08	0.41
26:14:1062:G:N2	26:14:1077:A:N3	2.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
33:61:117:GLU:HG2	33:61:117:GLU:H	1.47	0.41
50:H5:5:LYS:HB3	50:H5:5:LYS:HE3	1.64	0.41
41:C8:34:LYS:HA	41:C8:34:LYS:HE3	2.03	0.41
17:8A:14:LYS:HB2	17:8A:14:LYS:HE3	1.86	0.41
1:1G:1048:G:C8	1:1G:1048:G:H3'	2.55	0.41
28:11:101:GLU:HG3	28:11:102:LYS:N	2.35	0.41
26:1H:1357:U:C4	26:1H:1358:G:C5	3.08	0.41
55:Q8:4:MET:CB	55:Q8:59:LYS:HD2	2.51	0.41
26:14:2017:U:O2	52:J5:10:LYS:HB2	2.21	0.41
1:13:1348:U:N3	1:13:1374:A:H2	1.94	0.41
41:85:50:ARG:NH2	42:95:72:VAL:HG23	2.34	0.41
26:14:1141:U:H2'	26:14:1141:U:H6	1.59	0.41
26:1H:195:A:H4'	26:1H:251:A:O2'	2.21	0.41
27:1J:40:U:C2'	27:1J:45:A:H61	2.33	0.41
18:9I:53:ARG:HE	18:9I:59:SER:C	2.24	0.41
46:D5:111:VAL:HG12	46:D5:145:GLU:HB2	2.03	0.41
33:69:76:THR:HG21	33:69:140:LEU:HA	2.03	0.41
1:1G:468:A:O2'	16:7A:81:ARG:HA	2.20	0.41
26:14:2392:A:H8	36:35:61:ARG:HD2	1.86	0.41
31:49:104:GLU:HG2	51:I5:23:GLU:CG	2.50	0.41
26:1H:1137:G:H2'	26:1H:1138:G:C8	2.55	0.41
31:41:63:ILE:HG12	31:41:64:THR:N	2.36	0.41
26:14:2128:C:C4	26:14:2129:C:C4	3.09	0.41
26:1H:2283:C:C2	26:1H:2389:G:C2	3.08	0.41
24:3K:25:C:C4	24:3K:26:A:C8	3.09	0.41
26:14:153:C:H42	26:14:173:G:H1	1.67	0.41
1:13:418:C:O2'	1:13:540:G:H1'	2.21	0.41
36:35:144:GLU:HA	36:35:145:PRO:HD3	1.76	0.41
1:1G:18:C:H4'	1:1G:1078:U:O2	2.21	0.41
10:1A:7:LYS:O	10:1A:96:ILE:HB	2.21	0.41
10:1A:8:LEU:HG	10:1A:96:ILE:HG21	2.02	0.41
1:1G:1004:A:N6	1:1G:1025:U:HO2'	2.18	0.41
13:4I:49:THR:C	13:4I:51:ALA:N	2.74	0.41
32:51:8:PRO:HD2	32:51:69:ARG:HE	1.85	0.41
1:1G:437:U:C4	1:1G:438:G:C5	3.09	0.41
45:G8:8:LYS:HG2	45:G8:11:ASP:OD1	2.20	0.41
1:13:1256:A:H4'	1:13:1258:G:C4	2.56	0.41
1:13:1256:A:N3	1:13:1277:C:C4	2.89	0.41
26:14:1709:U:H2'	26:14:1710:C:H6	1.81	0.41
29:21:52:LEU:O	29:21:75:VAL:HG23	2.20	0.41
15:6A:56:LEU:O	15:6A:60:VAL:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:42:76:ILE:HG23	5:42:77:PRO:HD2	2.02	0.41
26:14:270(I):G:H2'	26:14:270(J):G:H8	1.86	0.41
32:59:137:ASP:CB	32:59:140:LYS:HB3	2.49	0.41
26:14:2851:A:H2'	26:14:2852:G:C8	2.55	0.41
26:14:2611:U:O2	52:J5:3:LYS:HE3	2.21	0.41
30:31:40:GLN:NE2	30:31:184:TYR:CB	2.83	0.41
1:1G:382:A:H2'	1:1G:383:A:H8	1.83	0.41
1:13:484:G:HO2'	1:13:485:G:P	2.43	0.41
32:59:120:GLY:O	32:59:121:ILE:HD13	2.21	0.41
40:B8:43:GLN:HG2	40:B8:44:ASP:N	2.35	0.41
27:1J:33:G:H1'	27:1J:50:G:N2	2.36	0.41
42:D8:20:LEU:HD21	42:D8:22:VAL:CG2	2.51	0.41
45:C5:62:GLU:CD	45:C5:63:LYS:H	2.23	0.41
49:K8:42:GLY:C	49:K8:44:LEU:H	2.24	0.41
26:14:1005:C:C2	26:14:1143:A:C5	3.09	0.41
31:41:12:TYR:CD1	31:41:16:ARG:HD3	2.56	0.41
26:1H:447:A:C2	26:1H:473:G:C8	3.09	0.41
27:1J:57:A:C2'	27:1J:58:A:H5'	2.51	0.41
1:13:1333:A:H2'	1:13:1334:G:O4'	2.20	0.41
32:51:42:ARG:HH11	32:51:42:ARG:HB2	1.86	0.41
47:I8:53:MET:CG	47:I8:59:LEU:HD23	2.51	0.41
26:14:1648:C:N4	26:14:2009:G:H1	2.19	0.41
5:4E:87:SER:HB3	5:4E:125:SER:O	2.20	0.41
43:E8:23:LEU:HD22	52:N8:25:LEU:HD12	2.02	0.41
6:52:14:LEU:HB2	6:52:18:GLN:OE1	2.19	0.41
1:1G:1010:G:C2	1:1G:1020:U:H1'	2.56	0.41
50:H5:6:VAL:O	50:H5:34:GLU:HA	2.20	0.41
1:13:1234:C:O2'	1:13:1235:U:H5'	2.21	0.41
26:1H:275:G:C8	26:1H:363:G:C5	3.08	0.41
26:14:693:C:H42	26:14:769:G:H1	1.69	0.41
17:8A:81:ARG:HG2	17:8A:84:LEU:HD12	2.03	0.41
26:1H:2003:G:H2'	26:1H:2004:G:O5'	2.20	0.41
14:5A:38:GLY:O	14:5A:39:LEU:HD23	2.21	0.41
26:14:1961:C:O2'	26:14:1962:C:H5'	2.21	0.41
26:14:121:G:C2	26:14:131:G:C4	3.09	0.41
26:1H:1761:C:H5''	26:1H:1762:A:OP2	2.21	0.41
30:31:46:ARG:HD2	30:31:46:ARG:HA	1.73	0.41
43:E8:90:ARG:HE	43:E8:90:ARG:HB3	1.45	0.41
42:95:20:LEU:HA	42:95:20:LEU:HD12	1.76	0.41
19:AI:71:LEU:HA	19:AI:71:LEU:HD23	1.88	0.41
46:H8:135:GLU:HG2	46:H8:135:GLU:H	1.70	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:AA:17:GLU:O	19:AA:21:GLU:HG2	2.21	0.41
26:1H:760:G:C2'	26:1H:761:A:H5'	2.51	0.41
26:1H:511:U:C5	26:1H:512:G:C5	3.09	0.41
26:1H:2704:C:H2'	26:1H:2705:A:C8	2.55	0.41
26:1H:960:A:C8	26:1H:962:G:C8	3.08	0.41
26:1H:962:G:H2'	26:1H:963:U:H6	1.86	0.41
26:1H:963:U:H2'	26:1H:964:C:C6	2.56	0.41
26:1H:818:G:H4'	26:1H:838:C:O3'	2.21	0.41
46:H8:7:ALA:HB3	46:H8:61:LEU:CB	2.51	0.41
7:62:95:ARG:NH2	7:62:99:LEU:HD11	2.35	0.41
4:3E:107:ARG:HA	4:3E:107:ARG:HD2	1.78	0.41
10:1I:57:LYS:CD	10:1I:60:ARG:HH12	2.31	0.41
26:14:531:C:C5	26:14:2035:G:C2	3.09	0.41
34:15:63:THR:HG23	34:15:63:THR:H	1.60	0.41
39:A8:36:TYR:N	39:A8:36:TYR:HD1	2.18	0.41
26:1H:1165:U:C2	26:1H:1166:C:C5	3.08	0.41
1:13:963:G:H5'	58:13:1942:HOH:O	2.20	0.41
26:14:1970:A:P	58:14:3458:HOH:O	2.74	0.41
36:35:60:MET:C	36:35:61:ARG:HG2	2.33	0.41
55:M5:32:LEU:HA	55:M5:32:LEU:HD12	1.74	0.41
26:14:137(A):G:H2'	26:14:139:G:N7	2.35	0.41
1:1G:1300:G:O2'	1:1G:1301:U:P	2.79	0.41
30:39:80:ALA:O	30:39:83:PHE:HB2	2.21	0.41
31:49:104:GLU:OE1	51:I5:23:GLU:HG2	2.21	0.41
52:N8:40:LYS:HG3	52:N8:47:PRO:HD2	2.02	0.41
34:58:65:LYS:HE3	34:58:65:LYS:HB2	1.50	0.41
9:82:65:VAL:HG21	9:82:73:GLN:HB3	2.03	0.41
24:3L:15:G:C6	24:3L:48:C:N3	2.89	0.41
24:3L:18:G:H2'	24:3L:57:G:N2	2.36	0.41
26:14:660:G:H21	36:35:12:ALA:CA	2.34	0.41
26:1H:1732:A:H2'	26:1H:1733:G:O4'	2.20	0.41
26:14:1568:G:H5'	28:19:60:ARG:HA	2.03	0.41
29:29:23:VAL:O	29:29:184:VAL:O	2.39	0.41
26:14:629:G:N1	26:14:634:C:N3	2.49	0.41
26:1H:1675:C:H2'	26:1H:1676:A:O4'	2.21	0.41
34:15:17:ASP:O	34:15:56:ASN:ND2	2.54	0.41
29:29:9:VAL:HG23	29:29:26:ILE:O	2.21	0.41
40:B8:108:ARG:O	40:B8:111:ARG:HG2	2.20	0.41
26:1H:993:G:H2'	26:1H:994:C:H6	1.86	0.41
26:1H:994:C:H5''	26:1H:995:C:OP1	2.21	0.41
26:1H:994:C:H2'	26:1H:994:C:O2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:C8:90:VAL:O	41:C8:92:ARG:N	2.54	0.41
26:1H:802:A:P	58:1H:4587:HOH:O	2.77	0.41
26:1H:1041:C:H2'	26:1H:1042:G:C8	2.56	0.41
1:1G:1002:G:H22	1:1G:1038:C:N4	2.18	0.41
38:98:52:ILE:HB	38:98:94:TYR:HD2	1.86	0.41
26:1H:174:C:H2'	26:1H:175:G:O4'	2.19	0.41
24:3L:9:A:N6	24:3L:22:G:C8	2.89	0.41
24:3L:8:U:O2'	24:3L:9:A:H5''	2.21	0.41
1:13:708:C:H2'	1:13:709:G:H8	1.86	0.41
53:O8:11:LEU:HD22	53:O8:26:ASN:HB3	2.03	0.41
55:M5:34:TRP:CD1	55:M5:35:GLN:N	2.86	0.41
7:6E:91:VAL:HG12	7:6E:95:ARG:HB3	2.02	0.41
1:13:198:G:C2	1:13:199:G:C8	3.09	0.41
16:7I:77:ALA:HB1	16:7I:79:VAL:HG23	2.03	0.41
12:3I:53:ARG:HB3	12:3I:69:TYR:HE1	1.86	0.41
26:14:740:U:H5''	26:14:1784:A:H3'	2.03	0.41
20:BI:43:LEU:HD23	20:BI:43:LEU:HA	1.92	0.41
26:1H:154:G:H2'	26:1H:155:C:C6	2.56	0.41
14:5A:41:ARG:O	14:5A:45:ARG:HB2	2.20	0.41
38:55:74:LYS:CE	38:55:77:ARG:HH21	2.31	0.41
26:14:1420:U:HO2'	26:14:1421:G:P	2.44	0.41
35:25:68:GLU:HB3	35:25:78:ARG:HB2	2.03	0.41
26:14:1599:C:C4	26:14:1600:C:C5	3.08	0.41
1:13:338:A:P	35:68:97:ARG:HH12	2.43	0.41
26:1H:931:G:C4	26:1H:933:A:C8	3.08	0.41
15:6A:58:MET:HB2	15:6A:58:MET:HE2	1.93	0.41
26:14:1409:C:H2'	26:14:1410:G:O4'	2.21	0.41
2:12:207:ALA:O	2:12:211:ILE:HG13	2.21	0.41
26:14:2893:G:H5'	26:14:2894:G:OP1	2.21	0.41
31:49:68:PRO:HB2	31:49:90:LEU:HD12	2.03	0.41
50:L8:4:LEU:O	50:L8:36:VAL:HA	2.21	0.41
2:12:5:ILE:HG12	2:12:6:THR:HG22	2.02	0.41
10:1I:8:LEU:HD22	10:1I:96:ILE:HG22	2.02	0.41
2:1E:20:GLU:HG3	2:1E:191:ASP:N	2.36	0.41
35:68:88:ASN:HD21	35:68:92:GLU:HG3	1.86	0.41
26:14:582:G:H2'	26:14:583:G:H8	1.86	0.41
31:41:61:ALA:HA	31:41:66:GLN:O	2.21	0.41
26:14:2558:C:C4	26:14:2559:C:C5	3.09	0.41
1:1G:728:A:H2'	1:1G:729:A:H8	1.83	0.41
26:1H:128:C:H2'	26:1H:129:C:C6	2.56	0.41
22:1L:23:A:H2'	22:1L:24:G:C8	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:12:219:VAL:HG23	2:12:222:ILE:HD11	2.03	0.41
1:1G:735:C:H2'	1:1G:736:C:C6	2.56	0.41
1:1G:1205:U:O2'	3:22:195:VAL:HG13	2.20	0.41
28:11:238:GLY:O	28:11:239:ARG:C	2.60	0.41
1:1G:4:U:H3'	1:1G:5:U:H5'	2.02	0.41
46:D5:155:LEU:O	46:D5:157:LEU:HD13	2.21	0.41
3:22:33:LEU:O	3:22:36:ASP:N	2.53	0.41
18:9I:26:LEU:HD12	18:9I:29:PHE:CE1	2.56	0.41
50:L8:28:LEU:HA	50:L8:33:GLN:OE1	2.21	0.41
26:1H:1465:G:C6	26:1H:1466:G:C5	3.09	0.41
26:14:234:C:H2'	26:14:235:U:H6	1.85	0.41
26:14:2850:A:C8	26:14:2869:G:O4'	2.73	0.41
34:15:127:ASP:HB3	34:15:128:HIS:H	1.59	0.41
24:3K:37:A:HO2'	24:3K:38:A:C4'	2.34	0.41
38:55:96:ARG:NH2	38:55:117:VAL:HG13	2.36	0.41
6:52:8:ILE:N	6:52:8:ILE:HD12	2.36	0.41
26:1H:117:G:C6	26:1H:119:A:C6	3.09	0.41
29:21:15:PHE:HB3	40:B8:81:PRO:HG3	2.02	0.41
33:61:99:GLU:HG2	33:61:103:ARG:NH2	2.36	0.41
11:2I:59:TYR:OH	11:2I:63:LEU:HD21	2.21	0.41
34:58:40:PRO:HB3	41:C8:67:ALA:HB3	2.03	0.41
26:1H:806:C:C2	26:1H:807:U:C5	3.09	0.41
6:5E:44:GLY:HA2	6:5E:59:TYR:CZ	2.56	0.41
29:21:31:CYS:HA	29:21:32:PRO:HD3	1.64	0.41
26:14:2355:C:H5''	26:14:2356:C:OP2	2.21	0.41
1:13:222:U:H2'	1:13:223:U:C6	2.55	0.41
28:19:37:LEU:HB2	28:19:38:LYS:HG2	2.02	0.41
1:1G:485:G:O2'	1:1G:486:U:H6	2.03	0.41
54:L5:6:GLN:HA	54:L5:7:PRO:HD3	1.81	0.41
1:13:923:A:OP1	5:4E:21:ALA:HB2	2.21	0.41
16:7I:17:TYR:CE2	16:7I:41:PRO:HG3	2.56	0.41
8:7E:81:HIS:HB2	8:7E:138:TRP:CZ3	2.56	0.41
8:7E:82:HIS:ND1	8:7E:138:TRP:CE2	2.89	0.41
20:BI:56:MET:HB3	20:BI:56:MET:HE2	1.73	0.41
8:72:78:GLN:H	8:72:78:GLN:HG2	1.47	0.41
9:82:53:VAL:C	9:82:55:ALA:H	2.24	0.41
26:14:2238:G:N3	26:14:2238:G:H2'	2.35	0.41
22:1K:2:C:H2'	22:1K:3:C:C5	2.56	0.41
15:6A:76:GLU:HA	15:6A:79:ARG:HH11	1.85	0.41
1:13:979:C:N4	58:13:1827:HOH:O	2.54	0.41
26:14:2101:G:H2'	26:14:2102:U:O4'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:8E:15:ALA:HB2	9:8E:65:VAL:HG23	2.02	0.41
26:1H:188:G:H1	26:1H:208:C:N4	2.19	0.41
26:1H:107:C:H2'	26:1H:108:U:H6	1.86	0.41
26:1H:17:G:H2'	26:1H:18:C:C6	2.56	0.41
2:1E:36:ARG:C	2:1E:38:GLY:H	2.24	0.41
26:1H:1001:A:C8	26:1H:1002:G:C8	3.09	0.41
30:31:197:ASP:N	30:31:197:ASP:OD1	2.54	0.41
26:14:2399:G:H2'	26:14:2400:G:O4'	2.21	0.41
26:1H:2356:C:H2'	26:1H:2357:U:O4'	2.20	0.41
30:39:53:THR:HG22	30:39:56:GLU:OE1	2.21	0.41
45:C5:67:LEU:HA	45:C5:67:LEU:HD12	1.67	0.41
1:1G:1104:G:O3'	2:12:111:ARG:NH2	2.53	0.41
1:1G:596:C:H2'	1:1G:597:G:H8	1.86	0.41
2:12:22:LYS:HE3	2:12:22:LYS:HB2	1.66	0.41
1:13:1091:U:C2	1:13:1095:U:N3	2.89	0.41
37:88:1:MET:HB3	37:88:2:LEU:H	1.67	0.41
28:19:264:LYS:HE2	28:19:266:SER:HB3	2.02	0.41
26:1H:234:C:C2	26:1H:235:U:C6	3.09	0.41
1:13:9:G:C2	1:13:26:A:C2	3.08	0.41
22:1K:34:G:H5''	22:1K:35:A:OP2	2.21	0.41
10:1I:33:GLN:O	10:1I:75:ILE:HG13	2.21	0.41
26:14:1716:U:O2'	26:14:1717:G:H5'	2.20	0.41
26:1H:1761:C:H42	26:1H:1762:A:H62	1.67	0.41
32:59:27:LYS:HD2	32:59:32:GLU:HG3	2.03	0.41
26:1H:2751:G:O2'	26:1H:2752:C:H5'	2.21	0.41
31:41:53:LEU:HG	31:41:90:LEU:HD21	2.02	0.41
26:14:2436:G:C5	26:14:2437:U:C5	3.09	0.41
1:1G:775:G:H2'	1:1G:776:G:O4'	2.20	0.41
1:13:1247:U:H3	1:13:1290:G:H1	1.67	0.41
17:8I:52:LYS:HD2	17:8I:55:ASP:OD1	2.21	0.41
48:J8:82:LEU:HD22	48:J8:82:LEU:H	1.86	0.41
28:19:271:ILE:HG12	28:19:271:ILE:O	2.21	0.41
12:3I:28:LYS:HD2	12:3I:28:LYS:HA	1.77	0.41
6:5E:42:GLU:O	6:5E:42:GLU:HG2	2.20	0.41
27:1J:76:G:H2'	27:1J:77:U:O4'	2.20	0.41
9:8E:13:ALA:HB2	9:8E:68:GLY:HA3	2.03	0.41
26:14:1628:G:H2'	26:14:1629:U:C6	2.56	0.41
7:6E:109:ASN:OD1	7:6E:119:ARG:NH2	2.49	0.41
26:14:540:G:C6	26:14:541:C:C4	3.09	0.41
26:1H:1870:C:H2'	26:1H:1871:A:O4'	2.21	0.41
26:14:2643:G:H2'	26:14:2644:G:O4'	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:2E:113:ALA:HB3	3:2E:114:PRO:HD3	2.02	0.41
45:C5:77:PRO:O	45:C5:78:ALA:HB3	2.21	0.41
36:35:27:HIS:HB3	36:35:32:THR:HG23	2.02	0.41
3:22:130:VAL:O	3:22:134:ILE:HG12	2.21	0.41
23:2K:19:G:C2	23:2K:59:A:C5	3.09	0.41
38:55:18:LEU:HD22	38:55:22:ARG:HG3	2.03	0.41
11:2A:40:ILE:HA	11:2A:40:ILE:HD13	1.68	0.41
38:55:44:LEU:HD23	38:55:44:LEU:HA	1.83	0.41
8:72:54:ASP:OD1	8:72:54:ASP:N	2.53	0.41
23:2K:41:C:O2'	23:2K:42:C:H5'	2.21	0.41
47:I8:41:ARG:NE	47:I8:41:ARG:HA	2.33	0.41
28:11:165:ILE:H	28:11:165:ILE:HG12	1.44	0.41
10:1A:50:ILE:H	10:1A:50:ILE:HD12	1.86	0.41
33:61:9:LEU:HA	33:61:9:LEU:HD12	1.86	0.41
26:14:536:A:H2'	26:14:537:C:C6	2.56	0.41
45:C5:15:VAL:HG12	45:C5:21:LYS:HA	2.03	0.41
32:59:92:ILE:HD11	32:59:160:LYS:NZ	2.36	0.41
26:1H:590:A:H2'	26:1H:591:C:C6	2.56	0.41
55:Q8:30:ARG:NH1	55:Q8:30:ARG:CG	2.76	0.41
46:H8:128:VAL:HG22	46:H8:129:SER:O	2.20	0.41
26:14:2720:U:C2	26:14:2721:A:C8	3.09	0.41
1:13:1124:G:H5'	10:1I:35:SER:HB2	2.02	0.41
42:95:87:HIS:HE1	42:95:89:GLN:HB2	1.83	0.41
26:1H:2702:U:C6	26:1H:2702:U:OP1	2.70	0.41
26:14:2637:U:C4	26:14:2638:G:C6	3.08	0.41
26:1H:2028:U:H2'	26:1H:2029:G:O4'	2.21	0.41
1:13:1029:G:H1'	1:13:1032(A):G:H22	1.86	0.41
30:31:78:ILE:HA	30:31:83:PHE:CE2	2.56	0.41
36:35:62:LEU:CD1	55:M5:30:ARG:HH21	2.34	0.41
26:1H:1483:G:H2'	26:1H:1484:G:H8	1.85	0.41
30:39:78:ILE:HA	30:39:83:PHE:CE2	2.52	0.41
33:61:4:ILE:HD11	33:61:44:LEU:HD12	2.03	0.41
26:14:2507:C:H2'	26:14:2508:G:O4'	2.21	0.41
22:1L:72:C:C2'	22:1L:73:A:H5''	2.51	0.41
24:3L:58:A:H1'	24:3L:60:U:C5	2.56	0.41
26:14:2298:A:H61	26:14:2318:G:H2'	1.86	0.41
26:14:2208:U:H4'	28:19:151:LYS:HG2	2.02	0.41
26:14:2116:G:C2	26:14:2117:A:N6	2.89	0.41
34:58:42:TRP:HA	34:58:48:MET:CE	2.51	0.41
23:2K:48:U:O2'	23:2K:49:C:OP2	2.26	0.41
26:14:1839:G:C8	26:14:1927:A:H1'	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:4A:29:ARG:HD3	13:4A:64:TRP:CE3	2.55	0.41
26:14:1056:G:H1'	26:14:1103:A:N6	2.36	0.41
51:M8:40:HIS:HA	51:M8:44:THR:O	2.22	0.41
1:1G:503:C:O5'	1:1G:503:C:H6	2.04	0.41
1:13:114:U:H2'	1:13:115:G:C8	2.56	0.41
16:7I:74:LEU:HD23	16:7I:74:LEU:HA	1.87	0.41
37:88:85:LYS:HG3	37:88:86:GLY:N	2.36	0.41
37:88:52:VAL:HG13	37:88:56:ARG:HG2	2.03	0.41
26:14:1188:U:H2'	26:14:1189:A:H8	1.85	0.41
2:12:19:HIS:HE1	2:12:206:ASP:HB2	1.86	0.41
1:1G:626:U:C2	1:1G:627:G:C8	3.09	0.41
26:1H:2788:C:P	29:21:61:ARG:HH22	2.43	0.41
19:AI:40:ILE:HG23	19:AI:67:VAL:O	2.21	0.41
9:82:10:ARG:HD3	9:82:75:ASP:HB3	2.03	0.41
1:13:502:G:C6	1:13:503:C:C4	3.09	0.41
16:7I:26:ARG:HE	16:7I:31:LYS:HB3	1.86	0.41
1:13:1415:G:C6	1:13:1486:G:C6	3.09	0.41
5:42:42:GLY:HA3	5:42:66:MET:HA	2.02	0.41
26:1H:943:U:OP2	58:1H:4508:HOH:O	2.21	0.41
1:1G:15:G:H2'	1:1G:16:A:H8	1.87	0.41
26:1H:638:G:H2'	26:1H:639:U:C6	2.55	0.41
1:13:1291:G:H2'	1:13:1292:U:C6	2.56	0.41
9:82:3:GLN:O	9:82:4:TYR:HD1	2.04	0.41
17:8I:48:GLU:O	17:8I:50:LYS:N	2.54	0.41
1:1G:1442:G:H2'	1:1G:1443:G:H5''	2.02	0.41
31:41:36:LYS:HG2	31:41:38:VAL:HG23	2.02	0.41
26:1H:51:G:H1'	26:1H:119:A:N1	2.36	0.41
27:1J:44:G:O2'	27:1J:48:A:N6	2.54	0.41
26:14:1537:C:H2'	26:14:1538:G:N9	2.36	0.41
25:4K:23:A:H3'	25:4K:24:A:H8	1.83	0.41
20:BA:13:LEU:HD12	20:BA:13:LEU:HA	1.70	0.41
1:13:958:A:C6	1:13:959:A:C6	3.09	0.41
27:1J:33:G:N3	27:1J:50:G:N2	2.68	0.41
1:13:587:G:C2	1:13:755:G:C5	3.09	0.41
26:1H:856:C:H5'	47:I8:27:GLU:OE2	2.21	0.41
46:H8:44:PHE:CD1	46:H8:44:PHE:C	2.94	0.41
26:14:277:C:H5''	26:14:278:A:C8	2.55	0.41
46:D5:103:ARG:NE	46:D5:136:PHE:HB3	2.36	0.41
43:A5:29:LEU:HD21	43:A5:33:ARG:HH21	1.85	0.41
37:45:70:PRO:HA	37:45:94:VAL:O	2.20	0.41
30:31:125:LEU:HA	30:31:125:LEU:HD23	1.80	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1171:G:H2'	1:13:1172:C:C6	2.55	0.41
26:1H:2473:U:O2'	26:1H:2474:C:H5'	2.21	0.41
39:A8:58:LEU:HD12	39:A8:68:GLN:OE1	2.21	0.41
28:19:149:PRO:O	28:19:150:LYS:HB2	2.21	0.41
26:14:494:G:OP1	43:A5:8:ARG:HD3	2.21	0.41
16:7I:83:GLU:HB3	16:7I:84:ALA:H	1.52	0.41
46:H8:69:THR:HA	46:H8:89:PHE:O	2.21	0.41
1:13:875:C:O2'	8:7E:14:ARG:NH1	2.54	0.41
15:6A:10:LYS:HD2	15:6A:10:LYS:HA	1.92	0.41
30:39:107:LYS:HD3	30:39:107:LYS:HA	1.56	0.41
46:D5:151:HIS:HB2	46:D5:154:ASP:CG	2.42	0.41
43:A5:57:ASN:HA	43:A5:61:ASN:HD22	1.85	0.41
36:78:49:ARG:NE	55:Q8:57:ARG:HG2	2.35	0.40
26:1H:568:U:OP1	36:78:36:LYS:HE3	2.21	0.40
2:1E:176:GLU:O	2:1E:180:LEU:HG	2.21	0.40
22:1L:18:G:O2'	22:1L:19:G:P	2.79	0.40
1:13:247:G:O6	1:13:278:G:C6	2.74	0.40
26:14:1142:U:H5''	26:14:1142(A):A:C8	2.56	0.40
26:1H:1556:C:H2'	26:1H:1557:C:H6	1.86	0.40
26:14:631:A:O2'	36:35:67:MET:HB3	2.21	0.40
32:51:6:ARG:HG2	32:51:66:GLY:CA	2.51	0.40
26:14:2207:C:O2	28:19:151:LYS:NZ	2.52	0.40
26:1H:1803:A:H61	26:1H:1814:G:H1'	1.85	0.40
1:1G:376:G:O2'	1:1G:377:G:H5'	2.21	0.40
26:14:674:G:H1'	30:39:74:ARG:HD3	2.03	0.40
15:6A:82:ILE:HD13	15:6A:88:ARG:HB2	2.01	0.40
1:1G:1306:A:C6	1:1G:1307:U:C2	3.09	0.40
1:13:1183:A:O2'	1:13:1184:G:OP1	2.32	0.40
49:G5:20:GLU:O	49:G5:24:LEU:HB2	2.21	0.40
26:1H:2075:U:C4	26:1H:2238:G:C6	3.09	0.40
40:B8:55:ASN:N	40:B8:59:THR:HB	2.35	0.40
24:3L:69:G:N1	24:3L:70:G:C6	2.89	0.40
48:J8:58:ILE:HG23	48:J8:87:PRO:HG3	2.04	0.40
23:2L:30:G:N2	23:2L:42:C:O2	2.46	0.40
39:65:15:ARG:HD2	39:65:25:ARG:HH21	1.85	0.40
1:13:1498:U:C4	25:4K:17:U:H5''	2.56	0.40
46:D5:7:ALA:O	46:D5:8:TYR:CD2	2.75	0.40
13:4I:40:ASN:HA	13:4I:41:PRO:HD3	1.96	0.40
33:61:57:ARG:O	33:61:61:ARG:HB2	2.21	0.40
1:13:726:C:H2'	1:13:727:G:H8	1.85	0.40
37:88:136:ALA:HB2	46:H8:52:SER:HB2	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:H8:25:PRO:O	46:H8:85:HIS:HA	2.21	0.40
1:1G:167:G:C4	1:1G:168:G:C8	3.09	0.40
15:6I:26:GLU:H	15:6I:26:GLU:HG2	1.47	0.40
26:14:2439:A:H4'	26:14:2440:C:O5'	2.20	0.40
39:A8:67:ARG:O	39:A8:71:ARG:HG3	2.21	0.40
33:61:63:ALA:O	33:61:67:ARG:HD3	2.21	0.40
30:39:170:LEU:HA	30:39:171:PRO:HD3	1.92	0.40
2:1E:97:TRP:CZ3	2:1E:172:ILE:HD12	2.57	0.40
21:1B:6:ARG:HD3	21:1B:15:ARG:NH2	2.36	0.40
26:14:686:G:C8	54:L5:7:PRO:HA	2.57	0.40
26:14:686:G:O6	54:L5:12:ARG:HG3	2.21	0.40
40:B8:62:THR:CG2	40:B8:75:ILE:HG12	2.50	0.40
26:1H:104:U:C5	26:1H:105:C:C5	3.09	0.40
26:1H:1523:U:H2'	26:1H:1524:G:O4'	2.21	0.40
1:1G:38:G:N2	1:1G:397:A:C4	2.90	0.40
37:45:69:PHE:HA	37:45:70:PRO:HD2	1.68	0.40
43:E8:29:LEU:O	43:E8:29:LEU:HD12	2.20	0.40
26:1H:754:C:H4'	26:1H:1272:A:N6	2.37	0.40
26:1H:2744:G:N2	32:51:143:GLN:OE1	2.53	0.40
21:1F:6:ARG:O	21:1F:12:LYS:HD2	2.21	0.40
26:1H:826:U:H2'	26:1H:828:U:O4'	2.21	0.40
23:2L:14:A:H2'	23:2L:15:G:H5'	2.02	0.40
1:1G:1084:G:H5'	1:1G:1102:A:OP2	2.21	0.40
46:D5:114:GLY:HA2	46:D5:174:VAL:HG11	2.04	0.40
4:32:39:PRO:O	4:32:44:GLY:HA3	2.21	0.40
26:14:430:G:H5''	26:14:431:U:OP2	2.21	0.40
6:5E:18:GLN:O	6:5E:21:LEU:HB2	2.21	0.40
9:82:83:ARG:O	9:82:87:GLN:HB2	2.20	0.40
34:58:111:PRO:HA	34:58:114:ARG:NH1	2.35	0.40
26:1H:858:U:O2	26:1H:2268:A:H2'	2.21	0.40
1:1G:1137:C:H1'	1:1G:1138:G:C2	2.56	0.40
32:59:94:TYR:CD1	32:59:94:TYR:N	2.89	0.40
20:BA:58:LYS:HD3	20:BA:58:LYS:O	2.21	0.40
43:E8:9:TYR:CD1	43:E8:102:HIS:HE1	2.39	0.40
26:1H:1225:C:O2'	42:D8:85:LYS:HA	2.21	0.40
26:14:675:A:C8	26:14:804:A:C6	3.09	0.40
26:14:1614:A:H2	58:14:3409:HOH:O	2.02	0.40
10:1I:5:ARG:HB2	10:1I:73:ASP:OD1	2.22	0.40
26:1H:1640:C:H2'	26:1H:1641:A:C8	2.57	0.40
26:1H:1784:A:H4'	26:1H:1785:A:O5'	2.22	0.40
26:14:1019:U:O2'	26:14:1021:A:H2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:1J:13:A:O2'	27:1J:14:U:H3'	2.20	0.40
26:1H:2035:G:P	58:1H:3793:HOH:O	2.74	0.40
1:1G:1141:C:C2	1:1G:1142:G:C8	3.09	0.40
36:35:39:LYS:CB	36:35:45:LEU:HD21	2.51	0.40
1:13:868:C:H2'	1:13:869:G:O4'	2.20	0.40
1:13:658:G:C6	1:13:659:U:C4	3.08	0.40
12:3I:113:ARG:O	12:3I:114:LYS:HD2	2.21	0.40
29:29:5:LEU:HB2	29:29:51:PHE:CG	2.56	0.40
46:H8:126:VAL:HA	46:H8:164:ALA:N	2.36	0.40
1:1G:1310:G:C5'	13:4A:77:ASN:HD21	2.27	0.40
1:13:465:A:N7	1:13:467:G:C5	2.89	0.40
26:14:2287:A:C4	26:14:2289:G:C8	3.09	0.40
1:13:272:C:C2	1:13:273:A:C8	3.08	0.40
29:29:26:ILE:HG23	29:29:26:ILE:HD12	1.74	0.40
1:13:1285:A:H4'	1:13:1286:A:O5'	2.21	0.40
48:J8:91:LYS:HA	48:J8:91:LYS:HZ3	1.86	0.40
10:1A:33:GLN:HB3	10:1A:75:ILE:HG12	2.04	0.40
38:98:96:ARG:NH2	38:98:118:GLU:H	2.20	0.40
32:51:88:LEU:HB3	32:51:130:ARG:HG2	2.03	0.40
1:1G:1329:A:H5''	13:4A:25:ILE:O	2.21	0.40
26:14:2327:A:H2'	26:14:2328:A:H8	1.81	0.40
53:O8:15:GLU:CG	53:O8:16:CYS:N	2.84	0.40
8:72:34:GLU:HB3	8:72:118:VAL:HG21	2.04	0.40
1:13:338:A:C6	1:13:339:C:C4	3.09	0.40
1:13:940:C:H2'	1:13:941:G:C8	2.56	0.40
1:1G:1055:A:N6	1:1G:1056:U:O2	2.55	0.40
10:1I:76:ASN:HA	10:1I:77:PRO:HD2	1.92	0.40
26:1H:637:A:O5'	36:78:116:GLY:HA3	2.22	0.40
55:M5:59:LYS:C	55:M5:60:LEU:HG	2.41	0.40
26:14:1668:A:H4'	26:14:1669:A:O5'	2.21	0.40
1:13:1503:A:O2'	25:4K:13:A:C6	2.67	0.40
1:1G:109:A:H2'	1:1G:326:G:N2	2.37	0.40
5:42:76:ILE:O	5:42:93:PRO:HB3	2.21	0.40
9:8E:20:ARG:HA	9:8E:21:PRO:HD3	1.95	0.40
26:14:144:C:C2	26:14:145:G:C8	3.09	0.40
34:15:16:ILE:HG21	34:15:26:LEU:HD11	2.03	0.40
26:14:2881:C:C4	26:14:2882:A:N7	2.89	0.40
42:95:75:PHE:HE1	42:95:81:TYR:CE1	2.38	0.40
42:95:80:GLN:CD	42:95:81:TYR:H	2.24	0.40
1:1G:526:C:C4	1:1G:527:G:H1'	2.56	0.40
6:5E:97:PHE:N	18:9I:30:ASP:OD1	2.53	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:827:U:H4'	26:14:828:U:O2	2.21	0.40
28:11:69:ARG:NH2	28:11:130:ALA:H	2.19	0.40
9:8E:10:ARG:NE	9:8E:105:ASP:OD2	2.54	0.40
26:14:2660:A:H2'	26:14:2661:G:O4'	2.21	0.40
26:1H:2244:U:H2'	26:1H:2245:U:O4'	2.21	0.40
26:1H:309:G:O2'	26:1H:329:G:C8	2.73	0.40
26:14:2607:G:H2'	26:14:2608:G:O4'	2.21	0.40
26:1H:447:A:C4	26:1H:473:G:N7	2.90	0.40
26:14:2187:G:H2'	26:14:2188:C:O4'	2.20	0.40
34:15:121:LYS:HB3	34:15:123:TYR:HE1	1.86	0.40
50:H5:17:LYS:NZ	50:H5:21:ALA:HB2	2.35	0.40
26:1H:15:G:C4	26:1H:16:G:C8	3.09	0.40
20:BA:56:MET:HE2	20:BA:56:MET:HB3	2.00	0.40
26:1H:46:C:OP2	26:1H:215:G:H2'	2.21	0.40
26:14:513:A:C2	26:14:514:A:C8	3.10	0.40
30:31:206:ILE:HG21	30:31:206:ILE:HD13	1.86	0.40
44:B5:26:TYR:OH	44:B5:88:LYS:HB2	2.21	0.40
16:7I:15:PRO:HD2	16:7I:42:ARG:HD3	2.03	0.40
41:85:91:ASP:OD1	41:85:96:ALA:HB2	2.22	0.40
28:19:264:LYS:HA	28:19:265:PRO:HD2	1.95	0.40
36:35:124:LYS:HA	36:35:143:GLY:O	2.22	0.40
1:13:1170:A:C8	1:13:1171:G:C8	3.09	0.40
11:2A:40:ILE:HG23	11:2A:75:TYR:CD2	2.56	0.40
26:14:2303:G:O2'	31:49:132:ASN:ND2	2.48	0.40
26:1H:1663:C:O2'	26:1H:2686:G:H4'	2.21	0.40
1:1G:328:C:H4'	1:1G:329:A:C5'	2.51	0.40
33:69:62:LYS:HE2	33:69:133:HIS:ND1	2.35	0.40
35:25:76:ALA:HB3	40:75:75:ILE:HD12	2.03	0.40
26:1H:1612:C:O3'	54:P8:5:TRP:HB3	2.21	0.40
27:1J:68:C:O5'	27:1J:68:C:H6	2.04	0.40
33:69:99:GLU:HG3	33:69:99:GLU:H	1.60	0.40
26:1H:14:A:O5'	26:1H:14:A:H8	2.02	0.40
31:41:181:ARG:HH11	31:41:181:ARG:HB3	1.86	0.40
48:J8:11:ARG:HH11	48:J8:11:ARG:HD2	1.75	0.40
26:1H:489:G:H2'	26:1H:489:G:N3	2.36	0.40
38:55:28:LEU:HD23	38:55:28:LEU:HA	1.78	0.40
26:14:1249:U:H2'	26:14:1249:U:O2	2.21	0.40
23:2L:38:A:H2'	23:2L:39:A:O4'	2.21	0.40
1:1G:407:G:C2	1:1G:436:C:C2	3.09	0.40
4:32:18:LYS:HG2	4:32:20:TYR:CD1	2.57	0.40
26:14:1109:C:H2'	26:14:1110:G:N9	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:13:1133:G:H2'	1:13:1134:G:C8	2.56	0.40
42:95:71:LEU:HA	42:95:71:LEU:HD13	1.77	0.40
37:88:77:LYS:HA	37:88:78:PRO:HD3	1.92	0.40
49:K8:51:ARG:NH1	49:K8:55:ARG:HH12	2.19	0.40
36:35:85:LEU:HA	36:35:88:LEU:CB	2.41	0.40
40:B8:9:LEU:HD23	40:B8:9:LEU:HA	1.94	0.40
1:1G:861:G:N2	1:1G:872:A:H2	2.18	0.40
26:14:2582:G:C2	26:14:2583:G:C8	3.09	0.40
26:14:2782:G:OP2	58:14:3788:HOH:O	2.22	0.40
55:Q8:33:ASN:HA	55:Q8:34:TRP:CD1	2.57	0.40
26:14:1299:G:OP1	58:14:4092:HOH:O	2.22	0.40
16:7A:50:LYS:HG2	16:7A:51:VAL:N	2.36	0.40
26:14:1480:G:C6	26:14:1482:U:N3	2.90	0.40
26:1H:2053:G:N2	26:1H:2054:A:C4	2.90	0.40
7:6E:95:ARG:HE	7:6E:95:ARG:HB3	1.69	0.40
1:13:291:C:C2	1:13:310:G:C2	3.10	0.40
12:3I:53:ARG:HG3	12:3I:93:LEU:HD21	2.02	0.40
26:14:1846:G:H5''	26:14:1847:A:OP2	2.22	0.40
1:13:435:C:H2'	1:13:436:C:C6	2.55	0.40
1:13:339:C:H2'	1:13:340:U:C6	2.56	0.40
42:95:15:GLU:O	42:95:96:ILE:HB	2.21	0.40
1:13:324:G:N2	1:13:326:G:H3'	2.37	0.40
3:2E:40:ARG:HG3	3:2E:40:ARG:NH1	2.36	0.40
3:2E:19:GLU:HG3	3:2E:54:ARG:CZ	2.51	0.40
29:29:171:GLU:OE2	29:29:185:LYS:HE2	2.21	0.40
1:1G:1226:C:H2'	13:4A:103:THR:HB	2.03	0.40
26:1H:2829:C:C2'	26:1H:2830:G:H5''	2.52	0.40
26:1H:230:U:OP2	26:1H:230:U:H6	2.03	0.40
46:D5:126:VAL:HG12	46:D5:163:LEU:HB3	2.03	0.40
1:13:721:G:N1	1:13:733:A:C2	2.89	0.40
26:14:353:G:H2'	26:14:354:G:H8	1.83	0.40
26:14:1467:C:H2'	26:14:1468:C:C6	2.52	0.40
26:1H:1110:G:O2'	26:1H:1111:A:C8	2.71	0.40
26:14:2493:U:H2'	26:14:2494:G:O4'	2.21	0.40
1:13:1312:G:H5'	19:AI:6:LYS:HD2	2.03	0.40
6:52:10:LEU:HB2	6:52:59:TYR:HB3	2.03	0.40
3:2E:8:ILE:HG22	3:2E:9:GLY:N	2.37	0.40
1:1G:983:A:N3	1:1G:983:A:H3'	2.35	0.40
28:19:181:GLU:HA	28:19:272:ALA:HB1	2.02	0.40
26:14:1826:G:H2'	26:14:1827:C:H6	1.85	0.40
8:72:17:THR:HG22	8:72:78:GLN:HE21	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:14:1190:G:O2'	26:14:1191:G:H5'	2.21	0.40
26:1H:2117:A:H2'	26:1H:2118:U:C6	2.56	0.40
27:1J:57:A:H8	27:1J:57:A:O5'	2.03	0.40
2:1E:200:ILE:O	2:1E:201:ILE:HD13	2.22	0.40
26:1H:749:C:C4	26:1H:1618:A:C2	3.09	0.40
26:14:384:U:H2'	26:14:385:C:C6	2.55	0.40
51:M8:15:ILE:HB	51:M8:32:TYR:HD1	1.86	0.40
34:58:10:GLU:HA	34:58:11:PRO:HD3	1.96	0.40
26:14:455:C:N3	26:14:473:G:H5'	2.36	0.40
15:6I:4:THR:OG1	15:6I:7:GLU:HB2	2.21	0.40
48:F5:97:LEU:O	48:F5:98:LEU:HD23	2.21	0.40
26:14:2679:A:H2'	26:14:2680:C:O4'	2.21	0.40
13:4I:77:ASN:OD1	51:M8:55:ARG:HD2	2.21	0.40
7:62:150:ALA:HB2	11:2A:50:TYR:OH	2.22	0.40
43:A5:12:ILE:HG13	43:A5:42:ARG:NH1	2.37	0.40
26:14:415:A:H2'	26:14:416:C:C6	2.57	0.40
26:14:2288:A:H8	26:14:2288:A:O5'	2.03	0.40
26:1H:792:G:OP1	26:1H:792:G:H3'	2.21	0.40
1:1G:1354:C:H6	1:1G:1354:C:O5'	2.04	0.40
28:11:133:LEU:HD23	28:11:133:LEU:HA	1.87	0.40
8:7E:75:ARG:HE	8:7E:75:ARG:HB2	1.48	0.40
3:2E:62:ASP:N	3:2E:62:ASP:OD1	2.53	0.40
20:BI:16:HIS:O	20:BI:19:SER:HB2	2.21	0.40
1:13:1111:A:N1	3:2E:177:THR:HG23	2.36	0.40
26:14:363:G:H2'	26:14:363(A):A:O4'	2.20	0.40
45:C5:75:ILE:O	45:C5:76:CYS:HB3	2.21	0.40
26:1H:374:A:C8	26:1H:375:C:C5	3.10	0.40
7:62:94:ARG:O	7:62:97:GLN:HB3	2.22	0.40
26:1H:910:A:C6	26:1H:911:A:C6	3.09	0.40
1:13:1006:C:H2'	1:13:1007:C:H5''	2.03	0.40
1:13:1024:G:H2'	1:13:1025:U:C6	2.56	0.40
26:1H:248:G:C2	26:1H:2431:U:H4'	2.56	0.40
26:14:2331:G:O3'	47:E5:43:THR:HG22	2.21	0.40
26:1H:252:G:P	36:78:50:ARG:HH12	2.44	0.40
26:14:2086:U:H2'	26:14:2087:G:C8	2.56	0.40
16:7I:82:GLN:H	16:7I:82:GLN:HG2	1.65	0.40
26:1H:1993:U:H4'	29:21:128:SER:HB2	2.03	0.40
23:2K:9:G:H21	23:2K:47:7MG:C5'	2.34	0.40
32:51:124:GLU:O	32:51:126:PRO:HD3	2.22	0.40
9:8E:112:LYS:HD3	9:8E:117:HIS:O	2.22	0.40
26:14:2133:G:O2'	26:14:2158:A:N6	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:45:85:LYS:HG2	37:45:86:GLY:N	2.32	0.40
1:13:1517:G:H1'	26:1H:1919:A:O3'	2.21	0.40
2:1E:88:ALA:HB2	2:1E:219:VAL:HG13	2.04	0.40
35:68:106:LEU:HD23	35:68:106:LEU:HA	1.67	0.40
4:32:8:VAL:HG12	4:32:21:LEU:HD12	2.03	0.40
29:29:88:GLY:HA2	29:29:89:ASP:HA	1.15	0.40
1:1G:1054:C:O2'	1:1G:1055:A:O5'	2.38	0.40
1:13:67:C:H2'	1:13:68:G:C8	2.57	0.40
26:14:1952:A:H5''	26:14:1953:A:OP2	2.21	0.40
35:25:16:ALA:HB2	35:25:52:VAL:CG1	2.51	0.40
41:C8:69:CYS:HB3	41:C8:74:LEU:CD1	2.50	0.40
26:1H:1831:G:C4	26:1H:1975:G:N2	2.89	0.40
4:3E:89:THR:N	4:3E:92:VAL:HG23	2.36	0.40
12:3A:69:TYR:CG	12:3A:90:VAL:HG21	2.56	0.40
3:2E:150:LYS:HD3	3:2E:152:ILE:HD11	2.04	0.40
26:14:1096:A:C5	26:14:1097:U:H1'	2.56	0.40
1:13:1396:A:O4'	1:13:1398:A:H1'	2.21	0.40
1:13:669:U:N3	1:13:670:G:N7	2.69	0.40
26:14:2495:G:H2'	26:14:2496:C:C6	2.55	0.40
26:14:2520:C:H41	26:14:2542:A:N6	2.19	0.40
20:BA:97:ALA:HA	20:BA:98:PRO:HD3	1.81	0.40
28:19:267:SER:C	28:19:269:PHE:H	2.23	0.40
2:1E:17:PHE:HB2	2:1E:42:ILE:HG22	2.03	0.40
26:14:370:G:H8	26:14:370:G:O5'	2.04	0.40
2:12:119:GLU:HA	2:12:122:PHE:CD2	2.57	0.40
1:13:883:C:O2'	1:13:884:U:H5'	2.20	0.40
37:88:17:LEU:HD13	37:88:39:PRO:HB2	2.03	0.40
31:41:16:ARG:NH1	31:41:31:VAL:HG22	2.37	0.40
26:14:2586:C:H2'	26:14:2587:A:H5'	2.03	0.40
36:78:135:LEU:HA	36:78:135:LEU:HD23	1.73	0.40
1:1G:25:C:H2'	1:1G:26:A:C8	2.55	0.40
1:1G:878:G:H5'	8:72:89:PRO:HG2	2.04	0.40
26:1H:2744:G:C8	26:1H:2755:C:C6	3.10	0.40
26:14:265:A:H1'	26:14:266:G:O4'	2.22	0.40
18:9I:73:ALA:HB3	18:9I:79:LEU:HD12	2.04	0.40
26:14:2124:G:H2'	26:14:2125:G:H5'	2.03	0.40
26:14:1562:A:H2'	26:14:1563:G:O4'	2.21	0.40
3:2E:195:VAL:C	3:2E:196:LEU:HD23	2.42	0.40
1:1G:641:U:O3'	1:1G:642:A:H8	2.05	0.40
4:32:199:ASN:C	4:32:201:GLN:H	2.25	0.40
1:1G:1192:C:N4	1:1G:1193:G:C2	2.90	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1G:233:C:H2'	1:1G:234:C:H6	1.87	0.40
35:25:119:PRO:HB2	40:75:68:TYR:CE2	2.56	0.40
20:BA:59:ALA:O	20:BA:62:LEU:N	2.55	0.40
2:1E:135:GLN:O	2:1E:139:LYS:HB2	2.22	0.40
26:1H:2870:C:H5''	38:98:65:LEU:HD21	2.02	0.40
35:25:7:TYR:HE1	35:25:20:MET:CE	2.35	0.40
26:1H:55:G:C2	26:1H:116:C:C2	3.09	0.40
3:2E:87:LEU:O	3:2E:90:GLU:N	2.54	0.40
32:59:125:VAL:HG22	32:59:126:PRO:HA	2.04	0.40
18:9A:19:LYS:HG3	18:9A:20:ALA:H	1.87	0.40
2:12:171:ALA:HA	2:12:174:VAL:HG23	2.04	0.40
1:13:12:U:O2'	1:13:526:C:H4'	2.22	0.40
1:1G:1413:A:H2'	1:1G:1414:U:O4'	2.21	0.40
1:13:1467:G:H8	1:13:1467:G:O5'	2.04	0.40
43:A5:65:LEU:HA	43:A5:65:LEU:HD23	1.94	0.40
41:85:57:PHE:HD1	41:85:57:PHE:HA	1.81	0.40
26:1H:1791:A:H4'	28:11:206:LEU:HB2	2.04	0.40
37:45:59:ARG:O	37:45:60:ARG:HG3	2.22	0.40
27:16:12:C:N3	47:18:74:ARG:NH1	2.70	0.40
1:13:277:C:H2'	1:13:278:G:H8	1.87	0.40
26:1H:880:G:H22	26:1H:897:C:N4	2.20	0.40
9:8E:97:LYS:HD2	9:8E:97:LYS:HA	1.91	0.40
26:1H:95:G:O2'	49:K8:48:HIS:HB3	2.22	0.40
49:G5:27:GLU:O	49:G5:31:GLU:HG3	2.22	0.40
1:1G:474:G:H2'	1:1G:475:G:C8	2.56	0.40
26:14:2298:A:N6	26:14:2318:G:H2'	2.36	0.40
26:14:2705:A:H2'	26:14:2706:G:O4'	2.21	0.40
39:A8:25:ARG:HH12	39:A8:42:ASP:CG	2.25	0.40
45:C5:42:VAL:O	45:C5:65:ALA:N	2.45	0.40
1:1G:520:A:OP2	12:3A:51:ALA:HB1	2.22	0.40
26:14:172:C:H2'	26:14:173:G:C8	2.54	0.40
1:1G:649:G:H2'	1:1G:650:G:C8	2.54	0.40
1:13:1118:C:O4'	1:13:1179:A:H1'	2.21	0.40
24:3K:21:A:H61	24:3K:46:G:H2'	1.86	0.40
4:32:153:ARG:HA	4:32:153:ARG:HD3	1.72	0.40
1:13:750:G:C2	1:13:751:U:C5	3.09	0.40
52:J5:4:HIS:O	52:J5:6:VAL:HG23	2.21	0.40
26:14:2766:G:C2	26:14:2767:C:C6	3.09	0.40
30:31:101:LEU:HA	30:31:101:LEU:HD22	1.64	0.40
26:14:1954:G:N3	26:14:2551:C:H5''	2.37	0.40
9:82:47:LEU:HB2	9:82:50:LEU:HD12	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:1H:123:G:O6	26:1H:128:C:N3	2.55	0.40
1:1G:1182:G:H5'	1:1G:1183:A:C5'	2.48	0.40
28:19:206:LEU:HD23	28:19:206:LEU:HA	1.80	0.40
26:14:2862:G:C6	26:14:2863:C:C4	3.09	0.40
44:F8:89:ILE:O	44:F8:93:GLU:HG2	2.22	0.40
9:82:99:LEU:HB3	9:82:101:PHE:HD1	1.82	0.40
3:22:155:GLY:O	3:22:157:ILE:HG13	2.22	0.40
36:78:122:PRO:HA	36:78:142:GLY:N	2.36	0.40
26:14:1519:G:C6	26:14:1520:U:C4	3.09	0.40
38:98:2:ARG:O	38:98:5:LYS:HB2	2.22	0.40
40:B8:85:LYS:HE3	40:B8:87:ASP:OD2	2.20	0.40
26:14:1432:C:H2'	26:14:1433:U:O4'	2.22	0.40
47:E5:27:GLU:HG3	47:E5:68:GLU:HA	2.03	0.40
26:1H:1519:G:C2'	26:1H:1520:U:H5'	2.52	0.40
26:14:29:U:O2'	26:14:30:G:H5'	2.21	0.40
33:69:111:PRO:O	33:69:112:LYS:C	2.60	0.40
1:13:814:A:C8	1:13:816:A:C4	3.09	0.40
26:14:1149:G:C2	26:14:1150:C:N3	2.89	0.40
8:7E:39:LEU:HB3	8:7E:45:ILE:CD1	2.52	0.40
26:1H:598:G:H2'	26:1H:599:G:O4'	2.21	0.40
26:1H:2335:A:N7	26:1H:2337:G:C5	2.90	0.40
26:14:827:U:O2	26:14:2246:G:H4'	2.21	0.40
29:29:37:ARG:HD2	29:29:80:GLU:OE2	2.21	0.40
26:14:2244:U:O5'	26:14:2244:U:H6	2.04	0.40
1:1G:1122:U:N3	1:1G:1123:A:N7	2.69	0.40
28:19:182:LEU:H	28:19:272:ALA:CB	2.34	0.40
2:12:101:MET:O	2:12:105:PHE:HB2	2.22	0.40
7:6E:45:ASP:OD1	7:6E:48:LYS:NZ	2.52	0.40
31:49:122:PRO:O	31:49:125:PHE:HD2	2.04	0.40
9:82:53:VAL:N	9:82:95:LYS:HZ1	2.20	0.40
26:1H:1087:G:O6	26:1H:1089:G:N2	2.54	0.40
33:69:58:LEU:O	33:69:61:ARG:N	2.54	0.40
26:1H:2082:A:H2'	26:1H:2083:G:O4'	2.20	0.40
46:H8:100:VAL:HG21	46:H8:134:PRO:HG2	2.03	0.40
26:1H:1972:A:H2'	26:1H:1973:G:C8	2.56	0.40
14:5I:47:LEU:HA	14:5I:47:LEU:HD23	1.84	0.40
26:14:1260:G:H2'	26:14:1261:C:C6	2.56	0.40
26:1H:2549:G:H5''	26:1H:2549:G:C8	2.57	0.40
46:D5:24:LEU:HA	46:D5:25:PRO:HD3	1.96	0.40
20:BA:84:LEU:HD23	20:BA:84:LEU:HA	1.92	0.40
26:1H:533:G:H5'	41:C8:24:TYR:CE1	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:M8:42:PHE:O	51:M8:43:TYR:HB3	2.21	0.40
46:D5:104:PHE:O	46:D5:105:VAL:HB	2.21	0.40
1:1G:147:G:C2	1:1G:148:G:C8	3.09	0.40
13:4A:4:ILE:HG12	13:4A:5:ALA:N	2.34	0.40
12:3A:71:PRO:O	12:3A:102:ARG:NH1	2.50	0.40
38:98:81:ASP:O	38:98:85:PRO:HG2	2.22	0.40
26:14:1032:A:N1	26:14:1122:G:O6	2.54	0.40
29:29:152:LYS:HG3	34:15:77:GLY:O	2.21	0.40
26:1H:270(O):U:H3	33:61:52:ARG:HH21	1.69	0.40
5:4E:37:ARG:HH12	5:4E:111:GLU:HB3	1.87	0.40
1:1G:1246:C:H2'	1:1G:1247:U:H6	1.86	0.40
26:1H:2726:U:O2	26:1H:2726:U:H5'	2.21	0.40
39:65:39:ILE:HA	39:65:39:ILE:HD13	1.91	0.40
42:D8:83:ARG:HD3	42:D8:83:ARG:HA	1.82	0.40
26:1H:336:C:H5''	45:G8:6:HIS:CD2	2.55	0.40
43:A5:75:TYR:CZ	43:A5:104:THR:HG21	2.56	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
49:K8:17:SER:OG	26:14:307:G:OP1[2_564]	1.99	0.21
26:1H:277:C:O2'	49:G5:49:LYS:NZ[2_564]	2.13	0.07

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	12	235/256 (92%)	194 (83%)	38 (16%)	3 (1%)	15	57
2	1E	235/256 (92%)	198 (84%)	35 (15%)	2 (1%)	21	65
3	22	204/239 (85%)	180 (88%)	24 (12%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	2E	203/239 (85%)	179 (88%)	23 (11%)	1 (0%)	34	76
4	32	206/209 (99%)	181 (88%)	24 (12%)	1 (0%)	34	76
4	3E	206/209 (99%)	190 (92%)	14 (7%)	2 (1%)	19	63
5	42	149/162 (92%)	141 (95%)	8 (5%)	0	100	100
5	4E	149/162 (92%)	139 (93%)	9 (6%)	1 (1%)	26	71
6	52	99/101 (98%)	95 (96%)	4 (4%)	0	100	100
6	5E	99/101 (98%)	91 (92%)	8 (8%)	0	100	100
7	62	153/156 (98%)	146 (95%)	6 (4%)	1 (1%)	26	71
7	6E	153/156 (98%)	142 (93%)	11 (7%)	0	100	100
8	72	136/138 (99%)	126 (93%)	8 (6%)	2 (2%)	13	53
8	7E	136/138 (99%)	124 (91%)	11 (8%)	1 (1%)	26	71
9	82	125/128 (98%)	111 (89%)	13 (10%)	1 (1%)	24	67
9	8E	125/128 (98%)	107 (86%)	18 (14%)	0	100	100
10	1A	97/105 (92%)	91 (94%)	6 (6%)	0	100	100
10	1I	97/105 (92%)	90 (93%)	7 (7%)	0	100	100
11	2A	117/129 (91%)	102 (87%)	12 (10%)	3 (3%)	7	38
11	2I	117/129 (91%)	103 (88%)	13 (11%)	1 (1%)	21	65
12	3A	123/132 (93%)	100 (81%)	18 (15%)	5 (4%)	3	25
12	3I	123/132 (93%)	105 (85%)	18 (15%)	0	100	100
13	4A	115/126 (91%)	94 (82%)	19 (16%)	2 (2%)	11	50
13	4I	116/126 (92%)	94 (81%)	22 (19%)	0	100	100
14	5A	56/61 (92%)	47 (84%)	8 (14%)	1 (2%)	11	49
14	5I	58/61 (95%)	47 (81%)	9 (16%)	2 (3%)	5	29
15	6A	86/89 (97%)	82 (95%)	4 (5%)	0	100	100
15	6I	86/89 (97%)	77 (90%)	9 (10%)	0	100	100
16	7A	82/88 (93%)	77 (94%)	5 (6%)	0	100	100
16	7I	82/88 (93%)	75 (92%)	6 (7%)	1 (1%)	16	58
17	8A	98/105 (93%)	91 (93%)	7 (7%)	0	100	100
17	8I	98/105 (93%)	93 (95%)	5 (5%)	0	100	100
18	9A	70/88 (80%)	65 (93%)	5 (7%)	0	100	100
18	9I	70/88 (80%)	62 (89%)	7 (10%)	1 (1%)	14	55

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
19	AA	76/93 (82%)	61 (80%)	13 (17%)	2 (3%)	7	38
19	AI	79/93 (85%)	65 (82%)	10 (13%)	4 (5%)	2	19
20	BA	97/106 (92%)	84 (87%)	11 (11%)	2 (2%)	9	44
20	BI	97/106 (92%)	82 (84%)	14 (14%)	1 (1%)	19	63
21	1B	23/27 (85%)	21 (91%)	2 (9%)	0	100	100
21	1F	23/27 (85%)	21 (91%)	2 (9%)	0	100	100
28	11	271/276 (98%)	252 (93%)	14 (5%)	5 (2%)	11	49
28	19	271/276 (98%)	252 (93%)	14 (5%)	5 (2%)	11	49
29	21	203/206 (98%)	161 (79%)	30 (15%)	12 (6%)	2	15
29	29	203/206 (98%)	156 (77%)	35 (17%)	12 (6%)	2	15
30	31	200/210 (95%)	183 (92%)	15 (8%)	2 (1%)	19	63
30	39	206/210 (98%)	159 (77%)	41 (20%)	6 (3%)	6	35
31	41	179/182 (98%)	159 (89%)	17 (10%)	3 (2%)	11	50
31	49	179/182 (98%)	158 (88%)	20 (11%)	1 (1%)	30	74
32	51	172/180 (96%)	148 (86%)	17 (10%)	7 (4%)	3	25
32	59	168/180 (93%)	129 (77%)	34 (20%)	5 (3%)	5	34
33	61	144/148 (97%)	117 (81%)	23 (16%)	4 (3%)	6	36
33	69	144/148 (97%)	119 (83%)	21 (15%)	4 (3%)	6	36
34	15	136/140 (97%)	124 (91%)	11 (8%)	1 (1%)	26	71
34	58	136/140 (97%)	116 (85%)	16 (12%)	4 (3%)	6	35
35	25	120/122 (98%)	111 (92%)	9 (8%)	0	100	100
35	68	120/122 (98%)	111 (92%)	8 (7%)	1 (1%)	24	67
36	35	148/150 (99%)	114 (77%)	28 (19%)	6 (4%)	3	25
36	78	148/150 (99%)	121 (82%)	21 (14%)	6 (4%)	3	25
37	45	139/141 (99%)	114 (82%)	23 (16%)	2 (1%)	14	55
37	88	134/141 (95%)	110 (82%)	18 (13%)	6 (4%)	3	22
38	55	115/118 (98%)	104 (90%)	9 (8%)	2 (2%)	11	50
38	98	116/118 (98%)	97 (84%)	18 (16%)	1 (1%)	21	65
39	65	109/112 (97%)	87 (80%)	18 (16%)	4 (4%)	4	27
39	A8	109/112 (97%)	90 (83%)	18 (16%)	1 (1%)	21	65
40	75	135/146 (92%)	119 (88%)	15 (11%)	1 (1%)	26	71

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
40	B8	135/146 (92%)	122 (90%)	13 (10%)	0	100	100
41	85	115/118 (98%)	99 (86%)	15 (13%)	1 (1%)	21	65
41	C8	115/118 (98%)	101 (88%)	10 (9%)	4 (4%)	4	29
42	95	99/101 (98%)	78 (79%)	16 (16%)	5 (5%)	2	19
42	D8	99/101 (98%)	94 (95%)	3 (3%)	2 (2%)	9	46
43	A5	111/113 (98%)	100 (90%)	11 (10%)	0	100	100
43	E8	111/113 (98%)	101 (91%)	10 (9%)	0	100	100
44	B5	91/96 (95%)	82 (90%)	8 (9%)	1 (1%)	17	61
44	F8	92/96 (96%)	84 (91%)	7 (8%)	1 (1%)	17	61
45	C5	102/110 (93%)	75 (74%)	22 (22%)	5 (5%)	3	20
45	G8	102/110 (93%)	82 (80%)	14 (14%)	6 (6%)	2	15
46	D5	177/206 (86%)	130 (73%)	37 (21%)	10 (6%)	2	17
46	H8	173/206 (84%)	136 (79%)	28 (16%)	9 (5%)	2	18
47	E5	75/85 (88%)	66 (88%)	8 (11%)	1 (1%)	15	57
47	I8	78/85 (92%)	65 (83%)	12 (15%)	1 (1%)	15	57
48	F5	95/98 (97%)	88 (93%)	6 (6%)	1 (1%)	17	61
48	J8	95/98 (97%)	86 (90%)	7 (7%)	2 (2%)	9	44
49	G5	64/72 (89%)	59 (92%)	3 (5%)	2 (3%)	5	32
49	K8	65/72 (90%)	60 (92%)	3 (5%)	2 (3%)	5	32
50	H5	57/60 (95%)	53 (93%)	4 (7%)	0	100	100
50	L8	55/60 (92%)	50 (91%)	4 (7%)	1 (2%)	11	49
51	I5	61/71 (86%)	34 (56%)	24 (39%)	3 (5%)	3	20
51	M8	64/71 (90%)	40 (62%)	21 (33%)	3 (5%)	3	21
52	J5	57/60 (95%)	50 (88%)	6 (10%)	1 (2%)	11	49
52	N8	56/60 (93%)	48 (86%)	6 (11%)	2 (4%)	4	28
53	K5	43/54 (80%)	26 (60%)	14 (33%)	3 (7%)	1	9
53	O8	43/54 (80%)	28 (65%)	13 (30%)	2 (5%)	3	21
54	L5	44/49 (90%)	42 (96%)	2 (4%)	0	100	100
54	P8	46/49 (94%)	45 (98%)	1 (2%)	0	100	100
55	M5	58/65 (89%)	47 (81%)	9 (16%)	2 (3%)	5	29
55	Q8	58/65 (89%)	31 (53%)	16 (28%)	11 (19%)	0	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	11340/12054 (94%)	9786 (86%)	1332 (12%)	222 (2%)	9	46

All (222) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	1E	237	ALA
18	9I	22	VAL
29	21	83	ASP
32	51	169	VAL
36	78	57	THR
41	C8	89	GLU
45	G8	54	LYS
49	K8	48	HIS
50	L8	54	VAL
51	M8	50	VAL
52	N8	41	PRO
55	Q8	7	HIS
55	Q8	44	LYS
55	Q8	50	LEU
55	Q8	55	ALA
9	82	118	LYS
19	AA	11	VAL
28	19	237	GLU
29	29	25	VAL
29	29	66	HIS
30	39	28	ILE
30	39	84	VAL
37	45	27	VAL
38	55	107	ASP
42	95	45	THR
46	D5	53	ILE
46	D5	165	VAL
46	D5	171	ILE
48	F5	30	VAL
51	I5	5	ILE
8	7E	86	ILE
14	5I	13	THR
29	21	60	ASN
29	21	78	LEU
31	41	96	ARG
32	51	168	PRO
37	88	66	ILE

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Mol	Chain	Res	Type
37	88	79	LEU
38	98	11	ASN
41	C8	93	LYS
42	D8	47	VAL
46	H8	60	GLU
46	H8	165	VAL
53	O8	17	LYS
55	Q8	8	LYS
55	Q8	36	LYS
55	Q8	49	VAL
55	Q8	51	ALA
4	32	32	ALA
11	2A	48	ILE
11	2A	100	ALA
12	3A	18	VAL
12	3A	26	ALA
14	5A	29	ARG
20	BA	73	HIS
28	19	33	LEU
29	29	51	PHE
29	29	81	ILE
29	29	90	THR
32	59	131	VAL
33	69	111	PRO
33	69	117	GLU
36	35	15	ARG
36	35	35	HIS
39	65	89	ARG
45	C5	17	SER
45	C5	29	GLU
46	D5	105	VAL
49	G5	47	ASN
49	G5	48	HIS
52	J5	57	VAL
55	M5	35	GLN
4	3E	31	CYS
4	3E	155	LEU
19	AI	7	LYS
19	AI	67	VAL
29	21	56	PRO
29	21	82	ARG
33	61	145	VAL

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Mol	Chain	Res	Type
34	58	127	ASP
34	58	128	HIS
36	78	36	LYS
37	88	2	LEU
39	A8	88	ASP
41	C8	88	ILE
44	F8	68	ARG
45	G8	53	PRO
46	H8	6	LYS
47	I8	83	PRO
48	J8	75	GLU
48	J8	76	ARG
52	N8	42	PRO
7	62	33	ASP
11	2A	101	SER
29	29	9	VAL
30	39	124	LEU
30	39	167	ALA
32	59	92	ILE
33	69	145	VAL
42	95	71	LEU
46	D5	7	ALA
47	E5	33	ALA
51	I5	26	SER
53	K5	17	LYS
14	5I	14	PRO
28	11	122	ASP
28	11	239	ARG
29	21	21	VAL
29	21	118	LYS
30	31	198	ALA
31	41	97	ASP
32	51	10	PRO
32	51	167	GLU
33	61	133	HIS
34	58	97	ARG
35	68	97	ARG
36	78	42	SER
37	88	20	ALA
45	G8	84	ARG
46	H8	59	LEU
49	K8	47	ASN

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Mol	Chain	Res	Type
55	Q8	31	HIS
20	BA	49	ALA
28	19	239	ARG
34	15	128	HIS
36	35	6	LEU
39	65	110	LEU
40	75	2	ASN
42	95	37	VAL
45	C5	92	ASN
46	D5	116	VAL
46	D5	158	PRO
46	D5	161	VAL
55	M5	31	HIS
28	11	3	VAL
31	41	5	VAL
32	51	12	PRO
34	58	22	THR
36	78	19	VAL
45	G8	76	CYS
45	G8	81	LYS
46	H8	81	ARG
51	M8	5	ILE
53	O8	18	ARG
2	12	73	THR
8	72	73	ASP
12	3A	19	ARG
29	29	26	ILE
29	29	62	PRO
30	39	25	PRO
32	59	168	PRO
36	35	108	LYS
39	65	87	PHE
41	85	93	LYS
42	95	44	LYS
45	C5	101	LYS
53	K5	42	TRP
2	1E	239	VAL
16	7I	51	VAL
30	31	24	LEU
33	61	12	LEU
37	88	3	MET
41	C8	90	VAL

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Mol	Chain	Res	Type
51	M8	34	GLU
55	Q8	47	LYS
2	12	7	VAL
12	3A	47	LYS
13	4A	95	GLY
29	29	63	LEU
29	29	77	ILE
38	55	3	HIS
39	65	111	GLU
51	I5	33	VAL
5	4E	115	VAL
11	2I	82	VAL
19	AI	41	VAL
28	11	240	ALA
29	21	55	ASN
29	21	72	VAL
32	51	92	ILE
33	61	131	LYS
36	78	95	VAL
45	G8	77	PRO
37	45	78	PRO
44	B5	51	VAL
46	D5	141	VAL
3	2E	13	GLY
29	21	22	PRO
42	D8	49	THR
55	Q8	58	ILE
2	12	39	ILE
8	72	100	ILE
12	3A	96	VAL
13	4A	84	ILE
28	19	3	VAL
30	39	89	VAL
31	49	5	VAL
28	11	123	ALA
36	78	7	ARG
46	H8	61	LEU
46	H8	62	PRO
46	H8	141	VAL
19	AA	67	VAL
29	29	52	LEU
33	69	144	VAL

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Mol	Chain	Res	Type
42	95	72	VAL
19	AI	9	VAL
20	BI	63	ILE
29	21	4	ILE
29	21	138	PRO
37	88	27	VAL
46	H8	53	ILE
28	19	118	VAL
29	29	91	VAL
32	59	4	ILE
32	59	167	GLU
36	35	7	ARG
36	35	34	GLY
53	K5	52	VAL
32	51	127	GLU
45	C5	76	CYS
46	D5	176	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	12	205/220 (93%)	169 (82%)	36 (18%)	2	11
2	1E	205/220 (93%)	152 (74%)	53 (26%)	0	2
3	22	160/188 (85%)	134 (84%)	26 (16%)	3	13
3	2E	159/188 (85%)	128 (80%)	31 (20%)	2	9
4	32	180/181 (99%)	144 (80%)	36 (20%)	1	8
4	3E	180/181 (99%)	138 (77%)	42 (23%)	1	4
5	42	116/123 (94%)	84 (72%)	32 (28%)	0	2
5	4E	116/123 (94%)	88 (76%)	28 (24%)	1	3
6	52	90/90 (100%)	71 (79%)	19 (21%)	1	7
6	5E	90/90 (100%)	78 (87%)	12 (13%)	5	22
7	62	126/127 (99%)	97 (77%)	29 (23%)	1	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	6E	126/127 (99%)	99 (79%)	27 (21%)	1	6
8	72	119/119 (100%)	91 (76%)	28 (24%)	1	4
8	7E	119/119 (100%)	95 (80%)	24 (20%)	1	7
9	82	98/99 (99%)	77 (79%)	21 (21%)	1	6
9	8E	98/99 (99%)	77 (79%)	21 (21%)	1	6
10	1A	89/92 (97%)	71 (80%)	18 (20%)	1	7
10	1I	89/92 (97%)	71 (80%)	18 (20%)	1	7
11	2A	90/99 (91%)	64 (71%)	26 (29%)	0	1
11	2I	90/99 (91%)	71 (79%)	19 (21%)	1	7
12	3A	104/109 (95%)	85 (82%)	19 (18%)	2	10
12	3I	104/109 (95%)	84 (81%)	20 (19%)	2	9
13	4A	94/101 (93%)	70 (74%)	24 (26%)	1	2
13	4I	94/101 (93%)	72 (77%)	22 (23%)	1	4
14	5A	48/50 (96%)	40 (83%)	8 (17%)	3	12
14	5I	49/50 (98%)	37 (76%)	12 (24%)	1	3
15	6A	79/80 (99%)	65 (82%)	14 (18%)	2	11
15	6I	79/80 (99%)	70 (89%)	9 (11%)	7	30
16	7A	72/74 (97%)	57 (79%)	15 (21%)	1	7
16	7I	72/74 (97%)	52 (72%)	20 (28%)	0	2
17	8A	95/97 (98%)	77 (81%)	18 (19%)	2	10
17	8I	95/97 (98%)	73 (77%)	22 (23%)	1	4
18	9A	63/77 (82%)	46 (73%)	17 (27%)	0	2
18	9I	63/77 (82%)	51 (81%)	12 (19%)	2	10
19	AA	67/80 (84%)	50 (75%)	17 (25%)	1	2
19	AI	70/80 (88%)	49 (70%)	21 (30%)	0	1
20	BA	76/82 (93%)	64 (84%)	12 (16%)	3	14
20	BI	76/82 (93%)	61 (80%)	15 (20%)	1	8
21	1B	20/22 (91%)	18 (90%)	2 (10%)	9	36
21	1F	20/22 (91%)	19 (95%)	1 (5%)	30	70
28	11	214/218 (98%)	167 (78%)	47 (22%)	1	6
28	19	214/218 (98%)	173 (81%)	41 (19%)	2	9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
29	21	165/166 (99%)	122 (74%)	43 (26%)	0	2
29	29	165/166 (99%)	125 (76%)	40 (24%)	1	3
30	31	161/166 (97%)	127 (79%)	34 (21%)	1	7
30	39	165/166 (99%)	124 (75%)	41 (25%)	1	3
31	41	155/156 (99%)	125 (81%)	30 (19%)	2	9
31	49	155/156 (99%)	127 (82%)	28 (18%)	2	10
32	51	145/148 (98%)	111 (77%)	34 (23%)	1	4
32	59	142/148 (96%)	105 (74%)	37 (26%)	0	2
33	61	122/124 (98%)	95 (78%)	27 (22%)	1	6
33	69	122/124 (98%)	89 (73%)	33 (27%)	0	2
34	15	117/119 (98%)	94 (80%)	23 (20%)	1	8
34	58	117/119 (98%)	90 (77%)	27 (23%)	1	4
35	25	100/100 (100%)	78 (78%)	22 (22%)	1	6
35	68	100/100 (100%)	89 (89%)	11 (11%)	8	31
36	35	116/116 (100%)	76 (66%)	40 (34%)	0	0
36	78	116/116 (100%)	85 (73%)	31 (27%)	0	2
37	45	111/111 (100%)	85 (77%)	26 (23%)	1	4
37	88	104/111 (94%)	78 (75%)	26 (25%)	1	3
38	55	100/101 (99%)	80 (80%)	20 (20%)	1	8
38	98	101/101 (100%)	73 (72%)	28 (28%)	0	2
39	65	87/88 (99%)	60 (69%)	27 (31%)	0	1
39	A8	87/88 (99%)	63 (72%)	24 (28%)	0	2
40	75	120/127 (94%)	81 (68%)	39 (32%)	0	0
40	B8	120/127 (94%)	89 (74%)	31 (26%)	0	2
41	85	93/94 (99%)	74 (80%)	19 (20%)	1	7
41	C8	93/94 (99%)	76 (82%)	17 (18%)	2	10
42	95	82/82 (100%)	64 (78%)	18 (22%)	1	6
42	D8	82/82 (100%)	57 (70%)	25 (30%)	0	1
43	A5	92/92 (100%)	69 (75%)	23 (25%)	1	3
43	E8	92/92 (100%)	66 (72%)	26 (28%)	0	1
44	B5	74/78 (95%)	50 (68%)	24 (32%)	0	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
44	F8	76/78 (97%)	60 (79%)	16 (21%)	1	7
45	C5	85/91 (93%)	59 (69%)	26 (31%)	0	1
45	G8	85/91 (93%)	59 (69%)	26 (31%)	0	1
46	D5	158/179 (88%)	119 (75%)	39 (25%)	1	3
46	H8	154/179 (86%)	128 (83%)	26 (17%)	2	12
47	E5	62/67 (92%)	51 (82%)	11 (18%)	2	11
47	I8	61/67 (91%)	45 (74%)	16 (26%)	0	2
48	F5	82/83 (99%)	67 (82%)	15 (18%)	2	10
48	J8	82/83 (99%)	61 (74%)	21 (26%)	0	2
49	G5	62/67 (92%)	46 (74%)	16 (26%)	0	2
49	K8	62/67 (92%)	42 (68%)	20 (32%)	0	0
50	H5	51/52 (98%)	38 (74%)	13 (26%)	1	2
50	L8	49/52 (94%)	33 (67%)	16 (33%)	0	0
51	I5	57/63 (90%)	46 (81%)	11 (19%)	2	9
51	M8	59/63 (94%)	42 (71%)	17 (29%)	0	1
52	J5	51/52 (98%)	40 (78%)	11 (22%)	1	6
52	N8	51/52 (98%)	39 (76%)	12 (24%)	1	4
53	K5	44/52 (85%)	36 (82%)	8 (18%)	2	10
53	O8	44/52 (85%)	30 (68%)	14 (32%)	0	1
54	L5	39/42 (93%)	31 (80%)	8 (20%)	1	7
54	P8	41/42 (98%)	31 (76%)	10 (24%)	1	3
55	M5	49/55 (89%)	37 (76%)	12 (24%)	1	3
55	Q8	50/55 (91%)	30 (60%)	20 (40%)	0	0
All	All	9568/9998 (96%)	7376 (77%)	2192 (23%)	1	5

All (2192) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	1E	8	LYS
2	1E	15	VAL
2	1E	17	PHE
2	1E	20	GLU
2	1E	21	ARG
2	1E	28	PHE

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Mol	Chain	Res	Type
2	1E	32	ILE
2	1E	37	ASN
2	1E	40	HIS
2	1E	64	ARG
2	1E	67	THR
2	1E	69	LEU
2	1E	71	VAL
2	1E	74	LYS
2	1E	82	ARG
2	1E	87	ARG
2	1E	96	ARG
2	1E	108	ILE
2	1E	113	HIS
2	1E	116	GLU
2	1E	118	LEU
2	1E	122	PHE
2	1E	125	PRO
2	1E	127	ILE
2	1E	130	ARG
2	1E	136	VAL
2	1E	144	ARG
2	1E	145	LEU
2	1E	146	GLN
2	1E	155	LEU
2	1E	157	ARG
2	1E	158	LEU
2	1E	160	ASP
2	1E	162	ILE
2	1E	163	PHE
2	1E	164	VAL
2	1E	170	GLU
2	1E	172	ILE
2	1E	175	ARG
2	1E	178	ARG
2	1E	187	LEU
2	1E	190	THR
2	1E	196	LEU
2	1E	200	ILE
2	1E	209	ARG
2	1E	213	LEU
2	1E	215	LEU
2	1E	216	SER

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Mol	Chain	Res	Type
2	1E	223	ILE
2	1E	226	ARG
2	1E	229	VAL
2	1E	231	GLU
2	1E	233	SER
3	2E	3	ASN
3	2E	4	LYS
3	2E	5	ILE
3	2E	8	ILE
3	2E	12	LEU
3	2E	17	ASP
3	2E	21	ARG
3	2E	26	LYS
3	2E	29	TYR
3	2E	31	HIS
3	2E	36	ASP
3	2E	56	ASP
3	2E	58	GLU
3	2E	62	ASP
3	2E	64	VAL
3	2E	84	ILE
3	2E	93	LYS
3	2E	95	THR
3	2E	98	ASN
3	2E	104	GLN
3	2E	108	ASN
3	2E	111	LEU
3	2E	136	GLN
3	2E	157	ILE
3	2E	161	GLU
3	2E	164	ARG
3	2E	167	TRP
3	2E	179	ARG
3	2E	188	LEU
3	2E	191	THR
3	2E	202	ILE
4	3E	3	ARG
4	3E	5	ILE
4	3E	10	ARG
4	3E	11	LEU
4	3E	13	ARG
4	3E	15	GLU

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Mol	Chain	Res	Type
4	3E	18	LYS
4	3E	24	GLU
4	3E	31	CYS
4	3E	45	GLN
4	3E	49	ARG
4	3E	50	ARG
4	3E	58	LEU
4	3E	59	ARG
4	3E	66	ARG
4	3E	71	SER
4	3E	83	SER
4	3E	86	LYS
4	3E	99	SER
4	3E	101	LEU
4	3E	106	TYR
4	3E	108	LEU
4	3E	122	ARG
4	3E	127	THR
4	3E	128	VAL
4	3E	132	ARG
4	3E	135	LEU
4	3E	138	TYR
4	3E	146	ILE
4	3E	151	LYS
4	3E	154	ASN
4	3E	179	GLU
4	3E	182	LYS
4	3E	184	LYS
4	3E	187	ARG
4	3E	188	LEU
4	3E	190	ASP
4	3E	191	ARG
4	3E	193	ASP
4	3E	200	GLU
4	3E	201	GLN
4	3E	209	ARG
5	4E	6	PHE
5	4E	11	ILE
5	4E	13	ILE
5	4E	15	ARG
5	4E	16	THR
5	4E	24	ARG

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Mol	Chain	Res	Type
5	4E	25	ARG
5	4E	26	PHE
5	4E	27	ARG
5	4E	31	LEU
5	4E	33	VAL
5	4E	41	VAL
5	4E	50	GLU
5	4E	51	VAL
5	4E	57	LYS
5	4E	65	ASN
5	4E	66	MET
5	4E	68	GLU
5	4E	72	GLN
5	4E	79	GLU
5	4E	80	ILE
5	4E	87	SER
5	4E	91	LEU
5	4E	116	THR
5	4E	126	ARG
5	4E	131	ILE
5	4E	151	LEU
5	4E	152	ARG
6	5E	19	LEU
6	5E	23	LYS
6	5E	27	GLN
6	5E	46	ARG
6	5E	55	ASP
6	5E	64	GLN
6	5E	65	VAL
6	5E	74	ASP
6	5E	75	LEU
6	5E	87	ARG
6	5E	89	MET
6	5E	94	GLN
7	6E	8	GLU
7	6E	9	VAL
7	6E	13	GLN
7	6E	22	LEU
7	6E	27	ILE
7	6E	31	MET
7	6E	36	LYS
7	6E	38	LEU

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Mol	Chain	Res	Type
7	6E	45	ASP
7	6E	48	LYS
7	6E	54	THR
7	6E	63	LYS
7	6E	66	VAL
7	6E	75	VAL
7	6E	80	VAL
7	6E	89	MET
7	6E	90	GLU
7	6E	91	VAL
7	6E	97	GLN
7	6E	98	SER
7	6E	104	LEU
7	6E	113	GLU
7	6E	124	LEU
7	6E	131	LYS
7	6E	146	GLU
7	6E	155	ARG
7	6E	156	TRP
8	7E	1	MET
8	7E	26	VAL
8	7E	29	SER
8	7E	33	GLU
8	7E	35	ILE
8	7E	39	LEU
8	7E	41	ARG
8	7E	45	ILE
8	7E	50	ARG
8	7E	52	ASP
8	7E	54	ASP
8	7E	63	LEU
8	7E	68	ARG
8	7E	80	ILE
8	7E	82	HIS
8	7E	83	ILE
8	7E	85	ARG
8	7E	88	LYS
8	7E	95	VAL
8	7E	102	ARG
8	7E	107	LEU
8	7E	112	LEU
8	7E	129	VAL

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Mol	Chain	Res	Type
8	7E	137	VAL
9	8E	9	ARG
9	8E	10	ARG
9	8E	16	ARG
9	8E	25	LYS
9	8E	27	THR
9	8E	31	GLN
9	8E	38	GLN
9	8E	40	LEU
9	8E	42	ARG
9	8E	47	LEU
9	8E	53	VAL
9	8E	54	ASP
9	8E	65	VAL
9	8E	79	LEU
9	8E	91	ASP
9	8E	95	LYS
9	8E	112	LYS
9	8E	117	HIS
9	8E	121	ARG
9	8E	124	GLN
9	8E	125	TYR
10	1I	5	ARG
10	1I	13	HIS
10	1I	16	LEU
10	1I	17	ASP
10	1I	19	SER
10	1I	24	VAL
10	1I	38	ILE
10	1I	48	THR
10	1I	49	VAL
10	1I	60	ARG
10	1I	62	HIS
10	1I	70	ARG
10	1I	76	ASN
10	1I	78	ASN
10	1I	84	GLN
10	1I	88	LEU
10	1I	92	THR
10	1I	96	ILE
11	2I	12	ARG
11	2I	13	GLN

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Mol	Chain	Res	Type
11	2I	14	VAL
11	2I	48	ILE
11	2I	63	LEU
11	2I	79	SER
11	2I	81	ASP
11	2I	84	VAL
11	2I	87	THR
11	2I	91	ARG
11	2I	93	GLN
11	2I	96	ARG
11	2I	103	LEU
11	2I	105	VAL
11	2I	106	LYS
11	2I	108	ILE
11	2I	109	VAL
11	2I	114	VAL
11	2I	116	HIS
12	3I	7	ILE
12	3I	11	VAL
12	3I	12	ARG
12	3I	18	VAL
12	3I	20	LYS
12	3I	28	LYS
12	3I	33	ARG
12	3I	38	THR
12	3I	55	VAL
12	3I	60	LEU
12	3I	61	THR
12	3I	62	SER
12	3I	64	TYR
12	3I	79	GLU
12	3I	82	VAL
12	3I	89	ARG
12	3I	96	VAL
12	3I	104	VAL
12	3I	114	LYS
12	3I	116	SER
13	4I	11	ARG
13	4I	14	ARG
13	4I	19	LEU
13	4I	32	GLU
13	4I	44	ARG

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Mol	Chain	Res	Type
13	4I	45	VAL
13	4I	48	LEU
13	4I	52	GLU
13	4I	56	LEU
13	4I	64	TRP
13	4I	67	GLU
13	4I	70	LEU
13	4I	83	ASP
13	4I	86	CYS
13	4I	88	ARG
13	4I	94	ARG
13	4I	105	THR
13	4I	106	ASN
13	4I	108	ARG
13	4I	110	ARG
13	4I	114	ARG
13	4I	115	LYS
14	5I	4	LYS
14	5I	8	GLU
14	5I	12	ARG
14	5I	17	LYS
14	5I	22	THR
14	5I	23	ARG
14	5I	26	ARG
14	5I	33	VAL
14	5I	40	CYS
14	5I	44	LEU
14	5I	50	LYS
14	5I	58	LYS
15	6I	6	GLU
15	6I	26	GLU
15	6I	39	LEU
15	6I	47	LYS
15	6I	66	LEU
15	6I	67	LEU
15	6I	68	ARG
15	6I	83	GLU
15	6I	87	ILE
16	7I	2	VAL
16	7I	4	ILE
16	7I	6	LEU
16	7I	19	ILE

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Mol	Chain	Res	Type
16	7I	20	VAL
16	7I	21	VAL
16	7I	22	THR
16	7I	25	ARG
16	7I	33	ILE
16	7I	45	THR
16	7I	47	ASP
16	7I	50	LYS
16	7I	54	GLU
16	7I	62	VAL
16	7I	67	THR
16	7I	69	THR
16	7I	72	ARG
16	7I	76	GLN
16	7I	82	GLN
16	7I	83	GLU
17	8I	6	LEU
17	8I	9	VAL
17	8I	12	SER
17	8I	27	PHE
17	8I	35	VAL
17	8I	38	ARG
17	8I	45	HIS
17	8I	48	GLU
17	8I	52	LYS
17	8I	60	ILE
17	8I	63	ARG
17	8I	67	LYS
17	8I	68	ARG
17	8I	74	LEU
17	8I	76	LEU
17	8I	79	SER
17	8I	87	LYS
17	8I	92	ARG
17	8I	97	SER
17	8I	99	SER
17	8I	100	LYS
17	8I	101	ARG
18	9I	18	ARG
18	9I	25	THR
18	9I	26	LEU
18	9I	29	PHE

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Mol	Chain	Res	Type
18	9I	31	LEU
18	9I	32	ARG
18	9I	33	ASP
18	9I	35	ARG
18	9I	46	GLU
18	9I	53	ARG
18	9I	58	LEU
18	9I	84	LYS
19	AI	5	LEU
19	AI	7	LYS
19	AI	12	ASP
19	AI	13	ASP
19	AI	15	LEU
19	AI	20	LEU
19	AI	21	GLU
19	AI	22	LEU
19	AI	29	ARG
19	AI	30	LEU
19	AI	31	ILE
19	AI	37	ARG
19	AI	43	GLU
19	AI	58	VAL
19	AI	60	VAL
19	AI	61	TYR
19	AI	64	GLU
19	AI	67	VAL
19	AI	77	THR
19	AI	78	ARG
19	AI	81	ARG
20	BI	9	ASN
20	BI	10	LEU
20	BI	11	SER
20	BI	13	LEU
20	BI	24	LEU
20	BI	30	LYS
20	BI	57	ARG
20	BI	64	ASP
20	BI	72	LEU
20	BI	73	HIS
20	BI	74	LYS
20	BI	75	ASN
20	BI	90	GLN

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Mol	Chain	Res	Type
20	BI	99	LEU
20	BI	100	ILE
21	1F	10	ARG
28	11	3	VAL
28	11	13	ARG
28	11	15	PHE
28	11	17	THR
28	11	23	GLU
28	11	28	GLU
28	11	30	GLU
28	11	34	VAL
28	11	35	LYS
28	11	37	LEU
28	11	46	GLN
28	11	58	HIS
28	11	59	LYS
28	11	61	LEU
28	11	64	ILE
28	11	65	ILE
28	11	68	LYS
28	11	94	LEU
28	11	95	LEU
28	11	105	ILE
28	11	106	ILE
28	11	111	LEU
28	11	113	VAL
28	11	131	LEU
28	11	136	ILE
28	11	141	VAL
28	11	142	VAL
28	11	150	LYS
28	11	155	LEU
28	11	162	SER
28	11	165	ILE
28	11	183	ARG
28	11	192	THR
28	11	193	VAL
28	11	200	ASP
28	11	212	SER
28	11	217	ARG
28	11	222	ARG
28	11	229	VAL

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Mol	Chain	Res	Type
28	11	239	ARG
28	11	242	ARG
28	11	257	LEU
28	11	259	THR
28	11	260	ARG
28	11	270	ILE
28	11	271	ILE
28	11	273	ARG
29	21	2	LYS
29	21	13	ARG
29	21	14	ILE
29	21	16	ARG
29	21	33	VAL
29	21	34	VAL
29	21	45	THR
29	21	47	VAL
29	21	48	GLN
29	21	55	ASN
29	21	59	VAL
29	21	63	LEU
29	21	66	HIS
29	21	72	VAL
29	21	78	LEU
29	21	87	GLU
29	21	89	ASP
29	21	92	THR
29	21	101	ARG
29	21	111	ARG
29	21	113	PHE
29	21	116	VAL
29	21	117	MET
29	21	118	LYS
29	21	119	ARG
29	21	136	ARG
29	21	138	PRO
29	21	144	ARG
29	21	146	THR
29	21	149	ARG
29	21	167	VAL
29	21	170	LEU
29	21	175	VAL
29	21	179	GLU

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Mol	Chain	Res	Type
29	21	180	ASN
29	21	185	LYS
29	21	188	VAL
29	21	195	LEU
29	21	196	VAL
29	21	197	ILE
29	21	200	GLU
29	21	202	LYS
29	21	203	LYS
30	31	8	GLN
30	31	9	ILE
30	31	12	LEU
30	31	15	SER
30	31	18	ARG
30	31	27	GLU
30	31	32	LEU
30	31	57	VAL
30	31	59	TYR
30	31	64	ILE
30	31	66	PRO
30	31	70	THR
30	31	77	ASP
30	31	78	ILE
30	31	88	VAL
30	31	101	LEU
30	31	103	LYS
30	31	106	ARG
30	31	116	ASP
30	31	117	ARG
30	31	119	ARG
30	31	132	VAL
30	31	145	GLU
30	31	148	LEU
30	31	158	THR
30	31	161	GLU
30	31	164	ARG
30	31	170	LEU
30	31	181	LEU
30	31	191	ARG
30	31	196	LEU
30	31	197	ASP
30	31	201	VAL

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Mol	Chain	Res	Type
30	31	203	GLN
31	41	3	LEU
31	41	4	ASP
31	41	10	LYS
31	41	13	GLU
31	41	14	GLU
31	41	20	ILE
31	41	21	ARG
31	41	26	GLN
31	41	28	VAL
31	41	31	VAL
31	41	34	LEU
31	41	43	LEU
31	41	45	GLU
31	41	62	LEU
31	41	63	ILE
31	41	67	LYS
31	41	74	LYS
31	41	76	SER
31	41	77	ILE
31	41	80	PHE
31	41	82	LEU
31	41	84	LYS
31	41	86	MET
31	41	94	LEU
31	41	101	ILE
31	41	135	LEU
31	41	155	MET
31	41	162	THR
31	41	166	ASP
31	41	174	GLU
32	51	2	SER
32	51	3	ARG
32	51	4	ILE
32	51	7	LEU
32	51	10	PRO
32	51	11	VAL
32	51	13	LYS
32	51	15	VAL
32	51	24	VAL
32	51	40	GLU
32	51	41	MET

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Mol	Chain	Res	Type
32	51	42	ARG
32	51	43	VAL
32	51	45	VAL
32	51	50	VAL
32	51	64	LEU
32	51	68	THR
32	51	71	LEU
32	51	72	ILE
32	51	77	LYS
32	51	80	SER
32	51	87	LEU
32	51	88	LEU
32	51	97	ARG
32	51	104	GLU
32	51	105	LEU
32	51	107	VAL
32	51	122	THR
32	51	129	THR
32	51	132	ARG
32	51	136	ILE
32	51	139	GLN
32	51	149	ARG
32	51	169	VAL
33	61	2	LYS
33	61	6	LEU
33	61	9	LEU
33	61	20	ASP
33	61	37	VAL
33	61	47	LEU
33	61	56	LYS
33	61	64	GLU
33	61	70	GLU
33	61	74	ASN
33	61	77	LEU
33	61	78	THR
33	61	82	ARG
33	61	85	GLU
33	61	86	THR
33	61	92	VAL
33	61	96	ASP
33	61	105	HIS
33	61	110	ASP

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Mol	Chain	Res	Type
33	61	116	LEU
33	61	117	GLU
33	61	131	LYS
33	61	135	GLU
33	61	139	GLN
33	61	140	LEU
33	61	142	VAL
33	61	143	SER
34	58	1	MET
34	58	2	LYS
34	58	5	VAL
34	58	7	LYS
34	58	12	ARG
34	58	14	VAL
34	58	28	THR
34	58	32	THR
34	58	34	LEU
34	58	38	HIS
34	58	43	THR
34	58	48	MET
34	58	50	ASP
34	58	58	ASP
34	58	60	ILE
34	58	61	ARG
34	58	65	LYS
34	58	67	LEU
34	58	79	PRO
34	58	87	LEU
34	58	90	MET
34	58	91	LEU
34	58	96	GLU
34	58	99	LEU
34	58	120	LEU
34	58	127	ASP
34	58	134	ARG
35	68	3	GLN
35	68	22	ILE
35	68	23	ARG
35	68	24	VAL
35	68	28	SER
35	68	66	LYS
35	68	88	ASN

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Mol	Chain	Res	Type
35	68	91	LEU
35	68	94	ARG
35	68	98	VAL
35	68	112	MET
36	78	10	PRO
36	78	13	ASN
36	78	15	ARG
36	78	16	ARG
36	78	18	ARG
36	78	19	VAL
36	78	30	THR
36	78	32	THR
36	78	41	ARG
36	78	45	LEU
36	78	46	LYS
36	78	49	ARG
36	78	50	ARG
36	78	55	ARG
36	78	57	THR
36	78	61	ARG
36	78	74	GLU
36	78	77	ARG
36	78	88	LEU
36	78	96	THR
36	78	99	LEU
36	78	100	LEU
36	78	105	LEU
36	78	106	LEU
36	78	112	LEU
36	78	119	GLU
36	78	126	VAL
36	78	135	LEU
36	78	138	LEU
36	78	144	GLU
36	78	147	LEU
37	88	1	MET
37	88	5	ARG
37	88	6	ARG
37	88	16	ARG
37	88	25	ASP
37	88	26	TYR
37	88	42	ILE

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Mol	Chain	Res	Type
37	88	45	GLN
37	88	58	PHE
37	88	59	ARG
37	88	67	ARG
37	88	76	LYS
37	88	77	LYS
37	88	82	ARG
37	88	85	LYS
37	88	87	LYS
37	88	103	MET
37	88	109	VAL
37	88	110	THR
37	88	112	GLU
37	88	115	MET
37	88	119	ARG
37	88	129	THR
37	88	134	ARG
37	88	139	GLU
37	88	141	GLN
38	98	4	LEU
38	98	6	SER
38	98	9	LYS
38	98	10	LEU
38	98	12	ARG
38	98	28	LEU
38	98	29	LEU
38	98	34	ILE
38	98	36	THR
38	98	44	LEU
38	98	45	ARG
38	98	48	VAL
38	98	59	ASP
38	98	65	LEU
38	98	67	LEU
38	98	71	GLN
38	98	75	LEU
38	98	79	LEU
38	98	82	GLU
38	98	88	ARG
38	98	91	GLN
38	98	94	TYR
38	98	96	ARG

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Mol	Chain	Res	Type
38	98	105	ARG
38	98	107	ASP
38	98	113	LEU
38	98	117	VAL
38	98	118	GLU
39	A8	3	ARG
39	A8	4	LEU
39	A8	8	GLU
39	A8	15	ARG
39	A8	17	ARG
39	A8	29	PHE
39	A8	32	LEU
39	A8	35	ILE
39	A8	36	TYR
39	A8	43	GLU
39	A8	44	LYS
39	A8	46	VAL
39	A8	48	LEU
39	A8	50	SER
39	A8	56	LEU
39	A8	58	LEU
39	A8	61	ASN
39	A8	73	LEU
39	A8	80	LEU
39	A8	83	LYS
39	A8	89	ARG
39	A8	98	VAL
39	A8	101	LEU
39	A8	106	ARG
40	B8	6	LEU
40	B8	10	VAL
40	B8	11	GLU
40	B8	16	ARG
40	B8	19	LEU
40	B8	21	GLU
40	B8	27	THR
40	B8	30	VAL
40	B8	35	LYS
40	B8	38	ASN
40	B8	39	ARG
40	B8	42	ILE
40	B8	48	ILE

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Mol	Chain	Res	Type
40	B8	50	ILE
40	B8	58	ASN
40	B8	59	THR
40	B8	64	ARG
40	B8	65	LYS
40	B8	82	LEU
40	B8	85	LYS
40	B8	86	ILE
40	B8	89	VAL
40	B8	95	ARG
40	B8	96	ARG
40	B8	98	LYS
40	B8	99	LEU
40	B8	106	SER
40	B8	108	ARG
40	B8	110	ILE
40	B8	111	ARG
40	B8	112	ARG
41	C8	3	ARG
41	C8	5	LYS
41	C8	17	ILE
41	C8	27	LEU
41	C8	34	LYS
41	C8	52	ARG
41	C8	60	LEU
41	C8	70	ARG
41	C8	74	LEU
41	C8	85	LYS
41	C8	89	GLU
41	C8	92	ARG
41	C8	94	ASN
41	C8	104	GLN
41	C8	108	GLU
41	C8	111	GLU
41	C8	112	ARG
42	D8	1	MET
42	D8	5	VAL
42	D8	6	LYS
42	D8	7	THR
42	D8	12	TYR
42	D8	15	GLU
42	D8	18	LEU

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Mol	Chain	Res	Type
42	D8	19	LYS
42	D8	20	LEU
42	D8	22	VAL
42	D8	24	LYS
42	D8	33	VAL
42	D8	35	LEU
42	D8	38	LEU
42	D8	40	LEU
42	D8	44	LYS
42	D8	47	VAL
42	D8	73	SER
42	D8	79	VAL
42	D8	82	ARG
42	D8	83	ARG
42	D8	85	LYS
42	D8	87	HIS
42	D8	88	ARG
42	D8	100	ARG
43	E8	1	MET
43	E8	11	ARG
43	E8	14	PRO
43	E8	15	ARG
43	E8	27	LYS
43	E8	51	LEU
43	E8	60	ASN
43	E8	62	HIS
43	E8	63	ASP
43	E8	66	GLU
43	E8	69	LEU
43	E8	70	TYR
43	E8	76	VAL
43	E8	78	GLU
43	E8	84	ARG
43	E8	86	LEU
43	E8	88	ARG
43	E8	90	ARG
43	E8	92	ARG
43	E8	96	ILE
43	E8	97	LYS
43	E8	103	ILE
43	E8	107	LEU
43	E8	109	GLU

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Mol	Chain	Res	Type
43	E8	111	HIS
43	E8	113	LYS
44	F8	2	LYS
44	F8	12	VAL
44	F8	14	SER
44	F8	15	GLU
44	F8	23	GLU
44	F8	48	LYS
44	F8	53	LYS
44	F8	54	VAL
44	F8	57	LEU
44	F8	68	ARG
44	F8	72	LYS
44	F8	76	ARG
44	F8	78	LYS
44	F8	80	ILE
44	F8	87	GLN
44	F8	88	LYS
45	G8	6	HIS
45	G8	7	VAL
45	G8	9	LYS
45	G8	21	LYS
45	G8	24	VAL
45	G8	26	LYS
45	G8	33	LYS
45	G8	38	ILE
45	G8	39	VAL
45	G8	40	GLU
45	G8	42	VAL
45	G8	44	ILE
45	G8	50	ARG
45	G8	51	VAL
45	G8	52	SER
45	G8	54	LYS
45	G8	57	GLN
45	G8	63	LYS
45	G8	64	GLU
45	G8	67	LEU
45	G8	79	CYS
45	G8	84	ARG
45	G8	85	VAL
45	G8	86	ARG

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Mol	Chain	Res	Type
45	G8	97	ARG
45	G8	99	CYS
46	H8	1	MET
46	H8	5	LEU
46	H8	19	ARG
46	H8	24	LEU
46	H8	42	VAL
46	H8	46	LYS
46	H8	53	ILE
46	H8	60	GLU
46	H8	61	LEU
46	H8	71	VAL
46	H8	76	LEU
46	H8	77	ASP
46	H8	80	ARG
46	H8	81	ARG
46	H8	86	VAL
46	H8	91	LEU
46	H8	94	GLU
46	H8	96	VAL
46	H8	105	VAL
46	H8	117	LEU
46	H8	119	GLU
46	H8	140	ASP
46	H8	144	LEU
46	H8	148	ASP
46	H8	150	LEU
46	H8	154	ASP
47	I8	11	ARG
47	I8	27	GLU
47	I8	29	GLN
47	I8	36	ILE
47	I8	37	LEU
47	I8	38	VAL
47	I8	40	GLN
47	I8	41	ARG
47	I8	46	LYS
47	I8	49	LYS
47	I8	53	MET
47	I8	55	ARG
47	I8	58	THR
47	I8	66	VAL

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Mol	Chain	Res	Type
47	I8	70	GLN
47	I8	74	ARG
48	J8	13	ILE
48	J8	21	ARG
48	J8	35	THR
48	J8	41	ARG
48	J8	46	LEU
48	J8	51	VAL
48	J8	57	GLU
48	J8	62	VAL
48	J8	72	GLU
48	J8	78	LYS
48	J8	80	LEU
48	J8	81	LYS
48	J8	82	LEU
48	J8	83	GLU
48	J8	86	SER
48	J8	90	ILE
48	J8	91	LYS
48	J8	93	GLU
48	J8	94	LEU
48	J8	97	LEU
48	J8	98	LEU
49	K8	4	SER
49	K8	5	GLU
49	K8	6	VAL
49	K8	16	LEU
49	K8	19	VAL
49	K8	24	LEU
49	K8	25	VAL
49	K8	41	ILE
49	K8	45	SER
49	K8	46	GLN
49	K8	47	ASN
49	K8	48	HIS
49	K8	49	LYS
49	K8	50	ILE
49	K8	52	ASP
49	K8	53	LEU
49	K8	54	LYS
49	K8	55	ARG
49	K8	62	THR

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Mol	Chain	Res	Type
49	K8	64	LEU
50	L8	3	ARG
50	L8	6	VAL
50	L8	8	LEU
50	L8	11	SER
50	L8	13	ILE
50	L8	17	LYS
50	L8	30	ARG
50	L8	31	LEU
50	L8	33	GLN
50	L8	36	VAL
50	L8	40	THR
50	L8	43	ILE
50	L8	44	ARG
50	L8	53	LEU
50	L8	57	GLU
50	L8	58	VAL
51	M8	6	HIS
51	M8	15	ILE
51	M8	16	CYS
51	M8	23	GLU
51	M8	26	SER
51	M8	36	CYS
51	M8	38	LYS
51	M8	39	CYS
51	M8	48	ARG
51	M8	50	VAL
51	M8	52	THR
51	M8	55	ARG
51	M8	59	PHE
51	M8	61	ARG
51	M8	62	ARG
51	M8	63	TYR
51	M8	65	ASP
52	N8	3	LYS
52	N8	9	LYS
52	N8	15	ARG
52	N8	16	ARG
52	N8	29	THR
52	N8	31	VAL
52	N8	40	LYS
52	N8	51	TYR

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Mol	Chain	Res	Type
52	N8	55	ARG
52	N8	56	LYS
52	N8	59	GLU
52	N8	60	VAL
53	O8	9	LEU
53	O8	10	LEU
53	O8	12	GLU
53	O8	17	LYS
53	O8	23	THR
53	O8	27	LYS
53	O8	30	THR
53	O8	32	ASN
53	O8	36	LEU
53	O8	37	ARG
53	O8	39	TYR
53	O8	44	ARG
53	O8	47	THR
53	O8	51	GLU
54	P8	4	THR
54	P8	8	ASN
54	P8	14	LYS
54	P8	19	ARG
54	P8	22	MET
54	P8	32	LYS
54	P8	41	ARG
54	P8	43	THR
54	P8	47	ARG
54	P8	48	LYS
55	Q8	6	THR
55	Q8	8	LYS
55	Q8	11	LYS
55	Q8	19	SER
55	Q8	21	LYS
55	Q8	26	LYS
55	Q8	30	ARG
55	Q8	31	HIS
55	Q8	32	LEU
55	Q8	41	ILE
55	Q8	42	ARG
55	Q8	43	GLN
55	Q8	46	ARG
55	Q8	47	LYS

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Mol	Chain	Res	Type
55	Q8	48	PHE
55	Q8	50	LEU
55	Q8	52	LYS
55	Q8	57	ARG
55	Q8	58	ILE
55	Q8	59	LYS
2	12	5	ILE
2	12	17	PHE
2	12	22	LYS
2	12	23	ARG
2	12	24	TRP
2	12	31	TYR
2	12	33	TYR
2	12	42	ILE
2	12	44	LEU
2	12	55	PHE
2	12	69	LEU
2	12	71	VAL
2	12	75	LYS
2	12	83	MET
2	12	96	ARG
2	12	107	THR
2	12	108	ILE
2	12	117	GLU
2	12	122	PHE
2	12	130	ARG
2	12	139	LYS
2	12	144	ARG
2	12	145	LEU
2	12	155	LEU
2	12	164	VAL
2	12	172	ILE
2	12	178	ARG
2	12	185	ILE
2	12	193	ASP
2	12	196	LEU
2	12	204	ASN
2	12	212	GLN
2	12	213	LEU
2	12	222	ILE
2	12	226	ARG
2	12	238	LEU

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Mol	Chain	Res	Type
3	22	4	LYS
3	22	5	ILE
3	22	16	ARG
3	22	21	ARG
3	22	29	TYR
3	22	32	LEU
3	22	43	LEU
3	22	47	LEU
3	22	52	LEU
3	22	67	THR
3	22	69	HIS
3	22	95	THR
3	22	119	ARG
3	22	120	VAL
3	22	124	ILE
3	22	128	PHE
3	22	139	GLN
3	22	140	ARG
3	22	142	MET
3	22	167	TRP
3	22	175	LEU
3	22	190	ARG
3	22	191	THR
3	22	193	TYR
3	22	195	VAL
3	22	202	ILE
4	32	3	ARG
4	32	5	ILE
4	32	8	VAL
4	32	18	LYS
4	32	24	GLU
4	32	30	LYS
4	32	36	ARG
4	32	50	ARG
4	32	59	ARG
4	32	61	LYS
4	32	71	SER
4	32	73	ARG
4	32	76	ARG
4	32	78	LEU
4	32	106	TYR
4	32	107	ARG

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Mol	Chain	Res	Type
4	32	108	LEU
4	32	119	GLN
4	32	122	ARG
4	32	127	THR
4	32	134	ASP
4	32	135	LEU
4	32	138	TYR
4	32	139	ARG
4	32	141	ARG
4	32	152	SER
4	32	159	ARG
4	32	162	LEU
4	32	163	GLU
4	32	169	LYS
4	32	187	ARG
4	32	191	ARG
4	32	192	GLU
4	32	199	ASN
4	32	200	GLU
4	32	209	ARG
5	42	8	GLU
5	42	12	LEU
5	42	14	ARG
5	42	16	THR
5	42	19	MET
5	42	25	ARG
5	42	40	ARG
5	42	41	VAL
5	42	43	LEU
5	42	45	PHE
5	42	47	LYS
5	42	51	VAL
5	42	53	LEU
5	42	56	GLN
5	42	68	GLU
5	42	72	GLN
5	42	73	ASN
5	42	78	HIS
5	42	79	GLU
5	42	82	VAL
5	42	83	GLU
5	42	87	SER

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Mol	Chain	Res	Type
5	42	90	VAL
5	42	91	LEU
5	42	101	ILE
5	42	115	VAL
5	42	116	THR
5	42	126	ARG
5	42	136	MET
5	42	137	GLU
5	42	144	THR
5	42	150	ARG
6	52	2	ARG
6	52	3	ARG
6	52	7	ASN
6	52	14	LEU
6	52	15	ASP
6	52	21	LEU
6	52	27	GLN
6	52	28	ARG
6	52	40	VAL
6	52	45	LEU
6	52	47	ARG
6	52	54	LYS
6	52	57	GLN
6	52	64	GLN
6	52	65	VAL
6	52	69	GLU
6	52	70	ASP
6	52	72	VAL
6	52	86	ARG
7	62	5	ARG
7	62	8	GLU
7	62	9	VAL
7	62	21	VAL
7	62	29	LYS
7	62	31	MET
7	62	45	ASP
7	62	51	GLN
7	62	52	GLU
7	62	54	THR
7	62	60	LYS
7	62	61	VAL
7	62	63	LYS

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Mol	Chain	Res	Type
7	62	70	LYS
7	62	72	ARG
7	62	73	MET
7	62	78	ARG
7	62	84	ASN
7	62	85	TYR
7	62	86	GLN
7	62	89	MET
7	62	90	GLU
7	62	91	VAL
7	62	104	LEU
7	62	114	ARG
7	62	137	LYS
7	62	144	MET
7	62	146	GLU
7	62	149	ARG
8	72	1	MET
8	72	2	LEU
8	72	12	ARG
8	72	21	LYS
8	72	25	ASP
8	72	35	ILE
8	72	39	LEU
8	72	54	ASP
8	72	69	ARG
8	72	73	ASP
8	72	77	GLU
8	72	78	GLN
8	72	81	HIS
8	72	82	HIS
8	72	83	ILE
8	72	84	ARG
8	72	86	ILE
8	72	88	LYS
8	72	91	ARG
8	72	92	ARG
8	72	95	VAL
8	72	97	VAL
8	72	99	GLU
8	72	102	ARG
8	72	112	LEU
8	72	114	THR

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Mol	Chain	Res	Type
8	72	115	SER
8	72	133	LEU
9	82	10	ARG
9	82	31	GLN
9	82	33	PHE
9	82	42	ARG
9	82	53	VAL
9	82	54	ASP
9	82	70	LYS
9	82	75	ASP
9	82	79	LEU
9	82	88	TYR
9	82	93	ARG
9	82	95	LYS
9	82	99	LEU
9	82	104	ARG
9	82	109	VAL
9	82	112	LYS
9	82	113	LYS
9	82	117	HIS
9	82	118	LYS
9	82	125	TYR
9	82	128	ARG
10	1A	16	LEU
10	1A	17	ASP
10	1A	29	ARG
10	1A	40	LEU
10	1A	43	ARG
10	1A	47	PHE
10	1A	48	THR
10	1A	55	LYS
10	1A	58	ASP
10	1A	59	SER
10	1A	62	HIS
10	1A	69	ASN
10	1A	70	ARG
10	1A	74	ILE
10	1A	79	ARG
10	1A	86	MET
10	1A	96	ILE
10	1A	99	LYS
11	2A	12	ARG

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Mol	Chain	Res	Type
11	2A	13	GLN
11	2A	14	VAL
11	2A	18	ARG
11	2A	24	SER
11	2A	29	ILE
11	2A	31	THR
11	2A	33	THR
11	2A	38	ASN
11	2A	57	THR
11	2A	63	LEU
11	2A	70	LYS
11	2A	82	VAL
11	2A	87	THR
11	2A	93	GLN
11	2A	95	ILE
11	2A	96	ARG
11	2A	103	LEU
11	2A	105	VAL
11	2A	106	LYS
11	2A	109	VAL
11	2A	116	HIS
11	2A	119	CYS
11	2A	120	ARG
11	2A	124	LYS
11	2A	127	LYS
12	3A	13	LYS
12	3A	19	ARG
12	3A	20	LYS
12	3A	28	LYS
12	3A	33	ARG
12	3A	40	VAL
12	3A	41	ARG
12	3A	42	THR
12	3A	50	SER
12	3A	54	LYS
12	3A	57	LYS
12	3A	64	TYR
12	3A	81	SER
12	3A	83	VAL
12	3A	85	ILE
12	3A	89	ARG
12	3A	96	VAL

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Mol	Chain	Res	Type
12	3A	111	LYS
12	3A	118	SER
13	4A	4	ILE
13	4A	8	GLU
13	4A	17	VAL
13	4A	32	GLU
13	4A	39	ILE
13	4A	47	ASP
13	4A	48	LEU
13	4A	64	TRP
13	4A	65	LYS
13	4A	66	LEU
13	4A	69	GLU
13	4A	70	LEU
13	4A	71	ARG
13	4A	82	MET
13	4A	83	ASP
13	4A	88	ARG
13	4A	92	HIS
13	4A	93	ARG
13	4A	94	ARG
13	4A	98	VAL
13	4A	103	THR
13	4A	108	ARG
13	4A	114	ARG
13	4A	115	LYS
14	5A	8	GLU
14	5A	12	ARG
14	5A	17	LYS
14	5A	22	THR
14	5A	23	ARG
14	5A	33	VAL
14	5A	37	PHE
14	5A	44	LEU
15	6A	3	ILE
15	6A	6	GLU
15	6A	22	THR
15	6A	31	LEU
15	6A	41	GLU
15	6A	47	LYS
15	6A	48	LYS
15	6A	51	HIS

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Mol	Chain	Res	Type
15	6A	58	MET
15	6A	82	ILE
15	6A	83	GLU
15	6A	84	LYS
15	6A	87	ILE
15	6A	88	ARG
16	7A	1	MET
16	7A	2	VAL
16	7A	5	ARG
16	7A	6	LEU
16	7A	8	ARG
16	7A	21	VAL
16	7A	25	ARG
16	7A	43	LYS
16	7A	51	VAL
16	7A	55	ARG
16	7A	65	GLN
16	7A	67	THR
16	7A	72	ARG
16	7A	74	LEU
16	7A	82	GLN
17	8A	6	LEU
17	8A	9	VAL
17	8A	25	ARG
17	8A	26	GLN
17	8A	35	VAL
17	8A	50	LYS
17	8A	53	LEU
17	8A	60	ILE
17	8A	62	SER
17	8A	63	ARG
17	8A	68	ARG
17	8A	70	ARG
17	8A	74	LEU
17	8A	79	SER
17	8A	91	ARG
17	8A	92	ARG
17	8A	100	LYS
17	8A	101	ARG
18	9A	21	LYS
18	9A	26	LEU
18	9A	29	PHE

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Mol	Chain	Res	Type
18	9A	31	LEU
18	9A	36	ASN
18	9A	42	ARG
18	9A	44	LEU
18	9A	47	THR
18	9A	53	ARG
18	9A	54	ARG
18	9A	58	LEU
18	9A	59	SER
18	9A	65	ILE
18	9A	82	THR
18	9A	84	LYS
18	9A	86	VAL
18	9A	87	ARG
19	AA	11	VAL
19	AA	13	ASP
19	AA	15	LEU
19	AA	22	LEU
19	AA	25	LYS
19	AA	30	LEU
19	AA	31	ILE
19	AA	33	THR
19	AA	34	TRP
19	AA	41	VAL
19	AA	44	MET
19	AA	49	ILE
19	AA	64	GLU
19	AA	66	MET
19	AA	79	THR
19	AA	81	ARG
19	AA	83	HIS
20	BA	11	SER
20	BA	13	LEU
20	BA	23	ARG
20	BA	56	MET
20	BA	68	LYS
20	BA	70	SER
20	BA	73	HIS
20	BA	74	LYS
20	BA	75	ASN
20	BA	84	LEU
20	BA	86	ARG

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Mol	Chain	Res	Type
20	BA	88	VAL
21	1B	9	ARG
21	1B	25	LYS
28	19	10	THR
28	19	13	ARG
28	19	24	ILE
28	19	25	THR
28	19	32	SER
28	19	38	LYS
28	19	43	ARG
28	19	49	ILE
28	19	61	LEU
28	19	64	ILE
28	19	65	ILE
28	19	68	LYS
28	19	69	ARG
28	19	73	VAL
28	19	78	LYS
28	19	88	ARG
28	19	91	ARG
28	19	94	LEU
28	19	99	ASP
28	19	105	ILE
28	19	138	VAL
28	19	141	VAL
28	19	154	LYS
28	19	169	GLU
28	19	181	GLU
28	19	182	LEU
28	19	192	THR
28	19	208	LYS
28	19	211	ARG
28	19	212	SER
28	19	217	ARG
28	19	222	ARG
28	19	239	ARG
28	19	242	ARG
28	19	244	ARG
28	19	255	LYS
28	19	260	ARG
28	19	263	ARG
28	19	266	SER

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Mol	Chain	Res	Type
28	19	268	ARG
28	19	271	ILE
29	29	1	MET
29	29	4	ILE
29	29	7	VAL
29	29	16	ARG
29	29	23	VAL
29	29	25	VAL
29	29	27	LEU
29	29	41	LYS
29	29	44	TYR
29	29	45	THR
29	29	48	GLN
29	29	58	ARG
29	29	59	VAL
29	29	66	HIS
29	29	67	PHE
29	29	75	VAL
29	29	76	ARG
29	29	79	ARG
29	29	82	ARG
29	29	87	GLU
29	29	90	THR
29	29	93	VAL
29	29	107	THR
29	29	113	PHE
29	29	116	VAL
29	29	117	MET
29	29	119	ARG
29	29	121	ASN
29	29	144	ARG
29	29	146	THR
29	29	154	LYS
29	29	170	LEU
29	29	175	VAL
29	29	178	GLU
29	29	179	GLU
29	29	181	LEU
29	29	188	VAL
29	29	200	GLU
29	29	201	THR
29	29	202	LYS

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Mol	Chain	Res	Type
30	39	2	LYS
30	39	7	TYR
30	39	8	GLN
30	39	11	VAL
30	39	12	LEU
30	39	13	SER
30	39	20	LEU
30	39	24	LEU
30	39	29	ASN
30	39	38	ARG
30	39	45	ARG
30	39	62	ARG
30	39	65	TRP
30	39	66	PRO
30	39	67	GLN
30	39	69	HIS
30	39	74	ARG
30	39	83	PHE
30	39	106	ARG
30	39	107	LYS
30	39	110	LEU
30	39	112	MET
30	39	123	LEU
30	39	124	LEU
30	39	125	LEU
30	39	127	GLU
30	39	140	LEU
30	39	152	GLU
30	39	153	SER
30	39	158	THR
30	39	169	ASN
30	39	181	LEU
30	39	191	ARG
30	39	192	LEU
30	39	193	VAL
30	39	194	MET
30	39	196	LEU
30	39	197	ASP
30	39	201	VAL
30	39	202	PHE
30	39	205	ARG
31	49	3	LEU

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Mol	Chain	Res	Type
31	49	7	LEU
31	49	13	GLU
31	49	19	LEU
31	49	20	ILE
31	49	28	VAL
31	49	52	ILE
31	49	58	GLN
31	49	64	THR
31	49	67	LYS
31	49	70	VAL
31	49	71	THR
31	49	82	LEU
31	49	83	ARG
31	49	84	LYS
31	49	91	ARG
31	49	96	ARG
31	49	109	VAL
31	49	116	ASP
31	49	118	ARG
31	49	130	ASN
31	49	137	GLU
31	49	139	LEU
31	49	144	ILE
31	49	153	ARG
31	49	157	ILE
31	49	159	VAL
31	49	162	THR
32	59	3	ARG
32	59	4	ILE
32	59	6	ARG
32	59	7	LEU
32	59	11	VAL
32	59	16	SER
32	59	37	VAL
32	59	41	MET
32	59	43	VAL
32	59	44	VAL
32	59	45	VAL
32	59	46	GLU
32	59	71	LEU
32	59	74	ASN
32	59	83	TYR

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Mol	Chain	Res	Type
32	59	85	LYS
32	59	86	GLU
32	59	89	ILE
32	59	99	VAL
32	59	103	LEU
32	59	105	LEU
32	59	122	THR
32	59	123	PHE
32	59	124	GLU
32	59	125	VAL
32	59	129	THR
32	59	131	VAL
32	59	132	ARG
32	59	137	ASP
32	59	139	GLN
32	59	143	GLN
32	59	144	VAL
32	59	147	ASN
32	59	152	ARG
32	59	160	LYS
32	59	164	TYR
32	59	171	LEU
33	69	2	LYS
33	69	7	GLU
33	69	9	LEU
33	69	10	GLU
33	69	35	LEU
33	69	37	VAL
33	69	43	ASN
33	69	44	LEU
33	69	51	ILE
33	69	56	LYS
33	69	58	LEU
33	69	64	GLU
33	69	67	ARG
33	69	68	LEU
33	69	73	GLU
33	69	75	LEU
33	69	77	LEU
33	69	78	THR
33	69	81	VAL
33	69	82	ARG

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Mol	Chain	Res	Type
33	69	87	LYS
33	69	88	ILE
33	69	91	SER
33	69	96	ASP
33	69	101	LEU
33	69	109	ILE
33	69	110	ASP
33	69	113	ARG
33	69	114	LEU
33	69	117	GLU
33	69	125	GLU
33	69	131	LYS
33	69	133	HIS
34	15	4	TYR
34	15	5	VAL
34	15	29	LYS
34	15	32	THR
34	15	33	LEU
34	15	34	LEU
34	15	35	ARG
34	15	38	HIS
34	15	43	THR
34	15	46	VAL
34	15	48	MET
34	15	56	ASN
34	15	59	LYS
34	15	63	THR
34	15	67	LEU
34	15	73	THR
34	15	74	ARG
34	15	85	ILE
34	15	93	THR
34	15	106	MET
34	15	112	LEU
34	15	130	HIS
34	15	137	LYS
35	25	5	GLN
35	25	22	ILE
35	25	24	VAL
35	25	28	SER
35	25	31	LYS
35	25	32	TYR

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Mol	Chain	Res	Type
35	25	38	VAL
35	25	47	ILE
35	25	49	ARG
35	25	53	LYS
35	25	58	VAL
35	25	70	LYS
35	25	80	ASP
35	25	87	ILE
35	25	94	ARG
35	25	96	THR
35	25	97	ARG
35	25	98	VAL
35	25	104	ARG
35	25	113	LYS
35	25	116	SER
35	25	117	LEU
36	35	1	MET
36	35	3	LEU
36	35	6	LEU
36	35	10	PRO
36	35	14	LYS
36	35	15	ARG
36	35	18	ARG
36	35	21	ARG
36	35	30	THR
36	35	36	LYS
36	35	41	ARG
36	35	45	LEU
36	35	46	LYS
36	35	50	ARG
36	35	61	ARG
36	35	62	LEU
36	35	65	ARG
36	35	67	MET
36	35	70	GLN
36	35	75	ILE
36	35	81	GLN
36	35	85	LEU
36	35	90	ARG
36	35	96	THR
36	35	98	GLU
36	35	105	LEU

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Mol	Chain	Res	Type
36	35	106	LEU
36	35	110	TYR
36	35	111	ARG
36	35	112	LEU
36	35	114	ILE
36	35	123	LEU
36	35	124	LYS
36	35	125	VAL
36	35	133	SER
36	35	135	LEU
36	35	138	LEU
36	35	144	GLU
36	35	146	VAL
36	35	147	LEU
37	45	5	ARG
37	45	6	ARG
37	45	10	ARG
37	45	16	ARG
37	45	18	LYS
37	45	21	THR
37	45	22	LYS
37	45	25	ASP
37	45	27	VAL
37	45	45	GLN
37	45	56	ARG
37	45	59	ARG
37	45	60	ARG
37	45	77	LYS
37	45	79	LEU
37	45	81	VAL
37	45	83	MET
37	45	91	GLU
37	45	103	MET
37	45	112	GLU
37	45	116	GLU
37	45	118	LEU
37	45	127	ILE
37	45	128	LYS
37	45	134	ARG
37	45	141	GLN
38	55	15	SER
38	55	18	LEU

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Mol	Chain	Res	Type
38	55	24	GLN
38	55	28	LEU
38	55	29	LEU
38	55	35	THR
38	55	44	LEU
38	55	57	ARG
38	55	65	LEU
38	55	67	LEU
38	55	74	LYS
38	55	75	LEU
38	55	76	VAL
38	55	79	LEU
38	55	80	PHE
38	55	81	ASP
38	55	82	GLU
38	55	91	GLN
38	55	95	THR
38	55	102	GLU
39	65	4	LEU
39	65	12	PHE
39	65	20	ARG
39	65	21	THR
39	65	23	ARG
39	65	30	ARG
39	65	36	TYR
39	65	38	GLN
39	65	42	ASP
39	65	52	SER
39	65	56	LEU
39	65	58	LEU
39	65	64	GLU
39	65	65	VAL
39	65	69	VAL
39	65	71	ARG
39	65	78	LEU
39	65	82	ILE
39	65	88	ASP
39	65	89	ARG
39	65	93	LYS
39	65	95	HIS
39	65	101	LEU
39	65	106	ARG

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Mol	Chain	Res	Type
39	65	107	GLU
39	65	110	LEU
39	65	112	PHE
40	75	6	LEU
40	75	7	ILE
40	75	8	LYS
40	75	9	LEU
40	75	10	VAL
40	75	11	GLU
40	75	13	ARG
40	75	19	LEU
40	75	21	GLU
40	75	23	ARG
40	75	24	PRO
40	75	27	THR
40	75	28	VAL
40	75	34	VAL
40	75	39	ARG
40	75	40	THR
40	75	41	ARG
40	75	52	ILE
40	75	54	ARG
40	75	55	ASN
40	75	60	THR
40	75	62	THR
40	75	63	VAL
40	75	64	ARG
40	75	66	VAL
40	75	67	SER
40	75	74	ARG
40	75	88	ILE
40	75	91	ARG
40	75	93	ARG
40	75	105	LEU
40	75	106	SER
40	75	107	ASP
40	75	112	ARG
40	75	120	ARG
40	75	124	ASP
40	75	125	ARG
40	75	129	ARG
40	75	132	LYS

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Mol	Chain	Res	Type
41	85	3	ARG
41	85	5	LYS
41	85	8	VAL
41	85	15	LYS
41	85	17	ILE
41	85	31	SER
41	85	34	LYS
41	85	55	ARG
41	85	57	PHE
41	85	65	ILE
41	85	71	GLN
41	85	74	LEU
41	85	83	LEU
41	85	88	ILE
41	85	92	ARG
41	85	95	LEU
41	85	97	ASP
41	85	100	VAL
41	85	105	VAL
42	95	13	ARG
42	95	15	GLU
42	95	19	LYS
42	95	26	ASP
42	95	32	THR
42	95	33	VAL
42	95	38	LEU
42	95	47	VAL
42	95	49	THR
42	95	62	LEU
42	95	66	ARG
42	95	70	ILE
42	95	74	LYS
42	95	83	ARG
42	95	85	LYS
42	95	89	GLN
42	95	91	TYR
42	95	95	LEU
43	A5	6	ILE
43	A5	11	ARG
43	A5	23	LEU
43	A5	28	SER
43	A5	35	ILE

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Mol	Chain	Res	Type
43	A5	37	ARG
43	A5	39	THR
43	A5	41	LYS
43	A5	51	LEU
43	A5	65	LEU
43	A5	67	ASP
43	A5	76	VAL
43	A5	86	LEU
43	A5	88	ARG
43	A5	90	ARG
43	A5	94	ASP
43	A5	95	ILE
43	A5	96	ILE
43	A5	97	LYS
43	A5	100	THR
43	A5	106	ILE
43	A5	107	LEU
43	A5	110	LYS
44	B5	3	THR
44	B5	12	VAL
44	B5	15	GLU
44	B5	23	GLU
44	B5	25	LYS
44	B5	27	THR
44	B5	30	VAL
44	B5	35	THR
44	B5	36	LYS
44	B5	37	THR
44	B5	40	LYS
44	B5	45	THR
44	B5	49	VAL
44	B5	52	VAL
44	B5	54	VAL
44	B5	60	ARG
44	B5	63	LYS
44	B5	65	ARG
44	B5	69	TYR
44	B5	70	LEU
44	B5	76	ARG
44	B5	78	LYS
44	B5	87	GLN
44	B5	88	LYS

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Mol	Chain	Res	Type
45	C5	14	LEU
45	C5	29	GLU
45	C5	31	LEU
45	C5	33	LYS
45	C5	37	VAL
45	C5	38	ILE
45	C5	43	ASN
45	C5	44	ILE
45	C5	47	LYS
45	C5	50	ARG
45	C5	51	VAL
45	C5	55	TYR
45	C5	60	PHE
45	C5	62	GLU
45	C5	68	HIS
45	C5	84	ARG
45	C5	85	VAL
45	C5	86	ARG
45	C5	87	LYS
45	C5	88	LYS
45	C5	89	PHE
45	C5	90	LEU
45	C5	92	ASN
45	C5	97	ARG
45	C5	99	CYS
45	C5	102	CYS
46	D5	2	GLU
46	D5	4	ARG
46	D5	5	LEU
46	D5	19	ARG
46	D5	24	LEU
46	D5	27	VAL
46	D5	28	MET
46	D5	31	ARG
46	D5	33	LEU
46	D5	36	LYS
46	D5	41	LEU
46	D5	44	PHE
46	D5	53	ILE
46	D5	56	VAL
46	D5	59	LEU
46	D5	63	ASP

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Mol	Chain	Res	Type
46	D5	70	LEU
46	D5	72	ARG
46	D5	73	GLN
46	D5	74	VAL
46	D5	76	LEU
46	D5	77	ASP
46	D5	84	GLU
46	D5	87	ASP
46	D5	88	PHE
46	D5	89	PHE
46	D5	94	GLU
46	D5	103	ARG
46	D5	107	THR
46	D5	117	LEU
46	D5	121	HIS
46	D5	122	ARG
46	D5	123	ASP
46	D5	136	PHE
46	D5	144	LEU
46	D5	161	VAL
46	D5	165	VAL
46	D5	168	GLU
46	D5	175	VAL
47	E5	10	THR
47	E5	12	ASN
47	E5	14	ARG
47	E5	19	LYS
47	E5	36	ILE
47	E5	38	VAL
47	E5	43	THR
47	E5	44	ARG
47	E5	55	ARG
47	E5	64	ASP
47	E5	70	GLN
48	F5	11	ARG
48	F5	13	ILE
48	F5	21	ARG
48	F5	25	LYS
48	F5	35	THR
48	F5	38	SER
48	F5	46	LEU
48	F5	56	GLN

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Mol	Chain	Res	Type
48	F5	62	VAL
48	F5	76	ARG
48	F5	78	LYS
48	F5	82	LEU
48	F5	83	GLU
48	F5	85	LEU
48	F5	90	ILE
49	G5	4	SER
49	G5	5	GLU
49	G5	14	ARG
49	G5	15	LYS
49	G5	16	LEU
49	G5	17	SER
49	G5	24	LEU
49	G5	32	LEU
49	G5	44	LEU
49	G5	46	GLN
49	G5	48	HIS
49	G5	49	LYS
49	G5	53	LEU
49	G5	55	ARG
49	G5	60	LEU
49	G5	66	GLU
50	H5	5	LYS
50	H5	7	LYS
50	H5	8	LEU
50	H5	13	ILE
50	H5	17	LYS
50	H5	18	ASP
50	H5	23	LEU
50	H5	35	ARG
50	H5	37	LEU
50	H5	40	THR
50	H5	44	ARG
50	H5	48	GLU
50	H5	55	ARG
51	I5	1	MET
51	I5	8	LYS
51	I5	18	CYS
51	I5	21	VAL
51	I5	22	ILE
51	I5	30	GLU

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Mol	Chain	Res	Type
51	I5	39	CYS
51	I5	44	THR
51	I5	59	PHE
51	I5	60	GLN
51	I5	61	ARG
52	J5	3	LYS
52	J5	4	HIS
52	J5	10	LYS
52	J5	15	ARG
52	J5	25	LEU
52	J5	29	THR
52	J5	33	CYS
52	J5	35	GLU
52	J5	36	CYS
52	J5	48	GLU
52	J5	55	ARG
53	K5	10	LEU
53	K5	12	GLU
53	K5	20	ASN
53	K5	27	LYS
53	K5	28	ARG
53	K5	29	ASN
53	K5	34	LEU
53	K5	47	THR
54	L5	1	MET
54	L5	2	LYS
54	L5	4	THR
54	L5	8	ASN
54	L5	14	LYS
54	L5	41	ARG
54	L5	42	LEU
54	L5	43	THR
55	M5	15	LYS
55	M5	23	VAL
55	M5	25	MET
55	M5	31	HIS
55	M5	33	ASN
55	M5	34	TRP
55	M5	37	SER
55	M5	46	ARG
55	M5	50	LEU
55	M5	52	LYS

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Mol	Chain	Res	Type
55	M5	58	ILE
55	M5	60	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
3	2E	6	HIS
3	2E	108	ASN
4	3E	43	HIS
6	5E	64	GLN
10	1I	84	GLN
16	7I	16	HIS
28	11	227	ASN
29	21	143	ASN
30	31	40	GLN
35	68	88	ASN
36	78	13	ASN
37	88	12	GLN
38	98	13	HIS
43	E8	62	HIS
44	F8	31	HIS
44	F8	82	GLN
47	I8	70	GLN
2	12	212	GLN
13	4A	77	ASN
13	4A	92	HIS
19	AA	23	ASN
32	59	74	ASN
34	15	38	HIS
37	45	113	GLN
38	55	91	GLN
45	C5	92	ASN
51	I5	6	HIS
51	I5	20	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	13	1495/1522 (98%)	358 (23%)	29 (1%)
1	1G	1495/1522 (98%)	366 (24%)	36 (2%)

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
22	1K	74/76 (97%)	36 (48%)	1 (1%)
22	1L	74/76 (97%)	32 (43%)	3 (4%)
23	2K	76/77 (98%)	21 (27%)	3 (3%)
23	2L	76/77 (98%)	19 (25%)	4 (5%)
24	3K	75/76 (98%)	39 (52%)	8 (10%)
24	3L	75/76 (98%)	41 (54%)	3 (4%)
25	4K	13/30 (43%)	6 (46%)	1 (7%)
25	4L	6/30 (20%)	2 (33%)	1 (16%)
26	14	2908/2917 (99%)	777 (26%)	51 (1%)
26	1H	2911/2917 (99%)	743 (25%)	61 (2%)
27	16	121/122 (99%)	21 (17%)	1 (0%)
27	1J	121/122 (99%)	35 (28%)	2 (1%)
All	All	9520/9640 (98%)	2496 (26%)	204 (2%)

All (2496) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	13	5	U
1	13	6	G
1	13	8	A
1	13	9	G
1	13	28	G
1	13	32	A
1	13	39	G
1	13	41	G
1	13	44	G
1	13	47	C
1	13	48	C
1	13	50	A
1	13	51	A
1	13	54	C
1	13	61	G
1	13	65	U
1	13	66	G
1	13	77	C
1	13	78	G
1	13	79	G
1	13	91	C
1	13	95	G
1	13	101	A
1	13	113	G
1	13	116	A

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Mol	Chain	Res	Type
1	13	117	G
1	13	121	C
1	13	131	C
1	13	138	G
1	13	144	G
1	13	147	G
1	13	151	A
1	13	161	A
1	13	163	C
1	13	169	C
1	13	172	A
1	13	173	U
1	13	174	C
1	13	175	C
1	13	189	U
1	13	190	G
1	13	191(A)	G
1	13	195	A
1	13	197	A
1	13	199	G
1	13	201	C
1	13	208	U
1	13	209	U
1	13	210	U
1	13	217	C
1	13	222	U
1	13	226	G
1	13	231	G
1	13	243	A
1	13	244	U
1	13	245	C
1	13	247	G
1	13	251	G
1	13	253	U
1	13	256	U
1	13	262	A
1	13	266	G
1	13	267	C
1	13	273	A
1	13	289	G
1	13	316	G
1	13	318	G

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Mol	Chain	Res	Type
1	13	319	G
1	13	321	A
1	13	324	G
1	13	328	C
1	13	329	A
1	13	330	C
1	13	332	G
1	13	341	C
1	13	342	C
1	13	344	A
1	13	345	C
1	13	346	G
1	13	347	G
1	13	351	G
1	13	352	C
1	13	353	A
1	13	354	G
1	13	357	G
1	13	367	U
1	13	372	C
1	13	388	G
1	13	390	C
1	13	392	G
1	13	397	A
1	13	398	C
1	13	406	G
1	13	412	A
1	13	413	G
1	13	414	A
1	13	415	A
1	13	418	C
1	13	419	C
1	13	421	U
1	13	423	G
1	13	424	G
1	13	428	G
1	13	429	U
1	13	430	A
1	13	439	A
1	13	451	A
1	13	452	A
1	13	454	C

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Mol	Chain	Res	Type
1	13	466	C
1	13	467	G
1	13	485	G
1	13	487	A
1	13	496	A
1	13	497	U
1	13	498	A
1	13	505	G
1	13	506	G
1	13	509	A
1	13	510	A
1	13	511	C
1	13	513	C
1	13	518	C
1	13	519	C
1	13	524	G
1	13	527	G
1	13	531	U
1	13	532	A
1	13	533	A
1	13	536	C
1	13	547	A
1	13	549	C
1	13	559	A
1	13	561	U
1	13	562	C
1	13	567	G
1	13	572	A
1	13	573	A
1	13	576	G
1	13	577	G
1	13	581	G
1	13	607	A
1	13	610	G
1	13	616	G
1	13	617	G
1	13	619	U
1	13	630	G
1	13	631	G
1	13	632	A
1	13	633	G
1	13	639	G

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Mol	Chain	Res	Type
1	13	650	G
1	13	655	A
1	13	661	G
1	13	665	A
1	13	666	G
1	13	704	A
1	13	723	U
1	13	724	G
1	13	734	G
1	13	747	C
1	13	748	C
1	13	749	C
1	13	750	G
1	13	753	A
1	13	755	G
1	13	763	G
1	13	777	A
1	13	786	G
1	13	792	A
1	13	793	U
1	13	794	A
1	13	805	C
1	13	812	C
1	13	813	U
1	13	817	C
1	13	818	G
1	13	828	A
1	13	841	U
1	13	842	C
1	13	843	U
1	13	848	C
1	13	853	G
1	13	859	A
1	13	864	A
1	13	870	U
1	13	874	G
1	13	882	C
1	13	884	U
1	13	885	G
1	13	887	G
1	13	902	G
1	13	914	A

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Mol	Chain	Res	Type
1	13	926	G
1	13	927	G
1	13	934	C
1	13	935	A
1	13	936	C
1	13	940	C
1	13	949	A
1	13	953	G
1	13	960	U
1	13	968	A
1	13	969	A
1	13	971	G
1	13	974	A
1	13	975	A
1	13	976	G
1	13	977	A
1	13	982	U
1	13	984	C
1	13	991	U
1	13	992	U
1	13	993	G
1	13	994	A
1	13	995	C
1	13	999	U
1	13	1004	A
1	13	1005	A
1	13	1006	C
1	13	1007	C
1	13	1008	C
1	13	1009	G
1	13	1017	G
1	13	1021	G
1	13	1024	G
1	13	1025	U
1	13	1028	C
1	13	1029	G
1	13	1032(A)	G
1	13	1032(B)	G
1	13	1033	G
1	13	1040	U
1	13	1042	G
1	13	1044	A

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Mol	Chain	Res	Type
1	13	1046	A
1	13	1055	A
1	13	1065	U
1	13	1066	C
1	13	1081	G
1	13	1092	A
1	13	1094	G
1	13	1095	U
1	13	1101	A
1	13	1123	A
1	13	1124	G
1	13	1125	U
1	13	1126	U
1	13	1127	G
1	13	1129	C
1	13	1130	A
1	13	1131	G
1	13	1136	U
1	13	1137	C
1	13	1138	G
1	13	1139	G
1	13	1143	G
1	13	1146	A
1	13	1152	A
1	13	1154	G
1	13	1155	G
1	13	1157	A
1	13	1158	C
1	13	1159	U
1	13	1160	G
1	13	1161	C
1	13	1164	G
1	13	1165	C
1	13	1177	G
1	13	1178	G
1	13	1181	G
1	13	1184	G
1	13	1189	C
1	13	1194	U
1	13	1196	U
1	13	1197	G
1	13	1201	A

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Mol	Chain	Res	Type
1	13	1202	G
1	13	1203	C
1	13	1212	U
1	13	1213	A
1	13	1225	A
1	13	1227	A
1	13	1238	A
1	13	1240	U
1	13	1246	C
1	13	1253	G
1	13	1256	A
1	13	1257	U
1	13	1258	G
1	13	1260	C
1	13	1261	A
1	13	1262	C
1	13	1270	C
1	13	1272	G
1	13	1273	G
1	13	1280	A
1	13	1281	U
1	13	1282	C
1	13	1285	A
1	13	1286	A
1	13	1287	A
1	13	1299	A
1	13	1300	G
1	13	1302	U
1	13	1305	G
1	13	1320	C
1	13	1322	C
1	13	1323	G
1	13	1326	C
1	13	1331	G
1	13	1335	C
1	13	1336	C
1	13	1337	G
1	13	1338	G
1	13	1340	A
1	13	1346	A
1	13	1347	G
1	13	1350	A

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Mol	Chain	Res	Type
1	13	1353	G
1	13	1360	A
1	13	1363	A
1	13	1364	U
1	13	1368	G
1	13	1370	G
1	13	1377	A
1	13	1396	A
1	13	1401	G
1	13	1402	C
1	13	1406	U
1	13	1419	G
1	13	1442	G
1	13	1443	G
1	13	1446	A
1	13	1451	A
1	13	1452	C
1	13	1453	G
1	13	1454	G
1	13	1455	G
1	13	1467	G
1	13	1469	G
1	13	1487	G
1	13	1492	A
1	13	1494	G
1	13	1495	U
1	13	1499	A
1	13	1503	A
1	13	1504	G
1	13	1505	G
1	13	1506	U
1	13	1507	A
1	13	1509	C
1	13	1514	C
1	13	1517	G
1	13	1519	A
1	13	1520	G
1	13	1529	G
1	13	1530	G
22	1K	3	C
22	1K	4	C
22	1K	9	A

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Mol	Chain	Res	Type
22	1K	10	G
22	1K	11	C
22	1K	16	U
22	1K	17	C
22	1K	18	G
22	1K	19	G
22	1K	20	U
22	1K	21	A
22	1K	22	G
22	1K	24	G
22	1K	26	A
22	1K	34	G
22	1K	41	C
22	1K	44	G
22	1K	45	U
22	1K	46	G
22	1K	47	U
22	1K	48	C
22	1K	49	C
22	1K	52	G
22	1K	56	C
22	1K	59	U
22	1K	62	C
22	1K	64	A
22	1K	66	U
22	1K	68	C
22	1K	69	G
22	1K	70	G
22	1K	72	C
22	1K	73	A
22	1K	74	C
22	1K	75	C
22	1K	76	A
23	2K	2	G
23	2K	6	G
23	2K	9	G
23	2K	13	C
23	2K	15	G
23	2K	18	C
23	2K	20	G
23	2K	21	U
23	2K	22	A

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Mol	Chain	Res	Type
23	2K	23	G
23	2K	35	C
23	2K	44	A
23	2K	47	7MG
23	2K	48	U
23	2K	49	C
23	2K	50	G
23	2K	53	G
23	2K	62	C
23	2K	68	C
23	2K	73	A
23	2K	77	A
24	3K	2	C
24	3K	3	C
24	3K	4	C
24	3K	5	G
24	3K	7	A
24	3K	9	A
24	3K	10	G
24	3K	13	C
24	3K	14	A
24	3K	17	C
24	3K	19	G
24	3K	20	U
24	3K	21	A
24	3K	22	G
24	3K	24	G
24	3K	26	A
24	3K	33	U
24	3K	34	G
24	3K	35	A
24	3K	36	A
24	3K	37	A
24	3K	45	U
24	3K	46	G
24	3K	47	U
24	3K	48	C
24	3K	49	C
24	3K	52	G
24	3K	55	U
24	3K	56	C
24	3K	58	A

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Mol	Chain	Res	Type
24	3K	59	U
24	3K	60	U
24	3K	61	C
24	3K	65	G
24	3K	69	G
24	3K	70	G
24	3K	72	C
24	3K	73	A
24	3K	76	A
25	4K	14	A
25	4K	15	A
25	4K	19	C
25	4K	22	A
25	4K	25	A
25	4K	26	A
26	1H	5	A
26	1H	7	G
26	1H	9	U
26	1H	12	U
26	1H	17	G
26	1H	27	G
26	1H	34	C
26	1H	37	C
26	1H	46	C
26	1H	51	G
26	1H	54	G
26	1H	56	A
26	1H	63	U
26	1H	71	A
26	1H	72	U
26	1H	74	A
26	1H	75	G
26	1H	85	G
26	1H	93	C
26	1H	95	G
26	1H	102	G
26	1H	118	A
26	1H	119	A
26	1H	120	U
26	1H	123	G
26	1H	125	G
26	1H	138	G

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Mol	Chain	Res	Type
26	1H	147	U
26	1H	155	C
26	1H	163	U
26	1H	164	U
26	1H	181	A
26	1H	188	G
26	1H	196	A
26	1H	197	A
26	1H	199	A
26	1H	200	U
26	1H	214	G
26	1H	215	G
26	1H	216	A
26	1H	222	A
26	1H	223	A
26	1H	228	A
26	1H	229	A
26	1H	230	U
26	1H	233	A
26	1H	235	U
26	1H	243	U
26	1H	244	A
26	1H	245	G
26	1H	248	G
26	1H	250	G
26	1H	252	G
26	1H	261	G
26	1H	269	U
26	1H	270(F)	U
26	1H	270(G)	C
26	1H	270(M)	U
26	1H	270(P)	C
26	1H	271(B)	G
26	1H	271(C)	U
26	1H	271	G
26	1H	274	G
26	1H	275	G
26	1H	278	A
26	1H	299	A
26	1H	308	G
26	1H	311	A
26	1H	323	G

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Mol	Chain	Res	Type
26	1H	324	A
26	1H	326	G
26	1H	329	G
26	1H	330	A
26	1H	342	G
26	1H	345	A
26	1H	346	A
26	1H	347	A
26	1H	352	G
26	1H	363	G
26	1H	364	C
26	1H	372	G
26	1H	382	G
26	1H	386	G
26	1H	389	G
26	1H	393	C
26	1H	405	U
26	1H	407	G
26	1H	411	G
26	1H	413	C
26	1H	428	A
26	1H	435	C
26	1H	436	C
26	1H	444	C
26	1H	445	C
26	1H	446	G
26	1H	447	A
26	1H	448	U
26	1H	451	C
26	1H	454	A
26	1H	455	C
26	1H	457	A
26	1H	460	A
26	1H	462	C
26	1H	467	G
26	1H	470	A
26	1H	471	A
26	1H	481	G
26	1H	482	A
26	1H	488	G
26	1H	501	A
26	1H	505	A

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Mol	Chain	Res	Type
26	1H	508	G
26	1H	509	C
26	1H	513	A
26	1H	529	A
26	1H	531	C
26	1H	532	A
26	1H	533	G
26	1H	545	G
26	1H	546	C
26	1H	549	G
26	1H	556	G
26	1H	563	G
26	1H	564	C
26	1H	567	A
26	1H	573	G
26	1H	574	C
26	1H	575	A
26	1H	583	G
26	1H	586	A
26	1H	587	C
26	1H	588	U
26	1H	603	A
26	1H	607	U
26	1H	609(A)	G
26	1H	614	U
26	1H	617	G
26	1H	621	A
26	1H	622	G
26	1H	627	A
26	1H	632	A
26	1H	637	A
26	1H	644	A
26	1H	645	C
26	1H	646	A
26	1H	654	A
26	1H	654(A)	A
26	1H	654(H)	G
26	1H	654(I)	C
26	1H	654(J)	A
26	1H	654(K)	C
26	1H	654(M)	C
26	1H	654(N)	G

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Mol	Chain	Res	Type
26	1H	654(O)	G
26	1H	654(T)	A
26	1H	664	C
26	1H	667	U
26	1H	669	G
26	1H	686	G
26	1H	699	A
26	1H	715	G
26	1H	717	G
26	1H	730	C
26	1H	731	C
26	1H	739	G
26	1H	752	A
26	1H	753	C
26	1H	762	U
26	1H	765	G
26	1H	776	G
26	1H	777	A
26	1H	782	A
26	1H	783	A
26	1H	784	A
26	1H	785	G
26	1H	790	C
26	1H	791	C
26	1H	792	G
26	1H	793	A
26	1H	801	G
26	1H	805	G
26	1H	812	C
26	1H	824	A
26	1H	827	U
26	1H	828	U
26	1H	836	G
26	1H	846	C
26	1H	847	U
26	1H	853	G
26	1H	859	G
26	1H	861	A
26	1H	863	A
26	1H	866	A
26	1H	879	G
26	1H	880	G

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Mol	Chain	Res	Type
26	1H	881	G
26	1H	882	G
26	1H	884	C
26	1H	885	C
26	1H	886	C
26	1H	887	A
26	1H	888	C
26	1H	890	A
26	1H	892	G
26	1H	893	C
26	1H	894	C
26	1H	895	U
26	1H	899	A
26	1H	900	A
26	1H	901	A
26	1H	902	C
26	1H	907	U
26	1H	910	A
26	1H	917	A
26	1H	918	A
26	1H	925	C
26	1H	926	A
26	1H	932	G
26	1H	938	G
26	1H	940	G
26	1H	941	A
26	1H	946	G
26	1H	947	G
26	1H	953	A
26	1H	958	U
26	1H	959	A
26	1H	961	C
26	1H	974	G
26	1H	974(A)	C
26	1H	980	A
26	1H	983	A
26	1H	990	A
26	1H	996	A
26	1H	997	G
26	1H	1003	G
26	1H	1005	C
26	1H	1008	C

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Mol	Chain	Res	Type
26	1H	1011	G
26	1H	1012	U
26	1H	1013	C
26	1H	1015	G
26	1H	1016	G
26	1H	1020	A
26	1H	1022	G
26	1H	1023	U
26	1H	1024	G
26	1H	1025	G
26	1H	1026	U
26	1H	1027	A
26	1H	1033	U
26	1H	1041	C
26	1H	1046	A
26	1H	1047	G
26	1H	1056	G
26	1H	1057	A
26	1H	1060	U
26	1H	1061	U
26	1H	1062	G
26	1H	1064	C
26	1H	1066	U
26	1H	1068	G
26	1H	1070	A
26	1H	1071	G
26	1H	1072	C
26	1H	1073	A
26	1H	1076	C
26	1H	1078	U
26	1H	1082	U
26	1H	1084	A
26	1H	1085	A
26	1H	1086	A
26	1H	1087	G
26	1H	1088	A
26	1H	1089	G
26	1H	1090	U
26	1H	1095	A
26	1H	1096	A
26	1H	1098	A
26	1H	1111	A

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Mol	Chain	Res	Type
26	1H	1112	G
26	1H	1122	G
26	1H	1126	A
26	1H	1129	A
26	1H	1130	U
26	1H	1131	G
26	1H	1133	U
26	1H	1135	C
26	1H	1136	G
26	1H	1139	G
26	1H	1142	U
26	1H	1142(A)	A
26	1H	1143	A
26	1H	1149	G
26	1H	1150	C
26	1H	1154	G
26	1H	1155	A
26	1H	1156	A
26	1H	1157	G
26	1H	1173	G
26	1H	1175	U
26	1H	1176	G
26	1H	1177	A
26	1H	1178	C
26	1H	1179	C
26	1H	1180	C
26	1H	1194	A
26	1H	1195	G
26	1H	1204	A
26	1H	1205	U
26	1H	1218	C
26	1H	1220	A
26	1H	1221	C
26	1H	1231	G
26	1H	1234	U
26	1H	1237	A
26	1H	1241	A
26	1H	1242	A
26	1H	1244	G
26	1H	1245	G
26	1H	1250	G
26	1H	1253	A

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Mol	Chain	Res	Type
26	1H	1255	U
26	1H	1256	G
26	1H	1265	A
26	1H	1269	A
26	1H	1271	G
26	1H	1272	A
26	1H	1273	U
26	1H	1274	A
26	1H	1275	A
26	1H	1280	G
26	1H	1287	A
26	1H	1292	U
26	1H	1293	C
26	1H	1298	C
26	1H	1300	U
26	1H	1301	A
26	1H	1302	A
26	1H	1313	U
26	1H	1314	C
26	1H	1319	G
26	1H	1329	U
26	1H	1332	G
26	1H	1334	G
26	1H	1338	G
26	1H	1342	A
26	1H	1345	C
26	1H	1348	G
26	1H	1349	A
26	1H	1351	C
26	1H	1352	U
26	1H	1358	G
26	1H	1359	A
26	1H	1360	A
26	1H	1365	A
26	1H	1368	G
26	1H	1369	G
26	1H	1370	C
26	1H	1378	A
26	1H	1379	A
26	1H	1380	G
26	1H	1383	C
26	1H	1384	A

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Mol	Chain	Res	Type
26	1H	1385	G
26	1H	1386	C
26	1H	1391	U
26	1H	1395	A
26	1H	1397	U
26	1H	1403	C
26	1H	1406	U
26	1H	1416	G
26	1H	1420	U
26	1H	1421	G
26	1H	1428	C
26	1H	1431	U
26	1H	1437	C
26	1H	1444(A)	A
26	1H	1449	A
26	1H	1449(A)	G
26	1H	1454	U
26	1H	1455	G
26	1H	1458	C
26	1H	1459	G
26	1H	1460	A
26	1H	1461	G
26	1H	1467	C
26	1H	1471	A
26	1H	1472	A
26	1H	1483	G
26	1H	1493	C
26	1H	1497	U
26	1H	1506	C
26	1H	1507	A
26	1H	1508	A
26	1H	1509	C
26	1H	1510	A
26	1H	1511	A
26	1H	1517	G
26	1H	1522	G
26	1H	1526	G
26	1H	1534	G
26	1H	1535	U
26	1H	1536	A
26	1H	1537	C
26	1H	1538	G

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Mol	Chain	Res	Type
26	1H	1540	G
26	1H	1543	A
26	1H	1545	A
26	1H	1548	C
26	1H	1554	A
26	1H	1558	A
26	1H	1559	G
26	1H	1560	G
26	1H	1566	A
26	1H	1567	A
26	1H	1569	A
26	1H	1578	U
26	1H	1579	A
26	1H	1580	A
26	1H	1586	A
26	1H	1587	A
26	1H	1592	C
26	1H	1594	G
26	1H	1608	A
26	1H	1609	A
26	1H	1610	A
26	1H	1617	C
26	1H	1634	A
26	1H	1640	C
26	1H	1647	G
26	1H	1648	C
26	1H	1651	G
26	1H	1664	A
26	1H	1666	G
26	1H	1674	G
26	1H	1684	C
26	1H	1685	C
26	1H	1694	C
26	1H	1695	G
26	1H	1699	G
26	1H	1707	G
26	1H	1728	G
26	1H	1729	A
26	1H	1731	G
26	1H	1743	G
26	1H	1756	G
26	1H	1762	A

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Mol	Chain	Res	Type
26	1H	1763	G
26	1H	1764	G
26	1H	1765	C
26	1H	1773	A
26	1H	1786	A
26	1H	1791	A
26	1H	1799	G
26	1H	1800	C
26	1H	1801	G
26	1H	1802	A
26	1H	1816	G
26	1H	1819	A
26	1H	1829	A
26	1H	1833	U
26	1H	1835	G
26	1H	1836	C
26	1H	1839	G
26	1H	1847	A
26	1H	1858	G
26	1H	1859	A
26	1H	1870	C
26	1H	1878	G
26	1H	1889	A
26	1H	1896	G
26	1H	1900	A
26	1H	1901	A
26	1H	1904	G
26	1H	1906	G
26	1H	1914	C
26	1H	1915	U
26	1H	1919	A
26	1H	1924	C
26	1H	1926	U
26	1H	1929	G
26	1H	1930	G
26	1H	1931	U
26	1H	1932	A
26	1H	1938	A
26	1H	1951	U
26	1H	1952	A
26	1H	1953	A
26	1H	1955	U

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Mol	Chain	Res	Type
26	1H	1960	A
26	1H	1963	U
26	1H	1967	C
26	1H	1969	A
26	1H	1970	A
26	1H	1971	A
26	1H	1972	A
26	1H	1976	U
26	1H	1982	C
26	1H	1983	C
26	1H	1993	U
26	1H	2023	G
26	1H	2030	A
26	1H	2031	A
26	1H	2032	G
26	1H	2033	A
26	1H	2036	C
26	1H	2043	C
26	1H	2047	U
26	1H	2048	G
26	1H	2049	G
26	1H	2051	A
26	1H	2054	A
26	1H	2055	C
26	1H	2056	G
26	1H	2058	A
26	1H	2060	A
26	1H	2061	G
26	1H	2062	A
26	1H	2063	C
26	1H	2069	G
26	1H	2070	G
26	1H	2087	G
26	1H	2098	U
26	1H	2108	C
26	1H	2110	G
26	1H	2111	C
26	1H	2112	G
26	1H	2113	U
26	1H	2114	A
26	1H	2115	G
26	1H	2116	G

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Mol	Chain	Res	Type
26	1H	2118	U
26	1H	2122	U
26	1H	2126	A
26	1H	2127	G
26	1H	2128	C
26	1H	2131	G
26	1H	2132	U
26	1H	2133	G
26	1H	2135	A
26	1H	2136	C
26	1H	2138	C
26	1H	2139	C
26	1H	2145	C
26	1H	2147	G
26	1H	2148	G
26	1H	2154	G
26	1H	2157	G
26	1H	2158	A
26	1H	2161	C
26	1H	2166	G
26	1H	2167	U
26	1H	2168	G
26	1H	2169	A
26	1H	2170	A
26	1H	2171	A
26	1H	2173	A
26	1H	2176	A
26	1H	2177	C
26	1H	2181	G
26	1H	2190	G
26	1H	2198	A
26	1H	2207	C
26	1H	2210	G
26	1H	2211	G
26	1H	2212	A
26	1H	2213	U
26	1H	2215	G
26	1H	2225	A
26	1H	2226	C
26	1H	2237	G
26	1H	2238	G
26	1H	2240	C

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Mol	Chain	Res	Type
26	1H	2264	C
26	1H	2267	A
26	1H	2271	G
26	1H	2273	A
26	1H	2275	C
26	1H	2278	A
26	1H	2280	G
26	1H	2283	C
26	1H	2285	C
26	1H	2286	A
26	1H	2287	A
26	1H	2288	A
26	1H	2294	C
26	1H	2298	A
26	1H	2305	A
26	1H	2307	G
26	1H	2308	G
26	1H	2310	A
26	1H	2312	U
26	1H	2314	C
26	1H	2320	A
26	1H	2321	G
26	1H	2325	G
26	1H	2327	A
26	1H	2334	G
26	1H	2336	A
26	1H	2340	G
26	1H	2343	C
26	1H	2346	A
26	1H	2347	C
26	1H	2350	C
26	1H	2352	A
26	1H	2360	A
26	1H	2361	A
26	1H	2372	G
26	1H	2377	A
26	1H	2379	G
26	1H	2383	G
26	1H	2385	C
26	1H	2388	A
26	1H	2389	G
26	1H	2392	A

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Mol	Chain	Res	Type
26	1H	2393	A
26	1H	2395	C
26	1H	2402	C
26	1H	2403	C
26	1H	2406	U
26	1H	2413	G
26	1H	2414	G
26	1H	2424	C
26	1H	2425	A
26	1H	2428	G
26	1H	2429	G
26	1H	2430	A
26	1H	2431	U
26	1H	2439	A
26	1H	2440	C
26	1H	2441	C
26	1H	2448	A
26	1H	2468	G
26	1H	2469	A
26	1H	2474	C
26	1H	2475	C
26	1H	2476	A
26	1H	2478	A
26	1H	2479	G
26	1H	2482	G
26	1H	2484	G
26	1H	2487	G
26	1H	2494	G
26	1H	2502	G
26	1H	2505	G
26	1H	2506	U
26	1H	2507	C
26	1H	2518	A
26	1H	2525	G
26	1H	2529	G
26	1H	2531	A
26	1H	2553	G
26	1H	2554	U
26	1H	2566	A
26	1H	2567	G
26	1H	2568	C
26	1H	2569	G

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Mol	Chain	Res	Type
26	1H	2570	G
26	1H	2573	C
26	1H	2574	G
26	1H	2579	C
26	1H	2582	G
26	1H	2583	G
26	1H	2584	U
26	1H	2601	C
26	1H	2602	A
26	1H	2609	U
26	1H	2611	U
26	1H	2612	C
26	1H	2615	U
26	1H	2629	A
26	1H	2630	G
26	1H	2632	A
26	1H	2636	U
26	1H	2637	U
26	1H	2641	G
26	1H	2646	C
26	1H	2654	A
26	1H	2657	A
26	1H	2665	A
26	1H	2666	C
26	1H	2672	G
26	1H	2679	A
26	1H	2689	U
26	1H	2690	C
26	1H	2691	C
26	1H	2700	C
26	1H	2701	C
26	1H	2702	U
26	1H	2703	C
26	1H	2705	A
26	1H	2707	G
26	1H	2712(A)	A
26	1H	2713	A
26	1H	2714	G
26	1H	2718	G
26	1H	2719	G
26	1H	2726	U
26	1H	2729	G

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Mol	Chain	Res	Type
26	1H	2733	A
26	1H	2744	G
26	1H	2749	A
26	1H	2750	A
26	1H	2752	C
26	1H	2757	A
26	1H	2764	A
26	1H	2765	A
26	1H	2766	G
26	1H	2778	A
26	1H	2779	U
26	1H	2781	A
26	1H	2787	C
26	1H	2789	C
26	1H	2790	A
26	1H	2791	C
26	1H	2793	G
26	1H	2794	C
26	1H	2795	G
26	1H	2797	U
26	1H	2798	C
26	1H	2799	A
26	1H	2801	A
26	1H	2802	G
26	1H	2808	U
26	1H	2813	A
26	1H	2818	G
26	1H	2820	A
26	1H	2821	A
26	1H	2830	G
26	1H	2833	G
26	1H	2834	G
26	1H	2835	A
26	1H	2838	G
26	1H	2849	U
26	1H	2850	A
26	1H	2851	A
26	1H	2872	G
26	1H	2891	G
26	1H	2892	A
26	1H	2893	G
26	1H	2894	G

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Mol	Chain	Res	Type
26	1H	2895	U
26	1H	2899	G
27	16	1	U
27	16	5	C
27	16	7	G
27	16	13	A
27	16	15	A
27	16	25	A
27	16	35	U
27	16	42	C
27	16	45	A
27	16	56	G
27	16	65	C
27	16	66	A
27	16	73	A
27	16	74	U
27	16	81	G
27	16	82	G
27	16	83	G
27	16	105	G
27	16	109	G
27	16	117	G
27	16	119	A
1	1G	5	U
1	1G	6	G
1	1G	9	G
1	1G	15	G
1	1G	21	G
1	1G	26	A
1	1G	32	A
1	1G	39	G
1	1G	41	G
1	1G	42	G
1	1G	44	G
1	1G	47	C
1	1G	48	C
1	1G	50	A
1	1G	51	A
1	1G	54	C
1	1G	65	U
1	1G	90	C
1	1G	91	C

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Mol	Chain	Res	Type
1	1G	95	G
1	1G	101	A
1	1G	115	G
1	1G	116	A
1	1G	121	C
1	1G	127	G
1	1G	131	C
1	1G	146	G
1	1G	162	A
1	1G	163	C
1	1G	167	G
1	1G	174	C
1	1G	182	U
1	1G	183	G
1	1G	188	U
1	1G	189	U
1	1G	190	G
1	1G	191(A)	G
1	1G	191(D)	U
1	1G	191(E)	G
1	1G	195	A
1	1G	196	A
1	1G	197	A
1	1G	198	G
1	1G	208	U
1	1G	209	U
1	1G	210	U
1	1G	216	G
1	1G	217	C
1	1G	231	G
1	1G	240	C
1	1G	247	G
1	1G	249	U
1	1G	250	A
1	1G	251	G
1	1G	266	G
1	1G	267	C
1	1G	281	G
1	1G	283	C
1	1G	289	G
1	1G	298	A
1	1G	299	G

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Mol	Chain	Res	Type
1	1G	316	G
1	1G	317	G
1	1G	318	G
1	1G	321	A
1	1G	328	C
1	1G	329	A
1	1G	332	G
1	1G	345	C
1	1G	346	G
1	1G	347	G
1	1G	349	A
1	1G	350	G
1	1G	351	G
1	1G	352	C
1	1G	353	A
1	1G	354	G
1	1G	356	A
1	1G	363	A
1	1G	366	C
1	1G	367	U
1	1G	372	C
1	1G	373	A
1	1G	388	G
1	1G	397	A
1	1G	398	C
1	1G	406	G
1	1G	411	A
1	1G	412	A
1	1G	413	G
1	1G	419	C
1	1G	422	C
1	1G	423	G
1	1G	424	G
1	1G	429	U
1	1G	430	A
1	1G	439	A
1	1G	442	C
1	1G	452	A
1	1G	465	A
1	1G	466	C
1	1G	467	G
1	1G	475	G

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Mol	Chain	Res	Type
1	1G	476	G
1	1G	482	A
1	1G	484	G
1	1G	485	G
1	1G	486	U
1	1G	496	A
1	1G	497	U
1	1G	498	A
1	1G	500	G
1	1G	505	G
1	1G	509	A
1	1G	510	A
1	1G	511	C
1	1G	518	C
1	1G	521	G
1	1G	524	G
1	1G	527	G
1	1G	528	C
1	1G	529	G
1	1G	530	G
1	1G	531	U
1	1G	532	A
1	1G	533	A
1	1G	534	U
1	1G	536	C
1	1G	544	G
1	1G	545	C
1	1G	547	A
1	1G	559	A
1	1G	560	U
1	1G	561	U
1	1G	562	C
1	1G	564	C
1	1G	566	G
1	1G	572	A
1	1G	573	A
1	1G	576	G
1	1G	577	G
1	1G	581	G
1	1G	596	C
1	1G	607	A
1	1G	614	A

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Mol	Chain	Res	Type
1	1G	618	C
1	1G	620	C
1	1G	630	G
1	1G	632	A
1	1G	633	G
1	1G	642	A
1	1G	651	C
1	1G	652	U
1	1G	653	A
1	1G	660	G
1	1G	665	A
1	1G	670	G
1	1G	687	A
1	1G	688	G
1	1G	702	A
1	1G	704	A
1	1G	710	G
1	1G	723	U
1	1G	724	G
1	1G	731	G
1	1G	742	G
1	1G	748	C
1	1G	764	C
1	1G	777	A
1	1G	778	G
1	1G	792	A
1	1G	793	U
1	1G	794	A
1	1G	805	C
1	1G	813	U
1	1G	817	C
1	1G	821	G
1	1G	828	A
1	1G	841	U
1	1G	842	C
1	1G	843	U
1	1G	848	C
1	1G	859	A
1	1G	871	U
1	1G	885	G
1	1G	889	A
1	1G	891	U

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Mol	Chain	Res	Type
1	1G	902	G
1	1G	914	A
1	1G	916	G
1	1G	926	G
1	1G	927	G
1	1G	934	C
1	1G	935	A
1	1G	936	C
1	1G	940	C
1	1G	958	A
1	1G	960	U
1	1G	961	U
1	1G	968	A
1	1G	969	A
1	1G	971	G
1	1G	972	C
1	1G	974	A
1	1G	975	A
1	1G	976	G
1	1G	977	A
1	1G	978	A
1	1G	980	C
1	1G	981	U
1	1G	982	U
1	1G	984	C
1	1G	991	U
1	1G	992	U
1	1G	993	G
1	1G	994	A
1	1G	1004	A
1	1G	1006	C
1	1G	1009	G
1	1G	1014	A
1	1G	1023	G
1	1G	1024	G
1	1G	1025	U
1	1G	1027	C
1	1G	1028	C
1	1G	1028(B)	C
1	1G	1029	G
1	1G	1031	G
1	1G	1032(A)	G

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Mol	Chain	Res	Type
1	1G	1033	G
1	1G	1036	G
1	1G	1038	C
1	1G	1040	U
1	1G	1046	A
1	1G	1047	G
1	1G	1050	G
1	1G	1053	G
1	1G	1054	C
1	1G	1055	A
1	1G	1064	G
1	1G	1066	C
1	1G	1081	G
1	1G	1082	G
1	1G	1085	U
1	1G	1086	U
1	1G	1092	A
1	1G	1094	G
1	1G	1095	U
1	1G	1101	A
1	1G	1113	C
1	1G	1118	C
1	1G	1124	G
1	1G	1125	U
1	1G	1127	G
1	1G	1128	C
1	1G	1129	C
1	1G	1130	A
1	1G	1131	G
1	1G	1133	G
1	1G	1136	U
1	1G	1137	C
1	1G	1138	G
1	1G	1139	G
1	1G	1141	C
1	1G	1145	C
1	1G	1146	A
1	1G	1147	C
1	1G	1154	G
1	1G	1155	G
1	1G	1157	A
1	1G	1158	C

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Mol	Chain	Res	Type
1	1G	1159	U
1	1G	1160	G
1	1G	1177	G
1	1G	1178	G
1	1G	1181	G
1	1G	1185	G
1	1G	1187	G
1	1G	1188	A
1	1G	1189	C
1	1G	1190	G
1	1G	1196	U
1	1G	1197	G
1	1G	1201	A
1	1G	1211	U
1	1G	1212	U
1	1G	1213	A
1	1G	1214	C
1	1G	1225	A
1	1G	1227	A
1	1G	1236	A
1	1G	1238	A
1	1G	1240	U
1	1G	1241	G
1	1G	1256	A
1	1G	1257	U
1	1G	1258	G
1	1G	1260	C
1	1G	1263	C
1	1G	1267	C
1	1G	1269	A
1	1G	1278	U
1	1G	1279	A
1	1G	1280	A
1	1G	1286	A
1	1G	1287	A
1	1G	1289	A
1	1G	1291	G
1	1G	1295	G
1	1G	1297	C
1	1G	1298	C
1	1G	1299	A
1	1G	1300	G

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Mol	Chain	Res	Type
1	1G	1301	U
1	1G	1303	C
1	1G	1305	G
1	1G	1313	U
1	1G	1317	C
1	1G	1319	A
1	1G	1320	C
1	1G	1322	C
1	1G	1323	G
1	1G	1331	G
1	1G	1335	C
1	1G	1336	C
1	1G	1337	G
1	1G	1346	A
1	1G	1347	G
1	1G	1353	G
1	1G	1359	C
1	1G	1362(A)	C
1	1G	1363	A
1	1G	1364	U
1	1G	1365	G
1	1G	1370	G
1	1G	1379	G
1	1G	1382	C
1	1G	1396	A
1	1G	1397	C
1	1G	1398	A
1	1G	1399	C
1	1G	1401	G
1	1G	1404	C
1	1G	1405	G
1	1G	1406	U
1	1G	1419	G
1	1G	1442	G
1	1G	1443	G
1	1G	1446	A
1	1G	1450	U
1	1G	1451	A
1	1G	1452	C
1	1G	1453	G
1	1G	1454	G
1	1G	1484	C

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Mol	Chain	Res	Type
1	1G	1492	A
1	1G	1499	A
1	1G	1502	A
1	1G	1503	A
1	1G	1504	G
1	1G	1506	U
1	1G	1510	U
1	1G	1517	G
1	1G	1520	G
1	1G	1529	G
1	1G	1530	G
22	1L	8	U
22	1L	9	A
22	1L	10	G
22	1L	11	C
22	1L	13	C
22	1L	16	U
22	1L	17	C
22	1L	18	G
22	1L	19	G
22	1L	21	A
22	1L	22	G
22	1L	25	C
22	1L	26	A
22	1L	27	G
22	1L	36	A
22	1L	46	G
22	1L	47	U
22	1L	48	C
22	1L	51	U
22	1L	56	C
22	1L	58	A
22	1L	59	U
22	1L	61	C
22	1L	64	A
22	1L	68	C
22	1L	69	G
22	1L	70	G
22	1L	72	C
22	1L	73	A
22	1L	74	C
22	1L	75	C

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Mol	Chain	Res	Type
22	1L	76	A
23	2L	2	G
23	2L	7	G
23	2L	8	4SU
23	2L	9	G
23	2L	16	C
23	2L	17	C
23	2L	19	G
23	2L	21	U
23	2L	22	A
23	2L	23	G
23	2L	27	G
23	2L	32	G
23	2L	47	7MG
23	2L	48	U
23	2L	49	C
23	2L	61	U
23	2L	68	C
23	2L	70	C
23	2L	77	A
24	3L	5	G
24	3L	6	G
24	3L	7	A
24	3L	9	A
24	3L	13	C
24	3L	16	U
24	3L	17	C
24	3L	18	G
24	3L	19	G
24	3L	20	U
24	3L	21	A
24	3L	22	G
24	3L	23	A
24	3L	24	G
24	3L	26	A
24	3L	27	G
24	3L	30	G
24	3L	31	A
24	3L	34	G
24	3L	36	A
24	3L	37	A
24	3L	38	A

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Mol	Chain	Res	Type
24	3L	39	U
24	3L	40	C
24	3L	41	C
24	3L	44	G
24	3L	46	G
24	3L	47	U
24	3L	48	C
24	3L	49	C
24	3L	53	G
24	3L	54	U
24	3L	58	A
24	3L	59	U
24	3L	61	C
24	3L	65	G
24	3L	66	U
24	3L	68	C
24	3L	72	C
24	3L	73	A
24	3L	76	A
25	4L	14	A
25	4L	15	A
26	14	4	C
26	14	5	A
26	14	6	A
26	14	9	U
26	14	14	A
26	14	15	G
26	14	34	C
26	14	35	G
26	14	36	G
26	14	41	C
26	14	46	C
26	14	49	A
26	14	50	U
26	14	54	G
26	14	55	G
26	14	58	G
26	14	70	G
26	14	71	A
26	14	72	U
26	14	74	A
26	14	75	G

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Mol	Chain	Res	Type
26	14	76	C
26	14	78	A
26	14	82	G
26	14	83	G
26	14	84	A
26	14	91	A
26	14	93	C
26	14	99	U
26	14	102	G
26	14	112	U
26	14	118	A
26	14	119	A
26	14	120	U
26	14	121	G
26	14	129	C
26	14	138	G
26	14	140	A
26	14	152	G
26	14	154	G
26	14	155	C
26	14	161	U
26	14	162	U
26	14	176	G
26	14	181	A
26	14	183	C
26	14	196	A
26	14	199	A
26	14	205	G
26	14	214	G
26	14	216	A
26	14	218	A
26	14	221	A
26	14	222	A
26	14	225	A
26	14	229	A
26	14	233	A
26	14	240	G
26	14	248	G
26	14	249	C
26	14	250	G
26	14	252	G
26	14	265	A

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Mol	Chain	Res	Type
26	14	266	G
26	14	270(G)	C
26	14	270(K)	C
26	14	270(L)	U
26	14	270(M)	U
26	14	270(O)	U
26	14	271(B)	G
26	14	271(C)	U
26	14	271	G
26	14	273(D)	C
26	14	273(F)	C
26	14	274	G
26	14	275	G
26	14	276	A
26	14	277	C
26	14	278	A
26	14	279	C
26	14	283	A
26	14	289	A
26	14	311	A
26	14	324	A
26	14	327	G
26	14	329	G
26	14	330	A
26	14	331	A
26	14	335	C
26	14	336	C
26	14	346	A
26	14	352	G
26	14	363	G
26	14	363(A)	A
26	14	363(E)	U
26	14	386	G
26	14	391	G
26	14	396	G
26	14	399	G
26	14	405	U
26	14	406	G
26	14	411	G
26	14	412	A
26	14	428	A
26	14	443	A

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Mol	Chain	Res	Type
26	14	444	C
26	14	447	A
26	14	448	U
26	14	451	C
26	14	454	A
26	14	455	C
26	14	457	A
26	14	459	U
26	14	460	A
26	14	470	A
26	14	471	A
26	14	479	A
26	14	481	G
26	14	498	G
26	14	501	A
26	14	504	U
26	14	505	A
26	14	509	C
26	14	510	C
26	14	513	A
26	14	528	A
26	14	529	A
26	14	531	C
26	14	532	A
26	14	533	G
26	14	537	C
26	14	549	G
26	14	556	G
26	14	563	G
26	14	573	G
26	14	575	A
26	14	587	C
26	14	599	G
26	14	602	G
26	14	603	A
26	14	607	U
26	14	609(A)	G
26	14	613	U
26	14	614	U
26	14	615	G
26	14	617	G
26	14	618	G

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Mol	Chain	Res	Type
26	14	618(A)	C
26	14	620	G
26	14	621	A
26	14	622	G
26	14	627	A
26	14	637	A
26	14	645	C
26	14	646	A
26	14	650	C
26	14	651	G
26	14	654	A
26	14	654(E)	C
26	14	654(G)	C
26	14	654(H)	G
26	14	654(I)	C
26	14	654(K)	C
26	14	654(L)	G
26	14	654(N)	G
26	14	654(T)	A
26	14	654(V)	A
26	14	667	U
26	14	669	G
26	14	672	C
26	14	676	A
26	14	686	G
26	14	689	A
26	14	699	A
26	14	715	G
26	14	717	G
26	14	720	C
26	14	722	A
26	14	730	C
26	14	748	G
26	14	750	A
26	14	752	A
26	14	753	C
26	14	757	U
26	14	758	C
26	14	762	U
26	14	765	G
26	14	769	G
26	14	771	G

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Mol	Chain	Res	Type
26	14	775	G
26	14	776	G
26	14	779	U
26	14	782	A
26	14	784	A
26	14	785	G
26	14	792	G
26	14	797	C
26	14	805	G
26	14	812	C
26	14	819	A
26	14	820	A
26	14	827	U
26	14	828	U
26	14	831	G
26	14	832	G
26	14	836	G
26	14	845	G
26	14	846	C
26	14	847	U
26	14	858	U
26	14	859	G
26	14	863	A
26	14	865	C
26	14	866	A
26	14	869	G
26	14	880	G
26	14	881	G
26	14	882	G
26	14	885	C
26	14	886	C
26	14	887	A
26	14	888	C
26	14	889	C
26	14	890	A
26	14	894	C
26	14	896	A
26	14	897	C
26	14	899	A
26	14	900	A
26	14	901	A
26	14	902	C

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Mol	Chain	Res	Type
26	14	903	C
26	14	904	C
26	14	905	U
26	14	910	A
26	14	914	C
26	14	915	C
26	14	917	A
26	14	918	A
26	14	919	G
26	14	932	G
26	14	933	A
26	14	938	G
26	14	941	A
26	14	945	A
26	14	946	G
26	14	952	G
26	14	957	A
26	14	958	U
26	14	959	A
26	14	961	C
26	14	968	G
26	14	972	G
26	14	974	G
26	14	980	A
26	14	983	A
26	14	986	C
26	14	989	G
26	14	990	A
26	14	991	C
26	14	996	A
26	14	1002	G
26	14	1005	C
26	14	1012	U
26	14	1013	C
26	14	1017	G
26	14	1020	A
26	14	1022	G
26	14	1023	U
26	14	1025	G
26	14	1026	U
26	14	1028	A
26	14	1034	G

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Mol	Chain	Res	Type
26	14	1035	U
26	14	1037	G
26	14	1038	C
26	14	1039	G
26	14	1044	G
26	14	1045	A
26	14	1046	A
26	14	1047	G
26	14	1048	A
26	14	1050	A
26	14	1051	G
26	14	1054	A
26	14	1056	G
26	14	1057	A
26	14	1058	U
26	14	1059	G
26	14	1060	U
26	14	1062	G
26	14	1065	U
26	14	1067	A
26	14	1068	G
26	14	1070	A
26	14	1073	A
26	14	1077	A
26	14	1079	C
26	14	1082	U
26	14	1083	U
26	14	1085	A
26	14	1086	A
26	14	1087	G
26	14	1088	A
26	14	1090	U
26	14	1091	G
26	14	1093	G
26	14	1094	U
26	14	1095	A
26	14	1096	A
26	14	1097	U
26	14	1098	A
26	14	1105	U
26	14	1110	G
26	14	1111	A

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Mol	Chain	Res	Type
26	14	1112	G
26	14	1122	G
26	14	1129	A
26	14	1130	U
26	14	1135	C
26	14	1136	G
26	14	1138	G
26	14	1139	G
26	14	1141	U
26	14	1142	U
26	14	1142(A)	A
26	14	1143	A
26	14	1144	G
26	14	1149	G
26	14	1155	A
26	14	1157	G
26	14	1160	G
26	14	1167	U
26	14	1170	G
26	14	1173	G
26	14	1174	A
26	14	1175	U
26	14	1177	A
26	14	1178	C
26	14	1183	G
26	14	1204	A
26	14	1205	U
26	14	1206	G
26	14	1210	A
26	14	1211	U
26	14	1212	G
26	14	1220	A
26	14	1229(A)	G
26	14	1237	A
26	14	1244	G
26	14	1247	A
26	14	1248	G
26	14	1249	U
26	14	1250	G
26	14	1253	A
26	14	1256	G
26	14	1267	U

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Mol	Chain	Res	Type
26	14	1271	G
26	14	1272	A
26	14	1275	A
26	14	1287	A
26	14	1298	C
26	14	1300	U
26	14	1301	A
26	14	1306	C
26	14	1313	U
26	14	1314	C
26	14	1316	U
26	14	1318	C
26	14	1319	G
26	14	1321	A
26	14	1329	U
26	14	1332	G
26	14	1338	G
26	14	1342	A
26	14	1345	C
26	14	1347	G
26	14	1349	A
26	14	1359	A
26	14	1360	A
26	14	1365	A
26	14	1368	G
26	14	1370	C
26	14	1378	A
26	14	1379	A
26	14	1383	C
26	14	1384	A
26	14	1385	G
26	14	1386	C
26	14	1395	A
26	14	1403	C
26	14	1406	U
26	14	1416	G
26	14	1417	C
26	14	1419	A
26	14	1420	U
26	14	1421	G
26	14	1427	A
26	14	1428	C

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Mol	Chain	Res	Type
26	14	1437	C
26	14	1444(A)	A
26	14	1445	C
26	14	1449	A
26	14	1449(A)	G
26	14	1455	G
26	14	1459	G
26	14	1460	A
26	14	1461	G
26	14	1467	C
26	14	1471	A
26	14	1475	G
26	14	1483	G
26	14	1490	A
26	14	1493	C
26	14	1494	A
26	14	1506	C
26	14	1508	A
26	14	1509	C
26	14	1510	A
26	14	1515	C
26	14	1522	G
26	14	1524	G
26	14	1528	A
26	14	1533	C
26	14	1535	U
26	14	1537	C
26	14	1538	G
26	14	1543	A
26	14	1554	A
26	14	1558	A
26	14	1559	G
26	14	1560	G
26	14	1569	A
26	14	1578	U
26	14	1583	A
26	14	1585	C
26	14	1586	A
26	14	1588	C
26	14	1589	C
26	14	1598	C
26	14	1608	A

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Mol	Chain	Res	Type
26	14	1609	A
26	14	1610	A
26	14	1614	A
26	14	1616	A
26	14	1617	C
26	14	1618	A
26	14	1622	G
26	14	1625	C
26	14	1631	A
26	14	1632	A
26	14	1636	C
26	14	1640	C
26	14	1641	A
26	14	1644	C
26	14	1647	G
26	14	1648	C
26	14	1654	A
26	14	1661	G
26	14	1669	A
26	14	1670	C
26	14	1671	U
26	14	1674	G
26	14	1678	G
26	14	1694	C
26	14	1697	G
26	14	1700	A
26	14	1701	A
26	14	1703	G
26	14	1725	G
26	14	1726	G
26	14	1728	G
26	14	1729	A
26	14	1730	U
26	14	1742	C
26	14	1750	G
26	14	1756	G
26	14	1761	C
26	14	1762	A
26	14	1763	G
26	14	1764	G
26	14	1773	A
26	14	1780	A

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Mol	Chain	Res	Type
26	14	1781	C
26	14	1782	C
26	14	1791	A
26	14	1800	C
26	14	1801	G
26	14	1802	A
26	14	1812	A
26	14	1816	G
26	14	1819	A
26	14	1820	U
26	14	1829	A
26	14	1830	C
26	14	1839	G
26	14	1847	A
26	14	1848	A
26	14	1858	G
26	14	1878	G
26	14	1884	A
26	14	1888	G
26	14	1889	A
26	14	1895	C
26	14	1906	G
26	14	1909	C
26	14	1917	U
26	14	1919	A
26	14	1929	G
26	14	1930	G
26	14	1933	G
26	14	1936	A
26	14	1937	A
26	14	1938	A
26	14	1955	U
26	14	1960	A
26	14	1963	U
26	14	1964	G
26	14	1967	C
26	14	1968	G
26	14	1969	A
26	14	1970	A
26	14	1971	A
26	14	1972	A
26	14	1985	G

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Mol	Chain	Res	Type
26	14	1993	U
26	14	1996	C
26	14	2018	G
26	14	2019	A
26	14	2023	G
26	14	2031	A
26	14	2032	G
26	14	2033	A
26	14	2039	C
26	14	2043	C
26	14	2048	G
26	14	2049	G
26	14	2053	G
26	14	2055	C
26	14	2056	G
26	14	2060	A
26	14	2061	G
26	14	2062	A
26	14	2063	C
26	14	2069	G
26	14	2082	A
26	14	2096	U
26	14	2100	G
26	14	2105	C
26	14	2108	C
26	14	2111	C
26	14	2112	G
26	14	2114	A
26	14	2117	A
26	14	2118	U
26	14	2119	A
26	14	2122	U
26	14	2125	G
26	14	2126	A
26	14	2127	G
26	14	2128	C
26	14	2131	G
26	14	2132	U
26	14	2133	G
26	14	2136	C
26	14	2137	C
26	14	2140	C

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Mol	Chain	Res	Type
26	14	2144	U
26	14	2145	C
26	14	2146	C
26	14	2147	G
26	14	2148	G
26	14	2153	G
26	14	2157	G
26	14	2158	A
26	14	2160	G
26	14	2171	A
26	14	2173	A
26	14	2174	C
26	14	2177	C
26	14	2183	C
26	14	2188	C
26	14	2189	U
26	14	2191	G
26	14	2192	G
26	14	2198	A
26	14	2210	G
26	14	2211	G
26	14	2212	A
26	14	2213	U
26	14	2215	G
26	14	2225	A
26	14	2226	C
26	14	2238	G
26	14	2239	G
26	14	2240	C
26	14	2251	G
26	14	2252	G
26	14	2261	C
26	14	2267	A
26	14	2268	A
26	14	2273	A
26	14	2275	C
26	14	2276	G
26	14	2278	A
26	14	2280	G
26	14	2283	C
26	14	2287	A
26	14	2288	A

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Mol	Chain	Res	Type
26	14	2289	G
26	14	2305	A
26	14	2307	G
26	14	2308	G
26	14	2309	A
26	14	2310	A
26	14	2311	A
26	14	2312	U
26	14	2318	G
26	14	2321	G
26	14	2324	C
26	14	2325	G
26	14	2327	A
26	14	2333	A
26	14	2334	G
26	14	2335	A
26	14	2336	A
26	14	2342	C
26	14	2343	C
26	14	2346	A
26	14	2347	C
26	14	2350	C
26	14	2352	A
26	14	2353	G
26	14	2354	G
26	14	2355	C
26	14	2383	G
26	14	2385	C
26	14	2388	A
26	14	2394	C
26	14	2396	G
26	14	2401	U
26	14	2402	C
26	14	2403	C
26	14	2406	U
26	14	2407	G
26	14	2410	G
26	14	2414	G
26	14	2422	A
26	14	2423	U
26	14	2425	A
26	14	2429	G

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Mol	Chain	Res	Type
26	14	2430	A
26	14	2432	A
26	14	2434	A
26	14	2435	A
26	14	2439	A
26	14	2440	C
26	14	2441	C
26	14	2445	G
26	14	2447	G
26	14	2448	A
26	14	2449	U
26	14	2465	C
26	14	2468	G
26	14	2469	A
26	14	2470	G
26	14	2474	C
26	14	2476	A
26	14	2477	C
26	14	2478	A
26	14	2480	C
26	14	2482	G
26	14	2483	C
26	14	2484	G
26	14	2487	G
26	14	2489	G
26	14	2496	C
26	14	2497	A
26	14	2502	G
26	14	2504	U
26	14	2505	G
26	14	2506	U
26	14	2507	C
26	14	2514	U
26	14	2518	A
26	14	2519	U
26	14	2527	C
26	14	2529	G
26	14	2532	G
26	14	2542	A
26	14	2543	G
26	14	2552	U
26	14	2554	U

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Mol	Chain	Res	Type
26	14	2555	U
26	14	2566	A
26	14	2567	G
26	14	2569	G
26	14	2573	C
26	14	2577	A
26	14	2585	U
26	14	2587	A
26	14	2601	C
26	14	2602	A
26	14	2603	G
26	14	2608	G
26	14	2609	U
26	14	2611	U
26	14	2612	C
26	14	2613	U
26	14	2615	U
26	14	2617	C
26	14	2630	G
26	14	2636	U
26	14	2646	C
26	14	2647	U
26	14	2654	A
26	14	2655	G
26	14	2663	G
26	14	2665	A
26	14	2667	C
26	14	2673	G
26	14	2679	A
26	14	2689	U
26	14	2690	C
26	14	2691	C
26	14	2700	C
26	14	2702	U
26	14	2703	C
26	14	2707	G
26	14	2712(A)	A
26	14	2713	A
26	14	2714	G
26	14	2718	G
26	14	2719	G
26	14	2726	U

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Mol	Chain	Res	Type
26	14	2729	G
26	14	2733	A
26	14	2736	G
26	14	2744	G
26	14	2748	A
26	14	2750	A
26	14	2751	G
26	14	2752	C
26	14	2754	U
26	14	2758	A
26	14	2762	G
26	14	2764	A
26	14	2765	A
26	14	2766	G
26	14	2777	G
26	14	2778	A
26	14	2779	U
26	14	2789	C
26	14	2790	A
26	14	2791	C
26	14	2795	G
26	14	2797	U
26	14	2798	C
26	14	2802	G
26	14	2808	U
26	14	2811	G
26	14	2818	G
26	14	2820	A
26	14	2821	A
26	14	2833	G
26	14	2834	G
26	14	2835	A
26	14	2849	U
26	14	2860	A
26	14	2861	G
26	14	2872	G
26	14	2873	A
26	14	2880	C
26	14	2886	G
26	14	2889	C
26	14	2894	G
26	14	2896	C

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Mol	Chain	Res	Type
27	1J	0	A
27	1J	2	C
27	1J	5	C
27	1J	7	G
27	1J	8	U
27	1J	12	C
27	1J	13	A
27	1J	15	A
27	1J	16	G
27	1J	19	G
27	1J	22	U
27	1J	24	G
27	1J	29	A
27	1J	30	C
27	1J	32	C
27	1J	41	U
27	1J	42	C
27	1J	44	G
27	1J	45	A
27	1J	47	C
27	1J	52	A
27	1J	53	A
27	1J	58	A
27	1J	73	A
27	1J	81	G
27	1J	88	C
27	1J	89	G
27	1J	89(A)	A
27	1J	90	C
27	1J	100	G
27	1J	101	A
27	1J	108	C
27	1J	109	G
27	1J	113	C
27	1J	114	G

All (204) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	13	5	U
1	13	50	A
1	13	115	G

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Mol	Chain	Res	Type
1	13	244	U
1	13	251	G
1	13	266	G
1	13	412	A
1	13	422	C
1	13	428	G
1	13	429	U
1	13	484	G
1	13	509	A
1	13	560	U
1	13	703	G
1	13	748	C
1	13	793	U
1	13	812	C
1	13	991	U
1	13	992	U
1	13	1027	C
1	13	1065	U
1	13	1126	U
1	13	1285	A
1	13	1302	U
1	13	1336	C
1	13	1452	C
1	13	1453	G
1	13	1498	U
1	13	1504	G
22	1K	10	G
23	2K	21	U
23	2K	47	7MG
23	2K	48	U
24	3K	2	C
24	3K	8	U
24	3K	18	G
24	3K	33	U
24	3K	36	A
24	3K	58	A
24	3K	60	U
24	3K	71	G
25	4K	25	A
26	1H	125	G
26	1H	196	A
26	1H	199	A

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Mol	Chain	Res	Type
26	1H	222	A
26	1H	229	A
26	1H	242	G
26	1H	271(B)	G
26	1H	404	C
26	1H	481	G
26	1H	508	G
26	1H	528	A
26	1H	587	C
26	1H	668	G
26	1H	685	A
26	1H	752	A
26	1H	764	A
26	1H	776	G
26	1H	800	A
26	1H	880	G
26	1H	974	G
26	1H	1022	G
26	1H	1026	U
26	1H	1060	U
26	1H	1084	A
26	1H	1085	A
26	1H	1110	G
26	1H	1178	C
26	1H	1241	A
26	1H	1252	G
26	1H	1301	A
26	1H	1312	U
26	1H	1378	A
26	1H	1396	U
26	1H	1420	U
26	1H	1508	A
26	1H	1558	A
26	1H	1608	A
26	1H	1609	A
26	1H	1694	C
26	1H	1699	G
26	1H	1799	G
26	1H	1800	C
26	1H	1900	A
26	1H	1980	G
26	1H	2060	A

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Mol	Chain	Res	Type
26	1H	2062	A
26	1H	2111	C
26	1H	2135	A
26	1H	2157	G
26	1H	2167	U
26	1H	2225	A
26	1H	2286	A
26	1H	2307	G
26	1H	2426	A
26	1H	2428	G
26	1H	2475	C
26	1H	2481	G
26	1H	2566	A
26	1H	2611	U
26	1H	2689	U
26	1H	2756	U
27	16	108	C
1	1G	115	G
1	1G	173	U
1	1G	197	A
1	1G	201	C
1	1G	250	A
1	1G	266	G
1	1G	327	A
1	1G	328	C
1	1G	345	C
1	1G	412	A
1	1G	429	U
1	1G	465	A
1	1G	485	G
1	1G	509	A
1	1G	528	C
1	1G	560	U
1	1G	632	A
1	1G	687	A
1	1G	723	U
1	1G	793	U
1	1G	812	C
1	1G	913	A
1	1G	992	U
1	1G	1054	C
1	1G	1064	G

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Mol	Chain	Res	Type
1	1G	1128	C
1	1G	1157	A
1	1G	1212	U
1	1G	1285	A
1	1G	1297	C
1	1G	1298	C
1	1G	1300	G
1	1G	1346	A
1	1G	1396	A
1	1G	1453	G
1	1G	1498	U
22	1L	10	G
22	1L	18	G
22	1L	73	A
23	2L	20	G
23	2L	21	U
23	2L	46	G
23	2L	48	U
24	3L	8	U
24	3L	58	A
24	3L	71	G
25	4L	13	A
26	14	34	C
26	14	49	A
26	14	123	G
26	14	128	C
26	14	196	A
26	14	197	A
26	14	278	A
26	14	310	A
26	14	385	C
26	14	503	A
26	14	528	A
26	14	685	A
26	14	686	G
26	14	752	A
26	14	764	A
26	14	783	A
26	14	886	C
26	14	960	A
26	14	974	G
26	14	990	A

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Mol	Chain	Res	Type
26	14	1085	A
26	14	1141	U
26	14	1378	A
26	14	1420	U
26	14	1460	A
26	14	1558	A
26	14	1608	A
26	14	1609	A
26	14	1647	G
26	14	1819	A
26	14	1984	G
26	14	2191	G
26	14	2210	G
26	14	2211	G
26	14	2225	A
26	14	2238	G
26	14	2275	C
26	14	2335	A
26	14	2406	U
26	14	2425	A
26	14	2439	A
26	14	2447	G
26	14	2506	U
26	14	2602	A
26	14	2629	A
26	14	2689	U
26	14	2726	U
26	14	2776	A
26	14	2778	A
26	14	2859	G
26	14	2893	G
27	1J	52	A
27	1J	88	C

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

16 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
22	PSU	1K	32	22	15,21,22	4.06	5 (33%)	16,30,33	2.07	3 (18%)
22	MIA	1K	37	22	22,31,32	0.93	2 (9%)	26,44,47	1.10	2 (7%)
22	PSU	1K	39	22	15,21,22	3.78	4 (26%)	16,30,33	2.11	2 (12%)
22	PSU	1L	32	22	15,21,22	3.67	5 (33%)	16,30,33	2.09	4 (25%)
22	MIA	1L	37	22	22,31,32	1.22	2 (9%)	26,44,47	2.63	6 (23%)
22	PSU	1L	39	22	15,21,22	4.06	5 (33%)	16,30,33	2.13	2 (12%)
23	OMC	2K	33	23	15,22,23	1.99	4 (26%)	20,31,34	1.87	2 (10%)
23	7MG	2K	47	23	20,26,27	3.33	6 (30%)	23,39,42	2.49	9 (39%)
23	5MU	2K	55	56,23	13,22,23	1.69	2 (15%)	16,32,35	1.40	2 (12%)
23	PSU	2K	56	23	15,21,22	4.16	4 (26%)	16,30,33	1.76	3 (18%)
23	4SU	2K	8	23	12,21,22	3.27	2 (16%)	15,30,33	0.58	0
23	OMC	2L	33	23	15,22,23	2.16	4 (26%)	20,31,34	1.98	3 (15%)
23	7MG	2L	47	23	20,26,27	3.37	5 (25%)	23,39,42	2.16	7 (30%)
23	5MU	2L	55	23	13,22,23	1.69	2 (15%)	16,32,35	1.36	1 (6%)
23	PSU	2L	56	23	15,21,22	3.97	5 (33%)	16,30,33	2.01	3 (18%)
23	4SU	2L	8	23	12,21,22	3.20	2 (16%)	15,30,33	0.79	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
22	PSU	1K	32	22	-	0/7/25/26	0/2/2/2
22	MIA	1K	37	22	-	0/11/33/34	0/3/3/3
22	PSU	1K	39	22	-	0/7/25/26	0/2/2/2
22	PSU	1L	32	22	-	0/7/25/26	0/2/2/2
22	MIA	1L	37	22	-	0/11/33/34	0/3/3/3
22	PSU	1L	39	22	-	0/7/25/26	0/2/2/2
23	OMC	2K	33	23	-	0/5/27/28	0/2/2/2
23	7MG	2K	47	23	-	0/7/37/38	0/3/3/3
23	5MU	2K	55	56,23	-	0/3/25/26	0/2/2/2
23	PSU	2K	56	23	-	0/7/25/26	0/2/2/2
23	4SU	2K	8	23	-	0/3/25/26	0/2/2/2
23	OMC	2L	33	23	-	0/5/27/28	0/2/2/2
23	7MG	2L	47	23	-	0/7/37/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
23	5MU	2L	55	23	-	0/3/25/26	0/2/2/2
23	PSU	2L	56	23	-	0/7/25/26	0/2/2/2
23	4SU	2L	8	23	-	0/3/25/26	0/2/2/2

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
23	2K	56	PSU	C5-C1'	-14.09	1.40	1.52
22	1K	32	PSU	C5-C1'	-13.59	1.40	1.52
22	1L	39	PSU	C5-C1'	-12.98	1.41	1.52
23	2L	56	PSU	C5-C1'	-12.76	1.41	1.52
22	1K	39	PSU	C5-C1'	-12.46	1.41	1.52
22	1L	32	PSU	C5-C1'	-11.38	1.42	1.52
23	2K	47	7MG	C5-C4	-5.96	1.23	1.39
23	2L	47	7MG	C5-C4	-5.78	1.23	1.39
23	2L	55	5MU	C4-N3	-3.18	1.27	1.33
23	2K	55	5MU	C4-N3	-2.70	1.28	1.33
22	1K	37	MIA	C4-N3	-2.12	1.32	1.35
22	1K	39	PSU	C2-N1	2.06	1.42	1.38
22	1K	32	PSU	C2-N3	2.11	1.42	1.38
22	1L	39	PSU	C2-N3	2.24	1.42	1.38
22	1K	37	MIA	C2-S10	2.29	1.77	1.75
22	1L	32	PSU	C2-N3	2.38	1.43	1.38
23	2L	56	PSU	C2-N3	2.40	1.43	1.38
22	1K	32	PSU	C2-N1	2.50	1.43	1.38
23	2K	47	7MG	C2-N1	2.50	1.40	1.35
23	2K	56	PSU	C2-N1	2.65	1.43	1.38
22	1L	37	MIA	C6-N1	2.82	1.36	1.33
22	1L	39	PSU	C2-N1	2.85	1.44	1.38
22	1L	32	PSU	C2-N1	2.88	1.44	1.38
23	2K	56	PSU	C4-N3	3.06	1.38	1.33
23	2L	56	PSU	C2-N1	3.16	1.44	1.38
23	2K	33	OMC	C4-N4	3.21	1.44	1.35
23	2L	33	OMC	C4-N4	3.38	1.44	1.35
23	2K	47	7MG	C2-N2	3.41	1.41	1.34
23	2K	33	OMC	C2-N3	3.64	1.45	1.38
23	2L	33	OMC	C2-N3	3.68	1.45	1.38
23	2K	33	OMC	C6-N1	3.69	1.40	1.35
23	2L	47	7MG	C2-N2	3.69	1.41	1.34
22	1K	32	PSU	C4-N3	3.75	1.39	1.33
22	1K	39	PSU	C4-N3	3.77	1.39	1.33
23	2K	33	OMC	C5-C4	4.11	1.50	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
23	2L	33	OMC	C5-C4	4.23	1.50	1.41
22	1L	32	PSU	C4-N3	4.31	1.40	1.33
23	2L	56	PSU	C4-N3	4.32	1.40	1.33
22	1L	37	MIA	C2-S10	4.37	1.79	1.75
23	2K	47	7MG	C6-C5	4.38	1.47	1.41
22	1L	39	PSU	C4-N3	4.44	1.41	1.33
23	2L	47	7MG	C8-N7	4.56	1.64	1.43
23	2K	47	7MG	C8-N7	4.67	1.65	1.43
23	2L	55	5MU	C2-N3	4.78	1.48	1.38
23	2L	33	OMC	C6-N1	4.89	1.42	1.35
23	2K	55	5MU	C2-N3	4.90	1.48	1.38
23	2L	47	7MG	C6-C5	5.01	1.48	1.41
22	1K	32	PSU	C6-N1	5.52	1.46	1.34
23	2K	56	PSU	C6-N1	5.53	1.46	1.34
22	1K	39	PSU	C6-N1	5.69	1.46	1.34
23	2L	56	PSU	C6-N1	5.93	1.46	1.34
22	1L	32	PSU	C6-N1	6.04	1.47	1.34
22	1L	39	PSU	C6-N1	6.26	1.47	1.34
23	2L	8	4SU	C6-N1	6.56	1.44	1.35
23	2K	8	4SU	C6-N1	7.03	1.44	1.35
23	2K	8	4SU	C5-C4	8.64	1.49	1.38
23	2L	8	4SU	C5-C4	8.80	1.49	1.38
23	2K	47	7MG	C4-N3	10.67	1.48	1.34
23	2L	47	7MG	C4-N3	11.06	1.48	1.34

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	2L	47	7MG	C5-C4-N3	-6.35	120.27	126.74
23	2K	47	7MG	C4-N9-C1'	-5.11	114.54	126.65
23	2L	47	7MG	N1-C2-N3	-3.80	119.30	125.51
23	2K	47	7MG	C5-C4-N3	-3.56	123.11	126.74
23	2K	47	7MG	N3-C4-N9	-3.53	122.42	126.98
23	2K	47	7MG	C5-C6-N1	-3.50	118.19	123.39
23	2K	47	7MG	N1-C2-N3	-3.27	120.17	125.51
23	2L	56	PSU	C5-C6-N1	-3.16	119.98	124.38
22	1K	37	MIA	C5-C6-N1	-2.71	117.83	120.58
22	1L	37	MIA	C12-N6-C6	-2.62	120.43	123.46
22	1L	37	MIA	N3-C2-N1	-2.53	122.17	126.84
23	2L	8	4SU	C5-C4-N3	-2.51	120.90	123.56
23	2K	56	PSU	C5-C6-N1	-2.46	120.95	124.38
22	1L	32	PSU	C5-C6-N1	-2.23	121.27	124.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	1L	37	MIA	C5-C6-N1	-2.20	118.35	120.58
22	1L	32	PSU	C5-C1'-C2'	-2.19	111.71	115.44
23	2L	47	7MG	C5-C6-N1	-2.12	120.24	123.39
22	1K	32	PSU	C5-C6-N1	-2.11	121.44	124.38
23	2K	55	5MU	C2'-C1'-N1	-2.10	107.83	113.46
23	2L	33	OMC	C5-C4-N3	-2.09	119.14	121.79
23	2K	56	PSU	O4'-C1'-C2'	2.09	106.95	104.69
23	2L	47	7MG	C2-N3-C4	2.15	120.61	114.50
23	2L	47	7MG	N2-C2-N1	2.16	120.76	117.20
22	1K	39	PSU	O4'-C1'-C2'	2.32	107.20	104.69
22	1K	32	PSU	O4'-C1'-C2'	2.44	107.33	104.69
22	1L	37	MIA	N6-C6-N1	2.50	121.52	118.55
22	1L	39	PSU	O4'-C1'-C2'	2.50	107.39	104.69
22	1L	32	PSU	O4'-C1'-C2'	2.61	107.51	104.69
23	2K	47	7MG	N2-C2-N1	2.67	121.61	117.20
22	1K	37	MIA	C2-N1-C6	2.70	120.56	113.13
23	2L	56	PSU	O4'-C1'-C2'	2.73	107.64	104.69
22	1L	37	MIA	C2-N1-C6	2.88	121.04	113.13
23	2K	33	OMC	N4-C4-N3	3.13	121.96	116.50
23	2L	47	7MG	C6-N1-C2	3.38	119.85	115.88
23	2K	47	7MG	C8-N9-C1'	3.43	132.74	122.43
23	2L	33	OMC	N4-C4-N3	3.51	122.63	116.50
23	2K	55	5MU	C4-N3-C2	4.18	118.65	115.16
23	2K	47	7MG	C6-N1-C2	4.30	120.92	115.88
23	2L	47	7MG	C5-C4-N9	4.43	113.41	106.25
23	2L	55	5MU	C4-N3-C2	4.48	118.89	115.16
23	2K	47	7MG	C5-C4-N9	5.09	114.47	106.25
23	2K	56	PSU	C4-N3-C2	5.18	119.48	115.16
23	2L	56	PSU	C4-N3-C2	6.40	120.50	115.16
22	1L	32	PSU	C4-N3-C2	6.95	120.96	115.16
22	1K	32	PSU	C4-N3-C2	7.06	121.05	115.16
23	2K	33	OMC	C6-C5-C4	7.08	120.21	117.44
23	2L	33	OMC	C6-C5-C4	7.28	120.29	117.44
22	1K	39	PSU	C4-N3-C2	7.38	121.31	115.16
22	1L	39	PSU	C4-N3-C2	7.47	121.39	115.16
22	1L	37	MIA	C11-S10-C2	11.85	110.67	102.31

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

11 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
22	1K	37	MIA	2	0
22	1L	39	PSU	1	0
23	2K	47	7MG	4	0
23	2K	55	5MU	2	0
23	2K	56	PSU	1	0
23	2K	8	4SU	2	0
23	2L	33	OMC	3	0
23	2L	47	7MG	4	0
23	2L	55	5MU	2	0
23	2L	56	PSU	1	0
23	2L	8	4SU	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1245 ligands modelled in this entry, 1245 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	13	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	13	1530:G	O3'	1531:A	P	3.03

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	13	1498/1522 (98%)	-0.66	0 100 100	61, 110, 194, 290	0
1	1G	1497/1522 (98%)	-0.67	2 (0%) 95 95	73, 129, 204, 306	0
2	12	237/256 (92%)	0.59	29 (12%) 5 3	151, 185, 205, 218	0
2	1E	237/256 (92%)	0.42	25 (10%) 8 4	119, 155, 183, 193	0
3	22	206/239 (86%)	1.05	38 (18%) 2 1	149, 166, 190, 202	0
3	2E	205/239 (85%)	0.48	18 (8%) 12 6	98, 122, 159, 171	0
4	32	208/209 (99%)	0.60	15 (7%) 18 9	107, 128, 153, 160	0
4	3E	208/209 (99%)	0.51	19 (9%) 11 5	92, 123, 146, 155	0
5	42	151/162 (93%)	0.67	14 (9%) 11 5	118, 139, 159, 195	0
5	4E	151/162 (93%)	0.55	14 (9%) 11 5	88, 109, 135, 174	0
6	52	101/101 (100%)	0.53	10 (9%) 9 4	92, 115, 133, 160	0
6	5E	101/101 (100%)	0.39	3 (2%) 54 37	87, 110, 132, 154	0
7	62	155/156 (99%)	0.44	19 (12%) 5 3	124, 143, 178, 199	0
7	6E	155/156 (99%)	0.07	14 (9%) 12 5	110, 129, 160, 179	0
8	72	138/138 (100%)	0.30	7 (5%) 32 17	115, 145, 160, 163	0
8	7E	138/138 (100%)	0.46	10 (7%) 18 9	93, 118, 128, 142	0
9	82	127/128 (99%)	0.55	6 (4%) 35 20	123, 171, 189, 197	0
9	8E	127/128 (99%)	-0.07	0 100 100	94, 149, 170, 180	0
10	1A	99/105 (94%)	0.61	10 (10%) 9 4	139, 174, 192, 197	0
10	1I	99/105 (94%)	0.70	11 (11%) 7 3	90, 148, 181, 187	0
11	2A	119/129 (92%)	1.75	45 (37%) 0 0	94, 125, 152, 183	0
11	2I	119/129 (92%)	1.36	26 (21%) 1 1	77, 113, 153, 182	0
12	3A	125/132 (94%)	1.18	30 (24%) 1 1	91, 114, 144, 175	0
12	3I	125/132 (94%)	0.40	6 (4%) 34 20	74, 84, 117, 179	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	4A	117/126 (92%)	0.41	12 (10%) 9 4	127, 171, 192, 200	0
13	4I	118/126 (93%)	0.06	3 (2%) 61 45	86, 127, 148, 156	0
14	5A	58/61 (95%)	0.90	12 (20%) 1 1	152, 166, 179, 184	0
14	5I	60/61 (98%)	-0.20	2 (3%) 50 33	92, 113, 128, 135	0
15	6A	88/89 (98%)	0.08	1 (1%) 82 71	96, 125, 143, 148	0
15	6I	88/89 (98%)	0.37	6 (6%) 20 10	82, 104, 122, 129	0
16	7A	84/88 (95%)	0.10	1 (1%) 81 69	100, 117, 142, 174	0
16	7I	84/88 (95%)	-0.42	0 100 100	110, 126, 158, 174	0
17	8A	100/105 (95%)	0.09	2 (2%) 68 52	105, 122, 138, 160	0
17	8I	100/105 (95%)	0.39	5 (5%) 32 17	97, 114, 125, 129	0
18	9A	72/88 (81%)	1.55	21 (29%) 1 0	106, 128, 167, 178	0
18	9I	72/88 (81%)	1.18	18 (25%) 1 1	90, 113, 143, 175	0
19	AA	78/93 (83%)	0.66	11 (14%) 4 2	160, 189, 203, 207	0
19	AI	81/93 (87%)	0.57	12 (14%) 3 2	103, 127, 156, 165	0
20	BA	99/106 (93%)	0.09	3 (3%) 54 37	97, 118, 150, 162	0
20	BI	99/106 (93%)	-0.45	0 100 100	115, 131, 165, 172	0
21	1B	25/27 (92%)	0.16	0 100 100	125, 153, 169, 189	0
21	1F	25/27 (92%)	-0.74	0 100 100	100, 112, 127, 154	0
22	1K	73/76 (96%)	2.25	31 (42%) 0 0	97, 221, 259, 264	0
22	1L	73/76 (96%)	3.31	38 (52%) 0 0	134, 254, 277, 285	0
23	2K	72/77 (93%)	-0.41	1 (1%) 78 64	73, 95, 128, 139	0
23	2L	72/77 (93%)	-0.26	1 (1%) 78 64	87, 121, 153, 165	0
24	3K	76/76 (100%)	-0.08	4 (5%) 30 16	77, 253, 279, 282	0
24	3L	76/76 (100%)	0.43	6 (7%) 15 8	89, 259, 278, 280	0
25	4K	14/30 (46%)	-0.25	0 100 100	76, 107, 151, 155	0
25	4L	6/30 (20%)	-0.25	0 100 100	100, 108, 158, 161	0
26	14	2909/2917 (99%)	-0.33	38 (1%) 79 67	53, 91, 249, 335	0
26	1H	2912/2917 (99%)	-0.34	30 (1%) 84 74	40, 77, 231, 309	0
27	16	122/122 (100%)	-0.71	1 (0%) 87 79	69, 98, 122, 202	0
27	1J	122/122 (100%)	-0.66	1 (0%) 87 79	89, 133, 162, 212	0
28	11	273/276 (98%)	0.13	1 (0%) 93 90	43, 66, 84, 98	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
28	19	273/276 (98%)	0.08	1 (0%) 93 90	54, 80, 97, 112	0
29	21	205/206 (99%)	0.57	24 (11%) 6 3	49, 94, 143, 161	0
29	29	205/206 (99%)	0.65	26 (12%) 5 2	64, 99, 146, 173	0
30	31	202/210 (96%)	-0.10	3 (1%) 76 62	47, 80, 119, 141	0
30	39	208/210 (99%)	0.66	25 (12%) 6 3	60, 109, 170, 198	0
31	41	181/182 (99%)	0.72	22 (12%) 5 3	83, 109, 145, 156	0
31	49	181/182 (99%)	1.92	78 (43%) 0 0	127, 154, 181, 199	0
32	51	174/180 (96%)	0.54	13 (7%) 17 9	85, 109, 129, 142	0
32	59	170/180 (94%)	1.30	46 (27%) 1 0	155, 205, 228, 248	0
33	61	146/148 (98%)	0.43	12 (8%) 14 7	80, 139, 157, 163	0
33	69	146/148 (98%)	0.81	25 (17%) 2 1	88, 135, 157, 168	0
34	15	138/140 (98%)	0.52	8 (5%) 26 13	85, 116, 149, 165	0
34	58	138/140 (98%)	0.19	6 (4%) 39 23	69, 95, 137, 156	0
35	25	122/122 (100%)	0.52	3 (2%) 61 45	73, 93, 108, 124	0
35	68	122/122 (100%)	0.32	2 (1%) 74 61	61, 80, 99, 112	0
36	35	150/150 (100%)	0.90	30 (20%) 1 1	62, 113, 145, 183	0
36	78	150/150 (100%)	0.28	10 (6%) 21 11	49, 85, 116, 166	0
37	45	141/141 (100%)	0.65	18 (12%) 5 2	81, 111, 145, 155	0
37	88	138/141 (97%)	0.10	3 (2%) 65 49	56, 83, 107, 138	0
38	55	117/118 (99%)	0.09	3 (2%) 59 44	64, 84, 99, 117	0
38	98	118/118 (100%)	0.28	2 (1%) 73 59	63, 87, 109, 120	0
39	65	111/112 (99%)	0.66	14 (12%) 5 3	98, 128, 148, 162	0
39	A8	111/112 (99%)	0.28	5 (4%) 37 21	75, 94, 123, 133	0
40	75	137/146 (93%)	0.12	4 (2%) 55 39	83, 101, 161, 195	0
40	B8	137/146 (93%)	-0.20	2 (1%) 76 62	76, 97, 154, 179	0
41	85	117/118 (99%)	0.44	7 (5%) 25 13	73, 104, 150, 174	0
41	C8	117/118 (99%)	0.04	2 (1%) 73 59	56, 82, 115, 134	0
42	95	101/101 (100%)	0.95	18 (17%) 2 1	71, 134, 155, 175	0
42	D8	101/101 (100%)	0.62	13 (12%) 5 2	61, 111, 141, 155	0
43	A5	113/113 (100%)	0.10	1 (0%) 85 77	66, 80, 117, 174	0
43	E8	113/113 (100%)	0.00	4 (3%) 48 30	58, 76, 112, 159	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
44	B5	93/96 (96%)	0.29	5 (5%) 29 15	71, 89, 113, 121	0
44	F8	94/96 (97%)	0.24	5 (5%) 30 16	56, 73, 97, 117	0
45	C5	104/110 (94%)	1.19	20 (19%) 2 1	89, 122, 160, 166	0
45	G8	104/110 (94%)	0.67	6 (5%) 26 13	75, 98, 142, 149	0
46	D5	179/206 (86%)	2.08	78 (43%) 0 0	124, 161, 240, 246	0
46	H8	175/206 (84%)	0.47	13 (7%) 17 9	86, 128, 212, 218	0
47	E5	77/85 (90%)	0.75	6 (7%) 16 8	77, 96, 116, 162	0
47	I8	80/85 (94%)	0.06	2 (2%) 61 45	58, 77, 114, 117	0
48	F5	97/98 (98%)	0.98	14 (14%) 3 2	63, 89, 143, 161	0
48	J8	97/98 (98%)	0.55	10 (10%) 9 4	54, 77, 128, 161	0
49	G5	66/72 (91%)	0.50	4 (6%) 25 12	89, 112, 130, 148	0
49	K8	67/72 (93%)	-0.00	1 (1%) 76 62	65, 81, 99, 132	0
50	H5	59/60 (98%)	0.83	5 (8%) 13 6	80, 109, 155, 173	0
50	L8	57/60 (95%)	0.59	3 (5%) 30 16	64, 83, 112, 120	0
51	I5	63/71 (88%)	3.84	45 (71%) 0 0	167, 206, 219, 221	0
51	M8	66/71 (92%)	1.62	21 (31%) 1 0	114, 161, 192, 200	0
52	J5	59/60 (98%)	0.17	3 (5%) 32 17	66, 91, 159, 212	0
52	N8	58/60 (96%)	0.71	7 (12%) 6 3	53, 100, 178, 183	0
53	K5	45/54 (83%)	6.26	42 (93%) 0 0	141, 181, 194, 201	0
53	O8	45/54 (83%)	4.74	42 (93%) 0 0	117, 148, 168, 171	0
54	L5	46/49 (93%)	-0.17	0 100 100	53, 63, 75, 87	0
54	P8	48/49 (97%)	-0.26	0 100 100	46, 51, 88, 113	0
55	M5	60/65 (92%)	0.29	2 (3%) 50 33	73, 86, 108, 133	0
55	Q8	60/65 (92%)	0.13	1 (1%) 73 59	61, 73, 104, 119	0
All	All	21056/21694 (97%)	0.14	1397 (6%) 22 11	40, 109, 201, 335	0

All (1397) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
26	14	2901	C	18.2
53	K5	13	CYS	17.4
53	K5	51	GLU	15.3
52	N8	60	VAL	14.5
22	1L	3	C	14.2

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Mol	Chain	Res	Type	RSRZ
26	1H	2900	A	14.1
22	1L	71	G	13.7
30	39	1	MET	13.2
26	1H	2901	C	13.1
53	K5	26	ASN	12.8
48	F5	98	LEU	12.4
26	14	2899	G	12.3
22	1L	72	C	12.0
51	I5	31	ILE	11.9
22	1K	17	C	11.4
53	K5	12	GLU	11.0
22	1L	70	G	11.0
53	K5	25	LYS	11.0
53	K5	42	TRP	10.9
7	62	82	GLY	10.6
46	D5	178	GLU	10.5
51	I5	52	THR	10.4
7	62	81	GLY	10.3
46	D5	147	GLY	10.3
22	1K	71	G	10.2
53	K5	50	ARG	10.2
26	14	2900	A	10.1
18	9A	88	LYS	10.0
11	2I	129	SER	10.0
26	14	2902	C	9.9
22	1L	76	A	9.8
53	K5	52	VAL	9.8
26	14	2797	U	9.7
22	1L	17	C	9.7
53	O8	13	CYS	9.5
7	6E	83	ALA	9.3
11	2I	128	ALA	9.3
22	1L	75	C	9.1
52	N8	59	GLU	9.1
53	O8	20	ASN	9.1
26	1H	4	C	9.1
53	K5	14	THR	9.0
31	49	138	GLN	8.9
12	3I	129	ALA	8.8
22	1K	3	C	8.8
51	I5	47	GLN	8.8
22	1K	70	G	8.8

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Mol	Chain	Res	Type	RSRZ
37	45	1	MET	8.7
48	F5	97	LEU	8.6
53	O8	42	TRP	8.6
26	1H	1	G	8.6
52	J5	59	GLU	8.5
46	D5	179	ASP	8.4
26	1H	2902	C	8.4
22	1K	76	A	8.4
52	J5	60	VAL	8.2
18	9A	87	ARG	8.2
42	95	45	THR	8.2
11	2I	11	LYS	8.1
17	8A	101	ARG	8.0
22	1L	45	U	7.9
7	62	80	VAL	7.9
53	K5	22	ALA	7.9
53	K5	53	LYS	7.9
22	1K	69	G	7.9
51	I5	44	THR	7.9
51	I5	40	HIS	7.8
46	D5	146	ILE	7.8
7	6E	82	GLY	7.8
53	K5	31	PRO	7.8
53	K5	39	TYR	7.8
22	1L	68	C	7.8
36	35	149	GLU	7.7
51	I5	30	GLU	7.7
26	14	654(K)	C	7.6
22	1K	16	U	7.5
30	39	208	GLY	7.5
51	I5	42	PHE	7.5
53	O8	18	ARG	7.4
53	O8	49	HIS	7.4
53	O8	50	ARG	7.4
46	D5	172	ALA	7.3
53	K5	9	LEU	7.3
46	D5	112	ARG	7.3
53	K5	24	GLU	7.3
22	1L	47	U	7.3
32	59	100	GLY	7.3
51	I5	29	PRO	7.3
53	K5	49	HIS	7.2

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Mol	Chain	Res	Type	RSRZ
43	A5	113	LYS	7.2
53	O8	53	LYS	7.2
51	I5	28	LYS	7.2
53	K5	21	TYR	7.2
26	14	2799	A	7.2
29	29	69	LYS	7.2
26	1H	654(I)	C	7.2
11	2I	12	ARG	7.2
29	21	90	THR	7.1
51	I5	43	TYR	7.1
31	49	139	LEU	7.1
7	6E	81	GLY	7.0
36	35	110	TYR	7.0
22	1L	2	C	7.0
22	1L	16	U	6.9
26	14	2795	G	6.9
22	1L	4	C	6.8
29	21	205	ALA	6.8
53	O8	43	CYS	6.8
26	14	2798	C	6.7
42	D8	1	MET	6.7
13	4A	4	ILE	6.7
18	9I	88	LYS	6.7
22	1K	73	A	6.6
12	3I	127	GLU	6.6
53	K5	41	PRO	6.6
11	2A	12	ARG	6.6
7	62	78	ARG	6.6
47	I8	85	ALA	6.5
52	N8	58	LEU	6.5
26	14	654(J)	A	6.5
51	I5	46	GLN	6.5
26	1H	2	G	6.5
51	I5	32	TYR	6.5
7	6E	154	TYR	6.4
12	3A	64	TYR	6.4
11	2A	128	ALA	6.4
46	D5	162	GLU	6.4
53	O8	22	ALA	6.4
53	K5	23	THR	6.4
46	D5	68	PRO	6.4
7	62	85	TYR	6.4

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Mol	Chain	Res	Type	RSRZ
53	K5	36	LEU	6.4
31	49	137	GLU	6.4
53	O8	15	GLU	6.3
51	I5	41	PRO	6.3
46	D5	117	LEU	6.3
53	O8	14	THR	6.3
19	AA	82	GLY	6.2
53	O8	47	THR	6.2
53	O8	48	VAL	6.2
46	H8	113	ALA	6.2
12	3I	128	ALA	6.2
26	1H	2899	G	6.2
26	14	2	G	6.2
50	H5	60	GLU	6.1
26	14	654(I)	C	6.1
26	14	654(L)	G	6.1
22	1L	56	C	6.1
53	K5	11	LEU	6.1
53	K5	16	CYS	6.1
36	78	150	ALA	6.0
51	I5	24	THR	6.0
26	14	4	C	6.0
29	21	72	VAL	6.0
22	1L	73	A	6.0
13	4A	2	ALA	6.0
22	1L	74	C	5.9
53	K5	30	THR	5.9
31	49	90	LEU	5.9
48	J8	98	LEU	5.9
32	59	99	VAL	5.9
46	D5	107	THR	5.9
3	22	53	ALA	5.8
47	E5	85	ALA	5.8
11	2A	50	TYR	5.8
53	K5	20	ASN	5.8
26	1H	654(J)	A	5.7
26	1H	654(K)	C	5.7
24	3L	34	G	5.7
45	C5	47	LYS	5.6
37	45	91	GLU	5.6
51	I5	55	ARG	5.6
36	35	148	LEU	5.6

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Mol	Chain	Res	Type	RSRZ
53	K5	40	CYS	5.6
33	69	4	ILE	5.6
46	D5	154	ASP	5.6
22	1L	1	G	5.6
53	O8	35	GLU	5.5
51	I5	10	VAL	5.5
46	D5	163	LEU	5.5
53	O8	34	LEU	5.5
26	14	3	U	5.5
46	D5	142	SER	5.4
11	2A	13	GLN	5.4
53	O8	44	ARG	5.4
31	49	178	PHE	5.4
31	49	133	LEU	5.4
31	49	39	ILE	5.4
29	21	88	GLY	5.3
53	O8	21	TYR	5.3
22	1K	4	C	5.3
7	6E	84	ASN	5.3
12	3A	128	ALA	5.3
46	D5	149	SER	5.3
43	E8	113	LYS	5.2
46	D5	176	PRO	5.2
31	49	48	GLU	5.2
51	I5	63	TYR	5.2
53	K5	35	GLU	5.2
26	1H	3	U	5.2
29	29	54	GLN	5.2
14	5A	39	LEU	5.2
31	49	155	MET	5.2
31	49	146	TYR	5.2
53	K5	18	ARG	5.1
53	O8	51	GLU	5.1
7	62	86	GLN	5.1
45	C5	58	GLY	5.1
42	D8	36	PRO	5.1
30	39	2	LYS	5.1
45	C5	49	VAL	5.1
29	21	89	ASP	5.1
10	1I	6	ILE	5.1
53	O8	26	ASN	5.1
26	14	2898	U	5.0

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Mol	Chain	Res	Type	RSRZ
31	49	177	GLY	5.0
22	1K	67	C	5.0
51	I5	7	PRO	5.0
40	B8	1	MET	5.0
22	1L	20	U	5.0
3	22	146	ALA	5.0
48	J8	92	LYS	5.0
53	O8	16	CYS	5.0
46	D5	141	VAL	5.0
53	K5	10	LEU	4.9
12	3A	68	ALA	4.9
51	M8	34	GLU	4.9
2	12	240	GLN	4.9
12	3A	32	PHE	4.9
46	D5	148	ASP	4.9
51	M8	31	ILE	4.9
37	45	90	VAL	4.9
51	I5	51	ASP	4.9
45	C5	59	GLY	4.9
31	49	37	VAL	4.8
7	62	153	HIS	4.8
46	D5	153	SER	4.8
11	2A	129	SER	4.8
48	F5	96	LYS	4.8
53	O8	36	LEU	4.8
11	2A	11	LYS	4.8
46	D5	159	PRO	4.8
36	35	150	ALA	4.8
53	K5	19	ARG	4.8
11	2A	49	GLY	4.8
36	78	149	GLU	4.8
46	D5	171	ILE	4.8
13	4A	6	GLY	4.8
14	5A	38	GLY	4.8
35	25	1	MET	4.8
11	2A	89	ALA	4.8
24	3L	6	G	4.8
36	35	106	LEU	4.8
32	59	169	VAL	4.7
51	I5	39	CYS	4.7
12	3A	69	TYR	4.7
29	29	70	ALA	4.7

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Mol	Chain	Res	Type	RSRZ
32	59	45	VAL	4.7
18	9A	26	LEU	4.7
31	49	182	LYS	4.7
46	D5	121	HIS	4.7
51	I5	54	GLY	4.7
31	49	89	GLY	4.7
44	F8	92	LEU	4.7
12	3A	129	ALA	4.7
22	1L	67	C	4.7
39	65	2	ALA	4.7
26	14	2794	C	4.7
13	4A	65	LYS	4.6
31	49	92	VAL	4.6
45	G8	106	LEU	4.6
3	22	143	GLU	4.6
31	49	175	LEU	4.6
13	4A	7	VAL	4.6
33	69	3	VAL	4.6
7	62	79	ARG	4.6
29	21	55	ASN	4.6
22	1K	72	C	4.6
33	69	1	MET	4.6
26	1H	2799	A	4.6
46	D5	111	VAL	4.6
26	14	2802	G	4.6
51	I5	45	GLY	4.5
52	N8	57	VAL	4.5
53	K5	34	LEU	4.5
24	3L	7	A	4.5
51	I5	11	PRO	4.5
51	M8	66	SER	4.5
53	O8	19	ARG	4.5
30	39	20	LEU	4.5
3	22	147	LYS	4.5
32	59	55	PRO	4.5
31	41	26	GLN	4.5
31	49	58	GLN	4.5
46	D5	116	VAL	4.5
48	J8	97	LEU	4.5
51	I5	36	CYS	4.5
46	D5	151	HIS	4.5
7	6E	85	TYR	4.4

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Mol	Chain	Res	Type	RSRZ
53	O8	52	VAL	4.4
19	AA	78	ARG	4.4
5	42	155	GLU	4.4
7	62	154	TYR	4.4
2	12	163	PHE	4.4
48	J8	93	GLU	4.4
6	5E	101	ALA	4.4
51	I5	5	ILE	4.4
22	1K	75	C	4.4
22	1L	48	C	4.4
11	2A	91	ARG	4.4
53	O8	23	THR	4.3
22	1K	13	C	4.3
18	9A	43	PHE	4.3
45	C5	29	GLU	4.3
22	1L	5	G	4.3
53	O8	45	LYS	4.3
11	2I	42	TRP	4.3
51	I5	13	ARG	4.3
22	1K	74	C	4.3
31	49	136	ARG	4.3
51	I5	27	THR	4.3
22	1K	68	C	4.3
29	29	76	ARG	4.3
22	1L	69	G	4.3
51	M8	52	THR	4.3
37	45	104	PHE	4.3
18	9A	17	SER	4.3
15	6I	89	GLY	4.3
33	61	113	ARG	4.3
7	6E	80	VAL	4.2
29	21	78	LEU	4.2
51	M8	32	TYR	4.2
36	78	106	LEU	4.2
33	69	16	GLY	4.2
53	O8	32	ASN	4.1
53	O8	31	PRO	4.1
29	29	73	GLU	4.1
48	F5	93	GLU	4.1
24	3K	17	C	4.1
22	1L	44	G	4.1
26	14	2790	A	4.1

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Mol	Chain	Res	Type	RSRZ
26	14	654(H)	G	4.1
32	59	4	ILE	4.1
49	K8	43	GLN	4.1
31	49	145	THR	4.1
46	D5	115	GLY	4.1
51	M8	22	ILE	4.1
32	59	49	VAL	4.1
18	9A	23	LYS	4.1
33	69	11	ASN	4.1
22	1L	57	G	4.0
26	1H	2795	G	4.0
31	49	75	LYS	4.0
7	62	83	ALA	4.0
33	69	85	GLU	4.0
51	I5	18	CYS	4.0
53	O8	25	LYS	4.0
53	O8	29	ASN	4.0
26	14	1	G	4.0
48	J8	96	LYS	4.0
2	1E	188	ALA	4.0
32	51	172	LYS	4.0
11	2A	21	ILE	4.0
31	49	63	ILE	4.0
3	22	198	VAL	4.0
31	49	34	LEU	4.0
46	D5	96	VAL	4.0
31	49	152	LEU	4.0
22	1K	20	U	4.0
26	1H	654(F)	C	4.0
51	I5	35	VAL	4.0
46	D5	155	LEU	4.0
22	1L	18	G	3.9
32	59	114	VAL	3.9
3	22	204	LEU	3.9
22	1K	12	U	3.9
11	2A	65	ALA	3.9
30	39	22	ALA	3.9
11	2A	90	GLY	3.9
45	C5	44	ILE	3.9
22	1L	65	G	3.9
18	9A	18	ARG	3.9
22	1K	66	U	3.9

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Mol	Chain	Res	Type	RSRZ
31	49	167	GLU	3.9
29	29	205	ALA	3.9
33	69	35	LEU	3.9
29	29	51	PHE	3.9
26	1H	2898	U	3.9
7	62	88	PRO	3.9
26	1H	654(H)	G	3.9
7	62	152	ALA	3.9
29	29	77	ILE	3.9
36	35	123	LEU	3.8
31	41	2	PRO	3.8
26	1H	2798	C	3.8
31	49	62	LEU	3.8
36	35	125	VAL	3.8
5	42	45	PHE	3.8
45	C5	92	ASN	3.8
12	3A	56	ALA	3.8
6	52	101	ALA	3.8
2	12	14	GLY	3.8
46	D5	2	GLU	3.8
7	6E	78	ARG	3.8
11	2A	59	TYR	3.8
12	3A	28	LYS	3.8
46	D5	168	GLU	3.8
29	21	204	ALA	3.8
31	49	82	LEU	3.8
46	D5	125	LEU	3.8
46	D5	150	LEU	3.8
37	45	92	GLY	3.8
41	85	91	ASP	3.8
14	5A	26	ARG	3.8
2	1E	187	LEU	3.8
51	M8	56	VAL	3.8
44	B5	92	LEU	3.7
3	22	103	VAL	3.7
6	52	38	GLU	3.7
53	K5	37	ARG	3.7
45	C5	60	PHE	3.7
32	51	3	ARG	3.7
5	4E	154	GLY	3.7
45	C5	45	VAL	3.7
3	22	206	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
32	59	25	LYS	3.7
12	3I	64	TYR	3.7
12	3A	55	VAL	3.7
4	32	179	GLU	3.7
48	F5	95	LEU	3.7
31	49	140	ILE	3.7
29	21	91	VAL	3.7
51	M8	55	ARG	3.7
7	6E	79	ARG	3.7
11	2A	92	GLU	3.7
31	49	135	LEU	3.6
46	D5	1	MET	3.6
26	14	2801	A	3.6
53	O8	41	PRO	3.6
53	O8	12	GLU	3.6
29	29	78	LEU	3.6
51	M8	40	HIS	3.6
22	1K	7	A	3.6
26	14	1536	A	3.6
2	12	165	VAL	3.6
29	21	76	ARG	3.6
2	12	155	LEU	3.6
5	42	151	LEU	3.6
46	D5	92	SER	3.6
19	AI	61	TYR	3.6
53	K5	45	LYS	3.6
14	5I	52	GLN	3.6
26	14	888	C	3.6
36	35	111	ARG	3.6
32	59	94	TYR	3.6
39	65	112	PHE	3.6
29	29	59	VAL	3.6
22	1L	58	A	3.6
22	1L	66	U	3.6
3	22	155	GLY	3.6
12	3A	60	LEU	3.6
32	59	46	GLU	3.6
42	95	1	MET	3.6
33	69	36	ALA	3.6
42	95	14	VAL	3.6
33	69	12	LEU	3.5
6	5E	46	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
42	95	91	TYR	3.5
29	29	96	PHE	3.5
53	O8	17	LYS	3.5
51	I5	50	VAL	3.5
3	22	132	ARG	3.5
19	AA	83	HIS	3.5
19	AA	84	GLY	3.5
45	C5	46	LYS	3.5
46	D5	132	ASN	3.5
11	2A	31	THR	3.5
40	75	6	LEU	3.5
50	H5	26	LEU	3.5
46	D5	173	ALA	3.5
36	35	91	PHE	3.5
31	49	166	ASP	3.5
32	51	171	LEU	3.5
31	49	142	PRO	3.5
34	15	51	PHE	3.5
29	21	199	ARG	3.5
53	K5	48	VAL	3.5
33	69	20	ASP	3.5
32	59	106	THR	3.5
46	D5	5	LEU	3.5
41	85	72	HIS	3.5
11	2A	48	ILE	3.5
7	6E	153	HIS	3.5
53	K5	46	HIS	3.5
31	41	114	ILE	3.4
32	59	103	LEU	3.4
33	69	34	GLY	3.4
51	I5	12	ALA	3.4
4	32	176	LEU	3.4
46	D5	55	HIS	3.4
46	D5	169	GLU	3.4
3	22	201	TYR	3.4
36	35	126	VAL	3.4
2	12	232	PRO	3.4
6	52	39	LYS	3.4
22	1K	1	G	3.4
13	4I	6	GLY	3.4
17	8I	17	LYS	3.4
46	D5	51	ALA	3.4

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Mol	Chain	Res	Type	RSRZ
46	D5	4	ARG	3.4
11	2I	82	VAL	3.4
13	4A	66	LEU	3.4
26	1H	5	A	3.4
36	35	145	PRO	3.4
51	I5	26	SER	3.4
31	41	135	LEU	3.4
11	2A	42	TRP	3.4
42	D8	37	VAL	3.4
26	14	654(O)	G	3.4
43	E8	111	HIS	3.4
11	2I	98	LEU	3.3
15	6I	88	ARG	3.3
3	22	64	VAL	3.3
30	31	6	VAL	3.3
12	3A	94	PRO	3.3
22	1K	47	U	3.3
33	69	17	GLN	3.3
33	69	18	VAL	3.3
51	I5	33	VAL	3.3
31	49	157	ILE	3.3
42	95	12	TYR	3.3
45	C5	50	ARG	3.3
24	3K	20	U	3.3
31	49	97	ASP	3.3
42	95	5	VAL	3.3
19	AI	48	THR	3.3
32	59	39	PRO	3.3
3	22	205	GLY	3.3
33	61	140	LEU	3.3
31	41	181	ARG	3.3
46	D5	156	LYS	3.3
53	K5	17	LYS	3.3
3	22	144	SER	3.3
5	4E	118	ILE	3.3
11	2A	43	SER	3.3
12	3A	127	GLU	3.3
45	C5	90	LEU	3.3
2	1E	160	ASP	3.3
46	D5	138	GLU	3.3
8	7E	24	THR	3.3
36	35	124	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
29	29	67	PHE	3.3
29	29	204	ALA	3.3
53	O8	9	LEU	3.3
39	A8	2	ALA	3.3
14	5A	58	LYS	3.3
18	9I	21	LYS	3.3
46	D5	109	ALA	3.3
4	3E	110	PHE	3.3
32	59	107	VAL	3.3
31	49	68	PRO	3.3
11	2I	21	ILE	3.3
42	95	44	LYS	3.3
10	1I	5	ARG	3.2
3	22	124	ILE	3.2
26	1H	2797	U	3.2
31	49	134	GLY	3.2
53	O8	33	LYS	3.2
36	35	128	HIS	3.2
42	95	16	PRO	3.2
40	75	99	LEU	3.2
10	1I	23	ILE	3.2
18	9A	86	VAL	3.2
46	D5	139	VAL	3.2
22	1L	59	U	3.2
37	45	105	GLU	3.2
45	C5	5	MET	3.2
53	O8	37	ARG	3.2
53	O8	40	CYS	3.2
37	45	33	GLY	3.2
45	C5	53	PRO	3.2
32	59	84	SER	3.2
37	45	103	MET	3.2
11	2I	43	SER	3.2
32	51	16	SER	3.2
13	4A	5	ALA	3.2
31	49	35	GLU	3.2
33	69	37	VAL	3.2
51	I5	34	GLU	3.2
31	41	182	LYS	3.2
3	2E	128	PHE	3.2
32	59	115	VAL	3.2
22	1L	23	A	3.2

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Mol	Chain	Res	Type	RSRZ
46	H8	148	ASP	3.2
13	4A	3	ARG	3.1
26	14	2897	U	3.1
3	2E	139	GLN	3.1
30	39	133	ASN	3.1
31	49	160	VAL	3.1
39	65	37	ALA	3.1
46	H8	173	ALA	3.1
53	K5	29	ASN	3.1
3	22	7	PRO	3.1
31	49	83	ARG	3.1
7	62	151	TYR	3.1
11	2I	13	GLN	3.1
31	49	108	ASN	3.1
3	22	10	PHE	3.1
29	21	69	LYS	3.1
5	4E	155	GLU	3.1
32	59	43	VAL	3.1
40	B8	106	SER	3.1
51	M8	13	ARG	3.1
42	D8	38	LEU	3.1
19	AI	40	ILE	3.1
5	4E	98	THR	3.1
18	9I	42	ARG	3.1
48	J8	70	VAL	3.1
11	2A	127	LYS	3.1
3	22	203	PHE	3.1
11	2A	108	ILE	3.1
2	12	157	ARG	3.1
31	49	147	ASP	3.1
31	49	141	PHE	3.1
51	M8	33	VAL	3.1
18	9I	31	LEU	3.1
48	J8	95	LEU	3.1
46	D5	133	ILE	3.1
10	1I	101	VAL	3.1
18	9I	19	LYS	3.1
46	D5	52	SER	3.1
3	22	177	THR	3.1
8	7E	137	VAL	3.1
14	5A	25	VAL	3.1
50	L8	57	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
31	49	176	LEU	3.1
39	65	56	LEU	3.1
42	95	36	PRO	3.1
31	49	181	ARG	3.1
13	4I	5	ALA	3.0
10	1I	10	GLY	3.0
46	D5	106	GLY	3.0
53	K5	33	LYS	3.0
31	49	94	LEU	3.0
32	59	87	LEU	3.0
49	G5	44	LEU	3.0
2	12	152	PHE	3.0
48	F5	92	LYS	3.0
31	49	74	LYS	3.0
10	1A	10	GLY	3.0
45	G8	42	VAL	3.0
39	65	83	LYS	3.0
46	D5	170	THR	3.0
4	32	23	GLY	3.0
26	1H	2897	U	3.0
12	3A	65	GLU	3.0
19	AA	47	HIS	3.0
53	O8	46	HIS	3.0
22	1L	13	C	3.0
39	65	108	GLY	3.0
22	1K	21	A	3.0
44	B5	69	TYR	3.0
19	AI	41	VAL	3.0
46	D5	56	VAL	3.0
3	2E	79	ARG	3.0
9	82	33	PHE	3.0
31	41	139	LEU	3.0
31	41	136	ARG	3.0
31	49	86	MET	3.0
37	45	102	VAL	3.0
3	22	142	MET	3.0
24	3L	17	C	3.0
52	N8	55	ARG	3.0
31	49	109	VAL	3.0
18	9A	84	LYS	3.0
30	39	172	TRP	3.0
46	D5	9	TYR	3.0

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Mol	Chain	Res	Type	RSRZ
4	3E	135	LEU	3.0
11	2I	96	ARG	3.0
12	3A	62	SER	3.0
31	49	41	GLN	3.0
36	35	130	PHE	3.0
46	H8	106	GLY	3.0
11	2A	83	ILE	3.0
3	2E	91	LEU	3.0
3	2E	82	GLU	3.0
3	22	200	ALA	2.9
2	12	70	PHE	2.9
13	4A	8	GLU	2.9
18	9A	42	ARG	2.9
24	3K	6	G	2.9
36	35	1	MET	2.9
46	D5	91	LEU	2.9
14	5A	37	PHE	2.9
29	21	79	ARG	2.9
40	75	1	MET	2.9
45	G8	64	GLU	2.9
2	12	5	ILE	2.9
2	12	102	LEU	2.9
31	49	149	VAL	2.9
36	35	138	LEU	2.9
4	32	49	ARG	2.9
31	49	102	PHE	2.9
7	62	73	MET	2.9
23	2L	1	C	2.9
29	29	181	LEU	2.9
11	2I	83	ILE	2.9
11	2A	66	LEU	2.9
26	14	1535	U	2.9
7	6E	86	GLN	2.9
51	M8	53	GLU	2.9
11	2A	61	ALA	2.9
36	35	92	GLU	2.9
3	22	145	GLY	2.9
44	F8	86	GLY	2.9
55	M5	34	TRP	2.9
42	95	93	GLU	2.9
30	39	207	GLY	2.9
26	1H	163	U	2.9

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Mol	Chain	Res	Type	RSRZ
2	1E	77	ALA	2.9
2	1E	11	LEU	2.9
4	32	69	GLY	2.9
14	5A	53	LEU	2.9
28	11	111	LEU	2.8
46	D5	32	HIS	2.8
47	E5	75	LEU	2.8
10	1A	47	PHE	2.8
3	22	170	GLN	2.8
29	21	4	ILE	2.8
26	1H	896	A	2.8
46	H8	172	ALA	2.8
46	H8	146	ILE	2.8
51	M8	65	ASP	2.8
53	O8	24	GLU	2.8
52	N8	53	ALA	2.8
31	41	138	GLN	2.8
45	C5	81	LYS	2.8
51	M8	42	PHE	2.8
31	49	174	GLU	2.8
42	D8	60	GLU	2.8
18	9I	87	ARG	2.8
7	6E	156	TRP	2.8
5	4E	6	PHE	2.8
10	1A	34	VAL	2.8
19	AA	41	VAL	2.8
37	45	65	PHE	2.8
51	I5	14	ILE	2.8
8	7E	46	LYS	2.8
19	AI	47	HIS	2.8
36	35	129	ALA	2.8
45	G8	65	ALA	2.8
11	2I	84	VAL	2.8
18	9A	22	VAL	2.8
22	1L	46	G	2.8
2	12	6	THR	2.8
46	D5	69	THR	2.8
31	49	116	ASP	2.8
30	39	191	ARG	2.8
51	I5	22	ILE	2.8
31	49	143	GLU	2.8
19	AI	71	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
10	1A	93	GLY	2.8
46	D5	108	PRO	2.8
46	D5	145	GLU	2.8
11	2I	19	ALA	2.8
4	3E	168	ARG	2.8
46	D5	177	PRO	2.8
32	59	72	ILE	2.7
39	A8	48	LEU	2.7
5	4E	73	ASN	2.7
34	58	136	GLU	2.7
32	59	83	TYR	2.7
33	69	21	VAL	2.7
3	2E	101	LEU	2.7
20	BA	104	LEU	2.7
11	2A	25	TYR	2.7
29	21	67	PHE	2.7
6	52	64	GLN	2.7
22	1K	6	G	2.7
31	41	72	ARG	2.7
33	61	121	LYS	2.7
6	52	61	LEU	2.7
33	69	5	LEU	2.7
2	12	71	VAL	2.7
7	62	147	ALA	2.7
14	5A	36	PHE	2.7
32	59	35	VAL	2.7
32	59	101	ARG	2.7
31	49	85	GLY	2.7
43	E8	112	GLY	2.7
5	42	94	ALA	2.7
45	C5	35	TYR	2.7
32	59	53	GLU	2.7
29	29	53	PRO	2.7
34	58	133	GLN	2.7
42	95	39	LEU	2.7
46	H8	107	THR	2.7
31	49	159	VAL	2.7
46	D5	8	TYR	2.7
8	72	1	MET	2.7
2	1E	31	TYR	2.7
36	35	127	ALA	2.7
12	3A	70	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
44	B5	89	ILE	2.7
18	9I	40	LEU	2.7
32	51	96	ALA	2.7
18	9I	29	PHE	2.7
35	68	52	VAL	2.7
36	35	118	GLY	2.7
22	1L	60	U	2.7
42	D8	39	LEU	2.7
48	F5	94	LEU	2.7
5	42	12	LEU	2.7
8	72	112	LEU	2.7
15	6A	31	LEU	2.7
2	1E	66	GLY	2.7
22	1L	6	G	2.7
31	49	15	VAL	2.7
39	65	60	GLY	2.7
46	D5	95	PRO	2.7
46	D5	3	TYR	2.6
12	3A	85	ILE	2.6
33	61	109	ILE	2.6
35	25	11	ALA	2.6
8	72	2	LEU	2.6
30	39	148	LEU	2.6
51	I5	9	LEU	2.6
2	1E	228	GLY	2.6
12	3A	59	ARG	2.6
42	95	4	ILE	2.6
11	2A	88	GLY	2.6
2	1E	76	GLN	2.6
4	3E	170	VAL	2.6
32	59	17	VAL	2.6
6	52	45	LEU	2.6
14	5A	34	TYR	2.6
36	35	134	ALA	2.6
44	F8	89	ILE	2.6
3	22	135	LYS	2.6
42	D8	44	LYS	2.6
46	D5	161	VAL	2.6
22	1L	55	U	2.6
31	49	179	PRO	2.6
29	21	51	PHE	2.6
12	3A	35	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
46	D5	97	GLU	2.6
2	1E	213	LEU	2.6
18	9I	79	LEU	2.6
23	2K	1	C	2.6
33	61	65	ALA	2.6
51	M8	28	LYS	2.6
3	22	187	ALA	2.6
4	32	17	VAL	2.6
12	3A	77	LEU	2.6
18	9A	85	LEU	2.6
20	BA	41	ILE	2.6
39	65	57	LYS	2.6
2	1E	96	ARG	2.6
3	22	184	TYR	2.6
46	D5	164	ALA	2.6
12	3I	65	GLU	2.6
51	I5	17	GLY	2.6
4	32	19	LEU	2.6
32	59	41	MET	2.6
12	3A	34	ARG	2.6
5	42	29	GLY	2.6
1	1G	1032	A	2.6
33	69	38	LEU	2.6
39	65	35	ILE	2.6
51	M8	5	ILE	2.6
2	12	116	GLU	2.6
4	3E	167	GLY	2.6
31	41	75	LYS	2.6
39	A8	111	GLU	2.6
12	3A	71	PRO	2.6
19	AI	62	ILE	2.6
19	AA	62	ILE	2.6
46	D5	57	ILE	2.6
3	2E	71	ALA	2.6
32	59	18	GLU	2.6
2	12	68	ILE	2.6
26	1H	1095	A	2.5
36	78	118	GLY	2.5
29	21	171	GLU	2.5
33	61	146	ALA	2.5
36	35	119	GLU	2.5
47	E5	72	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
2	1E	165	VAL	2.5
4	3E	102	ASP	2.5
35	68	122	LEU	2.5
46	H8	5	LEU	2.5
36	78	110	TYR	2.5
36	35	94	GLU	2.5
32	59	50	VAL	2.5
33	69	144	VAL	2.5
2	1E	27	LYS	2.5
5	42	6	PHE	2.5
34	15	48	MET	2.5
26	14	884	C	2.5
2	12	62	ALA	2.5
37	45	89	ASN	2.5
4	3E	90	GLY	2.5
8	7E	66	GLY	2.5
9	82	115	GLY	2.5
11	2I	14	VAL	2.5
26	1H	654(L)	G	2.5
33	61	107	VAL	2.5
10	1I	98	ILE	2.5
2	1E	14	GLY	2.5
2	12	197	VAL	2.5
3	22	60	ALA	2.5
44	F8	26	TYR	2.5
46	D5	110	GLY	2.5
5	42	43	LEU	2.5
18	9A	58	LEU	2.5
11	2I	81	ASP	2.5
32	59	168	PRO	2.5
32	59	96	ALA	2.5
33	69	2	LYS	2.5
46	D5	70	LEU	2.5
53	K5	47	THR	2.5
30	39	19	GLU	2.5
31	41	137	GLU	2.5
11	2A	86	GLY	2.5
14	5A	56	VAL	2.5
29	29	52	LEU	2.5
38	98	29	LEU	2.5
46	D5	50	GLN	2.5
12	3A	100	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
31	49	93	THR	2.5
3	2E	76	VAL	2.5
19	AA	80	TYR	2.5
30	39	11	VAL	2.5
39	65	85	VAL	2.5
51	I5	25	TYR	2.5
48	F5	60	PHE	2.5
2	12	229	VAL	2.5
4	3E	111	ALA	2.5
10	1I	94	VAL	2.5
11	2A	96	ARG	2.5
18	9I	22	VAL	2.5
29	29	182	LEU	2.5
31	49	70	VAL	2.5
11	2A	67	ASP	2.5
46	D5	93	ASP	2.5
2	1E	28	PHE	2.5
3	22	186	PHE	2.5
5	4E	40	ARG	2.5
22	1L	50	U	2.5
3	2E	201	TYR	2.5
29	29	49	LEU	2.5
32	51	87	LEU	2.5
39	65	24	LEU	2.5
48	F5	71	TYR	2.5
34	58	51	PHE	2.5
46	H8	25	PRO	2.5
3	2E	200	ALA	2.5
4	32	162	LEU	2.5
5	4E	74	GLY	2.5
37	88	86	GLY	2.5
53	K5	27	LYS	2.5
15	6I	87	ILE	2.5
26	14	883	G	2.5
53	K5	44	ARG	2.5
3	22	188	LEU	2.4
29	21	7	VAL	2.4
42	95	18	LEU	2.4
51	I5	19	GLY	2.4
30	39	27	GLU	2.4
2	12	230	VAL	2.4
5	4E	34	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
11	2A	82	VAL	2.4
26	1H	2803	C	2.4
34	58	130	HIS	2.4
42	D8	2	PHE	2.4
6	52	100	ASN	2.4
53	K5	43	CYS	2.4
2	1E	15	VAL	2.4
36	35	100	LEU	2.4
42	D8	42	GLY	2.4
46	D5	165	VAL	2.4
49	G5	60	LEU	2.4
6	52	35	ALA	2.4
11	2A	60	ALA	2.4
42	95	15	GLU	2.4
50	H5	5	LYS	2.4
55	M5	40	GLU	2.4
3	2E	66	VAL	2.4
10	1I	72	VAL	2.4
33	69	9	LEU	2.4
46	D5	131	ARG	2.4
12	3A	120	TYR	2.4
30	39	199	TRP	2.4
32	59	95	ARG	2.4
44	F8	87	GLN	2.4
8	72	133	LEU	2.4
3	22	202	ILE	2.4
10	1I	65	LEU	2.4
10	1A	101	VAL	2.4
12	3A	101	VAL	2.4
37	88	1	MET	2.4
42	D8	95	LEU	2.4
29	29	74	PRO	2.4
31	41	180	PHE	2.4
5	4E	81	GLU	2.4
19	AA	12	ASP	2.4
46	D5	59	LEU	2.4
29	29	60	ASN	2.4
3	2E	88	ARG	2.4
6	52	36	ARG	2.4
46	D5	38	TYR	2.4
32	59	89	ILE	2.4
9	82	127	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
29	29	75	VAL	2.4
47	E5	9	SER	2.4
51	M8	16	CYS	2.4
18	9I	43	PHE	2.4
37	45	83	MET	2.4
22	1L	7	A	2.4
30	31	133	ASN	2.4
38	55	102	GLU	2.4
8	7E	4	ASP	2.4
31	49	19	LEU	2.4
3	22	148	GLY	2.4
31	49	38	VAL	2.4
32	59	112	PRO	2.4
45	C5	2	ARG	2.4
37	45	80	GLU	2.4
28	19	5	LYS	2.4
29	21	49	LEU	2.4
30	39	125	LEU	2.4
30	39	181	LEU	2.4
50	L8	56	VAL	2.4
18	9I	18	ARG	2.4
31	49	80	PHE	2.4
37	45	60	ARG	2.4
2	12	4	GLU	2.4
16	7A	33	ILE	2.4
26	14	654(E)	C	2.4
18	9I	78	LEU	2.4
6	52	37	VAL	2.3
14	5I	51	GLY	2.3
42	95	101	GLY	2.3
48	F5	70	VAL	2.3
5	42	8	GLU	2.3
8	72	65	TYR	2.3
2	1E	115	LEU	2.3
4	32	108	LEU	2.3
7	6E	16	LEU	2.3
8	72	86	ILE	2.3
22	1K	10	G	2.3
30	39	10	PRO	2.3
30	39	14	PRO	2.3
33	61	126	TYR	2.3
34	58	15	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
53	O8	10	LEU	2.3
26	14	889	C	2.3
29	29	50	GLY	2.3
30	39	23	ASP	2.3
42	95	94	LEU	2.3
11	2A	84	VAL	2.3
10	1I	22	LYS	2.3
34	15	1	MET	2.3
2	1E	222	ILE	2.3
46	H8	70	LEU	2.3
50	H5	2	PRO	2.3
4	3E	3	ARG	2.3
4	3E	96	LEU	2.3
11	2I	108	ILE	2.3
11	2A	75	TYR	2.3
29	29	195	LEU	2.3
32	51	9	ILE	2.3
33	69	114	LEU	2.3
37	45	34	LEU	2.3
32	59	24	VAL	2.3
45	C5	34	LYS	2.3
13	4A	26	GLY	2.3
3	22	140	ARG	2.3
7	62	156	TRP	2.3
31	49	118	ARG	2.3
3	22	149	ALA	2.3
10	1A	65	LEU	2.3
31	49	150	ASP	2.3
11	2I	68	ALA	2.3
30	39	26	ALA	2.3
32	59	88	LEU	2.3
42	D8	3	ALA	2.3
46	H8	116	VAL	2.3
3	2E	78	GLY	2.3
4	3E	191	ARG	2.3
51	I5	6	HIS	2.3
12	3A	46	LYS	2.3
18	9A	31	LEU	2.3
33	69	6	LEU	2.3
31	49	158	ALA	2.3
32	51	17	VAL	2.3
5	42	14	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
31	49	72	ARG	2.3
32	59	170	ARG	2.3
2	1E	42	ILE	2.3
2	12	115	LEU	2.3
2	1E	197	VAL	2.3
5	4E	100	VAL	2.3
26	1H	887	A	2.3
31	49	132	ASN	2.3
31	41	80	PHE	2.3
32	51	123	PHE	2.3
33	69	14	ASP	2.3
36	35	90	ARG	2.3
48	F5	28	GLY	2.3
2	12	90	MET	2.3
3	22	199	LYS	2.3
26	14	1509	C	2.3
11	2A	30	VAL	2.3
18	9A	55	ARG	2.3
35	25	84	ALA	2.3
29	21	5	LEU	2.3
39	65	58	LEU	2.3
4	32	47	ARG	2.3
5	42	62	ALA	2.3
11	2I	30	VAL	2.3
15	6I	60	VAL	2.3
31	41	23	PHE	2.3
37	45	10	ARG	2.3
39	A8	112	PHE	2.3
41	C8	106	PHE	2.3
31	49	36	LYS	2.3
48	F5	91	LYS	2.3
6	5E	55	ASP	2.3
11	2A	63	LEU	2.3
18	9A	44	LEU	2.3
30	39	12	LEU	2.3
11	2A	54	ARG	2.3
42	D8	56	SER	2.3
32	59	52	VAL	2.3
46	D5	118	GLN	2.3
34	15	136	GLU	2.3
33	61	86	THR	2.3
2	1E	130	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
12	3A	27	LEU	2.3
5	42	133	TYR	2.3
1	1G	1033	G	2.2
8	7E	136	GLU	2.2
31	41	100	TRP	2.2
2	12	160	ASP	2.2
3	22	102	ASN	2.2
4	32	161	ASN	2.2
3	2E	72	LYS	2.2
22	1K	2	C	2.2
31	49	2	PRO	2.2
31	49	105	LYS	2.2
36	35	95	VAL	2.2
32	51	104	GLU	2.2
41	85	106	PHE	2.2
48	J8	51	VAL	2.2
5	4E	35	GLY	2.2
29	21	187	ALA	2.2
47	E5	76	GLY	2.2
11	2A	98	LEU	2.2
53	O8	11	LEU	2.2
4	3E	134	ASP	2.2
11	2A	32	ILE	2.2
31	41	88	ILE	2.2
19	AI	59	PRO	2.2
39	65	5	THR	2.2
29	21	48	GLN	2.2
51	M8	23	GLU	2.2
4	32	70	ILE	2.2
12	3I	33	ARG	2.2
19	AI	29	ARG	2.2
34	15	12	ARG	2.2
47	E5	55	ARG	2.2
3	22	180	ALA	2.2
19	AA	30	LEU	2.2
31	49	100	TRP	2.2
33	61	123	LEU	2.2
42	D8	94	LEU	2.2
7	62	149	ARG	2.2
29	29	176	ILE	2.2
32	59	51	ARG	2.2
41	85	90	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
53	K5	28	ARG	2.2
9	82	88	TYR	2.2
18	9A	57	GLY	2.2
51	I5	23	GLU	2.2
18	9I	32	ARG	2.2
18	9I	50	ILE	2.2
19	AI	60	VAL	2.2
45	G8	39	VAL	2.2
18	9A	46	GLU	2.2
46	D5	94	GLU	2.2
2	12	88	ALA	2.2
22	1K	15	G	2.2
26	14	1176	G	2.2
26	14	2793	G	2.2
36	35	81	GLN	2.2
45	C5	48	ALA	2.2
18	9A	76	LEU	2.2
4	32	110	PHE	2.2
31	49	23	PHE	2.2
5	4E	89	ILE	2.2
2	1E	129	GLU	2.2
5	42	115	VAL	2.2
46	D5	53	ILE	2.2
5	4E	97	GLY	2.2
50	H5	34	GLU	2.2
7	62	16	LEU	2.2
27	1J	88	C	2.2
29	21	32	PRO	2.2
11	2A	18	ARG	2.2
4	3E	140	VAL	2.2
17	8I	10	VAL	2.2
29	29	72	VAL	2.2
14	5A	59	ALA	2.2
20	BA	106	ALA	2.2
48	J8	80	LEU	2.2
36	78	107	LYS	2.2
31	41	178	PHE	2.2
10	1A	72	VAL	2.2
11	2A	114	VAL	2.2
22	1K	11	C	2.2
41	85	62	ILE	2.2
44	B5	3	THR	2.2

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Mol	Chain	Res	Type	RSRZ
18	9A	62	GLU	2.2
26	1H	654(P)	G	2.2
10	1A	59	SER	2.2
29	21	27	LEU	2.2
32	51	103	LEU	2.2
33	69	83	ALA	2.2
42	95	38	LEU	2.2
11	2I	110	ASP	2.2
27	16	1(M)	A	2.2
36	35	122	PRO	2.2
41	C8	117	GLN	2.2
2	12	214	ILE	2.2
10	1A	6	ILE	2.2
11	2I	31	THR	2.2
46	D5	137	ILE	2.2
26	1H	654(G)	C	2.2
34	15	116	LEU	2.2
46	H8	155	LEU	2.2
11	2I	29	ILE	2.2
37	45	68	ILE	2.2
41	85	89	GLU	2.2
18	9I	44	LEU	2.1
47	I8	7	LEU	2.1
11	2A	19	ALA	2.1
37	88	104	PHE	2.1
12	3A	72	GLY	2.1
22	1K	46	G	2.1
39	65	109	GLY	2.1
12	3A	93	LEU	2.1
52	J5	58	LEU	2.1
2	1E	93	VAL	2.1
8	7E	83	ILE	2.1
18	9I	80	PRO	2.1
26	14	654(F)	C	2.1
48	F5	69	LYS	2.1
36	78	135	LEU	2.1
12	3A	99	HIS	2.1
32	59	129	THR	2.1
51	M8	12	ALA	2.1
24	3L	44	G	2.1
19	AI	28	LYS	2.1
31	49	40	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
32	59	124	GLU	2.1
37	45	64	ILE	2.1
3	22	197	GLY	2.1
4	3E	2	GLY	2.1
18	9I	17	SER	2.1
26	14	882	G	2.1
31	49	87	PRO	2.1
46	H8	144	LEU	2.1
4	3E	185	PHE	2.1
11	2A	70	LYS	2.1
17	8I	101	ARG	2.1
24	3L	45	U	2.1
31	49	10	LYS	2.1
50	L8	3	ARG	2.1
32	59	111	HIS	2.1
10	1A	100	THR	2.1
32	59	40	GLU	2.1
46	D5	135	GLU	2.1
48	J8	90	ILE	2.1
51	I5	4	GLY	2.1
11	2A	94	ALA	2.1
2	1E	40	HIS	2.1
22	1K	50	U	2.1
26	14	885	C	2.1
26	14	893	C	2.1
31	49	88	ILE	2.1
41	85	110	VAL	2.1
53	O8	30	THR	2.1
36	78	112	LEU	2.1
51	M8	59	PHE	2.1
8	7E	25	ASP	2.1
31	49	73	ALA	2.1
32	51	18	GLU	2.1
31	41	63	ILE	2.1
3	22	101	LEU	2.1
4	3E	21	LEU	2.1
8	7E	133	LEU	2.1
11	2A	71	LYS	2.1
2	1E	202	PRO	2.1
11	2I	125	PHE	2.1
13	4A	97	PRO	2.1
3	2E	184	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
38	55	101	ALA	2.1
11	2A	109	VAL	2.1
17	8I	36	ILE	2.1
19	AA	45	VAL	2.1
34	58	53	VAL	2.1
36	78	138	LEU	2.1
38	55	70	LEU	2.1
31	41	146	TYR	2.1
44	B5	68	ARG	2.1
51	I5	8	LYS	2.1
8	7E	2	LEU	2.1
33	61	116	LEU	2.1
38	98	92	GLY	2.1
46	D5	160	GLY	2.1
52	N8	54	GLY	2.1
12	3A	67	THR	2.1
4	32	48	ALA	2.1
7	6E	155	ARG	2.1
32	51	128	PRO	2.1
17	8A	11	VAL	2.1
32	59	26	VAL	2.1
40	75	50	ILE	2.1
4	3E	196	LEU	2.1
32	59	105	LEU	2.1
36	78	130	PHE	2.1
49	G5	43	GLN	2.1
2	12	231	GLU	2.1
30	39	205	ARG	2.1
45	C5	33	LYS	2.1
3	2E	60	ALA	2.1
24	3K	16	U	2.1
30	39	97	TYR	2.1
15	6I	27	VAL	2.1
30	39	113	ALA	2.1
34	15	44	PRO	2.1
11	2I	95	ILE	2.1
19	AI	49	ILE	2.1
11	2A	17	GLY	2.1
14	5A	55	GLY	2.1
18	9A	27	GLY	2.1
22	1L	63	G	2.0
42	95	47	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
4	32	64	LEU	2.0
33	69	79	ILE	2.0
51	M8	64	GLY	2.0
3	2E	59	ARG	2.0
7	62	77	SER	2.0
31	49	115	ARG	2.0
46	D5	88	PHE	2.0
34	15	41	ASP	2.0
26	1H	277	C	2.0
2	12	112	VAL	2.0
8	72	61	VAL	2.0
4	3E	97	LEU	2.0
17	8I	50	LYS	2.0
26	1H	1534	G	2.0
9	82	93	ARG	2.0
9	82	26	VAL	2.0
30	39	128	ALA	2.0
33	61	120	ILE	2.0
5	42	28	PHE	2.0
31	41	102	PHE	2.0
22	1K	44	G	2.0
32	59	159	GLU	2.0
51	I5	53	GLU	2.0
11	2I	61	ALA	2.0
13	4I	4	ILE	2.0
31	41	94	LEU	2.0
43	E8	74	ALA	2.0
55	Q8	21	LYS	2.0
3	2E	81	GLY	2.0
32	59	123	PHE	2.0
10	1I	95	GLU	2.0
2	12	19	HIS	2.0
4	3E	169	LYS	2.0
39	A8	49	VAL	2.0
45	G8	37	VAL	2.0
11	2I	65	ALA	2.0
31	49	3	LEU	2.0
53	O8	28	ARG	2.0
22	1K	5	G	2.0
2	12	37	ASN	2.0
36	35	121	LYS	2.0
49	G5	9	GLN	2.0

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Mol	Chain	Res	Type	RSRZ
13	4A	60	VAL	2.0
15	6I	31	LEU	2.0
29	29	5	LEU	2.0
30	31	17	ARG	2.0
36	35	105	LEU	2.0
48	F5	74	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
23	7MG	2L	47	24/25	0.97	0.11	-	130,140,145,148	0
23	7MG	2K	47	24/25	0.95	0.11	-	97,107,122,130	0
23	5MU	2L	55	21/22	0.95	0.11	-	126,131,137,141	0
23	OMC	2L	33	21/22	0.95	0.14	-	105,111,114,116	0
23	4SU	2K	8	20/21	0.94	0.15	-	86,95,103,110	0
22	PSU	1K	32	20/21	0.92	0.12	-	112,119,129,131	0
23	PSU	2K	56	20/21	0.93	0.10	-	97,104,114,116	0
22	PSU	1L	39	20/21	0.86	0.20	-	113,140,153,154	0
22	MIA	1K	37	29/30	0.94	0.18	-	88,98,114,119	0
22	PSU	1K	39	20/21	0.93	0.15	-	90,106,111,114	0
23	PSU	2L	56	20/21	0.90	0.11	-	121,128,134,135	0
22	PSU	1L	32	20/21	0.91	0.16	-	139,147,158,158	0
23	4SU	2L	8	20/21	0.88	0.20	-	124,129,135,138	0
23	OMC	2K	33	21/22	0.97	0.16	-	79,84,93,100	0
22	MIA	1L	37	29/30	0.93	0.16	-	121,132,138,150	0
23	5MU	2K	55	21/22	0.95	0.12	-	99,108,119,128	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
56	MG	14	3112	1/1	0.90	0.87	62.97	90,90,90,90	0
56	MG	14	3142	1/1	0.83	0.65	61.53	73,73,73,73	0
56	MG	1H	3292	1/1	0.70	0.72	58.91	79,79,79,79	0
56	MG	1H	3117	1/1	0.94	0.56	44.29	63,63,63,63	0
56	MG	1H	3082	1/1	0.83	0.53	44.13	64,64,64,64	0
56	MG	14	3065	1/1	0.97	0.48	42.24	49,49,49,49	0
56	MG	13	1625	1/1	0.93	0.41	36.14	61,61,61,61	0
56	MG	14	3141	1/1	0.87	0.48	35.80	84,84,84,84	0
56	MG	1H	3110	1/1	0.90	0.41	35.17	64,64,64,64	0
56	MG	13	1655	1/1	0.92	0.37	34.77	80,80,80,80	0
56	MG	14	3070	1/1	0.95	0.52	31.66	59,59,59,59	0
56	MG	1G	1603	1/1	0.97	0.40	31.64	83,83,83,83	0
56	MG	14	3044	1/1	0.95	0.94	30.15	82,82,82,82	0
56	MG	14	3160	1/1	0.84	0.59	28.61	88,88,88,88	0
56	MG	1H	3122	1/1	0.84	0.51	26.27	78,78,78,78	0
56	MG	13	1651	1/1	0.90	0.45	25.93	73,73,73,73	0
56	MG	1G	1611	1/1	0.95	0.60	25.92	91,91,91,91	0
56	MG	1H	3316	1/1	0.92	0.55	25.71	64,64,64,64	0
56	MG	14	3234	1/1	0.87	0.64	25.19	80,80,80,80	0
56	MG	14	3036	1/1	0.71	0.60	25.14	68,68,68,68	0
56	MG	14	3079	1/1	0.96	0.55	24.96	69,69,69,69	0
56	MG	1H	3130	1/1	0.91	0.49	24.92	82,82,82,82	0
56	MG	14	3104	1/1	0.82	0.59	24.83	86,86,86,86	0
56	MG	1H	3001	1/1	0.97	0.43	24.36	45,45,45,45	0
56	MG	1H	3153	1/1	0.93	0.55	24.07	57,57,57,57	0
56	MG	14	3139	1/1	0.94	0.67	23.79	69,69,69,69	0
56	MG	14	3041	1/1	0.86	0.69	23.63	76,76,76,76	0
56	MG	1H	3165	1/1	0.83	0.30	23.60	68,68,68,68	0
56	MG	14	3110	1/1	0.77	0.51	22.76	88,88,88,88	0
56	MG	13	1615	1/1	0.84	0.31	22.56	103,103,103,103	0
56	MG	14	3023	1/1	0.91	0.38	21.42	66,66,66,66	0
56	MG	14	3253	1/1	0.74	0.47	21.16	86,86,86,86	0
56	MG	1G	1658	1/1	0.76	0.32	20.98	96,96,96,96	0
56	MG	1G	1659	1/1	0.84	0.32	20.85	110,110,110,110	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3126	1/1	0.95	0.46	20.85	88,88,88,88	0
56	MG	1H	3097	1/1	0.91	0.38	20.72	41,41,41,41	0
56	MG	14	3114	1/1	0.55	0.48	20.42	72,72,72,72	0
56	MG	14	3165	1/1	0.90	0.38	19.86	81,81,81,81	0
56	MG	14	3043	1/1	0.99	0.46	19.18	57,57,57,57	0
56	MG	1H	3129	1/1	0.89	0.24	19.11	77,77,77,77	0
56	MG	1H	3050	1/1	0.96	0.29	18.67	62,62,62,62	0
56	MG	14	3064	1/1	0.97	0.50	18.53	68,68,68,68	0
56	MG	13	1643	1/1	0.97	0.36	18.51	96,96,96,96	0
56	MG	14	3120	1/1	0.89	0.48	18.44	78,78,78,78	0
56	MG	14	3085	1/1	0.97	0.36	18.43	77,77,77,77	0
56	MG	14	3022	1/1	0.87	0.39	18.43	61,61,61,61	0
56	MG	14	3090	1/1	0.90	0.24	18.09	90,90,90,90	0
56	MG	1G	1618	1/1	0.89	0.28	17.26	85,85,85,85	0
56	MG	13	1610	1/1	0.75	0.40	17.06	72,72,72,72	0
56	MG	29	302	1/1	0.53	0.50	16.74	85,85,85,85	0
56	MG	1H	3281	1/1	0.78	0.31	16.67	60,60,60,60	0
56	MG	1H	3090	1/1	0.78	0.33	16.10	66,66,66,66	0
56	MG	14	3011	1/1	0.96	0.56	15.98	57,57,57,57	0
56	MG	1H	3046	1/1	0.95	0.34	15.97	80,80,80,80	0
56	MG	85	201	1/1	0.81	0.62	15.96	87,87,87,87	0
56	MG	14	3058	1/1	0.98	0.50	15.73	49,49,49,49	0
56	MG	14	3162	1/1	0.92	0.35	15.69	60,60,60,60	0
56	MG	13	1620	1/1	0.96	0.54	15.67	69,69,69,69	0
56	MG	1H	3007	1/1	0.97	0.36	15.60	40,40,40,40	0
56	MG	1G	1617	1/1	0.97	0.31	15.40	99,99,99,99	0
56	MG	1H	3068	1/1	0.84	0.29	15.02	67,67,67,67	0
56	MG	14	3053	1/1	0.92	0.35	14.72	55,55,55,55	0
56	MG	14	3035	1/1	0.99	0.47	14.56	71,71,71,71	0
56	MG	1G	1663	1/1	0.95	0.29	14.54	89,89,89,89	0
56	MG	1H	3022	1/1	0.83	0.52	14.39	78,78,78,78	0
56	MG	14	3071	1/1	0.85	0.21	14.38	88,88,88,88	0
56	MG	1H	3133	1/1	0.84	0.27	14.30	61,61,61,61	0
56	MG	1H	3023	1/1	0.89	0.27	14.14	57,57,57,57	0
56	MG	13	1607	1/1	0.91	0.36	14.12	78,78,78,78	0
56	MG	14	3040	1/1	0.97	0.43	13.94	49,49,49,49	0
56	MG	13	1603	1/1	0.98	0.36	13.82	74,74,74,74	0
56	MG	14	3108	1/1	0.88	0.43	13.73	64,64,64,64	0
56	MG	1H	3326	1/1	0.96	0.35	13.25	94,94,94,94	0
56	MG	13	1612	1/1	0.91	0.26	13.14	70,70,70,70	0
56	MG	1H	3029	1/1	0.99	0.31	12.89	65,65,65,65	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1641	1/1	0.89	0.26	12.84	90,90,90,90	0
56	MG	1H	3056	1/1	0.97	0.33	12.73	54,54,54,54	0
56	MG	13	1658	1/1	0.90	0.39	12.40	92,92,92,92	0
56	MG	1G	1602	1/1	0.98	0.45	12.38	72,72,72,72	0
56	MG	1H	3148	1/1	0.68	0.38	12.20	47,47,47,47	0
56	MG	1H	3047	1/1	0.77	0.25	12.16	71,71,71,71	0
56	MG	14	3034	1/1	0.94	0.26	12.12	64,64,64,64	0
56	MG	1H	3055	1/1	0.98	0.33	12.01	52,52,52,52	0
56	MG	1H	3241	1/1	0.92	0.42	11.89	70,70,70,70	0
56	MG	13	1605	1/1	0.96	0.34	11.88	81,81,81,81	0
56	MG	2L	101	1/1	0.99	0.37	11.77	88,88,88,88	0
56	MG	14	3167	1/1	0.79	0.42	11.76	71,71,71,71	0
56	MG	13	1677	1/1	0.72	0.34	11.69	109,109,109,109	0
56	MG	1G	1601	1/1	0.99	0.34	11.60	94,94,94,94	0
56	MG	14	3054	1/1	0.96	0.43	11.22	59,59,59,59	0
56	MG	13	1630	1/1	0.97	0.36	10.89	57,57,57,57	0
56	MG	1H	3075	1/1	0.86	0.30	10.69	53,53,53,53	0
56	MG	1H	3004	1/1	0.97	0.37	10.60	57,57,57,57	0
56	MG	14	3073	1/1	0.97	0.38	10.60	58,58,58,58	0
56	MG	1J	205	1/1	0.88	0.35	10.59	98,98,98,98	0
56	MG	14	3007	1/1	0.91	0.34	10.59	54,54,54,54	0
56	MG	1H	3020	1/1	0.94	0.36	10.54	56,56,56,56	0
56	MG	1H	3042	1/1	0.93	0.32	10.37	53,53,53,53	0
56	MG	14	3387	1/1	0.91	0.43	10.36	74,74,74,74	0
56	MG	14	3146	1/1	0.93	0.34	10.30	105,105,105,105	0
56	MG	1H	3312	1/1	0.62	0.26	10.21	82,82,82,82	0
56	MG	1H	3033	1/1	0.98	0.28	10.02	75,75,75,75	0
56	MG	1H	3028	1/1	0.80	0.24	9.76	77,77,77,77	0
56	MG	1H	3009	1/1	0.95	0.27	9.63	47,47,47,47	0
56	MG	1H	3057	1/1	0.96	0.32	9.62	40,40,40,40	0
56	MG	1G	1635	1/1	0.91	0.33	9.47	96,96,96,96	0
56	MG	1H	3217	1/1	0.95	0.32	9.47	61,61,61,61	0
56	MG	14	3004	1/1	0.98	0.31	9.43	53,53,53,53	0
56	MG	1G	1654	1/1	0.69	0.56	9.37	86,86,86,86	0
56	MG	1H	3524	1/1	0.94	0.32	9.27	49,49,49,49	0
56	MG	1H	3024	1/1	0.97	0.30	9.25	53,53,53,53	0
56	MG	13	1629	1/1	0.97	0.25	9.17	87,87,87,87	0
56	MG	1H	3300	1/1	0.60	0.26	9.13	90,90,90,90	0
56	MG	16	205	1/1	0.77	0.45	9.11	88,88,88,88	0
56	MG	13	1650	1/1	0.89	0.37	9.02	98,98,98,98	0
56	MG	14	3039	1/1	0.95	0.28	8.77	79,79,79,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3185	1/1	0.98	0.38	8.67	64,64,64,64	0
56	MG	14	3084	1/1	0.88	0.29	8.50	56,56,56,56	0
56	MG	1H	3031	1/1	0.94	0.25	8.41	63,63,63,63	0
56	MG	11	303	1/1	0.87	0.66	8.39	75,75,75,75	0
56	MG	1H	3083	1/1	0.85	0.31	7.95	55,55,55,55	0
56	MG	1H	3137	1/1	0.88	0.20	7.90	67,67,67,67	0
56	MG	1J	206	1/1	0.73	0.28	7.89	91,91,91,91	0
56	MG	1H	3096	1/1	0.89	0.32	7.85	70,70,70,70	0
56	MG	1H	3202	1/1	0.82	0.22	7.62	80,80,80,80	0
56	MG	1H	3120	1/1	0.66	0.18	7.53	68,68,68,68	0
56	MG	1H	3113	1/1	0.96	0.33	7.32	73,73,73,73	0
56	MG	13	1654	1/1	0.87	0.16	7.29	105,105,105,105	0
56	MG	14	3186	1/1	0.91	0.27	7.27	86,86,86,86	0
56	MG	1H	3053	1/1	0.96	0.27	7.26	48,48,48,48	0
56	MG	1H	3021	1/1	0.94	0.21	7.22	58,58,58,58	0
56	MG	3I	201	1/1	0.95	0.24	7.20	70,70,70,70	0
56	MG	14	3033	1/1	0.96	0.27	7.19	69,69,69,69	0
56	MG	1H	3201	1/1	0.95	0.22	7.14	79,79,79,79	0
56	MG	13	1601	1/1	0.98	0.26	7.04	83,83,83,83	0
56	MG	1H	3092	1/1	0.95	0.23	6.69	52,52,52,52	0
56	MG	13	1669	1/1	0.97	0.23	6.67	142,142,142,142	0
56	MG	1G	1639	1/1	0.74	0.39	6.64	99,99,99,99	0
56	MG	14	3046	1/1	0.93	0.30	6.59	65,65,65,65	0
56	MG	14	3100	1/1	0.86	0.31	6.53	70,70,70,70	0
56	MG	1H	3036	1/1	0.94	0.28	6.46	85,85,85,85	0
56	MG	13	1631	1/1	0.97	0.27	6.33	82,82,82,82	0
56	MG	1H	3041	1/1	0.92	0.29	6.19	57,57,57,57	0
56	MG	1H	3071	1/1	0.98	0.29	6.12	51,51,51,51	0
56	MG	98	201	1/1	0.84	0.51	5.83	99,99,99,99	0
56	MG	14	3150	1/1	0.93	0.25	5.64	64,64,64,64	0
56	MG	14	3050	1/1	0.99	0.28	5.64	81,81,81,81	0
56	MG	1H	3079	1/1	0.86	0.23	5.51	52,52,52,52	0
56	MG	14	3026	1/1	0.95	0.28	5.49	90,90,90,90	0
56	MG	1H	3255	1/1	0.75	0.33	5.47	68,68,68,68	0
56	MG	14	3227	1/1	0.92	0.28	5.47	103,103,103,103	0
56	MG	14	3155	1/1	0.93	0.19	5.30	82,82,82,82	0
56	MG	13	1614	1/1	0.94	0.29	5.20	97,97,97,97	0
56	MG	1G	1632	1/1	0.96	0.24	5.14	86,86,86,86	0
56	MG	1H	3167	1/1	0.97	0.19	5.10	62,62,62,62	0
56	MG	14	3246	1/1	0.89	0.30	4.99	98,98,98,98	0
56	MG	14	3111	1/1	0.90	0.31	4.95	86,86,86,86	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3087	1/1	0.86	0.23	4.92	72,72,72,72	0
56	MG	1H	3194	1/1	0.99	0.32	4.90	48,48,48,48	0
56	MG	14	3013	1/1	0.99	0.28	4.88	66,66,66,66	0
56	MG	1H	3318	1/1	0.71	0.18	4.86	86,86,86,86	0
56	MG	14	3200	1/1	0.86	0.17	4.77	82,82,82,82	0
56	MG	13	1671	1/1	0.69	0.28	4.69	110,110,110,110	0
56	MG	13	1681	1/1	0.73	0.57	4.61	83,83,83,83	0
56	MG	1H	3099	1/1	0.87	0.21	4.54	64,64,64,64	0
56	MG	1H	3164	1/1	0.91	0.26	4.37	70,70,70,70	0
56	MG	14	3091	1/1	0.97	0.27	4.36	54,54,54,54	0
57	ZN	3E	303	1/1	0.97	0.39	4.35	110,110,110,110	0
56	MG	1H	3234	1/1	0.86	0.23	4.14	74,74,74,74	0
56	MG	14	3248	1/1	0.98	0.21	4.11	85,85,85,85	0
56	MG	13	1679	1/1	0.82	0.18	4.07	95,95,95,95	0
56	MG	14	3178	1/1	0.86	0.18	4.06	95,95,95,95	0
56	MG	78	201	1/1	0.89	0.24	4.03	70,70,70,70	0
56	MG	13	1751	1/1	0.96	0.21	3.85	70,70,70,70	0
56	MG	14	3144	1/1	0.91	0.17	3.83	72,72,72,72	0
56	MG	2K	103	1/1	0.95	0.32	3.79	59,59,59,59	0
56	MG	11	301	1/1	0.87	0.40	3.77	48,48,48,48	0
56	MG	14	3018	1/1	0.93	0.27	3.67	68,68,68,68	0
56	MG	F5	101	1/1	0.93	0.47	3.60	74,74,74,74	0
56	MG	1H	3258	1/1	0.97	0.37	3.59	68,68,68,68	0
56	MG	1G	1633	1/1	0.93	0.33	3.50	92,92,92,92	0
56	MG	1G	1605	1/1	0.85	0.19	3.47	101,101,101,101	0
56	MG	1H	3297	1/1	0.60	0.21	3.39	80,80,80,80	0
56	MG	13	1699	1/1	0.77	0.24	3.37	98,98,98,98	0
56	MG	1H	3019	1/1	0.72	0.17	3.23	71,71,71,71	0
56	MG	1H	3072	1/1	0.87	0.23	3.19	46,46,46,46	0
56	MG	1H	3154	1/1	0.93	0.15	3.10	66,66,66,66	0
56	MG	13	1661	1/1	0.99	0.15	3.09	95,95,95,95	0
56	MG	14	3015	1/1	0.94	0.28	3.02	69,69,69,69	0
56	MG	1G	1641	1/1	0.44	0.17	2.90	121,121,121,121	0
56	MG	13	1619	1/1	0.58	0.22	2.84	82,82,82,82	0
56	MG	13	1668	1/1	0.66	0.26	2.67	111,111,111,111	0
56	MG	14	3136	1/1	0.91	0.25	2.60	60,60,60,60	0
56	MG	13	1608	1/1	0.93	0.18	2.56	75,75,75,75	0
56	MG	1H	3066	1/1	0.93	0.17	2.54	62,62,62,62	0
56	MG	1H	3088	1/1	0.98	0.24	2.52	61,61,61,61	0
56	MG	14	3148	1/1	0.75	0.46	2.45	75,75,75,75	0
56	MG	1H	3174	1/1	0.81	0.19	2.41	78,78,78,78	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1663	1/1	0.93	0.19	2.41	82,82,82,82	0
56	MG	13	1604	1/1	0.97	0.31	2.36	89,89,89,89	0
56	MG	1H	3309	1/1	0.94	0.18	2.23	87,87,87,87	0
56	MG	1H	3073	1/1	0.96	0.22	2.20	53,53,53,53	0
56	MG	31	301	1/1	0.90	0.29	2.17	62,62,62,62	0
57	ZN	G8	201	1/1	0.66	0.46	2.15	188,188,188,188	0
56	MG	1H	3197	1/1	0.89	0.27	2.09	68,68,68,68	0
56	MG	14	3152	1/1	0.88	0.18	2.07	61,61,61,61	0
56	MG	1G	1651	1/1	0.62	0.14	2.05	102,102,102,102	0
56	MG	1G	1648	1/1	0.72	0.21	1.94	115,115,115,115	0
57	ZN	32	301	1/1	0.98	0.35	1.71	123,123,123,123	0
56	MG	1H	3421	1/1	0.99	0.18	1.51	64,64,64,64	0
56	MG	1H	3018	1/1	0.98	0.23	1.49	83,83,83,83	0
56	MG	21	302	1/1	0.89	0.20	1.41	73,73,73,73	0
56	MG	14	3068	1/1	0.82	0.31	1.18	73,73,73,73	0
56	MG	88	201	1/1	0.93	0.26	1.15	79,79,79,79	0
56	MG	13	1690	1/1	0.92	0.15	1.11	96,96,96,96	0
56	MG	1H	3062	1/1	0.94	0.23	1.08	35,35,35,35	0
56	MG	14	3220	1/1	0.87	0.16	0.94	77,77,77,77	0
56	MG	C8	201	1/1	0.93	0.17	0.82	74,74,74,74	0
56	MG	1G	1612	1/1	0.60	0.18	0.82	94,94,94,94	0
56	MG	14	3159	1/1	0.95	0.20	0.81	60,60,60,60	0
56	MG	1G	1686	1/1	0.92	0.21	0.80	84,84,84,84	0
56	MG	1G	1655	1/1	0.67	0.22	0.75	124,124,124,124	0
56	MG	13	1695	1/1	0.67	0.12	0.70	88,88,88,88	0
56	MG	1H	3150	1/1	0.90	0.20	0.65	52,52,52,52	0
56	MG	16	208	1/1	0.91	0.13	0.57	70,70,70,70	0
56	MG	E8	201	1/1	0.96	0.31	0.48	79,79,79,79	0
56	MG	14	3288	1/1	0.95	0.19	0.48	62,62,62,62	0
56	MG	1G	1615	1/1	0.96	0.18	0.45	103,103,103,103	0
56	MG	11	302	1/1	0.84	0.26	0.42	82,82,82,82	0
56	MG	13	1700	1/1	0.85	0.16	0.36	100,100,100,100	0
56	MG	41	302	1/1	0.84	0.20	0.35	94,94,94,94	0
56	MG	14	3172	1/1	0.93	0.14	0.28	93,93,93,93	0
56	MG	1H	3064	1/1	0.96	0.20	0.26	45,45,45,45	0
56	MG	14	3135	1/1	0.92	0.20	0.22	54,54,54,54	0
56	MG	14	3042	1/1	0.96	0.23	0.22	48,48,48,48	0
56	MG	1J	201	1/1	0.94	0.16	0.21	114,114,114,114	0
56	MG	16	201	1/1	0.94	0.13	0.20	96,96,96,96	0
56	MG	14	3209	1/1	0.81	0.13	0.12	71,71,71,71	0
56	MG	14	3318	1/1	0.94	0.17	0.09	63,63,63,63	0
56	MG	1G	1646	1/1	0.95	0.19	0.08	84,84,84,84	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3358	1/1	0.98	0.12	0.06	65,65,65,65	0
56	MG	1H	3211	1/1	0.92	0.13	-0.07	88,88,88,88	0
56	MG	1H	3437	1/1	0.96	0.15	-0.09	82,82,82,82	0
56	MG	Q8	101	1/1	0.81	0.21	-0.13	53,53,53,53	0
56	MG	1H	3143	1/1	0.83	0.27	-0.14	68,68,68,68	0
56	MG	AI	101	1/1	0.71	0.14	-0.16	119,119,119,119	0
56	MG	1H	3314	1/1	0.90	0.17	-0.21	110,110,110,110	0
56	MG	1H	3465	1/1	0.94	0.12	-0.22	95,95,95,95	0
56	MG	1H	3346	1/1	0.96	0.18	-0.23	64,64,64,64	0
56	MG	1H	3262	1/1	0.96	0.17	-0.32	57,57,57,57	0
56	MG	14	3244	1/1	0.87	0.14	-0.36	80,80,80,80	0
56	MG	13	1660	1/1	0.90	0.15	-0.44	108,108,108,108	0
56	MG	3E	301	1/1	0.72	0.18	-0.55	128,128,128,128	0
56	MG	14	3158	1/1	0.95	0.17	-0.56	75,75,75,75	0
56	MG	1H	3015	1/1	0.99	0.16	-0.66	57,57,57,57	0
56	MG	14	3365	1/1	0.85	0.15	-0.70	87,87,87,87	0
56	MG	13	1735	1/1	0.98	0.13	-0.71	77,77,77,77	0
56	MG	1G	1638	1/1	0.94	0.11	-0.73	154,154,154,154	0
56	MG	1H	3406	1/1	0.89	0.15	-0.75	86,86,86,86	0
56	MG	14	3024	1/1	0.85	0.12	-0.87	79,79,79,79	0
56	MG	14	3364	1/1	0.95	0.14	-0.91	77,77,77,77	0
57	ZN	5I	101	1/1	0.99	0.13	-0.93	111,111,111,111	0
56	MG	13	1750	1/1	0.89	0.09	-0.97	97,97,97,97	0
56	MG	13	1702	1/1	0.36	0.14	-0.98	112,112,112,112	0
56	MG	13	1739	1/1	0.95	0.12	-0.99	118,118,118,118	0
56	MG	1H	3162	1/1	0.96	0.17	-1.01	55,55,55,55	0
57	ZN	5A	101	1/1	0.94	0.12	-1.03	160,160,160,160	0
56	MG	14	3357	1/1	0.97	0.14	-1.03	82,82,82,82	0
56	MG	1G	1660	1/1	0.98	0.07	-1.03	92,92,92,92	0
56	MG	1H	3481	1/1	0.97	0.06	-1.09	79,79,79,79	0
56	MG	13	1640	1/1	0.80	0.10	-1.10	95,95,95,95	0
56	MG	1G	1689	1/1	0.98	0.16	-1.19	80,80,80,80	0
56	MG	16	203	1/1	0.76	0.10	-1.24	91,91,91,91	0
56	MG	1H	3362	1/1	0.97	0.11	-1.25	72,72,72,72	0
56	MG	49	301	1/1	0.90	0.10	-1.26	116,116,116,116	0
56	MG	14	3367	1/1	0.81	0.17	-1.27	92,92,92,92	0
56	MG	3E	302	1/1	0.69	0.08	-1.28	125,125,125,125	0
56	MG	13	1724	1/1	0.90	0.14	-1.35	102,102,102,102	0
56	MG	1H	3388	1/1	0.95	0.15	-1.35	63,63,63,63	0
56	MG	1H	3369	1/1	0.96	0.14	-1.36	70,70,70,70	0
56	MG	14	3103	1/1	0.85	0.14	-1.43	52,52,52,52	0
56	MG	14	3290	1/1	0.97	0.16	-1.44	61,61,61,61	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3138	1/1	0.93	0.16	-1.46	57,57,57,57	0
56	MG	14	3366	1/1	0.94	0.15	-1.47	83,83,83,83	0
56	MG	14	3363	1/1	0.86	0.17	-1.55	96,96,96,96	0
56	MG	1H	3220	1/1	0.94	0.12	-1.58	61,61,61,61	0
56	MG	14	3293	1/1	0.96	0.13	-1.60	68,68,68,68	0
56	MG	14	3291	1/1	0.99	0.14	-1.64	54,54,54,54	0
56	MG	14	3388	1/1	0.99	0.12	-1.66	80,80,80,80	0
56	MG	1H	3393	1/1	0.98	0.16	-1.67	44,44,44,44	0
56	MG	1H	3392	1/1	0.98	0.19	-1.69	42,42,42,42	0
56	MG	1H	3351	1/1	0.94	0.15	-1.70	50,50,50,50	0
56	MG	14	3276	1/1	0.88	0.15	-1.72	64,64,64,64	0
56	MG	1H	3410	1/1	0.96	0.12	-1.73	62,62,62,62	0
56	MG	14	3377	1/1	0.93	0.10	-1.74	52,52,52,52	0
56	MG	1H	3395	1/1	0.99	0.14	-1.79	48,48,48,48	0
56	MG	1H	3382	1/1	0.97	0.13	-1.81	60,60,60,60	0
56	MG	1H	3529	1/1	0.85	0.12	-1.82	55,55,55,55	0
57	ZN	C5	202	1/1	0.65	0.12	-1.83	181,181,181,181	0
56	MG	41	301	1/1	0.87	0.09	-1.83	77,77,77,77	0
56	MG	14	3260	1/1	0.91	0.15	-1.84	62,62,62,62	0
56	MG	1H	3528	1/1	0.91	0.09	-1.85	77,77,77,77	0
56	MG	13	1717	1/1	0.96	0.07	-1.86	105,105,105,105	0
56	MG	1H	3526	1/1	0.96	0.15	-1.89	46,46,46,46	0
56	MG	13	1626	1/1	0.93	0.17	-1.96	65,65,65,65	0
56	MG	1H	3386	1/1	0.79	0.12	-1.96	58,58,58,58	0
56	MG	14	3296	1/1	0.99	0.17	-2.03	53,53,53,53	0
56	MG	14	3381	1/1	0.95	0.10	-2.10	79,79,79,79	0
56	MG	1H	3387	1/1	0.98	0.12	-2.10	57,57,57,57	0
56	MG	14	3315	1/1	0.95	0.13	-2.14	71,71,71,71	0
56	MG	1H	3515	1/1	0.96	0.15	-2.15	46,46,46,46	0
56	MG	1H	3342	1/1	0.99	0.14	-2.15	57,57,57,57	0
56	MG	1H	3430	1/1	0.91	0.11	-2.21	71,71,71,71	0
56	MG	16	213	1/1	0.95	0.10	-2.22	102,102,102,102	0
56	MG	1H	3370	1/1	0.97	0.15	-2.23	51,51,51,51	0
56	MG	1H	3512	1/1	0.95	0.15	-2.38	44,44,44,44	0
56	MG	1H	3447	1/1	0.96	0.12	-2.42	81,81,81,81	0
56	MG	14	3153	1/1	0.94	0.15	-2.53	60,60,60,60	0
56	MG	1H	3438	1/1	0.96	0.11	-2.56	52,52,52,52	0
56	MG	14	3328	1/1	0.73	0.15	-2.58	98,98,98,98	0
56	MG	14	3302	1/1	0.98	0.11	-2.59	73,73,73,73	0
56	MG	14	3372	1/1	0.88	0.15	-2.70	88,88,88,88	0
56	MG	1H	3402	1/1	0.95	0.16	-2.76	49,49,49,49	0
56	MG	1H	3352	1/1	0.98	0.14	-2.76	43,43,43,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1G	1692	1/1	0.93	0.07	-2.77	111,111,111,111	0
56	MG	1H	3077	1/1	0.96	0.10	-2.79	69,69,69,69	0
56	MG	14	3194	1/1	0.85	0.06	-2.80	140,140,140,140	0
56	MG	1H	3371	1/1	0.97	0.15	-2.83	53,53,53,53	0
56	MG	14	3333	1/1	0.94	0.09	-2.92	91,91,91,91	0
56	MG	1H	3114	1/1	0.92	0.12	-3.06	59,59,59,59	0
56	MG	14	3308	1/1	0.97	0.11	-3.14	68,68,68,68	0
56	MG	14	3317	1/1	0.92	0.10	-3.16	89,89,89,89	0
56	MG	14	3320	1/1	0.97	0.09	-3.21	72,72,72,72	0
56	MG	14	3261	1/1	0.92	0.14	-3.23	54,54,54,54	0
56	MG	14	3087	1/1	0.94	0.09	-3.24	59,59,59,59	0
56	MG	14	3198	1/1	0.94	0.06	-3.31	82,82,82,82	0
56	MG	1H	3525	1/1	0.97	0.10	-3.36	43,43,43,43	0
56	MG	1H	3408	1/1	0.98	0.13	-3.42	58,58,58,58	0
56	MG	14	3292	1/1	0.94	0.08	-3.49	90,90,90,90	0
56	MG	14	3386	1/1	0.97	0.09	-3.51	51,51,51,51	0
56	MG	1H	3521	1/1	0.96	0.13	-3.51	49,49,49,49	0
56	MG	1H	3147	1/1	0.98	0.15	-3.54	49,49,49,49	0
56	MG	14	3283	1/1	0.94	0.13	-3.55	57,57,57,57	0
56	MG	14	3030	1/1	0.87	0.07	-3.57	74,74,74,74	0
56	MG	1H	3527	1/1	0.87	0.10	-3.60	69,69,69,69	0
56	MG	1G	1690	1/1	0.89	0.08	-3.62	90,90,90,90	0
56	MG	14	3294	1/1	0.92	0.10	-3.77	69,69,69,69	0
56	MG	14	3378	1/1	0.85	0.12	-3.80	92,92,92,92	0
56	MG	1H	3116	1/1	0.87	0.13	-3.80	61,61,61,61	0
56	MG	13	1730	1/1	0.98	0.10	-3.82	75,75,75,75	0
56	MG	13	1728	1/1	0.94	0.09	-3.85	95,95,95,95	0
56	MG	1H	3364	1/1	0.92	0.09	-3.88	70,70,70,70	0
56	MG	14	3380	1/1	0.89	0.12	-3.98	58,58,58,58	0
56	MG	1H	3348	1/1	0.95	0.08	-4.01	60,60,60,60	0
56	MG	13	1720	1/1	0.98	0.08	-4.01	66,66,66,66	0
56	MG	14	3266	1/1	0.91	0.06	-4.11	69,69,69,69	0
56	MG	1G	1687	1/1	0.87	0.05	-4.11	121,121,121,121	0
56	MG	1H	3356	1/1	0.95	0.09	-4.12	67,67,67,67	0
56	MG	1H	3440	1/1	0.98	0.10	-4.23	63,63,63,63	0
56	MG	1G	1698	1/1	0.96	0.08	-4.31	91,91,91,91	0
56	MG	14	3307	1/1	0.95	0.09	-4.47	70,70,70,70	0
56	MG	13	1706	1/1	0.88	0.07	-4.52	108,108,108,108	0
56	MG	14	3270	1/1	0.98	0.13	-4.60	75,75,75,75	0
56	MG	14	3347	1/1	0.97	0.10	-4.65	75,75,75,75	0
56	MG	14	3338	1/1	0.95	0.06	-4.69	107,107,107,107	0
56	MG	1H	3350	1/1	0.94	0.13	-4.69	43,43,43,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3463	1/1	0.98	0.07	-4.76	82,82,82,82	0
56	MG	13	1713	1/1	0.87	0.12	-4.86	77,77,77,77	0
56	MG	1H	3368	1/1	0.96	0.06	-5.00	76,76,76,76	0
56	MG	14	3376	1/1	0.96	0.11	-5.08	65,65,65,65	0
56	MG	1H	3474	1/1	0.98	0.06	-5.12	91,91,91,91	0
56	MG	14	3362	1/1	0.94	0.06	-5.14	78,78,78,78	0
56	MG	1H	3341	1/1	0.98	0.13	-5.18	49,49,49,49	0
56	MG	13	1714	1/1	0.96	0.10	-5.20	62,62,62,62	0
56	MG	1H	3390	1/1	0.93	0.12	-5.32	55,55,55,55	0
56	MG	1H	3449	1/1	0.98	0.07	-5.65	66,66,66,66	0
56	MG	14	3325	1/1	0.93	0.05	-5.73	103,103,103,103	0
56	MG	1H	3472	1/1	0.97	0.07	-5.73	61,61,61,61	0
56	MG	1H	3337	1/1	0.89	0.09	-5.88	56,56,56,56	0
56	MG	14	3345	1/1	0.98	0.07	-5.95	72,72,72,72	0
56	MG	1H	3490	1/1	0.86	0.12	-5.96	92,92,92,92	0
56	MG	14	3272	1/1	0.95	0.12	-6.06	59,59,59,59	0
56	MG	13	1666	1/1	0.93	0.06	-6.16	99,99,99,99	0
56	MG	1H	3498	1/1	0.96	0.09	-6.21	64,64,64,64	0
56	MG	14	3375	1/1	0.92	0.10	-6.32	61,61,61,61	0
56	MG	14	3277	1/1	0.97	0.11	-6.41	67,67,67,67	0
56	MG	1H	3391	1/1	0.97	0.10	-6.74	53,53,53,53	0
56	MG	14	3295	1/1	0.92	0.09	-7.23	87,87,87,87	0
56	MG	1H	3339	1/1	0.97	0.11	-7.39	50,50,50,50	0
56	MG	1H	3436	1/1	0.97	0.07	-7.49	79,79,79,79	0
56	MG	1H	3355	1/1	0.97	0.08	-7.51	53,53,53,53	0
56	MG	1H	3495	1/1	0.93	0.07	-7.56	53,53,53,53	0
56	MG	14	3379	1/1	0.94	0.09	-7.80	60,60,60,60	0
56	MG	1H	3424	1/1	0.99	0.14	-7.91	49,49,49,49	0
56	MG	14	3262	1/1	0.97	0.10	-7.94	88,88,88,88	0
56	MG	14	3374	1/1	0.98	0.06	-8.07	57,57,57,57	0
56	MG	14	3361	1/1	0.73	0.07	-8.12	115,115,115,115	0
56	MG	1H	3448	1/1	0.98	0.04	-8.98	68,68,68,68	0
56	MG	14	3368	1/1	0.99	0.12	-9.69	53,53,53,53	0
56	MG	1H	3504	1/1	0.95	0.05	-9.72	75,75,75,75	0
56	MG	1H	3401	1/1	0.96	0.07	-9.83	50,50,50,50	0
56	MG	14	3369	1/1	0.97	0.05	-9.93	96,96,96,96	0
56	MG	1H	3505	1/1	0.95	0.07	-10.96	78,78,78,78	0
56	MG	14	3263	1/1	0.97	0.09	-12.38	60,60,60,60	0
56	MG	14	3348	1/1	0.93	0.05	-12.95	97,97,97,97	0
56	MG	1H	3383	1/1	0.99	0.05	-14.98	48,48,48,48	0
56	MG	1H	3336	1/1	0.95	0.09	-16.21	52,52,52,52	0
56	MG	1H	3485	1/1	0.96	0.07	-30.59	81,81,81,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1656	1/1	0.87	0.23	-	83,83,83,83	0
56	MG	1H	3354	1/1	0.96	0.09	-	61,61,61,61	0
56	MG	1G	1621	1/1	0.97	0.24	-	92,92,92,92	0
56	MG	1H	3166	1/1	0.96	0.43	-	60,60,60,60	0
56	MG	14	3196	1/1	0.75	0.27	-	91,91,91,91	0
56	MG	14	3344	1/1	0.98	0.09	-	85,85,85,85	0
56	MG	1H	3327	1/1	0.40	0.45	-	110,110,110,110	0
56	MG	1H	3310	1/1	0.91	0.30	-	82,82,82,82	0
56	MG	1H	3431	1/1	0.72	0.07	-	117,117,117,117	0
56	MG	14	3176	1/1	0.91	0.24	-	71,71,71,71	0
56	MG	16	210	1/1	0.89	0.28	-	97,97,97,97	0
56	MG	1H	3340	1/1	0.98	0.19	-	41,41,41,41	0
56	MG	1G	1653	1/1	0.76	0.32	-	91,91,91,91	0
56	MG	13	1683	1/1	0.78	0.31	-	78,78,78,78	0
56	MG	1H	3475	1/1	0.93	0.08	-	77,77,77,77	0
56	MG	1H	3308	1/1	0.87	0.56	-	81,81,81,81	0
56	MG	14	3188	1/1	0.81	0.24	-	83,83,83,83	0
56	MG	14	3083	1/1	0.95	0.42	-	93,93,93,93	0
56	MG	1H	3517	1/1	0.97	0.11	-	71,71,71,71	0
56	MG	14	3257	1/1	0.71	0.52	-	99,99,99,99	0
56	MG	1H	3299	1/1	0.87	0.48	-	93,93,93,93	0
56	MG	J8	102	1/1	0.74	0.31	-	66,66,66,66	0
56	MG	1H	3491	1/1	0.77	0.05	-	121,121,121,121	0
56	MG	16	204	1/1	0.87	0.43	-	86,86,86,86	0
56	MG	14	3113	1/1	0.77	0.62	-	96,96,96,96	0
56	MG	14	3115	1/1	0.89	0.38	-	86,86,86,86	0
56	MG	14	3062	1/1	0.97	0.24	-	69,69,69,69	0
56	MG	13	1748	1/1	0.92	0.06	-	156,156,156,156	0
56	MG	1H	3278	1/1	0.89	0.34	-	82,82,82,82	0
56	MG	13	1733	1/1	0.96	0.11	-	101,101,101,101	0
56	MG	1H	3032	1/1	0.97	0.36	-	46,46,46,46	0
56	MG	13	1687	1/1	0.81	0.90	-	79,79,79,79	0
56	MG	14	3239	1/1	0.74	0.26	-	88,88,88,88	0
56	MG	1H	3279	1/1	0.42	0.48	-	100,100,100,100	0
56	MG	1H	3094	1/1	0.86	0.33	-	76,76,76,76	0
56	MG	14	3217	1/1	0.74	0.30	-	88,88,88,88	0
56	MG	1H	3397	1/1	0.74	0.10	-	96,96,96,96	0
56	MG	13	1738	1/1	0.90	0.07	-	113,113,113,113	0
56	MG	13	1734	1/1	0.97	0.04	-	87,87,87,87	0
56	MG	1H	3231	1/1	0.89	0.20	-	63,63,63,63	0
56	MG	1H	3225	1/1	0.68	0.34	-	79,79,79,79	0
56	MG	1H	3179	1/1	0.77	0.27	-	59,59,59,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1G	1626	1/1	0.92	0.48	-	78,78,78,78	0
56	MG	1H	3226	1/1	0.73	0.51	-	97,97,97,97	0
56	MG	1H	3282	1/1	0.70	0.46	-	68,68,68,68	0
56	MG	14	3032	1/1	0.89	0.35	-	60,60,60,60	0
56	MG	1H	3132	1/1	0.87	0.70	-	74,74,74,74	0
56	MG	14	3319	1/1	0.96	0.05	-	88,88,88,88	0
56	MG	1H	3290	1/1	0.52	0.61	-	94,94,94,94	0
56	MG	14	3154	1/1	0.76	0.14	-	79,79,79,79	0
56	MG	13	1718	1/1	0.95	0.12	-	88,88,88,88	0
56	MG	14	3105	1/1	0.95	0.51	-	75,75,75,75	0
56	MG	1H	3115	1/1	0.92	0.24	-	59,59,59,59	0
56	MG	1H	3198	1/1	0.75	0.33	-	71,71,71,71	0
56	MG	14	3299	1/1	0.94	0.08	-	82,82,82,82	0
56	MG	1H	3466	1/1	0.83	0.07	-	106,106,106,106	0
56	MG	1H	3058	1/1	0.91	0.73	-	69,69,69,69	0
56	MG	1H	3061	1/1	0.97	0.42	-	61,61,61,61	0
56	MG	1H	3205	1/1	0.75	0.43	-	80,80,80,80	0
56	MG	1H	3123	1/1	0.85	0.67	-	73,73,73,73	0
56	MG	1H	3344	1/1	0.89	0.11	-	48,48,48,48	0
56	MG	14	3236	1/1	0.89	0.40	-	89,89,89,89	0
56	MG	14	3335	1/1	0.89	0.05	-	91,91,91,91	0
56	MG	13	1647	1/1	0.96	0.54	-	75,75,75,75	0
56	MG	13	1684	1/1	0.86	1.20	-	95,95,95,95	0
56	MG	1H	3059	1/1	0.49	0.71	-	63,63,63,63	0
56	MG	13	1634	1/1	0.90	0.28	-	80,80,80,80	0
56	MG	1H	3457	1/1	0.92	0.08	-	91,91,91,91	0
56	MG	1G	1661	1/1	0.81	1.07	-	87,87,87,87	0
56	MG	1H	3136	1/1	0.91	0.29	-	66,66,66,66	0
56	MG	14	3027	1/1	0.97	0.25	-	91,91,91,91	0
56	MG	1H	3051	1/1	0.93	0.35	-	72,72,72,72	0
56	MG	14	3181	1/1	0.73	0.96	-	86,86,86,86	0
56	MG	1G	1642	1/1	0.84	0.48	-	110,110,110,110	0
56	MG	1G	1684	1/1	0.70	0.15	-	113,113,113,113	0
56	MG	14	3219	1/1	0.52	1.15	-	89,89,89,89	0
56	MG	1H	3441	1/1	0.90	0.05	-	82,82,82,82	0
56	MG	1H	3516	1/1	0.90	0.06	-	100,100,100,100	0
56	MG	14	3097	1/1	0.89	0.56	-	80,80,80,80	0
56	MG	1H	3484	1/1	0.96	0.10	-	89,89,89,89	0
56	MG	1H	3043	1/1	0.96	0.23	-	50,50,50,50	0
56	MG	1H	3332	1/1	0.56	0.34	-	138,138,138,138	0
56	MG	14	3163	1/1	0.96	0.56	-	71,71,71,71	0
56	MG	14	3256	1/1	0.88	0.33	-	75,75,75,75	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1701	1/1	0.66	0.42	-	103,103,103,103	0
56	MG	1H	3376	1/1	0.99	0.06	-	70,70,70,70	0
56	MG	1H	3487	1/1	0.94	0.09	-	101,101,101,101	0
56	MG	14	3067	1/1	0.71	0.69	-	84,84,84,84	0
56	MG	14	3057	1/1	0.95	0.38	-	80,80,80,80	0
56	MG	1H	3215	1/1	0.87	0.65	-	71,71,71,71	0
56	MG	13	1726	1/1	0.92	0.10	-	91,91,91,91	0
56	MG	1H	3127	1/1	0.52	0.43	-	84,84,84,84	0
56	MG	14	3389	1/1	0.97	0.21	-	62,62,62,62	0
56	MG	13	1692	1/1	0.91	0.09	-	114,114,114,114	0
56	MG	1G	1688	1/1	0.97	0.12	-	73,73,73,73	0
56	MG	1H	3444	1/1	0.94	0.04	-	97,97,97,97	0
56	MG	1H	3317	1/1	0.91	0.37	-	66,66,66,66	0
56	MG	14	3038	1/1	0.86	0.26	-	74,74,74,74	0
56	MG	13	1709	1/1	0.71	0.27	-	127,127,127,127	0
56	MG	14	3166	1/1	0.92	0.19	-	112,112,112,112	0
56	MG	13	1743	1/1	0.68	0.05	-	156,156,156,156	0
56	MG	14	3355	1/1	0.97	0.10	-	97,97,97,97	0
56	MG	13	1715	1/1	0.91	0.09	-	116,116,116,116	0
56	MG	1H	3175	1/1	0.87	0.52	-	92,92,92,92	0
56	MG	1H	3329	1/1	0.86	0.24	-	99,99,99,99	0
56	MG	1H	3409	1/1	0.96	0.05	-	83,83,83,83	0
56	MG	1H	3125	1/1	0.96	0.62	-	77,77,77,77	0
56	MG	1H	3494	1/1	0.93	0.06	-	95,95,95,95	0
56	MG	1H	3293	1/1	0.79	0.52	-	82,82,82,82	0
56	MG	1H	3298	1/1	0.88	0.51	-	73,73,73,73	0
56	MG	1H	3380	1/1	0.81	0.07	-	82,82,82,82	0
56	MG	1H	3134	1/1	0.91	0.51	-	90,90,90,90	0
56	MG	1H	3121	1/1	0.80	0.61	-	88,88,88,88	0
56	MG	13	1633	1/1	0.92	0.53	-	73,73,73,73	0
56	MG	13	1613	1/1	0.89	0.21	-	83,83,83,83	0
56	MG	1H	3034	1/1	0.98	0.37	-	91,91,91,91	0
56	MG	14	3149	1/1	0.89	0.51	-	89,89,89,89	0
56	MG	1H	3285	1/1	0.64	0.49	-	89,89,89,89	0
56	MG	1H	3074	1/1	0.95	0.37	-	65,65,65,65	0
56	MG	14	3271	1/1	0.93	0.24	-	63,63,63,63	0
56	MG	13	1744	1/1	0.92	0.05	-	144,144,144,144	0
56	MG	14	3132	1/1	0.82	0.26	-	75,75,75,75	0
56	MG	1H	3138	1/1	0.82	0.46	-	69,69,69,69	0
56	MG	1G	1656	1/1	0.68	0.37	-	102,102,102,102	0
56	MG	1H	3266	1/1	0.78	0.47	-	75,75,75,75	0
56	MG	14	3216	1/1	0.85	0.55	-	70,70,70,70	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1609	1/1	0.85	0.28	-	94,94,94,94	0
56	MG	14	3251	1/1	0.89	0.16	-	102,102,102,102	0
56	MG	14	3099	1/1	0.84	0.30	-	68,68,68,68	0
56	MG	14	3323	1/1	0.96	0.10	-	96,96,96,96	0
56	MG	13	1606	1/1	0.92	0.32	-	80,80,80,80	0
56	MG	1G	1616	1/1	0.88	0.33	-	79,79,79,79	0
56	MG	14	3247	1/1	0.99	0.20	-	61,61,61,61	0
56	MG	1H	3170	1/1	0.83	0.49	-	86,86,86,86	0
56	MG	1H	3322	1/1	0.95	0.17	-	77,77,77,77	0
56	MG	1H	3460	1/1	0.98	0.07	-	63,63,63,63	0
56	MG	13	1675	1/1	0.89	0.22	-	92,92,92,92	0
56	MG	1H	3160	1/1	0.98	0.27	-	92,92,92,92	0
56	MG	14	3208	1/1	0.74	0.37	-	94,94,94,94	0
56	MG	1H	3156	1/1	0.89	0.40	-	78,78,78,78	0
56	MG	1G	1668	1/1	0.77	0.23	-	116,116,116,116	0
56	MG	14	3339	1/1	0.94	0.09	-	85,85,85,85	0
56	MG	1G	1694	1/1	0.96	0.05	-	104,104,104,104	0
56	MG	1G	1606	1/1	0.96	0.28	-	94,94,94,94	0
56	MG	1G	1670	1/1	0.65	0.43	-	98,98,98,98	0
56	MG	14	3051	1/1	0.94	0.72	-	71,71,71,71	0
56	MG	14	3197	1/1	0.43	0.42	-	94,94,94,94	0
56	MG	1H	3168	1/1	0.99	0.49	-	84,84,84,84	0
56	MG	14	3330	1/1	0.83	0.05	-	104,104,104,104	0
56	MG	1H	3473	1/1	0.68	0.11	-	92,92,92,92	0
56	MG	14	3232	1/1	0.83	0.97	-	109,109,109,109	0
56	MG	14	3322	1/1	0.93	0.06	-	101,101,101,101	0
56	MG	14	3350	1/1	0.98	0.06	-	84,84,84,84	0
56	MG	13	1716	1/1	0.93	0.05	-	89,89,89,89	0
56	MG	13	1644	1/1	0.91	0.49	-	77,77,77,77	0
56	MG	1H	3427	1/1	0.89	0.13	-	78,78,78,78	0
56	MG	29	301	1/1	0.99	0.38	-	49,49,49,49	0
56	MG	1G	1643	1/1	0.88	0.39	-	79,79,79,79	0
56	MG	14	3094	1/1	0.89	0.19	-	82,82,82,82	0
56	MG	14	3019	1/1	0.97	0.50	-	56,56,56,56	0
56	MG	1G	1682	1/1	0.72	0.66	-	102,102,102,102	0
56	MG	14	3093	1/1	0.81	0.47	-	63,63,63,63	0
56	MG	1H	3511	1/1	0.95	0.09	-	76,76,76,76	0
56	MG	1H	3212	1/1	0.88	0.68	-	92,92,92,92	0
56	MG	1K	101	1/1	0.96	0.34	-	88,88,88,88	0
56	MG	14	3224	1/1	0.87	0.38	-	90,90,90,90	0
56	MG	14	3356	1/1	0.93	0.09	-	85,85,85,85	0
56	MG	14	3195	1/1	0.79	0.98	-	96,96,96,96	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3118	1/1	0.83	0.46	-	92,92,92,92	0
56	MG	14	3342	1/1	0.98	0.10	-	72,72,72,72	0
56	MG	1H	3140	1/1	0.98	0.34	-	59,59,59,59	0
56	MG	14	3218	1/1	0.16	0.51	-	82,82,82,82	0
56	MG	14	3134	1/1	0.90	0.35	-	88,88,88,88	0
56	MG	13	1678	1/1	0.78	0.19	-	110,110,110,110	0
56	MG	14	3102	1/1	0.91	0.76	-	76,76,76,76	0
56	MG	14	3190	1/1	0.81	0.44	-	77,77,77,77	0
56	MG	13	1627	1/1	0.97	0.24	-	57,57,57,57	0
56	MG	14	3169	1/1	0.88	0.53	-	81,81,81,81	0
56	MG	1H	3413	1/1	0.96	0.08	-	71,71,71,71	0
56	MG	14	3164	1/1	0.89	0.20	-	74,74,74,74	0
56	MG	14	3264	1/1	0.97	0.10	-	53,53,53,53	0
56	MG	1G	1701	1/1	0.83	0.09	-	128,128,128,128	0
56	MG	13	1697	1/1	0.91	0.51	-	75,75,75,75	0
56	MG	1H	3095	1/1	0.75	0.32	-	78,78,78,78	0
56	MG	1H	3052	1/1	0.92	0.50	-	52,52,52,52	0
56	MG	1H	3443	1/1	0.91	0.12	-	76,76,76,76	0
56	MG	1H	3311	1/1	0.88	0.19	-	70,70,70,70	0
56	MG	1H	3014	1/1	0.96	0.42	-	57,57,57,57	0
56	MG	1H	3328	1/1	0.52	0.34	-	86,86,86,86	0
56	MG	1H	3275	1/1	0.77	0.23	-	94,94,94,94	0
56	MG	1H	3452	1/1	0.98	0.09	-	69,69,69,69	0
56	MG	14	3206	1/1	0.85	0.33	-	90,90,90,90	0
56	MG	1H	3508	1/1	0.81	0.08	-	96,96,96,96	0
56	MG	1H	3076	1/1	0.88	0.41	-	75,75,75,75	0
56	MG	1H	3477	1/1	0.96	0.11	-	82,82,82,82	0
56	MG	1H	3500	1/1	0.95	0.07	-	80,80,80,80	0
56	MG	14	3384	1/1	0.87	0.07	-	110,110,110,110	0
56	MG	1H	3086	1/1	0.93	0.47	-	67,67,67,67	0
56	MG	L8	101	1/1	0.85	0.26	-	72,72,72,72	0
56	MG	1H	3048	1/1	0.87	0.32	-	84,84,84,84	0
56	MG	1G	1680	1/1	0.74	1.13	-	107,107,107,107	0
56	MG	1H	3503	1/1	0.92	0.07	-	115,115,115,115	0
56	MG	14	3207	1/1	0.87	0.40	-	93,93,93,93	0
56	MG	14	3371	1/1	0.74	0.08	-	108,108,108,108	0
56	MG	1H	3288	1/1	0.95	0.14	-	60,60,60,60	0
56	MG	1G	1666	1/1	0.85	0.46	-	94,94,94,94	0
56	MG	1H	3171	1/1	0.95	0.36	-	101,101,101,101	0
56	MG	14	3351	1/1	0.96	0.03	-	99,99,99,99	0
56	MG	1H	3010	1/1	0.99	0.26	-	47,47,47,47	0
56	MG	J8	101	1/1	0.96	0.30	-	72,72,72,72	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1672	1/1	0.92	0.24	-	118,118,118,118	0
56	MG	1G	1673	1/1	0.83	0.38	-	138,138,138,138	0
56	MG	13	1676	1/1	0.86	0.10	-	104,104,104,104	0
56	MG	14	3222	1/1	0.95	0.35	-	87,87,87,87	0
56	MG	14	3382	1/1	0.91	0.06	-	85,85,85,85	0
56	MG	1H	3523	1/1	0.98	0.50	-	37,37,37,37	0
56	MG	1H	3320	1/1	0.90	0.31	-	86,86,86,86	0
56	MG	1H	3270	1/1	0.83	0.36	-	106,106,106,106	0
56	MG	14	3303	1/1	0.97	0.07	-	107,107,107,107	0
56	MG	1H	3119	1/1	0.95	1.25	-	86,86,86,86	0
56	MG	1H	3442	1/1	0.93	0.08	-	100,100,100,100	0
56	MG	1H	3151	1/1	0.86	0.23	-	71,71,71,71	0
56	MG	1H	3235	1/1	0.87	0.57	-	82,82,82,82	0
56	MG	14	3284	1/1	0.97	0.16	-	49,49,49,49	0
56	MG	1G	1604	1/1	0.96	0.20	-	95,95,95,95	0
56	MG	1H	3037	1/1	0.87	0.34	-	64,64,64,64	0
56	MG	13	1703	1/1	0.95	0.92	-	91,91,91,91	0
56	MG	1H	3294	1/1	0.51	0.24	-	122,122,122,122	0
56	MG	1H	3455	1/1	0.92	0.08	-	111,111,111,111	0
56	MG	13	1635	1/1	0.90	0.57	-	73,73,73,73	0
56	MG	14	3215	1/1	0.74	0.43	-	111,111,111,111	0
56	MG	13	1622	1/1	0.84	0.44	-	85,85,85,85	0
56	MG	14	3107	1/1	0.96	0.39	-	76,76,76,76	0
56	MG	14	3202	1/1	0.94	0.10	-	115,115,115,115	0
56	MG	13	1618	1/1	0.91	0.70	-	71,71,71,71	0
56	MG	1G	1676	1/1	0.92	0.15	-	157,157,157,157	0
56	MG	14	3126	1/1	0.87	0.37	-	71,71,71,71	0
56	MG	14	3297	1/1	0.92	0.10	-	107,107,107,107	0
56	MG	1H	3184	1/1	0.67	0.48	-	94,94,94,94	0
56	MG	13	1746	1/1	0.80	0.06	-	127,127,127,127	0
56	MG	14	3151	1/1	0.92	0.53	-	52,52,52,52	0
56	MG	1H	3398	1/1	0.99	0.13	-	63,63,63,63	0
56	MG	14	3316	1/1	0.90	0.10	-	82,82,82,82	0
56	MG	13	1725	1/1	0.90	0.14	-	90,90,90,90	0
56	MG	14	3242	1/1	0.93	0.53	-	78,78,78,78	0
56	MG	1H	3366	1/1	0.98	0.14	-	68,68,68,68	0
56	MG	1H	3210	1/1	0.95	0.17	-	95,95,95,95	0
56	MG	1H	3259	1/1	0.70	0.54	-	115,115,115,115	0
56	MG	14	3309	1/1	0.96	0.12	-	86,86,86,86	0
56	MG	1H	3081	1/1	0.98	0.37	-	63,63,63,63	0
56	MG	1H	3361	1/1	0.97	0.04	-	105,105,105,105	0
56	MG	1H	3263	1/1	0.85	0.40	-	82,82,82,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3014	1/1	0.96	0.17	-	81,81,81,81	0
56	MG	14	3255	1/1	0.54	0.67	-	94,94,94,94	0
56	MG	14	3287	1/1	0.92	0.08	-	99,99,99,99	0
56	MG	14	3229	1/1	0.87	0.31	-	82,82,82,82	0
56	MG	14	3326	1/1	0.92	0.05	-	110,110,110,110	0
56	MG	1H	3091	1/1	0.94	0.17	-	43,43,43,43	0
56	MG	13	1732	1/1	0.43	0.11	-	132,132,132,132	0
56	MG	14	3179	1/1	0.98	0.29	-	86,86,86,86	0
56	MG	13	1722	1/1	0.97	0.23	-	94,94,94,94	0
56	MG	1G	1681	1/1	0.83	0.32	-	106,106,106,106	0
56	MG	13	1648	1/1	0.60	0.47	-	89,89,89,89	0
56	MG	1H	3223	1/1	0.57	0.25	-	69,69,69,69	0
56	MG	14	3089	1/1	0.96	0.49	-	87,87,87,87	0
56	MG	1H	3419	1/1	0.91	0.11	-	53,53,53,53	0
56	MG	1H	3335	1/1	0.77	0.41	-	97,97,97,97	0
56	MG	1G	1703	1/1	0.91	0.07	-	124,124,124,124	0
56	MG	1H	3284	1/1	0.80	0.08	-	190,190,190,190	0
56	MG	14	3225	1/1	0.95	0.09	-	96,96,96,96	0
56	MG	14	3279	1/1	0.89	0.04	-	95,95,95,95	0
56	MG	13	1685	1/1	0.58	0.32	-	123,123,123,123	0
56	MG	1H	3185	1/1	0.93	0.79	-	68,68,68,68	0
56	MG	1H	3445	1/1	0.92	0.10	-	91,91,91,91	0
56	MG	1H	3192	1/1	0.96	0.36	-	81,81,81,81	0
56	MG	1H	3509	1/1	0.89	0.09	-	82,82,82,82	0
56	MG	14	3275	1/1	0.99	0.12	-	71,71,71,71	0
56	MG	1H	3405	1/1	0.95	0.11	-	75,75,75,75	0
56	MG	14	3267	1/1	0.96	0.11	-	67,67,67,67	0
56	MG	14	3137	1/1	0.94	0.36	-	53,53,53,53	0
56	MG	14	3008	1/1	0.96	0.48	-	58,58,58,58	0
56	MG	14	3235	1/1	0.56	0.25	-	76,76,76,76	0
56	MG	1H	3307	1/1	0.79	0.46	-	80,80,80,80	0
56	MG	14	3128	1/1	0.73	0.83	-	98,98,98,98	0
56	MG	13	1645	1/1	0.90	0.46	-	83,83,83,83	0
56	MG	1H	3085	1/1	0.94	0.43	-	74,74,74,74	0
56	MG	13	1638	1/1	0.90	0.63	-	107,107,107,107	0
56	MG	14	3180	1/1	0.90	0.59	-	93,93,93,93	0
56	MG	1H	3289	1/1	0.80	0.45	-	77,77,77,77	0
56	MG	14	3373	1/1	0.91	0.04	-	125,125,125,125	0
56	MG	13	1742	1/1	0.82	0.07	-	102,102,102,102	0
56	MG	1H	3428	1/1	0.80	0.16	-	100,100,100,100	0
56	MG	1G	1623	1/1	0.97	0.37	-	114,114,114,114	0
56	MG	1H	3343	1/1	0.97	0.14	-	48,48,48,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
56	MG	1H	3418	1/1	0.97	0.14	-	56,56,56,56	0
56	MG	14	3028	1/1	0.76	0.36	-	81,81,81,81	0
56	MG	14	3210	1/1	0.89	0.49	-	88,88,88,88	0
56	MG	1J	203	1/1	0.88	0.20	-	79,79,79,79	0
56	MG	14	3273	1/1	0.97	0.08	-	53,53,53,53	0
56	MG	1G	1644	1/1	0.87	0.50	-	74,74,74,74	0
56	MG	1H	3319	1/1	0.91	0.17	-	76,76,76,76	0
56	MG	16	202	1/1	0.91	0.40	-	66,66,66,66	0
56	MG	14	3334	1/1	0.80	0.09	-	116,116,116,116	0
56	MG	1H	3302	1/1	0.97	0.14	-	81,81,81,81	0
56	MG	14	3145	1/1	0.94	0.24	-	90,90,90,90	0
56	MG	1H	3249	1/1	0.91	0.47	-	89,89,89,89	0
56	MG	14	3193	1/1	0.74	0.44	-	74,74,74,74	0
56	MG	1H	3514	1/1	0.94	0.14	-	51,51,51,51	0
56	MG	1H	3372	1/1	0.96	0.14	-	65,65,65,65	0
56	MG	14	3233	1/1	0.91	0.49	-	112,112,112,112	0
56	MG	14	3300	1/1	0.98	0.13	-	47,47,47,47	0
56	MG	14	3121	1/1	0.93	0.78	-	79,79,79,79	0
56	MG	1H	3178	1/1	0.91	0.31	-	68,68,68,68	0
56	MG	1H	3248	1/1	0.74	0.15	-	81,81,81,81	0
56	MG	14	3092	1/1	0.89	0.39	-	68,68,68,68	0
56	MG	1G	1700	1/1	0.90	0.05	-	126,126,126,126	0
56	MG	1H	3103	1/1	0.92	0.33	-	76,76,76,76	0
56	MG	1H	3144	1/1	0.89	0.25	-	78,78,78,78	0
56	MG	1J	202	1/1	0.90	0.27	-	81,81,81,81	0
56	MG	1G	1702	1/1	0.95	0.07	-	146,146,146,146	0
56	MG	1H	3239	1/1	0.88	0.29	-	84,84,84,84	0
56	MG	13	1710	1/1	0.54	0.13	-	125,125,125,125	0
56	MG	1H	3367	1/1	0.92	0.06	-	83,83,83,83	0
56	MG	1H	3268	1/1	0.76	1.00	-	80,80,80,80	0
56	MG	14	3268	1/1	0.93	0.11	-	64,64,64,64	0
56	MG	1G	1622	1/1	0.71	0.34	-	90,90,90,90	0
56	MG	1H	3238	1/1	0.94	0.18	-	71,71,71,71	0
56	MG	1H	3108	1/1	0.82	0.37	-	58,58,58,58	0
56	MG	13	1694	1/1	0.71	0.50	-	95,95,95,95	0
56	MG	1H	3163	1/1	0.95	0.45	-	74,74,74,74	0
56	MG	14	3002	1/1	0.97	0.37	-	67,67,67,67	0
56	MG	1H	3489	1/1	0.74	0.14	-	86,86,86,86	0
56	MG	1H	3124	1/1	0.74	0.31	-	89,89,89,89	0
56	MG	6A	101	1/1	0.16	0.33	-	126,126,126,126	0
56	MG	1H	3013	1/1	0.91	0.30	-	53,53,53,53	0
56	MG	14	3096	1/1	0.63	0.66	-	78,78,78,78	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1711	1/1	0.79	0.16	-	109,109,109,109	0
56	MG	1H	3237	1/1	0.90	0.32	-	62,62,62,62	0
56	MG	1H	3416	1/1	0.98	0.06	-	69,69,69,69	0
56	MG	1H	3315	1/1	0.89	0.24	-	77,77,77,77	0
56	MG	13	1741	1/1	0.79	0.04	-	138,138,138,138	0
56	MG	1H	3323	1/1	0.86	0.37	-	87,87,87,87	0
56	MG	1H	3462	1/1	0.92	0.09	-	81,81,81,81	0
56	MG	14	3119	1/1	0.65	0.54	-	88,88,88,88	0
56	MG	1H	3389	1/1	0.98	0.09	-	51,51,51,51	0
56	MG	14	3117	1/1	0.76	0.44	-	84,84,84,84	0
56	MG	14	3265	1/1	0.98	0.11	-	61,61,61,61	0
56	MG	1J	208	1/1	0.91	0.05	-	109,109,109,109	0
56	MG	13	1673	1/1	0.93	0.50	-	83,83,83,83	0
56	MG	1G	1634	1/1	0.96	0.21	-	110,110,110,110	0
56	MG	1H	3423	1/1	0.95	0.06	-	93,93,93,93	0
56	MG	14	3080	1/1	0.96	0.35	-	54,54,54,54	0
56	MG	1H	3157	1/1	0.93	0.39	-	65,65,65,65	0
56	MG	1H	3467	1/1	0.90	0.04	-	86,86,86,86	0
56	MG	1H	3420	1/1	0.94	0.09	-	65,65,65,65	0
56	MG	14	3069	1/1	0.90	0.44	-	63,63,63,63	0
56	MG	1H	3334	1/1	0.88	0.20	-	77,77,77,77	0
56	MG	13	1611	1/1	0.94	0.26	-	70,70,70,70	0
56	MG	1G	1675	1/1	0.92	0.57	-	96,96,96,96	0
56	MG	16	207	1/1	0.80	0.36	-	87,87,87,87	0
56	MG	1H	3519	1/1	0.99	0.03	-	94,94,94,94	0
56	MG	1G	1652	1/1	0.77	0.62	-	86,86,86,86	0
56	MG	13	1686	1/1	0.89	0.22	-	106,106,106,106	0
56	MG	1H	3221	1/1	0.91	0.43	-	92,92,92,92	0
56	MG	14	3221	1/1	0.94	0.47	-	88,88,88,88	0
56	MG	1H	3017	1/1	0.97	0.27	-	43,43,43,43	0
56	MG	1H	3199	1/1	0.78	0.47	-	86,86,86,86	0
56	MG	2K	102	1/1	0.85	0.61	-	76,76,76,76	0
56	MG	1H	3216	1/1	0.68	0.58	-	68,68,68,68	0
56	MG	1H	3404	1/1	0.83	0.09	-	86,86,86,86	0
56	MG	13	1639	1/1	0.82	0.33	-	98,98,98,98	0
56	MG	1H	3461	1/1	0.85	0.09	-	105,105,105,105	0
56	MG	14	3238	1/1	0.94	0.33	-	77,77,77,77	0
56	MG	14	3332	1/1	0.97	0.08	-	88,88,88,88	0
56	MG	1H	3429	1/1	0.95	0.06	-	74,74,74,74	0
56	MG	1H	3152	1/1	0.92	0.40	-	53,53,53,53	0
56	MG	13	1657	1/1	0.82	0.06	-	87,87,87,87	0
56	MG	1H	3330	1/1	0.81	0.40	-	94,94,94,94	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3106	1/1	0.98	0.55	-	70,70,70,70	0
56	MG	1G	1667	1/1	0.88	0.29	-	97,97,97,97	0
56	MG	1H	3507	1/1	0.91	0.04	-	94,94,94,94	0
56	MG	14	3173	1/1	0.86	0.53	-	82,82,82,82	0
56	MG	14	3021	1/1	0.93	0.31	-	80,80,80,80	0
56	MG	1H	3169	1/1	0.93	0.15	-	74,74,74,74	0
56	MG	88	202	1/1	0.80	0.38	-	69,69,69,69	0
56	MG	16	206	1/1	0.77	0.29	-	76,76,76,76	0
56	MG	1H	3497	1/1	0.97	0.10	-	51,51,51,51	0
56	MG	1H	3464	1/1	0.89	0.07	-	111,111,111,111	0
56	MG	1H	3257	1/1	0.91	0.41	-	82,82,82,82	0
56	MG	14	3147	1/1	0.97	0.25	-	95,95,95,95	0
56	MG	1H	3111	1/1	0.96	0.28	-	100,100,100,100	0
56	MG	1H	3253	1/1	0.90	0.49	-	64,64,64,64	0
56	MG	14	3241	1/1	0.81	0.60	-	90,90,90,90	0
56	MG	1H	3233	1/1	0.85	1.00	-	75,75,75,75	0
56	MG	14	3140	1/1	0.96	0.28	-	65,65,65,65	0
56	MG	1H	3286	1/1	0.74	0.70	-	91,91,91,91	0
56	MG	14	3056	1/1	0.98	0.35	-	64,64,64,64	0
56	MG	13	1698	1/1	0.95	0.47	-	90,90,90,90	0
56	MG	1H	3331	1/1	0.81	0.73	-	74,74,74,74	0
56	MG	1H	3453	1/1	0.94	0.06	-	91,91,91,91	0
56	MG	1H	3357	1/1	0.98	0.10	-	52,52,52,52	0
56	MG	1H	3213	1/1	0.90	0.69	-	86,86,86,86	0
56	MG	1G	1691	1/1	0.95	0.05	-	119,119,119,119	0
56	MG	1H	3182	1/1	0.44	0.67	-	82,82,82,82	0
56	MG	13	1632	1/1	0.89	0.41	-	77,77,77,77	0
56	MG	1H	3276	1/1	0.88	0.33	-	81,81,81,81	0
56	MG	14	3048	1/1	0.97	0.40	-	65,65,65,65	0
56	MG	1H	3128	1/1	0.95	0.14	-	71,71,71,71	0
56	MG	1H	3422	1/1	0.96	0.09	-	83,83,83,83	0
56	MG	1H	3469	1/1	0.85	0.09	-	96,96,96,96	0
56	MG	1H	3027	1/1	0.91	0.56	-	64,64,64,64	0
56	MG	14	3047	1/1	0.97	0.30	-	59,59,59,59	0
56	MG	1H	3002	1/1	0.98	0.27	-	30,30,30,30	0
56	MG	14	3329	1/1	0.85	0.08	-	89,89,89,89	0
56	MG	1H	3193	1/1	0.93	0.41	-	67,67,67,67	0
56	MG	1H	3338	1/1	0.95	0.14	-	49,49,49,49	0
56	MG	1G	1610	1/1	0.94	0.09	-	98,98,98,98	0
56	MG	13	1712	1/1	0.73	0.29	-	114,114,114,114	0
56	MG	1H	3272	1/1	0.95	0.50	-	71,71,71,71	0
56	MG	14	3211	1/1	0.88	0.71	-	82,82,82,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3174	1/1	0.95	0.27	-	96,96,96,96	0
56	MG	13	1747	1/1	0.91	0.09	-	140,140,140,140	0
56	MG	1H	3476	1/1	0.97	0.07	-	65,65,65,65	0
56	MG	14	3189	1/1	0.73	0.86	-	86,86,86,86	0
56	MG	14	3010	1/1	0.99	0.33	-	58,58,58,58	0
56	MG	1H	3240	1/1	0.86	0.58	-	77,77,77,77	0
56	MG	14	3009	1/1	0.93	0.44	-	63,63,63,63	0
56	MG	1H	3003	1/1	0.98	0.32	-	63,63,63,63	0
56	MG	14	3301	1/1	0.96	0.06	-	81,81,81,81	0
56	MG	I8	102	1/1	0.95	0.07	-	67,67,67,67	0
56	MG	13	1745	1/1	0.90	0.05	-	119,119,119,119	0
56	MG	1H	3203	1/1	0.70	0.29	-	76,76,76,76	0
56	MG	1H	3407	1/1	0.98	0.08	-	84,84,84,84	0
56	MG	13	1731	1/1	0.97	0.07	-	69,69,69,69	0
56	MG	14	3131	1/1	0.89	0.58	-	93,93,93,93	0
56	MG	1H	3084	1/1	0.94	0.19	-	66,66,66,66	0
56	MG	1H	3492	1/1	0.92	0.11	-	108,108,108,108	0
56	MG	1H	3434	1/1	0.98	0.06	-	72,72,72,72	0
56	MG	1H	3209	1/1	0.96	0.42	-	60,60,60,60	0
56	MG	13	1617	1/1	0.90	0.42	-	70,70,70,70	0
56	MG	14	3231	1/1	0.83	0.41	-	77,77,77,77	0
56	MG	14	3269	1/1	0.93	0.15	-	90,90,90,90	0
56	MG	1H	3379	1/1	0.94	0.05	-	79,79,79,79	0
56	MG	16	211	1/1	0.94	0.07	-	84,84,84,84	0
56	MG	1G	1640	1/1	0.78	0.20	-	99,99,99,99	0
56	MG	1H	3107	1/1	0.89	0.19	-	84,84,84,84	0
56	MG	14	3066	1/1	0.96	0.52	-	65,65,65,65	0
56	MG	1H	3256	1/1	0.92	0.90	-	74,74,74,74	0
56	MG	1H	3306	1/1	0.96	0.31	-	81,81,81,81	0
56	MG	1H	3109	1/1	0.91	0.48	-	93,93,93,93	0
56	MG	14	3212	1/1	0.84	0.41	-	90,90,90,90	0
56	MG	14	3168	1/1	0.88	0.39	-	74,74,74,74	0
56	MG	4A	201	1/1	0.73	0.40	-	115,115,115,115	0
56	MG	16	209	1/1	0.62	0.27	-	90,90,90,90	0
56	MG	14	3012	1/1	0.97	0.47	-	49,49,49,49	0
56	MG	I8	101	1/1	0.90	0.05	-	88,88,88,88	0
56	MG	14	3031	1/1	0.88	0.20	-	73,73,73,73	0
56	MG	13	1723	1/1	0.95	0.12	-	105,105,105,105	0
56	MG	14	3324	1/1	0.85	0.10	-	107,107,107,107	0
56	MG	13	1636	1/1	0.71	0.14	-	93,93,93,93	0
56	MG	14	3171	1/1	0.76	0.90	-	91,91,91,91	0
56	MG	1H	3067	1/1	0.76	0.41	-	74,74,74,74	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3260	1/1	0.86	0.46	-	82,82,82,82	0
56	MG	1G	1647	1/1	0.93	0.31	-	102,102,102,102	0
56	MG	1H	3214	1/1	0.82	0.26	-	79,79,79,79	0
56	MG	14	3130	1/1	0.84	0.39	-	75,75,75,75	0
56	MG	14	3005	1/1	0.99	0.46	-	48,48,48,48	0
56	MG	5E	201	1/1	0.93	0.23	-	93,93,93,93	0
56	MG	2L	102	1/1	0.54	0.74	-	103,103,103,103	0
56	MG	1H	3025	1/1	0.92	0.18	-	86,86,86,86	0
56	MG	14	3226	1/1	0.84	0.36	-	83,83,83,83	0
56	MG	1H	3470	1/1	0.82	0.08	-	98,98,98,98	0
56	MG	1H	3065	1/1	0.97	0.66	-	51,51,51,51	0
56	MG	14	3327	1/1	0.94	0.10	-	120,120,120,120	0
56	MG	14	3360	1/1	0.77	0.27	-	112,112,112,112	0
56	MG	1H	3403	1/1	0.97	0.20	-	48,48,48,48	0
56	MG	C5	201	1/1	0.70	0.21	-	102,102,102,102	0
56	MG	1H	3177	1/1	0.96	0.23	-	62,62,62,62	0
56	MG	14	3183	1/1	0.96	0.45	-	83,83,83,83	0
56	MG	14	3109	1/1	0.59	0.84	-	85,85,85,85	0
56	MG	14	3204	1/1	0.66	0.26	-	126,126,126,126	0
56	MG	13	1729	1/1	0.97	0.08	-	111,111,111,111	0
56	MG	14	3314	1/1	0.97	0.08	-	64,64,64,64	0
56	MG	1H	3106	1/1	0.70	0.20	-	78,78,78,78	0
56	MG	1H	3187	1/1	0.97	0.40	-	89,89,89,89	0
56	MG	1G	1637	1/1	0.92	0.26	-	124,124,124,124	0
56	MG	1H	3384	1/1	0.96	0.11	-	62,62,62,62	0
56	MG	1G	1683	1/1	0.96	0.26	-	123,123,123,123	0
56	MG	14	3358	1/1	0.83	0.07	-	117,117,117,117	0
56	MG	1H	3269	1/1	0.78	0.40	-	99,99,99,99	0
56	MG	2I	301	1/1	0.97	0.38	-	55,55,55,55	0
56	MG	14	3184	1/1	0.92	0.35	-	53,53,53,53	0
56	MG	1J	209	1/1	0.94	0.07	-	94,94,94,94	0
56	MG	1H	3236	1/1	0.78	0.41	-	78,78,78,78	0
56	MG	1H	3191	1/1	0.95	0.22	-	59,59,59,59	0
56	MG	1H	3158	1/1	0.91	0.37	-	55,55,55,55	0
56	MG	1H	3181	1/1	0.95	0.23	-	99,99,99,99	0
56	MG	1G	1674	1/1	0.60	0.10	-	94,94,94,94	0
56	MG	14	3237	1/1	0.90	0.47	-	78,78,78,78	0
56	MG	13	1696	1/1	0.67	0.23	-	137,137,137,137	0
56	MG	14	3001	1/1	0.98	0.47	-	58,58,58,58	0
56	MG	13	1689	1/1	0.69	0.38	-	97,97,97,97	0
56	MG	14	3240	1/1	0.54	1.68	-	103,103,103,103	0
56	MG	1G	1695	1/1	0.93	0.04	-	103,103,103,103	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3146	1/1	0.95	0.20	-	42,42,42,42	0
56	MG	1H	3188	1/1	0.86	0.17	-	62,62,62,62	0
56	MG	1H	3040	1/1	0.96	0.23	-	47,47,47,47	0
56	MG	14	3254	1/1	0.84	0.57	-	81,81,81,81	0
56	MG	14	3086	1/1	0.87	0.25	-	53,53,53,53	0
56	MG	14	3230	1/1	0.92	0.17	-	103,103,103,103	0
56	MG	1G	1685	1/1	0.59	0.17	-	139,139,139,139	0
56	MG	1H	3070	1/1	0.86	0.53	-	74,74,74,74	0
56	MG	13	1688	1/1	0.78	0.15	-	102,102,102,102	0
56	MG	14	3143	1/1	0.93	0.75	-	70,70,70,70	0
56	MG	1H	3155	1/1	0.93	0.69	-	72,72,72,72	0
56	MG	13	1616	1/1	0.36	0.36	-	98,98,98,98	0
56	MG	14	3003	1/1	0.98	0.39	-	45,45,45,45	0
56	MG	1H	3426	1/1	0.95	0.08	-	53,53,53,53	0
56	MG	1H	3274	1/1	0.80	0.47	-	87,87,87,87	0
56	MG	1H	3439	1/1	0.94	0.08	-	76,76,76,76	0
56	MG	13	1708	1/1	0.53	0.14	-	116,116,116,116	0
56	MG	14	3076	1/1	0.91	0.30	-	79,79,79,79	0
56	MG	14	3129	1/1	0.94	0.44	-	76,76,76,76	0
56	MG	1H	3089	1/1	0.93	0.16	-	44,44,44,44	0
56	MG	1H	3353	1/1	0.99	0.15	-	61,61,61,61	0
56	MG	1H	3229	1/1	0.89	0.54	-	112,112,112,112	0
56	MG	1H	3267	1/1	0.65	0.26	-	84,84,84,84	0
56	MG	1H	3080	1/1	0.98	0.41	-	53,53,53,53	0
56	MG	13	1680	1/1	0.49	0.81	-	97,97,97,97	0
56	MG	1H	3347	1/1	0.97	0.16	-	59,59,59,59	0
56	MG	14	3156	1/1	0.87	0.23	-	80,80,80,80	0
56	MG	14	3258	1/1	0.58	0.28	-	90,90,90,90	0
56	MG	1H	3518	1/1	0.95	0.13	-	74,74,74,74	0
56	MG	1G	1629	1/1	0.92	0.39	-	89,89,89,89	0
56	MG	1H	3488	1/1	0.96	0.10	-	102,102,102,102	0
56	MG	1H	3207	1/1	0.89	0.35	-	79,79,79,79	0
56	MG	1H	3454	1/1	0.98	0.08	-	106,106,106,106	0
56	MG	1H	3030	1/1	0.95	0.28	-	70,70,70,70	0
56	MG	1H	3142	1/1	0.93	0.23	-	101,101,101,101	0
56	MG	1H	3200	1/1	0.90	0.17	-	98,98,98,98	0
56	MG	14	3052	1/1	0.83	0.78	-	74,74,74,74	0
56	MG	14	3063	1/1	0.94	0.27	-	62,62,62,62	0
56	MG	1H	3363	1/1	0.92	0.08	-	89,89,89,89	0
56	MG	1H	3324	1/1	0.46	0.17	-	93,93,93,93	0
56	MG	1H	3377	1/1	0.69	0.06	-	100,100,100,100	0
56	MG	14	3289	1/1	0.91	0.17	-	77,77,77,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3245	1/1	0.87	0.27	-	91,91,91,91	0
56	MG	14	3175	1/1	0.81	0.12	-	88,88,88,88	0
56	MG	1H	3291	1/1	0.66	0.36	-	71,71,71,71	0
56	MG	1H	3374	1/1	0.94	0.09	-	63,63,63,63	0
56	MG	14	3029	1/1	0.95	0.31	-	84,84,84,84	0
56	MG	1H	3486	1/1	0.95	0.09	-	61,61,61,61	0
56	MG	1G	1678	1/1	0.62	0.22	-	114,114,114,114	0
56	MG	14	3045	1/1	0.96	0.56	-	60,60,60,60	0
56	MG	1G	1669	1/1	0.40	0.65	-	103,103,103,103	0
56	MG	14	3187	1/1	0.83	0.71	-	62,62,62,62	0
56	MG	1H	3456	1/1	0.94	0.05	-	111,111,111,111	0
56	MG	1H	3172	1/1	0.92	0.51	-	70,70,70,70	0
56	MG	14	3213	1/1	0.93	0.55	-	93,93,93,93	0
56	MG	14	3346	1/1	0.97	0.11	-	80,80,80,80	0
56	MG	1H	3219	1/1	0.93	1.20	-	82,82,82,82	0
56	MG	14	3223	1/1	0.92	0.60	-	86,86,86,86	0
56	MG	1H	3035	1/1	0.93	0.26	-	86,86,86,86	0
56	MG	1H	3333	1/1	0.65	0.60	-	99,99,99,99	0
56	MG	1G	1696	1/1	0.90	0.05	-	98,98,98,98	0
56	MG	14	3203	1/1	0.81	0.92	-	111,111,111,111	0
56	MG	1H	3273	1/1	0.86	0.42	-	97,97,97,97	0
56	MG	16	212	1/1	0.95	0.06	-	79,79,79,79	0
56	MG	14	3285	1/1	0.95	0.12	-	57,57,57,57	0
56	MG	14	3282	1/1	0.83	0.08	-	92,92,92,92	0
56	MG	1H	3063	1/1	0.94	0.46	-	47,47,47,47	0
56	MG	1H	3045	1/1	0.96	0.35	-	75,75,75,75	0
56	MG	14	3016	1/1	0.95	0.30	-	68,68,68,68	0
56	MG	13	1719	1/1	0.96	0.14	-	64,64,64,64	0
56	MG	1H	3425	1/1	0.98	0.13	-	67,67,67,67	0
56	MG	1H	3078	1/1	0.94	0.17	-	60,60,60,60	0
56	MG	1H	3039	1/1	0.88	0.46	-	70,70,70,70	0
56	MG	14	3201	1/1	0.73	0.52	-	86,86,86,86	0
56	MG	14	3259	1/1	0.92	0.18	-	67,67,67,67	0
56	MG	1H	3232	1/1	0.78	0.15	-	90,90,90,90	0
56	MG	14	3082	1/1	0.71	0.36	-	97,97,97,97	0
56	MG	1H	3499	1/1	0.83	0.12	-	77,77,77,77	0
56	MG	1H	3501	1/1	0.96	0.12	-	70,70,70,70	0
56	MG	1H	3480	1/1	0.93	0.05	-	122,122,122,122	0
56	MG	14	3127	1/1	0.97	0.34	-	83,83,83,83	0
56	MG	14	3199	1/1	0.68	1.11	-	84,84,84,84	0
56	MG	13	1740	1/1	0.88	0.13	-	91,91,91,91	0
56	MG	1H	3131	1/1	0.92	0.45	-	79,79,79,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3141	1/1	0.98	0.13	-	53,53,53,53	0
56	MG	14	3352	1/1	0.96	0.08	-	106,106,106,106	0
56	MG	13	1667	1/1	0.81	0.64	-	87,87,87,87	0
56	MG	1G	1613	1/1	0.82	1.03	-	87,87,87,87	0
56	MG	13	1736	1/1	0.98	0.15	-	86,86,86,86	0
56	MG	1H	3038	1/1	0.91	0.26	-	76,76,76,76	0
56	MG	14	3359	1/1	0.87	0.09	-	97,97,97,97	0
56	MG	13	1727	1/1	0.90	0.13	-	117,117,117,117	0
56	MG	14	3077	1/1	0.96	0.43	-	78,78,78,78	0
56	MG	13	1659	1/1	0.95	0.12	-	107,107,107,107	0
56	MG	1H	3246	1/1	0.95	0.47	-	88,88,88,88	0
56	MG	14	3341	1/1	0.86	0.06	-	103,103,103,103	0
56	MG	1H	3247	1/1	0.79	0.38	-	65,65,65,65	0
56	MG	1G	1627	1/1	0.82	0.74	-	77,77,77,77	0
56	MG	14	3278	1/1	0.95	0.10	-	84,84,84,84	0
56	MG	1H	3250	1/1	0.93	0.25	-	90,90,90,90	0
56	MG	14	3310	1/1	0.97	0.05	-	78,78,78,78	0
56	MG	13	1749	1/1	0.85	0.07	-	101,101,101,101	0
56	MG	1H	3296	1/1	0.67	0.68	-	81,81,81,81	0
56	MG	1H	3261	1/1	0.93	0.22	-	58,58,58,58	0
56	MG	13	1670	1/1	0.80	0.40	-	92,92,92,92	0
56	MG	1H	3412	1/1	0.90	0.10	-	101,101,101,101	0
56	MG	1H	3471	1/1	0.96	0.08	-	94,94,94,94	0
56	MG	1H	3483	1/1	0.92	0.06	-	117,117,117,117	0
56	MG	1H	3176	1/1	0.74	0.65	-	90,90,90,90	0
56	MG	1G	1679	1/1	0.95	0.44	-	105,105,105,105	0
56	MG	14	3252	1/1	0.74	0.53	-	117,117,117,117	0
56	MG	14	3095	1/1	0.94	0.31	-	72,72,72,72	0
56	MG	11	304	1/1	0.73	0.70	-	68,68,68,68	0
56	MG	14	3370	1/1	0.97	0.10	-	88,88,88,88	0
56	MG	13	1674	1/1	0.72	0.43	-	95,95,95,95	0
56	MG	1G	1662	1/1	0.80	0.35	-	100,100,100,100	0
56	MG	1H	3008	1/1	0.97	0.31	-	56,56,56,56	0
56	MG	14	3286	1/1	0.98	0.14	-	58,58,58,58	0
56	MG	1H	3345	1/1	0.98	0.10	-	47,47,47,47	0
56	MG	14	3313	1/1	0.91	0.07	-	81,81,81,81	0
56	MG	14	3025	1/1	0.97	0.31	-	53,53,53,53	0
56	MG	13	1652	1/1	0.95	0.07	-	79,79,79,79	0
56	MG	1H	3411	1/1	0.97	0.03	-	101,101,101,101	0
56	MG	1G	1607	1/1	0.78	0.29	-	92,92,92,92	0
56	MG	14	3055	1/1	0.94	0.32	-	60,60,60,60	0
56	MG	14	3280	1/1	0.94	0.09	-	88,88,88,88	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1628	1/1	0.91	0.39	-	77,77,77,77	0
56	MG	13	1653	1/1	0.94	0.25	-	111,111,111,111	0
56	MG	13	1721	1/1	0.93	0.07	-	94,94,94,94	0
56	MG	1G	1636	1/1	0.71	0.31	-	94,94,94,94	0
56	MG	1G	1609	1/1	0.91	0.23	-	101,101,101,101	0
56	MG	1H	3101	1/1	0.91	0.27	-	80,80,80,80	0
56	MG	1H	3459	1/1	0.97	0.10	-	64,64,64,64	0
56	MG	14	3214	1/1	0.78	0.73	-	81,81,81,81	0
56	MG	14	3125	1/1	0.95	0.25	-	105,105,105,105	0
56	MG	1H	3098	1/1	0.88	0.38	-	74,74,74,74	0
56	MG	14	3306	1/1	0.97	0.17	-	49,49,49,49	0
56	MG	1G	1657	1/1	0.97	0.48	-	87,87,87,87	0
56	MG	1G	1664	1/1	0.80	0.51	-	88,88,88,88	0
56	MG	1G	1677	1/1	0.67	0.45	-	106,106,106,106	0
56	MG	13	1693	1/1	0.90	0.40	-	109,109,109,109	0
56	MG	1H	3190	1/1	0.90	0.43	-	88,88,88,88	0
56	MG	14	3353	1/1	0.94	0.05	-	99,99,99,99	0
56	MG	1H	3049	1/1	0.97	0.21	-	74,74,74,74	0
56	MG	1H	3230	1/1	0.60	0.45	-	66,66,66,66	0
56	MG	1G	1650	1/1	0.92	0.28	-	111,111,111,111	0
56	MG	14	3249	1/1	0.60	0.32	-	69,69,69,69	0
56	MG	78	202	1/1	0.69	0.32	-	78,78,78,78	0
56	MG	14	3304	1/1	0.97	0.08	-	91,91,91,91	0
56	MG	1H	3180	1/1	0.93	0.36	-	67,67,67,67	0
56	MG	1G	1619	1/1	0.64	0.21	-	130,130,130,130	0
56	MG	14	3177	1/1	0.86	0.40	-	80,80,80,80	0
56	MG	1G	1630	1/1	0.83	0.63	-	76,76,76,76	0
56	MG	14	3383	1/1	0.93	0.13	-	93,93,93,93	0
56	MG	1G	1671	1/1	0.86	0.37	-	96,96,96,96	0
56	MG	1H	3242	1/1	0.84	0.41	-	86,86,86,86	0
56	MG	1H	3159	1/1	0.88	0.20	-	65,65,65,65	0
56	MG	1H	3145	1/1	0.94	0.28	-	49,49,49,49	0
56	MG	14	3336	1/1	0.74	0.09	-	123,123,123,123	0
56	MG	13	1624	1/1	0.98	0.70	-	71,71,71,71	0
56	MG	1H	3254	1/1	0.93	0.23	-	84,84,84,84	0
56	MG	14	3245	1/1	0.72	0.40	-	84,84,84,84	0
56	MG	13	1623	1/1	0.78	0.42	-	85,85,85,85	0
56	MG	14	3321	1/1	0.86	0.13	-	97,97,97,97	0
56	MG	1H	3227	1/1	0.70	0.55	-	93,93,93,93	0
56	MG	13	1649	1/1	0.82	0.38	-	79,79,79,79	0
56	MG	1H	3415	1/1	0.97	0.07	-	100,100,100,100	0
56	MG	1H	3478	1/1	0.91	0.07	-	111,111,111,111	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3354	1/1	0.96	0.16	-	77,77,77,77	0
56	MG	14	3059	1/1	0.97	0.27	-	59,59,59,59	0
56	MG	75	201	1/1	0.37	0.34	-	108,108,108,108	0
56	MG	1G	1645	1/1	0.91	0.32	-	92,92,92,92	0
56	MG	1H	3251	1/1	0.96	0.20	-	96,96,96,96	0
56	MG	1H	3394	1/1	0.97	0.16	-	54,54,54,54	0
56	MG	1H	3399	1/1	0.96	0.09	-	67,67,67,67	0
56	MG	1H	3450	1/1	0.86	0.10	-	119,119,119,119	0
56	MG	1H	3183	1/1	0.73	0.21	-	76,76,76,76	0
56	MG	13	1637	1/1	0.96	0.62	-	59,59,59,59	0
56	MG	1H	3196	1/1	0.93	0.17	-	90,90,90,90	0
56	MG	13	1704	1/1	0.30	1.41	-	95,95,95,95	0
56	MG	1H	3482	1/1	0.98	0.16	-	66,66,66,66	0
56	MG	14	3081	1/1	0.93	0.46	-	71,71,71,71	0
56	MG	13	1737	1/1	0.77	0.05	-	122,122,122,122	0
56	MG	14	3182	1/1	0.81	0.32	-	152,152,152,152	0
56	MG	1H	3102	1/1	0.90	0.42	-	75,75,75,75	0
56	MG	1H	3139	1/1	0.95	0.20	-	65,65,65,65	0
56	MG	1H	3060	1/1	0.97	0.49	-	48,48,48,48	0
56	MG	1G	1697	1/1	0.84	0.10	-	108,108,108,108	0
56	MG	1H	3287	1/1	0.77	0.17	-	85,85,85,85	0
56	MG	1H	3321	1/1	0.90	0.17	-	72,72,72,72	0
56	MG	14	3017	1/1	0.96	0.28	-	85,85,85,85	0
56	MG	1H	3381	1/1	0.97	0.07	-	46,46,46,46	0
56	MG	1H	3400	1/1	0.95	0.12	-	68,68,68,68	0
56	MG	1H	3301	1/1	0.79	0.54	-	84,84,84,84	0
56	MG	1H	3044	1/1	0.95	0.39	-	75,75,75,75	0
56	MG	14	3124	1/1	0.86	0.36	-	81,81,81,81	0
56	MG	1H	3271	1/1	0.90	0.46	-	90,90,90,90	0
56	MG	14	3088	1/1	0.99	0.32	-	44,44,44,44	0
56	MG	1K	102	1/1	0.87	0.51	-	93,93,93,93	0
56	MG	13	1602	1/1	0.97	0.38	-	68,68,68,68	0
56	MG	1H	3252	1/1	0.79	0.38	-	83,83,83,83	0
56	MG	13	1642	1/1	0.93	0.24	-	89,89,89,89	0
56	MG	14	3122	1/1	0.84	0.30	-	81,81,81,81	0
56	MG	14	3205	1/1	0.94	0.34	-	91,91,91,91	0
56	MG	14	3037	1/1	0.61	0.38	-	103,103,103,103	0
56	MG	13	1665	1/1	0.77	0.40	-	86,86,86,86	0
56	MG	14	3020	1/1	0.92	0.20	-	72,72,72,72	0
56	MG	1H	3496	1/1	0.88	0.12	-	59,59,59,59	0
56	MG	14	3157	1/1	0.92	0.24	-	107,107,107,107	0
56	MG	1H	3493	1/1	0.67	0.09	-	113,113,113,113	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	13	1707	1/1	0.81	0.11	-	104,104,104,104	0
56	MG	1G	1699	1/1	0.84	0.11	-	122,122,122,122	0
56	MG	1G	1693	1/1	0.87	0.09	-	125,125,125,125	0
56	MG	1H	3277	1/1	0.95	0.47	-	84,84,84,84	0
56	MG	1H	3283	1/1	0.91	0.26	-	67,67,67,67	0
56	MG	1H	3435	1/1	0.93	0.08	-	95,95,95,95	0
56	MG	1H	3360	1/1	0.91	0.04	-	105,105,105,105	0
56	MG	1H	3359	1/1	0.94	0.06	-	102,102,102,102	0
56	MG	1H	3161	1/1	0.92	0.33	-	66,66,66,66	0
56	MG	14	3243	1/1	0.81	0.87	-	81,81,81,81	0
56	MG	13	1705	1/1	0.56	0.24	-	138,138,138,138	0
56	MG	1H	3446	1/1	0.96	0.05	-	84,84,84,84	0
56	MG	1H	3373	1/1	0.98	0.07	-	63,63,63,63	0
56	MG	1H	3100	1/1	0.95	0.39	-	71,71,71,71	0
56	MG	14	3331	1/1	0.88	0.10	-	65,65,65,65	0
56	MG	1J	204	1/1	0.84	0.70	-	100,100,100,100	0
56	MG	1H	3375	1/1	0.94	0.07	-	75,75,75,75	0
56	MG	1H	3243	1/1	0.78	0.62	-	88,88,88,88	0
56	MG	1H	3173	1/1	0.96	0.15	-	79,79,79,79	0
56	MG	1H	3208	1/1	0.96	0.22	-	71,71,71,71	0
56	MG	1H	3522	1/1	0.89	0.12	-	99,99,99,99	0
56	MG	1H	3149	1/1	0.84	0.37	-	77,77,77,77	0
56	MG	14	3074	1/1	0.94	0.49	-	63,63,63,63	0
56	MG	14	3192	1/1	0.68	0.58	-	85,85,85,85	0
56	MG	14	3228	1/1	0.74	0.06	-	192,192,192,192	0
56	MG	14	3337	1/1	0.94	0.06	-	83,83,83,83	0
56	MG	14	3123	1/1	0.82	0.35	-	82,82,82,82	0
56	MG	35	201	1/1	0.92	0.39	-	76,76,76,76	0
56	MG	14	3006	1/1	0.98	0.56	-	60,60,60,60	0
56	MG	1H	3206	1/1	0.93	0.16	-	77,77,77,77	0
56	MG	1G	1665	1/1	0.82	0.59	-	104,104,104,104	0
56	MG	1H	3349	1/1	0.95	0.10	-	70,70,70,70	0
56	MG	2K	101	1/1	0.81	0.34	-	86,86,86,86	0
56	MG	1H	3414	1/1	0.91	0.09	-	80,80,80,80	0
56	MG	1H	3396	1/1	0.98	0.04	-	79,79,79,79	0
56	MG	1H	3026	1/1	0.84	0.51	-	85,85,85,85	0
56	MG	1H	3118	1/1	0.83	0.34	-	93,93,93,93	0
56	MG	1H	3264	1/1	0.73	0.30	-	76,76,76,76	0
56	MG	14	3349	1/1	0.92	0.07	-	110,110,110,110	0
56	MG	14	3311	1/1	0.95	0.07	-	82,82,82,82	0
56	MG	13	1621	1/1	0.96	0.33	-	97,97,97,97	0
56	MG	1H	3105	1/1	0.83	0.75	-	68,68,68,68	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	14	3170	1/1	0.91	0.46	-	88,88,88,88	0
56	MG	1H	3513	1/1	0.95	0.14	-	58,58,58,58	0
56	MG	1G	1649	1/1	0.97	0.16	-	145,145,145,145	0
56	MG	14	3312	1/1	0.83	0.08	-	97,97,97,97	0
56	MG	14	3191	1/1	0.88	0.23	-	94,94,94,94	0
56	MG	1G	1672	1/1	0.96	0.35	-	100,100,100,100	0
56	MG	1H	3313	1/1	0.85	0.78	-	80,80,80,80	0
56	MG	14	3385	1/1	0.83	0.23	-	113,113,113,113	0
56	MG	1H	3506	1/1	0.97	0.06	-	88,88,88,88	0
56	MG	14	3101	1/1	0.98	0.16	-	55,55,55,55	0
56	MG	13	1682	1/1	0.81	0.50	-	92,92,92,92	0
56	MG	1H	3195	1/1	0.92	0.18	-	68,68,68,68	0
56	MG	1H	3520	1/1	0.90	0.07	-	109,109,109,109	0
56	MG	1G	1628	1/1	0.75	0.47	-	76,76,76,76	0
56	MG	14	3061	1/1	0.94	0.45	-	71,71,71,71	0
56	MG	14	3060	1/1	0.98	0.36	-	65,65,65,65	0
56	MG	1H	3204	1/1	0.85	0.53	-	76,76,76,76	0
56	MG	1H	3104	1/1	0.98	0.19	-	67,67,67,67	0
56	MG	1H	3304	1/1	0.68	0.28	-	79,79,79,79	0
56	MG	14	3340	1/1	0.85	0.10	-	83,83,83,83	0
56	MG	14	3049	1/1	0.89	0.35	-	71,71,71,71	0
56	MG	1H	3303	1/1	0.96	0.21	-	88,88,88,88	0
56	MG	1H	3433	1/1	0.91	0.17	-	77,77,77,77	0
56	MG	1H	3510	1/1	0.88	0.12	-	99,99,99,99	0
56	MG	1H	3468	1/1	0.92	0.04	-	109,109,109,109	0
56	MG	14	3281	1/1	0.97	0.11	-	79,79,79,79	0
56	MG	14	3298	1/1	0.98	0.10	-	59,59,59,59	0
56	MG	14	3116	1/1	0.77	0.18	-	77,77,77,77	0
56	MG	13	1691	1/1	0.94	0.26	-	101,101,101,101	0
56	MG	1H	3265	1/1	0.85	0.55	-	92,92,92,92	0
56	MG	1H	3479	1/1	0.95	0.11	-	67,67,67,67	0
56	MG	1H	3189	1/1	0.88	0.29	-	77,77,77,77	0
56	MG	1G	1614	1/1	0.97	0.64	-	78,78,78,78	0
56	MG	1H	3451	1/1	0.92	0.08	-	94,94,94,94	0
56	MG	1H	3224	1/1	0.91	0.55	-	89,89,89,89	0
56	MG	1H	3325	1/1	0.86	0.36	-	64,64,64,64	0
56	MG	1G	1631	1/1	0.94	0.34	-	89,89,89,89	0
56	MG	1H	3005	1/1	0.98	0.43	-	46,46,46,46	0
56	MG	14	3305	1/1	0.90	0.14	-	112,112,112,112	0
56	MG	14	3075	1/1	0.93	0.43	-	61,61,61,61	0
56	MG	13	1662	1/1	0.90	0.40	-	113,113,113,113	0
56	MG	1H	3006	1/1	0.90	0.43	-	55,55,55,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	MG	1H	3228	1/1	0.95	0.33	-	71,71,71,71	0
56	MG	1H	3218	1/1	0.83	0.40	-	76,76,76,76	0
56	MG	14	3078	1/1	0.98	0.23	-	68,68,68,68	0
56	MG	13	1664	1/1	0.90	0.38	-	71,71,71,71	0
56	MG	1H	3186	1/1	0.75	0.50	-	89,89,89,89	0
56	MG	14	3072	1/1	0.99	0.21	-	75,75,75,75	0
56	MG	1H	3385	1/1	0.98	0.12	-	66,66,66,66	0
56	MG	1H	3012	1/1	0.95	0.50	-	52,52,52,52	0
56	MG	1G	1608	1/1	0.88	0.18	-	90,90,90,90	0
56	MG	1H	3054	1/1	0.93	0.22	-	53,53,53,53	0
56	MG	1H	3222	1/1	0.92	0.40	-	77,77,77,77	0
56	MG	1J	207	1/1	0.86	0.20	-	109,109,109,109	0
56	MG	1H	3417	1/1	0.94	0.13	-	65,65,65,65	0
56	MG	1H	3011	1/1	0.98	0.39	-	54,54,54,54	0
56	MG	13	1646	1/1	0.71	0.26	-	100,100,100,100	0
56	MG	P8	101	1/1	0.79	0.40	-	70,70,70,70	0
56	MG	1H	3280	1/1	0.92	0.25	-	80,80,80,80	0
56	MG	1H	3432	1/1	0.99	0.03	-	83,83,83,83	0
56	MG	14	3250	1/1	0.95	0.37	-	108,108,108,108	0
56	MG	14	3161	1/1	0.93	0.58	-	70,70,70,70	0
56	MG	14	3133	1/1	0.89	0.50	-	82,82,82,82	0
56	MG	14	3274	1/1	0.88	0.08	-	99,99,99,99	0
56	MG	1H	3502	1/1	0.98	0.10	-	49,49,49,49	0
56	MG	1G	1620	1/1	0.57	0.36	-	111,111,111,111	0
56	MG	1G	1625	1/1	0.87	0.67	-	82,82,82,82	0
56	MG	1H	3458	1/1	0.96	0.14	-	109,109,109,109	0
56	MG	1H	3112	1/1	0.97	0.43	-	72,72,72,72	0
56	MG	1H	3305	1/1	0.92	0.27	-	87,87,87,87	0
56	MG	1H	3378	1/1	0.87	0.09	-	83,83,83,83	0
56	MG	1G	1624	1/1	0.48	0.55	-	85,85,85,85	0
56	MG	1H	3365	1/1	0.95	0.06	-	91,91,91,91	0
56	MG	1H	3295	1/1	0.82	0.24	-	94,94,94,94	0
56	MG	1H	3244	1/1	0.72	0.22	-	89,89,89,89	0
56	MG	1H	3016	1/1	0.98	0.40	-	61,61,61,61	0
56	MG	14	3098	1/1	0.89	0.38	-	53,53,53,53	0
56	MG	1H	3093	1/1	0.89	0.74	-	61,61,61,61	0
56	MG	1H	3135	1/1	0.82	0.48	-	68,68,68,68	0
56	MG	14	3343	1/1	0.96	0.04	-	100,100,100,100	0
56	MG	1H	3069	1/1	0.94	0.36	-	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.