



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 15, 2017 – 08:47 AM EST

PDB ID : 4WQF
Title : Crystal structure of the *Thermus thermophilus* 70S ribosome in complex with elongation factor G and fusidic acid in the post-translocational state
Authors : Lin, J.; Gagnon, M.G.; Steitz, T.A.
Deposited on : unknown
Resolution : 2.80 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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-20030345

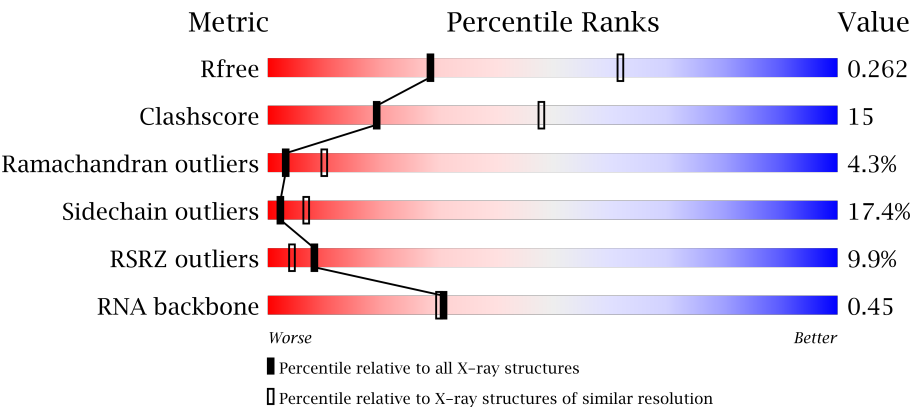
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)
RNA backbone	2435	1007 (3.10-2.50)

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	AA	2915	
1	CA	2915	
2	AB	121	
2	CB	121	





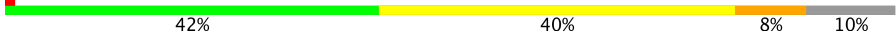




















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Mol	Chain	Length	Quality of chain
3	AC	228	
3	CC	228	
4	AD	276	
4	CD	276	
5	AE	206	
5	CE	206	
6	AF	210	
6	CF	210	
7	AG	182	
7	CG	182	
8	AH	180	
8	CH	180	
9	AK	173	
9	CK	173	
10	AL	147	
10	CL	147	
11	AN	140	
11	CN	140	
12	AO	122	
12	CO	122	
13	AP	150	
13	CP	150	
14	AQ	141	
14	CQ	141	
15	AR	118	

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Mol	Chain	Length	Quality of chain
15	CR	118	
16	AS	112	
16	CS	112	
17	AT	146	
17	CT	146	
18	AU	118	
18	CU	118	
19	AV	101	
19	CV	101	
20	AW	113	
20	CW	113	
21	AX	96	
21	CX	96	
22	AY	110	
22	CY	110	
23	AZ	206	
23	CZ	206	
24	A0	85	
24	C0	85	
25	A1	98	
25	C1	98	
26	A2	72	
26	C2	72	
27	A3	60	
27	C3	60	

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Mol	Chain	Length	Quality of chain
28	A4	71	
28	C4	71	
29	A5	60	
29	C5	60	
30	A6	54	
30	C6	54	
31	A7	49	
31	C7	49	
32	A8	65	
32	C8	65	
33	A9	37	
33	C9	37	
34	BA	1521	
34	DA	1521	
35	BB	256	
35	DB	256	
36	BC	239	
36	DC	239	
37	BD	209	
37	DD	209	
38	BE	162	
38	DE	162	
39	BF	101	
39	DF	101	
40	BG	156	

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Mol	Chain	Length	Quality of chain
40	DG	156	
41	BH	138	
41	DH	138	
42	BI	128	
42	DI	128	
43	BJ	105	
43	DJ	105	
44	BK	129	
44	DK	129	
45	BL	132	
45	DL	132	
46	BM	126	
46	DM	126	
47	BN	61	
47	DN	61	
48	BO	89	
48	DO	89	
49	BP	88	
49	DP	88	
50	BQ	105	
50	DQ	105	
51	BR	88	
51	DR	88	
52	BS	93	
52	DS	93	

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Mol	Chain	Length	Quality of chain
53	BT	106	
53	DT	106	
54	BU	27	
54	DU	27	
55	BV	18	
55	DV	18	
56	BW	76	
56	BY	76	
56	DW	76	
56	DY	76	
57	BZ	758	
57	DZ	758	

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
58	MG	A7	103	-	-	-	X
58	MG	A8	5001	-	-	-	X
58	MG	AA	3012	-	-	-	X
58	MG	AA	3014	-	-	-	X
58	MG	AA	3018	-	-	-	X
58	MG	AA	3020	-	-	-	X
58	MG	AA	3023	-	-	-	X
58	MG	AA	3034	-	-	-	X
58	MG	AA	3035	-	-	-	X
58	MG	AA	3036	-	-	-	X
58	MG	AA	3037	-	-	-	X
58	MG	AA	3039	-	-	-	X
58	MG	AA	3040	-	-	-	X
58	MG	AA	3043	-	-	-	X
58	MG	AA	3044	-	-	-	X
58	MG	AA	3048	-	-	-	X
58	MG	AA	3051	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
58	MG	AA	3059	-	-	-	X
58	MG	AA	3061	-	-	-	X
58	MG	AA	3081	-	-	-	X
58	MG	AA	3082	-	-	-	X
58	MG	AA	3101	-	-	-	X
58	MG	AA	3102	-	-	-	X
58	MG	AA	3109	-	-	-	X
58	MG	AA	3110	-	-	-	X
58	MG	AA	3112	-	-	-	X
58	MG	AA	3113	-	-	-	X
58	MG	AA	3116	-	-	-	X
58	MG	AA	3117	-	-	-	X
58	MG	AA	3120	-	-	-	X
58	MG	AA	3128	-	-	-	X
58	MG	AA	3130	-	-	-	X
58	MG	AA	3132	-	-	-	X
58	MG	AA	3133	-	-	-	X
58	MG	AA	3134	-	-	-	X
58	MG	AA	3135	-	-	-	X
58	MG	AA	3138	-	-	-	X
58	MG	AA	3142	-	-	-	X
58	MG	AA	3150	-	-	-	X
58	MG	AA	3162	-	-	-	X
58	MG	AA	3168	-	-	-	X
58	MG	AA	3171	-	-	-	X
58	MG	AA	3173	-	-	-	X
58	MG	AA	3174	-	-	-	X
58	MG	AA	3179	-	-	-	X
58	MG	AA	3183	-	-	-	X
58	MG	AA	3184	-	-	-	X
58	MG	AA	3185	-	-	-	X
58	MG	AA	3187	-	-	-	X
58	MG	AA	3190	-	-	-	X
58	MG	AA	3196	-	-	-	X
58	MG	AA	3206	-	-	-	X
58	MG	AA	3210	-	-	-	X
58	MG	AA	3211	-	-	-	X
58	MG	AA	3212	-	-	-	X
58	MG	AA	3221	-	-	-	X
58	MG	AA	3223	-	-	-	X
58	MG	AA	3231	-	-	-	X
58	MG	AA	3240	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
58	MG	AA	3247	-	-	-	X
58	MG	AA	3249	-	-	-	X
58	MG	AA	3250	-	-	-	X
58	MG	AA	3253	-	-	-	X
58	MG	AA	3257	-	-	-	X
58	MG	AA	3267	-	-	-	X
58	MG	AA	3272	-	-	-	X
58	MG	AA	3282	-	-	-	X
58	MG	AA	3297	-	-	-	X
58	MG	AA	3301	-	-	-	X
58	MG	AA	3309	-	-	-	X
58	MG	AA	3311	-	-	-	X
58	MG	AA	3314	-	-	-	X
58	MG	AA	3316	-	-	-	X
58	MG	AA	3329	-	-	-	X
58	MG	AA	3331	-	-	-	X
58	MG	AA	3354	-	-	-	X
58	MG	AA	3357	-	-	-	X
58	MG	AA	3372	-	-	-	X
58	MG	AA	3381	-	-	-	X
58	MG	AA	3388	-	-	-	X
58	MG	AA	3395	-	-	-	X
58	MG	AA	3400	-	-	-	X
58	MG	AA	3410	-	-	-	X
58	MG	AA	3418	-	-	-	X
58	MG	AA	3420	-	-	-	X
58	MG	AA	3436	-	-	-	X
58	MG	AA	3439	-	-	-	X
58	MG	AA	3440	-	-	-	X
58	MG	AA	3442	-	-	-	X
58	MG	AA	3443	-	-	-	X
58	MG	AA	3453	-	-	-	X
58	MG	AA	3456	-	-	-	X
58	MG	AA	3462	-	-	-	X
58	MG	AA	3485	-	-	-	X
58	MG	AA	3488	-	-	-	X
58	MG	AA	3499	-	-	-	X
58	MG	AA	3506	-	-	-	X
58	MG	AA	3507	-	-	-	X
58	MG	AA	3508	-	-	-	X
58	MG	AA	3525	-	-	-	X
58	MG	AA	3551	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
58	MG	AA	3559	-	-	-	X
58	MG	AA	3564	-	-	-	X
58	MG	AA	3565	-	-	-	X
58	MG	AA	3581	-	-	-	X
58	MG	AA	3589	-	-	-	X
58	MG	AA	3596	-	-	-	X
58	MG	AA	3602	-	-	-	X
58	MG	AA	3604	-	-	-	X
58	MG	AA	3606	-	-	-	X
58	MG	AA	3617	-	-	-	X
58	MG	AA	3620	-	-	-	X
58	MG	AA	3621	-	-	-	X
58	MG	AA	3623	-	-	-	X
58	MG	AA	3663	-	-	-	X
58	MG	AA	3698	-	-	-	X
58	MG	AA	3702	-	-	-	X
58	MG	AA	3704	-	-	-	X
58	MG	AA	3706	-	-	-	X
58	MG	AA	3708	-	-	-	X
58	MG	AA	3711	-	-	-	X
58	MG	AA	3717	-	-	-	X
58	MG	AA	3726	-	-	-	X
58	MG	AA	3735	-	-	-	X
58	MG	AA	3736	-	-	-	X
58	MG	AA	3739	-	-	-	X
58	MG	AA	3741	-	-	-	X
58	MG	AA	3768	-	-	-	X
58	MG	AA	3770	-	-	-	X
58	MG	AA	3771	-	-	-	X
58	MG	AA	3773	-	-	-	X
58	MG	AA	3791	-	-	-	X
58	MG	AA	3793	-	-	-	X
58	MG	AA	3798	-	-	-	X
58	MG	AA	3805	-	-	-	X
58	MG	AA	3806	-	-	-	X
58	MG	AA	3811	-	-	-	X
58	MG	AA	3812	-	-	-	X
58	MG	AA	3815	-	-	-	X
58	MG	AA	3816	-	-	-	X
58	MG	AA	3817	-	-	-	X
58	MG	AA	3819	-	-	-	X
58	MG	AA	3820	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
58	MG	AA	3822	-	-	-	X
58	MG	AA	3823	-	-	-	X
58	MG	AA	3824	-	-	-	X
58	MG	AA	3827	-	-	-	X
58	MG	AA	3828	-	-	-	X
58	MG	AA	3829	-	-	-	X
58	MG	AA	3830	-	-	-	X
58	MG	AA	3831	-	-	-	X
58	MG	AA	3832	-	-	-	X
58	MG	AA	3833	-	-	-	X
58	MG	AA	3835	-	-	-	X
58	MG	AB	3003	-	-	-	X
58	MG	AB	3008	-	-	-	X
58	MG	AB	3023	-	-	-	X
58	MG	AD	301	-	-	-	X
58	MG	AD	302	-	-	-	X
58	MG	AD	304	-	-	-	X
58	MG	AD	305	-	-	-	X
58	MG	AD	308	-	-	-	X
58	MG	AD	309	-	-	-	X
58	MG	AD	310	-	-	-	X
58	MG	AE	304	-	-	-	X
58	MG	AF	302	-	-	-	X
58	MG	AF	303	-	-	-	X
58	MG	AH	3001	-	-	-	X
58	MG	AH	3002	-	-	-	X
58	MG	AN	3001	-	-	-	X
58	MG	AQ	202	-	-	-	X
58	MG	AU	201	-	-	-	X
58	MG	AU	202	-	-	-	X
58	MG	AU	203	-	-	-	X
58	MG	AV	202	-	-	-	X
58	MG	AW	3003	-	-	-	X
58	MG	AX	101	-	-	-	X
58	MG	BA	1615	-	-	-	X
58	MG	BA	1616	-	-	-	X
58	MG	BA	1623	-	-	-	X
58	MG	BA	1626	-	-	-	X
58	MG	BA	1629	-	-	-	X
58	MG	BA	1630	-	-	-	X
58	MG	BA	1648	-	-	-	X
58	MG	BA	1657	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
58	MG	BA	1664	-	-	-	X
58	MG	BA	1671	-	-	-	X
58	MG	BA	1678	-	-	-	X
58	MG	BA	1683	-	-	-	X
58	MG	BA	1686	-	-	-	X
58	MG	BA	1690	-	-	-	X
58	MG	BA	1693	-	-	-	X
58	MG	BA	1695	-	-	-	X
58	MG	BA	1711	-	-	-	X
58	MG	BA	1721	-	-	-	X
58	MG	BA	1723	-	-	-	X
58	MG	BA	1738	-	-	-	X
58	MG	BA	1740	-	-	-	X
58	MG	BA	1755	-	-	-	X
58	MG	BA	1756	-	-	-	X
58	MG	BA	1763	-	-	-	X
58	MG	BA	1783	-	-	-	X
58	MG	BA	1801	-	-	-	X
58	MG	BA	1811	-	-	-	X
58	MG	BT	3001	-	-	-	X
58	MG	C3	3001	-	-	-	X
58	MG	C7	101	-	-	-	X
58	MG	CA	3002	-	-	-	X
58	MG	CA	3011	-	-	-	X
58	MG	CA	3013	-	-	-	X
58	MG	CA	3014	-	-	-	X
58	MG	CA	3023	-	-	-	X
58	MG	CA	3027	-	-	-	X
58	MG	CA	3028	-	-	-	X
58	MG	CA	3030	-	-	-	X
58	MG	CA	3035	-	-	-	X
58	MG	CA	3037	-	-	-	X
58	MG	CA	3038	-	-	-	X
58	MG	CA	3041	-	-	-	X
58	MG	CA	3043	-	-	-	X
58	MG	CA	3073	-	-	-	X
58	MG	CA	3084	-	-	-	X
58	MG	CA	3086	-	-	-	X
58	MG	CA	3088	-	-	-	X
58	MG	CA	3091	-	-	-	X
58	MG	CA	3106	-	-	-	X
58	MG	CA	3109	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
58	MG	CA	3110	-	-	-	X
58	MG	CA	3114	-	-	-	X
58	MG	CA	3119	-	-	-	X
58	MG	CA	3124	-	-	-	X
58	MG	CA	3127	-	-	-	X
58	MG	CA	3133	-	-	-	X
58	MG	CA	3137	-	-	-	X
58	MG	CA	3146	-	-	-	X
58	MG	CA	3159	-	-	-	X
58	MG	CA	3163	-	-	-	X
58	MG	CA	3166	-	-	-	X
58	MG	CA	3168	-	-	-	X
58	MG	CA	3169	-	-	-	X
58	MG	CA	3177	-	-	-	X
58	MG	CA	3178	-	-	-	X
58	MG	CA	3182	-	-	-	X
58	MG	CA	3185	-	-	-	X
58	MG	CA	3190	-	-	-	X
58	MG	CA	3201	-	-	-	X
58	MG	CA	3212	-	-	-	X
58	MG	CA	3213	-	-	-	X
58	MG	CA	3217	-	-	-	X
58	MG	CA	3218	-	-	-	X
58	MG	CA	3221	-	-	-	X
58	MG	CA	3223	-	-	-	X
58	MG	CA	3226	-	-	-	X
58	MG	CA	3229	-	-	-	X
58	MG	CA	3230	-	-	-	X
58	MG	CA	3251	-	-	-	X
58	MG	CA	3263	-	-	-	X
58	MG	CA	3266	-	-	-	X
58	MG	CA	3276	-	-	-	X
58	MG	CA	3281	-	-	-	X
58	MG	CA	3290	-	-	-	X
58	MG	CA	3302	-	-	-	X
58	MG	CA	3309	-	-	-	X
58	MG	CA	3313	-	-	-	X
58	MG	CA	3314	-	-	-	X
58	MG	CA	3322	-	-	-	X
58	MG	CA	3324	-	-	-	X
58	MG	CA	3326	-	-	-	X
58	MG	CA	3328	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
58	MG	CA	3330	-	-	-	X
58	MG	CA	3340	-	-	-	X
58	MG	CA	3348	-	-	-	X
58	MG	CA	3353	-	-	-	X
58	MG	CA	3358	-	-	-	X
58	MG	CA	3361	-	-	-	X
58	MG	CA	3364	-	-	-	X
58	MG	CA	3375	-	-	-	X
58	MG	CA	3383	-	-	-	X
58	MG	CA	3392	-	-	-	X
58	MG	CA	3396	-	-	-	X
58	MG	CA	3409	-	-	-	X
58	MG	CA	3420	-	-	-	X
58	MG	CA	3428	-	-	-	X
58	MG	CA	3432	-	-	-	X
58	MG	CA	3441	-	-	-	X
58	MG	CA	3452	-	-	-	X
58	MG	CA	3455	-	-	-	X
58	MG	CA	3458	-	-	-	X
58	MG	CA	3463	-	-	-	X
58	MG	CA	3467	-	-	-	X
58	MG	CA	3476	-	-	-	X
58	MG	CA	3489	-	-	-	X
58	MG	CA	3490	-	-	-	X
58	MG	CA	3491	-	-	-	X
58	MG	CA	3498	-	-	-	X
58	MG	CA	3500	-	-	-	X
58	MG	CA	3530	-	-	-	X
58	MG	CA	3542	-	-	-	X
58	MG	CA	3555	-	-	-	X
58	MG	CA	3567	-	-	-	X
58	MG	CA	3588	-	-	-	X
58	MG	CA	3596	-	-	-	X
58	MG	CA	3597	-	-	-	X
58	MG	CA	3603	-	-	-	X
58	MG	CA	3607	-	-	-	X
58	MG	CA	3618	-	-	-	X
58	MG	CA	3619	-	-	-	X
58	MG	CA	3626	-	-	-	X
58	MG	CA	3650	-	-	-	X
58	MG	CA	3654	-	-	-	X
58	MG	CA	3660	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
58	MG	CA	3661	-	-	-	X
58	MG	CD	303	-	-	-	X
58	MG	CE	301	-	-	-	X
58	MG	CE	303	-	-	-	X
58	MG	CF	301	-	-	-	X
58	MG	CF	303	-	-	-	X
58	MG	CQ	201	-	-	-	X
58	MG	CU	201	-	-	-	X
58	MG	CV	202	-	-	-	X
58	MG	DA	1606	-	-	-	X
58	MG	DA	1610	-	-	-	X
58	MG	DA	1618	-	-	-	X
58	MG	DA	1636	-	-	-	X
58	MG	DA	1638	-	-	-	X
58	MG	DA	1647	-	-	-	X
58	MG	DA	1649	-	-	-	X
58	MG	DA	1651	-	-	-	X
58	MG	DA	1658	-	-	-	X
58	MG	DA	1668	-	-	-	X
58	MG	DA	1669	-	-	-	X
58	MG	DA	1672	-	-	-	X
58	MG	DA	1680	-	-	-	X
58	MG	DA	1684	-	-	-	X
58	MG	DA	1689	-	-	-	X
58	MG	DA	1694	-	-	-	X
58	MG	DA	1697	-	-	-	X
58	MG	DA	1722	-	-	-	X
58	MG	DA	1743	-	-	-	X
58	MG	DA	1768	-	-	-	X
58	MG	DF	3001	-	-	-	X
58	MG	DT	3001	-	-	-	X
60	SF4	DD	501	-	-	X	-
62	GDP	DZ	704	-	-	X	-

2 Entry composition

There are 63 unique types of molecules in this entry. The entry contains 310279 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 23S Ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	2852	Total	C	N	O	P	0	0	0
			61426	27339	11489	19747	2851			
1	CA	2848	Total	C	N	O	P	0	0	0
			61337	27299	11470	19721	2847			

- Molecule 2 is a RNA chain called 5S Ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	AB	120	Total	C	N	O	P	0	0	0
			2573	1146	476	832	119			
2	CB	120	Total	C	N	O	P	0	0	0
			2573	1146	476	832	119			

- Molecule 3 is a protein called 50S ribosomal protein L1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	AC	137	Total	C	N	O	S	0	0	0
			1063	669	201	192	1			
3	CC	137	Total	C	N	O	S	0	0	0
			1063	669	201	192	1			

- Molecule 4 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	AD	275	Total	C	N	O	S	0	0	0
			2136	1349	423	361	3			
4	CD	275	Total	C	N	O	S	0	0	0
			2142	1352	426	361	3			

- Molecule 5 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	AE	204	Total	C	N	O	S	0	0	0
			1559	985	298	270	6			
5	CE	204	Total	C	N	O	S	0	0	0
			1559	985	298	270	6			

- Molecule 6 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	AF	203	Total	C	N	O	S	0	0	1
			1584	1009	298	275	2			
6	CF	203	Total	C	N	O	S	0	0	1
			1580	1007	297	274	2			

- Molecule 7 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	AG	181	Total	C	N	O	S	0	0	0
			1425	914	256	251	4			
7	CG	181	Total	C	N	O	S	0	0	0
			1424	911	258	251	4			

- Molecule 8 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	AH	174	Total	C	N	O	S	0	0	0
			1330	845	248	236	1			
8	CH	174	Total	C	N	O	S	0	0	0
			1330	845	248	236	1			

- Molecule 9 is a protein called 50S ribosomal protein L10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	AK	130	Total	C	N	O		0	0	0
			641	381	130	130				
9	CK	130	Total	C	N	O		0	0	0
			641	381	130	130				

- Molecule 10 is a protein called 50S ribosomal protein L11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	AL	139	Total	C	N	O	S	0	0	0
			1025	653	181	186	5			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	CL	139	Total	C	N	O	S	0	0	0
			1025	653	181	186	5			

- Molecule 11 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	AN	140	Total	C	N	O	S	0	0	0
			1117	719	207	187	4			
11	CN	140	Total	C	N	O	S	0	0	0
			1117	719	207	187	4			

- Molecule 12 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	AO	122	Total	C	N	O	S	0	0	0
			933	588	171	170	4			
12	CO	122	Total	C	N	O	S	0	0	0
			933	588	171	170	4			

- Molecule 13 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	AP	149	Total	C	N	O	S	0	0	0
			1139	709	231	196	3			
13	CP	149	Total	C	N	O	S	0	0	0
			1135	706	230	196	3			

- Molecule 14 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	AQ	141	Total	C	N	O	S	0	0	0
			1122	715	212	188	7			
14	CQ	141	Total	C	N	O	S	0	0	0
			1122	715	212	188	7			

- Molecule 15 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	AR	118	Total	C	N	O	S	0	0	0
			968	604	203	160	1			
15	CR	118	Total	C	N	O	S	0	0	0
			968	604	203	160	1			

- Molecule 16 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
16	AS	110	Total	C	N	O	0	0	0
			877	553	175	149			
16	CS	110	Total	C	N	O	0	0	0
			870	549	173	148			

- Molecule 17 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	AT	131	Total	C	N	O	S	0	0	0
			1091	680	225	185	1			
17	CT	131	Total	C	N	O	S	0	0	0
			1083	675	224	183	1			

- Molecule 18 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
18	AU	116	Total	C	N	O	S	0	0	0
			959	608	201	149	1			
18	CU	116	Total	C	N	O	S	0	0	0
			959	608	201	149	1			

- Molecule 19 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	AV	101	Total	C	N	O	S	0	0	0
			771	495	140	135	1			
19	CV	101	Total	C	N	O	S	0	0	0
			771	495	140	135	1			

- Molecule 20 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	AW	112	Total	C	N	O	S	0	0	0
			886	557	174	153	2			
20	CW	112	Total	C	N	O	S	0	0	0
			886	557	174	153	2			

- Molecule 21 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
21	AX	95	Total	C	N	O	S	0	0	0
			750	488	135	126	1			
21	CX	95	Total	C	N	O	S	0	0	0
			750	488	135	126	1			

- Molecule 22 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	AY	107	Total	C	N	O	S	0	0	0
			806	517	152	131	6			
22	CY	107	Total	C	N	O	S	0	0	0
			806	517	152	131	6			

- Molecule 23 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
23	AZ	185	Total	C	N	O	S	0	0	0
			1451	927	258	264	2			
23	CZ	185	Total	C	N	O	S	0	0	0
			1451	927	258	264	2			

- Molecule 24 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
24	A0	83	Total	C	N	O	S	0	0	0
			653	404	139	109	1			
24	C0	83	Total	C	N	O	S	0	0	0
			653	404	139	109	1			

- Molecule 25 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
25	A1	97	Total	C	N	O	S	0	0	0
			755	475	148	131	1			
25	C1	97	Total	C	N	O	S	0	0	0
			755	475	148	131	1			

- Molecule 26 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	A2	70	Total	C	N	O	S	0	0	0
			588	365	118	103	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
26	C2	70	Total	C	N	O	S	0	0	0
			588	365	118	103	2			

- Molecule 27 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
27	A3	59	Total	C	N	O		0	0	0
			469	298	90	81				
27	C3	59	Total	C	N	O		0	0	0
			464	296	90	78				

- Molecule 28 is a protein called 50S ribosomal protein L31.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
28	A4	69	Total	C	N	O	S	0	0	0
			558	352	102	99	5			
28	C4	69	Total	C	N	O	S	0	0	0
			532	339	97	91	5			

- Molecule 29 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
29	A5	59	Total	C	N	O	S	0	0	0
			455	285	89	76	5			
29	C5	59	Total	C	N	O	S	0	0	0
			455	285	89	76	5			

- Molecule 30 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
30	A6	53	Total	C	N	O	S	0	0	0
			453	281	91	77	4			
30	C6	53	Total	C	N	O	S	0	0	0
			449	279	91	75	4			

- Molecule 31 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
31	A7	48	Total	C	N	O	S	0	0	0
			418	257	104	55	2			
31	C7	48	Total	C	N	O	S	0	0	0
			418	257	104	55	2			

- Molecule 32 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
32	A8	64	Total	C	N	O	S	0	0	0
			517	331	102	82	2			
32	C8	64	Total	C	N	O	S	0	0	0
			517	331	102	82	2			

- Molecule 33 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
33	A9	37	Total	C	N	O	S	0	0	0
			307	188	68	47	4			
33	C9	37	Total	C	N	O	S	0	0	0
			307	188	68	47	4			

- Molecule 34 is a RNA chain called 16S Ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
34	BA	1495	Total	C	N	O	P	0	0	0
			32141	14304	5958	10384	1495			
34	DA	1501	Total	C	N	O	P	0	0	0
			32268	14361	5980	10426	1501			

- Molecule 35 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
35	BB	231	Total	C	N	O	S	0	0	0
			1846	1179	331	331	5			
35	DB	231	Total	C	N	O	S	0	0	0
			1825	1167	326	327	5			

- Molecule 36 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
36	BC	206	Total	C	N	O	S	0	0	0
			1552	976	302	273	1			
36	DC	206	Total	C	N	O	S	0	0	0
			1544	970	300	273	1			

- Molecule 37 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
37	BD	208	Total	C	N	O	S	0	0	0
			1659	1040	326	286	7			
37	DD	208	Total	C	N	O	S	0	0	0
			1678	1052	333	286	7			

- Molecule 38 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
38	BE	148	Total	C	N	O	S	0	0	0
			1129	714	213	198	4			
38	DE	148	Total	C	N	O	S	0	0	0
			1133	716	214	199	4			

- Molecule 39 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
39	BF	100	Total	C	N	O	S	0	0	0
			812	514	146	149	3			
39	DF	100	Total	C	N	O	S	0	0	0
			820	518	147	152	3			

- Molecule 40 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
40	BG	155	Total	C	N	O	S	0	0	0
			1231	766	243	216	6			
40	DG	155	Total	C	N	O	S	0	0	0
			1235	769	244	216	6			

- Molecule 41 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
41	BH	137	Total	C	N	O	S	0	0	0
			1088	689	206	191	2			
41	DH	137	Total	C	N	O	S	0	0	0
			1088	689	206	191	2			

- Molecule 42 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
42	BI	127	Total	C	N	O		0	0	0
			986	626	193	167				

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
42	DI	127	Total	C	N	O	0	0	0
			978	619	190	169			

- Molecule 43 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
43	BJ	97	Total	C	N	O	0	0	0
			709	440	138	131			
43	DJ	96	Total	C	N	O	0	0	0
			714	445	138	131			

- Molecule 44 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
44	BK	114	Total	C	N	O	S	0	0	0
			833	519	156	155	3			
44	DK	114	Total	C	N	O	S	0	0	0
			833	519	156	155	3			

- Molecule 45 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
45	BL	122	Total	C	N	O	S	0	0	0
			930	585	185	159	1			
45	DL	122	Total	C	N	O	S	0	0	0
			930	585	185	159	1			

- Molecule 46 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
46	BM	117	Total	C	N	O	S	0	0	0
			923	570	191	160	2			
46	DM	122	Total	C	N	O	S	0	0	0
			950	586	197	165	2			

- Molecule 47 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
47	BN	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			
47	DN	60	Total	C	N	O	S	0	0	0
			492	312	104	72	4			

- Molecule 48 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
48	BO	88	Total	C	N	O	S	0	0	0
			728	456	144	126	2			
48	DO	88	Total	C	N	O	S	0	0	0
			728	456	144	126	2			

- Molecule 49 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
49	BP	82	Total	C	N	O	S	0	0	0
			681	433	134	113	1			
49	DP	82	Total	C	N	O	S	0	0	0
			677	430	133	113	1			

- Molecule 50 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
50	BQ	99	Total	C	N	O	S	0	0	0
			823	528	151	142	2			
50	DQ	99	Total	C	N	O	S	0	0	0
			823	528	151	142	2			

- Molecule 51 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
51	BR	68	Total	C	N	O	0	0	0
			555	355	108	92			
51	DR	68	Total	C	N	O	0	0	0
			555	355	108	92			

- Molecule 52 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
52	BS	84	Total	C	N	O	S	0	0	0
			661	423	122	114	2			
52	DS	83	Total	C	N	O	S	0	0	0
			646	412	119	113	2			

- Molecule 53 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
53	BT	96	Total	C	N	O	S	0	0	0
			728	446	156	124	2			
53	DT	96	Total	C	N	O	S	0	0	0
			731	449	156	124	2			

- Molecule 54 is a protein called 30S ribosomal protein Thx.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
54	BU	23	Total	C	N	O		0	0	0
			199	122	48	29				
54	DU	23	Total	C	N	O		0	0	0
			199	122	48	29				

- Molecule 55 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
55	BV	7	Total	C	N	O	P	0	0	0
			148	67	27	47	7			
55	DV	6	Total	C	N	O	P	0	0	0
			123	57	22	39	5			

- Molecule 56 is a RNA chain called P-site tRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
56	BW	76	Total	C	N	O	P	S	0	0
			1631	731	290	532	76	2		
56	BY	74	Total	C	N	O	P	S	0	0
			1581	707	285	515	73	1		
56	DW	76	Total	C	N	O	P	S	0	0
			1631	731	290	532	76	2		
56	DY	73	Total	C	N	O	P	S	0	0
			1561	698	283	507	72	1		

- Molecule 57 is a protein called 50S ribosomal protein L9,Elongation factor G.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
57	BZ	730	Total	C	N	O	S	0	0	0
			5690	3616	980	1075	19			
57	DZ	730	Total	C	N	O	S	0	0	0
			5690	3616	980	1075	19			

- Molecule 58 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
58	AP	2	Total 2 Mg 2	0	0
58	CR	1	Total 1 Mg 1	0	0
58	BA	211	Total 211 Mg 211	0	0
58	CA	664	Total 664 Mg 664	0	0
58	C8	1	Total 1 Mg 1	0	0
58	C5	1	Total 1 Mg 1	0	0
58	AB	23	Total 23 Mg 23	0	0
58	BL	2	Total 2 Mg 2	0	0
58	CV	2	Total 2 Mg 2	0	0
58	A6	1	Total 1 Mg 1	0	0
58	BE	1	Total 1 Mg 1	0	0
58	AW	4	Total 4 Mg 4	0	0
58	AN	3	Total 3 Mg 3	0	0
58	DZ	2	Total 2 Mg 2	0	0
58	AX	1	Total 1 Mg 1	0	0
58	CN	1	Total 1 Mg 1	0	0
58	A2	2	Total 2 Mg 2	0	0
58	CY	1	Total 1 Mg 1	0	0
58	DD	1	Total 1 Mg 1	0	0
58	BB	1	Total 1 Mg 1	0	0
58	BT	1	Total 1 Mg 1	0	0
58	AE	4	Total 4 Mg 4	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
58	BM	2	Total 2	Mg 2	0	0
58	CU	1	Total 1	Mg 1	0	0
58	BF	1	Total 1	Mg 1	0	0
58	AV	3	Total 3	Mg 3	0	0
58	DA	168	Total 168	Mg 168	0	0
58	CB	13	Total 13	Mg 13	0	0
58	AA	835	Total 835	Mg 835	0	0
58	CQ	5	Total 5	Mg 5	0	0
58	A5	1	Total 1	Mg 1	0	0
58	AR	1	Total 1	Mg 1	0	0
58	CG	1	Total 1	Mg 1	0	0
58	DK	1	Total 1	Mg 1	0	0
58	DF	1	Total 1	Mg 1	0	0
58	AD	10	Total 10	Mg 10	0	0
58	BN	1	Total 1	Mg 1	0	0
58	DJ	1	Total 1	Mg 1	0	0
58	C7	1	Total 1	Mg 1	0	0
58	C3	1	Total 1	Mg 1	0	0
58	AZ	2	Total 2	Mg 2	0	0
58	A4	1	Total 1	Mg 1	0	0
58	BK	1	Total 1	Mg 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
58	AU	3	Total 3	Mg 3	0	0
58	DW	3	Total 3	Mg 3	0	0
58	A9	1	Total 1	Mg 1	0	0
58	CF	5	Total 5	Mg 5	0	0
58	BV	1	Total 1	Mg 1	0	0
58	A0	3	Total 3	Mg 3	0	0
58	AG	2	Total 2	Mg 2	0	0
58	DE	2	Total 2	Mg 2	0	0
58	AQ	3	Total 3	Mg 3	0	0
58	CE	6	Total 6	Mg 6	0	0
58	AH	2	Total 2	Mg 2	0	0
58	BZ	2	Total 2	Mg 2	0	0
58	CO	2	Total 2	Mg 2	0	0
58	CP	3	Total 3	Mg 3	0	0
58	CW	1	Total 1	Mg 1	0	0
58	A7	3	Total 3	Mg 3	0	0
58	CD	3	Total 3	Mg 3	0	0
58	BD	1	Total 1	Mg 1	0	0
58	DT	1	Total 1	Mg 1	0	0
58	A8	2	Total 2	Mg 2	0	0
58	AO	1	Total 1	Mg 1	0	0

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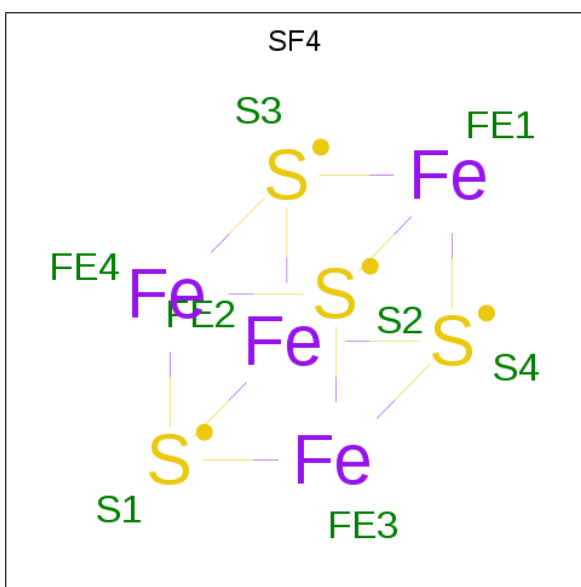
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
58	BW	3	Total 3	Mg 3	0	0
58	AY	1	Total 1	Mg 1	0	0
58	AF	5	Total 5	Mg 5	0	0

- Molecule 59 is ZINC ION (three-letter code: ZN) (formula: Zn).

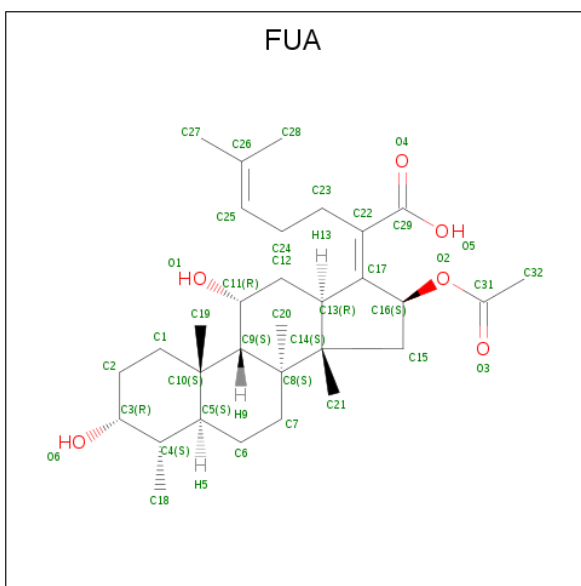
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
59	AY	1	Total 1	Zn 1	0	0
59	BN	1	Total 1	Zn 1	0	0
59	C4	1	Total 1	Zn 1	0	0
59	C5	1	Total 1	Zn 1	0	0
59	C6	1	Total 1	Zn 1	0	0
59	A6	1	Total 1	Zn 1	0	0
59	C9	1	Total 1	Zn 1	0	0
59	DN	1	Total 1	Zn 1	0	0
59	A4	1	Total 1	Zn 1	0	0
59	A5	1	Total 1	Zn 1	0	0
59	A9	1	Total 1	Zn 1	0	0
59	CY	1	Total 1	Zn 1	0	0

- Molecule 60 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe₄S₄).



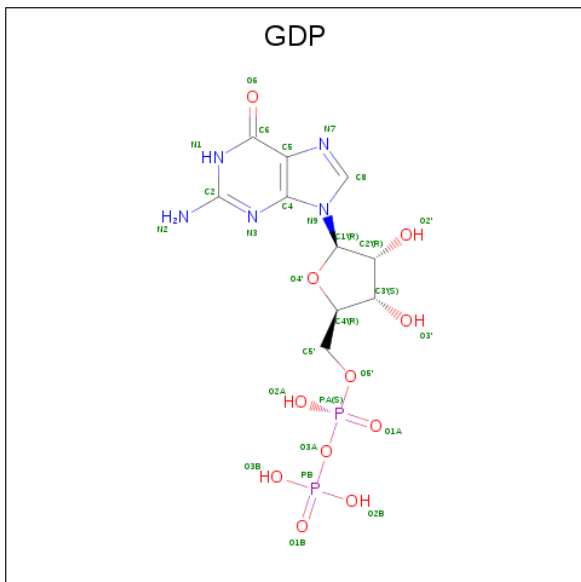
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
60	BD	1	Total	Fe	S	0	0
			8	4	4		
60	DD	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 61 is FUSIDIC ACID (three-letter code: FUA) (formula: $C_{31}H_{48}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
61	BZ	1	Total	C	O	0	0
			37	31	6		
61	DZ	1	Total	C	O	0	0
			37	31	6		

- Molecule 62 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
62	BZ	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
62	DZ	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

- Molecule 63 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
63	AA	1408	Total	O	0	0
			1408	1408		
63	AB	36	Total	O	0	0
			36	36		
63	AD	15	Total	O	0	0
			15	15		
63	AE	19	Total	O	0	0
			19	19		
63	AF	8	Total	O	0	0
			8	8		
63	AG	3	Total	O	0	0
			3	3		
63	AH	1	Total	O	0	0
			1	1		
63	AN	2	Total	O	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
63	AO	3	Total 3	O 3	0	0
63	AP	15	Total 15	O 15	0	0
63	AQ	3	Total 3	O 3	0	0
63	AR	3	Total 3	O 3	0	0
63	AS	1	Total 1	O 1	0	0
63	AT	2	Total 2	O 2	0	0
63	AU	6	Total 6	O 6	0	0
63	AW	1	Total 1	O 1	0	0
63	AX	2	Total 2	O 2	0	0
63	AZ	1	Total 1	O 1	0	0
63	A0	7	Total 7	O 7	0	0
63	A1	3	Total 3	O 3	0	0
63	A3	1	Total 1	O 1	0	0
63	A5	2	Total 2	O 2	0	0
63	A6	1	Total 1	O 1	0	0
63	A7	3	Total 3	O 3	0	0
63	A8	11	Total 11	O 11	0	0
63	BA	205	Total 205	O 205	0	0
63	BD	3	Total 3	O 3	0	0
63	BE	3	Total 3	O 3	0	0
63	BJ	1	Total 1	O 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
63	BL	2	Total	O	0	0
			2	2		
63	BM	1	Total	O	0	0
			1	1		
63	BO	1	Total	O	0	0
			1	1		
63	BV	2	Total	O	0	0
			2	2		
63	BW	1	Total	O	0	0
			1	1		
63	BZ	3	Total	O	0	0
			3	3		
63	CA	981	Total	O	0	0
			981	981		
63	CB	9	Total	O	0	0
			9	9		
63	CD	15	Total	O	0	0
			15	15		
63	CE	9	Total	O	0	0
			9	9		
63	CF	6	Total	O	0	0
			6	6		
63	CP	13	Total	O	0	0
			13	13		
63	CQ	1	Total	O	0	0
			1	1		
63	CT	3	Total	O	0	0
			3	3		
63	CU	4	Total	O	0	0
			4	4		
63	CV	1	Total	O	0	0
			1	1		
63	CW	1	Total	O	0	0
			1	1		
63	CX	1	Total	O	0	0
			1	1		
63	CY	1	Total	O	0	0
			1	1		
63	C0	5	Total	O	0	0
			5	5		
63	C1	3	Total	O	0	0
			3	3		

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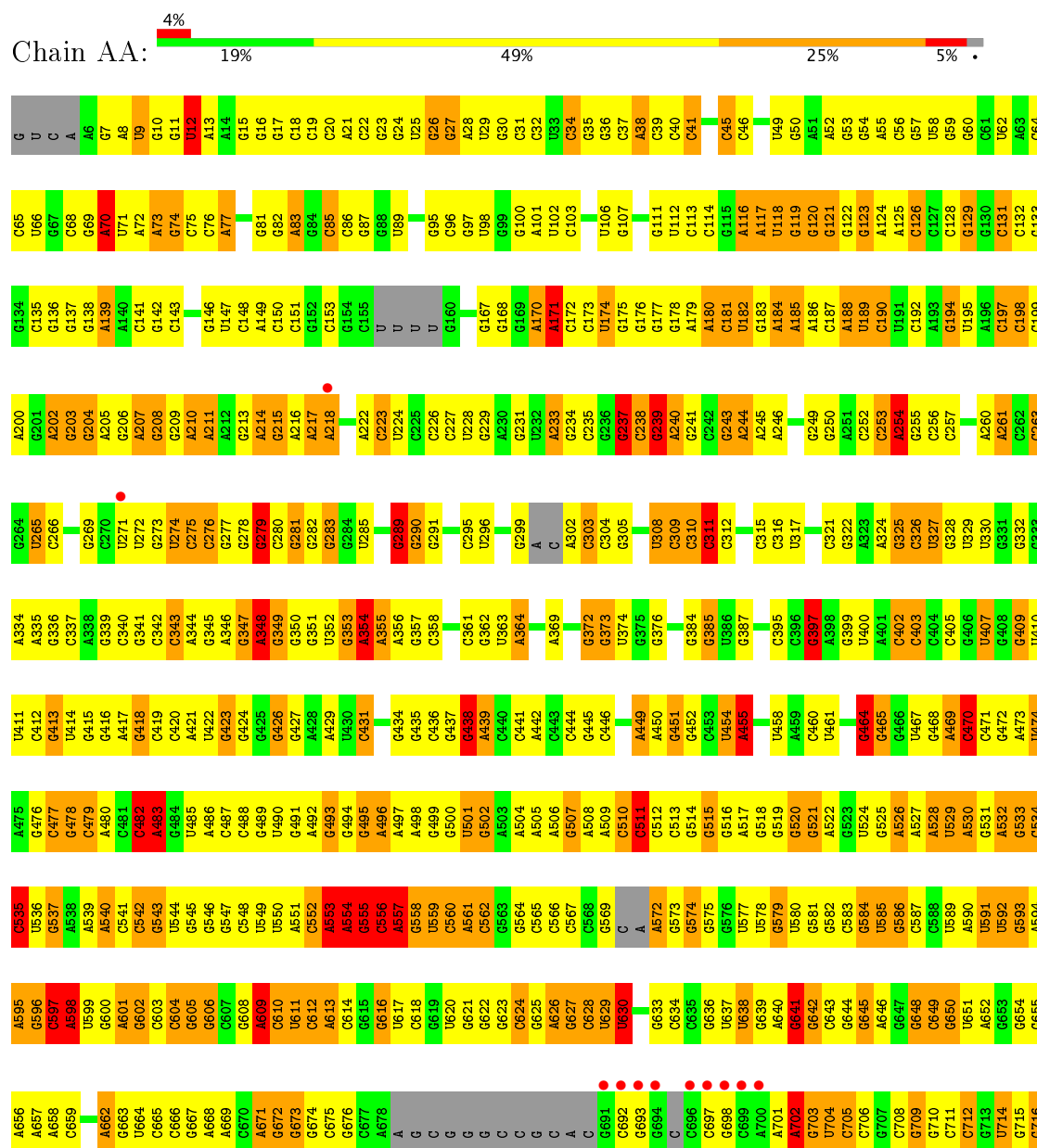
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
63	C3	1	Total 1	O 1	0	0
63	C5	1	Total 1	O 1	0	0
63	C7	3	Total 3	O 3	0	0
63	C8	3	Total 3	O 3	0	0
63	DA	153	Total 153	O 153	0	0
63	DE	2	Total 2	O 2	0	0
63	DH	1	Total 1	O 1	0	0
63	DJ	1	Total 1	O 1	0	0
63	DK	2	Total 2	O 2	0	0
63	DL	1	Total 1	O 1	0	0
63	DP	1	Total 1	O 1	0	0
63	DT	1	Total 1	O 1	0	0
63	DY	1	Total 1	O 1	0	0
63	DZ	2	Total 2	O 2	0	0

3 Residue-property plots

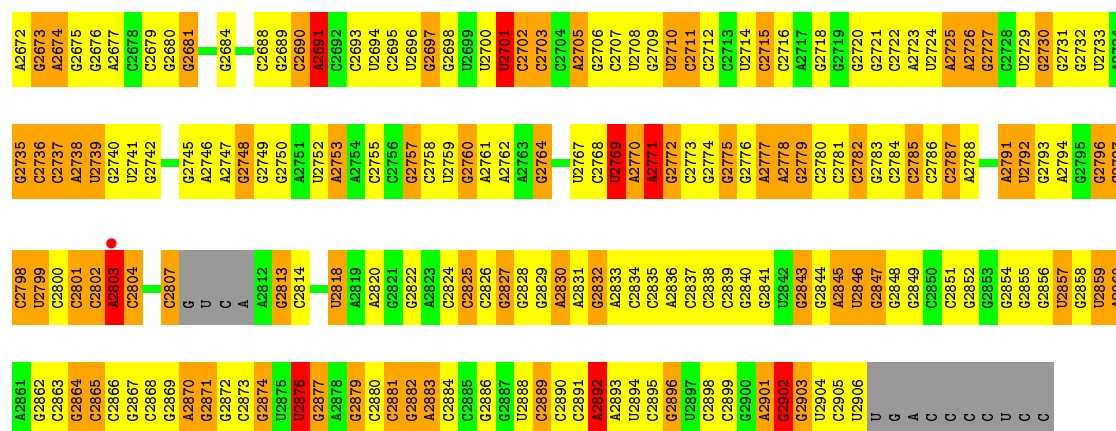
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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 ($RSRZ > 2$). 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: 23S Ribosomal RNA

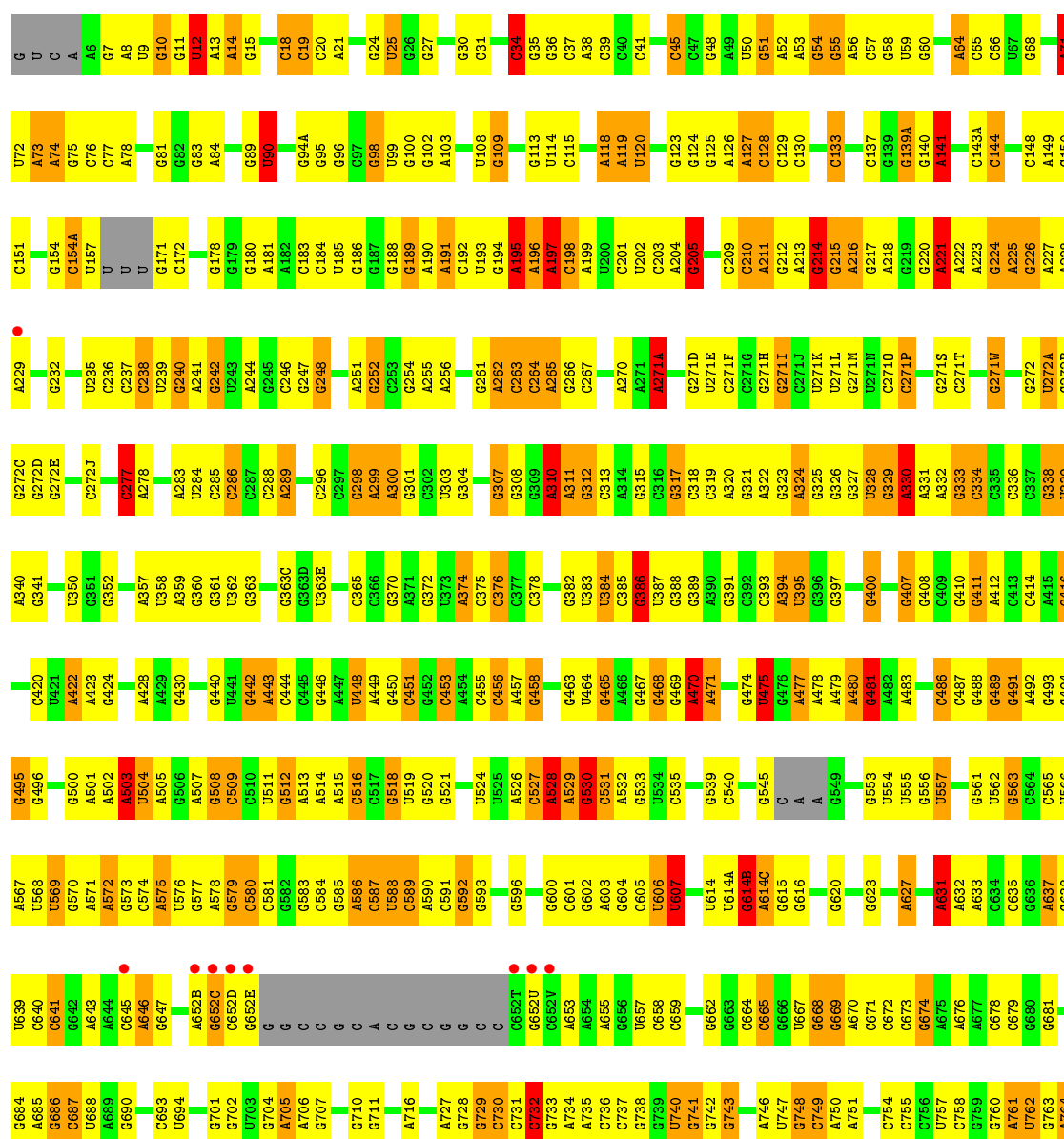




U2608	U2609	U2610	U2611	U2612	U2613	U2614	U2615	U2616	U2617	U2618	U2619	U2620	U2621	U2622	U2623	U2624	U2625	U2626	U2627	U2628	U2629	U2630	U2631	U2632	U2633	U2634	U2635	U2636	U2637	U2638	U2639	U2640	U2641	U2642	U2643	U2644	U2645	U2646	U2647																															
C2488	C2489	C2490	C2491	C2492	C2493	C2494	C2495	C2496	C2497	C2498	C2499	C2500	C2501	C2502	C2503	C2504	C2505	C2506	C2507	C2508	C2509	C2510	C2511	C2512	C2513	C2514	C2515	C2516	C2517	C2518	C2519	C2520	C2521	C2522	C2523	C2524	C2525	C2526	C2527	C2528	C2529	C2530	C2531	C2532	C2533	C2534	C2535	C2536	C2537	C2538	C2539	C2540	C2541	C2542	C2543	C2544	C2545	C2546	C2547											
G2488	G2489	G2490	G2491	G2492	G2493	G2494	G2495	G2496	G2497	G2498	G2499	G2500	G2501	G2502	G2503	G2504	G2505	G2506	G2507	G2508	G2509	G2510	G2511	G2512	G2513	G2514	G2515	G2516	G2517	G2518	G2519	G2520	G2521	G2522	G2523	G2524	G2525	G2526	G2527	G2528	G2529	G2530	G2531	G2532	G2533	G2534	G2535	G2536	G2537	G2538	G2539	G2540	G2541	G2542	G2543	G2544	G2545	G2546	G2547											
C2428	C2429	C2430	C2431	C2432	C2433	C2434	C2435	C2436	C2437	C2438	C2439	C2440	C2441	C2442	C2443	C2444	C2445	C2446	C2447	C2448	C2449	C2450	C2451	C2452	C2453	C2454	C2455	C2456	C2457	C2458	C2459	C2460	C2461	C2462	C2463	C2464	C2465	C2466	C2467	C2468	C2469	C2470	C2471	C2472	C2473	C2474	C2475	C2476	C2477	C2478	C2479	C2480	C2481	C2482	C2483	C2484	C2485	C2486	C2487											
C2367	C2368	C2369	C2370	C2371	C2372	C2373	C2374	C2375	C2376	C2377	C2378	C2379	C2380	C2381	C2382	C2383	C2384	C2385	C2386	C2387	C2388	C2389	C2390	C2391	C2392	C2393	C2394	C2395	C2396	C2397	C2398	C2399	C2400	C2401	C2402	C2403	C2404	C2405	C2406	C2407	C2408	C2409	C2410	C2411	C2412	C2413	C2414	C2415	C2416	C2417	C2418	C2419	C2420	C2421	C2422	C2423	C2424	C2425	C2426	C2427										
C2304	C2305	C2306	C2307	C2308	C2309	C2310	C2311	C2312	C2313	C2314	C2315	C2316	C2317	C2318	C2319	C2320	C2321	C2322	C2323	C2324	C2325	C2326	C2327	C2328	C2329	C2330	C2331	C2332	C2333	C2334	C2335	C2336	C2337	C2338	C2339	C2340	C2341	C2342	C2343	C2344	C2345	C2346	C2347	C2348	C2349	C2350	C2351	C2352	C2353	C2354	C2355	C2356	C2357	C2358	C2359	C2360	C2361	C2362	C2363	C2364	C2365	C2366								
G2242	G2243	G2244	G2245	G2246	G2247	G2248	G2249	G2250	G2251	G2252	G2253	G2254	G2255	G2256	G2257	G2258	G2259	G2260	G2261	G2262	G2263	G2264	G2265	G2266	G2267	G2268	G2269	G2270	G2271	G2272	G2273	G2274	G2275	G2276	G2277	G2278	G2279	G2280	G2281	G2282	G2283	G2284	G2285	G2286	G2287	G2288	G2289	G2290	G2291	G2292	G2293	G2294	G2295	G2296	G2297	G2298	G2299	G2300	G2301	G2302	G2303									
C2179	C2180	C2181	C2182	C2183	C2184	C2185	C2186	C2187	C2188	C2189	C2190	C2191	C2192	C2193	C2194	C2195	C2196	C2197	C2198	C2199	C2200	C2201	C2202	C2203	C2204	C2205	C2206	C2207	C2208	C2209	C2210	C2211	C2212	C2213	C2214	C2215	C2216	C2217	C2218	C2219	C2220	C2221	C2222	C2223	C2224	C2225	C2226	C2227	C2228	C2229	C2230	C2231	C2232	C2233	C2234	C2235	C2236	C2237	C2238	C2239	C2240	C2241								
U2119	U2120	U2121	U2122	U2123	U2124	U2125	U2126	U2127	U2128	U2129	U2130	U2131	U2132	U2133	U2134	U2135	U2136	U2137	U2138	U2139	U2140	U2141	U2142	U2143	U2144	U2145	U2146	U2147	U2148	U2149	U2150	U2151	U2152	U2153	U2154	U2155	U2156	U2157	U2158	U2159	U2160	U2161	U2162	U2163	U2164	U2165	U2166	U2167	U2168	U2169	U2170	U2171	U2172	U2173	U2174	U2175	U2176	U2177	U2178											
G2059	G2060	G2061	G2062	G2063	G2064	G2065	G2066	G2067	G2068	G2069	G2070	G2071	G2072	G2073	G2074	G2075	G2076	G2077	G2078	G2079	G2080	G2081	G2082	G2083	G2084	G2085	G2086	G2087	G2088	G2089	G2090	G2091	G2092	G2093	G2094	G2095	G2096	G2097	G2098	G2099	G2100	G2101	G2102	G2103	G2104	G2105	G2106	G2107	G2108	G2109	G2110	G2111	G2112	G2113	G2114	G2115	G2116	G2117	G2118											
U1998	U1999	U2000	U2001	U2002	U2003	U2004	U2005	U2006	U2007	U2008	U2009	U2010	U2011	U2012	U2013	U2014	U2015	U2016	U2017	U2018	U2019	U2020	U2021	U2022	U2023	U2024	U2025	U2026	U2027	U2028	U2029	U2030	U2031	U2032	U2033	U2034	U2035	U2036	U2037	U2038	U2039	U2040	U2041	U2042	U2043	U2044	U2045	U2046	U2047	U2048	U2049	U2050	U2051	U2052	U2053	U2054	U2055	U2056	U2057	U2058										
C1936	C1937	C1938	C1939	C1940	C1941	C1942	C1943	C1944	C1945	C1946	C1947	C1948	C1949	C1950	C1951	C1952	C1953	C1954	C1955	C1956	C1957	C1958	C1959	C1960	C1961	C1962	C1963	C1964	C1965	C1966	C1967	C1968	C1969	C1970	C1971	C1972	C1973	C1974	C1975	C1976	C1977	C1978	C1979	C1980	C1981	C1982	C1983	C1984	C1985	C1986	C1987	C1988	C1989	C1990	C1991	C1992	C1993	C1994	C1995											
U1865	U1866	U1867	U1868	U1869	U1870	U1871	U1872	U1873	U1874	U1875	U1876	U1877	U1878	U1879	U1880	U1881	U1882	U1883	U1884	U1885	U1886	U1887	U1888	U1889	U1890	U1891	U1892	U1893	U1894	U1895	U1896	U1897	U1898	U1899	U1900	U1901	U1902	U1903	U1904	U1905	U1906	U1907	U1908	U1909	U1910	U1911	U1912	U1913	U1914	U1915	U1916	U1917	U1918	U1919	U1920	U1921	U1922	U1923	U1924	U1925	U1926	U1927	U1928	U1929	U1930	U1931	U1932	U1933	U1934	
C1805	C1806	C1807	C1808	C1809	C1810	C1811	C1812	C1813	C1814	C1815	C1816	C1817	C1818	C1819	C1820	C1821	C1822	C1823	C1824	C1825	C1826	C1827	C1828	C1829	C1830	C1831	C1832	C1833	C1834	C1835	C1836	C1837	C1838	C1839	C1840	C1841	C1842	C1843	C1844	C1845	C1846	C1847	C1848	C1849	C1850	C1851	C1852	C1853	C1854	C1855	C1856	C1857	C1858	C1859	C1860	C1861	C1862	C1863	C1864											
U1734	U1735	U1736	U1737	U1738	U1739	U1740	U1741	U1742	U1743	U1744	U1745	U1746	U1747	U1748	U1749	U1750	U1751	U1752	U1753	U1754	U1755	U1756	U1757	U1758	U1759	U1760	U1761	U1762	U1763	U1764	U1765	U1766	U1767	U1768	U1769	U1770	U1771	U1772	U1773	U1774	U1775	U1776	U1777	U1778	U1779	U1780	U1781	U1782	U1783	U1784	U1785	U1786	U1787	U1788	U1789	U1790	U1791	U1792	U1793	U1794	U1795	U1796	U1797	U1798	U1799	U1800	U1801	U1802	U1803	U1804

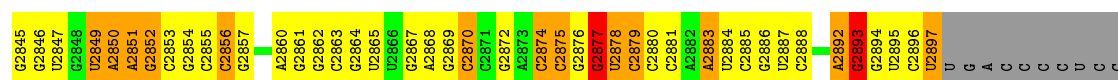


• Molecule 1: 23S Ribosomal RNA



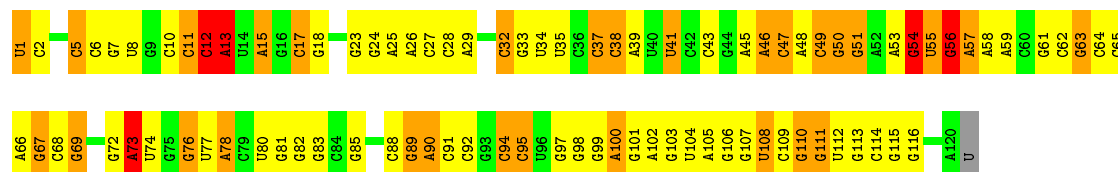
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G1674	G1606	G1538	G1459	C1386	G1318	C1250	A	C1110	G874	C904	U836	U767
C1675	A1608	G1539	A1460	C1387	C1320	C1251	C1178	G1111	C975	U905	U839	G769
G1678	A1609	G1540	G1461	C1388	A1321	A1253	C1179	A1112	C976	G906	U839	C772
U1679	A1610	G1541	G1462	C1389	A1322	A1254	G1181	U1113	C976	C908	C840	U773
G1680	C1611	C1542	C1463	U1390	U1323	U1255	A1182	U1114	C976	C909	A841	A774
G1681	C1612	C1543	C1464	U1391	G1324	C1256	C1185	G1115	A981	A910	G842	G775
G1682	G1613	A1544	G1465	U1394	C1327	C1257	C1186	C1117	A981	A911	G843	G776
G1683	A1614	A1545	G1466	A1395	G1328	G1260	U1188	G1118	A981	A912	G844	A777
G1684	C1615	C1546	C1467	U1396	U1329	U1261	A1189	C	A981	A913	G845	G778
C1685	A1616	C1547	G1468	U1397	C1330	U1262	G1190	C	A981	A914	G846	U779
C1686	C1617	C1548	A1469	U1397	A1331	G1263	G1191	C	A981	A915	G847	G780
U1687	A1618	G1470	G1471	U1396	G1332	U1264	G1195	C	A981	A916	G848	A781
U1688	G1619	A1471	A1472	U1397	C1333	G1265	C1196	C	A981	A917	G849	A782
A1689	G1620	G1473	G1473	U1397	C1334	U1266	C1197	C	A981	A918	G850	A783
U1692	G1621	A1477	A1477	U1397	U1335	U1267	U1198	G	A981	A919	G851	A784
U1693	G1622	A1478	A1478	U1397	A1336	A1268	U1199	C	A981	A920	G852	G785
C1694	C1623	G1479	G1479	U1401	G1337	A1269	C1200	A	A981	A921	G853	C786
U1695	G1624	G1479	G1479	C1402	G1338	C1270	C1201	A	A981	A922	G854	U787
G1696	G1625	G1480	G1480	C1403	G1339	A1271	C1202	A	A981	A923	G855	A788
G1697	G1626	U1481	U1481	C1404	U1340	A1272	C1203	U	A981	A924	G856	A789
G1698	G1627	G1482	G1482	U1405	U1341	A1273	G1204	U	A981	A925	G857	C790
A1698	G1628	G1482	G1482	U1406	A1342	A1274	U1205	A	A981	A926	G858	G791
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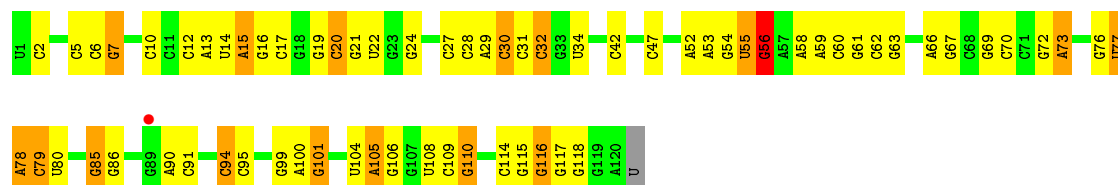
• Molecule 2: 5S Ribosomal RNA

Chain AB: 25% 46% 24% . .



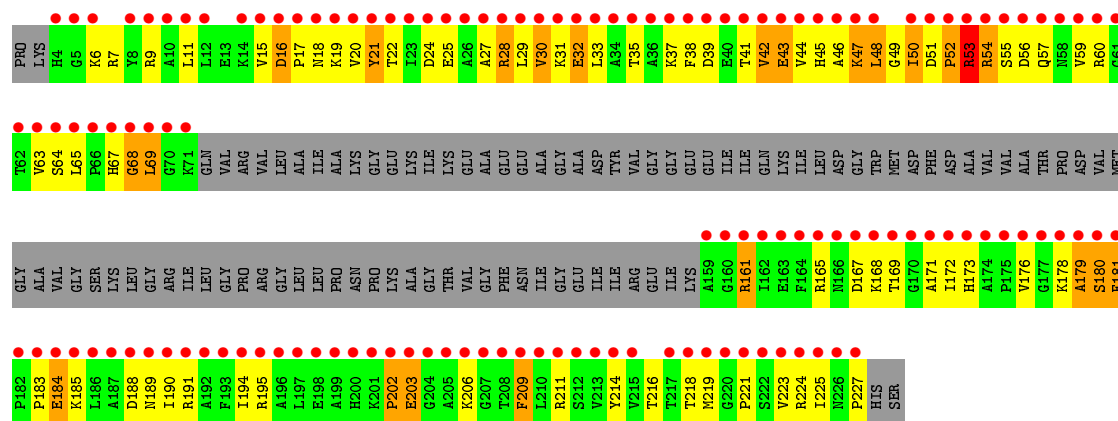
• Molecule 2: 5S Ribosomal RNA

Chain CB: 44% 41% 13% . .



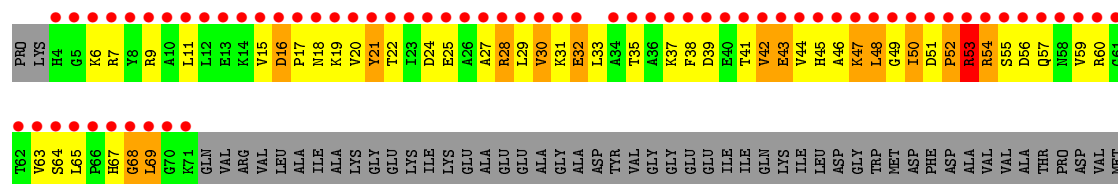
• Molecule 3: 50S ribosomal protein L1

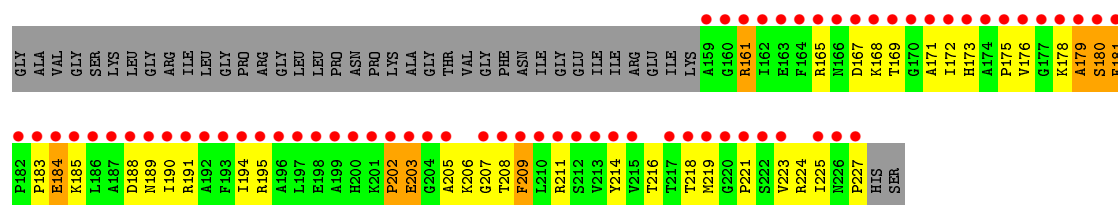
Chain AC: 22% 58% 28% 10% 40%



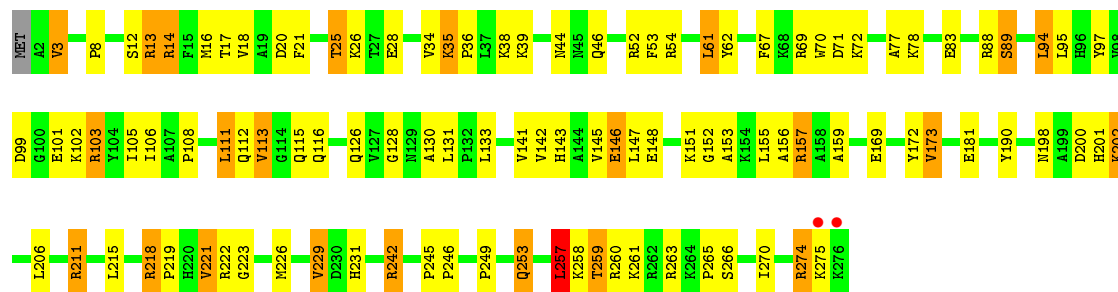
• Molecule 3: 50S ribosomal protein L1

Chain CC: 21% 58% 29% 10% 40%

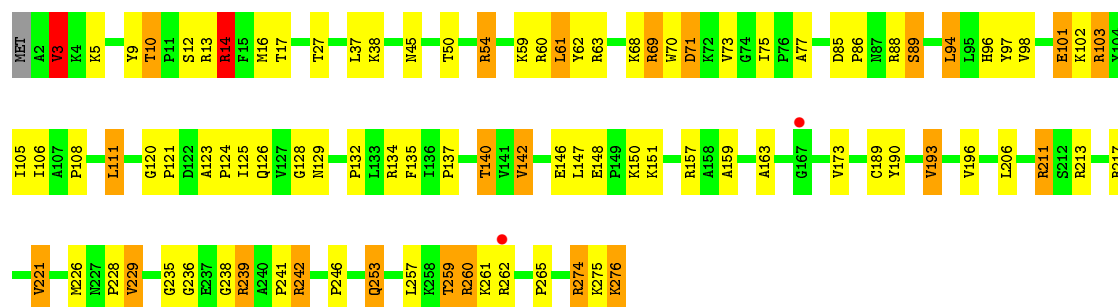




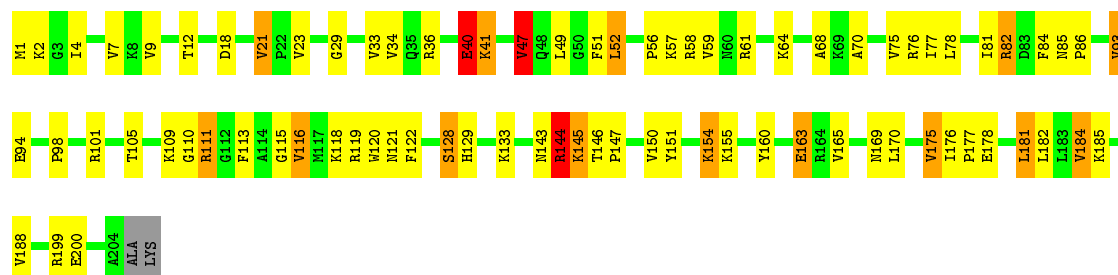
• Molecule 4: 50S ribosomal protein L2



• Molecule 4: 50S ribosomal protein L2

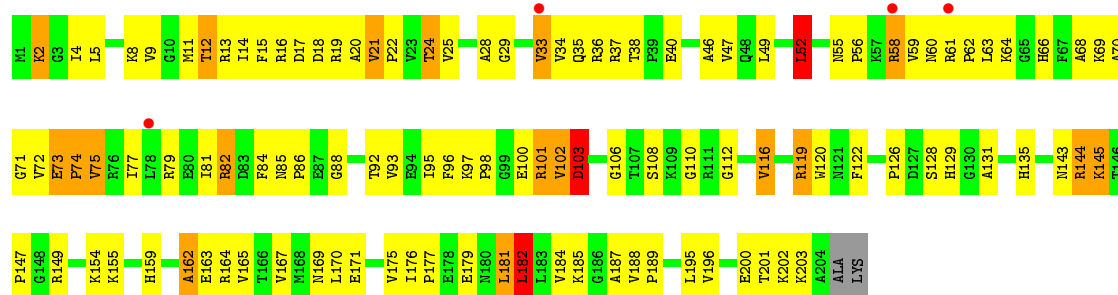


• Molecule 5: 50S ribosomal protein L3



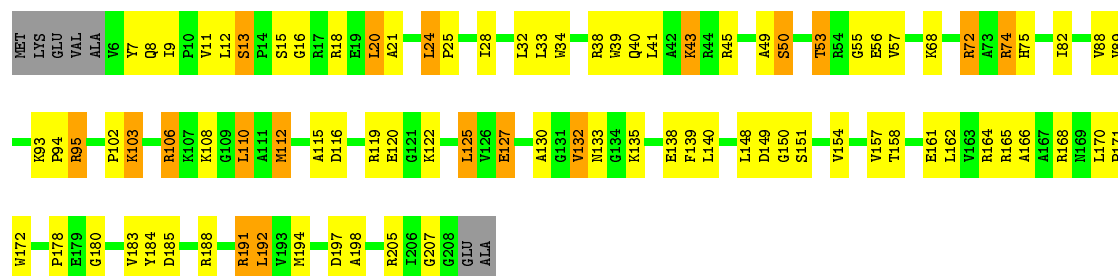
• Molecule 5: 50S ribosomal protein L3





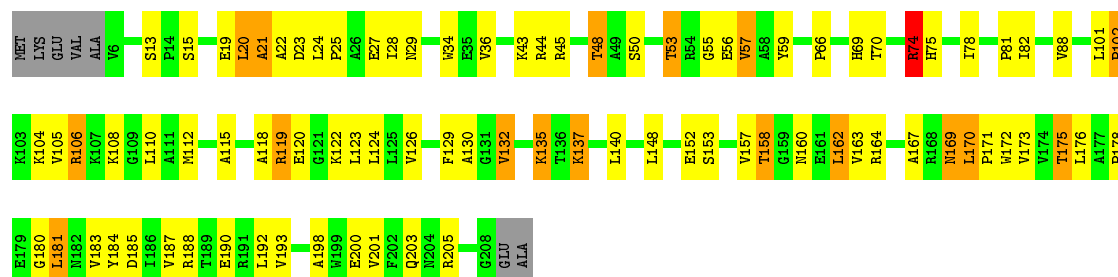
- Molecule 6: 50S ribosomal protein L4

Chain AF: 55% 33% 9% .



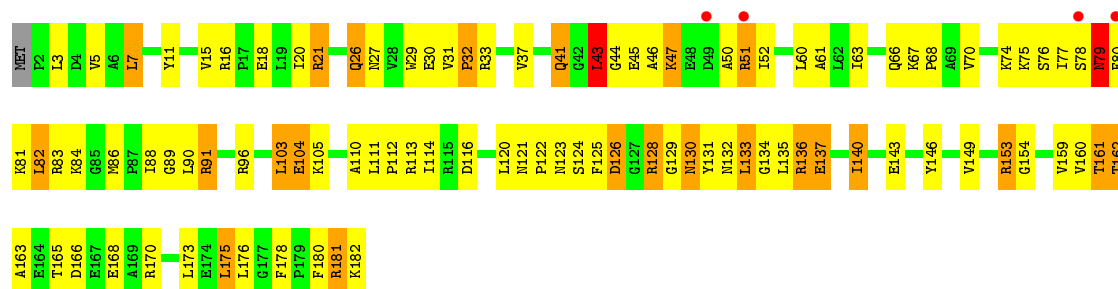
- Molecule 6: 50S ribosomal protein L4

Chain CF: 55% 33% 8% .

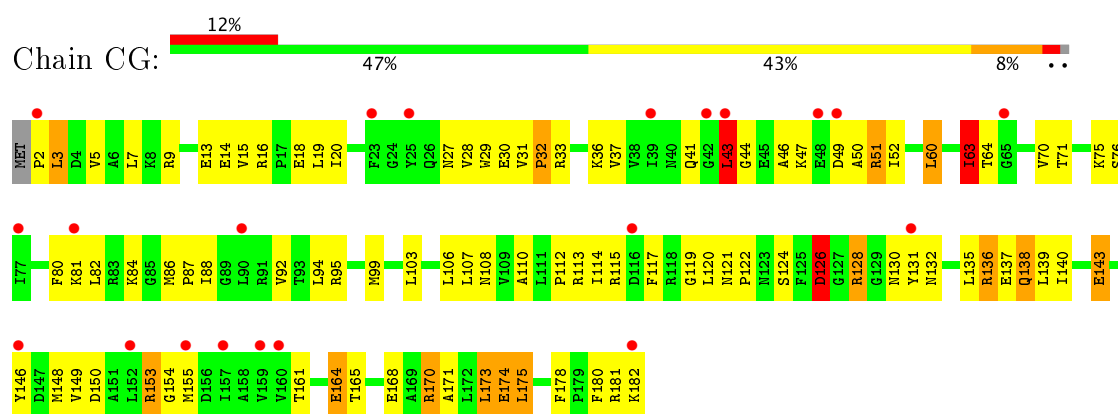


- Molecule 7: 50S ribosomal protein L5

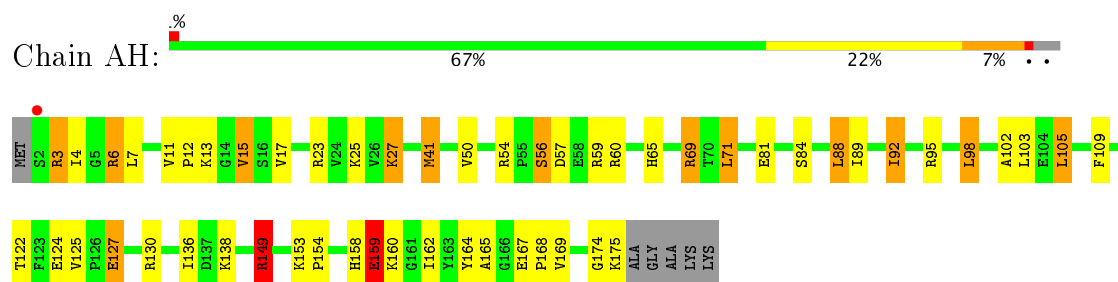
Chain AG: 2% 46% 40% 13% ..



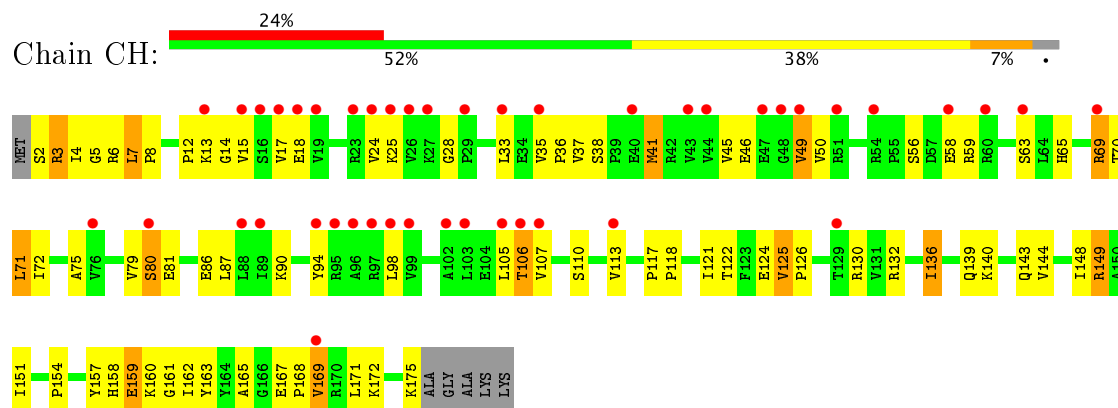
- Molecule 7: 50S ribosomal protein L5



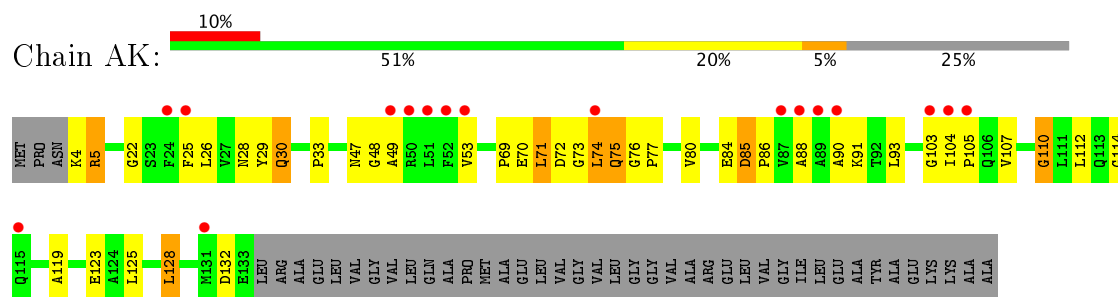
• Molecule 8: 50S ribosomal protein L6



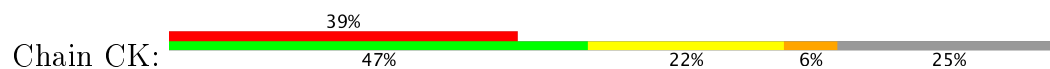
• Molecule 8: 50S ribosomal protein L6

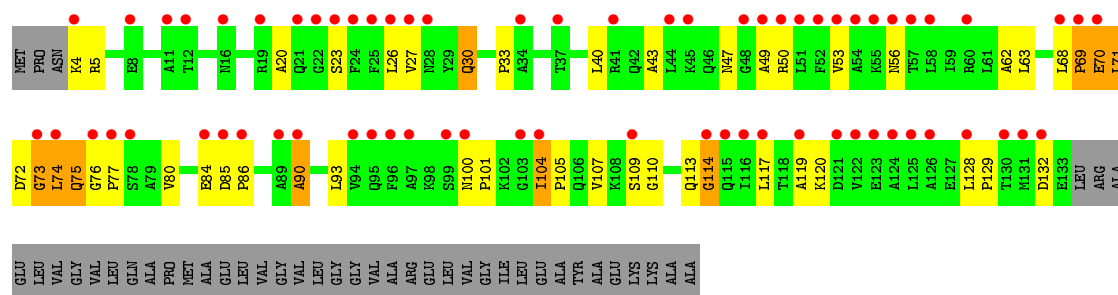


• Molecule 9: 50S ribosomal protein L10

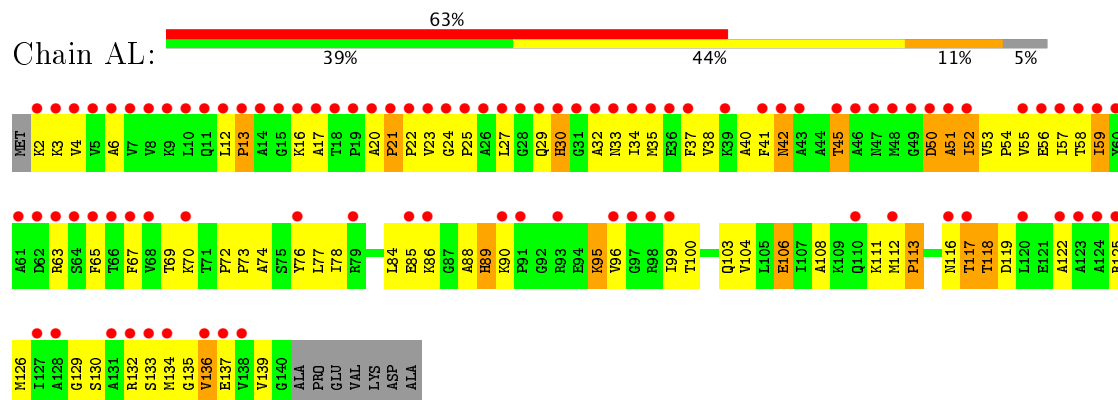


• Molecule 9: 50S ribosomal protein L10

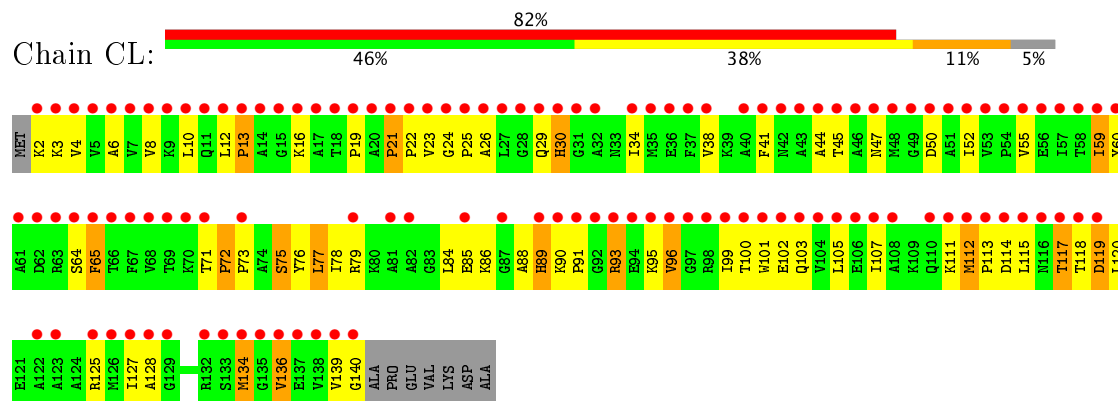




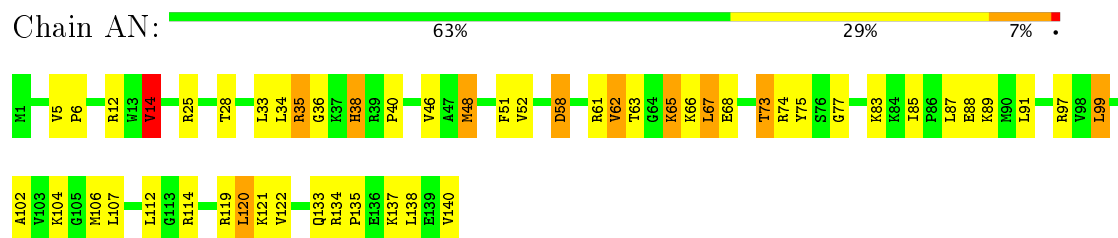
• Molecule 10: 50S ribosomal protein L11



• Molecule 10: 50S ribosomal protein L11

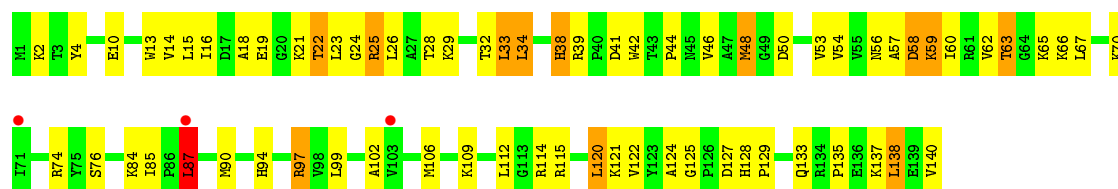


• Molecule 11: 50S ribosomal protein L13

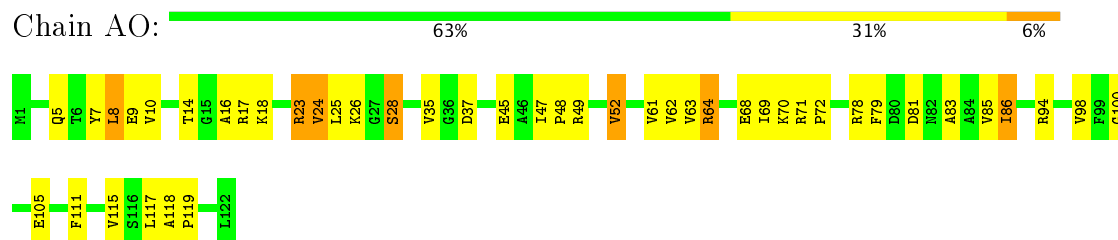


• Molecule 11: 50S ribosomal protein L13

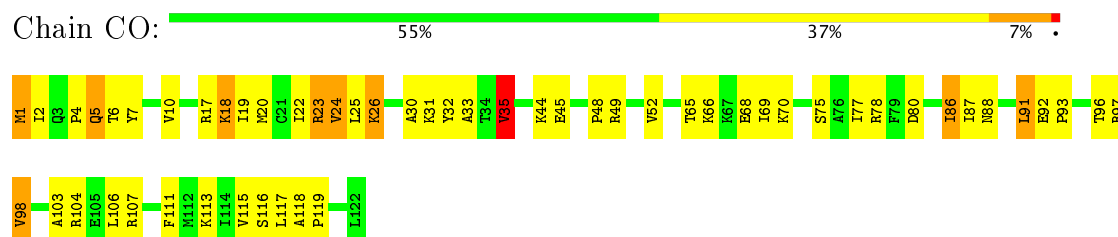




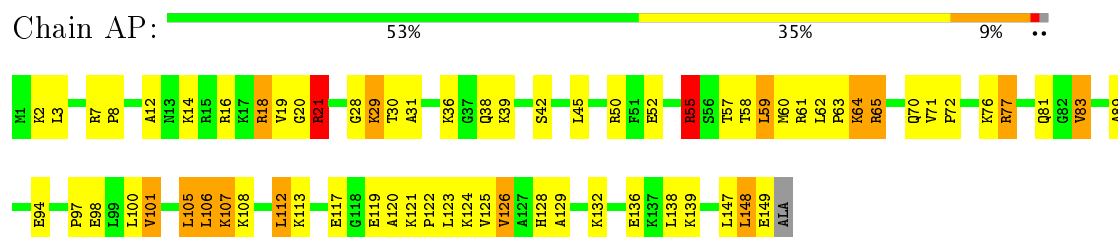
• Molecule 12: 50S ribosomal protein L14



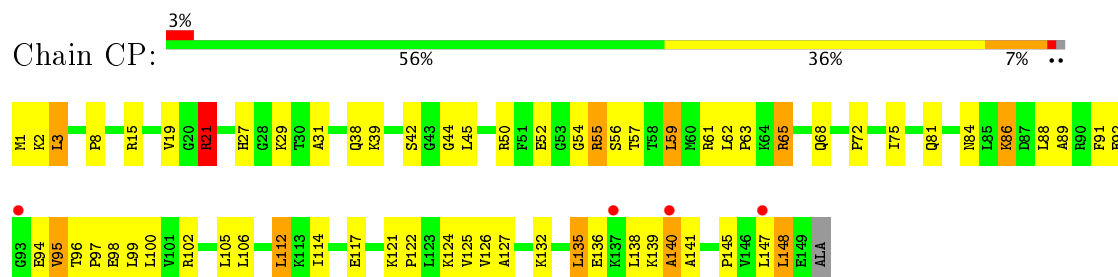
• Molecule 12: 50S ribosomal protein L14



• Molecule 13: 50S ribosomal protein L15

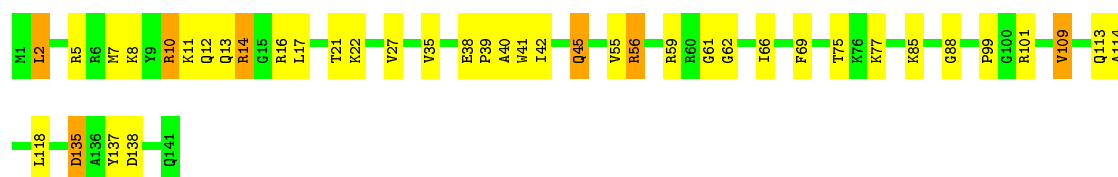


• Molecule 13: 50S ribosomal protein L15

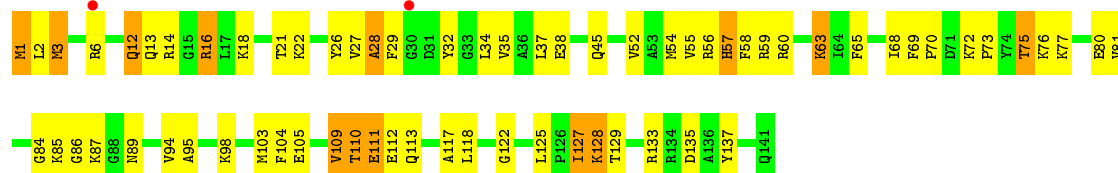


• Molecule 14: 50S ribosomal protein L16

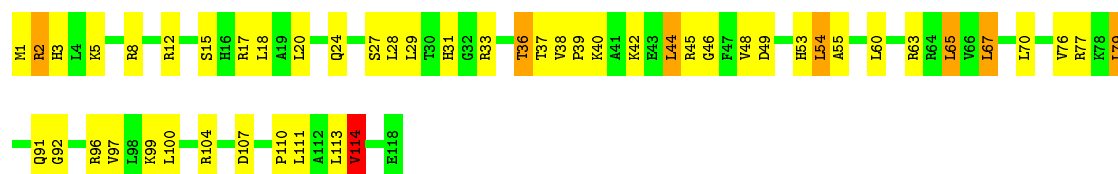




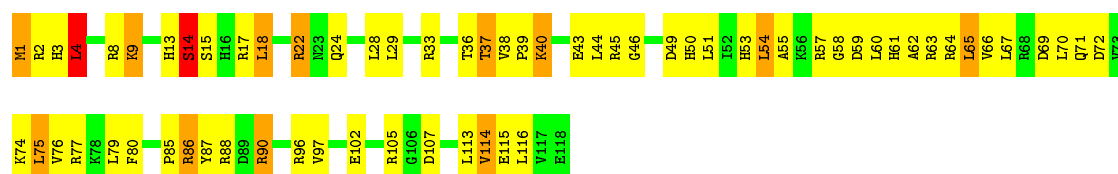
- Molecule 14: 50S ribosomal protein L16



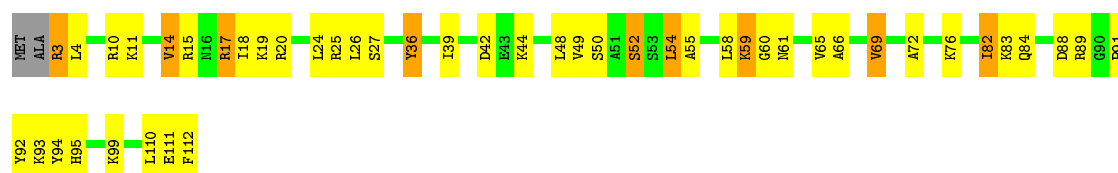
- Molecule 15: 50S ribosomal protein L17



- Molecule 15: 50S ribosomal protein L17

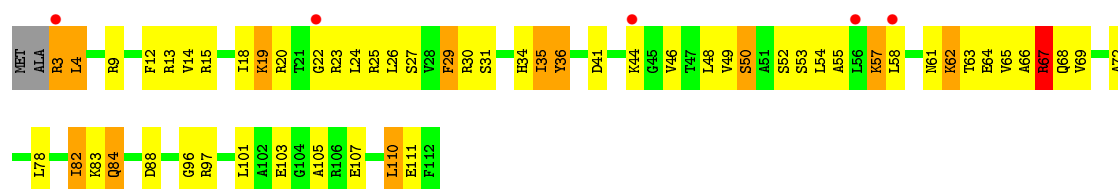


- Molecule 16: 50S ribosomal protein L18

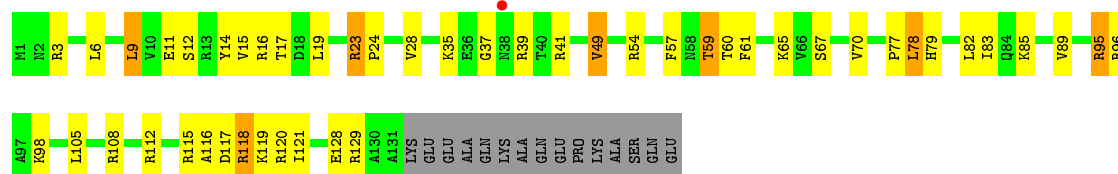


- Molecule 16: 50S ribosomal protein L18

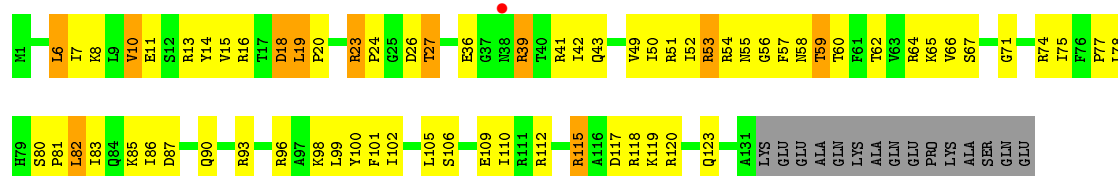
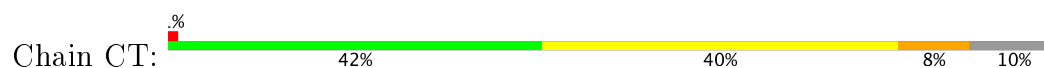




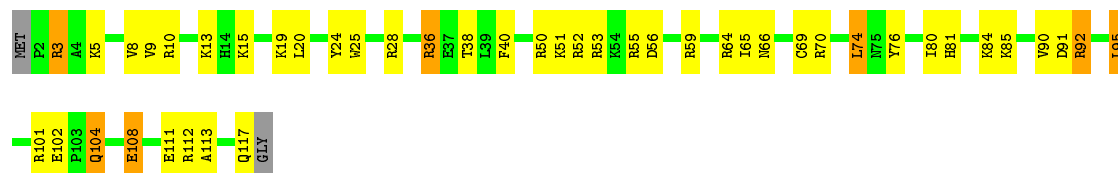
- Molecule 17: 50S ribosomal protein L19



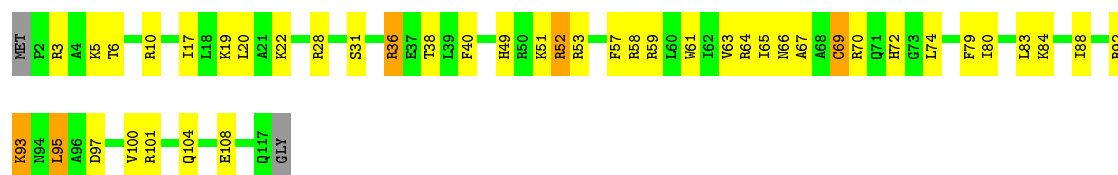
- Molecule 17: 50S ribosomal protein L19



- Molecule 18: 50S ribosomal protein L20

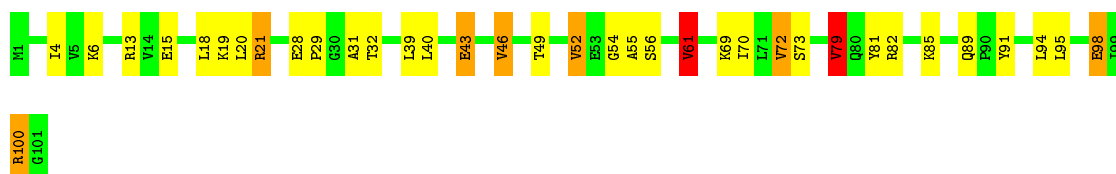


- Molecule 18: 50S ribosomal protein L20

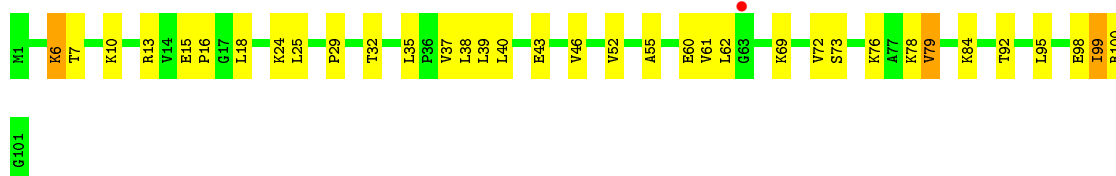


- Molecule 19: 50S ribosomal protein L21

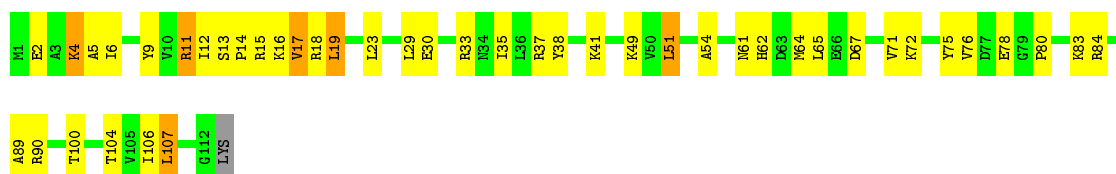




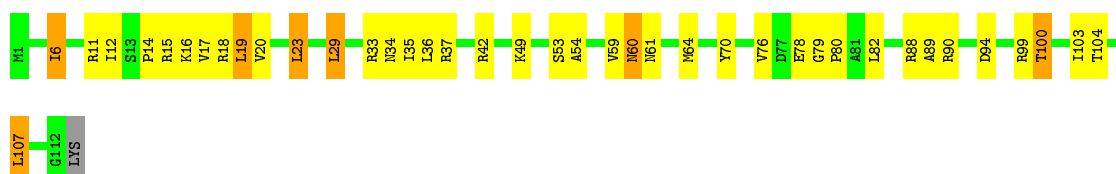
- Molecule 19: 50S ribosomal protein L21



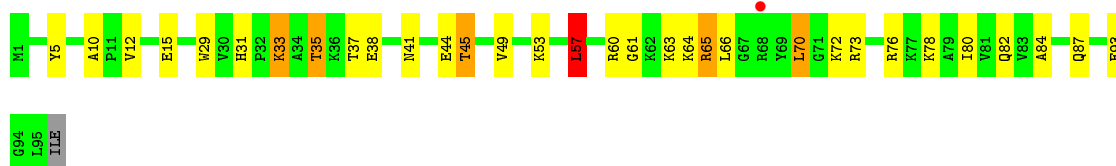
- Molecule 20: 50S ribosomal protein L22



- Molecule 20: 50S ribosomal protein L22



- Molecule 21: 50S ribosomal protein L23

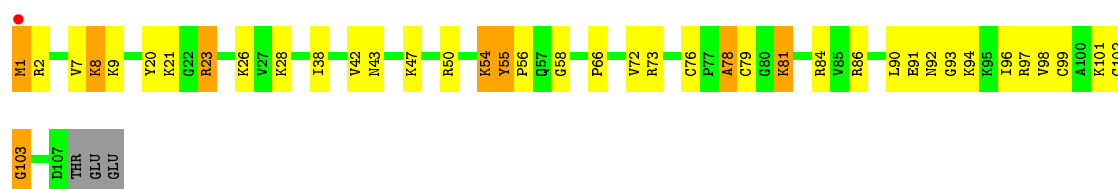


- Molecule 21: 50S ribosomal protein L23

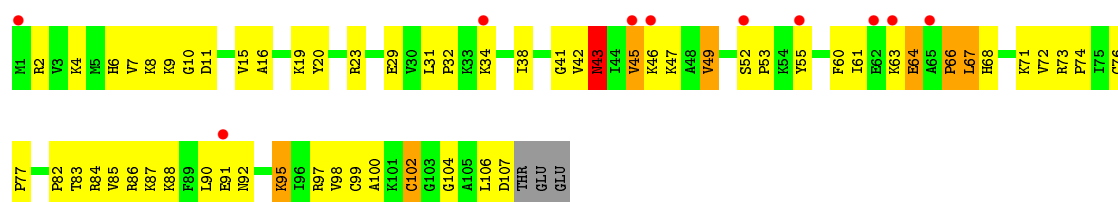




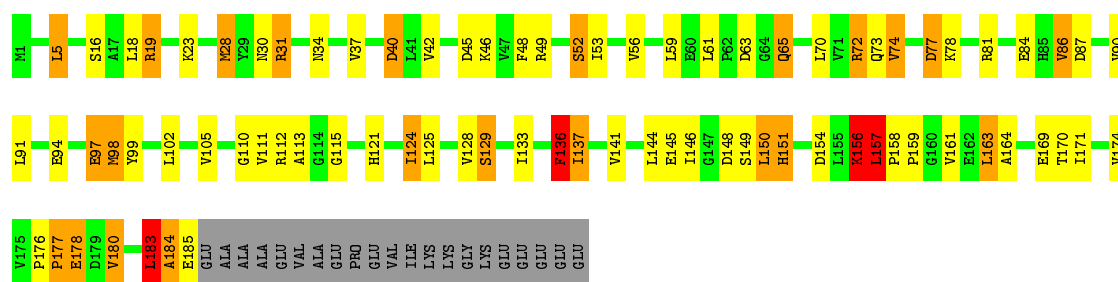
- Molecule 22: 50S ribosomal protein L24



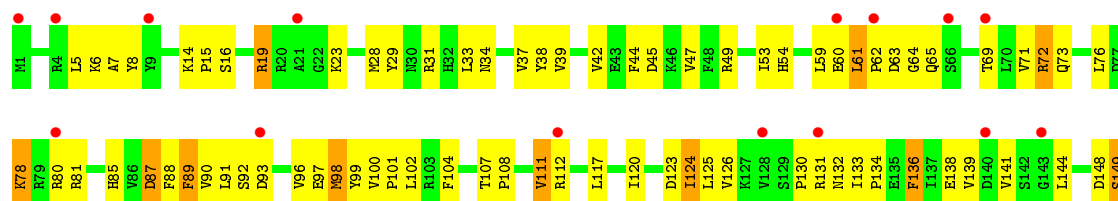
- Molecule 22: 50S ribosomal protein L24

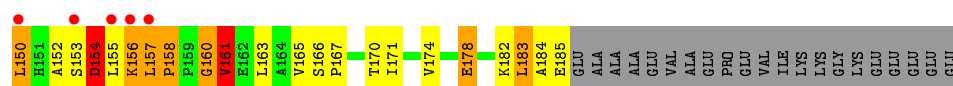


- Molecule 23: 50S ribosomal protein L25



- Molecule 23: 50S ribosomal protein L25





- Molecule 24: 50S ribosomal protein L27



- Molecule 24: 50S ribosomal protein L27



- Molecule 25: 50S ribosomal protein L28



- Molecule 25: 50S ribosomal protein L28



- Molecule 26: 50S ribosomal protein L29



- Molecule 26: 50S ribosomal protein L29



- Molecule 27: 50S ribosomal protein L30

Chain A3: 



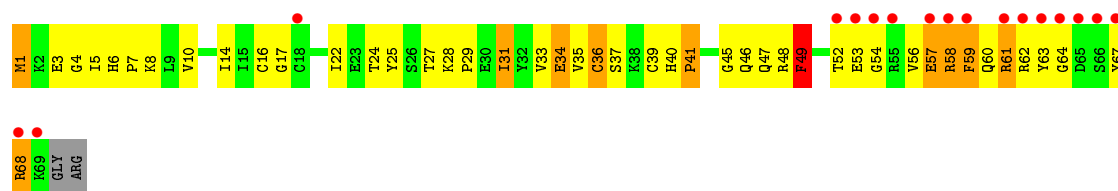
- Molecule 27: 50S ribosomal protein L30

Chain C3: 



- Molecule 28: 50S ribosomal protein L31

Chain A4: 



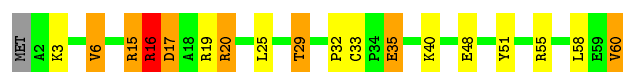
- Molecule 28: 50S ribosomal protein L31

Chain C4: 



- Molecule 29: 50S ribosomal protein L32

Chain A5: 



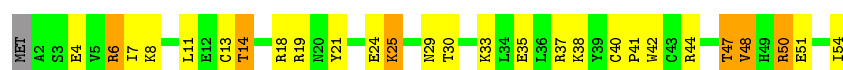
- Molecule 29: 50S ribosomal protein L32

Chain C5: 

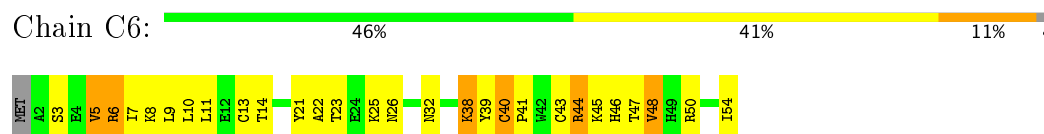


- Molecule 30: 50S ribosomal protein L33

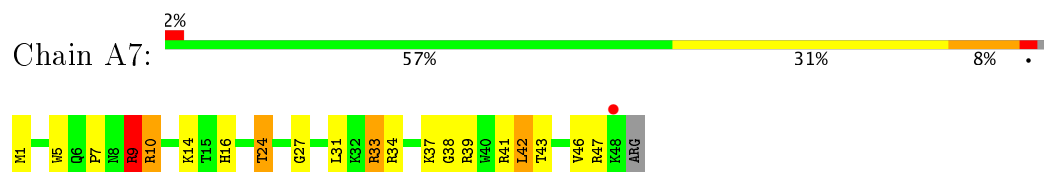
Chain A6: 



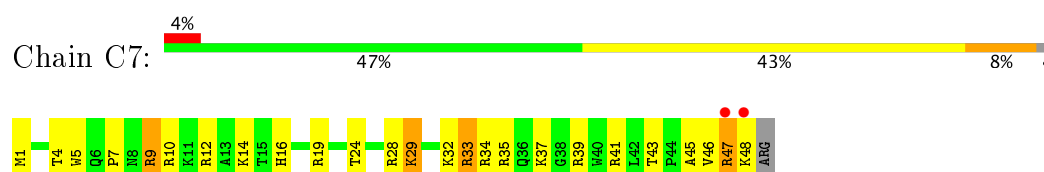
- Molecule 30: 50S ribosomal protein L33



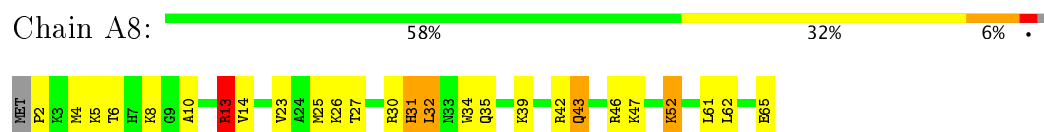
- Molecule 31: 50S ribosomal protein L34



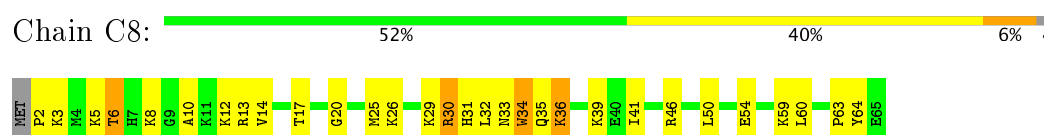
- Molecule 31: 50S ribosomal protein L34



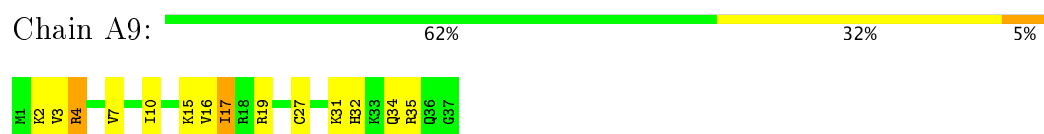
- Molecule 32: 50S ribosomal protein L35



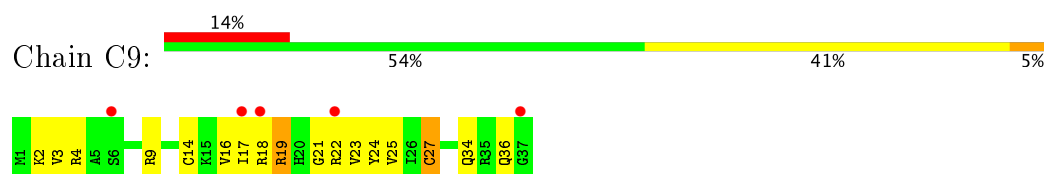
- Molecule 32: 50S ribosomal protein L35



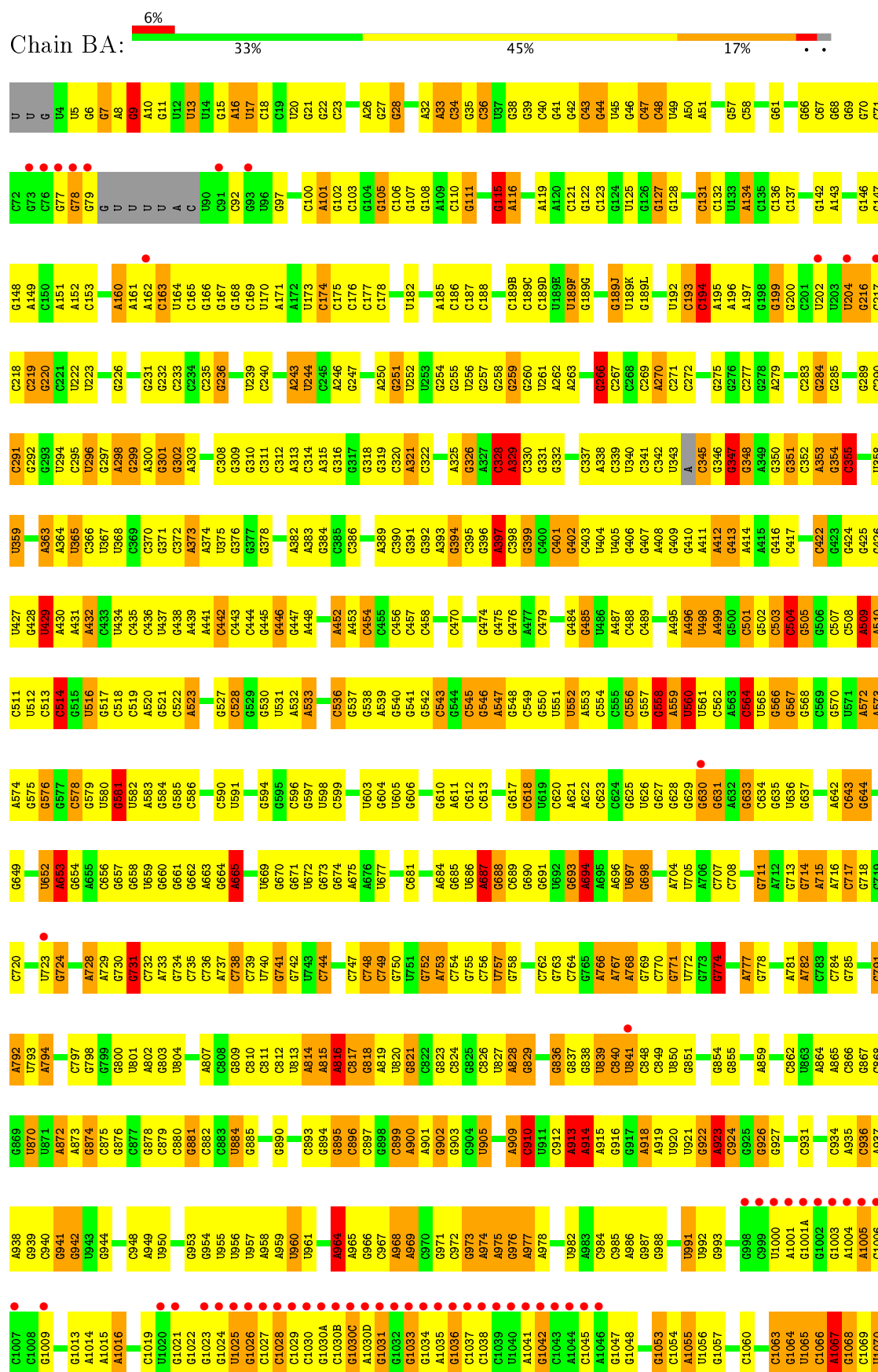
- Molecule 33: 50S ribosomal protein L36

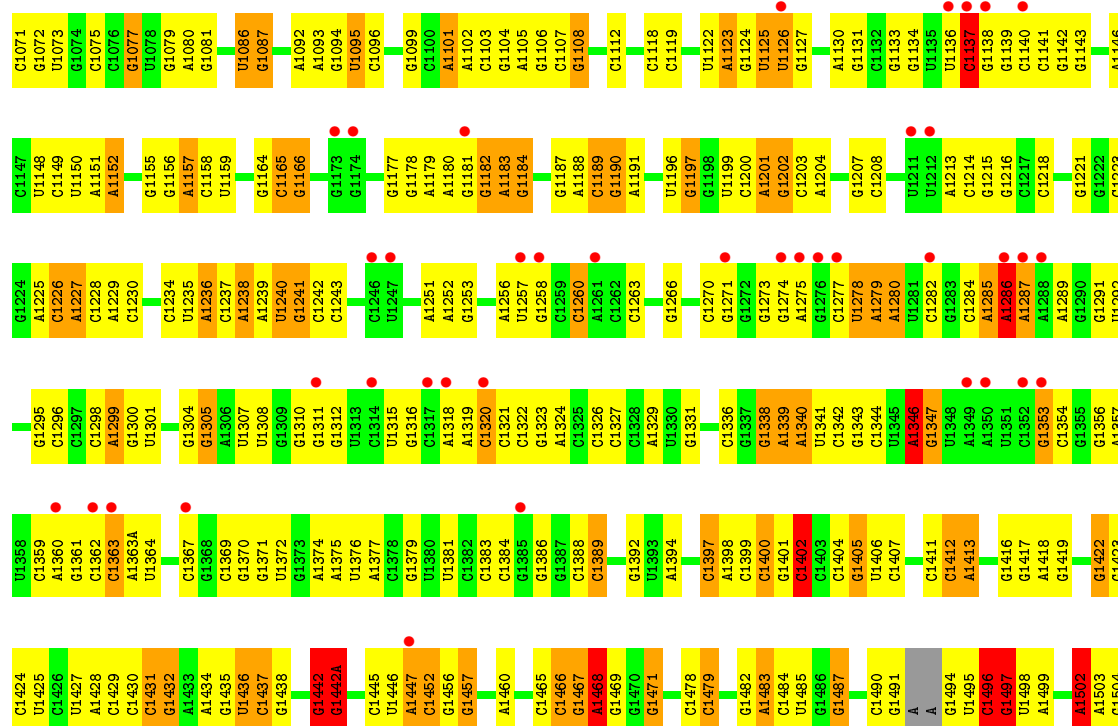


- Molecule 33: 50S ribosomal protein L36

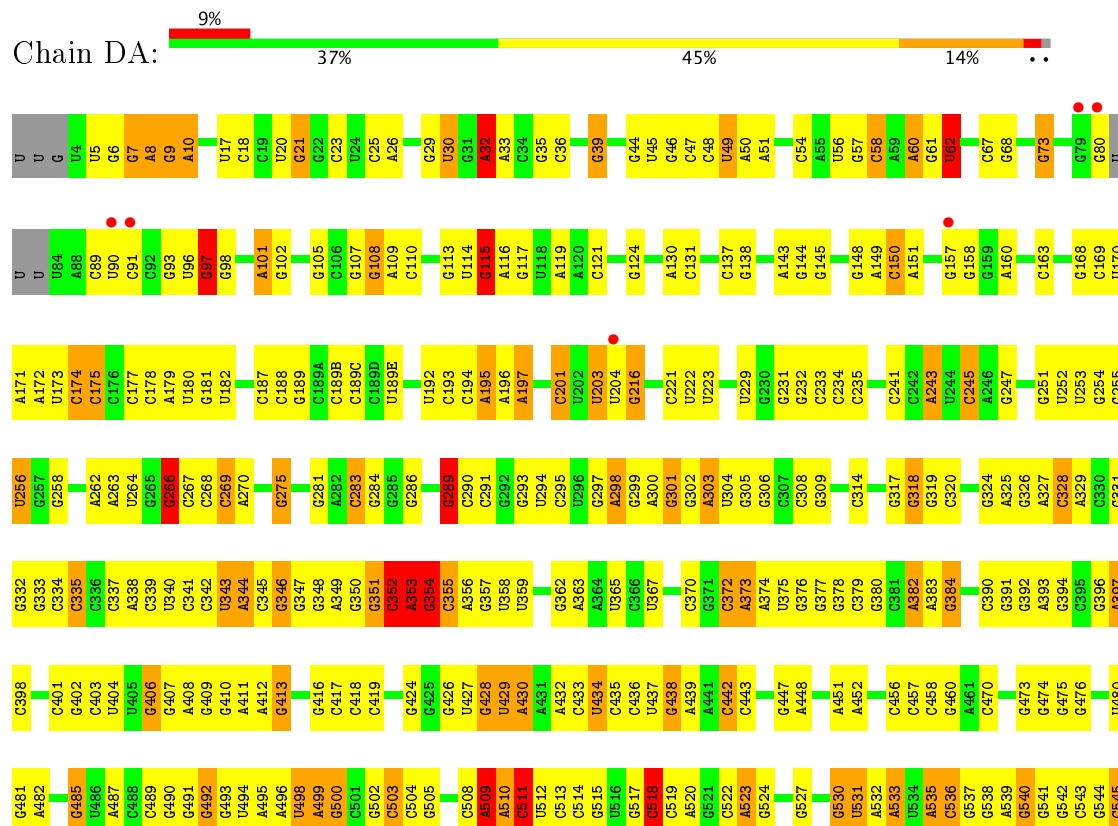


- Molecule 34: 16S Ribosomal RNA



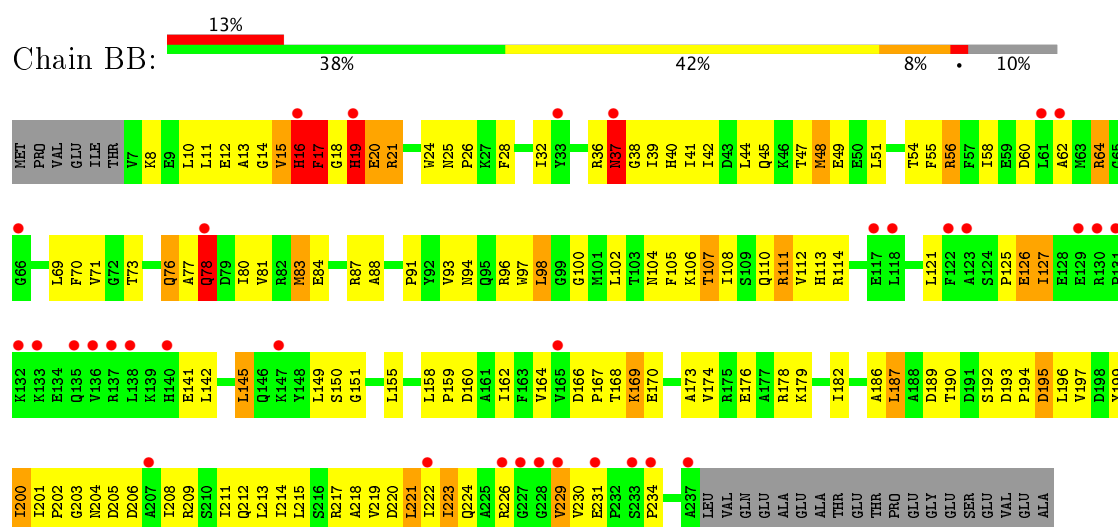


- Molecule 34: 16S Ribosomal RNA

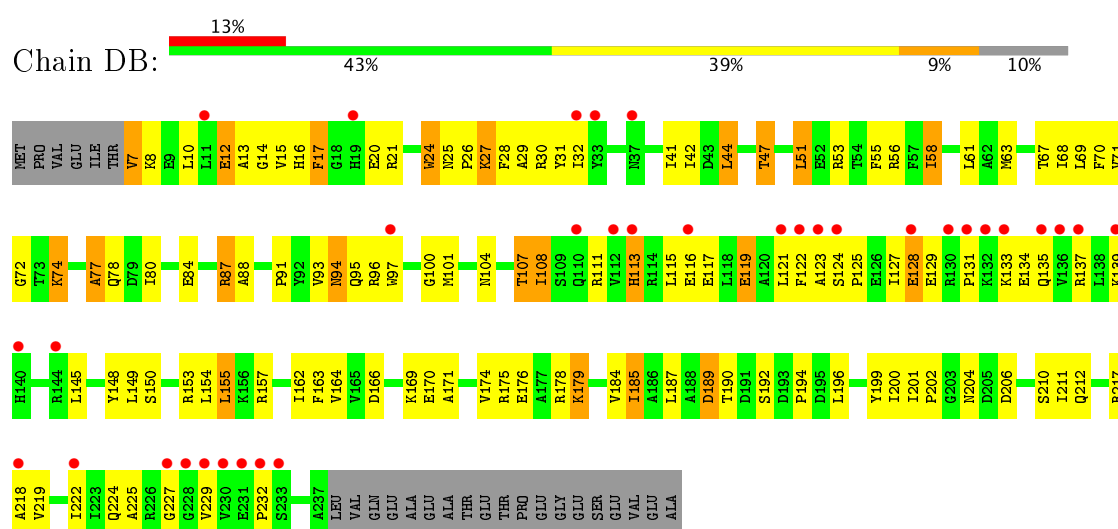


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U	A1468	C1383	C1317	G1254	G1127	G1063	A1003	A937	C866	G785	A707	G625	G548
C	G1469	G1384	A1319	G1255	C1128	G1064	G1004	A938	G867	G786	C708	A938	G550
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		G1386	C1321	G1258	G1131	C1066	C1006	G940	A873	U788	G710	G630	U552
		G1387	C1322	C1259	C1132	A1067	C1007	G941	A874	U789	G713	G631	
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		U1393	C1324	G1197	G1134	G1071	G1010	U943	C876	U792	A716	G633	C556
		A1394	C1325	G1198	U1135	G1072	G1011		C877	A793	C717	C634	G557
			G1326	A1201	U1136	U1073	U1012	U952	C878	A794	G718	G638	G558
			C1327	C1202	C1137	G1074	G1013	G953	C879	A795	C719	G639	A559
			C1328	C1203	G1138	C1075	A1014	G954	C880	C796	C720	A640	U560
				A1204	G1139	G1076	A1015	U955	G881	C797	G721	U641	U561
				U1205	C1140	U1077	A1016	U956	C882	G798	A722	A642	C562
				A1268	C1141	G1078	G1017		C883	C799	U723		A563
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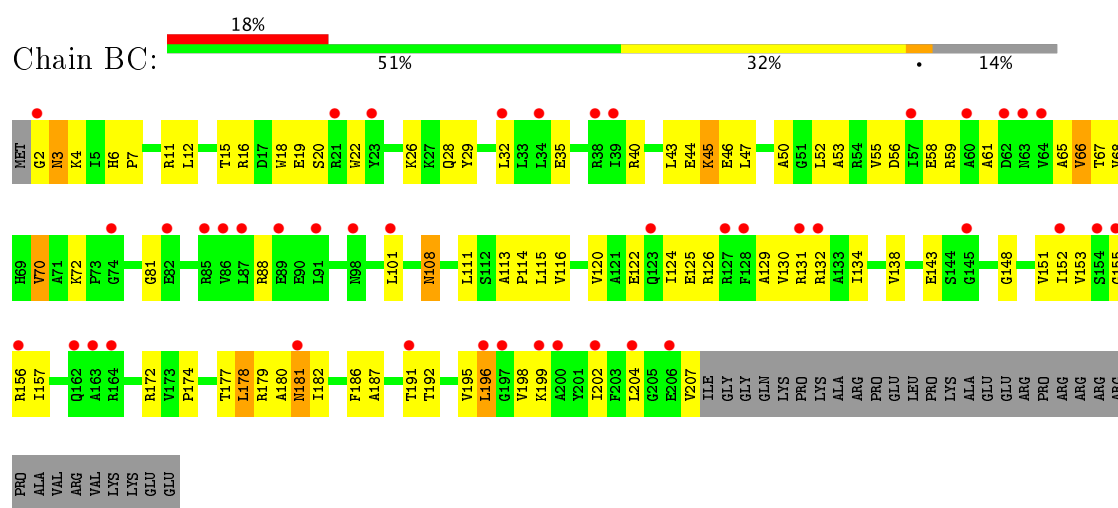
• Molecule 35: 30S ribosomal protein S2



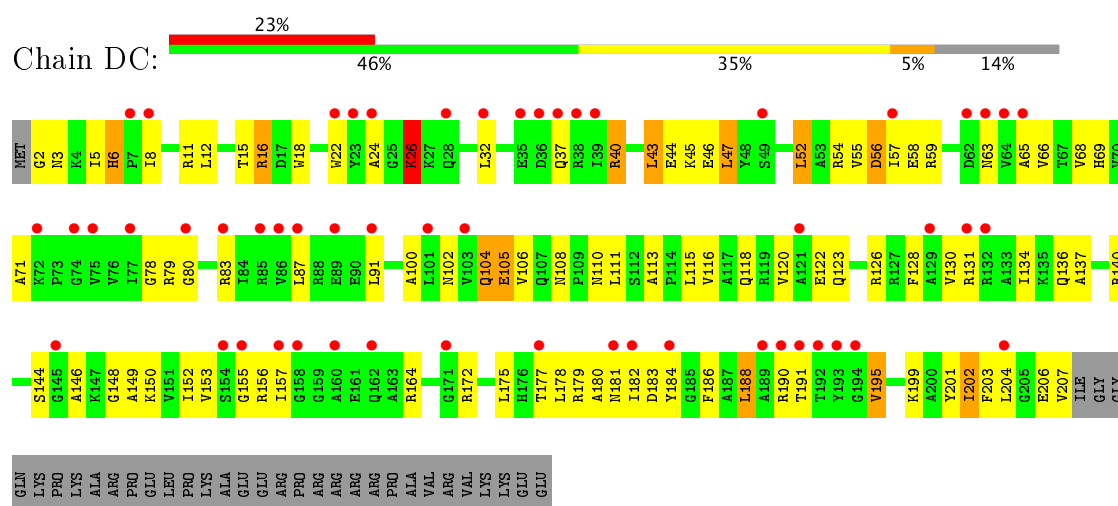
• Molecule 35: 30S ribosomal protein S2



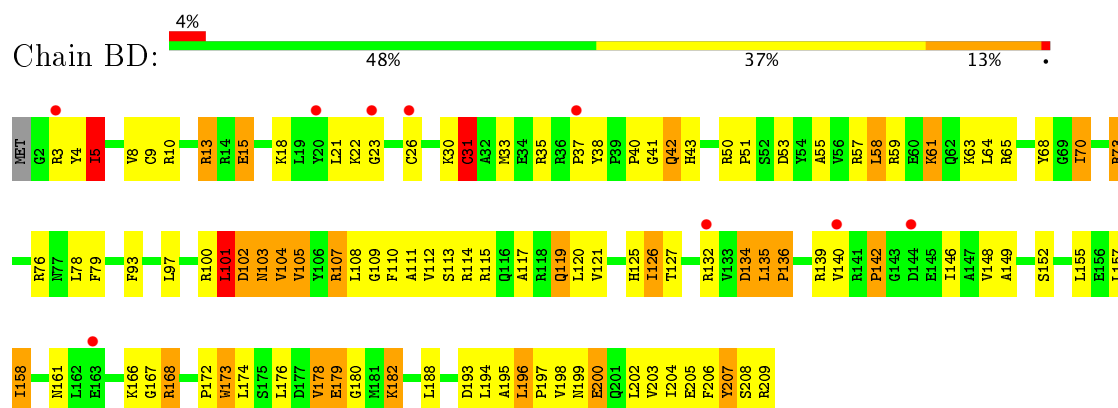
• Molecule 36: 30S ribosomal protein S3



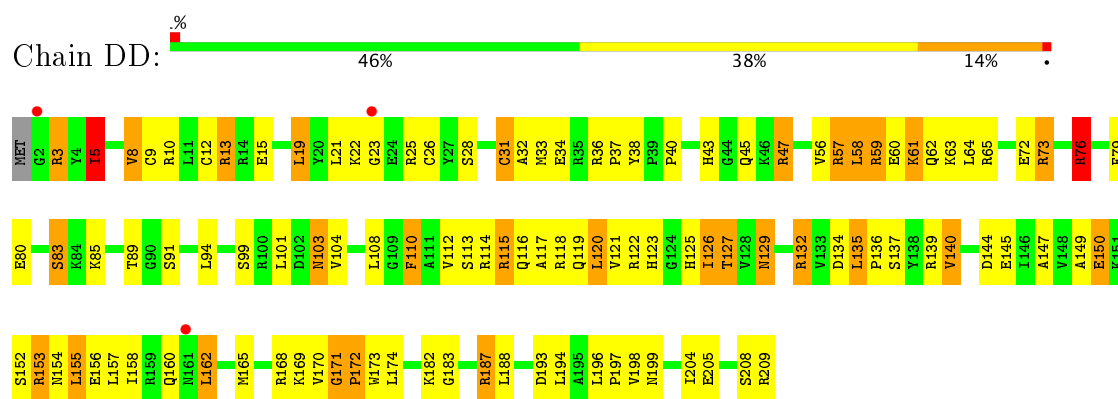
• Molecule 36: 30S ribosomal protein S3



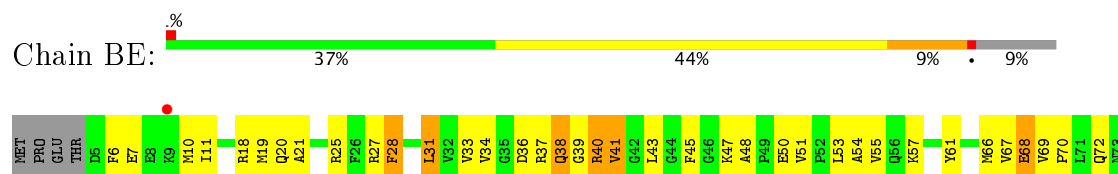
• Molecule 37: 30S ribosomal protein S4

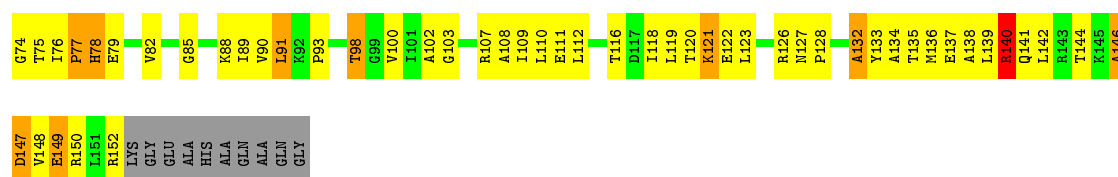


• Molecule 37: 30S ribosomal protein S4

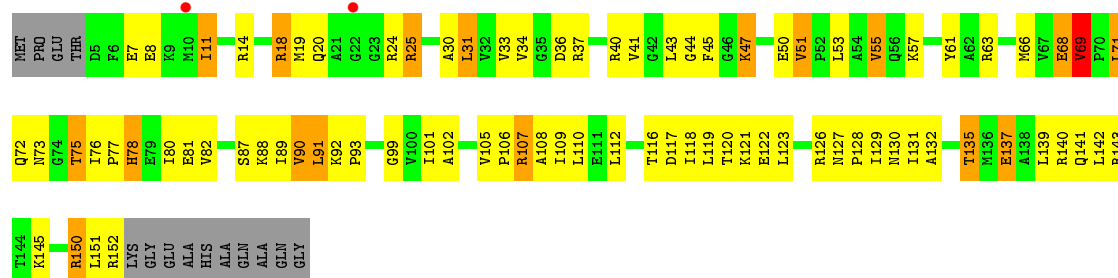


• Molecule 38: 30S ribosomal protein S5

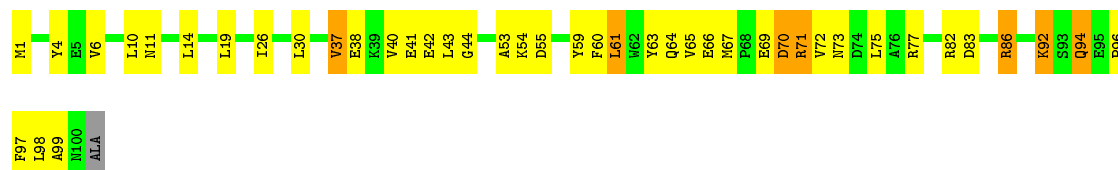




• Molecule 38: 30S ribosomal protein S5



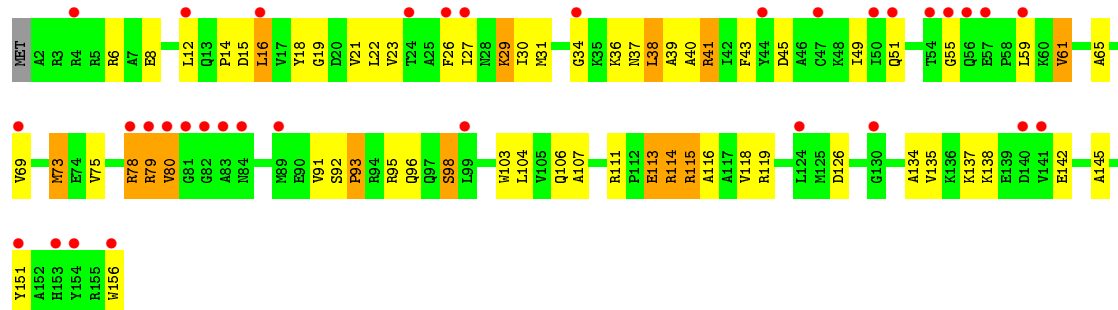
• Molecule 39: 30S ribosomal protein S6



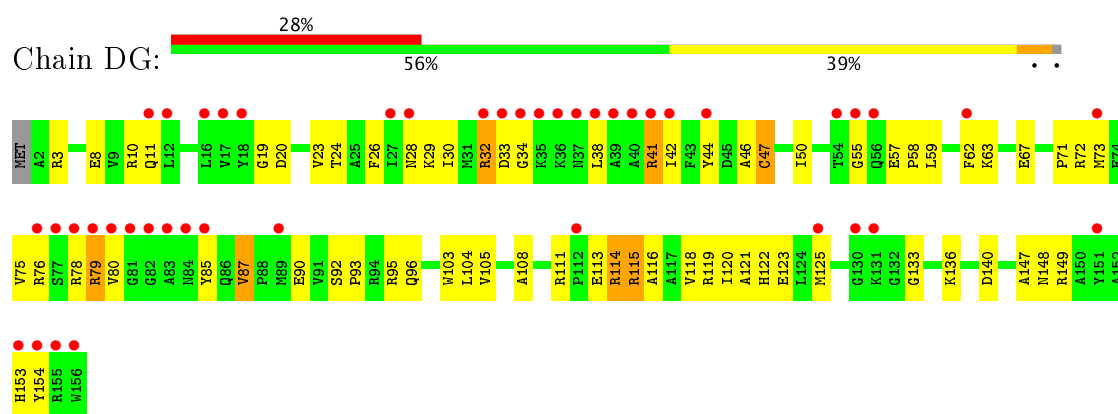
• Molecule 39: 30S ribosomal protein S6



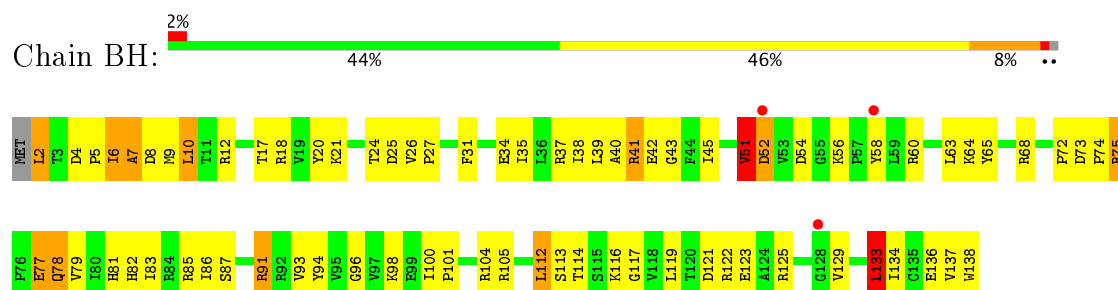
• Molecule 40: 30S ribosomal protein S7



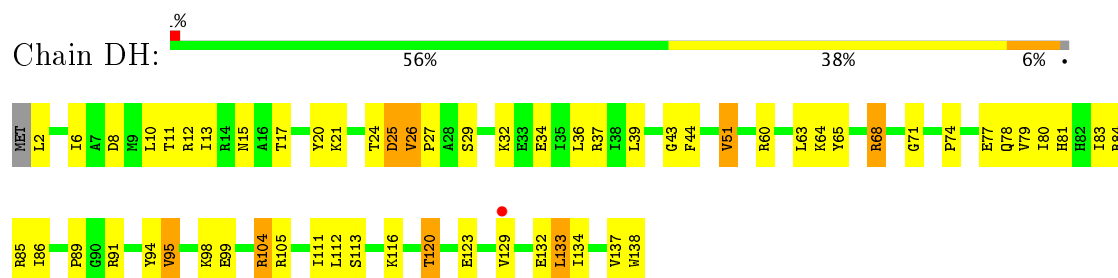
• Molecule 40: 30S ribosomal protein S7



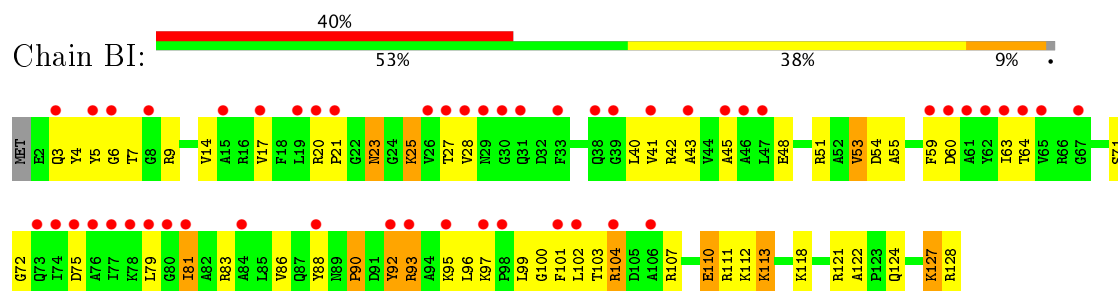
- Molecule 41: 30S ribosomal protein S8



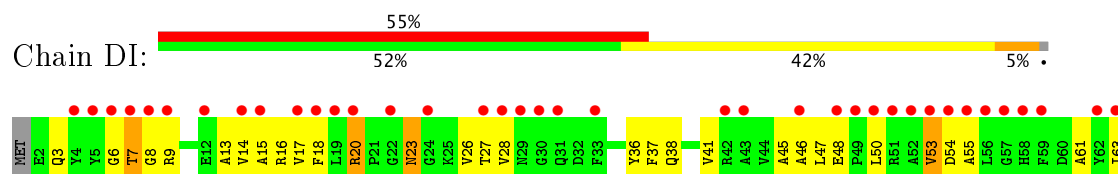
- Molecule 41: 30S ribosomal protein S8

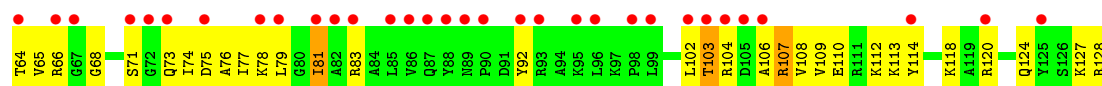


- Molecule 42: 30S ribosomal protein S9

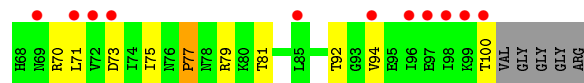
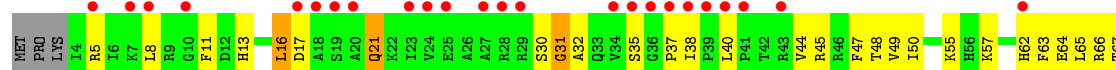


- Molecule 42: 30S ribosomal protein S9

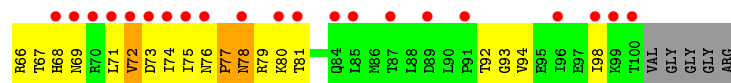
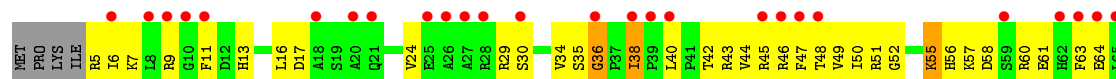




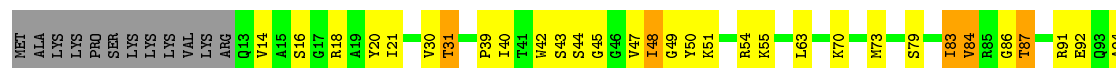
• Molecule 43: 30S ribosomal protein S10



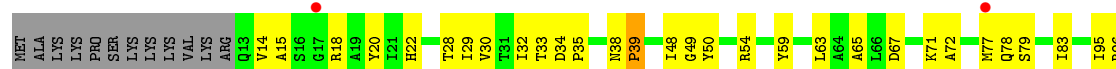
• Molecule 43: 30S ribosomal protein S10



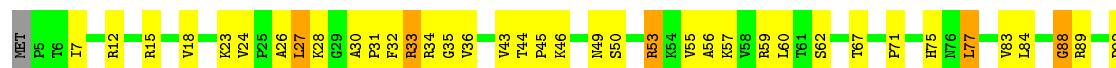
• Molecule 44: 30S ribosomal protein S11



• Molecule 44: 30S ribosomal protein S11

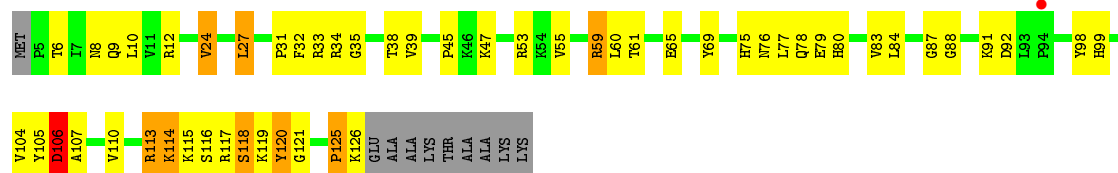


• Molecule 45: 30S ribosomal protein S12

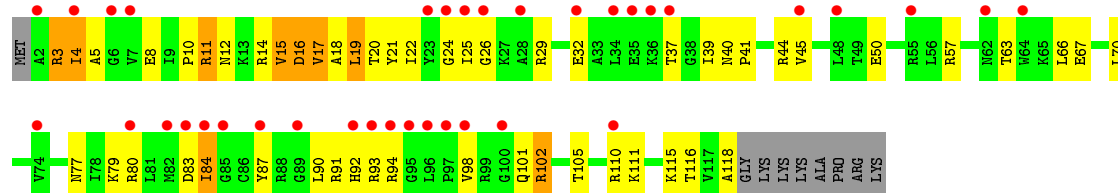




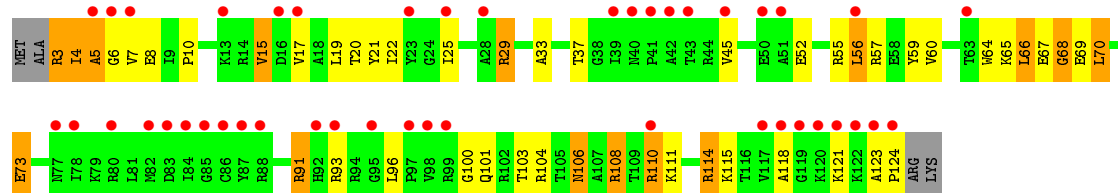
- Molecule 45: 30S ribosomal protein S12



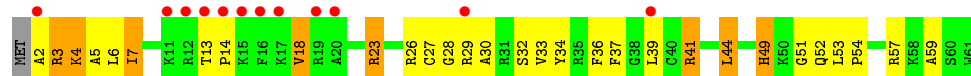
- Molecule 46: 30S ribosomal protein S13



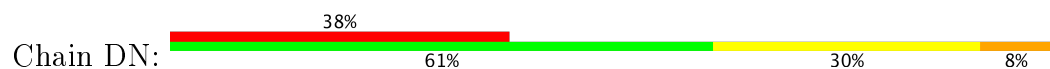
- Molecule 46: 30S ribosomal protein S13



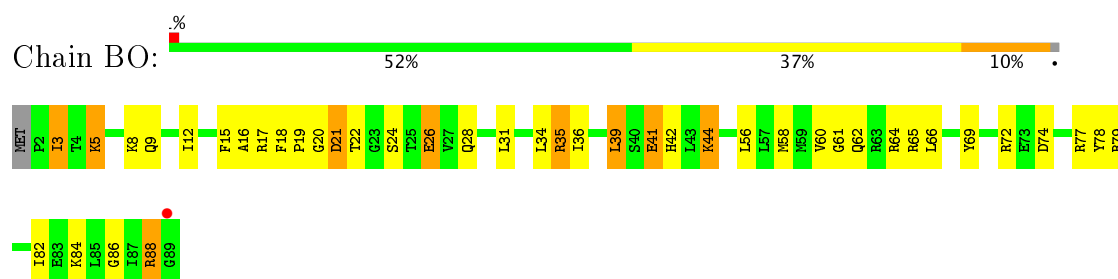
- Molecule 47: 30S ribosomal protein S14 type Z



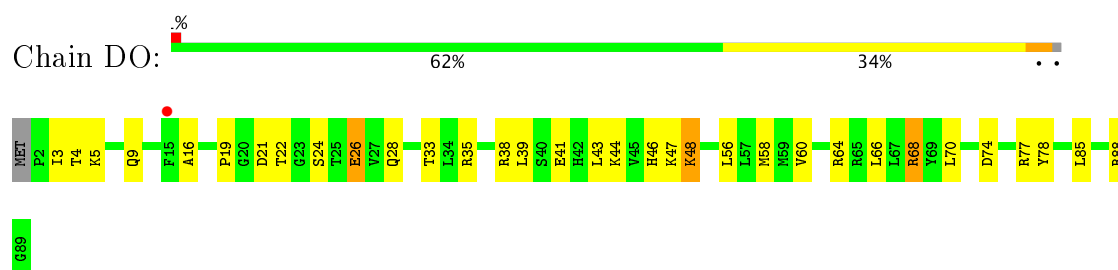
- Molecule 47: 30S ribosomal protein S14 type Z



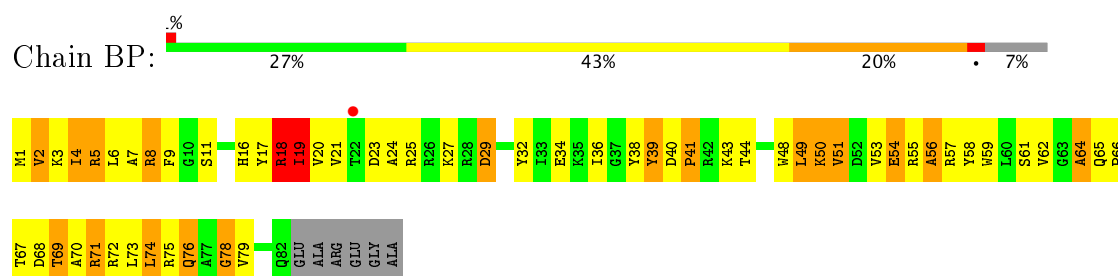
- Molecule 48: 30S ribosomal protein S15



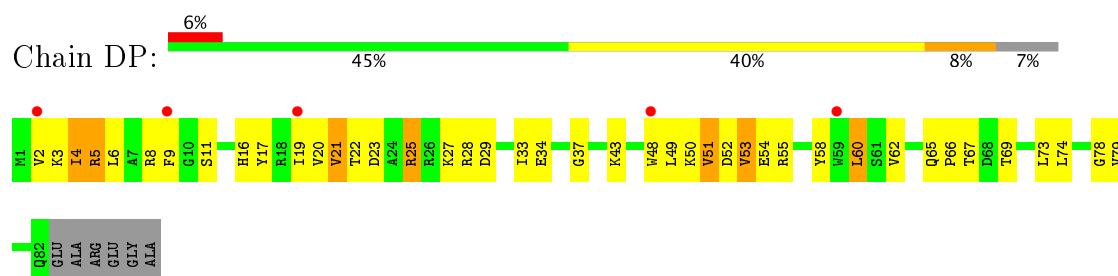
- Molecule 48: 30S ribosomal protein S15



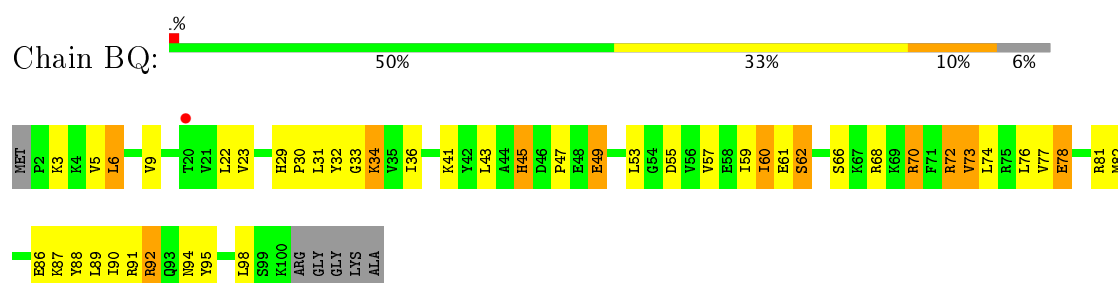
- Molecule 49: 30S ribosomal protein S16



- Molecule 49: 30S ribosomal protein S16

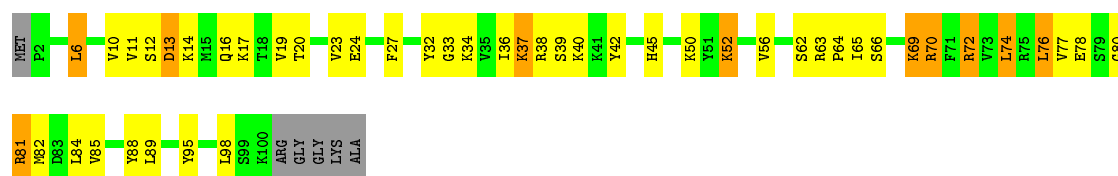


- Molecule 50: 30S ribosomal protein S17



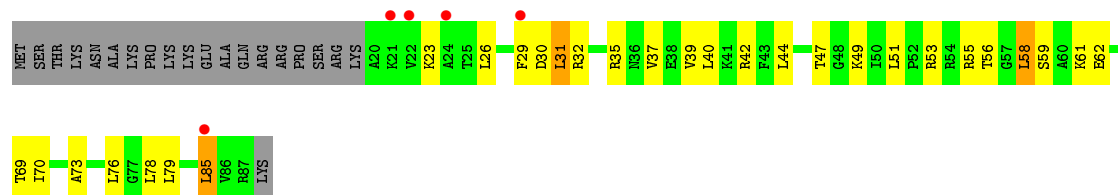
- Molecule 50: 30S ribosomal protein S17

Chain DQ: 



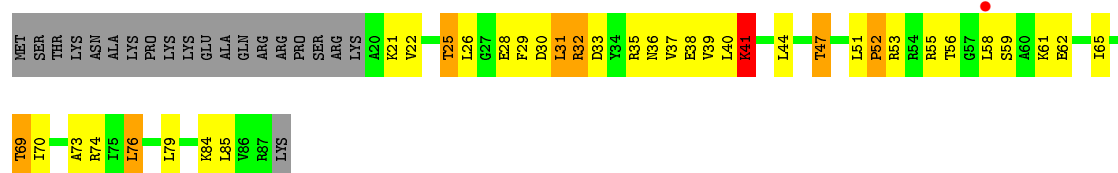
- Molecule 51: 30S ribosomal protein S18

Chain BR: 



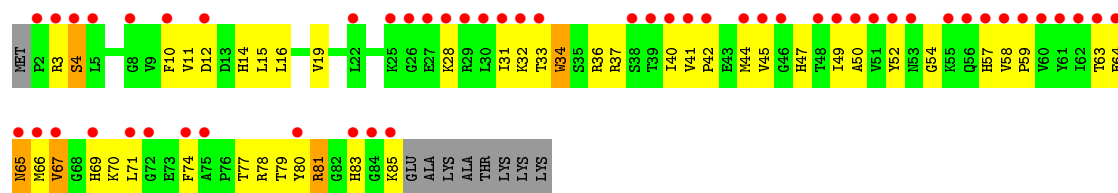
- Molecule 51: 30S ribosomal protein S18

Chain DR: 



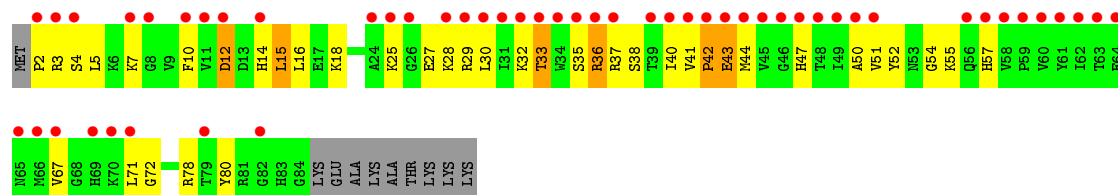
- Molecule 52: 30S ribosomal protein S19

Chain BS: 

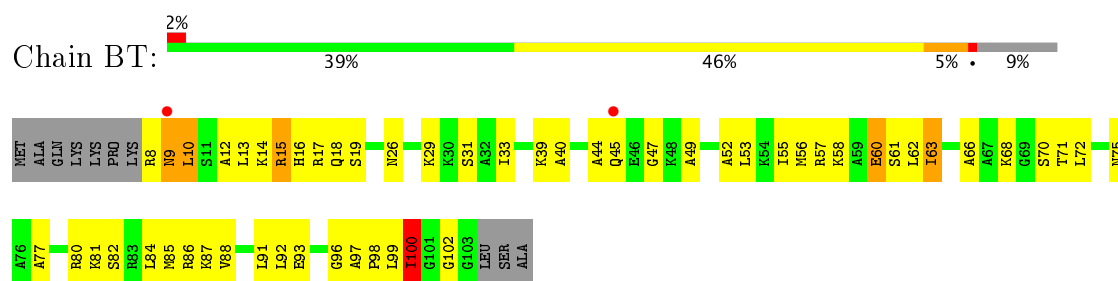


- Molecule 52: 30S ribosomal protein S19

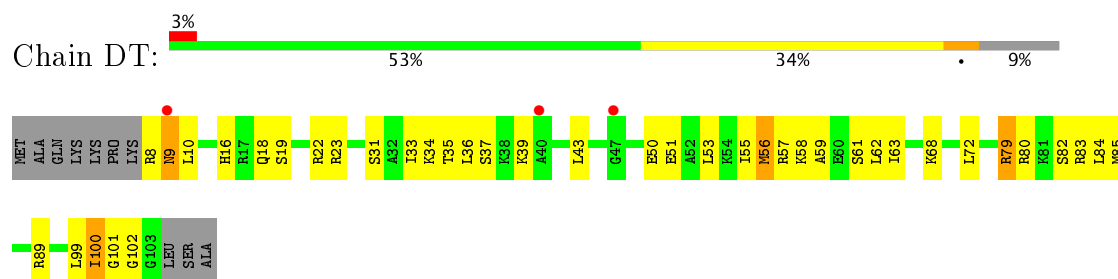
Chain DS: 



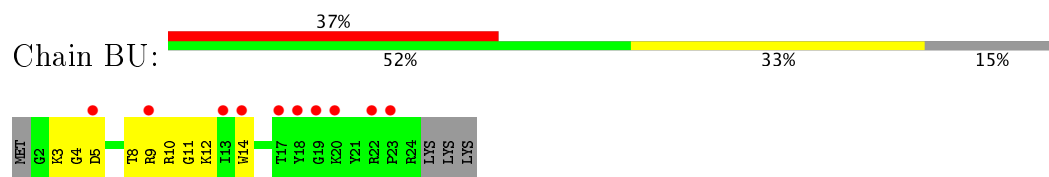
- Molecule 53: 30S ribosomal protein S20



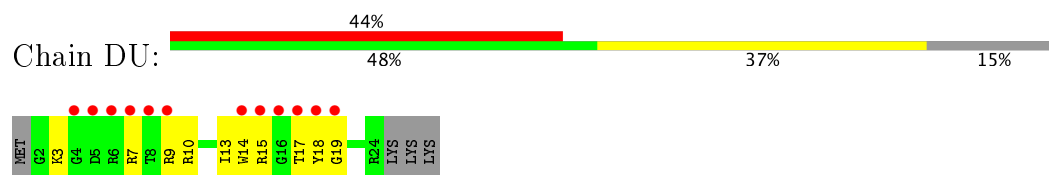
- Molecule 53: 30S ribosomal protein S20



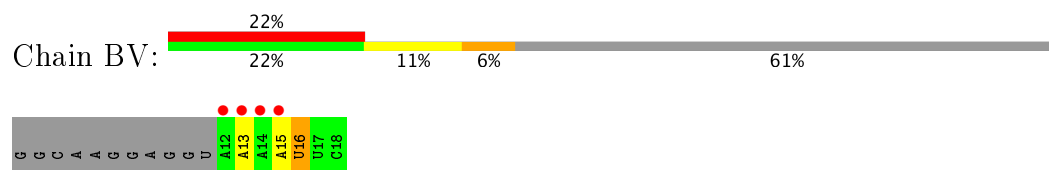
- Molecule 54: 30S ribosomal protein Thx



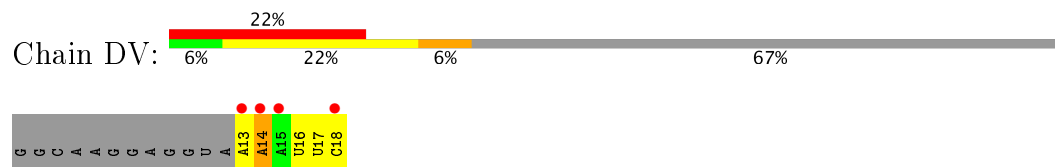
- Molecule 54: 30S ribosomal protein Thx



- Molecule 55: mRNA

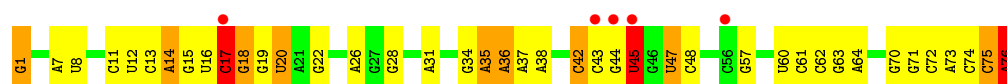


- Molecule 55: mRNA

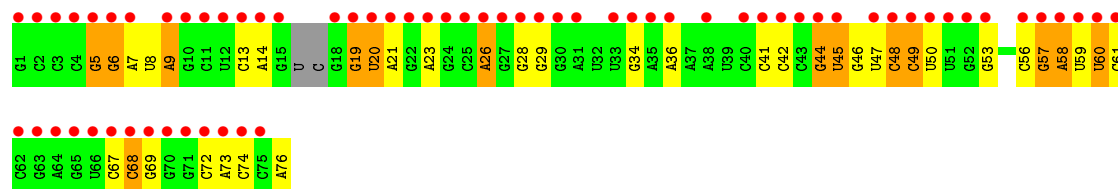


- Molecule 56: P-site tRNA

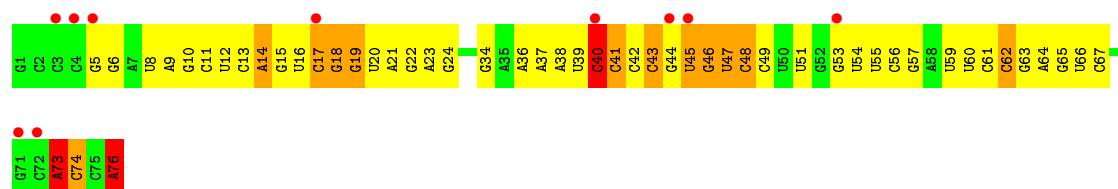




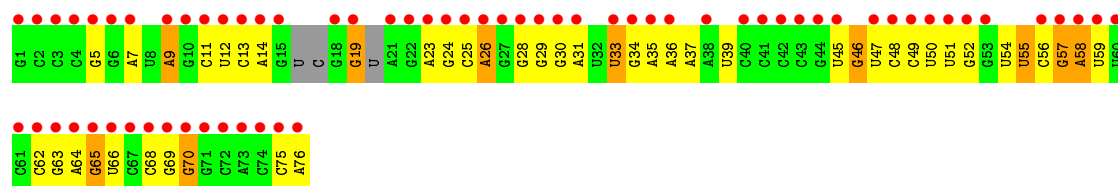
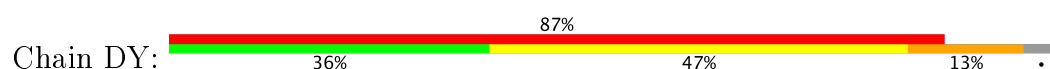
● Molecule 56: P-site tRNA



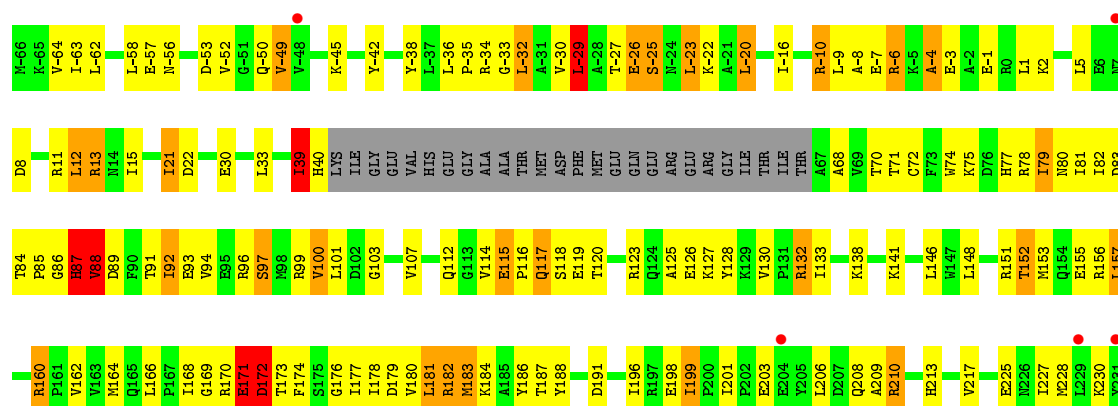
● Molecule 56: P-site tRNA

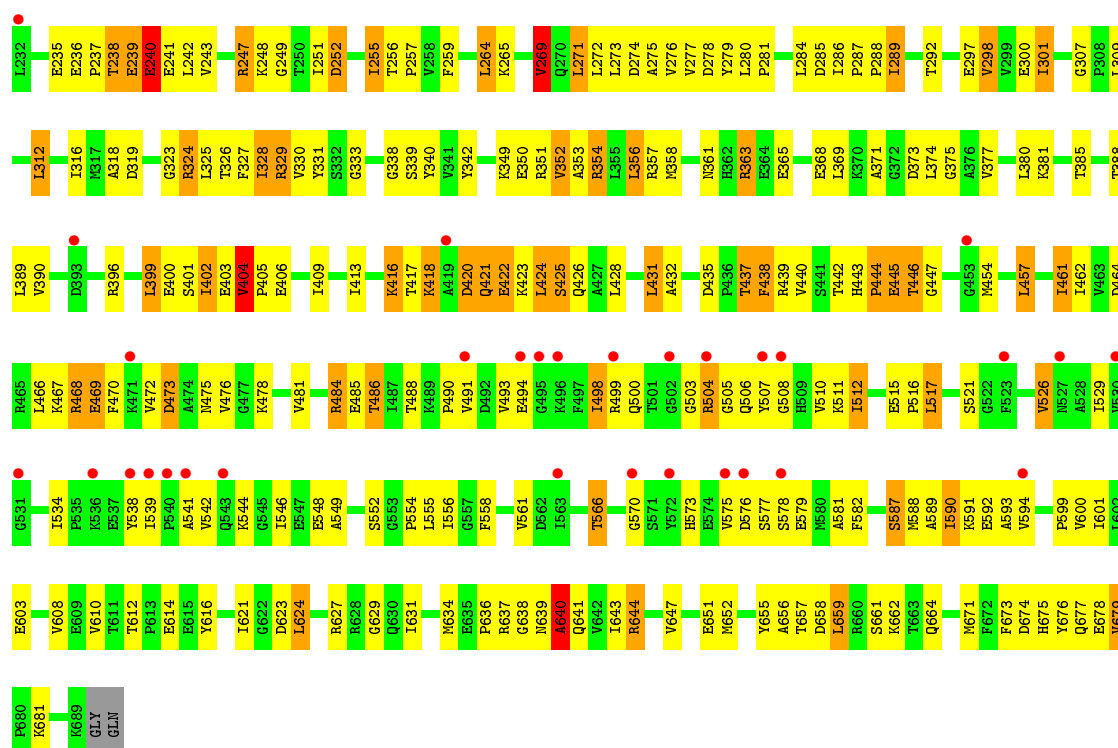


● Molecule 56: P-site tRNA

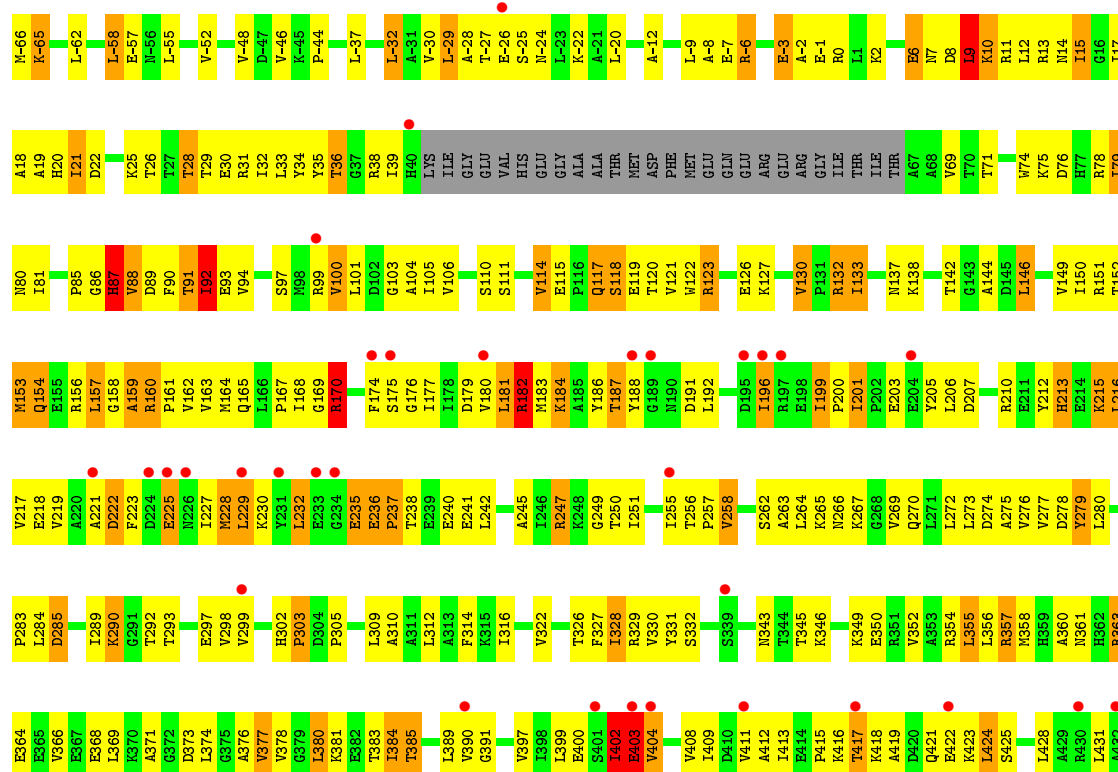


● Molecule 57: 50S ribosomal protein L9,Elongation factor G





- Molecule 57: 50S ribosomal protein L9,Elongation factor G



G638	L567	E493	Y507	L567	E493	Y507	L567	G638
M639	Y568	E494	G508	Y568	E494	G508	M639	M639
A640	D569	D435	H509	D569	D435	H509	A640	A640
Q641	G570	P436	V510	G570	P436	V510	Q641	Q641
V642	S571	T437	K511	S571	T437	K511	V642	V642
I643	Y572	F438	I512	Y572	F438	I512	I643	I643
R644	H573	R439	K513	H573	R439	K513	R644	R644
A645	E574	V440	V514	E574	V440	V514	A645	A645
F646	V575	P444	E515	V575	P444	E515	F646	F646
P647	D576	E445	P516	D576	E445	P516	P647	P647
P648	S577	T446	L517	S577	T446	L517	P648	P648
L649	S578		P518	S578		P518	L649	L649
A650	E579		R519	E579		R519	A650	A650
E651	M580	I451	G520	M580	I451	G520	E651	E651
M652	A581	S452	S521	M652	A581	S452	M652	M652
F653	F582	G453	G522	F653	F582	G453	F653	F653
G654		M454	F523	G654		M454	G654	G654
Y655		G455	E524	Y655		G455	Y655	Y655
A656	A589	E456	F525	A656	A589	E456	A656	A656
T657	I590	L457	V526	T657	I590	L457	T657	T657
D658		H458	M527	D658		H458	D658	D658
L659	V594	L459	A528	L659	V594	L459	L659	L659
R660	Q595	E460	K536	R660	Q595	E460	R660	R660
S661	K596	I461	I529	S661	K596	I461	S661	S661
K662	G597		V530	K662	G597		K662	K662
	D598		G531		D598			
	P599		G532		P599			
F669	V600	L466	V533	F669	V600	L466	F669	F669
V670	I601	R467	I534	V670	I601	R467	V670	V670
M671	L602	R468	P535	M671	L602	R468	M671	M671
F672		K471	K536	F672		K471	F672	F672
F673		V472	E537	F673		V472	F673	F673
D674	I605	V476	V538	D674	I605	V476	D674	D674
H675	M606		I539	H675	M606		H675	H675
Y676	R607		P540	Y676	R607		Y676	Y676
Q677	V608		A541	Q677	V608		Q677	Q677
E678	E609		V542	E678	E609		E678	E678
	V610		O543		V610			
	T611		K544		T611			
K681	T612		G545	K681	T612		K681	K681
Q682	P613		I546	Q682	P613		Q682	Q682
V683	E614		E547	V683	E614		V683	V683
Q684	E615		E548	Q684	E615		Q684	Q684
E685			A549	E685			E685	E685
	D619		M550		D619			
	I620		Q551		I620			
K689	G622		S552	K689	G622		K689	K689
GLY	D623		G553	GLY	D623		GLY	GLY
GLN	L624		P554	GLN	L624		GLN	GLN
	N625		I555		N625			
	A626		I556		A626			
			G557					
	G629		P558		G629			
	Q630		P559		Q630			
	I631		V560		I631			
	L632		V561		L632			
	G633		D562		G633			
	M634		I563		M634			
	E635		K564		E635			
	P636		V565		P636			
	R637		T566		R637			

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	210.45Å 449.00Å 625.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.74 – 2.80 49.74 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.74-2.80) 99.1 (49.74-2.79)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 2.77Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8.2_1309)	Depositor
R, R_{free}	0.200 , 0.255 0.211 , 0.262	Depositor DCC
R_{free} test set	71854 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	54.2	Xtriage
Anisotropy	0.399	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 67.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	310279	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.54% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: 5MU, GDP, ZN, MIA, SF4, MG, FUA, 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	AA	1.51	654/68792 (1.0%)	2.24	4887/107377 (4.6%)
1	CA	1.01	71/68691 (0.1%)	1.68	1752/107219 (1.6%)
2	AB	1.21	5/2878 (0.2%)	2.01	141/4490 (3.1%)
2	CB	0.67	0/2878	1.30	18/4490 (0.4%)
3	AC	0.34	0/1083	0.65	0/1460
3	CC	0.34	0/1083	0.65	0/1460
4	AD	1.00	4/2186 (0.2%)	1.08	8/2944 (0.3%)
4	CD	0.76	0/2192	0.92	2/2951 (0.1%)
5	AE	1.03	5/1592 (0.3%)	1.08	4/2149 (0.2%)
5	CE	0.68	0/1592	0.85	1/2149 (0.0%)
6	AF	0.96	2/1619 (0.1%)	1.07	3/2193 (0.1%)
6	CF	0.64	0/1615	0.85	1/2188 (0.0%)
7	AG	0.55	0/1450	0.77	0/1959
7	CG	0.40	0/1449	0.63	0/1958
8	AH	0.84	0/1356	0.96	3/1834 (0.2%)
8	CH	0.42	0/1356	0.64	0/1834
9	AK	0.42	0/640	0.76	1/889 (0.1%)
9	CK	0.29	0/640	0.62	0/889
10	AL	0.34	0/1044	0.58	0/1416
10	CL	0.31	0/1044	0.53	0/1416
11	AN	1.06	2/1144 (0.2%)	1.09	3/1543 (0.2%)
11	CN	0.55	0/1144	0.76	0/1543
12	AO	1.00	0/943	1.09	2/1269 (0.2%)
12	CO	0.71	0/943	0.82	0/1269
13	AP	0.89	0/1156	1.08	9/1537 (0.6%)
13	CP	0.58	0/1152	0.86	1/1533 (0.1%)
14	AQ	0.98	0/1143	1.05	4/1527 (0.3%)
14	CQ	0.62	0/1143	0.77	0/1527
15	AR	0.98	0/982	1.15	4/1312 (0.3%)
15	CR	0.62	0/982	0.95	2/1312 (0.2%)
16	AS	0.76	0/887	0.96	0/1180
16	CS	0.56	0/880	0.83	2/1172 (0.2%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
17	AT	0.87	0/1105	1.04	2/1477 (0.1%)
17	CT	0.68	0/1097	0.89	0/1468
18	AU	1.18	2/977 (0.2%)	1.19	5/1301 (0.4%)
18	CU	0.67	0/977	0.78	0/1301
19	AV	1.02	1/782 (0.1%)	1.09	1/1049 (0.1%)
19	CV	0.54	0/782	0.74	0/1049
20	AW	1.21	1/897 (0.1%)	1.19	3/1205 (0.2%)
20	CW	0.77	0/897	0.91	0/1205
21	AX	0.98	1/764 (0.1%)	0.99	2/1025 (0.2%)
21	CX	0.70	0/764	0.78	1/1025 (0.1%)
22	AY	0.92	0/819	1.00	1/1095 (0.1%)
22	CY	0.62	0/819	0.77	0/1095
23	AZ	0.77	0/1483	1.00	3/2017 (0.1%)
23	CZ	0.47	0/1483	0.71	0/2017
24	A0	0.97	0/662	1.03	0/881
24	C0	0.60	0/662	0.77	0/881
25	A1	0.94	0/762	1.04	3/1014 (0.3%)
25	C1	0.70	0/762	0.86	0/1014
26	A2	0.88	0/590	0.91	0/781
26	C2	0.60	0/590	0.75	0/781
27	A3	0.99	0/474	1.09	2/635 (0.3%)
27	C3	0.51	0/469	0.78	0/630
28	A4	0.47	0/571	0.74	0/768
28	C4	0.35	0/545	0.59	0/737
29	A5	1.22	3/469 (0.6%)	1.22	4/635 (0.6%)
29	C5	0.75	0/469	0.89	2/635 (0.3%)
30	A6	0.89	0/460	1.02	1/613 (0.2%)
30	C6	0.67	0/456	0.76	0/608
31	A7	1.11	0/426	1.21	4/561 (0.7%)
31	C7	0.86	0/426	1.03	2/561 (0.4%)
32	A8	1.00	0/525	1.04	3/691 (0.4%)
32	C8	0.72	0/525	0.85	0/691
33	A9	1.09	0/310	1.04	0/407
33	C9	0.57	0/310	0.70	0/407
34	BA	0.79	7/35976 (0.0%)	1.44	496/56145 (0.9%)
34	DA	0.70	2/36119 (0.0%)	1.31	266/56370 (0.5%)
35	BB	0.47	0/1881	0.72	0/2542
35	DB	0.39	0/1860	0.64	0/2518
36	BC	0.38	0/1576	0.57	0/2130
36	DC	0.34	0/1568	0.54	0/2122
37	BD	0.52	0/1689	0.77	0/2267
37	DD	0.51	0/1708	0.77	1/2289 (0.0%)
38	BE	0.62	0/1145	0.84	0/1543

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
38	DE	0.54	0/1149	0.78	1/1548 (0.1%)
39	BF	0.54	0/825	0.75	0/1118
39	DF	0.54	0/833	0.77	2/1128 (0.2%)
40	BG	0.40	0/1250	0.60	0/1679
40	DG	0.33	0/1254	0.52	0/1683
41	BH	0.58	0/1108	0.80	1/1494 (0.1%)
41	DH	0.48	0/1108	0.70	0/1494
42	BI	0.38	0/1005	0.61	0/1350
42	DI	0.33	0/997	0.58	0/1343
43	BJ	0.38	0/722	0.59	0/982
43	DJ	0.34	0/727	0.57	0/988
44	BK	0.55	0/848	0.75	0/1149
44	DK	0.51	0/848	0.70	0/1149
45	BL	0.74	0/946	0.88	1/1274 (0.1%)
45	DL	0.58	0/946	0.81	0/1274
46	BM	0.39	0/933	0.62	0/1253
46	DM	0.34	0/961	0.55	0/1291
47	BN	0.39	0/501	0.68	1/664 (0.2%)
47	DN	0.35	0/501	0.56	0/664
48	BO	0.56	0/739	0.81	0/985
48	DO	0.53	0/739	0.72	0/985
49	BP	0.55	0/697	0.79	1/939 (0.1%)
49	DP	0.53	0/693	0.71	0/935
50	BQ	0.63	0/836	0.78	0/1117
50	DQ	0.57	0/836	0.73	0/1117
51	BR	0.55	0/560	0.80	1/746 (0.1%)
51	DR	0.51	0/560	0.71	0/746
52	BS	0.35	0/676	0.57	0/911
52	DS	0.31	0/661	0.61	0/893
53	BT	0.49	0/730	0.74	0/965
53	DT	0.49	0/733	0.71	0/969
54	BU	0.38	0/203	0.65	0/266
54	DU	0.31	0/203	0.57	0/266
55	BV	0.71	0/165	1.15	1/254 (0.4%)
55	DV	0.60	0/137	1.05	0/211
56	BW	0.89	1/1650 (0.1%)	1.64	41/2569 (1.6%)
56	BY	0.42	0/1602	0.95	1/2493 (0.0%)
56	DW	0.70	0/1650	1.36	20/2569 (0.8%)
56	DY	0.35	0/1579	0.86	0/2455
57	BZ	0.58	0/5792	0.81	4/7844 (0.1%)
57	DZ	0.49	0/5792	0.72	4/7844 (0.1%)
All	All	0.99	761/330005 (0.2%)	1.56	7729/491779 (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
1	CA	0	1
17	CT	0	1
19	AV	0	1
21	AX	0	1
21	CX	0	1
23	AZ	0	1
24	A0	0	1
25	A1	0	1
28	A4	0	1
53	BT	0	1
53	DT	0	1
57	BZ	0	1
57	DZ	0	3
All	All	0	15

All (761) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	1067	A	N9-C4	-18.69	1.26	1.37
1	AA	1188	A	N9-C4	-16.21	1.28	1.37
1	AA	990	A	N9-C4	-15.62	1.28	1.37
1	AA	354	A	N9-C4	-13.34	1.29	1.37
1	AA	1988	A	N9-C4	-12.57	1.30	1.37
1	AA	2299	A	N9-C4	-12.41	1.30	1.37
1	AA	555	G	C2-N3	-11.33	1.23	1.32
1	AA	553	A	N9-C8	11.14	1.46	1.37
1	AA	553	A	C5-C6	-11.12	1.31	1.41
1	CA	528	A	N9-C4	-10.51	1.31	1.37
1	AA	978	A	N9-C4	-10.47	1.31	1.37
4	AD	28	GLU	CG-CD	9.95	1.66	1.51
1	AA	254	A	N9-C4	-9.80	1.31	1.37
1	AA	354	A	N9-C8	9.77	1.45	1.37
1	AA	2290	A	N3-C4	-9.74	1.29	1.34
1	AA	2285	A	C5-C6	-9.55	1.32	1.41
1	AA	1342	G	N1-C2	-9.43	1.30	1.37
1	AA	978	A	C5-C6	-9.31	1.32	1.41
1	CA	330	A	N9-C4	-9.22	1.32	1.37
1	AA	1001	G	C6-O6	9.16	1.32	1.24
1	CA	1204	A	N9-C4	-9.14	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	978	A	N3-C4	-9.14	1.29	1.34
1	AA	2043	C	N1-C6	-9.00	1.31	1.37
1	AA	586	G	C6-O6	8.99	1.32	1.24
1	AA	1272	A	N9-C4	-8.98	1.32	1.37
1	AA	550	U	C2-O2	-8.92	1.14	1.22
1	AA	851	A	N9-C4	-8.70	1.32	1.37
1	AA	2553	A	C6-N1	-8.70	1.29	1.35
1	AA	990	A	C5-C6	-8.59	1.33	1.41
1	CA	1142(A)	A	N9-C4	-8.58	1.32	1.37
1	AA	2331	G	N9-C4	-8.55	1.31	1.38
1	AA	555	G	N9-C8	8.52	1.43	1.37
1	AA	553	A	C5-C4	8.46	1.44	1.38
1	AA	1988	A	N3-C4	-8.45	1.29	1.34
1	AA	829	A	N7-C5	-8.45	1.34	1.39
1	AA	1249	A	C5-C6	-8.42	1.33	1.41
1	AA	2037	A	N3-C4	-8.32	1.29	1.34
1	AA	558	G	C5-C4	-8.23	1.32	1.38
1	AA	1067	A	C5-C6	-8.23	1.33	1.41
1	AA	254	A	C5-C6	-8.22	1.33	1.41
1	AA	553	A	N3-C4	-8.20	1.29	1.34
1	AA	1189	A	N3-C4	-8.16	1.29	1.34
1	AA	1314	A	N3-C4	-8.14	1.29	1.34
1	AA	2446	A	N7-C5	-8.07	1.34	1.39
1	AA	2269	U	C4-O4	-7.98	1.17	1.23
1	AA	2278	A	C5-C4	-7.98	1.33	1.38
1	AA	354	A	C5-C6	-7.97	1.33	1.41
1	AA	990	A	C5-C4	7.96	1.44	1.38
1	AA	550	U	C2-N3	-7.91	1.32	1.37
4	AD	28	GLU	CB-CG	7.86	1.67	1.52
1	AA	1745	A	C5-C6	-7.86	1.33	1.41
1	AA	1293	A	N3-C4	-7.85	1.30	1.34
1	AA	590	A	C5-C4	-7.84	1.33	1.38
1	AA	1076	G	C5-C4	-7.83	1.32	1.38
1	AA	710	G	N3-C4	-7.83	1.29	1.35
1	AA	1249	A	N7-C5	-7.80	1.34	1.39
1	CA	1791	A	N3-C4	-7.77	1.30	1.34
1	AA	2045	G	C5-C6	-7.73	1.34	1.42
1	AA	1249	A	N3-C4	-7.71	1.30	1.34
1	AA	726	C	N1-C6	-7.68	1.32	1.37
1	AA	2331	G	N3-C4	-7.68	1.30	1.35
1	AA	2657	G	N9-C4	-7.67	1.31	1.38
1	AA	1054	C	C4-N4	-7.67	1.27	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	CA	945	A	N9-C4	-7.64	1.33	1.37
1	AA	2298	A	C5-C4	7.64	1.44	1.38
1	AA	1311	A	N7-C5	-7.64	1.34	1.39
1	AA	1075	A	N3-C4	-7.63	1.30	1.34
1	AA	2777	A	N9-C4	-7.63	1.33	1.37
1	AA	470	C	N1-C2	-7.62	1.32	1.40
1	AA	2636	G	N9-C8	-7.61	1.32	1.37
1	AA	1076	G	C5-C6	-7.55	1.34	1.42
1	AA	2601	A	N3-C4	-7.53	1.30	1.34
1	AA	1302	G	C5-C4	-7.52	1.33	1.38
1	AA	1019	G	N7-C5	-7.51	1.34	1.39
1	AA	2775	G	C6-N1	-7.50	1.34	1.39
1	AA	2331	G	C2-N3	-7.49	1.26	1.32
1	AA	1274	G	C6-N1	-7.46	1.34	1.39
1	AA	2757	G	C2-N3	-7.46	1.26	1.32
1	CA	2441	C	N1-C6	-7.45	1.32	1.37
1	AA	896	A	N9-C4	-7.44	1.33	1.37
1	AA	2518	U	N1-C2	7.44	1.45	1.38
1	AA	2636	G	C5-C4	-7.43	1.33	1.38
1	AA	2083	G	N7-C5	-7.40	1.34	1.39
1	AA	2654	G	C6-N1	-7.39	1.34	1.39
1	AA	231	G	C6-N1	-7.37	1.34	1.39
1	AA	1188	A	C2-N3	-7.36	1.26	1.33
1	AA	1661	C	N1-C6	-7.36	1.32	1.37
1	AA	1659	G	C6-O6	-7.33	1.17	1.24
1	AA	559	U	C2-N3	-7.33	1.32	1.37
1	AA	2036	A	C6-N1	7.32	1.40	1.35
1	AA	831	A	C5-C4	-7.32	1.33	1.38
1	AA	1600	A	N7-C5	-7.31	1.34	1.39
1	AA	2500	A	N9-C4	-7.29	1.33	1.37
1	AA	1707	C	N3-C4	-7.29	1.28	1.33
1	AA	1745	A	N3-C4	-7.27	1.30	1.34
1	AA	856	G	C6-O6	-7.23	1.17	1.24
1	AA	593	G	C5-C4	-7.21	1.33	1.38
1	AA	2601	A	N9-C4	-7.21	1.33	1.37
1	AA	2451	A	N7-C5	-7.21	1.34	1.39
1	AA	2785	C	N1-C6	-7.19	1.32	1.37
1	AA	865	G	N3-C4	-7.14	1.30	1.35
1	AA	2278	A	N3-C4	-7.13	1.30	1.34
1	AA	1605	A	N7-C5	-7.06	1.35	1.39
1	AA	990	A	N1-C2	7.05	1.40	1.34
1	AA	598	A	C5-C4	-7.04	1.33	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	2000	A	C5-C4	-7.02	1.33	1.38
1	AA	591	U	C4-O4	-7.01	1.18	1.23
1	AA	356	A	N9-C4	-7.00	1.33	1.37
1	CA	256	A	N7-C5	-6.97	1.35	1.39
1	CA	1378	A	N9-C4	-6.93	1.33	1.37
1	AA	1026	A	N9-C4	-6.92	1.33	1.37
1	AA	799	A	N9-C4	-6.91	1.33	1.37
1	AA	2300	A	N9-C4	6.90	1.42	1.37
1	AA	2063	U	C2-O2	-6.90	1.16	1.22
1	AA	598	A	N7-C5	-6.89	1.35	1.39
1	AA	782	A	N3-C4	-6.88	1.30	1.34
1	AA	2858	G	C5-C6	-6.88	1.35	1.42
1	AA	729	G	C6-N1	-6.88	1.34	1.39
1	AA	2067	C	N3-C4	6.85	1.38	1.33
1	CA	1558	A	N9-C4	-6.85	1.33	1.37
1	AA	2820	A	N3-C4	-6.84	1.30	1.34
1	AA	53	G	N3-C4	-6.83	1.30	1.35
1	AA	990	A	N3-C4	-6.81	1.30	1.34
34	DA	1513	A	N9-C4	-6.80	1.33	1.37
1	AA	2876	U	C4-C5	6.79	1.49	1.43
1	CA	2225	A	N9-C4	-6.78	1.33	1.37
1	AA	1605	A	C5-C6	-6.76	1.34	1.41
1	AA	590	A	N1-C2	-6.75	1.28	1.34
1	AA	1241	C	N1-C6	-6.75	1.33	1.37
1	AA	872	C	N1-C6	-6.75	1.33	1.37
1	AA	1019	G	N9-C8	-6.75	1.33	1.37
1	AA	1050	C	C2-O2	-6.75	1.18	1.24
1	AA	1809	U	N1-C2	-6.74	1.32	1.38
1	AA	2778	A	N9-C8	-6.73	1.32	1.37
1	AA	1067	A	N9-C8	6.73	1.43	1.37
1	AA	2757	G	C5-C4	-6.73	1.33	1.38
1	AA	710	G	C5-C4	-6.72	1.33	1.38
4	AD	39	LYS	CE-NZ	6.72	1.65	1.49
1	AA	1261	G	C5-C4	-6.72	1.33	1.38
1	AA	553	A	N7-C5	-6.72	1.35	1.39
1	AA	978	A	N9-C8	6.71	1.43	1.37
1	AA	26	G	C6-O6	-6.71	1.18	1.24
1	AA	1711	A	N9-C4	-6.70	1.33	1.37
1	AA	1020	C	C2-N3	-6.70	1.30	1.35
1	AA	1605	A	N3-C4	-6.70	1.30	1.34
1	AA	2612	A	C6-N1	-6.70	1.30	1.35
1	AA	2627	U	C4-O4	-6.69	1.18	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	2641	A	C5-C6	-6.69	1.35	1.41
1	AA	1959	A	N9-C4	-6.69	1.33	1.37
1	AA	1725	G	N7-C5	-6.68	1.35	1.39
1	AA	835	A	N9-C4	-6.67	1.33	1.37
1	AA	1261	G	C5-C6	-6.67	1.35	1.42
1	AA	1711	A	N3-C4	-6.67	1.30	1.34
1	AA	171	A	N9-C4	-6.65	1.33	1.37
1	AA	2466	G	N3-C4	-6.65	1.30	1.35
1	CA	211	A	N9-C4	-6.64	1.33	1.37
1	AA	592	U	C2-N3	-6.63	1.33	1.37
1	AA	876	A	N7-C5	-6.62	1.35	1.39
1	AA	2060	G	N1-C2	-6.62	1.32	1.37
1	AA	2466	G	N1-C2	-6.61	1.32	1.37
1	AA	1243	U	C2-N3	-6.61	1.33	1.37
1	AA	2285	A	C6-N6	-6.61	1.28	1.33
1	AA	543	G	N9-C8	-6.60	1.33	1.37
1	AA	1745	A	N9-C4	-6.59	1.33	1.37
1	AA	978	A	N7-C5	-6.59	1.35	1.39
1	AA	1613	A	C6-N1	-6.59	1.30	1.35
1	CA	465	G	C6-N1	-6.58	1.34	1.39
1	AA	139	A	N9-C4	-6.58	1.33	1.37
1	AA	874	U	N1-C2	-6.58	1.32	1.38
1	AA	2013	U	C2-N3	-6.58	1.33	1.37
1	AA	2691	A	N3-C4	-6.57	1.30	1.34
1	AA	2778	A	N7-C5	-6.56	1.35	1.39
1	AA	732	A	N9-C4	-6.56	1.33	1.37
1	AA	2609	G	C5-C6	-6.55	1.35	1.42
1	AA	2239	A	N7-C5	-6.54	1.35	1.39
1	AA	1702	A	N9-C4	-6.54	1.33	1.37
1	AA	2641	A	N9-C8	6.53	1.43	1.37
1	AA	617	U	C2-N3	-6.52	1.33	1.37
1	AA	1200	G	C6-N1	-6.52	1.34	1.39
1	CA	2824	C	N1-C6	-6.52	1.33	1.37
1	AA	168	G	N3-C4	-6.51	1.30	1.35
1	CA	2287	A	N9-C4	-6.48	1.33	1.37
1	AA	1244	U	C2-N3	-6.46	1.33	1.37
1	AA	531	G	N1-C2	-6.46	1.32	1.37
1	AA	1993	A	N7-C5	-6.45	1.35	1.39
1	AA	597	C	N3-C4	-6.45	1.29	1.33
1	AA	2515	A	N3-C4	6.43	1.38	1.34
1	AA	2609	G	C5-C4	-6.43	1.33	1.38
1	AA	2840	G	N7-C5	-6.43	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	1061	G	N7-C5	-6.43	1.35	1.39
1	AA	2093	A	C5-C4	-6.42	1.34	1.38
1	AA	2737	C	N1-C6	-6.41	1.33	1.37
1	AA	2068	G	N7-C5	-6.40	1.35	1.39
1	AA	841	G	N3-C4	6.40	1.40	1.35
1	AA	1307	C	C4-N4	-6.40	1.28	1.33
1	AA	996	C	N3-C4	-6.39	1.29	1.33
1	AA	2278	A	C6-N1	-6.38	1.31	1.35
1	CA	37	C	N1-C6	-6.38	1.33	1.37
1	AA	126	C	N1-C6	-6.38	1.33	1.37
1	AA	716	G	C5-C4	-6.38	1.33	1.38
1	AA	1000	C	C2-O2	-6.38	1.18	1.24
1	AA	2389	A	N9-C4	-6.38	1.34	1.37
1	AA	2530	A	N3-C4	-6.37	1.31	1.34
1	AA	352	U	C4-O4	-6.36	1.18	1.23
1	AA	2037	A	N7-C5	-6.36	1.35	1.39
1	AA	2475	C	C2-O2	-6.35	1.18	1.24
1	AA	1657	C	N1-C6	-6.34	1.33	1.37
21	AX	15	GLU	CG-CD	6.34	1.61	1.51
6	AF	103	LYS	CE-NZ	6.34	1.65	1.49
1	AA	597	C	C4-C5	-6.34	1.37	1.43
1	AA	2584	A	N9-C4	-6.34	1.34	1.37
1	AA	553	A	N9-C4	-6.34	1.34	1.37
1	AA	2093	A	N7-C5	-6.33	1.35	1.39
1	AA	470	C	C2-N3	-6.33	1.30	1.35
1	AA	593	G	N7-C5	-6.33	1.35	1.39
1	AA	1721	G	N7-C5	-6.32	1.35	1.39
1	AA	1283	A	N9-C8	-6.32	1.32	1.37
1	AA	521	G	N9-C8	-6.31	1.33	1.37
1	AA	16	G	N7-C5	-6.29	1.35	1.39
1	AA	654	G	N7-C5	-6.29	1.35	1.39
1	AA	2693	C	N1-C6	-6.28	1.33	1.37
1	AA	416	G	C5-C4	-6.27	1.33	1.38
1	AA	2428	C	N1-C2	-6.27	1.33	1.40
1	AA	709	G	C2-N3	-6.26	1.27	1.32
1	AA	555	G	N3-C4	-6.25	1.31	1.35
1	AA	896	A	N3-C4	-6.25	1.31	1.34
34	BA	1483	A	N9-C4	-6.24	1.34	1.37
1	AA	775	G	N1-C2	-6.23	1.32	1.37
1	AA	2617	U	C2-N3	-6.22	1.33	1.37
1	AA	608	G	C5-C4	-6.22	1.33	1.38
1	AA	1382	A	C5-C4	-6.21	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	738	C	C2-O2	-6.20	1.18	1.24
1	CA	2764	A	N9-C4	-6.20	1.34	1.37
1	AA	753	A	N9-C4	-6.20	1.34	1.37
1	AA	2893	A	N9-C4	6.19	1.41	1.37
1	AA	2590	G	N7-C5	-6.19	1.35	1.39
1	AA	2794	A	N3-C4	-6.19	1.31	1.34
1	AA	1067	A	N3-C4	-6.19	1.31	1.34
1	AA	2619	G	C5-C4	-6.19	1.34	1.38
1	AA	2277	U	N1-C6	-6.19	1.32	1.38
1	AA	1388	A	N3-C4	-6.17	1.31	1.34
1	AA	549	U	C4-O4	-6.17	1.18	1.23
1	AA	2299	A	C5-C6	-6.17	1.35	1.41
1	AA	1020	C	C2-O2	-6.17	1.18	1.24
1	AA	52	A	N9-C4	-6.16	1.34	1.37
1	AA	1329	G	N7-C5	-6.16	1.35	1.39
1	AA	543	G	N7-C5	-6.15	1.35	1.39
1	AA	650	G	N7-C5	-6.15	1.35	1.39
1	AA	2830	A	N3-C4	-6.15	1.31	1.34
1	AA	1707	C	N1-C6	-6.14	1.33	1.37
1	AA	2371	C	N1-C6	-6.14	1.33	1.37
1	AA	1303	C	N1-C6	-6.14	1.33	1.37
1	AA	725	C	N3-C4	-6.13	1.29	1.33
1	AA	354	A	C5-C4	6.12	1.43	1.38
1	AA	2581	G	N3-C4	-6.12	1.31	1.35
1	AA	2092	G	N3-C4	-6.11	1.31	1.35
1	CA	2060	A	N9-C4	-6.11	1.34	1.37
11	AN	104	LYS	CE-NZ	6.11	1.64	1.49
1	AA	1733	C	N1-C2	6.11	1.46	1.40
1	AA	2278	A	C5-C6	-6.10	1.35	1.41
1	AA	2003	A	C6-N1	-6.10	1.31	1.35
1	AA	2003	A	N3-C4	-6.09	1.31	1.34
1	AA	1054	C	N3-C4	-6.09	1.29	1.33
1	AA	552	C	N1-C6	-6.08	1.33	1.37
1	AA	1066	A	N7-C5	-6.08	1.35	1.39
1	AA	541	C	N3-C4	-6.08	1.29	1.33
1	AA	1066	A	C5-C6	-6.07	1.35	1.41
1	AA	2029	C	N1-C6	-6.07	1.33	1.37
1	AA	255	G	C5-C4	-6.07	1.34	1.38
1	AA	1282	G	C5-C4	-6.07	1.34	1.38
1	AA	884	C	N3-C4	-6.06	1.29	1.33
5	AE	163	GLU	CG-CD	6.06	1.61	1.51
1	AA	552	C	N3-C4	-6.05	1.29	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	2466	G	C6-N1	-6.05	1.35	1.39
1	AA	2291	G	N7-C5	-6.04	1.35	1.39
1	AA	2514	G	C5-C4	-6.04	1.34	1.38
1	AA	613	A	N7-C5	-6.03	1.35	1.39
1	AA	1081	U	C2-N3	-6.03	1.33	1.37
1	AA	1845	G	C5-C6	-6.02	1.36	1.42
1	AA	1035	G	C2-N3	-6.02	1.27	1.32
1	AA	2035	A	N9-C4	-6.02	1.34	1.37
1	AA	791	G	N7-C5	-6.02	1.35	1.39
20	AW	30	GLU	CG-CD	6.01	1.60	1.51
1	AA	2044	U	N1-C6	-6.01	1.32	1.38
1	AA	2294	G	N1-C2	-6.01	1.32	1.37
1	AA	473	A	C6-N1	-6.01	1.31	1.35
1	AA	2605	U	C2-N3	-6.00	1.33	1.37
1	AA	1067	A	N7-C5	-6.00	1.35	1.39
1	AA	1274	G	N9-C4	-6.00	1.33	1.38
1	AA	1055	A	N9-C4	-5.99	1.34	1.37
1	AA	2405	A	N7-C5	-5.99	1.35	1.39
1	AA	2834	C	N1-C6	-5.99	1.33	1.37
1	AA	2562	G	N7-C5	-5.98	1.35	1.39
1	AA	605	G	C6-N1	-5.98	1.35	1.39
1	AA	1048	G	C2-N2	-5.98	1.28	1.34
1	AA	2034	G	C6-O6	-5.98	1.18	1.24
1	AA	595	A	N7-C5	-5.97	1.35	1.39
1	AA	1795	G	N1-C2	-5.97	1.32	1.37
1	AA	1345	G	N7-C5	-5.97	1.35	1.39
1	AA	1686	U	C4-O4	-5.96	1.18	1.23
29	A5	6	VAL	CB-CG2	-5.96	1.40	1.52
1	AA	2054	G	C6-N1	-5.96	1.35	1.39
1	AA	1395	A	N3-C4	5.95	1.38	1.34
1	AA	1029	A	C8-N7	-5.95	1.27	1.31
1	CA	1822	G	N3-C4	-5.95	1.31	1.35
1	AA	2019	G	C2-N3	5.94	1.37	1.32
1	CA	515	A	N3-C4	-5.94	1.31	1.34
1	AA	1249	A	N9-C8	5.93	1.42	1.37
1	AA	555	G	C6-N1	-5.93	1.35	1.39
1	AA	1302	G	N7-C5	-5.93	1.35	1.39
1	AA	2565	G	N9-C8	-5.93	1.33	1.37
1	AA	1041	C	N1-C6	-5.93	1.33	1.37
1	CA	2042	A	N9-C4	-5.92	1.34	1.37
1	AA	2650	G	N7-C5	-5.92	1.35	1.39
1	AA	597	C	N1-C2	-5.92	1.34	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	180	A	N9-C8	-5.91	1.33	1.37
1	AA	1711	A	C5-C6	-5.91	1.35	1.41
1	AA	119	G	N3-C4	-5.91	1.31	1.35
1	AA	470	C	C2-O2	-5.91	1.19	1.24
1	AA	561	A	C5-C4	-5.91	1.34	1.38
1	AA	1659	G	C5-C6	-5.91	1.36	1.42
1	AA	2084	A	C5-C6	5.91	1.46	1.41
1	AA	2828	G	C6-N1	-5.91	1.35	1.39
1	AA	1237	G	C5-C4	-5.90	1.34	1.38
1	AA	630	U	N1-C2	-5.90	1.33	1.38
1	AA	1441	A	C5-C6	-5.90	1.35	1.41
1	AA	780	G	N7-C5	-5.89	1.35	1.39
1	AA	1472	G	C5-C6	-5.89	1.36	1.42
1	AA	1343	C	N1-C6	-5.88	1.33	1.37
1	CA	310	A	N9-C4	-5.88	1.34	1.37
1	AA	1234	A	N3-C4	-5.87	1.31	1.34
1	AA	1255	A	C5-C4	-5.87	1.34	1.38
1	CA	1378	A	N3-C4	-5.87	1.31	1.34
1	AA	2711	C	N1-C6	-5.86	1.33	1.37
1	AA	2040	G	N9-C8	-5.86	1.33	1.37
1	AA	183	G	C6-O6	5.86	1.29	1.24
1	AA	2476	C	C4-N4	-5.85	1.28	1.33
1	AA	842	C	C2-O2	-5.85	1.19	1.24
1	AA	2505	U	C2-O2	-5.85	1.17	1.22
1	CA	2438	U	C2-N3	-5.85	1.33	1.37
1	AA	598	A	C8-N7	-5.83	1.27	1.31
34	BA	1523	G	N3-C4	-5.83	1.31	1.35
1	AA	710	G	C2-N3	-5.83	1.28	1.32
1	AA	826	U	C5-C6	-5.83	1.28	1.34
1	AA	825	G	N7-C5	-5.83	1.35	1.39
1	CA	2058	A	N3-C4	-5.82	1.31	1.34
1	AA	1332	A	N7-C5	-5.82	1.35	1.39
1	AA	990	A	N7-C5	-5.82	1.35	1.39
34	BA	1511	G	C5-C4	-5.82	1.34	1.38
1	AA	1378	G	N3-C4	-5.82	1.31	1.35
1	AA	2525	G	N3-C4	5.82	1.39	1.35
1	AA	2605	U	N3-C4	-5.81	1.33	1.38
1	AA	2525	G	C8-N7	-5.81	1.27	1.30
1	AA	2490	A	N9-C4	-5.81	1.34	1.37
1	CA	1791	A	C5-C4	-5.80	1.34	1.38
1	AA	1045	U	C4-O4	-5.80	1.19	1.23
1	AA	499	G	C6-N1	-5.80	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	2278	A	N1-C2	-5.80	1.29	1.34
1	AA	2820	A	N7-C5	-5.80	1.35	1.39
1	AA	889	G	N3-C4	-5.79	1.31	1.35
1	AA	1332	A	N3-C4	-5.79	1.31	1.34
1	AA	702	A	N9-C4	-5.79	1.34	1.37
1	CA	1781	C	N3-C4	5.79	1.38	1.33
1	AA	2383	G	C5-C4	-5.78	1.34	1.38
1	AA	2442	A	C6-N1	-5.77	1.31	1.35
1	AA	1287	A	N9-C8	-5.77	1.33	1.37
1	AA	855	G	N7-C5	-5.77	1.35	1.39
1	AA	2493	G	N7-C5	-5.76	1.35	1.39
1	AA	1679	A	N3-C4	-5.76	1.31	1.34
1	AA	1049	G	C8-N7	-5.75	1.27	1.30
1	CA	1237	A	N9-C4	-5.75	1.34	1.37
1	AA	2068	G	C6-O6	-5.75	1.19	1.24
1	CA	2585	U	C2-N3	5.73	1.41	1.37
1	AA	866	A	C5-C4	-5.73	1.34	1.38
1	AA	789	G	N9-C4	-5.72	1.33	1.38
1	AA	2553	A	N3-C4	-5.72	1.31	1.34
1	AA	2582	G	C2-N3	-5.72	1.28	1.32
1	AA	1994	A	C6-N1	-5.72	1.31	1.35
1	AA	2400	A	N3-C4	-5.72	1.31	1.34
18	AU	15	LYS	CE-NZ	5.71	1.63	1.49
1	AA	1694	G	C2-N3	-5.71	1.28	1.32
1	AA	590	A	C5-C6	-5.71	1.35	1.41
1	AA	2454	C	N1-C2	-5.71	1.34	1.40
1	AA	254	A	N7-C5	-5.70	1.35	1.39
1	AA	2039	U	N3-C4	-5.69	1.33	1.38
1	AA	1812	C	N3-C4	5.69	1.38	1.33
1	AA	421	A	C5-C4	-5.69	1.34	1.38
1	AA	781	A	N7-C5	-5.68	1.35	1.39
1	AA	2331	G	N9-C8	5.68	1.41	1.37
1	AA	240	A	N3-C4	5.68	1.38	1.34
1	AA	46	C	C2-N3	-5.68	1.31	1.35
1	AA	608	G	N3-C4	-5.67	1.31	1.35
1	AA	1082	G	C5-C4	-5.67	1.34	1.38
1	AA	553	A	N1-C2	5.66	1.39	1.34
1	AA	979	G	N7-C5	-5.66	1.35	1.39
1	AA	73	A	N3-C4	-5.66	1.31	1.34
1	AA	1659	G	N7-C5	-5.66	1.35	1.39
56	BW	47	U	N1-C2	5.66	1.43	1.38
1	AA	492	A	C6-N6	-5.65	1.29	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	2586	G	C6-O6	-5.65	1.19	1.24
1	AA	821	A	N9-C4	-5.64	1.34	1.37
1	AA	2643	G	N7-C5	-5.64	1.35	1.39
1	AA	849	A	N3-C4	-5.64	1.31	1.34
1	AA	1305	G	C6-N1	-5.64	1.35	1.39
1	AA	1711	A	N7-C5	-5.64	1.35	1.39
1	AA	1035	G	C5-C4	-5.63	1.34	1.38
1	AA	1709	C	C4-N4	-5.63	1.28	1.33
5	AE	184	VAL	CB-CG2	-5.63	1.41	1.52
1	CA	2448	A	N9-C4	-5.63	1.34	1.37
5	AE	40	GLU	CG-CD	5.62	1.60	1.51
34	BA	782	A	N3-C4	-5.62	1.31	1.34
1	AA	517	A	N7-C5	-5.62	1.35	1.39
1	AA	354	A	N3-C4	-5.62	1.31	1.34
1	AA	1617	A	C5-C6	-5.62	1.35	1.41
1	AA	494	G	C6-N1	-5.61	1.35	1.39
1	AA	622	G	N9-C8	-5.61	1.33	1.37
1	AA	1814	A	N3-C4	-5.61	1.31	1.34
1	AA	1048	G	N1-C2	-5.61	1.33	1.37
1	AA	2701	U	C3'-O3'	5.60	1.50	1.42
1	AA	487	C	N3-C4	-5.60	1.30	1.33
1	AA	602	G	C2-N3	-5.60	1.28	1.32
1	AA	1027	A	C5-C4	-5.60	1.34	1.38
1	CA	785	G	C6-N1	-5.60	1.35	1.39
1	AA	789	G	N9-C8	-5.59	1.33	1.37
1	AA	531	G	N3-C4	-5.59	1.31	1.35
1	AA	493	G	C6-N1	-5.59	1.35	1.39
1	AA	1255	A	C8-N7	-5.59	1.27	1.31
1	AA	1661	C	C2-N3	-5.59	1.31	1.35
1	AA	188	A	C6-N1	5.58	1.39	1.35
2	AB	26	A	N7-C5	-5.58	1.35	1.39
1	AA	340	C	N3-C4	-5.58	1.30	1.33
1	AA	1834	A	N3-C4	-5.58	1.31	1.34
1	CA	2577	A	N7-C5	-5.58	1.35	1.39
1	AA	2782	C	N1-C6	-5.58	1.33	1.37
1	CA	586	A	N3-C4	-5.57	1.31	1.34
1	AA	1072	U	N3-C4	5.57	1.43	1.38
1	AA	2444	A	N7-C5	-5.57	1.35	1.39
1	AA	2620	G	C6-O6	-5.57	1.19	1.24
1	AA	1741	C	N1-C6	-5.56	1.33	1.37
1	CA	563	G	N9-C4	-5.56	1.33	1.38
1	AA	438	G	N9-C4	5.56	1.42	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	1696	G	N7-C5	-5.56	1.35	1.39
1	AA	2834	C	N3-C4	-5.55	1.30	1.33
1	AA	1745	A	N9-C8	5.54	1.42	1.37
1	AA	499	G	N1-C2	-5.54	1.33	1.37
1	AA	1033	G	N9-C8	-5.54	1.33	1.37
1	AA	2881	C	N1-C2	-5.53	1.34	1.40
1	CA	1776	G	C8-N7	-5.53	1.27	1.30
1	AA	2387	G	C5-C4	-5.53	1.34	1.38
1	AA	1067	A	N1-C2	5.53	1.39	1.34
1	AA	2281	A	N9-C4	-5.53	1.34	1.37
1	AA	2003	A	C5-C6	-5.52	1.36	1.41
1	AA	55	A	C6-N1	-5.52	1.31	1.35
1	AA	31	C	N3-C4	-5.52	1.30	1.33
1	AA	1293	A	N9-C4	-5.51	1.34	1.37
1	AA	2340	A	N7-C5	-5.51	1.35	1.39
1	AA	987	G	C5-C4	-5.51	1.34	1.38
1	AA	2545	A	N9-C4	-5.51	1.34	1.37
1	AA	2691	A	N9-C4	-5.51	1.34	1.37
1	AA	517	A	C8-N7	-5.50	1.27	1.31
1	AA	1251	G	C5-C4	-5.50	1.34	1.38
1	AA	1832	G	N7-C5	5.50	1.42	1.39
1	AA	2094	G	C5-C4	-5.49	1.34	1.38
1	AA	990	A	C2-N3	5.49	1.38	1.33
1	AA	878	G	N7-C5	5.49	1.42	1.39
1	AA	2459	G	C6-N1	-5.48	1.35	1.39
1	AA	29	U	C4-O4	-5.48	1.19	1.23
1	AA	200	A	N3-C4	-5.48	1.31	1.34
1	AA	831	A	N9-C8	-5.48	1.33	1.37
1	AA	2065	C	N1-C6	-5.48	1.33	1.37
1	AA	541	C	C2-N3	-5.48	1.31	1.35
1	AA	593	G	C5-C6	-5.48	1.36	1.42
1	AA	2024	G	C6-N1	-5.47	1.35	1.39
1	AA	1853	G	C6-N1	-5.47	1.35	1.39
34	BA	1483	A	N9-C8	-5.47	1.33	1.37
1	AA	1073	A	N9-C8	-5.47	1.33	1.37
1	AA	2470	G	N3-C4	-5.47	1.31	1.35
1	AA	113	C	N1-C6	-5.46	1.33	1.37
1	AA	1078	A	C5-C6	-5.46	1.36	1.41
1	CA	1652	A	N9-C4	-5.46	1.34	1.37
1	AA	2087	C	N3-C4	-5.46	1.30	1.33
1	AA	810	G	N3-C4	-5.45	1.31	1.35
1	AA	1824	C	N3-C4	-5.45	1.30	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	CA	2435	A	N9-C4	-5.45	1.34	1.37
1	AA	2707	C	N3-C4	-5.44	1.30	1.33
1	AA	369	A	C5-C6	-5.44	1.36	1.41
1	AA	593	G	C6-O6	-5.44	1.19	1.24
1	AA	1423	G	N1-C2	-5.44	1.33	1.37
1	AA	2298	A	N3-C4	-5.44	1.31	1.34
1	AA	2386	C	N1-C6	-5.44	1.33	1.37
1	AA	511	C	C2-O2	-5.43	1.19	1.24
1	AA	2035	A	N9-C8	-5.43	1.33	1.37
1	AA	2092	G	C5-C4	-5.43	1.34	1.38
1	AA	2546	A	N9-C4	-5.43	1.34	1.37
1	AA	775	G	C6-N1	-5.43	1.35	1.39
1	AA	853	C	C4-C5	5.42	1.47	1.43
1	CA	2437	U	C2-N3	-5.42	1.33	1.37
1	AA	606	G	C8-N7	5.42	1.34	1.30
1	AA	2014	G	C3'-C2'	5.42	1.58	1.52
1	AA	2799	U	N3-C4	-5.42	1.33	1.38
1	AA	1371	G	C6-N1	-5.41	1.35	1.39
1	AA	2881	C	N1-C6	-5.41	1.33	1.37
1	AA	1049	G	C2-N2	-5.41	1.29	1.34
1	CA	1653	G	N7-C5	-5.41	1.36	1.39
1	AA	1374	G	C2-N2	-5.41	1.29	1.34
1	AA	2620	G	C5-C6	-5.41	1.36	1.42
1	AA	1707	C	C4-C5	-5.40	1.38	1.43
1	AA	2468	C	C2-O2	5.40	1.29	1.24
5	AE	165	VAL	CB-CG1	-5.40	1.41	1.52
1	CA	1791	A	C6-N1	-5.40	1.31	1.35
1	AA	1291	G	C6-N1	-5.39	1.35	1.39
2	AB	76	G	C6-N1	-5.39	1.35	1.39
1	AA	1022	C	C4-N4	-5.39	1.29	1.33
34	BA	1397	C	C2-N3	5.39	1.40	1.35
1	AA	2639	G	C5-C6	-5.38	1.36	1.42
1	AA	792	G	C2-N3	-5.38	1.28	1.32
1	AA	1274	G	N3-C4	-5.38	1.31	1.35
1	AA	2300	A	N3-C4	5.38	1.38	1.34
1	AA	2775	G	N1-C2	-5.38	1.33	1.37
29	A5	35	GLU	CD-OE1	5.38	1.31	1.25
1	AA	1026	A	C6-N6	-5.38	1.29	1.33
1	AA	26	G	C5-C6	-5.37	1.36	1.42
1	AA	2583	C	C4-C5	-5.37	1.38	1.43
1	AA	1312	G	N7-C5	-5.37	1.36	1.39
1	AA	1911	A	N9-C4	-5.37	1.34	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	1959	A	N3-C4	-5.37	1.31	1.34
1	AA	2731	G	N3-C4	-5.37	1.31	1.35
1	CA	783	A	C6-N1	-5.37	1.31	1.35
1	CA	1698	A	N7-C5	-5.37	1.36	1.39
1	AA	2005	C	N1-C6	-5.36	1.33	1.37
1	CA	2542	A	N7-C5	-5.36	1.36	1.39
1	AA	725	C	C2-N3	-5.36	1.31	1.35
1	AA	1313	U	C2-O2	-5.36	1.17	1.22
1	AA	1665	G	C5-C4	-5.36	1.34	1.38
1	AA	1994	A	N3-C4	-5.36	1.31	1.34
5	AE	151	TYR	CB-CG	-5.36	1.43	1.51
1	AA	173	C	N1-C6	-5.35	1.33	1.37
1	AA	1048	G	C2-N3	-5.35	1.28	1.32
1	AA	787	U	N3-C4	-5.34	1.33	1.38
1	AA	2559	U	N1-C6	-5.34	1.33	1.38
1	AA	580	U	C4-O4	-5.34	1.19	1.23
1	AA	984	G	N7-C5	-5.33	1.36	1.39
6	AF	88	VAL	CB-CG1	-5.33	1.41	1.52
1	AA	355	A	C5-C4	-5.33	1.35	1.38
1	CA	330	A	C5-C6	-5.33	1.36	1.41
1	CA	1303	G	C6-N1	-5.33	1.35	1.39
1	AA	26	G	N7-C5	-5.33	1.36	1.39
1	AA	1259	A	N9-C8	-5.33	1.33	1.37
1	AA	1277	G	N3-C4	-5.33	1.31	1.35
1	AA	1804	A	N9-C4	-5.33	1.34	1.37
1	AA	2860	A	N7-C5	-5.33	1.36	1.39
1	CA	2062	A	N9-C4	5.33	1.41	1.37
1	AA	2046	G	C6-O6	-5.32	1.19	1.24
1	AA	2638	C	N1-C6	-5.32	1.33	1.37
1	AA	787	U	C2-N3	-5.32	1.34	1.37
1	AA	2449	U	C2-O2	-5.32	1.17	1.22
1	AA	2455	C	N1-C6	-5.31	1.33	1.37
34	BA	1397	C	N1-C6	5.31	1.40	1.37
1	AA	795	G	N1-C2	-5.31	1.33	1.37
1	CA	144	C	N3-C4	-5.31	1.30	1.33
1	AA	129	G	C5-C6	-5.31	1.37	1.42
1	AA	2068	G	C5-C4	-5.31	1.34	1.38
1	AA	2849	G	C6-N1	-5.31	1.35	1.39
1	CA	2286	A	C5-C6	-5.30	1.36	1.41
1	AA	956	A	N9-C8	-5.30	1.33	1.37
1	AA	1054	C	N1-C6	-5.30	1.33	1.37
1	AA	178	G	N7-C5	-5.30	1.36	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	471	C	N1-C6	-5.30	1.33	1.37
1	CA	2444	G	C2-N3	5.30	1.36	1.32
1	AA	120	G	N3-C4	-5.29	1.31	1.35
1	AA	887	C	N3-C4	-5.29	1.30	1.33
1	AA	2019	G	C6-O6	-5.29	1.19	1.24
1	CA	197	A	N9-C4	-5.29	1.34	1.37
1	AA	2295	C	N3-C4	5.29	1.37	1.33
1	AA	2041	A	C6-N1	-5.28	1.31	1.35
4	AD	39	LYS	CD-CE	5.28	1.64	1.51
1	AA	2520	G	N1-C2	-5.28	1.33	1.37
1	AA	880	U	O3'-P	-5.28	1.54	1.61
1	AA	2291	G	C8-N7	-5.28	1.27	1.30
1	AA	2611	G	N1-C2	-5.28	1.33	1.37
1	CA	563	G	C2-N3	-5.28	1.28	1.32
1	AA	2073	A	C5-C4	-5.27	1.35	1.38
1	AA	2597	U	C2-O2	5.27	1.27	1.22
1	AA	2239	A	N3-C4	-5.27	1.31	1.34
1	AA	2431	U	C2-N3	-5.26	1.34	1.37
1	AA	710	G	C6-N1	-5.26	1.35	1.39
18	AU	9	VAL	CB-CG1	-5.26	1.41	1.52
1	AA	55	A	C5-C6	-5.26	1.36	1.41
1	AA	836	A	N7-C5	-5.26	1.36	1.39
1	AA	2616	U	C2-N3	-5.26	1.34	1.37
1	AA	2525	G	C2-N3	5.25	1.36	1.32
1	CA	126	A	N3-C4	-5.25	1.31	1.34
1	CA	785	G	N1-C2	-5.25	1.33	1.37
1	AA	872	C	N3-C4	-5.25	1.30	1.33
1	CA	1204	A	N3-C4	-5.25	1.31	1.34
1	AA	1322	A	C6-N1	-5.25	1.31	1.35
1	AA	2060	G	C6-N1	-5.25	1.35	1.39
1	AA	1985	U	C2-N3	5.25	1.41	1.37
1	AA	2115	G	N3-C4	-5.25	1.31	1.35
1	AA	2440	G	C6-N1	-5.25	1.35	1.39
1	AA	808	A	N7-C5	-5.24	1.36	1.39
1	AA	1470	G	C6-N1	-5.24	1.35	1.39
1	AA	1846	A	N3-C4	-5.24	1.31	1.34
1	AA	2525	G	C6-N1	5.24	1.43	1.39
1	AA	753	A	C5-C6	-5.24	1.36	1.41
1	AA	2502	G	N1-C2	-5.24	1.33	1.37
1	AA	2402	U	C4-O4	-5.23	1.19	1.23
1	AA	2660	C	C2-O2	-5.23	1.19	1.24
1	AA	2582	G	C8-N7	-5.23	1.27	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	CA	2570	G	N3-C4	-5.23	1.31	1.35
1	AA	1240	G	C5-C4	-5.22	1.34	1.38
1	AA	2285	A	N7-C5	-5.22	1.36	1.39
1	AA	2641	A	C5-C4	5.22	1.42	1.38
1	AA	1830	G	N3-C4	-5.22	1.31	1.35
1	CA	270	A	N9-C4	-5.21	1.34	1.37
1	AA	210	A	N7-C5	-5.21	1.36	1.39
1	AA	795	G	C6-N1	-5.21	1.35	1.39
1	AA	2639	G	C6-O6	-5.21	1.19	1.24
1	AA	353	G	C8-N7	-5.20	1.27	1.30
19	AV	61	VAL	CB-CG1	-5.20	1.42	1.52
1	CA	768	G	N7-C5	-5.20	1.36	1.39
1	AA	821	A	C8-N7	-5.20	1.27	1.31
1	AA	172	C	N3-C4	-5.19	1.30	1.33
1	AA	2545	A	N9-C8	-5.19	1.33	1.37
1	AA	1833	A	N7-C5	-5.19	1.36	1.39
1	CA	1027	A	N9-C4	-5.18	1.34	1.37
1	AA	1700	G	C6-N1	-5.17	1.35	1.39
1	AA	2402	U	C2-O2	-5.17	1.17	1.22
2	AB	98	G	N9-C8	-5.17	1.34	1.37
1	AA	2460	A	N7-C5	-5.17	1.36	1.39
1	AA	2605	U	C4-O4	-5.17	1.19	1.23
1	AA	2652	G	C2-N3	-5.17	1.28	1.32
1	AA	243	G	N7-C5	-5.17	1.36	1.39
1	AA	2556	G	N9-C4	-5.17	1.33	1.38
1	AA	1076	G	C6-N1	-5.17	1.35	1.39
1	AA	992	G	P-O5'	-5.16	1.54	1.59
1	AA	1377	A	C6-N1	-5.16	1.31	1.35
1	AA	2609	G	N3-C4	-5.16	1.31	1.35
1	AA	1616	A	C6-N1	-5.16	1.31	1.35
1	CA	2542	A	C5-C6	-5.16	1.36	1.41
1	AA	1401	G	C6-N1	-5.16	1.35	1.39
1	AA	835	A	C6-N1	-5.15	1.31	1.35
1	AA	255	G	N9-C8	-5.15	1.34	1.37
1	AA	593	G	N1-C2	-5.15	1.33	1.37
1	AA	2748	G	C6-N1	-5.15	1.35	1.39
1	AA	1878	A	N9-C4	5.15	1.41	1.37
1	AA	1979	C	N1-C6	-5.15	1.34	1.37
1	AA	1375	U	C2-O2	-5.15	1.17	1.22
1	AA	2621	U	P-O5'	-5.15	1.54	1.59
29	A5	35	GLU	CD-OE2	5.15	1.31	1.25
1	AA	2283	G	N1-C2	-5.15	1.33	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	CA	2258	C	N1-C6	-5.14	1.34	1.37
1	AA	54	G	N1-C2	-5.14	1.33	1.37
1	AA	2619	G	C5-C6	-5.14	1.37	1.42
1	AA	194	G	C5-C6	-5.14	1.37	1.42
1	AA	1324	A	N3-C4	-5.14	1.31	1.34
1	AA	2239	A	C5-C4	-5.14	1.35	1.38
1	CA	2241	A	N9-C4	-5.14	1.34	1.37
1	AA	1282	G	C5-C6	-5.14	1.37	1.42
1	AA	1715	A	C6-N6	-5.13	1.29	1.33
1	AA	2084	A	N3-C4	5.13	1.38	1.34
1	AA	511	C	N1-C6	5.13	1.40	1.37
1	AA	608	G	C2-N3	-5.13	1.28	1.32
1	AA	847	A	N3-C4	-5.13	1.31	1.34
1	AA	1371	G	N1-C2	-5.13	1.33	1.37
1	AA	2463	A	C5-C4	-5.12	1.35	1.38
2	AB	98	G	N7-C5	-5.12	1.36	1.39
1	AA	2502	G	C6-N1	-5.12	1.35	1.39
11	AN	65	LYS	CE-NZ	5.12	1.61	1.49
1	CA	1890	A	C5-C6	-5.12	1.36	1.41
1	AA	522	A	N7-C5	-5.12	1.36	1.39
1	AA	2080	A	N7-C5	-5.12	1.36	1.39
1	AA	476	G	N9-C8	-5.12	1.34	1.37
1	AA	751	G	C5-C4	-5.12	1.34	1.38
1	AA	2576	A	N3-C4	-5.12	1.31	1.34
1	AA	726	C	C5-C6	-5.12	1.30	1.34
1	AA	2298	A	N9-C8	5.12	1.41	1.37
1	AA	561	A	N3-C4	-5.11	1.31	1.34
1	CA	1890	A	N9-C4	-5.11	1.34	1.37
1	AA	1723	A	N3-C4	-5.11	1.31	1.34
1	AA	1833	A	N3-C4	-5.11	1.31	1.34
1	AA	2092	G	N9-C4	-5.10	1.33	1.38
1	AA	2863	C	C2-N3	-5.10	1.31	1.35
1	AA	531	G	C5-C4	-5.09	1.34	1.38
1	AA	2057	G	N9-C8	-5.09	1.34	1.37
1	CA	180	G	N3-C4	5.09	1.39	1.35
1	AA	1600	A	N9-C8	-5.09	1.33	1.37
1	AA	340	C	C4-C5	-5.09	1.38	1.43
1	AA	1382	A	N3-C4	-5.09	1.31	1.34
1	AA	1814	A	C5-C4	-5.09	1.35	1.38
1	AA	2514	G	C6-N1	-5.09	1.35	1.39
1	CA	679	C	N1-C2	-5.08	1.35	1.40
1	AA	1067	A	C2-N3	-5.08	1.28	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	805	C	N1-C6	-5.08	1.34	1.37
1	AA	1833	A	N9-C4	-5.08	1.34	1.37
1	AA	1974	A	C5-C6	-5.08	1.36	1.41
1	AA	586	G	C2-N3	-5.08	1.28	1.32
1	AA	1260	G	N7-C5	-5.08	1.36	1.39
1	AA	2400	A	C6-N1	-5.08	1.31	1.35
1	AA	28	A	C2-N3	-5.07	1.28	1.33
1	AA	736	A	N9-C8	-5.07	1.33	1.37
1	AA	893	C	N3-C4	-5.07	1.30	1.33
1	AA	730	C	N1-C6	-5.06	1.34	1.37
1	AA	846	G	N7-C5	-5.06	1.36	1.39
1	AA	2024	G	C8-N7	-5.06	1.27	1.30
1	AA	1029	A	N9-C4	-5.06	1.34	1.37
1	AA	2409	G	N1-C2	-5.06	1.33	1.37
1	AA	2487	C	N1-C6	-5.06	1.34	1.37
1	AA	474	U	N3-C4	-5.06	1.33	1.38
1	AA	585	U	C4-O4	-5.06	1.19	1.23
1	AA	1309	U	C2-N3	-5.06	1.34	1.37
1	AA	1546	G	C5-C4	-5.06	1.34	1.38
1	AA	2569	G	C5-C6	-5.06	1.37	1.42
1	CA	2576	G	C5-C4	-5.06	1.34	1.38
1	AA	2576	A	C6-N1	-5.06	1.32	1.35
1	AA	1201	A	N9-C4	-5.05	1.34	1.37
1	AA	2070	G	N1-C2	-5.05	1.33	1.37
1	AA	1243	U	C2-O2	-5.05	1.17	1.22
1	AA	1648	U	C4-O4	-5.05	1.19	1.23
1	AA	2637	G	N9-C8	-5.05	1.34	1.37
1	AA	1742	G	C5-C6	-5.05	1.37	1.42
1	AA	1829	U	C2-N3	-5.05	1.34	1.37
1	AA	2556	G	N3-C4	-5.05	1.31	1.35
1	AA	883	G	N1-C2	-5.05	1.33	1.37
1	AA	894	U	C4-C5	5.05	1.48	1.43
1	AA	2011	G	C6-O6	5.05	1.28	1.24
1	AA	641	G	C6-N1	-5.04	1.36	1.39
1	AA	1455	C	C2-O2	-5.04	1.20	1.24
1	AA	1857	G	N3-C4	-5.04	1.31	1.35
1	CA	1945	G	N7-C5	-5.04	1.36	1.39
1	CA	2581	G	N7-C5	-5.04	1.36	1.39
1	AA	2033	U	N3-C4	-5.04	1.33	1.38
1	AA	851	A	N3-C4	-5.04	1.31	1.34
1	AA	2528	G	C2-N2	-5.04	1.29	1.34
1	AA	1270	C	C5-C6	-5.04	1.30	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	1824	C	C4-N4	-5.04	1.29	1.33
1	AA	585	U	C2-O2	-5.03	1.17	1.22
1	AA	1334	U	P-OP2	-5.03	1.40	1.49
1	AA	2039	U	C2-N3	-5.03	1.34	1.37
1	CA	801	G	N7-C5	-5.03	1.36	1.39
1	AA	186	A	C6-N1	-5.03	1.32	1.35
1	AA	474	U	C2-O2	-5.03	1.17	1.22
1	AA	1240	G	C2-N3	-5.03	1.28	1.32
1	AA	1684	A	N9-C4	-5.02	1.34	1.37
1	AA	1321	A	C6-N1	-5.02	1.32	1.35
1	AA	27	G	P-O5'	-5.02	1.54	1.59
1	AA	584	G	C5-C4	-5.02	1.34	1.38
1	AA	2902	G	N9-C8	5.02	1.41	1.37
34	DA	295	C	N1-C6	-5.02	1.34	1.37
1	AA	1658	C	C4-N4	-5.01	1.29	1.33
2	AB	82	G	C6-N1	-5.01	1.36	1.39
1	AA	1359	U	N1-C6	-5.01	1.33	1.38
1	AA	2041	A	N7-C5	-5.01	1.36	1.39
1	CA	2556	C	N1-C6	-5.01	1.34	1.37
1	AA	1822	A	N9-C4	-5.01	1.34	1.37
1	AA	2084	A	C6-N1	5.01	1.39	1.35
1	AA	648	G	N1-C2	-5.01	1.33	1.37
1	AA	1231	G	N3-C4	5.01	1.39	1.35
1	AA	211	A	C5-C4	-5.00	1.35	1.38
1	AA	472	G	N7-C5	5.00	1.42	1.39
1	CA	330	A	N9-C8	5.00	1.41	1.37

All (7729) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1067	A	C2-N3-C4	-25.69	97.76	110.60
1	AA	1701	A	O5'-P-OP2	-25.21	80.44	110.70
1	AA	553	A	N1-C6-N6	25.19	133.71	118.60
1	AA	990	A	C5-N7-C8	-25.10	91.35	103.90
1	AA	553	A	C5-N7-C8	-23.43	92.19	103.90
1	AA	990	A	N1-C6-N6	22.68	132.21	118.60
1	AA	553	A	C6-C5-N7	-21.21	117.45	132.30
1	AA	354	A	C2-N3-C4	-21.08	100.06	110.60
1	AA	474	U	O5'-P-OP2	-21.05	85.44	110.70
1	AA	990	A	C6-C5-N7	-20.78	117.75	132.30
1	AA	553	A	N7-C8-N9	20.56	124.08	113.80
1	AA	1067	A	N3-C4-C5	20.20	140.94	126.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1605	A	C2-N3-C4	-19.70	100.75	110.60
1	AA	990	A	C2-N3-C4	-19.47	100.86	110.60
1	AA	553	A	C4-C5-N7	19.43	120.42	110.70
1	AA	990	A	C4-C5-N7	19.21	120.30	110.70
1	AA	1188	A	N3-C4-C5	19.00	140.10	126.80
1	AA	1807	G	O5'-P-OP2	-18.88	88.05	110.70
1	AA	2331	G	N3-C4-N9	-18.76	114.75	126.00
1	AA	553	A	C2-N3-C4	-18.63	101.28	110.60
1	AA	978	A	C5-N7-C8	-18.59	94.61	103.90
1	AA	990	A	N7-C8-N9	18.58	123.09	113.80
1	AA	1745	A	N1-C6-N6	18.46	129.68	118.60
1	AA	2299	A	C2-N3-C4	-18.38	101.41	110.60
1	AA	2045	G	O5'-P-OP1	-18.35	88.68	110.70
1	AA	2083	G	O5'-P-OP2	-18.25	88.80	110.70
1	AA	1811	A	O5'-P-OP2	-17.61	89.56	110.70
1	AA	254	A	C2-N3-C4	-17.57	101.82	110.60
1	AA	1188	A	C2-N3-C4	-17.51	101.84	110.60
1	AA	1067	A	N3-C4-N9	-17.42	113.46	127.40
1	AA	1605	A	N1-C2-N3	17.39	138.00	129.30
1	AA	2222	C	O5'-P-OP2	-17.26	89.99	110.70
1	AA	592	U	C5-C6-N1	-17.21	114.10	122.70
1	AA	1707	C	O5'-P-OP2	-17.14	90.13	110.70
1	AA	553	A	C8-N9-C4	-17.03	98.99	105.80
1	CA	1353	A	O5'-P-OP2	-16.66	90.71	105.70
1	AA	991	G	O5'-P-OP1	-16.59	90.77	105.70
1	AA	1188	A	N3-C4-N9	-16.57	114.14	127.40
1	AA	1745	A	C6-C5-N7	-16.48	120.76	132.30
1	AA	1067	A	C5-C6-N1	-16.45	109.47	117.70
1	AA	2045	G	C5-C6-O6	-16.34	118.80	128.60
1	AA	354	A	C5-N7-C8	-16.31	95.74	103.90
1	AA	2298	A	C2-N3-C4	-16.16	102.52	110.60
1	AA	978	A	C4-C5-N7	16.14	118.77	110.70
1	AA	555	G	N3-C4-N9	-16.14	116.31	126.00
1	AA	354	A	N3-C4-C5	16.10	138.07	126.80
1	AA	2418	U	O5'-P-OP1	-15.94	91.36	105.70
1	AA	1067	A	C5-N7-C8	-15.75	96.03	103.90
1	AA	1249	A	C6-C5-N7	-15.61	121.37	132.30
1	AA	254	A	N1-C6-N6	15.58	127.95	118.60
1	AA	1813	C	O5'-P-OP1	-15.46	91.78	105.70
1	AA	2331	G	N3-C4-C5	15.45	136.32	128.60
1	CA	2608	G	O5'-P-OP2	-15.30	91.93	105.70
1	AA	1745	A	C5-N7-C8	-15.21	96.30	103.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	528	A	C2-N3-C4	-15.14	103.03	110.60
1	AA	2298	A	N7-C8-N9	15.13	121.36	113.80
1	AA	1745	A	C4-C5-N7	15.05	118.22	110.70
1	AA	2619	G	C5-C6-O6	-14.92	119.65	128.60
1	AA	1249	A	C2-N3-C4	-14.82	103.19	110.60
1	AA	612	C	O5'-P-OP2	-14.81	92.37	105.70
1	AA	555	G	N3-C4-C5	14.65	135.93	128.60
1	AA	1812	C	C6-N1-C2	14.64	126.16	120.30
1	AA	1486	G	O5'-P-OP2	-14.55	92.60	105.70
34	BA	797	C	O5'-P-OP1	-14.55	92.60	105.70
1	AA	1686	U	O5'-P-OP2	-14.55	92.61	105.70
1	AA	990	A	N1-C2-N3	14.40	136.50	129.30
1	AA	2298	A	C5-N7-C8	-14.30	96.75	103.90
1	AA	1694	G	O5'-P-OP1	-14.28	92.85	105.70
1	CA	1998	G	O5'-P-OP2	-14.22	92.90	105.70
1	AA	1155	C	C5-C6-N1	14.20	128.10	121.00
1	AA	2065	C	O5'-P-OP1	-14.14	92.97	105.70
1	AA	990	A	C5-C6-N6	-14.14	112.39	123.70
1	AA	560	C	C5-C6-N1	-14.08	113.96	121.00
1	AA	2556	G	C5-C6-O6	-14.04	120.17	128.60
1	AA	1815	A	O5'-P-OP2	-14.03	93.07	105.70
1	CA	330	A	C2-N3-C4	-14.00	103.60	110.60
1	AA	204	G	O5'-P-OP1	-13.98	93.11	105.70
1	AA	1266	C	C6-N1-C2	-13.98	114.71	120.30
1	AA	2298	A	C6-C5-N7	-13.82	122.63	132.30
1	AA	1274	G	C2-N3-C4	-13.79	105.00	111.90
1	AA	354	A	N3-C4-N9	-13.72	116.42	127.40
1	AA	2515	A	N1-C2-N3	-13.66	122.47	129.30
1	AA	1249	A	N1-C6-N6	13.63	126.78	118.60
1	AA	2632	C	N1-C2-O2	-13.61	110.73	118.90
1	AA	553	A	N1-C2-N3	13.47	136.04	129.30
1	CA	1204	A	C2-N3-C4	-13.39	103.91	110.60
1	AA	2285	A	O5'-P-OP2	-13.38	93.66	105.70
1	AA	978	A	C2-N3-C4	-13.36	103.92	110.60
1	AA	553	A	O4'-C1'-N9	-13.24	97.61	108.20
1	AA	1612	C	O5'-P-OP2	-13.18	93.84	105.70
1	AA	1847	G	O5'-P-OP1	-13.11	93.90	105.70
1	AA	543	G	O5'-P-OP2	-13.09	93.92	105.70
1	AA	254	A	C5-N7-C8	-13.08	97.36	103.90
1	AA	18	C	O5'-P-OP2	-13.04	93.97	105.70
1	AA	1001	G	N1-C6-O6	13.03	127.72	119.90
1	CA	1204	A	N1-C6-N6	12.99	126.40	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2636	G	N7-C8-N9	-12.99	106.60	113.10
1	AA	2641	A	C5-N7-C8	-12.98	97.41	103.90
1	AA	553	A	C5-C6-N6	-12.97	113.33	123.70
1	AA	1472	G	C5-C6-O6	-12.97	120.82	128.60
1	AA	2269	U	C2-N3-C4	-12.93	119.24	127.00
1	AA	978	A	N1-C6-N6	12.90	126.34	118.60
1	AA	2298	A	N1-C6-N6	12.90	126.34	118.60
1	AA	1342	G	N1-C6-O6	-12.88	112.17	119.90
1	AA	1745	A	C2-N3-C4	-12.88	104.16	110.60
1	AA	354	A	C4-C5-N7	12.86	117.13	110.70
1	AA	539	A	O5'-P-OP2	-12.86	94.13	105.70
1	AA	894	U	N1-C2-N3	12.85	122.61	114.90
1	AA	1188	A	C6-N1-C2	12.85	126.31	118.60
1	AA	1188	A	C5-C6-N1	-12.85	111.28	117.70
1	AA	1249	A	N1-C2-N3	12.85	135.72	129.30
1	AA	2400	A	N9-C4-C5	12.80	110.92	105.80
1	AA	2627	U	O5'-P-OP1	-12.76	94.22	105.70
1	AA	978	A	N7-C8-N9	12.72	120.16	113.80
1	AA	2641	A	C6-C5-N7	-12.72	123.39	132.30
1	AA	812	G	O5'-P-OP2	-12.71	94.26	105.70
1	AA	2299	A	C5-N7-C8	-12.69	97.55	103.90
1	AA	2298	A	N1-C2-N3	12.63	135.62	129.30
1	AA	1605	A	C6-C5-N7	-12.62	123.46	132.30
1	AA	2368	C	C5-C4-N4	-12.62	111.37	120.20
1	AA	1032	C	N1-C2-O2	-12.62	111.33	118.90
1	AA	1543	U	C5-C4-O4	12.60	133.46	125.90
1	AA	2641	A	N1-C6-N6	12.59	126.15	118.60
1	AA	113	C	C6-N1-C2	12.57	125.33	120.30
1	AA	726	C	C2-N3-C4	-12.57	113.62	119.90
1	AA	2299	A	N3-C4-C5	12.54	135.58	126.80
1	AA	1926	G	N1-C6-O6	-12.54	112.38	119.90
1	AA	2539	C	C6-N1-C2	12.51	125.30	120.30
1	CA	512	G	O4'-C1'-N9	12.49	118.19	108.20
1	AA	989	G	C5-C6-O6	-12.46	121.12	128.60
1	AA	2452	C	C6-N1-C2	12.46	125.28	120.30
1	AA	2858	G	O5'-P-OP2	-12.44	94.50	105.70
1	AA	2636	G	C8-N9-C4	12.41	111.37	106.40
1	AA	2045	G	C4-C5-N7	12.41	115.77	110.80
1	AA	2641	A	C4-C5-N7	12.37	116.89	110.70
1	AA	1211	U	C2-N3-C4	-12.37	119.58	127.00
1	AA	1709	C	C2-N3-C4	-12.35	113.72	119.90
1	AA	2566	U	O5'-P-OP1	-12.34	94.59	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1422	C	O5'-P-OP1	-12.32	94.61	105.70
1	AA	2876	U	C5-C4-O4	12.29	133.27	125.90
1	AA	992	G	O5'-P-OP1	-12.26	94.67	105.70
1	AA	2298	A	C8-N9-C4	-12.25	100.90	105.80
1	AA	1429	C	O5'-P-OP1	-12.24	94.69	105.70
1	AA	884	C	C5-C6-N1	-12.23	114.88	121.00
1	CA	528	A	N3-C4-N9	-12.20	117.64	127.40
1	AA	840	A	O5'-P-OP2	-12.17	94.75	105.70
1	AA	1249	A	C4-C5-N7	12.16	116.78	110.70
1	AA	1605	A	N1-C6-N6	12.16	125.90	118.60
1	AA	2107	C	N3-C4-C5	12.13	126.75	121.90
1	CA	2435	A	O5'-P-OP1	-12.12	94.80	105.70
1	AA	978	A	N3-C4-C5	12.10	135.27	126.80
1	AA	1824	C	N3-C4-C5	12.09	126.74	121.90
1	AA	709	G	O5'-P-OP1	-12.08	94.83	105.70
1	AA	2015	U	O5'-P-OP1	-12.08	94.83	105.70
1	AA	1188	A	C5-N7-C8	-12.08	97.86	103.90
2	AB	73	A	O5'-P-OP2	-12.07	94.84	105.70
1	AA	1329	G	C6-C5-N7	-12.05	123.17	130.40
34	BA	1521	G	O5'-P-OP1	-12.05	94.86	105.70
1	AA	593	G	C5-C6-O6	-12.05	121.37	128.60
1	AA	2532	C	N1-C2-O2	-12.04	111.67	118.90
1	AA	2065	C	O5'-P-OP2	12.02	125.13	110.70
1	AA	1185	C	O5'-P-OP1	-12.02	94.89	105.70
1	AA	355	A	O5'-P-OP1	-11.98	94.92	105.70
1	AA	555	G	C5-C6-O6	11.97	135.78	128.60
1	AA	2426	G	O5'-P-OP1	11.97	125.07	110.70
1	AA	2556	G	N1-C6-O6	11.96	127.08	119.90
1	AA	254	A	C4-C5-N7	11.96	116.68	110.70
1	AA	139	A	C5-N7-C8	-11.95	97.92	103.90
1	AA	733	G	O5'-P-OP1	-11.94	94.95	105.70
1	AA	2053	A	O5'-P-OP1	-11.91	94.98	105.70
1	AA	325	G	C8-N9-C4	11.89	111.16	106.40
1	CA	330	A	C5-N7-C8	-11.88	97.96	103.90
1	AA	1249	A	C5-N7-C8	-11.85	97.97	103.90
1	AA	805	C	C6-N1-C2	11.81	125.03	120.30
1	AA	2331	G	N3-C2-N2	-11.80	111.64	119.90
1	AA	2550	C	C6-N1-C2	11.80	125.02	120.30
1	AA	2578	A	O5'-P-OP2	-11.79	95.09	105.70
1	AA	2367	C	N3-C4-C5	11.78	126.61	121.90
1	AA	2660	C	C2-N3-C4	-11.78	114.01	119.90
1	AA	1961	U	C4-C5-C6	-11.77	112.64	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1067	A	C4-C5-N7	11.75	116.58	110.70
1	AA	38	A	N1-C6-N6	-11.74	111.56	118.60
1	AA	1721	G	C6-N1-C2	-11.74	118.06	125.10
1	AA	1001	G	C5-C6-N1	-11.69	105.65	111.50
1	AA	641	G	O5'-P-OP2	-11.67	95.20	105.70
1	AA	2386	C	C6-N1-C2	11.66	124.96	120.30
1	AA	2331	G	C8-N9-C4	-11.63	101.75	106.40
1	CA	2193	G	OP1-P-O3'	-11.60	79.68	105.20
1	AA	126	C	O5'-P-OP1	-11.59	95.27	105.70
1	AA	1742	G	O5'-P-OP1	-11.55	95.30	105.70
1	AA	1691	C	N1-C2-O2	11.55	125.83	118.90
1	AA	254	A	C6-C5-N7	-11.53	124.23	132.30
1	AA	1745	A	C5-C6-N6	-11.52	114.48	123.70
1	AA	1745	A	N1-C2-N3	11.51	135.06	129.30
1	CA	763	G	O5'-P-OP1	-11.51	95.34	105.70
1	AA	716	G	C5-C6-O6	-11.45	121.73	128.60
1	AA	1694	G	N3-C2-N2	-11.44	111.89	119.90
1	CA	528	A	N3-C4-C5	11.43	134.80	126.80
1	CA	1673	U	O5'-P-OP1	-11.43	95.42	105.70
1	AA	2639	G	C8-N9-C4	11.42	110.97	106.40
1	AA	1709	C	N3-C4-C5	11.40	126.46	121.90
1	AA	2641	A	C2-N3-C4	-11.37	104.91	110.60
1	AA	1472	G	C4-C5-N7	11.37	115.35	110.80
34	DA	354	G	O5'-P-OP2	-11.35	95.49	105.70
1	AA	894	U	C5-C6-N1	-11.33	117.03	122.70
1	CA	1698	A	N1-C2-N3	11.31	134.96	129.30
1	AA	2018	C	C2-N3-C4	-11.30	114.25	119.90
1	AA	45	C	O5'-P-OP1	-11.29	95.54	105.70
1	AA	2018	C	N3-C4-C5	11.27	126.41	121.90
1	AA	618	C	O5'-P-OP2	-11.26	95.56	105.70
1	AA	1720	U	C5-C6-N1	-11.23	117.09	122.70
1	AA	1244	U	O5'-P-OP2	-11.21	95.61	105.70
1	CA	1269	A	O5'-P-OP2	-11.21	95.61	105.70
1	AA	1721	G	C5-C6-O6	-11.20	121.88	128.60
1	AA	2299	A	C5-C6-N1	-11.19	112.10	117.70
1	AA	2447	A	O5'-P-OP1	-11.17	95.64	105.70
1	AA	471	C	N1-C2-O2	-11.17	112.20	118.90
1	AA	2331	G	C5-N7-C8	-11.16	98.72	104.30
1	AA	1282	G	C8-N9-C4	11.12	110.85	106.40
1	AA	1080	G	OP1-P-OP2	11.12	136.28	119.60
1	CA	1362	C	C6-N1-C2	-11.08	115.87	120.30
1	AA	2876	U	C4-C5-C6	11.08	126.34	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	254	A	C5-C6-N1	-11.07	112.16	117.70
1	AA	847	A	N1-C6-N6	-11.06	111.96	118.60
1	AA	552	C	C5-C6-N1	-11.06	115.47	121.00
1	AA	2876	U	N1-C2-N3	11.06	121.54	114.90
1	CA	679	C	N1-C2-O2	-11.05	112.27	118.90
1	AA	139	A	N7-C8-N9	11.04	119.32	113.80
1	CA	1558	A	C2-N3-C4	-11.03	105.08	110.60
1	AA	438	G	N3-C4-C5	-11.02	123.09	128.60
1	AA	470	C	O5'-P-OP1	11.02	123.92	110.70
1	AA	2876	U	C5-C6-N1	-11.00	117.20	122.70
1	AA	2397	C	N3-C4-C5	10.97	126.29	121.90
1	AA	1067	A	C6-N1-C2	10.97	125.18	118.60
1	AA	2251	G	O5'-P-OP1	-10.96	95.84	105.70
1	AA	978	A	C5-C6-N1	-10.95	112.22	117.70
1	CA	2573	C	O5'-P-OP2	-10.95	95.84	105.70
1	AA	1371	G	O4'-C1'-N9	10.91	116.93	108.20
34	DA	1045	C	OP1-P-O3'	-10.89	81.23	105.20
1	AA	2426	G	O5'-P-OP2	-10.89	95.90	105.70
1	AA	1001	G	C2-N3-C4	-10.89	106.46	111.90
1	AA	1074	A	C5-C6-N6	-10.88	115.00	123.70
1	AA	726	C	C5-C4-N4	-10.86	112.60	120.20
1	AA	2045	G	N1-C6-O6	10.84	126.41	119.90
1	AA	2609	G	C5-C6-O6	-10.84	122.09	128.60
1	CA	330	A	C4-C5-N7	10.84	116.12	110.70
1	AA	1995	G	O5'-P-OP2	-10.82	95.96	105.70
1	CA	2261	C	O5'-P-OP2	-10.82	95.96	105.70
1	AA	2367	C	C2-N3-C4	-10.81	114.49	119.90
1	AA	1724	A	N1-C2-N3	10.81	134.70	129.30
1	AA	16	G	N1-C2-N3	10.80	130.38	123.90
1	AA	842	C	N3-C2-O2	-10.80	114.34	121.90
1	AA	537	G	O4'-C1'-N9	10.79	116.83	108.20
1	AA	552	C	C4-C5-C6	10.79	122.79	117.40
1	CA	1653	G	C8-N9-C4	-10.79	102.08	106.40
1	AA	1721	G	C5-C6-N1	10.78	116.89	111.50
1	AA	1013	G	N1-C6-O6	-10.77	113.44	119.90
1	AA	2400	A	N1-C6-N6	-10.77	112.14	118.60
1	AA	786	G	C5-C6-N1	10.76	116.88	111.50
1	AA	1237	G	C5-N7-C8	10.76	109.68	104.30
1	AA	614	C	N1-C2-O2	-10.75	112.45	118.90
1	CA	2286	A	N1-C6-N6	10.74	125.05	118.60
1	AA	978	A	C6-C5-N7	-10.74	124.78	132.30
1	AA	555	G	C2-N3-C4	-10.73	106.54	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1411	A	N1-C6-N6	10.72	125.03	118.60
34	DA	438	G	O5'-P-OP2	-10.71	96.06	105.70
1	AA	2019	G	O5'-P-OP2	-10.71	96.06	105.70
1	CA	1790	C	C6-N1-C2	-10.71	116.02	120.30
1	CA	2588	G	C8-N9-C4	10.70	110.68	106.40
1	AA	50	G	O5'-P-OP2	-10.70	96.07	105.70
1	AA	552	C	C2-N3-C4	-10.70	114.55	119.90
1	CA	1272	A	O5'-P-OP2	-10.69	96.08	105.70
1	AA	586	G	C5-C6-N1	-10.68	106.16	111.50
1	CA	1204	A	C5-N7-C8	-10.68	98.56	103.90
1	CA	2824	C	C6-N1-C2	10.68	124.57	120.30
1	CA	1698	A	C6-C5-N7	-10.67	124.83	132.30
1	CA	2004	G	O5'-P-OP2	-10.65	96.11	105.70
1	CA	450	G	N1-C6-O6	-10.63	113.52	119.90
1	AA	2295	C	O5'-P-OP2	-10.63	96.14	105.70
1	CA	330	A	N3-C4-C5	10.63	134.24	126.80
1	AA	2620	G	C5-C6-O6	-10.60	122.24	128.60
1	AA	884	C	C2-N3-C4	-10.60	114.60	119.90
1	AA	2884	C	N1-C2-O2	-10.59	112.55	118.90
1	AA	844	C	N1-C2-O2	-10.58	112.55	118.90
1	AA	2452	C	C5-C6-N1	-10.58	115.71	121.00
1	AA	1371	G	N1-C6-O6	-10.55	113.57	119.90
1	AA	2298	A	C5-C6-N1	-10.55	112.43	117.70
2	AB	82	G	C5-C6-O6	10.54	134.93	128.60
1	AA	1317	G	C5-C6-O6	-10.54	122.28	128.60
1	AA	357	G	C8-N9-C4	-10.53	102.19	106.40
34	DA	898	G	C5-C6-O6	-10.53	122.28	128.60
1	AA	2702	C	N1-C2-O2	-10.53	112.58	118.90
1	AA	2066	C	C2-N3-C4	-10.52	114.64	119.90
1	AA	1050	C	N1-C2-N3	10.52	126.56	119.20
1	AA	1189	A	C8-N9-C4	-10.52	101.59	105.80
1	AA	1239	A	N9-C4-C5	10.51	110.00	105.80
1	AA	1360	C	C2-N3-C4	-10.50	114.65	119.90
1	AA	1040	C	C6-N1-C2	10.49	124.50	120.30
1	CA	1941	C	O5'-P-OP1	-10.49	96.26	105.70
1	AA	2368	C	N3-C4-N4	10.48	125.34	118.00
1	AA	2520	G	O5'-P-OP1	-10.48	96.27	105.70
1	AA	1042	A	O5'-P-OP2	10.47	123.27	110.70
1	AA	1720	U	C2-N3-C4	-10.47	120.72	127.00
1	AA	2750	G	N1-C6-O6	-10.47	113.62	119.90
1	AA	1186	U	C2-N3-C4	-10.46	120.72	127.00
1	AA	591	U	N3-C4-C5	10.46	120.88	114.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	524	U	C5-C6-N1	-10.46	117.47	122.70
1	AA	1020	C	C5-C4-N4	10.46	127.52	120.20
1	AA	1186	U	N1-C2-N3	10.46	121.17	114.90
1	AA	2269	U	C5-C6-N1	-10.45	117.47	122.70
34	BA	816	A	O5'-P-OP1	10.42	123.21	110.70
1	AA	2269	U	N3-C4-C5	10.42	120.85	114.60
1	AA	2757	G	O5'-P-OP1	-10.39	96.35	105.70
1	CA	2234	G	C6-C5-N7	-10.39	124.17	130.40
2	AB	103	G	C2-N3-C4	-10.37	106.72	111.90
1	AA	181	C	C4-C5-C6	10.37	122.58	117.40
34	BA	817	C	O5'-P-OP1	-10.36	96.37	105.70
1	AA	2402	U	C5-C6-N1	-10.36	117.52	122.70
1	AA	2639	G	N9-C4-C5	-10.36	101.26	105.40
1	AA	1239	A	C8-N9-C4	-10.36	101.66	105.80
1	AA	847	A	N9-C4-C5	10.35	109.94	105.80
1	AA	2041	A	C2-N3-C4	-10.34	105.43	110.60
1	AA	496	A	O5'-P-OP1	-10.34	96.40	105.70
1	AA	719	C	O5'-P-OP2	-10.34	96.40	105.70
1	AA	2602	A	C8-N9-C4	10.33	109.93	105.80
34	BA	567	G	O5'-P-OP1	-10.33	96.40	105.70
1	CA	262	A	O5'-P-OP2	-10.33	96.40	105.70
34	BA	509	A	C8-N9-C4	-10.33	101.67	105.80
1	CA	2544	G	N1-C6-O6	10.30	126.08	119.90
1	AA	1082	G	N7-C8-N9	-10.30	107.95	113.10
1	CA	562	U	N3-C2-O2	-10.30	114.99	122.20
1	AA	1332	A	O5'-P-OP2	-10.28	96.45	105.70
1	AA	1397	C	N3-C4-N4	-10.28	110.80	118.00
1	AA	1189	A	N1-C2-N3	10.27	134.44	129.30
1	AA	841	G	C8-N9-C4	10.27	110.51	106.40
1	AA	1237	G	N7-C8-N9	-10.27	107.97	113.10
1	AA	980	C	C2-N3-C4	-10.26	114.77	119.90
1	AA	2063	U	N3-C2-O2	-10.26	115.02	122.20
1	AA	412	C	N1-C2-O2	-10.25	112.75	118.90
1	AA	2299	A	N3-C4-N9	-10.25	119.20	127.40
1	AA	553	A	C4-C5-C6	10.24	122.12	117.00
1	AA	1318	A	O5'-P-OP2	-10.24	96.49	105.70
34	BA	1502	A	C5-N7-C8	-10.24	98.78	103.90
1	AA	2505	U	N1-C2-N3	10.24	121.04	114.90
1	AA	1423	G	N3-C4-C5	-10.22	123.49	128.60
1	AA	2460	A	C6-N1-C2	-10.22	112.47	118.60
1	AA	1691	C	N3-C2-O2	-10.21	114.75	121.90
1	AA	975	U	N1-C2-O2	-10.21	115.65	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	792	G	O5'-P-OP2	-10.21	96.52	105.70
1	AA	1371	G	C2-N3-C4	10.20	117.00	111.90
1	AA	1458	A	O5'-P-OP2	10.20	122.94	110.70
1	AA	743	G	C8-N9-C4	10.19	110.47	106.40
1	AA	1067	A	N1-C6-N6	10.18	124.71	118.60
34	DA	900	A	O5'-P-OP1	-10.17	96.54	105.70
1	AA	1745	A	N7-C8-N9	10.13	118.86	113.80
1	AA	1342	G	N3-C2-N2	10.12	126.99	119.90
1	AA	832	G	C5-C6-O6	10.11	134.66	128.60
1	CA	2583	G	O5'-P-OP2	-10.10	96.61	105.70
1	CA	2080	G	O5'-P-OP2	-10.09	96.62	105.70
1	AA	2641	A	N1-C2-N3	10.09	134.34	129.30
1	AA	2775	G	C5-C6-O6	10.08	134.65	128.60
1	CA	1021	A	C2-N3-C4	-10.08	105.56	110.60
1	AA	2455	C	C2-N3-C4	-10.08	114.86	119.90
1	AA	1297	C	O5'-P-OP2	-10.07	96.64	105.70
1	AA	2019	G	C6-N1-C2	-10.07	119.06	125.10
18	AU	50	ARG	NE-CZ-NH1	-10.07	115.27	120.30
1	AA	1377	A	N1-C6-N6	-10.06	112.56	118.60
1	AA	2608	U	C2-N3-C4	-10.06	120.96	127.00
1	CA	1623	G	N3-C2-N2	-10.06	112.86	119.90
1	AA	2518	U	OP2-P-O3'	10.06	127.32	105.20
1	AA	2737	C	N3-C4-C5	-10.04	117.88	121.90
1	AA	254	A	N3-C4-C5	10.04	133.83	126.80
1	AA	2791	A	C2-N3-C4	-10.04	105.58	110.60
1	CA	205	G	C8-N9-C4	10.04	110.42	106.40
1	AA	1076	G	C5-C6-O6	-10.04	122.58	128.60
1	AA	2515	A	C8-N9-C4	10.03	109.81	105.80
1	AA	738	C	N3-C2-O2	-10.03	114.88	121.90
1	AA	778	C	C2-N3-C4	-10.02	114.89	119.90
1	AA	1033	G	OP1-P-OP2	-10.01	104.59	119.60
1	AA	904	C	O5'-P-OP1	-9.98	96.71	105.70
1	CA	2063	C	O5'-P-OP2	-9.96	96.73	105.70
1	AA	592	U	C6-N1-C2	9.96	126.97	121.00
1	AA	1048	G	O5'-P-OP2	-9.96	96.74	105.70
1	AA	1316	C	C6-N1-C2	9.96	124.28	120.30
1	AA	2497	G	C8-N9-C4	9.95	110.38	106.40
1	AA	1188	A	C4-C5-C6	-9.94	112.03	117.00
1	AA	2466	G	N1-C2-N2	-9.93	107.26	116.20
1	AA	139	A	C2-N3-C4	-9.93	105.64	110.60
1	AA	1018	A	N7-C8-N9	9.92	118.76	113.80
1	AA	2481	A	O5'-P-OP2	-9.92	96.77	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2053	A	O5'-P-OP2	9.92	122.60	110.70
1	AA	2579	G	N1-C6-O6	-9.92	113.95	119.90
1	AA	524	U	C5-C4-O4	9.91	131.85	125.90
1	AA	2693	C	C4-C5-C6	9.90	122.35	117.40
1	AA	2526	U	C5-C4-O4	9.90	131.84	125.90
1	CA	2503	A	N1-C2-N3	-9.89	124.35	129.30
1	AA	2559	U	N3-C4-O4	9.89	126.32	119.40
1	AA	990	A	C4-C5-C6	9.89	121.94	117.00
1	AA	1249	A	O4'-C1'-N9	9.88	116.10	108.20
1	AA	2048	C	O5'-P-OP2	-9.88	96.81	105.70
1	AA	2070	G	N3-C2-N2	9.88	126.81	119.90
2	AB	73	A	O5'-P-OP1	9.87	122.55	110.70
1	AA	82	G	C8-N9-C4	9.87	110.35	106.40
1	AA	986	A	O5'-P-OP1	-9.87	96.82	105.70
1	AA	2619	G	N1-C6-O6	9.87	125.82	119.90
1	AA	2641	A	N7-C8-N9	9.86	118.73	113.80
1	AA	2514	G	N7-C8-N9	-9.86	108.17	113.10
1	AA	1972	G	O5'-P-OP1	-9.86	96.83	105.70
1	AA	2298	A	C4-C5-N7	9.86	115.63	110.70
1	AA	796	C	C2-N3-C4	-9.85	114.98	119.90
2	AB	102	A	N1-C2-N3	9.85	134.22	129.30
1	AA	438	G	O5'-P-OP1	9.84	122.51	110.70
34	BA	756	C	C6-N1-C2	9.84	124.23	120.30
1	AA	832	G	N1-C6-O6	-9.83	114.00	119.90
1	CA	2618	G	O5'-P-OP2	-9.83	96.85	105.70
2	AB	82	G	N1-C6-O6	-9.83	114.00	119.90
1	AA	2383	G	C5-C6-N1	9.82	116.41	111.50
1	AA	2515	A	N9-C4-C5	-9.82	101.87	105.80
1	AA	73	A	N1-C2-N3	9.82	134.21	129.30
1	AA	1031	C	C6-N1-C2	-9.82	116.37	120.30
18	AU	50	ARG	NE-CZ-NH2	9.82	125.21	120.30
1	CA	945	A	C2-N3-C4	-9.81	105.70	110.60
1	CA	1363	C	O5'-P-OP2	-9.80	96.88	105.70
1	AA	1421	C	O5'-P-OP1	-9.80	96.88	105.70
1	AA	1006	C	C2-N3-C4	-9.79	115.01	119.90
1	AA	883	G	C2-N3-C4	9.78	116.79	111.90
2	AB	13	A	O5'-P-OP1	-9.78	96.90	105.70
1	AA	739	C	OP1-P-OP2	-9.77	104.94	119.60
1	AA	1605	A	C4-C5-C6	9.77	121.89	117.00
1	CA	2824	C	O5'-P-OP2	-9.76	96.91	105.70
1	AA	2723	A	O5'-P-OP2	-9.76	96.92	105.70
34	BA	1502	A	C2-N3-C4	-9.75	105.72	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1038	C	C2-N3-C4	-9.75	115.03	119.90
1	AA	240	A	C8-N9-C4	9.74	109.70	105.80
1	AA	2533	C	N3-C4-C5	9.74	125.80	121.90
1	AA	593	G	C5-C6-N1	9.73	116.37	111.50
1	CA	1284	A	N1-C6-N6	9.73	124.44	118.60
34	BA	868	C	O5'-P-OP1	-9.72	96.95	105.70
1	AA	843	C	C2-N3-C4	-9.72	115.04	119.90
1	CA	330	A	N1-C6-N6	9.71	124.43	118.60
1	AA	513	C	O5'-P-OP1	-9.71	96.97	105.70
34	BA	819	A	O5'-P-OP1	-9.70	96.97	105.70
34	DA	914	A	O5'-P-OP1	-9.70	96.97	105.70
1	AA	593	G	C4-C5-N7	9.69	114.67	110.80
34	BA	550	G	O5'-P-OP1	-9.69	96.98	105.70
1	AA	604	C	O5'-P-OP2	-9.68	96.98	105.70
1	AA	1249	A	C4-C5-C6	9.68	121.84	117.00
1	CA	1946	U	C5-C4-O4	-9.67	120.10	125.90
1	AA	2525	G	C5-C6-O6	-9.66	122.80	128.60
1	AA	794	U	O5'-P-OP2	-9.66	97.01	105.70
1	AA	993	G	C5-C6-N1	9.65	116.32	111.50
1	AA	1307	C	C5-C6-N1	-9.64	116.18	121.00
1	CA	2708	G	C8-N9-C4	9.63	110.25	106.40
1	AA	1274	G	N1-C2-N2	-9.63	107.53	116.20
1	AA	575	G	N1-C6-O6	-9.62	114.12	119.90
1	AA	1157	A	O4'-C1'-N9	9.62	115.90	108.20
1	AA	989	G	N3-C4-N9	9.62	131.77	126.00
1	AA	1186	U	C5-C6-N1	-9.62	117.89	122.70
1	AA	978	A	C8-N9-C4	-9.61	101.95	105.80
1	AA	2556	G	N3-C2-N2	-9.61	113.17	119.90
1	AA	246	A	O5'-P-OP2	-9.60	97.06	105.70
1	CA	1698	A	N7-C8-N9	9.60	118.60	113.80
1	AA	716	G	C5-C6-N1	9.59	116.30	111.50
1	AA	1255	A	P-O3'-C3'	9.59	131.21	119.70
56	BW	17	C	C2-N1-C1'	9.59	129.35	118.80
1	CA	221	A	O5'-P-OP1	-9.59	97.07	105.70
1	AA	1812	C	C5-C4-N4	-9.59	113.49	120.20
1	AA	2466	G	N1-C2-N3	9.59	129.65	123.90
1	AA	2041	A	C5-C6-N6	9.58	131.36	123.70
1	AA	500	G	C8-N9-C4	-9.58	102.57	106.40
1	AA	2775	G	N1-C6-O6	-9.58	114.16	119.90
1	AA	2636	G	C5-N7-C8	9.57	109.09	104.30
34	BA	1502	A	C6-C5-N7	-9.57	125.60	132.30
1	AA	2441	G	O5'-P-OP1	9.57	122.18	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	778	C	C5-C6-N1	-9.56	116.22	121.00
1	AA	1724	A	C2-N3-C4	-9.56	105.82	110.60
1	AA	2508	C	O5'-P-OP2	-9.56	97.09	105.70
56	BW	36	A	C6-N1-C2	9.55	124.33	118.60
1	AA	560	C	C2-N3-C4	-9.55	115.13	119.90
1	AA	1082	G	C5-C6-O6	9.54	134.33	128.60
1	CA	1142(A)	A	C2-N3-C4	-9.55	105.83	110.60
1	AA	1097	G	N1-C6-O6	9.53	125.62	119.90
1	CA	2229	C	C6-N1-C2	-9.53	116.49	120.30
1	AA	894	U	C2-N3-C4	-9.53	121.28	127.00
1	AA	1023	G	O5'-P-OP2	-9.53	97.13	105.70
1	AA	600	G	C5-C6-O6	-9.52	122.89	128.60
1	CA	1698	A	C4-C5-C6	9.52	121.76	117.00
1	AA	1455	C	C2-N3-C4	-9.52	115.14	119.90
1	AA	1784	G	N3-C2-N2	9.52	126.56	119.90
1	AA	2033	U	C5-C4-O4	9.52	131.61	125.90
34	BA	1502	A	N1-C2-N3	9.51	134.06	129.30
1	AA	1680	G	O5'-P-OP1	-9.51	97.14	105.70
1	AA	1728	G	C4-C5-N7	9.50	114.60	110.80
1	AA	1961	U	N3-C4-C5	9.50	120.30	114.60
1	AA	2581	G	O5'-P-OP1	-9.49	97.16	105.70
1	AA	1038	C	N3-C4-C5	9.49	125.70	121.90
1	AA	139	A	C8-N9-C4	-9.48	102.01	105.80
1	AA	1082	G	C5-N7-C8	9.48	109.04	104.30
1	AA	2298	A	C4-C5-C6	9.48	121.74	117.00
1	AA	553	A	C5-C6-N1	-9.48	112.96	117.70
1	AA	45	C	N1-C2-O2	-9.48	113.21	118.90
1	AA	1328	U	C5-C6-N1	-9.47	117.96	122.70
1	AA	1282	G	N9-C4-C5	-9.47	101.61	105.40
1	AA	2533	C	C2-N3-C4	-9.46	115.17	119.90
1	AA	776	G	C8-N9-C4	-9.46	102.61	106.40
1	AA	714	U	C5-C6-N1	-9.46	117.97	122.70
1	AA	2610	A	OP2-P-O3'	9.46	126.00	105.20
1	AA	796	C	C5-C6-N1	-9.45	116.27	121.00
1	AA	1270	C	C5-C6-N1	-9.44	116.28	121.00
1	AA	1812	C	N3-C4-C5	9.44	125.68	121.90
2	AB	46	A	O5'-P-OP1	-9.44	97.20	105.70
1	AA	2464	C	N1-C2-O2	-9.44	113.24	118.90
1	AA	2302	G	O5'-P-OP2	9.43	122.02	110.70
1	AA	823	G	C5-N7-C8	9.43	109.01	104.30
1	AA	1301	U	C5-C6-N1	9.42	127.41	122.70
1	CA	1204	A	C4-C5-N7	9.42	115.41	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	26	G	C5-C6-O6	-9.41	122.95	128.60
1	AA	1261	G	C5-C6-O6	-9.41	122.95	128.60
1	AA	1189	A	N9-C4-C5	9.41	109.56	105.80
1	CA	133	C	C6-N1-C2	9.41	124.06	120.30
1	AA	147	U	C6-N1-C2	9.40	126.64	121.00
1	AA	2107	C	C6-N1-C2	9.40	124.06	120.30
1	AA	2057	G	C5-C6-O6	-9.40	122.96	128.60
1	AA	174	U	O5'-P-OP2	-9.40	97.24	105.70
1	AA	2609	G	N1-C6-O6	9.39	125.54	119.90
1	AA	436	C	O5'-P-OP2	-9.39	97.25	105.70
1	AA	884	C	C4-C5-C6	9.39	122.09	117.40
1	CA	1902	C	C6-N1-C2	-9.38	116.55	120.30
1	AA	1350	C	O5'-P-OP2	-9.38	97.26	105.70
1	AA	723	A	C5-C6-N1	-9.38	113.01	117.70
1	AA	1659	G	C4-C5-N7	9.38	114.55	110.80
1	AA	643	C	N1-C2-O2	-9.37	113.28	118.90
1	AA	2019	G	C5-C6-N1	9.37	116.18	111.50
1	CA	2584	U	C5-C4-O4	-9.36	120.28	125.90
1	AA	493	G	N1-C6-O6	-9.36	114.28	119.90
1	AA	290	G	C8-N9-C4	9.36	110.14	106.40
1	AA	2281	A	O5'-P-OP1	-9.36	97.28	105.70
1	AA	894	U	C5-C4-O4	9.35	131.51	125.90
34	BA	1496	C	O5'-P-OP2	-9.35	97.29	105.70
1	AA	2418	U	O5'-P-OP2	9.34	121.91	110.70
1	AA	1811	A	C8-N9-C4	-9.34	102.07	105.80
34	BA	235	C	O5'-P-OP1	-9.34	97.30	105.70
1	CA	2548	G	C5-C6-O6	-9.34	123.00	128.60
1	AA	1249	A	N7-C8-N9	9.33	118.46	113.80
1	AA	751	G	O4'-C1'-N9	9.32	115.66	108.20
1	AA	2797	C	C2-N3-C4	-9.32	115.24	119.90
1	CA	2818	G	C8-N9-C4	9.32	110.13	106.40
1	CA	2286	A	C6-C5-N7	-9.31	125.78	132.30
1	AA	787	U	C6-N1-C2	-9.30	115.42	121.00
1	AA	315	C	C6-N1-C2	9.30	124.02	120.30
34	BA	665	A	O5'-P-OP2	-9.30	97.33	105.70
1	AA	560	C	C4-C5-C6	9.30	122.05	117.40
1	AA	591	U	C4-C5-C6	-9.29	114.12	119.70
1	AA	1081	U	N3-C4-C5	9.29	120.18	114.60
1	AA	1821	C	OP1-P-O3'	9.29	125.64	105.20
1	AA	1342	G	C5-C6-O6	9.29	134.17	128.60
1	AA	1985	U	C5-C6-N1	9.28	127.34	122.70
1	CA	388	G	O5'-P-OP1	-9.28	97.35	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	858	U	N1-C2-O2	-9.28	116.31	122.80
1	AA	181	C	N1-C2-O2	-9.27	113.34	118.90
1	AA	2078	G	C5-C6-O6	9.27	134.16	128.60
1	AA	1307	C	C2-N3-C4	-9.27	115.27	119.90
1	AA	1058	U	C5-C6-N1	-9.26	118.07	122.70
1	AA	1020	C	N3-C2-O2	-9.26	115.42	121.90
1	AA	1317	G	N9-C4-C5	-9.25	101.70	105.40
1	AA	786	G	N3-C4-C5	-9.24	123.98	128.60
1	CA	1204	A	C5-C6-N1	-9.24	113.08	117.70
1	AA	2066	C	N1-C2-O2	-9.24	113.36	118.90
1	AA	2581	G	OP1-P-OP2	9.23	133.45	119.60
1	AA	1458	A	O5'-P-OP1	-9.23	97.39	105.70
1	AA	1659	G	C5-C6-N1	9.22	116.11	111.50
1	CA	2425	A	O5'-P-OP2	-9.22	97.40	105.70
1	AA	1053	C	C4-C5-C6	9.21	122.00	117.40
1	CA	1698	A	C2-N3-C4	-9.20	106.00	110.60
1	AA	2902	G	C5-N7-C8	-9.19	99.70	104.30
1	AA	2060	G	N1-C6-O6	-9.19	114.39	119.90
1	AA	2460	A	O5'-P-OP1	-9.19	97.43	105.70
1	AA	1053	C	C5-C6-N1	-9.19	116.41	121.00
1	AA	2711	C	N1-C2-O2	9.19	124.41	118.90
1	AA	625	G	N3-C2-N2	9.18	126.33	119.90
1	AA	1251	G	C2-N3-C4	9.18	116.49	111.90
1	AA	2110	G	N1-C6-O6	9.18	125.41	119.90
1	CA	520	G	O5'-P-OP2	-9.18	97.44	105.70
1	CA	1296	G	C5-C6-O6	9.18	134.11	128.60
1	AA	1299	A	C5-C6-N1	9.17	122.29	117.70
1	AA	1856	A	O5'-P-OP2	-9.17	97.44	105.70
1	AA	2617	U	O5'-P-OP2	-9.17	97.45	105.70
1	CA	1934	C	C6-N1-C2	9.17	123.97	120.30
1	AA	894	U	C2-N1-C1'	-9.17	106.70	117.70
1	AA	1720	U	N1-C2-O2	-9.17	116.38	122.80
1	AA	1964	C	OP1-P-OP2	-9.17	105.85	119.60
1	AA	2269	U	OP1-P-OP2	-9.17	105.85	119.60
1	CA	2234	G	C4-C5-N7	9.17	114.47	110.80
1	AA	2556	G	O5'-P-OP1	-9.16	97.45	105.70
34	BA	1502	A	N7-C8-N9	9.16	118.38	113.80
1	CA	1979	C	O5'-P-OP2	-9.16	97.45	105.70
34	BA	33	A	O5'-P-OP2	-9.16	97.45	105.70
1	AA	1851	U	C5-C6-N1	-9.16	118.12	122.70
1	AA	2100	C	N3-C4-C5	9.15	125.56	121.90
2	AB	103	G	N1-C6-O6	9.15	125.39	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1327	C	O5'-P-OP2	-9.15	97.47	105.70
1	AA	446	C	O5'-P-OP2	-9.15	97.47	105.70
1	AA	1006	C	O5'-P-OP2	-9.14	97.47	105.70
1	AA	1155	C	C6-N1-C2	-9.14	116.64	120.30
34	BA	914	A	O5'-P-OP1	-9.14	97.47	105.70
34	DA	1045	C	OP2-P-O3'	-9.14	85.08	105.20
1	AA	990	A	N9-C4-C5	-9.14	102.14	105.80
1	AA	416	G	C6-N1-C2	-9.13	119.62	125.10
1	AA	2902	G	N7-C8-N9	9.12	117.66	113.10
34	BA	365	U	C5-C6-N1	-9.12	118.14	122.70
1	AA	1001	G	N3-C2-N2	-9.12	113.52	119.90
1	AA	2368	C	N1-C2-O2	-9.12	113.43	118.90
1	AA	413	G	C6-C5-N7	-9.12	124.93	130.40
1	AA	2718	G	N7-C8-N9	-9.12	108.54	113.10
1	AA	601	A	N1-C6-N6	-9.11	113.13	118.60
1	AA	772	G	C8-N9-C4	9.11	110.04	106.40
1	CA	1638	C	O5'-P-OP2	-9.11	97.50	105.70
1	AA	649	C	O5'-P-OP1	-9.11	97.50	105.70
1	AA	1998	U	O5'-P-OP2	-9.09	97.52	105.70
1	AA	209	G	N7-C8-N9	-9.08	108.56	113.10
1	AA	1000	C	N1-C2-N3	9.08	125.56	119.20
1	AA	1312	G	C5-C6-N1	9.08	116.04	111.50
1	CA	659	C	C5-C6-N1	-9.08	116.46	121.00
1	CA	748	G	N1-C6-O6	-9.08	114.45	119.90
1	AA	550	U	N1-C2-N3	9.07	120.34	114.90
1	AA	2518	U	P-O3'-C3'	9.07	130.58	119.70
1	AA	560	C	C6-N1-C2	9.06	123.92	120.30
1	AA	2550	C	N3-C2-O2	9.05	128.24	121.90
1	CA	773	U	O5'-P-OP1	-9.05	97.55	105.70
1	CA	1957	C	N3-C4-N4	9.05	124.34	118.00
1	AA	1802	C	N1-C2-O2	-9.05	113.47	118.90
1	AA	2037	A	C4-C5-C6	9.05	121.53	117.00
1	AA	2895	C	C6-N1-C2	-9.05	116.68	120.30
1	AA	2515	A	C2-N3-C4	9.04	115.12	110.60
1	AA	2748	G	N1-C6-O6	-9.04	114.47	119.90
34	BA	394	G	O5'-P-OP1	-9.04	97.56	105.70
1	AA	786	G	C5-C6-O6	-9.04	123.17	128.60
1	AA	1645	C	C5-C6-N1	-9.04	116.48	121.00
34	BA	1064	G	N3-C4-C5	9.04	133.12	128.60
1	CA	1760	A	O5'-P-OP2	-9.04	97.57	105.70
1	AA	126	C	C5-C4-N4	-9.03	113.88	120.20
1	AA	1298	G	OP2-P-O3'	9.03	125.06	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2827	G	C5-C6-O6	-9.03	123.18	128.60
1	CA	2544	G	C5-C6-O6	-9.03	123.18	128.60
1	AA	2004	C	C4-C5-C6	9.02	121.91	117.40
1	AA	1331	G	C5-C6-O6	-9.02	123.19	128.60
1	AA	2046	G	C8-N9-C4	9.02	110.01	106.40
1	AA	2383	G	C5-C6-O6	-9.02	123.19	128.60
1	AA	730	C	C5-C6-N1	-9.01	116.49	121.00
1	CA	474	G	N1-C6-O6	-9.01	114.49	119.90
1	AA	592	U	C2-N3-C4	-9.01	121.60	127.00
1	AA	663	G	N1-C6-O6	-9.01	114.50	119.90
1	AA	1733	C	N3-C2-O2	-9.01	115.59	121.90
1	CA	2234	G	C5-C6-O6	-9.01	123.20	128.60
2	AB	41	U	C5-C6-N1	-9.00	118.20	122.70
1	CA	1947	C	C6-N1-C2	9.00	123.90	120.30
1	CA	1992	G	N1-C6-O6	-9.00	114.50	119.90
1	AA	786	G	C6-N1-C2	-9.00	119.70	125.10
1	AA	1474	C	O5'-P-OP1	-9.00	97.60	105.70
1	AA	1657	C	C2-N3-C4	-9.00	115.40	119.90
1	CA	945	A	C5-N7-C8	-8.99	99.40	103.90
1	AA	1428	G	C5-C6-O6	-8.99	123.21	128.60
1	AA	121	G	O5'-P-OP2	-8.98	97.61	105.70
1	AA	600	G	C5-C6-N1	8.98	115.99	111.50
1	AA	353	G	C5-C6-O6	-8.98	123.21	128.60
1	AA	1097	G	O5'-P-OP1	-8.98	97.62	105.70
1	CA	2874	C	C6-N1-C2	-8.98	116.71	120.30
1	AA	12	U	N3-C2-O2	-8.98	115.92	122.20
1	AA	1809	U	N1-C2-O2	-8.97	116.52	122.80
1	AA	2331	G	C2-N3-C4	-8.97	107.42	111.90
1	AA	354	A	N7-C8-N9	8.97	118.28	113.80
1	AA	1021	G	O5'-P-OP2	-8.96	97.63	105.70
1	AA	2460	A	C5-C6-N1	8.97	122.18	117.70
1	AA	2651	A	N1-C2-N3	8.97	133.78	129.30
1	AA	845	G	C4-C5-N7	-8.96	107.22	110.80
1	AA	2858	G	O4'-C1'-N9	8.96	115.37	108.20
1	AA	535	C	N1-C2-O2	8.96	124.27	118.90
1	AA	978	A	N3-C4-N9	-8.96	120.24	127.40
34	BA	365	U	C2-N1-C1'	-8.96	106.95	117.70
1	CA	1698	A	C8-N9-C4	-8.95	102.22	105.80
1	CA	2850	A	O5'-P-OP2	-8.95	97.65	105.70
1	AA	497	A	N1-C2-N3	8.94	133.77	129.30
1	AA	1543	U	N3-C4-O4	-8.94	113.14	119.40
1	CA	1296	G	N1-C6-O6	-8.94	114.54	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	324	A	C2-N3-C4	8.93	115.07	110.60
1	AA	2448	G	O5'-P-OP1	-8.93	97.66	105.70
1	CA	975	C	N1-C2-O2	8.93	124.26	118.90
1	AA	2248	C	N3-C4-C5	8.93	125.47	121.90
1	CA	1639	U	O5'-P-OP2	-8.93	97.67	105.70
1	AA	564	G	C8-N9-C4	8.92	109.97	106.40
1	AA	1010	C	C6-N1-C2	-8.92	116.73	120.30
1	AA	1745	A	O4'-C1'-N9	8.92	115.33	108.20
1	AA	609	A	N1-C2-N3	-8.92	124.84	129.30
1	AA	787	U	O5'-P-OP1	8.91	121.40	110.70
1	AA	1039	G	O5'-P-OP2	-8.91	97.68	105.70
1	AA	183	G	C4-C5-N7	-8.90	107.24	110.80
1	AA	841	G	N9-C4-C5	-8.90	101.84	105.40
1	AA	524	U	N3-C4-O4	-8.90	113.17	119.40
1	AA	1664	A	O5'-P-OP1	-8.90	97.69	105.70
34	BA	533	A	O5'-P-OP1	-8.90	97.69	105.70
1	AA	327	U	C5-C6-N1	-8.89	118.25	122.70
1	AA	2429	C	O5'-P-OP2	-8.89	97.70	105.70
1	AA	2497	G	N9-C4-C5	-8.89	101.84	105.40
1	CA	2046	G	C8-N9-C4	8.89	109.96	106.40
1	AA	1658	C	N3-C4-C5	8.88	125.45	121.90
1	AA	2071	G	N7-C8-N9	8.88	117.54	113.10
1	AA	1701	A	C8-N9-C4	8.88	109.35	105.80
1	CA	1284	A	N9-C4-C5	-8.87	102.25	105.80
1	AA	16	G	O5'-P-OP1	-8.87	97.72	105.70
1	AA	1270	C	C2-N3-C4	-8.87	115.47	119.90
34	DA	1415	G	O5'-P-OP2	-8.86	97.73	105.70
1	AA	790	G	N1-C6-O6	-8.85	114.59	119.90
1	AA	1316	C	C2-N3-C4	-8.85	115.47	119.90
1	AA	2409	G	N1-C6-O6	-8.84	114.59	119.90
1	AA	2107	C	C2-N3-C4	-8.84	115.48	119.90
1	CA	1901	A	C2-N3-C4	8.84	115.02	110.60
1	AA	1340	U	C5-C6-N1	-8.83	118.28	122.70
1	AA	2277	U	C4-C5-C6	8.83	125.00	119.70
1	AA	2874	G	O5'-P-OP2	-8.83	97.75	105.70
1	CA	575	A	O5'-P-OP2	8.83	121.30	110.70
34	BA	1502	A	N1-C6-N6	8.83	123.90	118.60
1	AA	195	U	N3-C4-O4	-8.83	113.22	119.40
1	AA	2786	C	N3-C4-C5	8.83	125.43	121.90
1	AA	775	G	N3-C2-N2	8.83	126.08	119.90
1	AA	599	U	N1-C2-N3	8.82	120.19	114.90
1	AA	2641	A	C5-C6-N6	-8.82	116.64	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2576	G	C2-N3-C4	8.82	116.31	111.90
1	AA	634	C	C6-N1-C2	8.82	123.83	120.30
1	AA	981	C	N1-C2-O2	-8.82	113.61	118.90
1	AA	174	U	C5-C6-N1	-8.82	118.29	122.70
1	CA	1698	A	O4'-C1'-N9	8.82	115.25	108.20
1	AA	184	A	P-O3'-C3'	8.82	130.28	119.70
1	AA	2014	G	N3-C4-C5	-8.82	124.19	128.60
1	AA	2331	G	N9-C4-C5	8.82	108.93	105.40
1	AA	2453	C	C2-N3-C4	-8.81	115.49	119.90
1	AA	1308	A	O5'-P-OP2	8.81	121.27	110.70
1	CA	2563	U	N3-C2-O2	-8.81	116.03	122.20
1	CA	2593	U	N3-C4-O4	-8.81	113.23	119.40
1	AA	2476	C	N3-C4-C5	8.80	125.42	121.90
1	AA	2470	G	N3-C2-N2	-8.80	113.74	119.90
1	AA	786	G	C2-N3-C4	8.80	116.30	111.90
1	AA	907	U	O5'-P-OP2	-8.80	97.78	105.70
1	AA	1078	A	N1-C6-N6	8.80	123.88	118.60
1	AA	2285	A	C4-C5-N7	8.80	115.10	110.70
1	AA	822	G	N9-C4-C5	8.79	108.92	105.40
1	AA	2548	G	C4-C5-N7	-8.79	107.28	110.80
1	AA	1860	A	O5'-P-OP2	-8.79	97.79	105.70
1	AA	2443	U	O5'-P-OP2	-8.79	97.79	105.70
1	AA	2631	C	C5-C6-N1	-8.79	116.61	121.00
1	AA	2638	C	C5-C6-N1	-8.79	116.61	121.00
1	AA	887	C	O5'-P-OP2	-8.78	97.80	105.70
1	AA	1694	G	C4-C5-N7	-8.78	107.29	110.80
1	AA	2023	A	N1-C6-N6	-8.78	113.33	118.60
1	AA	2068	G	C5-C6-N1	8.78	115.89	111.50
1	AA	1926	G	C5-C6-O6	8.78	133.87	128.60
1	AA	2791	A	N1-C2-N3	8.78	133.69	129.30
1	AA	2307	C	O5'-P-OP1	-8.77	97.80	105.70
1	CA	1345	C	C6-N1-C2	8.77	123.81	120.30
1	AA	2299	A	C6-N1-C2	8.77	123.86	118.60
1	AA	22	C	N1-C2-O2	8.77	124.16	118.90
1	CA	2490	G	C4-C5-N7	8.77	114.31	110.80
1	CA	2006	C	N1-C2-O2	-8.76	113.64	118.90
1	AA	2115	G	C8-N9-C4	-8.76	102.89	106.40
1	AA	2260	C	C5-C6-N1	-8.76	116.62	121.00
1	AA	2384	G	N1-C6-O6	8.76	125.15	119.90
1	AA	1065	U	C5-C6-N1	-8.75	118.32	122.70
1	AA	1249	A	C8-N9-C4	-8.75	102.30	105.80
1	AA	1824	C	N3-C4-N4	-8.75	111.87	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1994	A	O5'-P-OP1	-8.75	97.83	105.70
1	AA	772	G	N7-C8-N9	-8.75	108.73	113.10
1	AA	2586	G	C5-C6-N1	8.75	115.87	111.50
16	CS	67	ARG	NE-CZ-NH1	8.75	124.67	120.30
1	CA	2287	A	O5'-P-OP2	-8.74	97.83	105.70
1	AA	535	C	N3-C2-O2	-8.74	115.78	121.90
1	CA	2548	G	N1-C6-O6	8.74	125.14	119.90
1	AA	836	A	O5'-P-OP1	-8.73	97.84	105.70
1	CA	2420	C	O5'-P-OP1	-8.73	97.85	105.70
1	AA	1045	U	N3-C4-C5	8.72	119.83	114.60
1	CA	448	U	O5'-P-OP1	-8.72	97.85	105.70
1	CA	1284	A	C4-C5-N7	8.72	115.06	110.70
1	CA	1323	U	N3-C4-C5	8.71	119.83	114.60
1	AA	126	C	N3-C4-C5	8.71	125.38	121.90
1	AA	649	C	C2-N3-C4	-8.71	115.55	119.90
1	AA	412	C	N3-C2-O2	8.71	127.99	121.90
1	AA	2331	G	N7-C8-N9	8.71	117.45	113.10
34	BA	1502	A	C4-C5-N7	8.71	115.05	110.70
1	AA	1008	U	C2-N3-C4	-8.70	121.78	127.00
1	AA	2530	A	N1-C2-N3	8.70	133.65	129.30
1	AA	1685	C	N3-C4-C5	8.70	125.38	121.90
2	AB	108	U	O5'-P-OP2	-8.70	97.87	105.70
1	AA	112	U	OP1-P-OP2	-8.69	106.56	119.60
1	AA	1986	G	O5'-P-OP1	-8.69	97.88	105.70
2	AB	82	G	C4-C5-N7	-8.69	107.32	110.80
1	AA	1200	G	N3-C4-N9	8.68	131.21	126.00
34	BA	266	G	C6-C5-N7	-8.68	125.19	130.40
1	AA	2527	C	N3-C2-O2	8.68	127.98	121.90
1	AA	2471	A	C2-N3-C4	8.68	114.94	110.60
1	AA	2394	G	O5'-P-OP2	-8.67	97.90	105.70
1	AA	36	G	N1-C6-O6	-8.67	114.70	119.90
1	AA	353	G	O5'-P-OP2	-8.67	97.90	105.70
1	AA	2096	U	N1-C2-O2	-8.66	116.73	122.80
1	AA	1919	G	O5'-P-OP1	-8.66	97.91	105.70
1	AA	2657	G	N3-C4-C5	8.66	132.93	128.60
1	CA	1941	C	C6-N1-C2	-8.65	116.84	120.30
1	AA	2057	G	C8-N9-C4	8.65	109.86	106.40
1	AA	727	G	O5'-P-OP1	-8.65	97.92	105.70
1	CA	1937	A	O4'-C1'-N9	8.65	115.12	108.20
1	CA	2679	A	O5'-P-OP2	-8.65	97.92	105.70
1	AA	1052	C	N1-C2-O2	-8.64	113.71	118.90
1	AA	2657	G	C5-N7-C8	-8.64	99.98	104.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	354	A	C5-C6-N1	-8.64	113.38	117.70
1	CA	2599	G	C5-C6-O6	8.64	133.78	128.60
1	AA	418	G	N9-C4-C5	-8.64	101.94	105.40
1	AA	1342	G	C2-N3-C4	8.64	116.22	111.90
1	AA	1011	G	C8-N9-C4	-8.63	102.95	106.40
1	AA	2290	A	N9-C4-C5	8.64	109.25	105.80
1	AA	204	G	O5'-P-OP2	8.63	121.06	110.70
1	AA	1816	A	C4-C5-C6	8.63	121.31	117.00
6	AF	74	ARG	NE-CZ-NH1	8.63	124.61	120.30
34	DA	50	A	C8-N9-C4	-8.63	102.35	105.80
1	AA	2718	G	C8-N9-C4	8.63	109.85	106.40
1	AA	900	G	O5'-P-OP2	-8.62	97.94	105.70
1	AA	2487	C	N3-C4-C5	-8.62	118.45	121.90
1	AA	2041	A	C5-C6-N1	-8.62	113.39	117.70
1	AA	2527	C	C5-C4-N4	-8.62	114.17	120.20
1	CA	1654	A	O5'-P-OP2	8.62	121.04	110.70
1	AA	187	C	O5'-P-OP2	-8.62	97.94	105.70
1	AA	2690	C	C2-N3-C4	8.61	124.21	119.90
1	AA	254	A	O4'-C1'-N9	8.61	115.09	108.20
18	AU	3	ARG	NE-CZ-NH1	-8.61	116.00	120.30
34	BA	915	A	O5'-P-OP2	-8.60	97.96	105.70
34	BA	1530	G	N1-C6-O6	8.60	125.06	119.90
1	AA	1081	U	C2-N3-C4	-8.60	121.84	127.00
1	AA	2693	C	N3-C4-N4	8.60	124.02	118.00
1	AA	1503	G	OP1-P-OP2	-8.59	106.71	119.60
34	BA	1523	G	N1-C6-O6	-8.59	114.74	119.90
1	CA	1790	C	N3-C4-C5	-8.59	118.46	121.90
34	DA	562	C	O5'-P-OP1	-8.59	97.97	105.70
1	AA	2014	G	C2-N3-C4	8.59	116.19	111.90
1	AA	2377	G	O5'-P-OP2	-8.59	97.97	105.70
1	CA	1763	G	C8-N9-C4	8.59	109.83	106.40
34	BA	266	G	C4-C5-N7	8.58	114.23	110.80
1	CA	1428	C	O5'-P-OP2	8.57	120.99	110.70
1	AA	804	U	C2-N3-C4	-8.57	121.86	127.00
1	AA	1686	U	N1-C2-O2	8.57	128.80	122.80
1	AA	241	G	C5-C6-O6	-8.56	123.46	128.60
1	AA	2833	A	C2-N3-C4	-8.56	106.32	110.60
1	CA	2540	C	N3-C4-C5	8.56	125.33	121.90
1	AA	710	G	N3-C2-N2	-8.56	113.91	119.90
1	AA	1282	G	N7-C8-N9	-8.55	108.82	113.10
1	AA	2833	A	N1-C2-N3	8.55	133.58	129.30
1	CA	141	A	C5-N7-C8	-8.55	99.62	103.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1334	G	N1-C6-O6	-8.55	114.77	119.90
1	AA	1081	U	N3-C4-O4	-8.55	113.42	119.40
1	AA	2881	C	C6-N1-C2	8.55	123.72	120.30
1	AA	853	C	O5'-P-OP2	8.54	120.95	110.70
1	AA	1472	G	C5-C6-N1	8.54	115.77	111.50
1	AA	2457	G	C5-C6-O6	8.54	133.73	128.60
1	AA	1728	G	N3-C4-C5	8.54	132.87	128.60
1	AA	1050	C	C4-C5-C6	8.54	121.67	117.40
1	AA	1329	G	C4-C5-C6	8.54	123.92	118.80
1	AA	2299	A	C4-C5-N7	8.53	114.97	110.70
1	AA	2554	A	C6-N1-C2	8.53	123.72	118.60
1	AA	552	C	C5-C4-N4	8.53	126.17	120.20
1	AA	35	G	C8-N9-C4	8.52	109.81	106.40
1	AA	2290	A	OP1-P-OP2	-8.52	106.81	119.60
1	AA	2516	U	O5'-P-OP2	-8.52	98.03	105.70
34	BA	1509	C	C6-N1-C2	8.52	123.71	120.30
1	CA	141	A	N7-C8-N9	8.52	118.06	113.80
1	AA	1733	C	N1-C2-O2	8.52	124.01	118.90
1	AA	2220	A	OP1-P-O3'	8.51	123.92	105.20
1	AA	2797	C	C5-C6-N1	-8.51	116.75	121.00
1	AA	1041	C	C6-N1-C2	8.51	123.70	120.30
1	AA	2505	U	N3-C2-O2	-8.51	116.25	122.20
1	AA	2529	C	C2-N3-C4	-8.51	115.65	119.90
1	AA	2876	U	N3-C2-O2	-8.51	116.25	122.20
1	CA	2023	G	O5'-P-OP1	-8.51	98.05	105.70
1	AA	718	C	C5-C4-N4	8.50	126.15	120.20
1	AA	1921	G	N7-C8-N9	8.50	117.35	113.10
2	AB	67	G	O5'-P-OP1	-8.50	98.05	105.70
1	CA	1996	C	C6-N1-C2	8.50	123.70	120.30
1	AA	842	C	N1-C2-O2	8.49	124.00	118.90
1	AA	1067	A	N7-C8-N9	8.49	118.05	113.80
1	CA	795	C	O5'-P-OP2	-8.49	98.06	105.70
1	AA	66	U	O5'-P-OP1	-8.49	98.06	105.70
1	CA	83	G	O5'-P-OP2	-8.49	98.06	105.70
1	CA	2540	C	C6-N1-C2	8.49	123.70	120.30
1	AA	1050	C	C6-N1-C2	-8.48	116.91	120.30
1	CA	963	U	OP1-P-OP2	-8.48	106.87	119.60
1	AA	2244	U	N3-C4-O4	-8.48	113.46	119.40
1	AA	1097	G	C5-C6-O6	-8.48	123.51	128.60
1	AA	27	G	O5'-P-OP2	-8.47	98.08	105.70
1	AA	794	U	OP1-P-OP2	8.47	132.30	119.60
1	AA	860	U	N3-C4-O4	-8.47	113.47	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1921	G	C8-N9-C4	-8.47	103.01	106.40
1	AA	2285	A	C5-C6-N6	-8.47	116.93	123.70
34	BA	365	U	C5-C4-O4	8.47	130.98	125.90
1	AA	1718	U	C5-C6-N1	-8.46	118.47	122.70
1	AA	2643	G	C8-N9-C4	-8.46	103.02	106.40
1	CA	1996	C	O5'-P-OP1	-8.46	98.09	105.70
1	CA	2394	C	N1-C2-O2	-8.46	113.83	118.90
1	AA	1656	A	C5-C6-N1	8.46	121.93	117.70
1	CA	2570	G	N9-C4-C5	8.45	108.78	105.40
1	AA	1261	G	N1-C6-O6	8.45	124.97	119.90
1	AA	2896	G	O5'-P-OP1	-8.45	98.09	105.70
1	AA	2896	G	C2-N3-C4	-8.45	107.67	111.90
34	BA	1523	G	N9-C4-C5	8.45	108.78	105.40
1	AA	853	C	OP1-P-OP2	-8.44	106.94	119.60
1	AA	1274	G	N3-C4-C5	8.44	132.82	128.60
1	CA	575	A	O5'-P-OP1	-8.44	98.11	105.70
34	DA	266	G	C6-C5-N7	-8.44	125.34	130.40
1	AA	873	U	OP2-P-O3'	8.43	123.75	105.20
1	AA	1855	G	OP2-P-O3'	8.43	123.74	105.20
1	AA	111	G	C5-C6-O6	-8.43	123.54	128.60
1	AA	1211	U	C5-C6-N1	-8.43	118.49	122.70
1	AA	2257	U	O5'-P-OP2	-8.42	98.12	105.70
56	BW	47	U	N3-C2-O2	-8.42	116.30	122.20
1	CA	2503	A	C2-N3-C4	8.42	114.81	110.60
34	DA	923	A	C8-N9-C4	-8.42	102.43	105.80
1	AA	100	G	N9-C4-C5	-8.42	102.03	105.40
1	AA	906	G	N1-C6-O6	8.42	124.95	119.90
1	AA	1237	G	C4-C5-N7	-8.42	107.43	110.80
1	CA	2581	G	C8-N9-C4	-8.42	103.03	106.40
1	AA	1843	A	OP1-P-OP2	8.41	132.22	119.60
1	AA	1018	A	C8-N9-C4	-8.40	102.44	105.80
1	AA	1361	C	N1-C2-O2	-8.39	113.86	118.90
1	AA	2453	C	C5-C6-N1	-8.39	116.80	121.00
1	AA	1003	U	C6-N1-C2	-8.39	115.97	121.00
1	AA	1040	C	C5-C6-N1	-8.39	116.81	121.00
1	AA	1371	G	N9-C4-C5	8.39	108.75	105.40
13	AP	18	ARG	NE-CZ-NH2	-8.39	116.11	120.30
1	AA	2046	G	N7-C8-N9	-8.38	108.91	113.10
1	CA	2409	G	N1-C6-O6	8.38	124.93	119.90
1	AA	1429	C	O5'-P-OP2	8.38	120.75	110.70
1	AA	624	C	N3-C2-O2	-8.37	116.04	121.90
1	AA	918	U	C5-C4-O4	-8.37	120.88	125.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2110	G	C5-C6-O6	-8.37	123.58	128.60
1	AA	1611	C	O5'-P-OP2	-8.37	98.17	105.70
34	DA	266	G	C4-C5-N7	8.37	114.15	110.80
1	AA	1710	C	C4-C5-C6	8.37	121.58	117.40
1	AA	606	G	N9-C4-C5	8.36	108.74	105.40
34	BA	328	C	O5'-P-OP1	-8.36	98.18	105.70
1	AA	2828	G	C5-C6-O6	8.35	133.61	128.60
1	AA	2691	A	N9-C4-C5	8.35	109.14	105.80
1	AA	1745	A	N9-C4-C5	-8.34	102.46	105.80
1	AA	2244	U	O5'-P-OP2	-8.34	98.19	105.70
1	AA	209	G	C5-N7-C8	8.34	108.47	104.30
1	AA	1700	G	N1-C6-O6	-8.34	114.90	119.90
1	AA	2018	C	N1-C2-O2	-8.33	113.90	118.90
1	AA	2078	G	C8-N9-C4	-8.33	103.07	106.40
1	AA	2294	G	O5'-P-OP1	-8.33	98.20	105.70
34	BA	767	A	N1-C6-N6	-8.33	113.60	118.60
1	CA	2498	C	N1-C2-O2	-8.33	113.90	118.90
1	AA	1074	A	C5-C6-N1	8.33	121.86	117.70
1	AA	1266	C	N3-C2-O2	-8.33	116.07	121.90
1	AA	1813	C	O5'-P-OP2	8.33	120.69	110.70
1	AA	738	C	C6-N1-C2	-8.33	116.97	120.30
1	AA	2607	G	C5-C6-O6	8.32	133.59	128.60
29	A5	15	ARG	NE-CZ-NH1	-8.32	116.14	120.30
1	AA	2041	A	C5-N7-C8	8.32	108.06	103.90
1	AA	856	G	N1-C6-O6	-8.32	114.91	119.90
34	BA	1528	U	O5'-P-OP2	-8.32	98.21	105.70
1	AA	802	C	C5-C6-N1	-8.32	116.84	121.00
1	AA	1829	U	N3-C4-O4	-8.32	113.58	119.40
1	CA	2353	G	O5'-P-OP2	-8.32	98.21	105.70
1	CA	270	A	C8-N9-C4	8.32	109.13	105.80
1	AA	1274	G	C5-C6-O6	8.31	133.59	128.60
1	CA	1204	A	C6-C5-N7	-8.31	126.48	132.30
1	AA	464	G	C5-C6-O6	-8.31	123.61	128.60
1	AA	2749	G	C8-N9-C4	8.31	109.72	106.40
1	AA	1317	G	N3-C4-N9	8.30	130.98	126.00
1	AA	1316	C	C5-C6-N1	-8.30	116.85	121.00
1	AA	2619	G	C6-N1-C2	-8.30	120.12	125.10
1	AA	1197	G	O5'-P-OP2	-8.29	98.23	105.70
1	AA	1617	A	C8-N9-C4	8.29	109.12	105.80
1	AA	2029	C	N1-C2-O2	-8.29	113.93	118.90
1	AA	431	C	N1-C2-O2	8.29	123.87	118.90
1	AA	2304	C	N3-C4-C5	-8.28	118.59	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2894	U	N1-C2-O2	-8.28	117.01	122.80
34	BA	1498	U	C2-N3-C4	-8.27	122.04	127.00
1	AA	793	A	O4'-C1'-N9	8.27	114.82	108.20
1	AA	2325	C	OP2-P-O3'	8.27	123.39	105.20
1	AA	516	G	O5'-P-OP1	-8.27	98.26	105.70
1	AA	2056	U	OP1-P-OP2	8.27	132.00	119.60
1	AA	1379	C	N3-C4-C5	8.27	125.21	121.90
1	AA	1721	G	N3-C4-N9	8.27	130.96	126.00
1	CA	204	A	C5-C6-N1	8.27	121.83	117.70
1	AA	1082	G	N1-C6-O6	-8.26	114.94	119.90
1	AA	1097	G	C4-C5-N7	8.26	114.10	110.80
1	AA	2827	G	C5-C6-N1	8.26	115.63	111.50
1	AA	2657	G	N3-C4-N9	-8.26	121.05	126.00
1	AA	1033	G	C4-C5-N7	-8.25	107.50	110.80
1	AA	2250	G	OP1-P-OP2	8.25	131.98	119.60
1	CA	2646	C	O5'-P-OP2	-8.25	98.27	105.70
1	AA	2707	C	N3-C2-O2	-8.25	116.12	121.90
1	AA	2003	A	C2-N3-C4	-8.25	106.48	110.60
1	CA	614	U	N3-C2-O2	-8.24	116.43	122.20
1	CA	2708	G	N9-C4-C5	-8.24	102.10	105.40
1	AA	539	A	O5'-P-OP1	-8.24	98.28	105.70
1	AA	2103	C	C6-N1-C2	-8.24	117.00	120.30
1	CA	2489	G	C5-C6-O6	-8.23	123.66	128.60
34	BA	1484	C	O5'-P-OP1	-8.23	98.29	105.70
4	CD	14	ARG	NE-CZ-NH2	-8.23	116.19	120.30
1	AA	1045	U	C4-C5-C6	-8.23	114.76	119.70
1	AA	1445	C	N1-C2-O2	-8.23	113.96	118.90
1	AA	1658	C	C5-C6-N1	-8.23	116.89	121.00
1	CA	1638	C	N3-C4-C5	8.23	125.19	121.90
1	AA	172	C	C2-N3-C4	-8.22	115.79	119.90
1	AA	511	C	N3-C2-O2	-8.22	116.15	121.90
1	CA	811	U	C6-N1-C2	8.22	125.93	121.00
1	CA	987	G	O5'-P-OP2	8.22	120.56	110.70
1	CA	811	U	N3-C4-C5	8.21	119.53	114.60
1	AA	564	G	O5'-P-OP1	-8.21	98.31	105.70
1	AA	1661	C	N1-C2-O2	8.21	123.83	118.90
1	AA	2776	G	N1-C6-O6	8.21	124.83	119.90
1	AA	1849	U	O5'-P-OP2	-8.21	98.31	105.70
1	AA	327	U	C2-N3-C4	-8.21	122.08	127.00
1	AA	2605	U	N3-C2-O2	-8.21	116.45	122.20
1	CA	2489	G	N1-C6-O6	8.21	124.83	119.90
1	AA	2282	G	C5-C6-O6	-8.20	123.68	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2713	A	N1-C6-N6	-8.19	113.68	118.60
1	CA	2626	C	C6-N1-C2	8.19	123.58	120.30
1	AA	962	G	O5'-P-OP2	8.19	120.53	110.70
1	AA	2393	C	C2-N3-C4	-8.19	115.81	119.90
1	AA	2376	C	C6-N1-C2	8.19	123.58	120.30
1	AA	787	U	N3-C2-O2	-8.18	116.47	122.20
1	AA	1816	A	C6-C5-N7	-8.18	126.57	132.30
1	AA	2078	G	N3-C2-N2	8.18	125.63	119.90
1	AA	2331	G	N1-C2-N2	8.18	123.56	116.20
1	AA	2019	G	C5-C6-O6	-8.17	123.70	128.60
1	AA	178	G	O5'-P-OP1	-8.17	98.35	105.70
34	BA	509	A	N7-C8-N9	8.17	117.88	113.80
1	AA	959	U	N1-C2-O2	-8.16	117.09	122.80
34	DA	619	U	C2-N1-C1'	-8.16	107.91	117.70
1	CA	2079	U	C4-C5-C6	8.16	124.60	119.70
1	AA	479	C	C5-C6-N1	-8.16	116.92	121.00
1	AA	35	G	N7-C8-N9	-8.16	109.02	113.10
34	BA	355	C	N3-C2-O2	-8.16	116.19	121.90
2	AB	104	U	N3-C4-C5	8.15	119.49	114.60
1	AA	714	U	C2-N3-C4	-8.15	122.11	127.00
1	AA	2619	G	C6-C5-N7	-8.15	125.51	130.40
1	AA	1423	G	C4-C5-N7	-8.15	107.54	110.80
1	CA	1334	G	C5-C6-N1	8.15	115.57	111.50
1	AA	2609	G	C4-C5-N7	8.14	114.06	110.80
1	AA	400	U	C5-C4-O4	-8.14	121.02	125.90
1	AA	241	G	N3-C2-N2	-8.14	114.20	119.90
1	AA	552	C	N3-C4-N4	-8.13	112.31	118.00
1	AA	565	C	N3-C4-C5	8.13	125.15	121.90
1	AA	1462	G	O4'-C1'-N9	8.13	114.71	108.20
1	AA	1657	C	N1-C2-O2	-8.13	114.02	118.90
1	AA	872	C	C5-C6-N1	-8.13	116.93	121.00
1	AA	1420	G	O5'-P-OP2	8.13	120.46	110.70
1	AA	1697	G	C5-C6-N1	8.13	115.56	111.50
1	AA	2351	G	C2-N3-C4	-8.13	107.83	111.90
1	AA	1239	A	N1-C6-N6	-8.13	113.72	118.60
1	AA	540	A	O5'-P-OP1	-8.13	98.39	105.70
1	AA	1620	G	C8-N9-C4	8.13	109.65	106.40
34	BA	1512	U	O5'-P-OP1	8.13	120.45	110.70
1	CA	1204	A	N3-C4-C5	8.13	132.49	126.80
1	AA	172	C	N3-C4-C5	8.12	125.15	121.90
1	CA	1694	C	N1-C2-O2	8.12	123.77	118.90
1	CA	1807	G	C8-N9-C4	8.12	109.65	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	38	A	C5-N7-C8	8.11	107.96	103.90
1	AA	978	A	O4'-C1'-N9	8.12	114.69	108.20
1	AA	2546	A	C5-C6-N6	-8.12	117.21	123.70
31	A7	9	ARG	NE-CZ-NH1	8.12	124.36	120.30
1	CA	1565	C	N1-C2-O2	-8.11	114.03	118.90
1	CA	1692	U	C2-N3-C4	-8.12	122.13	127.00
1	AA	1658	C	C2-N3-C4	-8.11	115.84	119.90
1	AA	2644	A	O5'-P-OP1	-8.11	98.41	105.70
1	AA	2009	G	N3-C2-N2	-8.10	114.23	119.90
1	AA	2548	G	N1-C6-O6	-8.10	115.04	119.90
1	CA	2856	C	C6-N1-C2	-8.10	117.06	120.30
1	AA	1620	G	N7-C8-N9	-8.10	109.05	113.10
34	DA	1103	C	C6-N1-C2	-8.10	117.06	120.30
1	AA	1034	A	C2-N3-C4	8.10	114.65	110.60
34	DA	1046	A	OP1-P-OP2	8.10	131.74	119.60
1	AA	2693	C	C5-C4-N4	-8.09	114.53	120.20
1	AA	1439	A	N1-C2-N3	-8.09	125.25	129.30
1	AA	1053	C	C2-N3-C4	-8.09	115.86	119.90
1	AA	1317	G	OP1-P-OP2	-8.09	107.47	119.60
1	AA	1360	C	C5-C6-N1	-8.09	116.96	121.00
1	AA	1605	A	O4'-C1'-N9	8.09	114.67	108.20
1	AA	577	U	O5'-P-OP1	8.08	120.40	110.70
1	AA	351	G	C5-C6-O6	8.08	133.45	128.60
1	AA	1008	U	N3-C4-C5	8.07	119.44	114.60
1	AA	978	A	C6-N1-C2	8.07	123.44	118.60
1	AA	1431	G	O4'-C1'-N9	8.07	114.66	108.20
1	CA	277	C	N1-C2-O2	8.07	123.74	118.90
1	AA	2386	C	C5-C4-N4	-8.06	114.56	120.20
1	AA	2355	C	N1-C2-O2	-8.06	114.07	118.90
1	AA	605	G	C5-C6-O6	8.05	133.43	128.60
1	AA	902	G	O5'-P-OP2	-8.05	98.45	105.70
1	AA	1710	C	N1-C2-O2	-8.05	114.07	118.90
1	AA	2514	G	C8-N9-C4	8.05	109.62	106.40
1	AA	739	C	O5'-P-OP2	8.05	120.36	110.70
1	AA	1277	G	C8-N9-C4	-8.05	103.18	106.40
1	CA	204	A	N1-C6-N6	-8.05	113.77	118.60
1	AA	2448	G	C8-N9-C4	-8.04	103.18	106.40
2	AB	41	U	C5-C4-O4	8.04	130.73	125.90
1	AA	1423	G	N1-C6-O6	-8.04	115.07	119.90
1	AA	2044	U	OP1-P-OP2	-8.04	107.54	119.60
1	AA	2354	C	O5'-P-OP2	8.04	120.35	110.70
56	BW	17	C	N1-C2-O2	8.03	123.72	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1745	A	C4-C5-C6	8.03	121.01	117.00
1	AA	2260	C	C2-N3-C4	-8.03	115.89	119.90
34	BA	266	G	C5-N7-C8	-8.03	100.29	104.30
1	CA	1692	U	C5-C6-N1	-8.03	118.69	122.70
1	AA	807	G	OP1-P-OP2	-8.02	107.57	119.60
1	CA	765	G	C8-N9-C4	-8.02	103.19	106.40
1	CA	1571	A	N1-C6-N6	-8.02	113.79	118.60
1	AA	423	G	N3-C2-N2	-8.02	114.29	119.90
1	AA	555	G	N1-C6-O6	-8.02	115.09	119.90
1	AA	968	U	C5-C4-O4	-8.02	121.09	125.90
1	AA	336	G	N7-C8-N9	8.01	117.11	113.10
1	AA	805	C	C5-C6-N1	-8.01	116.99	121.00
1	AA	2037	A	C5-C6-N1	-8.01	113.69	117.70
1	AA	2833	A	C8-N9-C4	8.01	109.00	105.80
34	BA	1417	G	C5-C6-N1	8.01	115.51	111.50
1	CA	1397	U	C6-N1-C2	-8.01	116.19	121.00
1	AA	831	A	O4'-C1'-N9	8.01	114.61	108.20
1	AA	1046	A	C5-C6-N1	-8.01	113.69	117.70
1	AA	1050	C	N3-C2-O2	-8.01	116.29	121.90
1	CA	148	C	C5-C6-N1	-8.01	117.00	121.00
1	CA	2196	C	O5'-P-OP2	-8.00	98.50	105.70
34	BA	365	U	N1-C2-N3	8.00	119.70	114.90
1	CA	946	G	O5'-P-OP1	-8.00	98.50	105.70
1	CA	1323	U	O5'-P-OP2	-8.00	98.50	105.70
1	AA	2902	G	C8-N9-C4	-7.99	103.20	106.40
1	AA	911	G	O5'-P-OP2	-7.99	98.51	105.70
1	AA	2800	C	C6-N1-C2	-7.99	117.10	120.30
1	AA	844	C	C2-N3-C4	-7.99	115.91	119.90
1	AA	421	A	N1-C6-N6	-7.98	113.81	118.60
1	AA	1720	U	C2-N1-C1'	-7.98	108.12	117.70
1	AA	2871	G	C5-C6-O6	-7.98	123.81	128.60
1	AA	184	A	OP2-P-O3'	7.98	122.75	105.20
1	CA	1972	A	O5'-P-OP2	-7.97	98.52	105.70
1	AA	2609	G	OP1-P-OP2	-7.97	107.64	119.60
1	AA	53	G	N9-C4-C5	7.97	108.59	105.40
1	AA	1258	A	N7-C8-N9	7.97	117.78	113.80
1	AA	725	C	N3-C4-N4	-7.97	112.42	118.00
1	AA	885	C	N3-C4-C5	7.97	125.09	121.90
34	BA	1384	C	C6-N1-C2	-7.97	117.11	120.30
1	CA	528	A	C5-C6-N1	-7.97	113.72	117.70
1	AA	990	A	O5'-P-OP1	-7.97	98.53	105.70
1	AA	1826	C	O5'-P-OP1	-7.96	98.53	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2384	G	C8-N9-C4	7.96	109.58	106.40
2	AB	104	U	C6-N1-C2	7.96	125.78	121.00
1	CA	2006	C	N3-C2-O2	7.96	127.47	121.90
1	AA	1297	C	N3-C4-C5	7.96	125.08	121.90
1	AA	511	C	C6-N1-C2	-7.96	117.12	120.30
1	AA	1371	G	N3-C4-C5	-7.96	124.62	128.60
1	CA	2582	G	O5'-P-OP2	-7.96	98.54	105.70
1	AA	975	U	C6-N1-C2	-7.95	116.23	121.00
1	AA	2585	C	N1-C2-O2	-7.95	114.13	118.90
1	AA	716	G	OP1-P-OP2	-7.95	107.68	119.60
1	AA	718	C	C6-N1-C2	-7.95	117.12	120.30
1	AA	894	U	N3-C4-O4	-7.95	113.84	119.40
1	CA	2619	C	C6-N1-C2	7.94	123.48	120.30
1	AA	241	G	C8-N9-C4	-7.94	103.22	106.40
1	CA	2330	G	N1-C6-O6	7.94	124.67	119.90
1	AA	324	A	C8-N9-C4	7.94	108.98	105.80
1	AA	1646	C	C2-N3-C4	-7.94	115.93	119.90
34	BA	813	U	O5'-P-OP1	7.93	120.22	110.70
2	AB	81	G	N3-C2-N2	-7.93	114.35	119.90
1	AA	98	U	C5-C4-O4	7.93	130.66	125.90
1	AA	1862	G	OP2-P-O3'	7.93	122.64	105.20
34	BA	868	C	O5'-P-OP2	7.93	120.22	110.70
34	BA	770	C	OP1-P-OP2	-7.93	107.71	119.60
1	AA	858	U	N1-C2-N3	7.93	119.66	114.90
1	AA	426	G	C2-N3-C4	-7.92	107.94	111.90
1	AA	893	C	C2-N3-C4	-7.92	115.94	119.90
34	BA	804	U	N3-C4-O4	-7.92	113.85	119.40
1	AA	1329	G	N1-C6-O6	7.92	124.65	119.90
1	CA	2586	C	N3-C2-O2	-7.92	116.35	121.90
34	BA	345	C	N1-C2-O2	7.92	123.65	118.90
1	AA	992	G	N9-C4-C5	7.92	108.57	105.40
1	AA	1741	C	C5-C6-N1	-7.91	117.04	121.00
1	CA	962	G	N1-C6-O6	7.91	124.65	119.90
1	AA	2749	G	N7-C8-N9	-7.91	109.14	113.10
34	BA	1530	G	N3-C4-C5	7.91	132.55	128.60
1	AA	616	G	C4-C5-N7	-7.91	107.64	110.80
1	CA	2330	G	C5-C6-O6	-7.91	123.86	128.60
1	AA	1000	C	N3-C4-C5	-7.90	118.74	121.90
1	AA	1605	A	C4-C5-N7	7.90	114.65	110.70
1	AA	578	U	C5-C4-O4	-7.89	121.17	125.90
1	AA	831	A	N7-C8-N9	-7.89	109.85	113.80
1	AA	1188	A	C4-C5-N7	7.89	114.65	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2279	A	O5'-P-OP1	-7.89	98.60	105.70
1	AA	147	U	C5-C6-N1	-7.89	118.75	122.70
1	AA	2783	G	C8-N9-C4	7.89	109.56	106.40
2	AB	99	G	N3-C2-N2	7.89	125.42	119.90
1	AA	797	A	O5'-P-OP2	-7.89	98.60	105.70
1	CA	1199	U	O5'-P-OP1	-7.89	98.60	105.70
1	AA	738	C	N1-C2-N3	7.88	124.72	119.20
1	AA	1656	A	C5-C6-N6	-7.88	117.39	123.70
20	AW	11	ARG	NE-CZ-NH1	7.88	124.24	120.30
34	DA	503	C	C6-N1-C2	-7.88	117.15	120.30
1	AA	418	G	C8-N9-C4	7.88	109.55	106.40
1	AA	1694	G	N1-C2-N2	7.88	123.29	116.20
1	AA	2067	C	C6-N1-C2	7.88	123.45	120.30
1	CA	1558	A	N1-C6-N6	7.88	123.33	118.60
1	AA	770	G	C5-C6-N1	7.87	115.44	111.50
1	AA	2559	U	C4-C5-C6	7.87	124.42	119.70
1	AA	1020	C	N1-C2-N3	7.87	124.71	119.20
1	AA	1740	U	C5-C6-N1	-7.87	118.77	122.70
1	AA	1318	A	O4'-C1'-N9	7.87	114.49	108.20
1	AA	2047	C	C5-C6-N1	-7.87	117.07	121.00
1	AA	74	G	C5-C6-O6	-7.86	123.88	128.60
1	CA	189	G	O5'-P-OP2	-7.86	98.62	105.70
1	AA	552	C	N1-C2-N3	7.86	124.70	119.20
1	AA	1815	A	OP1-P-OP2	7.86	131.39	119.60
34	BA	769	G	O5'-P-OP1	7.86	120.13	110.70
1	CA	2599	G	N1-C6-O6	-7.86	115.18	119.90
1	AA	894	U	C6-N1-C1'	7.86	132.20	121.20
1	AA	2348	A	N9-C4-C5	-7.86	102.66	105.80
1	AA	46	C	C2-N3-C4	-7.85	115.97	119.90
1	AA	757	G	N1-C6-O6	7.85	124.61	119.90
1	AA	1249	A	C5-C6-N6	-7.85	117.42	123.70
1	AA	1843	A	C5-C6-N1	7.84	121.62	117.70
1	AA	2605	U	N3-C4-O4	-7.84	113.91	119.40
1	AA	918	U	N3-C2-O2	7.84	127.69	122.20
1	AA	1231	G	N3-C2-N2	7.84	125.39	119.90
56	BW	36	A	C5-C6-N1	-7.84	113.78	117.70
1	AA	2045	G	C5-N7-C8	-7.84	100.38	104.30
34	BA	1436	U	O5'-P-OP2	-7.84	98.64	105.70
56	BW	47	U	C2-N1-C1'	7.84	127.11	117.70
1	AA	593	G	C2-N3-C4	7.84	115.82	111.90
1	AA	718	C	O5'-P-OP1	-7.84	98.65	105.70
1	AA	1784	G	N1-C2-N2	-7.84	109.15	116.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1026	A	C2-N3-C4	-7.83	106.68	110.60
1	AA	525	G	O5'-P-OP2	-7.83	98.65	105.70
1	AA	545	G	C5-N7-C8	7.83	108.21	104.30
1	AA	2301	G	C4-C5-N7	7.83	113.93	110.80
1	CA	2287	A	C2-N3-C4	-7.83	106.69	110.60
1	CA	195	A	P-O3'-C3'	7.83	129.09	119.70
1	AA	1505	C	C5-C6-N1	-7.82	117.09	121.00
17	AT	95	ARG	NE-CZ-NH1	-7.82	116.39	120.30
1	CA	2013	A	C4-C5-C6	7.82	120.91	117.00
1	AA	1709	C	C5-C6-N1	-7.82	117.09	121.00
1	AA	1742	G	OP1-P-OP2	7.82	131.33	119.60
1	AA	114	C	C5-C6-N1	-7.81	117.09	121.00
1	AA	1830	G	N9-C4-C5	7.81	108.52	105.40
1	AA	2025	G	C2-N3-C4	7.81	115.80	111.90
1	AA	1287	A	O5'-P-OP2	7.81	120.07	110.70
1	CA	2261	C	O5'-P-OP1	7.81	120.07	110.70
1	AA	788	G	OP1-P-OP2	-7.80	107.89	119.60
1	AA	1683	C	C4-C5-C6	7.80	121.30	117.40
1	CA	2503	A	C5-C6-N6	-7.80	117.46	123.70
1	AA	744	C	C5-C6-N1	-7.80	117.10	121.00
1	AA	1952	G	O5'-P-OP2	-7.80	98.68	105.70
1	AA	1832	G	OP1-P-OP2	7.80	131.30	119.60
1	AA	2407	C	N3-C4-C5	7.80	125.02	121.90
1	AA	1377	A	C5-C6-N6	7.80	129.94	123.70
1	AA	630	U	O5'-P-OP1	-7.79	98.69	105.70
1	AA	714	U	N1-C2-O2	-7.79	117.35	122.80
1	AA	2014	G	C8-N9-C4	-7.79	103.28	106.40
1	CA	277	C	C2-N1-C1'	7.79	127.37	118.80
1	CA	2043	C	O5'-P-OP2	-7.79	98.69	105.70
34	DA	912	C	C6-N1-C2	7.79	123.42	120.30
1	AA	2070	G	C8-N9-C4	7.79	109.52	106.40
1	CA	2554	U	O5'-P-OP2	7.79	120.04	110.70
1	AA	1645	C	C4-C5-C6	7.78	121.29	117.40
1	AA	1817	A	OP1-P-O3'	7.78	122.32	105.20
1	AA	2221	A	O5'-P-OP2	7.78	120.04	110.70
1	AA	2351	G	N1-C2-N3	7.78	128.57	123.90
34	BA	365	U	C6-N1-C1'	7.78	132.09	121.20
1	CA	1369	G	C8-N9-C4	-7.78	103.29	106.40
1	AA	1011	G	C5-C6-O6	7.77	133.26	128.60
1	AA	1749	G	C4-C5-N7	-7.77	107.69	110.80
1	AA	1661	C	N3-C2-O2	-7.77	116.46	121.90
1	AA	254	A	N7-C8-N9	7.77	117.68	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	414	C	O5'-P-OP2	-7.77	98.71	105.70
1	CA	1142(A)	A	N3-C4-C5	7.77	132.24	126.80
1	AA	1020	C	OP1-P-OP2	-7.77	107.95	119.60
1	AA	534	C	N1-C2-O2	7.76	123.56	118.90
1	AA	579	G	N1-C6-O6	-7.76	115.24	119.90
1	AA	976	G	O5'-P-OP1	-7.76	98.71	105.70
1	AA	2537	G	N1-C2-N2	-7.76	109.21	116.20
1	AA	2702	C	N3-C4-N4	7.76	123.43	118.00
1	AA	1052	C	C2-N1-C1'	-7.76	110.26	118.80
1	AA	1317	G	C4-C5-N7	7.76	113.91	110.80
34	BA	1484	C	C6-N1-C2	7.76	123.41	120.30
1	CA	2695	C	C6-N1-C2	7.76	123.40	120.30
1	AA	2873	C	N1-C2-O2	-7.76	114.25	118.90
1	AA	1022	C	N3-C4-C5	7.75	125.00	121.90
1	AA	783	C	C5-C4-N4	-7.75	114.77	120.20
1	AA	23	G	C4-C5-N7	-7.75	107.70	110.80
1	AA	1741	C	C2-N3-C4	-7.75	116.02	119.90
1	CA	856	C	C6-N1-C2	-7.75	117.20	120.30
1	AA	1425	A	N1-C6-N6	7.75	123.25	118.60
1	AA	1646	C	C5-C6-N1	-7.75	117.13	121.00
1	AA	2902	G	C4-C5-N7	7.75	113.90	110.80
1	CA	679	C	N3-C2-O2	7.75	127.33	121.90
1	AA	1543	U	N1-C2-N3	7.75	119.55	114.90
34	BA	1030(B)	C	N1-C2-O2	7.75	123.55	118.90
34	BA	1030(B)	C	C2-N1-C1'	7.74	127.32	118.80
34	BA	1417	G	C5-C6-O6	-7.74	123.95	128.60
1	AA	2063	U	N1-C2-N3	7.74	119.55	114.90
34	BA	290	C	C6-N1-C2	7.74	123.40	120.30
1	AA	325	G	N7-C8-N9	-7.74	109.23	113.10
1	AA	992	G	C4-C5-N7	-7.74	107.70	110.80
1	CA	792	G	O4'-C1'-N9	-7.74	102.01	108.20
1	AA	2083	G	C8-N9-C4	-7.74	103.31	106.40
1	AA	1020	C	N3-C4-N4	-7.74	112.58	118.00
1	CA	732	C	N1-C2-O2	7.73	123.54	118.90
1	AA	847	A	C8-N9-C4	-7.73	102.71	105.80
1	AA	802	C	C4-C5-C6	7.73	121.27	117.40
1	CA	2193	G	OP2-P-O3'	-7.73	88.19	105.20
1	AA	2098	U	C5-C6-N1	-7.73	118.84	122.70
1	AA	1849	U	C5-C6-N1	-7.73	118.84	122.70
1	AA	2660	C	N1-C2-N3	7.73	124.61	119.20
1	AA	2470	G	C6-N1-C2	-7.72	120.47	125.10
1	AA	1665	G	C5-C6-N1	7.72	115.36	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1851	U	N3-C2-O2	-7.72	116.80	122.20
1	AA	622	G	C8-N9-C4	7.72	109.49	106.40
1	AA	343	C	N3-C4-C5	-7.72	118.81	121.90
1	CA	1988	C	N1-C2-O2	7.72	123.53	118.90
1	AA	241	G	N1-C6-O6	7.72	124.53	119.90
1	AA	1418	U	N3-C4-O4	7.72	124.80	119.40
1	AA	22	C	N3-C4-N4	-7.71	112.60	118.00
1	AA	1097	G	O5'-P-OP2	7.71	119.96	110.70
1	CA	265	A	C5-N7-C8	-7.71	100.04	103.90
1	AA	23	G	C5-N7-C8	7.71	108.16	104.30
1	AA	2506	G	C5-C6-O6	7.71	133.23	128.60
1	CA	2286	A	C4-C5-N7	7.71	114.56	110.70
1	CA	2581	G	O4'-C1'-N9	7.71	114.37	108.20
1	AA	1987	C	C2-N1-C1'	-7.71	110.32	118.80
2	AB	54	G	N7-C8-N9	7.71	116.95	113.10
1	AA	892	G	O4'-C1'-N9	7.71	114.37	108.20
1	CA	2570	G	C4-C5-N7	-7.71	107.72	110.80
1	CA	2234	G	N1-C6-O6	7.71	124.52	119.90
1	AA	907	U	N3-C4-O4	-7.70	114.01	119.40
1	AA	1816	A	C6-N1-C2	-7.70	113.98	118.60
1	CA	1963	U	C2-N1-C1'	7.70	126.94	117.70
1	AA	1274	G	N3-C2-N2	7.70	125.29	119.90
1	AA	2828	G	N1-C6-O6	-7.70	115.28	119.90
1	AA	1710	C	C5-C6-N1	-7.70	117.15	121.00
1	AA	1275	G	O5'-P-OP2	-7.69	98.78	105.70
1	AA	2684	G	C5-C6-O6	-7.69	123.99	128.60
1	CA	1284	A	C5-C6-N6	-7.69	117.55	123.70
1	AA	2277	U	C5-C6-N1	-7.69	118.86	122.70
1	AA	1837	C	O5'-P-OP1	-7.69	98.78	105.70
1	AA	2785	C	C2-N3-C4	-7.69	116.06	119.90
1	CA	2824	C	C5-C6-N1	-7.69	117.16	121.00
1	AA	605	G	N1-C6-O6	-7.68	115.29	119.90
1	AA	2783	G	N9-C4-C5	-7.68	102.33	105.40
1	AA	1358	U	C5-C6-N1	-7.68	118.86	122.70
34	BA	756	C	N3-C4-C5	7.68	124.97	121.90
34	BA	1064	G	N3-C4-N9	-7.68	121.39	126.00
1	AA	1071	G	O5'-P-OP1	-7.68	98.79	105.70
1	AA	2279	A	C2-N3-C4	7.68	114.44	110.60
1	AA	231	G	C5-C6-O6	7.68	133.21	128.60
1	AA	1341	C	C5-C6-N1	-7.68	117.16	121.00
1	AA	2545	A	C8-N9-C4	7.68	108.87	105.80
1	AA	1811	A	N7-C8-N9	7.67	117.64	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1961	U	N3-C4-O4	-7.67	114.03	119.40
1	AA	2265	G	O5'-P-OP1	7.67	119.91	110.70
1	CA	1955	U	N1-C2-O2	-7.67	117.43	122.80
1	AA	500	G	N9-C4-C5	7.67	108.47	105.40
1	AA	625	G	N9-C4-C5	-7.67	102.33	105.40
1	AA	1802	C	C2-N3-C4	-7.67	116.07	119.90
1	CA	2732	G	N1-C6-O6	-7.67	115.30	119.90
1	AA	49	U	N1-C2-O2	7.67	128.17	122.80
56	DW	17	C	C2-N1-C1'	7.67	127.23	118.80
1	CA	2856	C	C5-C6-N1	7.67	124.83	121.00
1	AA	530	A	OP1-P-O3'	7.66	122.06	105.20
1	AA	822	G	O5'-P-OP2	-7.66	98.80	105.70
1	AA	1605	A	C5-C6-N1	-7.66	113.87	117.70
2	AB	38	C	N1-C2-O2	7.66	123.50	118.90
1	AA	726	C	OP1-P-OP2	-7.66	108.11	119.60
1	AA	1703	C	O5'-P-OP1	7.66	119.89	110.70
1	AA	1859	G	C4-C5-N7	-7.66	107.74	110.80
1	CA	1953	A	O5'-P-OP1	-7.66	98.81	105.70
1	AA	787	U	C5-C4-O4	7.66	130.50	125.90
1	AA	2054	G	N7-C8-N9	-7.66	109.27	113.10
1	AA	1864	U	C5-C6-N1	-7.65	118.87	122.70
1	CA	765	G	N7-C8-N9	7.65	116.93	113.10
1	AA	1439	A	C2-N3-C4	7.65	114.43	110.60
1	CA	1288	U	O5'-P-OP1	-7.65	98.81	105.70
1	AA	1464	G	O5'-P-OP1	-7.65	98.81	105.70
34	BA	345	C	C6-N1-C2	-7.65	117.24	120.30
1	CA	2067	G	N9-C4-C5	7.65	108.46	105.40
1	AA	731	G	C5-C6-O6	-7.65	124.01	128.60
34	BA	296	U	C5-C6-N1	-7.65	118.88	122.70
1	CA	1605	C	C4-C5-C6	7.65	121.22	117.40
1	AA	126	C	C2-N3-C4	-7.65	116.08	119.90
1	CA	945	A	N3-C4-C5	7.65	132.15	126.80
1	AA	2058	C	O5'-P-OP1	-7.64	98.82	105.70
1	AA	2653	G	O5'-P-OP1	-7.64	98.82	105.70
1	AA	555	G	C4-C5-C6	-7.64	114.22	118.80
1	AA	1478	C	N1-C2-O2	-7.64	114.31	118.90
1	AA	1665	G	O5'-P-OP2	-7.64	98.82	105.70
1	AA	2889	C	O5'-P-OP2	-7.64	98.82	105.70
1	AA	2548	G	C5-C6-O6	7.64	133.19	128.60
1	AA	2701	U	C6-N1-C2	-7.64	116.42	121.00
1	AA	192	C	N3-C2-O2	7.64	127.25	121.90
1	AA	2263	G	OP1-P-OP2	-7.64	108.14	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	96	C	O5'-P-OP2	-7.64	98.83	105.70
1	AA	2737	C	C4-C5-C6	7.63	121.22	117.40
1	AA	550	U	C5-C6-N1	-7.63	118.88	122.70
1	AA	1682	G	C5-C6-O6	-7.63	124.02	128.60
1	AA	782	A	O5'-P-OP1	-7.63	98.83	105.70
34	BA	355	C	N1-C2-O2	7.63	123.48	118.90
1	CA	1314	C	O5'-P-OP1	-7.63	98.83	105.70
1	AA	1445	C	N3-C2-O2	7.63	127.24	121.90
1	AA	2736	C	N3-C4-C5	7.63	124.95	121.90
1	AA	195	U	C5-C6-N1	-7.62	118.89	122.70
1	AA	1016	C	N3-C4-C5	-7.62	118.85	121.90
1	AA	1022	C	OP1-P-OP2	-7.62	108.17	119.60
1	AA	2505	U	C5-C4-O4	7.62	130.47	125.90
1	AA	883	G	N3-C4-C5	-7.62	124.79	128.60
2	CB	56	G	N3-C4-C5	-7.62	124.79	128.60
1	AA	1659	G	C5-C6-O6	-7.62	124.03	128.60
1	CA	1352	U	C5-C6-N1	-7.62	118.89	122.70
34	BA	266	G	N1-C6-O6	7.62	124.47	119.90
1	AA	1784	G	N9-C4-C5	-7.61	102.36	105.40
1	AA	2014	G	P-O3'-C3'	7.61	128.83	119.70
1	AA	2359	C	C6-N1-C2	-7.61	117.26	120.30
1	AA	2901	A	C8-N9-C4	7.61	108.84	105.80
34	BA	1517	G	O5'-P-OP2	-7.61	98.85	105.70
34	DA	509	A	C8-N9-C4	-7.61	102.76	105.80
1	AA	1303	C	N3-C4-C5	-7.61	118.86	121.90
1	AA	2014	G	C2'-C3'-O3'	7.61	126.23	109.50
1	AA	2066	C	N1-C2-N3	7.61	124.53	119.20
1	AA	2776	G	C6-C5-N7	-7.61	125.84	130.40
34	BA	924	C	N3-C2-O2	-7.61	116.58	121.90
1	AA	256	C	C6-N1-C2	7.61	123.34	120.30
1	AA	1743	G	O5'-P-OP2	-7.61	98.85	105.70
1	AA	2504	U	N1-C2-N3	7.61	119.46	114.90
1	AA	823	G	N7-C8-N9	-7.60	109.30	113.10
1	CA	1288	U	N3-C2-O2	-7.60	116.88	122.20
1	CA	1653	G	O5'-P-OP1	-7.60	98.86	105.70
1	AA	1322	A	C5-C6-N1	7.60	121.50	117.70
1	CA	1983	C	C6-N1-C2	7.60	123.34	120.30
1	AA	2383	G	N3-C4-N9	7.59	130.56	126.00
1	CA	1947	C	C5-C6-N1	-7.59	117.20	121.00
1	AA	2502	G	O5'-P-OP2	-7.59	98.87	105.70
34	BA	1523	G	C5-C6-O6	7.59	133.15	128.60
1	AA	775	G	N1-C6-O6	-7.58	115.35	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1329	G	C8-N9-C1'	-7.58	117.14	127.00
1	AA	1695	C	O5'-P-OP1	-7.58	98.88	105.70
1	AA	1830	G	C5-C6-O6	7.58	133.15	128.60
1	AA	1192	C	C6-N1-C2	7.58	123.33	120.30
34	BA	924	C	C6-N1-C2	-7.58	117.27	120.30
1	CA	12	U	C2-N1-C1'	7.58	126.80	117.70
1	CA	1792	G	C8-N9-C4	7.58	109.43	106.40
1	AA	1427	G	C6-N1-C2	-7.58	120.55	125.10
1	AA	2115	G	N9-C4-C5	7.58	108.43	105.40
1	AA	901	G	OP1-P-OP2	-7.57	108.24	119.60
1	AA	1262	C	O5'-P-OP2	7.57	119.79	110.70
1	CA	794	G	O5'-P-OP2	-7.57	98.88	105.70
1	AA	2347	A	O5'-P-OP1	-7.57	98.89	105.70
1	CA	2008	C	N3-C4-C5	7.57	124.93	121.90
1	CA	1266	G	C8-N9-C4	7.57	109.43	106.40
1	AA	980	C	C5-C6-N1	-7.57	117.22	121.00
1	AA	2021	C	N3-C4-C5	7.57	124.93	121.90
1	AA	2707	C	N1-C2-O2	7.57	123.44	118.90
1	CA	1291	C	C6-N1-C2	7.57	123.33	120.30
1	AA	721	G	C4-C5-N7	7.56	113.83	110.80
56	BW	47	U	N1-C2-O2	7.56	128.09	122.80
1	AA	2460	A	O5'-P-OP2	7.56	119.77	110.70
1	AA	1844	G	C8-N9-C4	7.56	109.42	106.40
1	AA	2162	C	N1-C2-O2	7.56	123.43	118.90
1	CA	2012	G	OP2-P-O3'	7.56	121.83	105.20
1	CA	2248	C	N1-C2-O2	-7.56	114.37	118.90
1	AA	1074	A	N1-C6-N6	7.56	123.13	118.60
1	AA	1362	U	O5'-P-OP1	-7.56	98.90	105.70
1	AA	888	A	C2-N3-C4	-7.55	106.82	110.60
1	AA	1251	G	N1-C2-N3	-7.55	119.37	123.90
1	AA	2779	G	C8-N9-C4	-7.55	103.38	106.40
1	AA	2797	C	N3-C4-C5	7.55	124.92	121.90
1	AA	2841	G	N1-C6-O6	-7.55	115.37	119.90
1	AA	1281	G	C6-N1-C2	-7.55	120.57	125.10
1	CA	2269	A	N1-C6-N6	7.55	123.13	118.60
1	AA	240	A	N9-C4-C5	-7.54	102.78	105.80
12	AO	64	ARG	NE-CZ-NH2	-7.54	116.53	120.30
1	AA	888	A	N1-C2-N3	7.54	133.07	129.30
1	AA	121	G	C6-N1-C2	-7.54	120.58	125.10
2	AB	102	A	C2-N3-C4	-7.54	106.83	110.60
1	AA	2062	C	N1-C2-O2	7.54	123.42	118.90
1	AA	1966	U	O5'-P-OP1	-7.54	98.92	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2611	G	C5-C6-N1	7.54	115.27	111.50
1	AA	2277	U	N1-C2-O2	-7.53	117.53	122.80
1	AA	883	G	OP2-P-O3'	7.53	121.77	105.20
1	AA	1018	A	C5-N7-C8	-7.53	100.14	103.90
1	AA	1273	G	C4-C5-N7	-7.53	107.79	110.80
1	AA	2029	C	C6-N1-C2	-7.53	117.29	120.30
34	BA	1513	A	C2-N3-C4	7.53	114.36	110.60
1	AA	1852	A	O5'-P-OP2	7.53	119.73	110.70
1	CA	395	U	O5'-P-OP2	-7.53	98.92	105.70
1	CA	730	C	C6-N1-C2	7.53	123.31	120.30
34	BA	345	C	N3-C2-O2	-7.53	116.63	121.90
1	AA	1007	G	N3-C2-N2	-7.53	114.63	119.90
1	AA	198	C	OP1-P-O3'	-7.52	88.65	105.20
1	AA	2020	G	C8-N9-C4	7.52	109.41	106.40
1	AA	637	U	C5-C4-O4	7.52	130.41	125.90
1	AA	1371	G	C8-N9-C4	-7.52	103.39	106.40
1	AA	2028	C	N3-C4-C5	7.52	124.91	121.90
1	AA	2525	G	N1-C6-O6	7.52	124.41	119.90
1	AA	121	G	C5-C6-N1	7.52	115.26	111.50
1	AA	822	G	O5'-P-OP1	-7.52	98.94	105.70
1	CA	2073	C	N3-C4-C5	7.52	124.91	121.90
1	AA	1645	C	O5'-P-OP2	-7.51	98.94	105.70
1	AA	1688	A	C8-N9-C4	-7.51	102.79	105.80
1	AA	2828	G	N1-C2-N2	-7.51	109.44	116.20
1	AA	1057	G	OP2-P-O3'	7.51	121.72	105.20
1	AA	1831	C	C2-N3-C4	-7.51	116.14	119.90
34	BA	123	C	C6-N1-C2	7.51	123.30	120.30
34	DA	758	G	C5-C6-O6	7.51	133.10	128.60
1	AA	1329	G	C4-N9-C1'	7.51	136.26	126.50
20	AW	17	VAL	CB-CA-C	-7.51	97.13	111.40
1	AA	1710	C	C2-N3-C4	-7.51	116.15	119.90
1	AA	2356	U	O5'-P-OP2	-7.50	98.95	105.70
1	CA	205	G	N7-C8-N9	-7.50	109.35	113.10
1	CA	308	G	O5'-P-OP2	-7.50	98.95	105.70
1	AA	2012	C	N1-C2-O2	-7.50	114.40	118.90
34	BA	924	C	N3-C4-C5	-7.50	118.90	121.90
1	CA	2566	A	O5'-P-OP2	-7.50	98.95	105.70
1	CA	2012	G	C8-N9-C4	-7.50	103.40	106.40
1	AA	2030	C	O5'-P-OP2	-7.50	98.95	105.70
1	CA	1807	G	N7-C8-N9	-7.50	109.35	113.10
56	BW	12	U	O5'-P-OP2	-7.50	98.95	105.70
1	CA	332	A	OP2-P-O3'	7.50	121.69	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2699	C	N1-C2-O2	7.50	123.40	118.90
1	AA	886	U	N3-C4-O4	-7.50	114.15	119.40
1	AA	2362	C	C2-N3-C4	-7.50	116.15	119.90
1	CA	2033	A	O5'-P-OP1	-7.49	98.96	105.70
1	AA	1453	C	N1-C2-O2	-7.49	114.41	118.90
1	AA	2301	G	C5-C6-N1	7.49	115.25	111.50
1	CA	631	A	O5'-P-OP2	-7.49	98.96	105.70
34	DA	890	G	N3-C4-C5	-7.49	124.86	128.60
34	BA	1495	U	O5'-P-OP2	-7.49	98.96	105.70
1	CA	2013	A	N1-C2-N3	7.49	133.04	129.30
1	AA	2612	A	N1-C6-N6	-7.48	114.11	118.60
1	AA	2083	G	C6-N1-C2	-7.48	120.61	125.10
1	AA	2299	A	N7-C8-N9	7.48	117.54	113.80
1	AA	733	G	OP1-P-OP2	7.48	130.81	119.60
1	AA	743	G	N7-C8-N9	-7.48	109.36	113.10
1	AA	778	C	C4-C5-C6	7.47	121.14	117.40
1	CA	2238	G	OP1-P-OP2	7.47	130.81	119.60
1	AA	2709	G	C8-N9-C4	7.47	109.39	106.40
1	AA	1071	G	N1-C2-N3	7.47	128.38	123.90
1	AA	2616	U	C5-C6-N1	-7.47	118.97	122.70
1	CA	687	C	C4-C5-C6	-7.47	113.67	117.40
1	CA	2549	G	C5-C6-O6	-7.47	124.12	128.60
1	AA	209	G	C4-C5-N7	-7.47	107.81	110.80
1	AA	663	G	N1-C2-N2	-7.47	109.48	116.20
1	AA	1000	C	C4-C5-C6	7.47	121.13	117.40
1	AA	2725	A	O5'-P-OP1	-7.47	98.98	105.70
1	CA	446	G	N3-C2-N2	-7.47	114.67	119.90
1	AA	74	G	N1-C6-O6	7.46	124.38	119.90
1	AA	139	A	C4-C5-N7	7.46	114.43	110.70
1	AA	586	G	C6-N1-C2	7.46	129.58	125.10
1	AA	844	C	C5-C4-N4	-7.46	114.98	120.20
1	AA	2067	C	C2-N3-C4	-7.46	116.17	119.90
1	AA	2697	G	C5-N7-C8	7.46	108.03	104.30
1	AA	976	G	C6-N1-C2	-7.46	120.62	125.10
1	AA	2888	U	N1-C2-N3	7.46	119.38	114.90
1	CA	2439	A	O5'-P-OP2	-7.46	98.99	105.70
1	AA	1658	C	C6-N1-C2	7.46	123.28	120.30
1	AA	2221	A	O5'-P-OP1	-7.46	98.99	105.70
1	AA	2855	G	N3-C4-C5	-7.46	124.87	128.60
1	AA	2858	G	O5'-P-OP1	7.46	119.65	110.70
1	CA	1658	C	O5'-P-OP2	-7.46	98.99	105.70
1	AA	2288	G	N1-C6-O6	-7.46	115.43	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1957	C	C5-C4-N4	-7.46	114.98	120.20
1	AA	1208	G	C5-N7-C8	7.46	108.03	104.30
1	AA	2075	G	N9-C4-C5	7.45	108.38	105.40
1	AA	2579	G	C6-C5-N7	7.45	134.87	130.40
1	AA	1665	G	C6-N1-C2	-7.45	120.63	125.10
1	AA	2506	G	C6-N1-C2	7.45	129.57	125.10
1	AA	2539	C	N3-C4-C5	7.45	124.88	121.90
1	AA	2775	G	N3-C2-N2	7.45	125.11	119.90
1	AA	362	G	C5-C6-O6	7.45	133.07	128.60
1	AA	1827	U	N3-C4-O4	7.45	124.61	119.40
1	AA	783	C	N3-C4-C5	7.45	124.88	121.90
1	CA	1646	C	N3-C2-O2	-7.45	116.69	121.90
1	AA	1807	G	N9-C4-C5	-7.45	102.42	105.40
1	CA	702	G	C5-N7-C8	7.45	108.02	104.30
1	AA	1077	G	C5-C6-N1	7.44	115.22	111.50
1	AA	2450	U	N1-C2-O2	-7.44	117.59	122.80
4	CD	229	VAL	CB-CA-C	-7.44	97.26	111.40
1	AA	1655	A	C5-C6-N6	-7.44	117.75	123.70
1	AA	2693	C	C2-N3-C4	-7.44	116.18	119.90
1	AA	2777	A	O5'-P-OP1	-7.44	99.01	105.70
1	AA	2249	G	N3-C2-N2	7.43	125.10	119.90
1	CA	150	C	N1-C2-O2	7.43	123.36	118.90
1	CA	25	U	N3-C2-O2	7.43	127.40	122.20
1	AA	2537	G	N3-C2-N2	7.43	125.10	119.90
1	AA	197	C	C5-C6-N1	-7.43	117.29	121.00
1	AA	2400	A	N1-C2-N3	7.43	133.01	129.30
1	AA	1410	G	N3-C4-C5	-7.42	124.89	128.60
1	CA	827	U	C6-N1-C2	7.42	125.45	121.00
1	CA	959	A	C8-N9-C4	-7.42	102.83	105.80
1	AA	2638	C	C6-N1-C2	7.42	123.27	120.30
1	AA	2718	G	N3-C2-N2	7.42	125.09	119.90
1	CA	2611	U	N1-C2-O2	-7.42	117.61	122.80
1	AA	1011	G	N1-C6-O6	-7.42	115.45	119.90
2	AB	97	G	N3-C2-N2	-7.42	114.71	119.90
1	CA	2580	U	C5-C6-N1	-7.42	118.99	122.70
34	DA	769	G	N1-C6-O6	7.42	124.35	119.90
1	CA	2406	U	O5'-P-OP2	7.42	119.60	110.70
1	CA	2638	G	O5'-P-OP1	-7.41	99.03	105.70
1	CA	563	G	N3-C2-N2	-7.41	114.71	119.90
1	CA	2779	U	N1-C2-O2	7.41	127.99	122.80
34	DA	912	C	C5-C6-N1	-7.41	117.30	121.00
1	AA	96	C	N1-C2-O2	7.40	123.34	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	181	C	C5-C6-N1	-7.40	117.30	121.00
1	AA	1384	G	OP1-P-OP2	-7.40	108.49	119.60
1	AA	1817	A	C8-N9-C4	7.40	108.76	105.80
1	CA	767	U	N3-C2-O2	-7.40	117.02	122.20
1	CA	1142(A)	A	C5-N7-C8	-7.40	100.20	103.90
1	AA	2757	G	N1-C2-N2	7.40	122.86	116.20
1	CA	2442	C	N3-C2-O2	-7.40	116.72	121.90
1	AA	496	A	C8-N9-C4	-7.40	102.84	105.80
1	AA	2697	G	C2-N3-C4	7.40	115.60	111.90
34	BA	885	G	O5'-P-OP2	-7.40	99.04	105.70
1	AA	805	C	N1-C2-O2	7.40	123.34	118.90
1	CA	2586	C	N3-C4-C5	7.40	124.86	121.90
1	AA	2341	G	N7-C8-N9	-7.39	109.40	113.10
1	AA	2400	A	C8-N9-C4	-7.39	102.84	105.80
1	CA	1978	A	OP2-P-O3'	7.39	121.47	105.20
1	AA	726	C	C6-N1-C2	7.39	123.26	120.30
1	AA	1742	G	C4-C5-N7	7.39	113.76	110.80
1	CA	1698	A	C5-N7-C8	-7.39	100.20	103.90
1	AA	2033	U	C4-C5-C6	7.39	124.13	119.70
1	AA	2455	C	N1-C2-N3	7.39	124.37	119.20
1	AA	995	G	C5-C6-N1	7.39	115.19	111.50
1	AA	2591	C	C5-C4-N4	-7.39	115.03	120.20
1	CA	2060	A	N1-C6-N6	-7.39	114.17	118.60
1	AA	2608	U	N1-C2-N3	7.39	119.33	114.90
1	CA	37	C	N3-C4-N4	7.39	123.17	118.00
1	AA	1454	C	N1-C2-O2	-7.39	114.47	118.90
1	AA	1686	U	N3-C2-O2	-7.39	117.03	122.20
34	BA	578	C	C6-N1-C2	-7.39	117.34	120.30
1	AA	534	C	C2-N3-C4	7.38	123.59	119.90
1	AA	2398	C	N3-C4-C5	-7.38	118.95	121.90
1	AA	2634	C	C2-N3-C4	-7.38	116.21	119.90
2	AB	77	U	C2-N3-C4	-7.38	122.57	127.00
1	AA	2077	C	O5'-P-OP2	-7.38	99.06	105.70
1	AA	1244	U	N3-C2-O2	-7.38	117.03	122.20
1	AA	2533	C	N1-C2-O2	-7.38	114.47	118.90
1	CA	997	G	O5'-P-OP1	-7.38	99.06	105.70
1	CA	2409	G	C6-C5-N7	-7.38	125.97	130.40
1	AA	597	C	C6-N1-C2	-7.38	117.35	120.30
34	BA	28	G	C5-C6-O6	-7.38	124.17	128.60
1	AA	1394	G	C5-C6-O6	-7.37	124.17	128.60
1	AA	2562	G	C6-C5-N7	-7.37	125.98	130.40
1	AA	240	A	N7-C8-N9	-7.37	110.11	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1653	G	N3-C4-C5	-7.37	124.92	128.60
1	CA	2582	G	N1-C2-N2	-7.37	109.57	116.20
1	AA	761	U	C5-C4-O4	-7.37	121.48	125.90
1	AA	1315	A	N1-C2-N3	-7.37	125.61	129.30
1	AA	2615	G	C8-N9-C4	7.37	109.35	106.40
34	BA	1513	A	N1-C2-N3	-7.37	125.62	129.30
1	AA	1987	C	N3-C2-O2	7.37	127.06	121.90
1	AA	2034	G	O5'-P-OP2	-7.36	99.07	105.70
1	CA	2049	G	O5'-P-OP2	-7.36	99.07	105.70
1	CA	2877	G	O5'-P-OP1	7.36	119.53	110.70
1	AA	1723	A	O5'-P-OP2	-7.36	99.08	105.70
34	BA	119	A	O5'-P-OP2	-7.36	99.08	105.70
1	CA	941	A	O5'-P-OP2	-7.36	99.08	105.70
1	CA	2554	U	O5'-P-OP1	-7.36	99.08	105.70
1	CA	265	A	N7-C8-N9	7.36	117.48	113.80
1	AA	16	G	C6-N1-C2	-7.36	120.69	125.10
1	AA	1804	A	C8-N9-C4	7.36	108.74	105.80
1	AA	2532	C	C5-C6-N1	-7.36	117.32	121.00
34	BA	1505	G	C5-C6-O6	7.36	133.01	128.60
1	AA	776	G	C4-N9-C1'	7.36	136.06	126.50
1	AA	1656	A	O5'-P-OP1	-7.36	99.08	105.70
1	AA	841	G	N7-C8-N9	-7.35	109.42	113.10
1	CA	2611	U	O5'-P-OP1	-7.35	99.08	105.70
1	AA	874	U	O5'-P-OP2	-7.35	99.08	105.70
1	AA	1698	G	C8-N9-C4	-7.35	103.46	106.40
1	AA	60	G	N3-C2-N2	-7.34	114.76	119.90
1	AA	2450	U	C5-C6-N1	-7.34	119.03	122.70
1	AA	19	C	N1-C2-O2	-7.34	114.49	118.90
1	AA	2369	U	N3-C2-O2	-7.34	117.06	122.20
1	AA	581	G	C8-N9-C4	7.34	109.33	106.40
1	AA	716	G	C8-N9-C4	7.34	109.34	106.40
1	AA	906	G	O4'-C1'-N9	-7.34	102.33	108.20
1	AA	2303	U	N3-C4-C5	-7.34	110.20	114.60
1	CA	2441	C	O5'-P-OP1	-7.34	99.09	105.70
1	AA	19	C	C2-N3-C4	-7.34	116.23	119.90
1	AA	2085	C	C6-N1-C2	7.34	123.23	120.30
2	AB	103	G	C6-C5-N7	-7.34	126.00	130.40
1	AA	516	G	C5-C6-O6	7.34	133.00	128.60
1	AA	651	U	O5'-P-OP2	-7.34	99.10	105.70
1	AA	721	G	C5-N7-C8	-7.34	100.63	104.30
1	AA	1621	C	OP2-P-O3'	7.33	121.33	105.20
1	AA	2858	G	C4-C5-N7	7.33	113.73	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	18	C	N3-C4-C5	-7.33	118.97	121.90
1	AA	1742	G	C5-C6-O6	-7.33	124.20	128.60
1	AA	2573	A	O5'-P-OP1	-7.33	99.10	105.70
2	AB	49	C	N1-C2-O2	-7.33	114.50	118.90
2	AB	101	G	O5'-P-OP2	-7.33	99.10	105.70
1	AA	1055	A	C8-N9-C4	7.33	108.73	105.80
1	AA	601	A	C5-C6-N6	7.33	129.56	123.70
1	CA	1497	U	O4'-C1'-N1	7.33	114.06	108.20
1	AA	2539	C	C5-C4-N4	-7.33	115.07	120.20
34	BA	770	C	N3-C4-N4	-7.33	112.87	118.00
1	AA	609	A	C2-N3-C4	7.32	114.26	110.60
1	AA	1033	G	C2-N3-C4	7.32	115.56	111.90
34	BA	803	G	C2-N3-C4	-7.32	108.24	111.90
1	CA	186	G	O5'-P-OP1	7.32	119.49	110.70
1	AA	1640	G	C5-C6-O6	-7.32	124.21	128.60
1	AA	2690	C	C5-C6-N1	7.32	124.66	121.00
34	BA	836	G	C5-C6-O6	-7.32	124.21	128.60
1	AA	1329	G	N3-C4-N9	7.32	130.39	126.00
1	CA	1774	C	O5'-P-OP2	-7.32	99.11	105.70
1	CA	2512	C	C2-N3-C4	-7.32	116.24	119.90
1	AA	1605	A	C5-N7-C8	-7.32	100.24	103.90
34	BA	875	C	O5'-P-OP2	-7.32	99.11	105.70
1	CA	2058	A	C2-N3-C4	-7.32	106.94	110.60
1	CA	1812	A	C2-N3-C4	7.31	114.26	110.60
1	AA	860	U	N3-C4-C5	7.31	118.99	114.60
1	AA	2535	G	C5-C6-O6	7.31	132.99	128.60
2	AB	104	U	C5-C4-O4	-7.31	121.52	125.90
1	AA	1875	C	C6-N1-C2	-7.31	117.38	120.30
1	AA	544	U	N1-C2-N3	7.30	119.28	114.90
1	AA	918	U	N1-C2-O2	-7.30	117.69	122.80
1	AA	1986	G	C4-C5-N7	7.30	113.72	110.80
1	AA	121	G	C5-C6-O6	-7.30	124.22	128.60
1	AA	839	G	C5-C6-N1	7.30	115.15	111.50
1	AA	1632	A	N1-C2-N3	7.30	132.95	129.30
2	AB	69	G	C8-N9-C4	7.30	109.32	106.40
1	AA	2702	C	C5-C4-N4	-7.30	115.09	120.20
1	AA	2799	U	N3-C4-O4	-7.30	114.29	119.40
1	CA	670	A	O4'-C1'-N9	-7.30	102.36	108.20
1	AA	2162	C	C2-N1-C1'	7.29	126.82	118.80
1	AA	340	C	N3-C4-C5	7.29	124.82	121.90
1	AA	1979	C	N3-C4-C5	-7.29	118.98	121.90
1	AA	2818	U	N3-C4-C5	-7.29	110.22	114.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	708	C	C2-N3-C4	-7.29	116.25	119.90
1	AA	2858	G	C5-C6-O6	-7.29	124.22	128.60
1	AA	1261	G	C4-C5-N7	7.29	113.72	110.80
1	AA	2448	G	C5-C6-O6	7.29	132.97	128.60
1	CA	1698	A	C4-N9-C1'	7.29	139.42	126.30
1	AA	972	A	C8-N9-C4	-7.29	102.89	105.80
1	AA	348	A	C8-N9-C4	7.28	108.71	105.80
1	AA	2423	A	C8-N9-C4	7.28	108.71	105.80
1	AA	181	C	C2-N3-C4	-7.28	116.26	119.90
1	CA	34	C	C6-N1-C2	-7.28	117.39	120.30
1	AA	497	A	C2-N3-C4	-7.28	106.96	110.60
1	AA	1835	C	OP2-P-O3'	7.28	121.22	105.20
34	BA	365	U	C4-C5-C6	7.28	124.07	119.70
1	AA	2535	G	N1-C6-O6	-7.28	115.53	119.90
1	AA	860	U	C2-N3-C4	-7.27	122.64	127.00
1	AA	1312	G	N1-C6-O6	-7.27	115.53	119.90
1	CA	732	C	N3-C2-O2	-7.27	116.81	121.90
1	AA	725	C	C6-N1-C2	-7.27	117.39	120.30
1	AA	452	G	C8-N9-C4	7.27	109.31	106.40
1	AA	795	G	O5'-P-OP1	-7.27	99.16	105.70
1	AA	1612	C	C6-N1-C2	7.27	123.21	120.30
1	AA	2641	A	C8-N9-C4	-7.27	102.89	105.80
1	AA	793	A	C8-N9-C4	-7.27	102.89	105.80
1	CA	1780	A	O5'-P-OP2	-7.27	99.16	105.70
1	AA	585	U	C5-C6-N1	-7.27	119.07	122.70
1	AA	2302	G	OP1-P-OP2	-7.27	108.70	119.60
1	AA	100	G	C8-N9-C4	7.26	109.31	106.40
1	AA	137	G	C8-N9-C4	-7.26	103.50	106.40
1	AA	1076	G	C4-C5-N7	7.26	113.71	110.80
1	AA	1691	C	C2-N1-C1'	7.26	126.79	118.80
1	CA	1328	G	O5'-P-OP2	7.26	119.42	110.70
1	AA	1926	G	N3-C4-C5	-7.26	124.97	128.60
1	AA	2589	A	C8-N9-C4	7.26	108.70	105.80
56	BW	73	A	O4'-C1'-N9	7.26	114.01	108.20
1	CA	2258	C	C5-C6-N1	-7.26	117.37	121.00
1	AA	129	G	N3-C2-N2	7.26	124.98	119.90
1	AA	1704	C	C5-C6-N1	-7.26	117.37	121.00
1	AA	2282	G	C5-C6-N1	7.26	115.13	111.50
34	BA	816	A	OP1-P-O3'	7.26	121.16	105.20
1	CA	1611	C	OP1-P-OP2	-7.26	108.72	119.60
1	CA	1842	G	O5'-P-OP2	-7.25	99.17	105.70
1	AA	283	G	OP1-P-O3'	7.25	121.16	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1711	A	N1-C6-N6	7.25	122.95	118.60
1	CA	2063	C	N1-C2-O2	-7.25	114.55	118.90
1	AA	2639	G	C4-C5-N7	7.25	113.70	110.80
1	AA	1378	G	N3-C4-C5	-7.25	124.97	128.60
2	AB	38	C	N3-C2-O2	-7.25	116.83	121.90
1	AA	518	G	O5'-P-OP2	-7.25	99.18	105.70
1	AA	1296	G	N1-C6-O6	-7.25	115.55	119.90
1	CA	242	G	C8-N9-C4	7.25	109.30	106.40
1	CA	1960	A	OP1-P-OP2	7.25	130.47	119.60
1	AA	2039	U	C5-C4-O4	7.25	130.25	125.90
1	CA	572	A	N1-C6-N6	7.25	122.95	118.60
1	CA	2006	C	O5'-P-OP2	-7.25	99.18	105.70
1	AA	2252	C	OP1-P-OP2	-7.25	108.73	119.60
1	CA	238	C	O5'-P-OP1	-7.24	99.18	105.70
1	CA	2382	G	O5'-P-OP2	-7.24	99.18	105.70
1	AA	1314	A	N1-C2-N3	7.24	132.92	129.30
1	AA	1986	G	N9-C4-C5	-7.24	102.50	105.40
1	AA	2437	A	C5-N7-C8	-7.24	100.28	103.90
1	AA	1951	G	O5'-P-OP1	-7.24	99.18	105.70
1	AA	2044	U	N3-C4-O4	7.24	124.47	119.40
1	CA	330	A	N9-C4-C5	-7.24	102.90	105.80
1	AA	2040	G	N1-C2-N3	-7.24	119.56	123.90
1	AA	2045	G	O5'-P-OP2	7.24	119.39	110.70
2	AB	7	G	N1-C6-O6	7.24	124.24	119.90
1	AA	256	C	N3-C4-C5	7.24	124.80	121.90
1	AA	1711	A	C2-N3-C4	-7.24	106.98	110.60
1	AA	2025	G	C4-C5-N7	-7.24	107.91	110.80
1	AA	1397	C	C5-C4-N4	7.23	125.26	120.20
1	AA	2108	U	O5'-P-OP2	-7.23	99.19	105.70
1	AA	2576	A	N1-C2-N3	7.23	132.92	129.30
1	AA	2701	U	OP1-P-O3'	7.23	121.11	105.20
1	AA	1308	A	O5'-P-OP1	-7.23	99.19	105.70
1	AA	2636	G	C4-C5-N7	-7.23	107.91	110.80
1	CA	1776	G	N3-C4-N9	7.23	130.34	126.00
34	DA	266	G	N1-C6-O6	7.23	124.24	119.90
1	AA	751	G	C5-C6-N1	7.23	115.11	111.50
1	AA	1987	C	C6-N1-C2	7.23	123.19	120.30
1	AA	227	C	C6-N1-C2	7.22	123.19	120.30
1	AA	905	U	N3-C4-C5	7.22	118.94	114.60
1	AA	1514	C	O5'-P-OP2	7.22	119.37	110.70
1	CA	774	A	O5'-P-OP2	-7.22	99.20	105.70
1	AA	734	C	N3-C4-N4	-7.22	112.94	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1208	G	C4-C5-N7	-7.22	107.91	110.80
1	AA	74	G	C4-C5-N7	7.22	113.69	110.80
34	BA	1511	G	N7-C8-N9	-7.22	109.49	113.10
1	CA	246	C	O5'-P-OP1	-7.22	99.20	105.70
1	AA	2871	G	N1-C6-O6	7.22	124.23	119.90
1	AA	2636	G	C6-N1-C2	-7.22	120.77	125.10
1	AA	1371	G	C5-C6-O6	7.21	132.93	128.60
1	CA	2561	A	OP1-P-OP2	7.21	130.42	119.60
1	AA	324	A	N1-C2-N3	-7.21	125.69	129.30
1	AA	2601	A	N1-C6-N6	-7.21	114.27	118.60
1	AA	2750	G	N9-C4-C5	7.21	108.28	105.40
1	AA	2892	A	OP1-P-OP2	7.21	130.42	119.60
1	CA	1768	U	C5-C4-O4	7.21	130.23	125.90
1	AA	45	C	N3-C2-O2	7.21	126.94	121.90
1	CA	73	A	N1-C6-N6	-7.21	114.27	118.60
1	CA	448	U	N3-C2-O2	-7.21	117.15	122.20
1	CA	1286	A	O5'-P-OP2	-7.21	99.21	105.70
1	CA	1626	G	C5-C6-O6	-7.21	124.28	128.60
34	DA	1414	U	C5-C6-N1	-7.21	119.10	122.70
1	AA	2548	G	C5-N7-C8	7.21	107.90	104.30
2	AB	97	G	N1-C2-N2	7.21	122.69	116.20
34	BA	1523	G	OP1-P-OP2	7.21	130.41	119.60
1	AA	2491	G	O5'-P-OP2	-7.21	99.22	105.70
1	AA	2638	C	C2-N3-C4	-7.21	116.30	119.90
1	AA	2602	A	C2-N3-C4	-7.20	107.00	110.60
1	AA	2833	A	C4-C5-C6	7.20	120.60	117.00
1	AA	823	G	N3-C4-C5	-7.20	125.00	128.60
1	AA	1042	A	C5-C6-N1	-7.20	114.10	117.70
34	BA	514	C	N1-C2-O2	-7.20	114.58	118.90
1	AA	260	A	O5'-P-OP2	-7.20	99.22	105.70
1	AA	545	G	C4-C5-N7	-7.20	107.92	110.80
1	AA	591	U	C5-C4-O4	-7.20	121.58	125.90
1	AA	2001	C	N3-C4-C5	7.20	124.78	121.90
1	AA	614	C	C2-N3-C4	-7.20	116.30	119.90
1	AA	1464	G	O5'-P-OP2	7.20	119.33	110.70
1	AA	975	U	N3-C4-O4	7.19	124.44	119.40
1	AA	541	C	N3-C4-N4	-7.19	112.97	118.00
1	AA	1837	C	C5-C6-N1	-7.19	117.40	121.00
1	AA	2566	U	O5'-P-OP2	7.19	119.33	110.70
1	CA	2607	G	N1-C2-N2	-7.19	109.73	116.20
1	AA	2535	G	N1-C2-N2	-7.19	109.73	116.20
1	AA	753	A	C2-N3-C4	-7.19	107.01	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2696	U	C2-N3-C4	-7.19	122.69	127.00
1	CA	2286	A	C2-N3-C4	-7.19	107.01	110.60
1	CA	1313	U	C2-N1-C1'	7.19	126.32	117.70
1	CA	1451	C	N1-C2-O2	-7.19	114.59	118.90
1	AA	1701	A	OP1-P-OP2	7.18	130.38	119.60
1	AA	2400	A	C5-C6-N6	7.18	129.45	123.70
1	AA	2655	G	C8-N9-C4	-7.18	103.53	106.40
1	AA	2078	G	N7-C8-N9	7.18	116.69	113.10
1	AA	2824	C	N1-C2-O2	-7.18	114.59	118.90
1	CA	562	U	N1-C2-N3	7.18	119.21	114.90
1	AA	823	G	N3-C4-N9	7.18	130.31	126.00
34	BA	836	G	N1-C6-O6	7.18	124.21	119.90
1	AA	1065	U	C2-N3-C4	-7.18	122.69	127.00
1	AA	2484	G	C5-N7-C8	7.18	107.89	104.30
1	AA	2779	G	O5'-P-OP2	-7.18	99.24	105.70
1	AA	1563	G	N1-C6-O6	7.18	124.21	119.90
1	AA	30	G	N1-C6-O6	-7.17	115.59	119.90
1	CA	1947	C	OP1-P-OP2	-7.17	108.84	119.60
1	AA	36	G	C4-C5-N7	-7.17	107.93	110.80
1	AA	575	G	C5-C6-O6	7.17	132.90	128.60
1	CA	59	U	OP2-P-O3'	7.17	120.98	105.20
1	CA	762	U	C5-C4-O4	-7.17	121.60	125.90
1	AA	555	G	C8-N9-C1'	7.17	136.32	127.00
1	AA	910	A	C8-N9-C4	-7.17	102.93	105.80
1	CA	474	G	C5-C6-O6	7.17	132.90	128.60
1	AA	1695	C	O5'-P-OP2	7.17	119.30	110.70
1	CA	2540	C	O5'-P-OP2	-7.17	99.25	105.70
1	AA	627	G	N1-C6-O6	-7.17	115.60	119.90
1	AA	2285	A	N1-C6-N6	7.16	122.90	118.60
1	CA	1698	A	N1-C6-N6	7.16	122.90	118.60
1	AA	1699	A	C8-N9-C4	-7.16	102.94	105.80
34	DA	295	C	C6-N1-C2	7.16	123.16	120.30
1	AA	715	G	C4-C5-N7	7.16	113.66	110.80
1	AA	989	G	N1-C6-O6	7.16	124.19	119.90
1	AA	1650	C	C2-N3-C4	-7.16	116.32	119.90
1	AA	2758	C	N1-C2-O2	-7.16	114.61	118.90
1	AA	2727	G	N1-C2-N3	-7.16	119.61	123.90
34	BA	519	C	C6-N1-C2	7.16	123.16	120.30
1	AA	1483	C	C6-N1-C2	-7.16	117.44	120.30
1	CA	1791	A	OP1-P-OP2	-7.15	108.87	119.60
1	AA	1704	C	OP2-P-O3'	7.15	120.93	105.20
1	CA	265	A	O4'-C1'-N9	7.15	113.92	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	659	C	C6-N1-C2	7.15	123.16	120.30
1	CA	1611	C	C6-N1-C2	7.15	123.16	120.30
1	AA	2711	C	N3-C4-C5	7.14	124.76	121.90
56	BW	1	G	C8-N9-C4	-7.14	103.54	106.40
1	AA	1032	C	N3-C2-O2	7.14	126.90	121.90
1	AA	2641	A	O4'-C1'-N9	7.14	113.91	108.20
1	AA	726	C	N3-C4-C5	7.14	124.76	121.90
1	AA	1312	G	N1-C2-N2	-7.14	109.78	116.20
1	AA	1839	U	O5'-P-OP1	-7.14	99.27	105.70
1	AA	2633	A	N1-C6-N6	-7.14	114.32	118.60
1	CA	1323	U	C6-N1-C2	7.14	125.28	121.00
1	CA	2818	G	N7-C8-N9	-7.14	109.53	113.10
1	AA	2455	C	O5'-P-OP1	-7.14	99.28	105.70
1	AA	721	G	C5-C6-O6	-7.14	124.32	128.60
1	AA	419	C	O5'-P-OP1	-7.13	99.28	105.70
1	AA	486	A	C8-N9-C4	7.13	108.65	105.80
1	AA	1060	U	OP2-P-O3'	7.13	120.89	105.20
1	AA	85	C	N3-C4-N4	-7.13	113.01	118.00
1	AA	710	G	N1-C6-O6	7.13	124.18	119.90
1	AA	1858	C	N1-C2-O2	7.13	123.18	118.90
1	CA	678	C	C6-N1-C2	7.13	123.15	120.30
2	AB	56	G	O5'-P-OP2	-7.13	99.29	105.70
1	CA	945	A	O4'-C1'-N9	7.13	113.90	108.20
1	AA	38	A	C4-C5-N7	-7.12	107.14	110.70
1	AA	1543	U	O5'-P-OP2	-7.12	99.29	105.70
1	AA	597	C	C5-C4-N4	7.12	125.19	120.20
1	CA	659	C	C2-N3-C4	-7.12	116.34	119.90
1	AA	975	U	N1-C2-N3	7.12	119.17	114.90
1	CA	918	A	O5'-P-OP1	-7.12	99.29	105.70
1	AA	1059	C	N1-C2-O2	-7.12	114.63	118.90
56	BW	47	U	C6-N1-C2	-7.12	116.73	121.00
1	CA	903	C	C6-N1-C2	7.12	123.15	120.30
1	CA	2046	G	N7-C8-N9	-7.12	109.54	113.10
1	AA	529	U	O5'-P-OP2	-7.11	99.30	105.70
1	AA	702	A	N3-C4-N9	-7.11	121.71	127.40
1	CA	1293	C	N3-C4-C5	7.11	124.75	121.90
1	AA	803	C	N1-C2-O2	-7.11	114.63	118.90
1	AA	1357	G	N7-C8-N9	7.11	116.66	113.10
1	AA	2757	G	N3-C4-C5	7.11	132.16	128.60
1	CA	593	G	N1-C6-O6	7.11	124.17	119.90
1	CA	408	G	C8-N9-C4	7.11	109.24	106.40
1	AA	1041	C	C5-C4-N4	-7.11	115.23	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	582	G	C8-N9-C4	-7.10	103.56	106.40
34	BA	899	C	C6-N1-C2	7.10	123.14	120.30
1	AA	1080	G	O5'-P-OP2	-7.10	99.31	105.70
1	CA	976	C	C6-N1-C2	7.10	123.14	120.30
1	AA	26	G	C5-C6-N1	7.10	115.05	111.50
1	AA	674	G	N3-C4-C5	-7.10	125.05	128.60
1	AA	953	U	OP2-P-O3'	7.10	120.81	105.20
1	CA	2003	G	OP2-P-O3'	7.10	120.82	105.20
1	AA	474	U	N3-C2-O2	-7.09	117.23	122.20
1	AA	640	A	C8-N9-C4	-7.09	102.96	105.80
1	AA	987	G	N7-C8-N9	-7.09	109.55	113.10
1	AA	990	A	O4'-C1'-N9	7.09	113.88	108.20
1	CA	945	A	C4-C5-N7	7.09	114.25	110.70
34	BA	1505	G	C4-C5-N7	-7.09	107.96	110.80
1	AA	2527	C	N1-C2-O2	-7.09	114.64	118.90
2	AB	72	G	C5-C6-O6	-7.09	124.35	128.60
1	AA	345	G	C5-C6-O6	-7.09	124.35	128.60
34	BA	900	A	O5'-P-OP2	7.09	119.20	110.70
1	CA	2576	G	N3-C4-N9	7.09	130.25	126.00
1	AA	111	G	N1-C6-O6	7.08	124.15	119.90
29	C5	16	ARG	NE-CZ-NH1	7.08	123.84	120.30
1	AA	107	G	N1-C6-O6	-7.08	115.65	119.90
1	AA	1747	A	O5'-P-OP1	-7.08	99.33	105.70
1	AA	1979	C	C4-C5-C6	7.08	120.94	117.40
1	AA	354	A	N1-C6-N6	7.08	122.85	118.60
1	CA	690	G	N3-C4-N9	7.08	130.25	126.00
1	CA	1899	G	OP1-P-O3'	7.08	120.77	105.20
1	AA	1698	G	N7-C8-N9	7.08	116.64	113.10
1	AA	2064	A	C2-N3-C4	-7.07	107.06	110.60
1	CA	2062	A	C2-N3-C4	7.07	114.14	110.60
34	DA	266	G	C5-C6-O6	-7.07	124.36	128.60
1	AA	1784	G	C8-N9-C4	7.07	109.23	106.40
1	AA	871	A	C2-N3-C4	7.07	114.14	110.60
1	AA	2065	C	N3-C4-C5	-7.07	119.07	121.90
1	AA	2576	A	C6-N1-C2	-7.07	114.36	118.60
1	AA	206	G	N3-C2-N2	-7.07	114.95	119.90
1	AA	552	C	O4'-C1'-N1	7.07	113.85	108.20
1	AA	1010	C	C5-C4-N4	7.07	125.15	120.20
1	AA	2000	A	C2-N3-C4	7.07	114.13	110.60
34	DA	353	A	N1-C6-N6	7.07	122.84	118.60
1	AA	901	G	C5-C6-O6	7.07	132.84	128.60
1	CA	600	G	C8-N9-C4	-7.07	103.57	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1807	G	N3-C4-N9	7.06	130.24	126.00
1	AA	1053	C	O5'-P-OP2	-7.06	99.34	105.70
1	AA	46	C	N3-C4-C5	7.06	124.72	121.90
1	AA	860	U	N3-C2-O2	-7.06	117.26	122.20
1	AA	2631	C	C4-C5-C6	7.06	120.93	117.40
2	AB	91	C	N3-C4-C5	7.06	124.72	121.90
1	AA	539	A	OP1-P-OP2	7.06	130.19	119.60
34	BA	1064	G	C4-N9-C1'	-7.06	117.33	126.50
1	AA	1926	G	C4-C5-N7	-7.05	107.98	110.80
1	AA	354	A	C6-N1-C2	7.05	122.83	118.60
1	AA	71	U	N1-C2-O2	-7.05	117.86	122.80
1	AA	89	U	C5-C4-O4	7.05	130.13	125.90
1	AA	343	C	N3-C2-O2	-7.05	116.97	121.90
1	AA	1402	G	N1-C6-O6	-7.05	115.67	119.90
1	AA	2592	U	C2-N3-C4	-7.05	122.77	127.00
1	AA	1715	A	N1-C6-N6	-7.05	114.37	118.60
1	AA	1718	U	C6-N1-C2	7.05	125.23	121.00
1	AA	2058	C	C2-N3-C4	-7.05	116.38	119.90
1	AA	822	G	OP1-P-O3'	7.04	120.70	105.20
1	CA	1558	A	C5-C6-N1	-7.04	114.18	117.70
1	AA	2409	G	C5-C6-N1	7.04	115.02	111.50
1	CA	141	A	C6-C5-N7	-7.04	127.37	132.30
1	CA	270	A	C2-N3-C4	-7.04	107.08	110.60
1	CA	2038	G	C5-C6-O6	7.04	132.82	128.60
1	AA	600	G	N3-C4-N9	7.04	130.22	126.00
1	AA	1281	G	C5-C6-O6	-7.04	124.38	128.60
1	AA	2601	A	N9-C4-C5	7.04	108.61	105.80
1	AA	347	G	O4'-C1'-N9	7.03	113.83	108.20
1	AA	778	C	N1-C2-N3	7.03	124.12	119.20
1	AA	2532	C	O5'-P-OP2	-7.03	99.37	105.70
34	BA	1422	G	O5'-P-OP1	-7.03	99.37	105.70
1	AA	753	A	N1-C6-N6	7.03	122.82	118.60
1	AA	1617	A	N1-C6-N6	7.03	122.82	118.60
1	AA	2781	C	N3-C4-C5	7.03	124.71	121.90
1	AA	38	A	N7-C8-N9	-7.03	110.28	113.80
1	AA	2048	C	OP1-P-OP2	7.03	130.15	119.60
1	CA	2258	C	N1-C2-O2	-7.03	114.68	118.90
1	AA	2750	G	C5-C6-O6	7.03	132.82	128.60
1	CA	1795	C	C6-N1-C2	7.03	123.11	120.30
1	AA	1057	G	O5'-P-OP2	-7.03	99.38	105.70
1	AA	176	G	C8-N9-C4	-7.03	103.59	106.40
1	AA	352	U	C2-N3-C4	-7.02	122.79	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2619	G	C4-C5-N7	7.02	113.61	110.80
1	AA	2836	A	OP1-P-OP2	-7.02	109.06	119.60
1	AA	1243	U	N3-C2-O2	-7.02	117.28	122.20
2	AB	12	C	N3-C4-C5	7.02	124.71	121.90
1	CA	2805	G	O4'-C1'-N9	7.02	113.82	108.20
1	AA	1805	C	C2-N3-C4	-7.02	116.39	119.90
1	AA	2684	G	N1-C6-O6	7.02	124.11	119.90
34	BA	1516	G	O5'-P-OP2	-7.02	99.38	105.70
1	AA	2833	A	N1-C6-N6	7.02	122.81	118.60
1	AA	491	G	N7-C8-N9	7.01	116.61	113.10
1	AA	1258	A	N1-C2-N3	7.01	132.81	129.30
1	AA	2368	C	C2-N3-C4	-7.01	116.39	119.90
1	AA	2865	C	OP2-P-O3'	7.01	120.62	105.20
1	CA	446	G	N1-C6-O6	7.01	124.11	119.90
34	DA	1501	C	C6-N1-C2	7.01	123.11	120.30
1	AA	1257	G	O5'-P-OP2	-7.01	99.39	105.70
1	CA	2437	U	OP2-P-O3'	7.01	120.62	105.20
1	AA	1475	G	C8-N9-C4	-7.01	103.60	106.40
1	AA	2060	G	C5-C6-O6	7.01	132.81	128.60
34	BA	1506	U	N3-C4-O4	7.01	124.31	119.40
2	AB	81	G	N9-C4-C5	7.00	108.20	105.40
1	AA	1921	G	C5-N7-C8	-7.00	100.80	104.30
1	AA	586	G	C4-C5-N7	-7.00	108.00	110.80
34	BA	643	C	O5'-P-OP1	-7.00	99.40	105.70
1	AA	1255	A	C2-N3-C4	7.00	114.10	110.60
1	AA	1752	G	N1-C6-O6	-7.00	115.70	119.90
1	AA	535	C	OP1-P-OP2	7.00	130.10	119.60
1	AA	1872	U	N1-C2-N3	7.00	119.10	114.90
1	CA	1812	A	C5-C6-N1	7.00	121.20	117.70
1	CA	2708	G	N3-C2-N2	7.00	124.80	119.90
1	AA	1595	C	N1-C2-O2	-7.00	114.70	118.90
1	AA	2236	G	C5-N7-C8	-7.00	100.80	104.30
1	AA	73	A	C2-N3-C4	-7.00	107.10	110.60
1	AA	2103	C	N1-C2-N3	6.99	124.10	119.20
1	CA	450	G	N9-C4-C5	6.99	108.20	105.40
1	CA	1776	G	N3-C4-C5	-6.99	125.10	128.60
1	AA	1666	G	C4-C5-N7	-6.99	108.00	110.80
1	AA	2450	U	C2-N3-C4	-6.99	122.81	127.00
1	AA	1255	A	OP2-P-O3'	6.99	120.58	105.20
1	CA	1314	C	C2-N1-C1'	6.99	126.49	118.80
1	AA	188	A	O5'-P-OP1	-6.99	99.41	105.70
1	AA	1038	C	C5-C6-N1	-6.99	117.50	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2067	G	C4-C5-N7	-6.99	108.00	110.80
1	AA	2025	G	C6-C5-N7	6.99	134.59	130.40
1	CA	1623	G	N1-C2-N2	6.99	122.49	116.20
56	DW	76	A	N1-C6-N6	6.99	122.79	118.60
1	AA	10	G	C5-C6-O6	6.98	132.79	128.60
1	AA	77	A	N1-C2-N3	6.98	132.79	129.30
1	AA	2609	G	C8-N9-C4	6.98	109.19	106.40
34	BA	1402	C	N3-C4-C5	-6.98	119.11	121.90
1	CA	241	A	O5'-P-OP2	-6.98	99.42	105.70
1	AA	1058	U	C6-N1-C2	6.98	125.19	121.00
34	BA	575	G	O4'-C1'-N9	-6.98	102.62	108.20
1	AA	451	G	C8-N9-C4	6.98	109.19	106.40
1	AA	1026	A	C8-N9-C4	6.98	108.59	105.80
1	AA	1670	G	N1-C6-O6	-6.98	115.71	119.90
1	CA	198	C	O5'-P-OP1	-6.98	99.42	105.70
1	AA	235	C	C5-C6-N1	-6.98	117.51	121.00
1	AA	2073	A	C5-N7-C8	6.98	107.39	103.90
1	AA	1017	G	C8-N9-C4	-6.98	103.61	106.40
1	AA	179	A	OP1-P-OP2	6.97	130.06	119.60
1	AA	804	U	N3-C4-C5	6.97	118.78	114.60
1	AA	1721	G	N3-C4-C5	-6.97	125.11	128.60
34	BA	605	U	C5-C4-O4	6.97	130.08	125.90
1	AA	1003	U	N3-C4-C5	-6.97	110.42	114.60
1	CA	933	A	O4'-C1'-N9	6.97	113.78	108.20
1	AA	865	G	N7-C8-N9	-6.97	109.61	113.10
1	AA	1307	C	C4-C5-C6	6.97	120.89	117.40
1	AA	1806	U	N3-C4-O4	-6.97	114.52	119.40
1	AA	1816	A	N1-C2-N3	6.97	132.78	129.30
1	CA	563	G	C5-N7-C8	-6.97	100.82	104.30
1	AA	2064	A	N7-C8-N9	-6.96	110.32	113.80
1	CA	797	C	C6-N1-C2	-6.96	117.51	120.30
2	AB	100	A	OP1-P-OP2	6.96	130.04	119.60
1	CA	2674	G	O5'-P-OP2	-6.96	99.43	105.70
1	AA	1264	G	O5'-P-OP2	-6.96	99.43	105.70
1	AA	1695	C	C5-C4-N4	6.96	125.07	120.20
1	AA	2040	G	C6-N1-C2	6.96	129.28	125.10
2	AB	54	G	C5-N7-C8	-6.96	100.82	104.30
34	BA	912	C	N1-C2-O2	-6.96	114.72	118.90
1	AA	1027	A	N7-C8-N9	-6.96	110.32	113.80
1	AA	1289	G	N1-C6-O6	-6.96	115.72	119.90
56	BW	35	A	OP2-P-O3'	6.96	120.51	105.20
1	CA	1339	G	N1-C6-O6	-6.96	115.72	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1836	C	O5'-P-OP1	6.96	119.05	110.70
1	AA	2497	G	N3-C2-N2	6.96	124.77	119.90
2	AB	81	G	N1-C2-N3	6.96	128.07	123.90
1	CA	1949	G	N1-C6-O6	6.96	124.07	119.90
1	AA	586	G	C5-C6-O6	6.96	132.77	128.60
1	AA	716	G	OP2-P-O3'	6.96	120.50	105.20
2	AB	68	C	C6-N1-C2	6.96	123.08	120.30
1	CA	783	A	O5'-P-OP2	-6.96	99.44	105.70
1	CA	1554	A	N1-C6-N6	-6.95	114.43	118.60
1	AA	1698	G	C5-C6-N1	6.95	114.98	111.50
1	AA	598	A	N7-C8-N9	-6.95	110.33	113.80
1	AA	797	A	C5-N7-C8	-6.95	100.42	103.90
1	CA	148	C	C6-N1-C2	6.95	123.08	120.30
1	AA	2015	U	N1-C2-N3	6.95	119.07	114.90
1	AA	2053	A	C8-N9-C4	-6.95	103.02	105.80
1	CA	383	U	O4'-C1'-N1	6.95	113.76	108.20
1	AA	38	A	C5-C6-N1	6.95	121.17	117.70
1	AA	411	U	N3-C4-O4	-6.95	114.54	119.40
1	AA	736	A	O5'-P-OP2	-6.95	99.45	105.70
34	BA	800	G	O5'-P-OP2	-6.95	99.45	105.70
1	AA	328	G	C5-C6-O6	-6.94	124.43	128.60
1	AA	841	G	C5-C6-N1	6.94	114.97	111.50
1	AA	1815	A	C8-N9-C4	6.94	108.58	105.80
1	CA	2256	G	O5'-P-OP2	-6.94	99.45	105.70
1	AA	851	A	C2-N3-C4	-6.94	107.13	110.60
1	AA	1988	A	N3-C4-C5	6.94	131.66	126.80
1	AA	1276	C	C6-N1-C2	-6.94	117.53	120.30
1	AA	1360	C	N1-C2-N3	6.94	124.06	119.20
1	AA	2580	C	C6-N1-C2	6.94	123.08	120.30
1	AA	2511	C	C5-C6-N1	6.94	124.47	121.00
1	AA	2594	G	OP1-P-OP2	-6.93	109.20	119.60
1	CA	665	C	C5-C6-N1	6.93	124.47	121.00
1	CA	2503	A	O5'-P-OP1	-6.93	99.46	105.70
1	AA	1737	A	N9-C4-C5	6.93	108.57	105.80
1	AA	1832	G	O5'-P-OP1	-6.93	99.46	105.70
1	CA	571	A	C2-N3-C4	6.93	114.07	110.60
34	DA	58	C	C6-N1-C2	-6.93	117.53	120.30
34	DA	898	G	C5-C6-N1	6.93	114.97	111.50
1	AA	1806	U	O5'-P-OP2	-6.93	99.46	105.70
34	DA	882	C	N3-C2-O2	-6.93	117.05	121.90
1	AA	2522	C	O5'-P-OP2	-6.93	99.46	105.70
1	AA	2650	G	C6-C5-N7	-6.93	126.24	130.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2385	C	N1-C2-O2	-6.93	114.74	118.90
1	CA	1963	U	N1-C2-O2	6.93	127.65	122.80
34	DA	677	U	O5'-P-OP1	-6.93	99.47	105.70
1	AA	663	G	N3-C2-N2	6.92	124.75	119.90
1	AA	1085	G	C8-N9-C4	6.92	109.17	106.40
1	AA	2712	C	OP1-P-OP2	6.92	129.99	119.60
1	CA	2498	C	N3-C2-O2	6.92	126.75	121.90
34	DA	289	G	C5-C6-O6	-6.92	124.45	128.60
1	CA	2625	G	C5-C6-O6	-6.92	124.45	128.60
1	CA	2827	C	C6-N1-C2	6.92	123.07	120.30
1	AA	1405	A	C4-C5-C6	-6.92	113.54	117.00
1	AA	1312	G	C6-N1-C2	-6.92	120.95	125.10
1	AA	2348	A	C8-N9-C4	6.92	108.57	105.80
1	AA	2703	C	OP1-P-OP2	-6.92	109.22	119.60
34	BA	1498	U	N3-C4-C5	6.92	118.75	114.60
1	AA	199	C	C6-N1-C2	6.92	123.07	120.30
1	AA	1666	G	N9-C4-C5	6.92	108.17	105.40
1	AA	2434	A	N1-C2-N3	6.92	132.76	129.30
1	AA	2890	C	O5'-P-OP2	-6.92	99.47	105.70
1	AA	478	G	N1-C2-N2	6.92	122.42	116.20
1	AA	1316	C	N3-C4-C5	6.92	124.67	121.90
1	AA	2601	A	N3-C4-N9	-6.92	121.87	127.40
1	AA	557	A	N1-C2-N3	6.91	132.76	129.30
1	AA	1277	G	N9-C4-C5	6.91	108.17	105.40
1	AA	1411	A	C2-N3-C4	-6.91	107.14	110.60
1	AA	627	G	C4-C5-N7	-6.91	108.03	110.80
1	AA	1029	A	C8-N9-C4	6.91	108.56	105.80
1	AA	1398	U	O5'-P-OP1	-6.91	99.48	105.70
1	AA	2787	C	C5-C6-N1	-6.91	117.54	121.00
1	AA	2797	C	C5-C4-N4	-6.91	115.36	120.20
1	AA	351	G	N1-C2-N2	-6.91	109.98	116.20
1	AA	555	G	C8-N9-C4	-6.91	103.64	106.40
2	AB	41	U	C4-C5-C6	6.91	123.84	119.70
34	DA	780	A	C8-N9-C4	-6.91	103.04	105.80
2	AB	13	A	OP1-P-OP2	6.91	129.96	119.60
1	CA	1353	A	OP1-P-OP2	6.91	129.96	119.60
1	AA	2457	G	N1-C6-O6	-6.91	115.76	119.90
1	AA	2546	A	C5-N7-C8	-6.91	100.45	103.90
1	AA	182	U	O5'-P-OP1	6.90	118.98	110.70
1	AA	2609	G	OP2-P-O3'	6.90	120.39	105.20
1	CA	442	G	N3-C2-N2	-6.90	115.07	119.90
1	AA	29	U	N1-C2-O2	-6.90	117.97	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1455	C	N1-C2-N3	6.90	124.03	119.20
1	CA	567	A	C5-C6-N6	-6.90	118.18	123.70
1	AA	1052	C	N3-C2-O2	6.90	126.73	121.90
1	CA	515	A	N1-C2-N3	6.90	132.75	129.30
1	AA	455	A	C5'-C4'-C3'	-6.89	104.97	116.00
1	AA	2556	G	C5-N7-C8	-6.89	100.85	104.30
1	AA	2718	G	C5-N7-C8	6.89	107.75	104.30
1	AA	737	G	N1-C6-O6	-6.89	115.77	119.90
1	AA	1006	C	C2-N1-C1'	-6.89	111.22	118.80
1	AA	1660	A	OP2-P-O3'	6.89	120.36	105.20
1	AA	2458	G	C2-N3-C4	6.89	115.35	111.90
1	AA	545	G	N7-C8-N9	-6.89	109.66	113.10
1	CA	749	C	C6-N1-C2	6.89	123.06	120.30
1	AA	1047	A	O5'-P-OP1	-6.89	99.50	105.70
1	AA	728	G	C8-N9-C4	6.88	109.15	106.40
1	AA	1788	U	C5-C6-N1	-6.88	119.26	122.70
1	AA	912	C	N3-C4-C5	-6.88	119.15	121.90
1	AA	1155	C	C4-C5-C6	-6.88	113.96	117.40
1	AA	2108	U	C5-C6-N1	-6.88	119.26	122.70
34	BA	893	C	OP1-P-OP2	-6.88	109.28	119.60
1	AA	415	G	O5'-P-OP2	-6.88	99.51	105.70
1	AA	2252	C	C5-C4-N4	-6.88	115.38	120.20
1	AA	2279	A	O4'-C1'-N9	-6.88	102.70	108.20
1	AA	106	U	C2-N3-C4	-6.88	122.87	127.00
1	AA	708	C	C5-C6-N1	-6.88	117.56	121.00
1	AA	728	G	N7-C8-N9	-6.88	109.66	113.10
1	CA	2674	G	C5-C6-O6	6.88	132.73	128.60
1	AA	1428	G	N1-C6-O6	6.88	124.03	119.90
1	AA	1746	G	O5'-P-OP1	6.88	118.95	110.70
1	AA	53	G	C8-N9-C4	-6.88	103.65	106.40
1	AA	1157	A	C5-N7-C8	-6.88	100.46	103.90
1	AA	1342	G	N1-C2-N2	-6.88	110.01	116.20
1	AA	1024	G	C4-C5-N7	-6.87	108.05	110.80
1	AA	2639	G	N3-C2-N2	6.87	124.71	119.90
1	AA	2888	U	N3-C2-O2	-6.87	117.39	122.20
1	CA	2552	U	O5'-P-OP2	-6.87	99.52	105.70
1	AA	478	G	N3-C4-N9	-6.87	121.88	126.00
1	AA	2604	G	O5'-P-OP1	-6.87	99.52	105.70
1	AA	2694	U	O5'-P-OP2	-6.87	99.52	105.70
1	CA	2442	C	C6-N1-C2	-6.87	117.55	120.30
1	AA	357	G	N7-C8-N9	6.87	116.53	113.10
1	AA	712	C	N3-C2-O2	-6.87	117.09	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	120	G	C6-N1-C2	-6.87	120.98	125.10
1	AA	803	C	C6-N1-C2	-6.87	117.55	120.30
1	CA	748	G	C5-C6-O6	6.87	132.72	128.60
1	CA	811	U	C5-C6-N1	-6.87	119.27	122.70
1	CA	2732	G	C4-C5-N7	-6.87	108.05	110.80
1	AA	2780	C	C2-N3-C4	-6.87	116.47	119.90
15	AR	114	VAL	CB-CA-C	-6.87	98.36	111.40
1	AA	98	U	N3-C4-O4	-6.86	114.59	119.40
1	AA	1749	G	C6-C5-N7	6.86	134.52	130.40
1	AA	846	G	N3-C4-C5	-6.86	125.17	128.60
1	AA	734	C	C5-C4-N4	6.86	125.00	120.20
1	AA	2078	G	N1-C6-O6	-6.86	115.78	119.90
1	AA	2465	A	OP1-P-OP2	-6.86	109.31	119.60
1	AA	1755	C	C6-N1-C2	6.86	123.04	120.30
1	AA	2446	A	C8-N9-C4	-6.86	103.06	105.80
1	AA	315	C	C5-C6-N1	-6.86	117.57	121.00
1	AA	621	G	O5'-P-OP2	-6.86	99.53	105.70
1	CA	1823	G	C5-C6-O6	6.86	132.72	128.60
1	AA	896	A	C2-N3-C4	-6.86	107.17	110.60
1	AA	2632	C	N1-C2-N3	6.86	124.00	119.20
1	CA	2572	A	C5-N7-C8	6.86	107.33	103.90
1	AA	610	C	C2-N3-C4	-6.85	116.47	119.90
1	AA	1922	A	C2-N3-C4	6.85	114.03	110.60
1	AA	50	G	N3-C4-C5	-6.85	125.17	128.60
1	AA	734	C	N3-C2-O2	-6.85	117.10	121.90
1	AA	1424	A	O5'-P-OP2	-6.85	99.53	105.70
1	AA	730	C	C2-N3-C4	-6.85	116.47	119.90
1	AA	1266	C	N3-C4-C5	-6.85	119.16	121.90
1	AA	2001	C	N1-C2-O2	-6.85	114.79	118.90
1	AA	2562	G	O5'-P-OP1	6.85	118.92	110.70
1	AA	520	G	C4-C5-N7	-6.85	108.06	110.80
1	AA	2546	A	C6-N1-C2	-6.85	114.49	118.60
2	AB	46	A	OP2-P-O3'	6.85	120.26	105.20
1	CA	2617	C	C6-N1-C2	6.85	123.04	120.30
1	AA	989	G	N3-C4-C5	-6.84	125.18	128.60
1	AA	1852	A	C8-N9-C4	-6.84	103.06	105.80
1	AA	2500	A	N1-C6-N6	-6.84	114.50	118.60
1	AA	2651	A	C6-N1-C2	-6.84	114.50	118.60
1	AA	2752	U	O5'-P-OP2	6.84	118.91	110.70
2	AB	32	C	C6-N1-C2	6.84	123.04	120.30
1	AA	637	U	N3-C2-O2	-6.84	117.41	122.20
1	AA	1231	G	N1-C6-O6	-6.84	115.80	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1543	U	C6-N1-C2	-6.84	116.90	121.00
1	AA	464	G	N1-C6-O6	6.84	124.00	119.90
1	AA	776	G	N7-C8-N9	6.84	116.52	113.10
1	AA	2033	U	N3-C4-C5	-6.83	110.50	114.60
1	AA	27	G	O4'-C1'-N9	6.83	113.67	108.20
1	AA	1266	C	N1-C2-N3	6.83	123.98	119.20
1	CA	1812	A	O5'-P-OP1	-6.83	99.55	105.70
1	AA	69	G	OP1-P-OP2	-6.83	109.36	119.60
1	AA	183	G	N9-C4-C5	6.83	108.13	105.40
1	AA	542	C	N3-C4-C5	6.83	124.63	121.90
1	AA	1316	C	OP2-P-O3'	6.83	120.23	105.20
1	AA	2282	G	O5'-P-OP1	-6.83	99.55	105.70
34	BA	811	C	C6-N1-C2	6.83	123.03	120.30
1	AA	129	G	N3-C4-N9	6.83	130.10	126.00
1	AA	148	C	N3-C4-N4	-6.83	113.22	118.00
34	DA	812	C	O5'-P-OP2	-6.83	99.56	105.70
1	AA	2828	G	C2-N3-C4	-6.82	108.49	111.90
1	CA	1554	A	N9-C4-C5	6.82	108.53	105.80
2	CB	79	C	O5'-P-OP2	-6.82	99.56	105.70
56	DW	73	A	C5-N7-C8	-6.82	100.49	103.90
1	AA	82	G	N1-C6-O6	6.82	123.99	119.90
1	CA	2732	G	N9-C4-C5	6.82	108.13	105.40
1	AA	511	C	N3-C4-N4	-6.82	113.23	118.00
1	AA	1683	C	OP1-P-OP2	-6.82	109.37	119.60
1	CA	141	A	N1-C6-N6	6.82	122.69	118.60
1	CA	530	G	N3-C4-N9	-6.82	121.91	126.00
1	AA	254	A	N9-C4-C5	-6.81	103.08	105.80
2	AB	82	G	N9-C4-C5	6.81	108.12	105.40
34	BA	239	U	N3-C2-O2	-6.81	117.43	122.20
1	CA	511	U	O5'-P-OP1	-6.81	99.57	105.70
1	AA	796	C	C6-N1-C2	6.81	123.02	120.30
1	AA	2655	G	C5-C6-O6	6.81	132.69	128.60
1	AA	2718	G	N1-C2-N2	-6.81	110.07	116.20
1	CA	790	C	C5-C4-N4	-6.81	115.43	120.20
1	AA	514	G	N1-C6-O6	-6.81	115.81	119.90
1	AA	663	G	N3-C4-C5	-6.81	125.20	128.60
1	AA	1744	G	C8-N9-C4	6.81	109.12	106.40
1	CA	372	G	O4'-C1'-N9	6.81	113.65	108.20
1	CA	693	C	N3-C4-C5	6.81	124.62	121.90
1	AA	1012	C	C6-N1-C2	-6.81	117.58	120.30
1	AA	2448	G	N1-C6-O6	-6.81	115.82	119.90
1	AA	726	C	C5-C6-N1	-6.80	117.60	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2634	C	N3-C4-C5	6.80	124.62	121.90
1	CA	2194	G	OP1-P-OP2	6.80	129.81	119.60
1	AA	353	G	N9-C4-C5	-6.80	102.68	105.40
2	AB	101	G	C8-N9-C4	6.80	109.12	106.40
1	AA	1749	G	N3-C4-N9	-6.80	121.92	126.00
1	AA	2071	G	C8-N9-C4	-6.80	103.68	106.40
34	DA	692	U	N3-C4-O4	6.80	124.16	119.40
1	AA	1341	C	OP1-P-OP2	6.80	129.80	119.60
1	AA	1796	C	OP2-P-O3'	6.80	120.16	105.20
1	AA	2443	U	C5-C6-N1	-6.80	119.30	122.70
1	CA	2618	G	O5'-P-OP1	6.80	118.86	110.70
1	AA	716	G	C6-N1-C2	-6.80	121.02	125.10
1	AA	410	U	O4'-C1'-N1	6.80	113.64	108.20
1	AA	798	A	C6-N1-C2	-6.80	114.52	118.60
1	AA	1206	G	OP2-P-O3'	6.80	120.15	105.20
2	AB	94	C	C6-N1-C2	-6.80	117.58	120.30
34	BA	1412	C	C5-C6-N1	-6.80	117.60	121.00
1	AA	1212	C	N3-C4-N4	6.79	122.76	118.00
1	AA	780	G	C5-N7-C8	6.79	107.70	104.30
1	AA	1200	G	N3-C4-C5	-6.79	125.20	128.60
1	AA	200	A	O5'-P-OP2	-6.79	99.59	105.70
1	AA	1009	C	N1-C2-O2	6.79	122.97	118.90
1	AA	2054	G	C5-N7-C8	6.79	107.69	104.30
1	CA	2512	C	C5-C6-N1	-6.79	117.61	121.00
1	AA	290	G	N3-C4-C5	6.79	131.99	128.60
1	AA	1919	G	C2-N3-C4	-6.79	108.51	111.90
1	AA	2594	G	C5-C6-O6	-6.79	124.53	128.60
1	AA	227	C	N3-C4-C5	6.79	124.61	121.90
1	AA	413	G	C4-C5-N7	6.79	113.52	110.80
1	CA	450	G	C5-C6-N1	6.79	114.89	111.50
1	AA	865	G	C4-C5-N7	-6.79	108.08	110.80
1	AA	1078	A	N9-C4-C5	-6.79	103.09	105.80
1	AA	2250	G	C5-N7-C8	-6.79	100.91	104.30
1	AA	2525	G	N9-C4-C5	-6.79	102.69	105.40
1	AA	1273	G	N1-C2-N3	6.78	127.97	123.90
1	AA	1505	C	C6-N1-C2	6.78	123.01	120.30
34	BA	1530	G	N3-C2-N2	-6.78	115.15	119.90
1	AA	895	G	N1-C6-O6	-6.78	115.83	119.90
1	AA	2480	G	C5-C6-N1	6.78	114.89	111.50
1	CA	2607	G	N3-C4-N9	6.78	130.07	126.00
1	AA	233	A	N1-C6-N6	6.78	122.67	118.60
1	AA	2435	U	OP1-P-OP2	6.78	129.77	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	AH	149	ARG	NE-CZ-NH1	-6.78	116.91	120.30
1	CA	1211	U	N1-C2-O2	6.78	127.55	122.80
1	CA	1324	G	C8-N9-C4	6.78	109.11	106.40
1	AA	796	C	C4-C5-C6	6.78	120.79	117.40
1	AA	2736	C	C5-C6-N1	-6.78	117.61	121.00
1	AA	719	C	N1-C2-O2	-6.78	114.83	118.90
1	CA	393	C	O5'-P-OP1	-6.78	99.60	105.70
1	CA	2490	G	N3-C4-C5	6.78	131.99	128.60
1	AA	867	A	C5-N7-C8	6.77	107.29	103.90
1	AA	1061	G	C6-C5-N7	-6.77	126.34	130.40
1	AA	1616	A	N1-C6-N6	-6.77	114.54	118.60
1	AA	2518	U	N3-C2-O2	-6.77	117.46	122.20
1	AA	2660	C	C4-C5-C6	6.77	120.79	117.40
1	AA	57	G	C5-C6-N1	6.77	114.89	111.50
1	AA	384	G	C8-N9-C4	6.77	109.11	106.40
1	AA	718	C	N3-C4-N4	-6.77	113.26	118.00
1	AA	2264	G	C8-N9-C1'	6.77	135.80	127.00
1	AA	1579	C	C6-N1-C2	-6.77	117.59	120.30
1	AA	1987	C	N1-C2-O2	-6.77	114.84	118.90
1	CA	2597	G	O5'-P-OP2	-6.77	99.61	105.70
1	AA	1237	G	N1-C6-O6	-6.77	115.84	119.90
1	AA	60	G	N3-C4-N9	-6.76	121.94	126.00
1	AA	1859	G	C5-C6-N1	-6.76	108.12	111.50
1	AA	2852	G	C2-N3-C4	-6.76	108.52	111.90
1	CA	150	C	N3-C4-C5	6.76	124.61	121.90
1	CA	1617	C	C6-N1-C2	-6.76	117.59	120.30
1	AA	907	U	C5-C6-N1	-6.76	119.32	122.70
1	AA	200	A	N1-C2-N3	6.76	132.68	129.30
1	AA	984	G	OP2-P-O3'	6.76	120.07	105.20
1	AA	2718	G	N1-C6-O6	-6.76	115.84	119.90
1	CA	2287	A	C5-N7-C8	-6.76	100.52	103.90
1	AA	60	G	N3-C4-C5	6.76	131.98	128.60
1	AA	1317	G	C6-C5-N7	-6.76	126.34	130.40
1	AA	1627	A	C8-N9-C4	-6.76	103.10	105.80
1	CA	2494	G	O5'-P-OP2	-6.76	99.62	105.70
1	AA	1208	G	N7-C8-N9	-6.76	109.72	113.10
1	AA	1331	G	N9-C4-C5	-6.76	102.70	105.40
1	AA	2115	G	N3-C4-N9	-6.76	121.95	126.00
1	AA	1411	A	N9-C4-C5	-6.75	103.10	105.80
1	AA	2833	A	N7-C8-N9	-6.75	110.42	113.80
2	AB	85	G	N3-C4-C5	-6.75	125.22	128.60
1	AA	887	C	C2-N3-C4	-6.75	116.52	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1736	A	OP2-P-O3'	6.75	120.06	105.20
34	DA	266	G	C5-N7-C8	-6.75	100.92	104.30
1	AA	1033	G	C5-N7-C8	6.75	107.67	104.30
1	AA	2540	U	N3-C2-O2	6.75	126.92	122.20
1	AA	2650	G	C5-C6-O6	-6.75	124.55	128.60
1	AA	620	U	N3-C2-O2	-6.75	117.48	122.20
1	AA	2697	G	N7-C8-N9	-6.75	109.73	113.10
1	AA	2466	G	C6-N1-C2	-6.75	121.05	125.10
1	AA	1200	G	N1-C2-N2	-6.75	110.13	116.20
1	AA	1864	U	C4-C5-C6	6.75	123.75	119.70
1	AA	738	C	OP1-P-OP2	-6.74	109.48	119.60
1	AA	1232	G	N1-C6-O6	-6.74	115.85	119.90
1	AA	1715	A	O5'-P-OP2	-6.74	99.63	105.70
1	AA	566	C	C6-N1-C2	-6.74	117.60	120.30
1	AA	1563	G	C5-C6-O6	-6.74	124.56	128.60
34	BA	913	A	P-O3'-C3'	6.74	127.79	119.70
1	CA	120	U	O5'-P-OP1	-6.74	99.63	105.70
1	CA	1142	U	C5-C6-N1	6.74	126.07	122.70
1	AA	1373	C	N3-C4-C5	-6.74	119.20	121.90
1	AA	2329	C	O5'-P-OP1	-6.74	99.64	105.70
1	AA	2470	G	C5-C6-O6	-6.74	124.56	128.60
1	AA	2898	C	C2-N3-C4	-6.74	116.53	119.90
56	BW	17	C	C6-N1-C1'	-6.74	112.71	120.80
1	CA	1812	A	OP1-P-OP2	6.74	129.71	119.60
1	AA	24	G	OP2-P-O3'	6.74	120.02	105.20
1	AA	525	G	C8-N9-C4	6.74	109.09	106.40
1	AA	1000	C	N1-C2-O2	-6.74	114.86	118.90
1	AA	1885	A	C6-N1-C2	-6.74	114.56	118.60
1	AA	1247	C	C5-C6-N1	-6.73	117.63	121.00
1	AA	1749	G	N9-C4-C5	6.73	108.09	105.40
56	BW	36	A	C2-N3-C4	-6.73	107.23	110.60
1	AA	881	C	C2-N3-C4	-6.73	116.53	119.90
1	CA	843	G	C8-N9-C4	6.73	109.09	106.40
34	DA	819	A	C2-N3-C4	-6.73	107.23	110.60
1	AA	339	G	O5'-P-OP2	-6.73	99.64	105.70
1	AA	990	A	C8-N9-C4	-6.73	103.11	105.80
1	AA	1097	G	N9-C4-C5	-6.73	102.71	105.40
1	AA	1843	A	O5'-P-OP1	-6.73	99.64	105.70
1	AA	2040	G	N3-C2-N2	6.73	124.61	119.90
1	AA	2572	C	C6-N1-C2	-6.73	117.61	120.30
1	CA	2520	C	C6-N1-C2	6.73	122.99	120.30
1	AA	1046	A	N1-C6-N6	6.73	122.64	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1381	U	N1-C2-O2	-6.73	118.09	122.80
34	BA	899	C	N1-C2-O2	-6.73	114.86	118.90
1	AA	2455	C	C5-C6-N1	-6.73	117.64	121.00
1	AA	239	G	C5-C6-O6	-6.72	124.56	128.60
1	AA	845	G	C5-N7-C8	6.72	107.66	104.30
1	AA	1339	C	N3-C4-C5	6.72	124.59	121.90
34	BA	326	G	C5-C6-O6	-6.72	124.57	128.60
1	AA	2762	A	C6-N1-C2	-6.72	114.57	118.60
1	CA	2360	A	C6-N1-C2	-6.72	114.57	118.60
1	CA	2409	G	C5-C6-O6	-6.72	124.57	128.60
1	AA	517	A	C5-N7-C8	6.72	107.26	103.90
1	AA	564	G	OP2-P-O3'	6.72	119.98	105.20
1	AA	718	C	OP1-P-OP2	6.72	129.68	119.60
1	AA	1784	G	C4-C5-N7	6.72	113.49	110.80
34	BA	880	C	N3-C4-C5	6.72	124.59	121.90
1	CA	109	G	N1-C6-O6	-6.72	115.87	119.90
1	CA	2390	U	O5'-P-OP1	-6.72	99.65	105.70
1	AA	1874	C	N3-C4-C5	6.72	124.59	121.90
1	AA	363	U	N3-C4-C5	6.72	118.63	114.60
1	AA	2757	G	N3-C2-N2	-6.72	115.20	119.90
1	CA	2361	A	N1-C6-N6	6.72	122.63	118.60
1	AA	285	U	O4'-C1'-N1	6.71	113.57	108.20
1	AA	1857	G	C8-N9-C4	-6.71	103.71	106.40
1	AA	2546	A	C4-C5-N7	6.71	114.06	110.70
34	BA	1527	C	C2-N3-C4	-6.71	116.54	119.90
2	CB	70	C	C6-N1-C2	-6.71	117.61	120.30
1	AA	562	C	C2-N3-C4	-6.71	116.54	119.90
1	AA	1378	G	C6-N1-C2	-6.71	121.07	125.10
1	AA	1749	G	C5-C6-O6	6.71	132.63	128.60
1	AA	2513	C	C5-C6-N1	-6.71	117.64	121.00
1	AA	551	A	OP1-P-O3'	-6.71	90.44	105.20
1	AA	1343	C	C5-C6-N1	-6.71	117.64	121.00
1	CA	34	C	N1-C2-O2	6.71	122.93	118.90
1	AA	1453	C	C5-C4-N4	-6.71	115.50	120.20
1	AA	1837	C	C4-C5-C6	6.71	120.75	117.40
1	AA	2301	G	C5-C6-O6	-6.71	124.57	128.60
1	AA	2383	G	C2-N3-C4	6.71	115.25	111.90
34	BA	1442	G	C2-N3-C4	-6.71	108.55	111.90
1	AA	2049	G	OP1-P-OP2	6.71	129.66	119.60
1	AA	2113	U	C5-C6-N1	-6.71	119.35	122.70
1	AA	2399	U	N1-C2-O2	-6.71	118.10	122.80
1	CA	261	G	C8-N9-C4	6.71	109.08	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	135	C	N3-C4-C5	6.71	124.58	121.90
1	AA	625	G	C6-N1-C2	6.71	129.12	125.10
1	AA	1924	C	OP2-P-O3'	6.71	119.95	105.20
1	CA	2608	G	N7-C8-N9	-6.71	109.75	113.10
34	BA	894	G	N9-C4-C5	6.71	108.08	105.40
34	DA	822	C	O5'-P-OP1	-6.71	99.67	105.70
1	AA	1627	A	N7-C8-N9	6.70	117.15	113.80
1	AA	2608	U	C5-C6-N1	-6.70	119.35	122.70
2	AB	56	G	N3-C4-C5	-6.70	125.25	128.60
1	CA	2576	G	C5-C6-N1	6.70	114.85	111.50
1	AA	1092	A	N1-C6-N6	-6.70	114.58	118.60
1	AA	1542	A	C8-N9-C4	-6.70	103.12	105.80
1	AA	2260	C	C6-N1-C2	6.70	122.98	120.30
1	CA	1950	G	OP1-P-OP2	6.70	129.65	119.60
1	CA	2569	G	C5-C6-O6	-6.70	124.58	128.60
1	CA	133	C	N3-C4-C5	6.70	124.58	121.90
1	AA	1042	A	O5'-P-OP1	-6.70	99.67	105.70
1	AA	1074	A	C6-N1-C2	-6.70	114.58	118.60
1	AA	1341	C	C2-N3-C4	-6.70	116.55	119.90
1	AA	1474	C	C5-C6-N1	-6.69	117.65	121.00
1	AA	2221	A	OP2-P-O3'	6.69	119.92	105.20
1	AA	666	C	N3-C4-C5	-6.69	119.22	121.90
1	AA	905	U	N3-C4-O4	-6.69	114.72	119.40
1	AA	2488	A	C4-C5-N7	6.69	114.05	110.70
34	BA	918	A	N1-C6-N6	-6.69	114.59	118.60
1	CA	456	C	C6-N1-C2	6.69	122.98	120.30
1	CA	1914	C	C6-N1-C2	-6.69	117.62	120.30
1	AA	1648	U	O5'-P-OP2	6.69	118.73	110.70
1	AA	139	A	OP1-P-OP2	-6.69	109.57	119.60
1	AA	495	G	O5'-P-OP1	6.69	118.73	110.70
1	AA	1030	A	C8-N9-C4	6.69	108.47	105.80
1	AA	2391	G	N3-C2-N2	-6.69	115.22	119.90
1	CA	2586	C	O5'-P-OP2	-6.69	99.68	105.70
1	AA	2269	U	N3-C4-O4	-6.69	114.72	119.40
1	AA	987	G	C5-N7-C8	6.68	107.64	104.30
1	AA	1341	C	N3-C2-O2	-6.68	117.22	121.90
1	AA	2304	C	OP2-P-O3'	6.68	119.91	105.20
1	AA	2344	U	C5-C6-N1	-6.68	119.36	122.70
1	AA	2515	A	C6-N1-C2	6.68	122.61	118.60
1	CA	2510	C	C6-N1-C2	-6.68	117.63	120.30
1	CA	962	G	C5-C6-O6	-6.68	124.59	128.60
1	AA	1275	G	O5'-P-OP1	6.68	118.72	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1486	G	N7-C8-N9	-6.68	109.76	113.10
1	AA	725	C	N3-C2-O2	-6.68	117.23	121.90
1	AA	822	G	N1-C6-O6	-6.68	115.89	119.90
1	AA	2532	C	C2-N3-C4	-6.68	116.56	119.90
34	DA	286	G	C8-N9-C4	6.68	109.07	106.40
1	AA	1802	C	N1-C2-N3	6.67	123.87	119.20
1	AA	2375	C	N3-C4-N4	-6.67	113.33	118.00
1	AA	2528	G	C4-C5-N7	-6.67	108.13	110.80
34	BA	574	A	N1-C6-N6	6.67	122.61	118.60
1	CA	37	C	C5-C4-N4	-6.67	115.53	120.20
1	CA	150	C	N3-C2-O2	-6.67	117.23	121.90
1	AA	976	G	N3-C4-C5	-6.67	125.26	128.60
34	BA	731	G	C8-N9-C4	-6.67	103.73	106.40
2	AB	37	C	OP2-P-O3'	6.67	119.88	105.20
2	AB	61	G	C5-C6-N1	-6.67	108.16	111.50
1	AA	2522	C	N3-C2-O2	-6.67	117.23	121.90
1	AA	2399	U	C5-C6-N1	-6.67	119.37	122.70
1	AA	2761	A	OP2-P-O3'	6.67	119.87	105.20
1	CA	1636	C	C6-N1-C2	6.67	122.97	120.30
1	CA	2629	A	O4'-C1'-N9	6.67	113.53	108.20
34	DA	813	U	C6-N1-C2	6.67	125.00	121.00
1	AA	1766	G	C5-N7-C8	-6.67	100.97	104.30
1	CA	1565	C	C5-C6-N1	-6.67	117.67	121.00
1	AA	103	C	OP2-P-O3'	6.67	119.86	105.20
1	AA	75	C	O5'-P-OP2	-6.66	99.70	105.70
1	AA	867	A	N7-C8-N9	-6.66	110.47	113.80
1	AA	1614	A	O5'-P-OP1	6.66	118.70	110.70
1	AA	2065	C	C6-N1-C2	-6.66	117.63	120.30
1	CA	1332	G	O5'-P-OP2	-6.66	99.70	105.70
1	AA	214	A	O5'-P-OP2	-6.66	99.71	105.70
1	AA	787	U	N1-C2-N3	6.66	118.90	114.90
1	AA	913	A	N7-C8-N9	6.66	117.13	113.80
1	AA	2494	G	N1-C6-O6	6.66	123.90	119.90
1	AA	2609	G	N9-C4-C5	-6.66	102.74	105.40
1	CA	1558	A	C5-N7-C8	-6.66	100.57	103.90
34	BA	501	C	C6-N1-C2	-6.66	117.64	120.30
1	AA	2741	U	N1-C2-N3	6.66	118.89	114.90
1	CA	220	G	C5-C6-O6	-6.66	124.61	128.60
1	AA	174	U	C4-C5-C6	6.66	123.69	119.70
1	AA	415	G	O5'-P-OP1	6.66	118.69	110.70
1	AA	2335	G	C5-C6-O6	-6.66	124.61	128.60
1	CA	384	U	C5-C6-N1	-6.66	119.37	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	827	U	C5-C6-N1	-6.66	119.37	122.70
1	AA	514	G	C5-C6-O6	6.65	132.59	128.60
1	AA	549	U	C2-N3-C4	-6.65	123.01	127.00
1	AA	663	G	N3-C4-N9	6.65	129.99	126.00
1	AA	2450	U	N3-C2-O2	6.65	126.86	122.20
1	AA	2588	G	OP1-P-OP2	-6.65	109.62	119.60
1	AA	2794	A	N7-C8-N9	-6.65	110.47	113.80
1	AA	702	A	C2-N3-C4	-6.65	107.28	110.60
1	AA	2639	G	N3-C4-C5	6.65	131.93	128.60
1	AA	2647	C	C2-N3-C4	-6.65	116.58	119.90
1	CA	383	U	C4-C5-C6	6.65	123.69	119.70
1	CA	109	G	O5'-P-OP2	-6.65	99.72	105.70
1	CA	1792	G	O5'-P-OP2	-6.65	99.72	105.70
1	AA	1334	U	O5'-P-OP2	-6.65	99.72	105.70
2	AB	12	C	C2-N3-C4	-6.65	116.58	119.90
1	AA	1038	C	N3-C4-N4	-6.65	113.35	118.00
1	AA	2012	C	C5-C4-N4	-6.65	115.55	120.20
1	AA	540	A	N1-C6-N6	-6.64	114.61	118.60
1	AA	1811	A	O5'-P-OP1	6.64	118.67	110.70
1	AA	2442	A	C6-N1-C2	6.64	122.59	118.60
1	AA	2741	U	C6-N1-C2	-6.64	117.01	121.00
1	CA	741	G	C5-C6-O6	6.64	132.59	128.60
1	AA	436	C	C2-N3-C4	6.64	123.22	119.90
1	AA	719	C	N3-C4-C5	6.64	124.56	121.90
1	AA	2028	C	C5-C4-N4	-6.64	115.55	120.20
1	AA	2290	A	O5'-P-OP1	6.64	118.67	110.70
1	AA	2380	C	C6-N1-C2	6.64	122.96	120.30
1	AA	2550	C	C5-C4-N4	-6.64	115.55	120.20
1	CA	141	A	C2-N3-C4	-6.64	107.28	110.60
1	CA	2195	C	OP1-P-O3'	6.64	119.81	105.20
1	AA	726	C	N3-C4-N4	6.64	122.64	118.00
1	AA	1100	A	N1-C6-N6	6.64	122.58	118.60
1	AA	1922	A	C8-N9-C4	-6.64	103.14	105.80
1	AA	2530	A	C2-N3-C4	-6.63	107.28	110.60
1	AA	1050	C	C2-N3-C4	-6.63	116.58	119.90
1	AA	1449	C	O5'-P-OP1	-6.63	99.73	105.70
34	BA	345	C	C2-N1-C1'	6.63	126.09	118.80
1	CA	1333	C	C5-C4-N4	-6.63	115.56	120.20
1	AA	1331	G	N3-C4-N9	6.63	129.98	126.00
1	AA	1857	G	N9-C4-C5	6.63	108.05	105.40
1	AA	593	G	C5-N7-C8	-6.63	100.98	104.30
1	CA	798	G	C5-C6-O6	6.63	132.58	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	97	G	N9-C4-C5	6.63	108.05	105.40
1	AA	132	C	O5'-P-OP1	6.63	118.65	110.70
1	AA	1252	C	C2-N3-C4	-6.63	116.59	119.90
1	AA	1665	G	N3-C4-N9	6.63	129.98	126.00
1	AA	2295	C	N3-C4-N4	6.63	122.64	118.00
1	AA	2785	C	N3-C2-O2	-6.63	117.26	121.90
1	AA	423	G	O5'-P-OP1	-6.62	99.74	105.70
1	AA	2366	G	C8-N9-C1'	-6.62	118.39	127.00
34	BA	567	G	C4-C5-N7	-6.62	108.15	110.80
1	AA	630	U	C6-N1-C2	6.62	124.97	121.00
1	AA	1204	C	C6-N1-C2	-6.62	117.65	120.30
1	CA	572	A	O5'-P-OP2	-6.62	99.74	105.70
1	CA	961	C	C6-N1-C2	6.62	122.95	120.30
1	CA	1829	A	N1-C6-N6	6.62	122.57	118.60
1	CA	2043	C	N3-C4-C5	6.62	124.55	121.90
1	AA	712	C	N3-C4-N4	-6.62	113.37	118.00
1	AA	2659	U	N3-C4-C5	6.62	118.57	114.60
1	AA	114	C	C2-N3-C4	-6.62	116.59	119.90
1	AA	1725	G	C6-N1-C2	-6.62	121.13	125.10
1	AA	2220	A	O4'-C1'-N9	6.62	113.50	108.20
1	AA	2702	C	N3-C2-O2	6.62	126.53	121.90
34	BA	373	A	OP1-P-O3'	6.62	119.76	105.20
1	AA	46	C	C5-C6-N1	-6.62	117.69	121.00
1	AA	2387	G	C5-C6-N1	6.62	114.81	111.50
1	CA	1997	G	O5'-P-OP2	-6.62	99.75	105.70
34	DA	442	C	C6-N1-C2	-6.62	117.65	120.30
1	AA	1021	G	C6-N1-C2	6.61	129.07	125.10
1	AA	2414	C	O5'-P-OP2	-6.61	99.75	105.70
1	AA	2724	U	OP1-P-OP2	6.61	129.52	119.60
1	CA	2618	G	C6-N1-C2	-6.61	121.13	125.10
1	CA	769	G	C6-C5-N7	-6.61	126.44	130.40
1	CA	2673	G	C5-C6-O6	6.61	132.57	128.60
1	AA	1040	C	C2-N3-C4	-6.61	116.60	119.90
1	AA	1421	C	O5'-P-OP2	6.61	118.63	110.70
1	AA	581	G	N7-C8-N9	-6.61	109.80	113.10
1	CA	563	G	N3-C4-N9	-6.61	122.04	126.00
15	CR	64	ARG	NE-CZ-NH1	6.61	123.60	120.30
1	AA	418	G	N1-C6-O6	6.60	123.86	119.90
1	AA	580	U	OP2-P-O3'	6.60	119.73	105.20
1	AA	859	C	C2-N3-C4	-6.60	116.60	119.90
1	AA	1474	C	N1-C2-N3	6.60	123.82	119.20
1	CA	2490	G	C5-N7-C8	-6.60	101.00	104.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	549	U	N3-C4-C5	6.60	118.56	114.60
1	AA	2379	G	N3-C4-C5	-6.60	125.30	128.60
1	CA	1901	A	C5-C6-N1	6.60	121.00	117.70
1	CA	2500	U	C5-C6-N1	-6.60	119.40	122.70
1	AA	1477	U	C5-C6-N1	6.60	126.00	122.70
1	AA	1741	C	C6-N1-C2	6.60	122.94	120.30
34	DA	640	A	C8-N9-C4	-6.60	103.16	105.80
1	AA	29	U	C2-N3-C4	-6.59	123.04	127.00
1	AA	844	C	N3-C4-N4	6.59	122.62	118.00
1	AA	464	G	C6-C5-N7	-6.59	126.44	130.40
1	AA	2876	U	N3-C4-O4	-6.59	114.78	119.40
1	AA	555	G	N3-C2-N2	-6.59	115.29	119.90
1	AA	2307	C	O5'-P-OP2	6.59	118.61	110.70
1	AA	2838	C	C6-N1-C2	6.59	122.94	120.30
1	CA	1653	G	OP1-P-OP2	6.59	129.49	119.60
1	AA	839	G	C5-C6-O6	-6.59	124.65	128.60
1	AA	998	A	C2-N3-C4	-6.59	107.31	110.60
1	AA	1847	G	N1-C2-N2	6.59	122.13	116.20
34	BA	1524	C	N1-C2-N3	6.59	123.81	119.20
1	CA	2029	G	C5-C6-O6	-6.59	124.65	128.60
1	AA	1193	C	C2-N1-C1'	-6.59	111.55	118.80
1	AA	2591	C	N3-C4-N4	6.59	122.61	118.00
1	AA	1411	A	C6-C5-N7	-6.59	127.69	132.30
1	AA	2662	U	C5-C4-O4	6.59	129.85	125.90
1	AA	2826	C	N1-C2-O2	6.59	122.85	118.90
30	A6	19	ARG	NE-CZ-NH1	-6.59	117.01	120.30
1	AA	1724	A	C6-N1-C2	-6.58	114.65	118.60
1	CA	1300	U	P-O3'-C3'	6.58	127.60	119.70
1	AA	627	G	N3-C4-C5	-6.58	125.31	128.60
1	AA	1330	A	OP1-P-OP2	6.58	129.47	119.60
1	AA	2070	G	N7-C8-N9	-6.58	109.81	113.10
1	AA	2445	A	N1-C6-N6	6.58	122.55	118.60
1	CA	2827	C	N3-C2-O2	6.58	126.51	121.90
34	DA	1484	C	C6-N1-C2	6.58	122.93	120.30
1	AA	1053	C	N1-C2-O2	-6.58	114.95	118.90
1	AA	1356	G	O5'-P-OP1	-6.58	99.78	105.70
1	AA	2417	G	C4-C5-N7	6.58	113.43	110.80
1	AA	211	A	N9-C4-C5	6.58	108.43	105.80
1	AA	1021	G	OP1-P-OP2	6.58	129.46	119.60
34	BA	770	C	C5-C4-N4	6.58	124.80	120.20
1	CA	2238	G	O5'-P-OP1	-6.58	99.78	105.70
1	AA	993	G	O5'-P-OP1	-6.57	99.78	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2490	G	N9-C4-C5	-6.57	102.77	105.40
1	CA	2580	U	C6-N1-C2	6.57	124.94	121.00
1	AA	2386	C	C2-N3-C4	-6.57	116.61	119.90
1	AA	2440	G	N1-C2-N2	-6.57	110.28	116.20
34	BA	1412	C	C6-N1-C2	6.57	122.93	120.30
1	CA	1949	G	C5-C6-O6	-6.57	124.66	128.60
1	AA	901	G	C4-C5-N7	-6.57	108.17	110.80
1	AA	1757	C	O5'-P-OP1	6.57	118.58	110.70
1	AA	2045	G	C5-C6-N1	6.57	114.78	111.50
1	CA	600	G	N3-C4-C5	-6.57	125.32	128.60
1	AA	1034	A	N1-C2-N3	-6.57	126.02	129.30
1	CA	2832	U	C5-C6-N1	-6.57	119.42	122.70
1	AA	1063	G	C4-C5-N7	-6.56	108.17	110.80
1	AA	1457	C	N3-C4-C5	6.56	124.53	121.90
34	BA	363	A	O5'-P-OP2	-6.56	99.79	105.70
56	DW	17	C	C6-N1-C2	-6.56	117.67	120.30
1	AA	472	G	C5-C6-O6	-6.56	124.66	128.60
1	AA	1064	C	N1-C2-O2	6.56	122.84	118.90
1	AA	1345	G	O5'-P-OP2	6.56	118.58	110.70
1	CA	2503	A	N9-C4-C5	-6.56	103.17	105.80
1	CA	2763	G	C5-C6-O6	6.56	132.54	128.60
1	AA	70	A	C4-C5-C6	6.56	120.28	117.00
1	AA	107	G	C6-C5-N7	6.56	134.34	130.40
1	AA	555	G	C5-N7-C8	-6.56	101.02	104.30
1	AA	2471	A	N3-C4-C5	-6.56	122.21	126.80
1	CA	204	A	C6-N1-C2	-6.56	114.66	118.60
1	AA	1617	A	N9-C4-C5	-6.56	103.18	105.80
1	AA	1766	G	C4-C5-N7	6.56	113.42	110.80
34	BA	1530	G	C5-C6-O6	-6.56	124.67	128.60
1	CA	811	U	C2-N3-C4	-6.56	123.07	127.00
1	AA	327	U	C5-C4-O4	-6.56	121.97	125.90
1	AA	534	C	N1-C2-N3	-6.56	114.61	119.20
34	BA	115	G	P-O3'-C3'	6.56	127.57	119.70
1	CA	382	G	C2-N3-C4	6.56	115.18	111.90
1	AA	431	C	N3-C2-O2	-6.55	117.31	121.90
1	AA	861	C	O5'-P-OP1	6.55	118.56	110.70
1	AA	1040	C	N3-C4-C5	6.55	124.52	121.90
1	CA	2560	C	N1-C2-O2	6.55	122.83	118.90
1	CA	512	G	O5'-P-OP1	-6.55	99.80	105.70
1	AA	213	G	N9-C4-C5	6.55	108.02	105.40
1	AA	743	G	N9-C4-C5	-6.55	102.78	105.40
1	CA	1424	G	O5'-P-OP2	-6.55	99.80	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1682	G	C5-C6-N1	6.55	114.78	111.50
1	AA	1700	G	O4'-C1'-N9	-6.55	102.96	108.20
34	BA	964	A	C8-N9-C4	6.55	108.42	105.80
1	AA	611	U	O5'-P-OP2	-6.55	99.81	105.70
1	AA	1493	C	N3-C2-O2	-6.55	117.32	121.90
1	AA	1599	G	O5'-P-OP2	-6.55	99.81	105.70
1	AA	2822	G	C4-C5-N7	-6.55	108.18	110.80
2	AB	39	A	C8-N9-C4	6.55	108.42	105.80
34	BA	1383	C	C6-N1-C2	-6.55	117.68	120.30
34	DA	586	C	C5-C4-N4	-6.55	115.62	120.20
1	AA	211	A	C2-N3-C4	6.54	113.87	110.60
1	AA	1694	G	C6-C5-N7	6.54	134.33	130.40
1	CA	1296	G	O5'-P-OP1	-6.54	99.81	105.70
1	AA	100	G	C4-C5-N7	6.54	113.42	110.80
1	CA	264	C	C6-N1-C2	6.54	122.92	120.30
1	CA	1704	G	N1-C6-O6	6.54	123.83	119.90
1	AA	1439	A	N1-C6-N6	-6.54	114.67	118.60
1	AA	2857	U	C5-C6-N1	-6.54	119.43	122.70
1	CA	1745(A)	C	C6-N1-C2	-6.54	117.68	120.30
1	CA	2732	G	C5-C6-O6	6.54	132.53	128.60
1	AA	2011	G	C5-C6-N1	-6.54	108.23	111.50
1	AA	137	G	N3-C4-C5	-6.54	125.33	128.60
18	AU	3	ARG	NE-CZ-NH2	6.54	123.57	120.30
1	AA	1669	G	O5'-P-OP2	-6.54	99.82	105.70
1	AA	1978	U	N3-C4-C5	6.54	118.52	114.60
1	AA	597	C	N3-C4-N4	-6.54	113.43	118.00
1	AA	627	G	N9-C4-C5	6.54	108.01	105.40
1	AA	208	G	O5'-P-OP2	-6.53	99.82	105.70
1	AA	550	U	N3-C2-O2	-6.53	117.63	122.20
1	AA	816	G	N3-C2-N2	6.53	124.47	119.90
1	AA	1030	A	N7-C8-N9	-6.53	110.53	113.80
1	AA	1343	C	C2-N3-C4	-6.53	116.63	119.90
1	AA	243	G	C5-C6-N1	6.53	114.77	111.50
1	AA	436	C	O5'-P-OP1	6.53	118.54	110.70
1	AA	872	C	C6-N1-C2	6.53	122.91	120.30
1	AA	2468	C	C6-N1-C2	6.53	122.91	120.30
2	CB	85	G	O5'-P-OP2	-6.53	99.82	105.70
1	AA	112	U	C5-C4-O4	-6.53	121.98	125.90
1	AA	1024	G	N9-C4-C5	6.53	108.01	105.40
1	AA	2660	C	N1-C2-O2	-6.53	114.98	118.90
1	CA	195	A	OP2-P-O3'	6.53	119.56	105.20
1	CA	1902	C	C5-C6-N1	6.53	124.26	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2713	A	C5-C6-N6	6.53	128.92	123.70
34	DA	1513	A	C8-N9-C4	6.53	108.41	105.80
1	AA	426	G	N1-C2-N3	6.53	127.81	123.90
1	AA	1232	G	O5'-P-OP1	6.53	118.53	110.70
1	AA	1797	U	OP1-P-OP2	-6.53	109.81	119.60
1	AA	2367	C	C5-C6-N1	-6.53	117.74	121.00
1	AA	2454	C	N1-C2-O2	-6.53	114.98	118.90
1	AA	1008	U	C5-C6-N1	-6.52	119.44	122.70
1	AA	1024	G	N1-C6-O6	-6.52	115.99	119.90
1	AA	1210	G	O5'-P-OP2	-6.52	99.83	105.70
1	CA	139(A)	G	N3-C4-N9	6.52	129.91	126.00
1	CA	933	A	C5-N7-C8	-6.52	100.64	103.90
2	AB	62	C	O5'-P-OP2	-6.52	99.83	105.70
1	CA	914	C	N1-C2-O2	6.52	122.81	118.90
1	AA	715	G	OP2-P-O3'	6.52	119.54	105.20
1	AA	1718	U	C5-C4-O4	-6.52	121.99	125.90
1	AA	2092	G	OP1-P-OP2	6.52	129.38	119.60
34	DA	1505	G	N3-C4-C5	6.52	131.86	128.60
1	AA	1307	C	N3-C4-N4	-6.52	113.44	118.00
1	AA	2422	G	C8-N9-C4	6.52	109.01	106.40
1	AA	1390	G	N7-C8-N9	-6.51	109.84	113.10
1	AA	1443	U	C5-C6-N1	-6.51	119.44	122.70
1	AA	2612	A	C5-C6-N6	6.51	128.91	123.70
1	CA	1792	G	N9-C4-C5	-6.51	102.79	105.40
1	CA	2553	G	N3-C2-N2	6.51	124.46	119.90
1	AA	1735	U	N3-C4-O4	-6.51	114.84	119.40
1	AA	2272	C	O5'-P-OP1	-6.51	99.84	105.70
1	AA	455	A	N7-C8-N9	6.51	117.06	113.80
1	AA	1234	A	N1-C2-N3	6.51	132.56	129.30
1	AA	1700	G	N3-C4-C5	-6.51	125.34	128.60
1	AA	2513	C	OP2-P-O3'	6.51	119.52	105.20
1	AA	2733	U	C5-C6-N1	-6.51	119.44	122.70
1	CA	39	C	N3-C4-C5	6.51	124.50	121.90
34	DA	546	G	N1-C6-O6	-6.51	115.99	119.90
1	AA	704	U	C2-N3-C4	-6.51	123.09	127.00
1	CA	2222	G	N1-C6-O6	-6.51	116.00	119.90
1	CA	2501	C	C6-N1-C2	6.51	122.90	120.30
1	AA	1097	G	N3-C4-C5	6.51	131.85	128.60
34	BA	1405	G	N1-C6-O6	-6.51	116.00	119.90
1	AA	122	G	OP1-P-OP2	6.50	129.36	119.60
1	AA	907	U	C6-N1-C2	6.50	124.90	121.00
1	AA	16	G	C8-N9-C4	-6.50	103.80	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	420	C	N1-C2-O2	6.50	122.80	118.90
1	AA	2293	C	N1-C2-O2	-6.50	115.00	118.90
1	AA	2716	C	C5-C6-N1	6.50	124.25	121.00
1	AA	438	G	N3-C4-N9	6.50	129.90	126.00
1	AA	775	G	N1-C2-N2	-6.50	110.35	116.20
1	AA	912	C	N3-C4-N4	6.50	122.55	118.00
1	AA	1846	A	O5'-P-OP1	-6.50	99.85	105.70
1	AA	341	G	N1-C6-O6	-6.50	116.00	119.90
1	AA	592	U	C4-C5-C6	6.50	123.60	119.70
1	AA	2244	U	C5-C6-N1	-6.50	119.45	122.70
1	AA	2327	G	C8-N9-C4	6.50	109.00	106.40
1	AA	2107	C	C5-C4-N4	-6.50	115.65	120.20
1	AA	2294	G	C2-N3-C4	6.50	115.15	111.90
1	AA	2787	C	C4-C5-C6	6.50	120.65	117.40
1	AA	832	G	C6-C5-N7	6.49	134.30	130.40
1	AA	1573	G	N1-C6-O6	6.49	123.80	119.90
1	AA	1795	G	N3-C2-N2	6.49	124.45	119.90
1	AA	2000	A	N1-C2-N3	-6.49	126.05	129.30
1	AA	2466	G	N3-C4-C5	-6.49	125.35	128.60
1	CA	1154	G	N3-C4-N9	6.49	129.89	126.00
1	AA	354	A	C4-C5-C6	-6.49	113.76	117.00
1	AA	1474	C	C4-C5-C6	6.49	120.64	117.40
1	AA	1411	A	C5-C6-N6	-6.49	118.51	123.70
1	AA	1819	C	O5'-P-OP1	-6.49	99.86	105.70
1	CA	1021	A	N1-C2-N3	6.49	132.54	129.30
1	AA	827	G	C5-C6-O6	-6.48	124.71	128.60
1	AA	1783	C	N1-C2-O2	-6.48	115.01	118.90
1	AA	26	G	C4-C5-N7	6.48	113.39	110.80
1	AA	2604	G	O5'-P-OP2	6.48	118.48	110.70
1	AA	1068	G	N3-C4-N9	-6.48	122.11	126.00
1	AA	2294	G	O5'-P-OP2	6.48	118.48	110.70
1	AA	2712	C	O5'-P-OP2	-6.48	99.87	105.70
1	CA	2058	A	N1-C2-N3	6.48	132.54	129.30
1	AA	789	G	C2-N3-C4	-6.48	108.66	111.90
1	CA	741	G	C2-N3-C4	-6.48	108.66	111.90
1	AA	1805	C	N1-C2-O2	-6.48	115.01	118.90
1	AA	2000	A	N7-C8-N9	-6.48	110.56	113.80
1	AA	2400	A	C4-C5-N7	-6.48	107.46	110.70
1	AA	2478	C	C4-C5-C6	6.48	120.64	117.40
1	CA	801	G	N3-C4-C5	-6.48	125.36	128.60
1	CA	954	G	C8-N9-C4	-6.48	103.81	106.40
1	AA	1243	U	N1-C2-N3	6.47	118.78	114.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2374	G	N1-C2-N3	6.47	127.78	123.90
34	BA	854	G	C8-N9-C4	-6.47	103.81	106.40
1	AA	1194	A	N1-C6-N6	-6.47	114.72	118.60
1	AA	1307	C	C6-N1-C2	6.47	122.89	120.30
1	AA	2052	A	OP1-P-OP2	6.47	129.31	119.60
1	AA	2367	C	C6-N1-C2	6.47	122.89	120.30
1	CA	2249	U	N1-C2-O2	6.47	127.33	122.80
1	AA	1013	G	C5-C6-O6	6.47	132.48	128.60
1	AA	1700	G	P-O3'-C3'	6.47	127.47	119.70
1	AA	2773	C	OP2-P-O3'	6.47	119.44	105.20
1	AA	2792	U	N1-C2-N3	6.47	118.78	114.90
34	DA	1509	C	N3-C4-C5	6.47	124.49	121.90
1	AA	2024	G	OP2-P-O3'	-6.47	90.97	105.20
1	AA	1200	G	C6-C5-N7	-6.46	126.52	130.40
1	AA	1411	A	N1-C2-N3	6.46	132.53	129.30
1	AA	1481	G	O5'-P-OP1	6.46	118.46	110.70
1	AA	2227	G	N3-C4-C5	6.46	131.83	128.60
1	CA	383	U	N1-C2-N3	6.46	118.78	114.90
1	AA	422	U	O5'-P-OP1	-6.46	99.88	105.70
56	BW	45	U	N1-C2-O2	6.46	127.32	122.80
1	AA	420	C	N3-C4-C5	-6.46	119.31	121.90
1	AA	531	G	C2-N3-C4	6.46	115.13	111.90
1	AA	1948	U	N3-C2-O2	-6.46	117.68	122.20
1	AA	2677	A	N1-C6-N6	6.46	122.48	118.60
34	BA	1510	U	C6-N1-C2	6.46	124.88	121.00
1	CA	579	G	C5-C6-O6	-6.46	124.72	128.60
34	DA	498	U	O5'-P-OP2	-6.46	99.88	105.70
34	BA	1397	C	C5-C6-N1	6.46	124.23	121.00
34	DA	634	C	C6-N1-C2	-6.46	117.72	120.30
1	AA	989	G	C6-C5-N7	-6.46	126.53	130.40
1	CA	2083	G	C6-C5-N7	-6.46	126.53	130.40
1	CA	2275	C	OP1-P-O3'	6.46	119.41	105.20
1	AA	1187	U	C5-C4-O4	-6.46	122.03	125.90
34	BA	841	U	C5-C6-N1	6.46	125.93	122.70
1	AA	2640	C	N3-C4-C5	6.46	124.48	121.90
34	BA	1437	C	O5'-P-OP1	-6.46	99.89	105.70
34	BA	885	G	C5-C6-O6	-6.45	124.73	128.60
1	AA	598	A	C5-N7-C8	6.45	107.13	103.90
34	BA	1263	C	C6-N1-C2	6.45	122.88	120.30
1	CA	214	G	C8-N9-C4	6.45	108.98	106.40
1	CA	474	G	OP2-P-O3'	6.45	119.39	105.20
1	CA	2257	U	N3-C4-C5	6.45	118.47	114.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2610	A	OP1-P-OP2	-6.45	109.93	119.60
1	CA	2413	G	C8-N9-C4	6.45	108.98	106.40
1	AA	1443	U	O5'-P-OP2	6.45	118.44	110.70
1	AA	1645	C	C2-N3-C4	-6.45	116.68	119.90
1	AA	1876	G	C2-N3-C4	-6.45	108.68	111.90
34	DA	1502	A	C2-N3-C4	-6.45	107.38	110.60
1	AA	472	G	C4-C5-N7	6.45	113.38	110.80
1	AA	1985	U	OP1-P-O3'	6.45	119.38	105.20
1	CA	1142(A)	A	N3-C4-N9	-6.45	122.24	127.40
1	CA	845	G	O4'-C1'-N9	6.44	113.36	108.20
1	AA	1515	C	O5'-P-OP1	6.44	118.43	110.70
1	AA	139	A	O4'-C1'-N9	6.44	113.35	108.20
1	AA	2612	A	C6-C5-N7	6.44	136.81	132.30
1	AA	2740	G	C5-C6-O6	6.44	132.46	128.60
34	BA	1520	G	O5'-P-OP1	6.44	118.43	110.70
1	AA	791	G	C5-C6-O6	6.44	132.46	128.60
1	AA	2825	C	C6-N1-C2	-6.44	117.73	120.30
1	AA	831	A	C5-N7-C8	6.43	107.12	103.90
1	AA	2256	U	N1-C2-N3	6.43	118.76	114.90
1	AA	2538	G	O5'-P-OP1	6.43	118.42	110.70
1	AA	822	G	C8-N9-C4	-6.43	103.83	106.40
1	AA	839	G	C6-N1-C2	-6.43	121.24	125.10
1	AA	241	G	C6-C5-N7	-6.43	126.54	130.40
1	AA	821	A	OP2-P-O3'	6.43	119.34	105.20
1	AA	1031	C	C5-C6-N1	6.43	124.21	121.00
1	AA	2566	U	N1-C2-O2	-6.43	118.30	122.80
1	CA	1609	A	C8-N9-C4	6.43	108.37	105.80
1	CA	2628	C	N1-C2-O2	-6.43	115.04	118.90
1	AA	69	G	C4-C5-N7	6.43	113.37	110.80
1	AA	1410	G	N3-C4-N9	6.43	129.86	126.00
1	AA	477	C	C5-C6-N1	-6.43	117.79	121.00
1	AA	623	G	OP1-P-OP2	6.43	129.24	119.60
1	CA	2442	C	N1-C2-O2	6.43	122.76	118.90
1	AA	364	A	O5'-P-OP2	6.42	118.41	110.70
1	AA	472	G	C5-N7-C8	-6.42	101.09	104.30
1	AA	1960	A	C2-N3-C4	-6.42	107.39	110.60
1	AA	2528	G	N1-C6-O6	-6.42	116.05	119.90
1	AA	2894	U	N3-C2-O2	6.42	126.70	122.20
34	BA	1497	G	C8-N9-C4	6.42	108.97	106.40
1	AA	1331	G	C4-C5-N7	6.42	113.37	110.80
1	AA	1609	A	C2-N3-C4	-6.42	107.39	110.60
1	AA	1861	C	N3-C4-C5	6.42	124.47	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2092	G	O5'-P-OP2	-6.42	99.92	105.70
1	AA	2386	C	C5-C6-N1	-6.42	117.79	121.00
1	AA	2782	C	C2-N3-C4	-6.42	116.69	119.90
32	A8	13	ARG	NE-CZ-NH1	6.42	123.51	120.30
1	CA	1607	C	C6-N1-C2	-6.42	117.73	120.30
1	CA	2588	G	N3-C4-C5	6.42	131.81	128.60
1	AA	1745	A	C4-N9-C1'	6.42	137.86	126.30
1	AA	1749	G	N1-C6-O6	-6.42	116.05	119.90
1	AA	195	U	C5-C4-O4	6.42	129.75	125.90
1	AA	1998	U	OP1-P-O3'	6.42	119.32	105.20
1	CA	141	A	C4-C5-N7	6.42	113.91	110.70
1	CA	1635	G	OP1-P-O3'	6.42	119.32	105.20
34	DA	397	A	O5'-P-OP2	-6.42	99.92	105.70
1	AA	36	G	C5-C6-O6	6.42	132.45	128.60
1	AA	1276	C	O5'-P-OP1	6.42	118.40	110.70
1	AA	1473	A	O5'-P-OP2	-6.42	99.93	105.70
19	AV	13	ARG	NE-CZ-NH1	-6.42	117.09	120.30
1	AA	385	G	N1-C6-O6	6.42	123.75	119.90
1	AA	908	A	N1-C6-N6	-6.42	114.75	118.60
1	AA	733	G	N1-C2-N2	-6.41	110.43	116.20
1	AA	1064	C	C6-N1-C2	-6.41	117.73	120.30
1	AA	2012	C	C6-N1-C2	-6.41	117.73	120.30
1	CA	2574	G	C5-C6-O6	-6.41	124.75	128.60
1	AA	1983	C	N3-C4-C5	-6.41	119.33	121.90
1	CA	144	C	N3-C4-C5	-6.41	119.33	121.90
1	CA	60	G	C4-C5-N7	6.41	113.36	110.80
1	CA	1655	A	OP1-P-OP2	-6.41	109.99	119.60
1	CA	2610	C	N3-C2-O2	-6.41	117.41	121.90
1	CA	2767	C	C6-N1-C2	-6.41	117.74	120.30
1	AA	743	G	O5'-P-OP1	-6.41	99.93	105.70
1	AA	243	G	N1-C2-N2	-6.41	110.44	116.20
1	AA	976	G	N1-C2-N3	6.41	127.74	123.90
1	AA	1978	U	O5'-P-OP2	-6.41	99.94	105.70
1	AA	2485	U	O5'-P-OP1	-6.41	99.93	105.70
1	AA	2755	C	N3-C2-O2	-6.40	117.42	121.90
1	CA	1683	C	N3-C4-C5	-6.40	119.34	121.90
1	CA	2023	G	N3-C2-N2	-6.40	115.42	119.90
1	AA	1365	G	N7-C8-N9	6.40	116.30	113.10
1	AA	1756	U	C5-C6-N1	-6.40	119.50	122.70
1	AA	2075	G	C4-C5-N7	-6.40	108.24	110.80
1	AA	2504	U	C6-N1-C2	-6.40	117.16	121.00
1	CA	1321	A	N1-C6-N6	6.40	122.44	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	82	G	N9-C4-C5	-6.40	102.84	105.40
1	AA	1270	C	C6-N1-C2	6.40	122.86	120.30
1	AA	857	U	O5'-P-OP1	6.40	118.38	110.70
1	AA	1066	A	N1-C6-N6	6.40	122.44	118.60
1	AA	1683	C	O5'-P-OP2	6.40	118.38	110.70
1	CA	37	C	C2-N3-C4	-6.40	116.70	119.90
1	CA	716	A	N9-C4-C5	6.40	108.36	105.80
1	AA	1262	C	C2-N3-C4	-6.40	116.70	119.90
1	CA	252	G	C4-C5-N7	-6.40	108.24	110.80
1	CA	1968	G	N3-C2-N2	-6.40	115.42	119.90
1	AA	181	C	N1-C2-N3	6.39	123.68	119.20
1	AA	674	G	C2-N3-C4	6.39	115.10	111.90
1	AA	2440	G	N1-C6-O6	-6.39	116.06	119.90
1	AA	2530	A	C4-C5-C6	6.39	120.20	117.00
1	CA	1770	G	O5'-P-OP2	6.39	118.37	110.70
1	CA	1774	C	O5'-P-OP1	6.39	118.37	110.70
1	CA	2708	G	N3-C4-N9	6.39	129.84	126.00
1	AA	188	A	N9-C4-C5	6.39	108.36	105.80
1	AA	1830	G	C4-C5-N7	-6.39	108.24	110.80
1	AA	2285	A	C6-C5-N7	-6.39	127.83	132.30
2	AB	68	C	OP2-P-O3'	6.39	119.26	105.20
34	DA	1158	C	C6-N1-C2	-6.39	117.74	120.30
1	AA	1639	G	O5'-P-OP1	-6.39	99.95	105.70
1	AA	2020	G	N3-C2-N2	6.39	124.37	119.90
1	AA	2432	C	C2-N3-C4	-6.39	116.70	119.90
1	AA	354	A	C8-N9-C4	-6.39	103.24	105.80
1	AA	525	G	N7-C8-N9	-6.39	109.91	113.10
1	AA	767	C	C6-N1-C2	6.39	122.86	120.30
1	AA	2441	G	N3-C2-N2	-6.39	115.43	119.90
1	CA	772	C	C5-C6-N1	-6.39	117.81	121.00
1	AA	749	G	O5'-P-OP2	-6.39	99.95	105.70
1	AA	984	G	N3-C2-N2	6.39	124.37	119.90
1	AA	2014	G	N1-C6-O6	-6.39	116.07	119.90
1	AA	716	G	N9-C4-C5	-6.39	102.85	105.40
1	AA	1012	C	N3-C4-N4	-6.39	113.53	118.00
1	AA	1686	U	OP1-P-OP2	6.39	129.18	119.60
1	AA	1814	A	N9-C4-C5	6.39	108.35	105.80
1	CA	64	A	N1-C6-N6	6.39	122.43	118.60
1	CA	572	A	C5-C6-N6	-6.39	118.59	123.70
34	DA	1081	G	C8-N9-C4	6.39	108.95	106.40
1	AA	531	G	N1-C6-O6	-6.38	116.07	119.90
1	AA	785	G	C4-C5-N7	6.38	113.35	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2697	G	C4-C5-N7	-6.38	108.25	110.80
1	CA	1647	G	O5'-P-OP1	-6.38	99.95	105.70
34	DA	335	C	N3-C2-O2	-6.38	117.43	121.90
34	BA	552	U	O5'-P-OP2	-6.38	99.95	105.70
1	AA	208	G	N3-C4-N9	6.38	129.83	126.00
1	AA	893	C	N1-C2-N3	6.38	123.67	119.20
1	AA	2095	C	OP2-P-O3'	6.38	119.24	105.20
1	AA	2461	U	C5-C4-O4	-6.38	122.07	125.90
1	AA	2484	G	N7-C8-N9	-6.38	109.91	113.10
1	AA	2515	A	N1-C6-N6	6.38	122.43	118.60
1	CA	1798	U	N3-C4-O4	-6.38	114.93	119.40
1	AA	1281	G	N1-C2-N3	6.38	127.73	123.90
1	AA	1394	G	C5-C6-N1	6.38	114.69	111.50
1	AA	1441	A	C5-C6-N1	6.38	120.89	117.70
1	AA	2081	A	OP2-P-O3'	6.38	119.23	105.20
1	AA	2580	C	C5-C6-N1	-6.38	117.81	121.00
1	CA	1835	G	N3-C4-N9	6.38	129.83	126.00
1	CA	2245	U	C5-C4-O4	-6.38	122.07	125.90
34	DA	29	G	O5'-P-OP1	-6.38	99.96	105.70
1	AA	1416	C	C5-C4-N4	-6.38	115.74	120.20
1	AA	2536	G	OP2-P-O3'	6.38	119.23	105.20
1	CA	1565	C	C6-N1-C2	6.38	122.85	120.30
1	AA	905	U	O5'-P-OP2	-6.38	99.96	105.70
1	AA	2782	C	N3-C4-C5	6.38	124.45	121.90
1	AA	1663	C	N3-C4-C5	6.37	124.45	121.90
1	AA	2051	G	O5'-P-OP2	-6.37	99.97	105.70
34	BA	556	C	C6-N1-C2	-6.37	117.75	120.30
1	CA	1792	G	C8-N9-C1'	-6.37	118.72	127.00
1	CA	690	G	O5'-P-OP1	-6.37	99.97	105.70
1	AA	326	C	N3-C4-C5	-6.37	119.35	121.90
34	BA	1417	G	C6-N1-C2	-6.37	121.28	125.10
1	CA	2824	C	C2-N3-C4	-6.37	116.72	119.90
1	AA	592	U	C2-N1-C1'	-6.37	110.06	117.70
1	AA	998	A	N7-C8-N9	6.37	116.98	113.80
1	AA	1822	A	C4-C5-C6	-6.37	113.82	117.00
1	CA	2574	G	N1-C6-O6	6.37	123.72	119.90
34	DA	729	A	C8-N9-C4	-6.37	103.25	105.80
56	DW	6	G	C8-N9-C4	6.37	108.95	106.40
1	AA	461	U	O5'-P-OP2	-6.36	99.97	105.70
1	AA	1387	U	O5'-P-OP1	-6.36	99.97	105.70
1	AA	1857	G	C5-C6-O6	6.36	132.42	128.60
1	CA	2584	U	N3-C2-O2	6.36	126.66	122.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1036	A	OP2-P-O3'	6.36	119.19	105.20
1	AA	1365	G	C5-C6-O6	6.36	132.42	128.60
1	CA	2610	C	N1-C2-O2	6.36	122.72	118.90
1	AA	2660	C	C5-C6-N1	-6.36	117.82	121.00
1	AA	207	A	C5-C6-N1	-6.36	114.52	117.70
1	AA	458	U	N3-C2-O2	-6.36	117.75	122.20
1	AA	578	U	C2-N3-C4	-6.36	123.19	127.00
1	AA	1252	C	C5-C4-N4	-6.36	115.75	120.20
1	AA	1472	G	N9-C4-C5	-6.36	102.86	105.40
1	AA	2223	C	N3-C2-O2	-6.36	117.45	121.90
1	AA	2657	G	C4-C5-N7	6.36	113.34	110.80
1	CA	776	G	C5-C6-O6	6.36	132.41	128.60
1	AA	133	G	OP2-P-O3'	6.36	119.18	105.20
1	AA	587	C	N1-C2-O2	-6.36	115.09	118.90
1	AA	1334	U	N1-C2-N3	6.36	118.71	114.90
1	AA	2057	G	N1-C6-O6	6.36	123.71	119.90
1	AA	2849	G	N3-C2-N2	6.36	124.35	119.90
1	CA	2713	A	N9-C4-C5	6.36	108.34	105.80
1	AA	884	C	N1-C2-N3	6.35	123.65	119.20
1	AA	1317	G	C5-C6-N1	6.35	114.68	111.50
1	AA	455	A	C5-N7-C8	-6.35	100.72	103.90
1	AA	905	U	C2-N3-C4	-6.35	123.19	127.00
1	CA	735	A	C2-N3-C4	6.35	113.78	110.60
1	CA	2561	A	O4'-C1'-N9	6.35	113.28	108.20
1	AA	599	U	C6-N1-C2	-6.35	117.19	121.00
1	AA	905	U	OP2-P-O3'	6.35	119.17	105.20
1	AA	2264	G	C4-N9-C1'	-6.35	118.25	126.50
4	AD	229	VAL	CB-CA-C	-6.35	99.33	111.40
1	CA	115	C	C6-N1-C2	6.35	122.84	120.30
34	DA	511	C	C6-N1-C2	-6.35	117.76	120.30
1	AA	541	C	N3-C4-C5	6.35	124.44	121.90
1	AA	1349	G	O5'-P-OP1	-6.35	99.99	105.70
1	AA	1474	C	N1-C2-O2	-6.35	115.09	118.90
34	BA	801	U	N3-C4-O4	-6.35	114.96	119.40
34	DA	583	A	C8-N9-C4	6.35	108.34	105.80
1	AA	171	A	OP2-P-O3'	6.35	119.16	105.20
1	AA	1067	A	C6-C5-N7	-6.35	127.86	132.30
1	AA	959	U	N3-C2-O2	6.34	126.64	122.20
1	AA	1834	A	OP2-P-O3'	6.34	119.16	105.20
1	AA	2236	G	C4-C5-N7	6.34	113.34	110.80
13	AP	14	LYS	CD-CE-NZ	-6.34	97.11	111.70
25	A1	40	ARG	NE-CZ-NH1	6.34	123.47	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	540	A	C4-C5-N7	-6.34	107.53	110.70
1	AA	1601	A	C8-N9-C4	-6.34	103.26	105.80
34	DA	758	G	N9-C4-C5	6.34	107.94	105.40
1	AA	336	G	C8-N9-C4	-6.34	103.86	106.40
1	AA	1642	A	N1-C6-N6	-6.34	114.80	118.60
34	BA	218	C	C6-N1-C2	-6.34	117.76	120.30
1	AA	881	C	OP1-P-OP2	6.34	129.11	119.60
1	AA	1244	U	OP2-P-O3'	6.34	119.15	105.20
1	AA	1319	U	N3-C4-O4	-6.34	114.96	119.40
1	AA	1948	U	C5-C4-O4	6.34	129.70	125.90
1	AA	2333	G	O5'-P-OP1	6.34	118.31	110.70
2	AB	54	G	C8-N9-C4	-6.34	103.86	106.40
1	CA	733	G	N9-C4-C5	-6.34	102.86	105.40
1	CA	965	C	O5'-P-OP2	-6.34	99.99	105.70
1	CA	1792	G	N3-C4-N9	6.34	129.80	126.00
1	CA	2708	G	N7-C8-N9	-6.34	109.93	113.10
1	AA	898	U	C5-C6-N1	-6.34	119.53	122.70
1	AA	2703	C	C2-N3-C4	-6.34	116.73	119.90
1	AA	535	C	N3-C4-N4	-6.34	113.56	118.00
1	AA	2537	G	OP2-P-O3'	6.34	119.14	105.20
1	CA	512	G	OP1-P-OP2	6.34	129.10	119.60
1	CA	563	G	N3-C4-C5	6.34	131.77	128.60
1	AA	2501	G	N1-C6-O6	6.33	123.70	119.90
1	AA	2544	G	N3-C2-N2	-6.33	115.47	119.90
1	AA	1300	A	C8-N9-C4	-6.33	103.27	105.80
1	AA	2833	A	N9-C4-C5	-6.33	103.27	105.80
34	DA	7	G	N3-C4-N9	-6.33	122.20	126.00
1	AA	733	G	N9-C4-C5	-6.33	102.87	105.40
34	DA	266	G	P-O3'-C3'	6.33	127.30	119.70
1	AA	1038	C	N1-C2-N3	6.33	123.63	119.20
1	AA	1713	G	C5-C6-O6	-6.33	124.80	128.60
1	CA	90	U	N3-C2-O2	-6.33	117.77	122.20
1	CA	515	A	N9-C4-C5	6.33	108.33	105.80
1	CA	769	G	N1-C6-O6	6.33	123.70	119.90
1	CA	1670	C	O5'-P-OP1	-6.33	100.00	105.70
1	CA	1614	A	O5'-P-OP1	-6.33	100.01	105.70
1	CA	2051	A	O5'-P-OP2	-6.33	100.00	105.70
1	AA	1252	C	N3-C4-N4	6.33	122.43	118.00
1	AA	1393	G	C8-N9-C4	-6.33	103.87	106.40
1	AA	2741	U	N3-C2-O2	-6.33	117.77	122.20
1	CA	127	A	OP1-P-O3'	6.33	119.11	105.20
1	CA	312	G	C4-N9-C1'	6.33	134.72	126.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1363	C	C5-C6-N1	-6.33	117.84	121.00
34	DA	755	G	N3-C2-N2	-6.33	115.47	119.90
1	AA	59	G	C5-C6-N1	6.32	114.66	111.50
1	AA	1475	G	N9-C4-C5	6.32	107.93	105.40
8	AH	3	ARG	CG-CD-NE	-6.32	98.52	111.80
1	AA	2499	G	C8-N9-C4	6.32	108.93	106.40
1	AA	2448	G	N7-C8-N9	6.32	116.26	113.10
1	CA	2576	G	N7-C8-N9	-6.32	109.94	113.10
1	AA	608	G	N7-C8-N9	-6.32	109.94	113.10
1	AA	2549	U	N3-C2-O2	-6.32	117.78	122.20
34	BA	1514	C	C4-C5-C6	6.32	120.56	117.40
1	AA	1733	C	C2-N1-C1'	6.32	125.75	118.80
34	BA	576	G	C5-C6-O6	-6.32	124.81	128.60
1	AA	826	U	C5-C4-O4	-6.32	122.11	125.90
1	AA	1431	G	C5-C6-O6	-6.32	124.81	128.60
1	AA	2609	G	N3-C4-C5	6.32	131.76	128.60
1	AA	865	G	C5-N7-C8	6.31	107.46	104.30
1	AA	875	U	OP2-P-O3'	6.31	119.09	105.20
1	AA	1011	G	N9-C4-C5	6.31	107.92	105.40
1	AA	1211	U	N3-C4-C5	6.31	118.39	114.60
1	AA	1660	A	O5'-P-OP1	-6.31	100.02	105.70
1	AA	2446	A	C4-C5-C6	6.31	120.16	117.00
1	CA	1615	C	C4-C5-C6	6.31	120.56	117.40
1	AA	1696	G	N3-C4-C5	-6.31	125.44	128.60
1	AA	2888	U	C6-N1-C2	-6.31	117.22	121.00
34	BA	303	A	C8-N9-C4	6.31	108.32	105.80
1	AA	2083	G	N1-C2-N3	6.31	127.69	123.90
1	AA	2452	C	C2-N3-C4	-6.31	116.75	119.90
2	AB	73	A	C4-C5-C6	-6.31	113.85	117.00
1	CA	190	A	O5'-P-OP2	-6.31	100.02	105.70
1	CA	1599	C	C5-C6-N1	-6.31	117.85	121.00
1	AA	114	C	C6-N1-C2	6.30	122.82	120.30
1	AA	1355	G	OP1-P-OP2	-6.30	110.14	119.60
1	AA	1678	A	OP1-P-O3'	6.30	119.07	105.20
1	AA	2227	G	C4-N9-C1'	-6.30	118.30	126.50
34	BA	912	C	C5-C4-N4	-6.30	115.79	120.20
1	AA	1735	U	OP2-P-O3'	6.30	119.07	105.20
1	CA	2549	G	N9-C4-C5	-6.30	102.88	105.40
1	AA	1416	C	N3-C2-O2	6.30	126.31	121.90
1	CA	1559	G	C8-N9-C4	6.30	108.92	106.40
1	CA	1772	G	N9-C1'-C2'	-6.30	105.07	112.00
1	AA	1282	G	N3-C2-N2	6.30	124.31	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2540	U	N1-C2-O2	-6.30	118.39	122.80
1	CA	98	G	N1-C6-O6	-6.30	116.12	119.90
1	AA	132	C	OP1-P-OP2	-6.30	110.15	119.60
1	AA	2735	G	C5-C6-N1	6.30	114.65	111.50
34	BA	873	A	C2-N3-C4	6.30	113.75	110.60
34	DA	821	G	O5'-P-OP1	-6.30	100.03	105.70
1	AA	257	C	N3-C4-C5	6.29	124.42	121.90
1	AA	490	U	OP1-P-OP2	-6.29	110.16	119.60
1	AA	1690	G	OP1-P-OP2	6.29	129.04	119.60
1	AA	1966	U	N1-C2-N3	6.29	118.68	114.90
1	AA	215	G	O4'-C1'-N9	6.29	113.23	108.20
1	AA	1283	A	C4-C5-N7	-6.29	107.55	110.70
1	AA	2559	U	N1-C2-O2	-6.29	118.39	122.80
1	CA	1260	G	C8-N9-C4	6.29	108.92	106.40
1	AA	1457	C	C6-N1-C2	6.29	122.82	120.30
1	AA	2397	C	OP1-P-OP2	6.29	129.04	119.60
1	AA	2409	G	C6-C5-N7	6.29	134.17	130.40
1	CA	751	A	O5'-P-OP1	-6.29	100.04	105.70
1	AA	1068	G	N3-C2-N2	-6.29	115.50	119.90
1	AA	1998	U	C5-C4-O4	-6.29	122.13	125.90
1	AA	873	U	OP1-P-O3'	-6.29	91.37	105.20
1	AA	1922	A	N9-C4-C5	6.29	108.31	105.80
1	AA	2464	C	C6-N1-C2	-6.29	117.78	120.30
34	BA	36	C	C6-N1-C2	-6.29	117.78	120.30
1	CA	2081	C	O5'-P-OP2	-6.29	100.04	105.70
1	CA	2689	U	P-O3'-C3'	6.29	127.25	119.70
34	DA	325	A	N1-C6-N6	-6.29	114.83	118.60
1	CA	801	G	C8-N9-C4	-6.29	103.89	106.40
1	AA	1249	A	C4-N9-C1'	6.29	137.62	126.30
1	AA	2051	G	C5-C6-O6	-6.29	124.83	128.60
2	AB	81	G	N3-C4-N9	-6.29	122.23	126.00
1	CA	1267	U	C5-C6-N1	6.29	125.84	122.70
1	CA	2364	C	O5'-P-OP1	6.29	118.24	110.70
1	AA	1342	G	N3-C4-C5	-6.28	125.46	128.60
1	CA	2348	U	N3-C4-O4	-6.28	115.00	119.40
34	DA	619	U	C5-C6-N1	-6.28	119.56	122.70
2	AB	34	U	N3-C4-O4	6.28	123.80	119.40
34	BA	769	G	OP1-P-OP2	-6.28	110.18	119.60
1	CA	2507	C	N3-C4-C5	-6.28	119.39	121.90
1	AA	562	C	N3-C4-N4	-6.28	113.60	118.00
1	AA	634	C	C5-C6-N1	-6.28	117.86	121.00
1	AA	1186	U	N1-C2-O2	-6.28	118.40	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1605	C	C5-C6-N1	-6.28	117.86	121.00
1	AA	30	G	C5-C6-O6	6.28	132.37	128.60
1	AA	867	A	OP2-P-O3'	6.28	119.01	105.20
1	AA	2025	G	N9-C4-C5	6.28	107.91	105.40
34	BA	564	C	C6-N1-C2	-6.28	117.79	120.30
1	AA	439	A	C8-N9-C4	-6.28	103.29	105.80
1	AA	871	A	OP2-P-O3'	6.28	119.01	105.20
1	AA	2252	C	N3-C4-C5	6.28	124.41	121.90
34	BA	1425	U	N3-C4-C5	6.28	118.37	114.60
1	CA	989	G	C5-C6-N1	6.28	114.64	111.50
1	AA	2273	C	O5'-P-OP2	-6.28	100.05	105.70
1	AA	2298	A	C4-N9-C1'	6.28	137.59	126.30
2	AB	41	U	N1-C2-N3	6.28	118.67	114.90
1	CA	2013	A	OP1-P-OP2	-6.28	110.19	119.60
34	DA	863	U	C5-C4-O4	6.28	129.67	125.90
1	AA	1400	A	OP2-P-O3'	6.27	119.00	105.20
1	AA	1802	C	C4-C5-C6	6.27	120.54	117.40
1	CA	2079	U	N3-C4-O4	6.27	123.79	119.40
1	CA	2447	G	C8-N9-C1'	6.27	135.16	127.00
1	CA	2604	U	N3-C4-C5	6.27	118.36	114.60
1	AA	2248	C	C6-N1-C2	6.27	122.81	120.30
34	BA	1490	C	N1-C2-O2	6.27	122.66	118.90
1	CA	2519	U	O5'-P-OP1	-6.27	100.05	105.70
1	CA	2782	G	N1-C6-O6	6.27	123.66	119.90
1	CA	2525	G	OP2-P-O3'	6.27	119.00	105.20
1	AA	244	A	N1-C2-N3	6.27	132.43	129.30
1	AA	2748	G	C5-C6-O6	6.27	132.36	128.60
1	AA	2755	C	C6-N1-C2	-6.27	117.79	120.30
34	BA	194	C	C6-N1-C2	-6.27	117.79	120.30
1	CA	1558	A	N3-C4-C5	6.27	131.19	126.80
1	AA	10	G	N1-C6-O6	-6.26	116.14	119.90
1	AA	107	G	C4-C5-N7	-6.26	108.29	110.80
1	AA	416	G	C5-C6-N1	6.26	114.63	111.50
1	AA	1808	U	C5-C6-N1	-6.26	119.57	122.70
1	AA	1894	G	C8-N9-C4	6.26	108.91	106.40
1	AA	2107	C	C5-C6-N1	-6.26	117.87	121.00
1	AA	2279	A	OP1-P-OP2	6.26	129.00	119.60
1	CA	2069	G	N3-C4-N9	-6.26	122.24	126.00
1	AA	2460	A	C5-C6-N6	-6.26	118.69	123.70
1	AA	2495	C	C5-C6-N1	6.26	124.13	121.00
1	CA	1266	G	N7-C8-N9	-6.26	109.97	113.10
1	AA	1299	A	C4-C5-C6	-6.26	113.87	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2092	G	C6-N1-C2	-6.26	121.34	125.10
1	AA	2542	A	N1-C6-N6	6.26	122.36	118.60
55	BV	16	U	O4'-C1'-N1	-6.26	103.19	108.20
1	AA	843	C	N3-C4-C5	6.26	124.40	121.90
1	CA	1559	G	N3-C4-C5	6.26	131.73	128.60
34	DA	576	G	C8-N9-C4	6.26	108.90	106.40
1	AA	1319	U	N3-C4-C5	6.26	118.36	114.60
1	AA	1845	G	C4-C5-N7	6.26	113.30	110.80
1	AA	756	U	OP1-P-O3'	-6.26	91.44	105.20
1	AA	917	A	O5'-P-OP2	-6.26	100.07	105.70
1	AA	1132	A	N1-C6-N6	6.26	122.35	118.60
34	BA	665	A	O5'-P-OP1	-6.26	100.07	105.70
1	AA	2022	G	C2-N3-C4	6.25	115.03	111.90
1	AA	2643	G	C5-C6-O6	6.25	132.35	128.60
1	AA	990	A	C4-N9-C1'	6.25	137.56	126.30
1	AA	1665	G	N7-C8-N9	-6.25	109.97	113.10
1	AA	1806	U	N3-C2-O2	-6.25	117.82	122.20
1	AA	1872	U	N3-C2-O2	-6.25	117.82	122.20
1	AA	2277	U	C2-N3-C4	-6.25	123.25	127.00
2	AB	41	U	N3-C2-O2	-6.25	117.82	122.20
1	AA	1655	A	C6-N1-C2	-6.25	114.85	118.60
1	CA	2272	U	N3-C2-O2	-6.25	117.82	122.20
1	AA	2529	C	O5'-P-OP1	6.25	118.20	110.70
1	CA	1836	C	O5'-P-OP2	-6.25	100.08	105.70
1	AA	36	G	O5'-P-OP2	-6.25	100.08	105.70
1	AA	1276	C	N1-C2-N3	6.25	123.58	119.20
1	AA	2636	G	C5-C6-N1	6.25	114.62	111.50
1	CA	1299	G	N3-C4-N9	-6.25	122.25	126.00
1	CA	2851	A	C2-N3-C4	-6.25	107.48	110.60
34	DA	689	C	C6-N1-C2	-6.25	117.80	120.30
1	AA	998	A	N1-C2-N3	6.25	132.42	129.30
1	CA	25	U	C6-N1-C2	6.25	124.75	121.00
1	AA	704	U	C5-C6-N1	-6.24	119.58	122.70
1	AA	865	G	C5-C6-O6	6.24	132.35	128.60
1	AA	1082	G	C4-C5-N7	-6.24	108.30	110.80
29	A5	17	ASP	CB-CG-OD2	6.24	123.92	118.30
1	AA	315	C	C2-N3-C4	-6.24	116.78	119.90
1	AA	716	G	C4-C5-N7	6.24	113.30	110.80
1	AA	172	C	C5-C6-N1	-6.24	117.88	121.00
1	AA	866	A	C2-N3-C4	6.24	113.72	110.60
1	AA	2299	A	N1-C6-N6	6.24	122.34	118.60
14	AQ	135	ASP	CB-CA-C	-6.24	97.92	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	123	C	N3-C2-O2	6.24	126.27	121.90
1	CA	1294	U	C5-C6-N1	6.24	125.82	122.70
1	AA	995	G	C4-C5-N7	6.24	113.30	110.80
1	AA	1071	G	O5'-P-OP2	6.24	118.19	110.70
1	AA	1282	G	OP2-P-O3'	6.24	118.92	105.20
1	AA	2238	C	C6-N1-C2	6.24	122.80	120.30
1	AA	168	G	N1-C2-N3	6.24	127.64	123.90
34	DA	872	A	O4'-C1'-N9	6.24	113.19	108.20
1	AA	1725	G	C6-C5-N7	-6.24	126.66	130.40
1	CA	1807	G	C5-N7-C8	6.24	107.42	104.30
1	AA	413	G	N7-C8-N9	6.23	116.22	113.10
1	CA	1320	C	N1-C2-O2	-6.23	115.16	118.90
1	AA	2780	C	N3-C4-C5	6.23	124.39	121.90
11	AN	65	LYS	CD-CE-NZ	6.23	126.04	111.70
1	AA	70	A	O5'-P-OP2	-6.23	100.09	105.70
1	AA	1379	C	C6-N1-C2	6.23	122.79	120.30
1	AA	1701	A	N9-C4-C5	-6.23	103.31	105.80
1	AA	2036	A	N1-C6-N6	6.23	122.34	118.60
1	CA	2413	G	O5'-P-OP2	-6.23	100.09	105.70
1	AA	1239	A	C2-N3-C4	6.23	113.72	110.60
1	CA	19	C	N3-C2-O2	-6.23	117.54	121.90
34	DA	886	G	C8-N9-C4	6.23	108.89	106.40
1	AA	823	G	C4-C5-N7	-6.23	108.31	110.80
1	AA	2281	A	C2-N3-C4	-6.23	107.49	110.60
1	CA	144	C	C6-N1-C2	-6.23	117.81	120.30
34	BA	115	G	O5'-P-OP2	-6.23	100.10	105.70
1	AA	521	G	C8-N9-C4	6.22	108.89	106.40
1	AA	1819	C	C5-C6-N1	-6.22	117.89	121.00
1	AA	1046	A	N9-C4-C5	-6.22	103.31	105.80
1	AA	1210	G	C5-C6-O6	6.22	132.33	128.60
1	AA	1238	G	C6-N1-C2	-6.22	121.37	125.10
1	AA	1720	U	C6-N1-C2	6.22	124.73	121.00
1	AA	1925	G	OP2-P-O3'	6.22	118.89	105.20
2	AB	114	C	C6-N1-C2	6.22	122.79	120.30
12	AO	8	LEU	CA-CB-CG	6.22	129.61	115.30
1	AA	397	G	O4'-C1'-N9	-6.22	103.22	108.20
1	AA	912	C	N1-C2-O2	-6.22	115.17	118.90
1	AA	2066	C	C6-N1-C2	-6.22	117.81	120.30
1	CA	2695	C	C2-N1-C1'	-6.22	111.96	118.80
34	DA	365	U	C5-C6-N1	-6.22	119.59	122.70
34	DA	885	G	C8-N9-C4	6.22	108.89	106.40
1	CA	1775	U	N1-C2-O2	-6.22	118.45	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	508	A	O5'-P-OP2	-6.22	100.11	105.70
1	AA	641	G	N7-C8-N9	-6.22	109.99	113.10
1	AA	879	G	N1-C6-O6	-6.22	116.17	119.90
1	AA	1001	G	N3-C4-C5	6.22	131.71	128.60
1	AA	2083	G	N3-C4-C5	-6.22	125.49	128.60
1	AA	2240	G	C5-C6-O6	6.22	132.33	128.60
1	AA	500	G	N7-C8-N9	6.21	116.21	113.10
1	AA	734	C	O5'-P-OP2	-6.21	100.11	105.70
1	AA	2610	A	N9-C4-C5	-6.21	103.31	105.80
1	AA	2896	G	N1-C2-N3	6.21	127.63	123.90
1	AA	310	C	C5-C6-N1	-6.21	117.89	121.00
1	AA	989	G	C5-C6-N1	6.21	114.61	111.50
1	AA	2899	C	N3-C4-C5	6.21	124.39	121.90
1	AA	1015	C	C4-C5-C6	6.21	120.51	117.40
1	AA	1479	U	C5-C6-N1	-6.21	119.59	122.70
1	AA	1922	A	C5-C6-N1	6.21	120.81	117.70
34	BA	219	C	C6-N1-C2	-6.21	117.82	120.30
34	BA	910	C	O5'-P-OP2	-6.21	100.11	105.70
1	CA	2572	A	C4-C5-N7	-6.21	107.59	110.70
1	AA	793	A	N1-C2-N3	6.21	132.41	129.30
1	AA	1003	U	C2-N3-C4	6.21	130.72	127.00
1	CA	915	C	C6-N1-C2	-6.21	117.82	120.30
1	AA	625	G	C8-N9-C4	6.21	108.88	106.40
1	CA	53	A	C5-N7-C8	6.21	107.00	103.90
1	AA	76	C	N3-C4-C5	6.21	124.38	121.90
1	AA	139	A	N3-C4-C5	6.21	131.14	126.80
1	AA	249	G	C4-C5-N7	-6.21	108.32	110.80
1	AA	351	G	N1-C6-O6	-6.20	116.18	119.90
1	AA	802	C	N1-C2-O2	-6.20	115.18	118.90
1	AA	1370	G	OP1-P-O3'	6.20	118.85	105.20
1	AA	1453	C	N3-C4-C5	6.20	124.38	121.90
1	CA	334	C	N3-C2-O2	-6.20	117.56	121.90
1	CA	1795	C	N3-C4-C5	6.20	124.38	121.90
1	AA	426	G	O5'-P-OP1	6.20	118.14	110.70
1	AA	1013	G	N9-C4-C5	6.20	107.88	105.40
1	AA	1362	U	N3-C4-C5	-6.20	110.88	114.60
1	AA	2070	G	N1-C2-N2	-6.20	110.62	116.20
34	BA	365	U	N1-C2-O2	-6.20	118.46	122.80
34	BA	553	A	N1-C6-N6	6.20	122.32	118.60
34	DA	898	G	C8-N9-C4	6.20	108.88	106.40
1	AA	209	G	C8-N9-C4	6.20	108.88	106.40
1	AA	637	U	N3-C4-O4	-6.20	115.06	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1835	G	C8-N9-C1'	-6.20	118.94	127.00
39	DF	79	LEU	CA-CB-CG	6.20	129.56	115.30
1	AA	511	C	N1-C2-N3	6.20	123.54	119.20
1	AA	543	G	O5'-P-OP1	6.20	118.14	110.70
1	CA	2326	C	O5'-P-OP1	-6.20	100.12	105.70
1	AA	2716	C	C6-N1-C2	-6.20	117.82	120.30
1	CA	489	G	OP2-P-O3'	6.20	118.83	105.20
1	AA	637	U	N1-C2-N3	6.19	118.62	114.90
1	AA	2226	C	C6-N1-C2	-6.19	117.82	120.30
1	AA	2560	G	C2-N3-C4	6.19	115.00	111.90
1	AA	2635	G	N1-C2-N2	-6.19	110.62	116.20
1	CA	1367	A	N1-C6-N6	-6.19	114.88	118.60
1	AA	57	G	C6-N1-C2	-6.19	121.39	125.10
1	AA	592	U	N1-C2-O2	-6.19	118.47	122.80
1	AA	2294	G	N1-C2-N3	-6.19	120.19	123.90
34	BA	553	A	O5'-P-OP1	6.19	118.13	110.70
1	AA	498	A	C5-C6-N1	-6.19	114.61	117.70
1	AA	1157	A	C4-C5-N7	6.19	113.80	110.70
1	AA	2470	G	O5'-P-OP2	-6.19	100.13	105.70
1	CA	1957	C	N1-C2-O2	-6.19	115.19	118.90
1	AA	12	U	C2-N1-C1'	6.19	125.13	117.70
1	AA	815	G	N9-C4-C5	6.19	107.88	105.40
1	AA	1659	G	C6-N1-C2	-6.19	121.39	125.10
1	AA	1860	A	C5-N7-C8	6.19	106.99	103.90
1	AA	2585	C	C2-N1-C1'	-6.19	111.99	118.80
1	CA	408	G	N9-C4-C5	-6.19	102.92	105.40
1	CA	1694	C	C6-N1-C1'	-6.19	113.38	120.80
1	AA	147	U	N3-C4-C5	6.19	118.31	114.60
1	AA	564	G	N7-C8-N9	-6.19	110.01	113.10
1	AA	910	A	C2-N3-C4	6.19	113.69	110.60
1	CA	1636	C	C5-C6-N1	-6.19	117.91	121.00
1	AA	717	A	C5-N7-C8	-6.18	100.81	103.90
1	AA	1317	G	C8-N9-C4	6.18	108.87	106.40
1	AA	1507	A	O4'-C1'-N9	6.18	113.15	108.20
1	AA	2270	C	N3-C4-C5	6.18	124.37	121.90
1	AA	2434	A	C4-C5-C6	6.18	120.09	117.00
1	AA	2626	A	C8-N9-C4	-6.18	103.33	105.80
1	CA	989	G	N1-C6-O6	-6.18	116.19	119.90
1	CA	1945	G	OP2-P-O3'	6.18	118.81	105.20
1	CA	2249	U	N3-C4-O4	-6.18	115.07	119.40
1	AA	2544	G	N1-C6-O6	6.18	123.61	119.90
1	CA	2893	G	C4-N9-C1'	6.18	134.54	126.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1826	C	OP1-P-OP2	6.18	128.87	119.60
1	AA	22	C	C6-N1-C2	6.18	122.77	120.30
1	AA	860	U	N1-C2-N3	6.18	118.61	114.90
1	AA	1056	A	N9-C4-C5	6.18	108.27	105.80
1	CA	1674	G	OP1-P-OP2	6.18	128.87	119.60
34	BA	550	G	N3-C2-N2	-6.18	115.58	119.90
1	CA	193	U	N3-C4-O4	6.18	123.72	119.40
1	CA	2272	U	N1-C2-O2	6.18	127.12	122.80
1	AA	819	C	N1-C2-O2	-6.18	115.19	118.90
1	AA	1788	U	C6-N1-C2	6.18	124.71	121.00
1	AA	2876	U	C2-N3-C4	-6.18	123.29	127.00
1	CA	2033	A	O4'-C1'-N9	6.18	113.14	108.20
1	AA	53	G	N3-C4-C5	-6.17	125.51	128.60
1	AA	176	G	N1-C2-N3	6.17	127.61	123.90
1	CA	2593	U	N3-C4-C5	6.17	118.30	114.60
1	AA	409	G	N3-C2-N2	6.17	124.22	119.90
1	AA	421	A	N9-C4-C5	6.17	108.27	105.80
1	AA	465	G	C4-C5-N7	6.17	113.27	110.80
1	AA	1699	A	N7-C8-N9	6.17	116.89	113.80
1	AA	1894	G	O5'-P-OP2	-6.17	100.14	105.70
6	AF	45	ARG	NE-CZ-NH2	6.17	123.39	120.30
1	CA	769	G	C4-C5-N7	6.17	113.27	110.80
1	CA	1790	C	P-O3'-C3'	6.17	127.11	119.70
1	CA	1796	U	C5-C6-N1	-6.17	119.61	122.70
1	CA	2356	C	N1-C2-O2	-6.17	115.20	118.90
1	AA	829	A	C6-N1-C2	-6.17	114.90	118.60
1	AA	1964	C	N1-C2-O2	-6.17	115.20	118.90
1	AA	2029	C	OP2-P-O3'	6.17	118.77	105.20
1	AA	2162	C	N3-C2-O2	-6.17	117.58	121.90
1	AA	2559	U	C5-C4-O4	-6.17	122.20	125.90
1	CA	600	G	N9-C4-C5	6.17	107.87	105.40
1	CA	778	G	N1-C6-O6	6.17	123.60	119.90
56	BW	73	A	N1-C6-N6	6.17	122.30	118.60
1	AA	210	A	N1-C6-N6	6.17	122.30	118.60
1	AA	575	G	C6-C5-N7	6.17	134.10	130.40
1	AA	1258	A	C8-N9-C4	-6.17	103.33	105.80
1	CA	2722	G	N3-C4-C5	-6.17	125.52	128.60
1	CA	1284	A	C5-N7-C8	-6.16	100.82	103.90
1	AA	2252	C	N3-C2-O2	6.16	126.21	121.90
1	AA	2697	G	N1-C6-O6	-6.16	116.20	119.90
1	CA	1331	A	C8-N9-C4	6.16	108.27	105.80
1	AA	1050	C	C5-C6-N1	-6.16	117.92	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2239	A	N9-C4-C5	6.16	108.26	105.80
1	CA	786	C	OP1-P-OP2	-6.16	110.36	119.60
1	CA	1154	G	C6-C5-N7	-6.16	126.70	130.40
1	AA	129	G	C4-C5-N7	6.16	113.26	110.80
1	AA	195	U	C2-N3-C4	-6.16	123.31	127.00
1	AA	2071	G	C5-N7-C8	-6.16	101.22	104.30
1	CA	2598	A	C8-N9-C4	6.16	108.26	105.80
1	AA	1021	G	C5-C6-O6	6.16	132.29	128.60
1	AA	1814	A	O4'-C1'-N9	-6.16	103.28	108.20
1	AA	2058	C	N3-C4-C5	6.16	124.36	121.90
1	CA	496	G	C2-N3-C4	-6.16	108.82	111.90
1	CA	330	A	C5-C6-N1	-6.15	114.62	117.70
1	AA	49	U	OP2-P-O3'	6.15	118.74	105.20
1	AA	2518	U	OP1-P-O3'	-6.15	91.67	105.20
1	AA	2610	A	N1-C6-N6	6.15	122.29	118.60
1	AA	2703	C	N3-C4-C5	6.15	124.36	121.90
34	BA	1484	C	N3-C2-O2	6.15	126.21	121.90
34	BA	1505	G	C6-C5-N7	6.15	134.09	130.40
1	CA	1648	C	N1-C2-O2	-6.15	115.21	118.90
1	AA	327	U	C6-N1-C2	6.15	124.69	121.00
1	AA	1009	C	C6-N1-C2	-6.15	117.84	120.30
1	AA	1258	A	C5-N7-C8	-6.15	100.83	103.90
1	CA	2242	G	N9-C4-C5	-6.15	102.94	105.40
1	AA	429	A	N1-C6-N6	-6.15	114.91	118.60
1	AA	1365	G	N1-C6-O6	-6.15	116.21	119.90
1	AA	1682	G	C6-N1-C2	-6.15	121.41	125.10
1	AA	1805	C	N1-C2-N3	6.15	123.50	119.20
1	AA	1849	U	C2-N3-C4	-6.15	123.31	127.00
32	A8	13	ARG	NE-CZ-NH2	-6.15	117.23	120.30
1	CA	2253	G	N1-C6-O6	6.15	123.59	119.90
34	DA	518	C	N1-C2-O2	6.15	122.59	118.90
1	CA	2673	G	OP1-P-OP2	6.15	128.82	119.60
1	AA	31	C	C6-N1-C2	-6.14	117.84	120.30
1	AA	39	C	C2-N3-C4	-6.14	116.83	119.90
1	AA	85	C	C5-C4-N4	6.14	124.50	120.20
1	AA	1007	G	C5-C6-O6	-6.14	124.91	128.60
1	AA	1376	C	N3-C2-O2	6.14	126.20	121.90
1	AA	1728	G	C5-N7-C8	-6.14	101.23	104.30
1	AA	2341	G	N1-C6-O6	-6.14	116.21	119.90
2	AB	55	U	C5-C6-N1	-6.14	119.63	122.70
1	CA	1914	C	N3-C2-O2	-6.14	117.60	121.90
1	CA	2318	G	C8-N9-C4	6.14	108.86	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2429	G	N3-C2-N2	-6.14	115.60	119.90
1	AA	237	G	C2-N3-C4	6.14	114.97	111.90
1	AA	1783	C	N3-C2-O2	6.14	126.20	121.90
1	AA	1989	C	OP1-P-O3'	6.14	118.71	105.20
1	AA	2287	C	C5'-C4'-O4'	-6.14	101.73	109.10
1	AA	2802	C	C2-N1-C1'	-6.14	112.04	118.80
1	CA	777	A	N1-C2-N3	6.14	132.37	129.30
1	CA	1350	C	N1-C2-O2	-6.14	115.21	118.90
1	AA	2015	U	C5-C4-O4	6.14	129.59	125.90
1	CA	2499	C	N3-C4-N4	6.14	122.30	118.00
1	AA	369	A	N1-C6-N6	6.14	122.28	118.60
1	AA	465	G	N1-C6-O6	6.14	123.58	119.90
1	AA	2073	A	C4-C5-N7	-6.14	107.63	110.70
1	AA	2266	C	C6-N1-C2	6.14	122.76	120.30
1	CA	141	A	C8-N9-C4	-6.14	103.34	105.80
1	CA	2512	C	C4-C5-C6	6.14	120.47	117.40
34	DA	720	C	C6-N1-C2	-6.14	117.84	120.30
1	AA	972	A	N9-C4-C5	6.14	108.25	105.80
1	AA	2060	G	N3-C2-N2	6.14	124.20	119.90
1	AA	2076	A	OP2-P-O3'	6.14	118.70	105.20
56	BW	17	C	C5-C6-N1	6.14	124.07	121.00
1	CA	2499	C	N1-C2-O2	-6.14	115.22	118.90
2	CB	30	C	C6-N1-C2	-6.14	117.84	120.30
1	AA	1866	G	O5'-P-OP1	6.14	118.06	110.70
34	BA	562	C	N3-C4-C5	6.14	124.36	121.90
1	CA	1430	C	N3-C2-O2	6.14	126.19	121.90
1	CA	1814	G	C5-C6-O6	-6.14	124.92	128.60
1	CA	2443	C	C6-N1-C2	6.14	122.75	120.30
1	CA	2874	C	N3-C4-C5	-6.14	119.44	121.90
1	AA	1656	A	C6-N1-C2	-6.13	114.92	118.60
1	AA	2602	A	N7-C8-N9	-6.13	110.73	113.80
2	AB	102	A	O5'-P-OP2	6.13	118.06	110.70
20	AW	11	ARG	NE-CZ-NH2	-6.13	117.23	120.30
34	BA	1412	C	C2-N3-C4	-6.13	116.83	119.90
34	DA	691	G	N1-C6-O6	6.13	123.58	119.90
34	DA	890	G	N3-C4-N9	6.13	129.68	126.00
34	BA	902	G	O5'-P-OP2	-6.13	100.18	105.70
34	BA	909	A	OP1-P-OP2	6.13	128.80	119.60
1	AA	753	A	C4-C5-N7	6.13	113.77	110.70
1	AA	772	G	C5-N7-C8	6.13	107.36	104.30
1	AA	1265	A	OP2-P-O3'	6.13	118.69	105.20
1	AA	2828	G	N1-C2-N3	6.13	127.58	123.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	675	C	N3-C4-N4	-6.13	113.71	118.00
1	CA	1683	C	C6-N1-C2	-6.13	117.85	120.30
1	AA	322	G	C5-N7-C8	6.13	107.36	104.30
1	AA	845	G	C6-C5-N7	6.13	134.08	130.40
34	DA	1505	G	N3-C4-N9	-6.13	122.32	126.00
1	AA	460	C	C2-N3-C4	-6.13	116.84	119.90
1	AA	628	C	OP1-P-OP2	-6.13	110.41	119.60
1	AA	1659	G	O5'-P-OP1	6.13	118.05	110.70
1	AA	1795	G	N1-C2-N2	-6.13	110.69	116.20
1	AA	1859	G	C5-C6-O6	6.13	132.28	128.60
1	AA	2602	A	N1-C2-N3	6.13	132.36	129.30
1	CA	277	C	N3-C2-O2	-6.13	117.61	121.90
1	CA	1329	U	OP1-P-O3'	6.13	118.68	105.20
1	AA	150	C	C5-C6-N1	-6.12	117.94	121.00
1	AA	451	G	N7-C8-N9	-6.12	110.04	113.10
1	AA	2796	G	OP2-P-O3'	6.12	118.67	105.20
34	BA	1064	G	C8-N9-C1'	6.12	134.96	127.00
1	AA	543	G	N3-C2-N2	6.12	124.19	119.90
1	AA	845	G	C5-C6-O6	6.12	132.27	128.60
1	AA	345	G	N9-C4-C5	-6.12	102.95	105.40
1	AA	1486	G	C8-N9-C4	6.12	108.85	106.40
1	CA	12	U	N1-C2-O2	6.12	127.08	122.80
1	AA	667	G	O5'-P-OP2	-6.12	100.19	105.70
1	AA	1812	C	C2-N3-C4	-6.12	116.84	119.90
1	AA	2162	C	C6-N1-C2	-6.12	117.85	120.30
1	CA	1314	C	C5-C4-N4	-6.12	115.92	120.20
1	AA	1423	G	N9-C4-C5	6.12	107.85	105.40
1	AA	2277	U	N1-C2-N3	6.12	118.57	114.90
1	AA	2525	G	O5'-P-OP2	-6.12	100.20	105.70
1	CA	205	G	OP1-P-OP2	6.12	128.77	119.60
1	CA	531	C	O5'-P-OP1	-6.12	100.20	105.70
1	CA	1396	U	N1-C2-O2	6.12	127.08	122.80
1	CA	1963	U	N3-C2-O2	-6.12	117.92	122.20
1	CA	2828	C	N3-C4-C5	6.12	124.35	121.90
1	AA	1824	C	OP2-P-O3'	6.11	118.65	105.20
1	AA	2014	G	OP1-P-OP2	-6.11	110.43	119.60
1	CA	2575	C	C4-C5-C6	6.11	120.46	117.40
1	AA	491	G	N1-C2-N3	6.11	127.57	123.90
1	AA	23	G	C5-C6-N1	-6.11	108.44	111.50
1	AA	2295	C	N3-C4-C5	-6.11	119.45	121.90
1	AA	2611	G	N3-C2-N2	6.11	124.18	119.90
1	AA	2696	U	N1-C2-N3	6.11	118.57	114.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2769	U	OP1-P-O3'	6.11	118.64	105.20
56	BW	76	A	N1-C6-N6	6.11	122.27	118.60
1	CA	312	G	N7-C8-N9	6.11	116.16	113.10
1	CA	481	G	O5'-P-OP2	-6.11	100.20	105.70
1	CA	2683	C	N3-C4-C5	-6.11	119.46	121.90
1	AA	745	C	N1-C2-O2	-6.11	115.23	118.90
1	AA	1660	A	OP1-P-O3'	-6.11	91.76	105.20
1	AA	1817	A	OP2-P-O3'	-6.11	91.76	105.20
1	AA	2096	U	C4-C5-C6	6.11	123.37	119.70
34	BA	893	C	N3-C4-N4	6.11	122.28	118.00
1	AA	357	G	N3-C4-C5	-6.11	125.55	128.60
1	AA	549	U	C5-C6-N1	-6.11	119.65	122.70
1	AA	2312	G	C8-N9-C4	-6.11	103.96	106.40
1	CA	2611	U	N1-C2-N3	6.11	118.56	114.90
1	AA	2654	G	C5-C6-O6	6.11	132.26	128.60
1	AA	2834	C	N3-C4-N4	6.11	122.27	118.00
1	CA	307	G	C2-N3-C4	6.11	114.95	111.90
1	AA	1357	G	C8-N9-C4	-6.10	103.96	106.40
1	AA	83	A	O4'-C1'-N9	6.10	113.08	108.20
1	AA	616	G	C5-N7-C8	6.10	107.35	104.30
1	AA	1440	U	OP1-P-OP2	-6.10	110.45	119.60
1	AA	1770	A	N1-C6-N6	6.10	122.26	118.60
1	AA	2383	G	C4-C5-N7	6.10	113.24	110.80
1	CA	64	A	C5-C6-N6	-6.10	118.82	123.70
1	AA	850	U	C2-N3-C4	-6.10	123.34	127.00
1	AA	1357	G	O5'-P-OP2	-6.10	100.21	105.70
1	AA	1988	A	C5-N7-C8	-6.10	100.85	103.90
1	CA	503	A	O4'-C1'-N9	6.10	113.08	108.20
57	DZ	229	LEU	CA-CB-CG	6.10	129.33	115.30
1	AA	554	A	C5-N7-C8	-6.10	100.85	103.90
1	AA	865	G	O5'-P-OP2	6.10	118.02	110.70
1	AA	1704	C	OP1-P-O3'	-6.10	91.78	105.20
1	AA	2402	U	C2-N3-C4	-6.10	123.34	127.00
1	AA	2867	G	OP1-P-OP2	6.10	128.75	119.60
2	AB	35	U	N3-C2-O2	-6.10	117.93	122.20
34	BA	579	G	N1-C6-O6	6.10	123.56	119.90
56	BW	45	U	O4'-C1'-N1	6.10	113.08	108.20
1	CA	1330	C	N1-C2-O2	-6.10	115.24	118.90
1	AA	597	C	N1-C2-O2	-6.10	115.24	118.90
1	AA	1065	U	C6-N1-C2	6.10	124.66	121.00
1	AA	1082	G	C8-N9-C4	6.10	108.84	106.40
1	AA	1652	G	N1-C6-O6	-6.10	116.24	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2710	U	N1-C2-O2	-6.10	118.53	122.80
1	AA	2867	G	C8-N9-C4	6.10	108.84	106.40
1	CA	1405	U	O5'-P-OP2	-6.10	100.21	105.70
1	CA	2713	A	C8-N9-C4	-6.10	103.36	105.80
1	AA	1628	G	C4-N9-C1'	6.10	134.43	126.50
34	BA	297	G	N1-C6-O6	6.10	123.56	119.90
1	AA	2428	C	OP2-P-O3'	6.09	118.61	105.20
1	AA	2342	G	C5-C6-N1	6.09	114.55	111.50
1	AA	2386	C	N3-C4-C5	6.09	124.34	121.90
1	AA	2566	U	N1-C2-N3	6.09	118.56	114.90
34	BA	23	C	OP1-P-OP2	-6.09	110.46	119.60
1	AA	399	G	O4'-C1'-N9	6.09	113.07	108.20
1	AA	791	G	N3-C2-N2	6.09	124.16	119.90
1	AA	2250	G	N7-C8-N9	6.09	116.14	113.10
1	AA	2372	A	N9-C4-C5	6.09	108.23	105.80
1	AA	2411	G	N1-C2-N2	-6.09	110.72	116.20
1	AA	2884	C	C6-N1-C2	-6.09	117.86	120.30
1	CA	1936	A	N9-C4-C5	-6.09	103.36	105.80
1	CA	2604	U	N1-C2-O2	6.09	127.06	122.80
1	AA	2380	C	C5-C6-N1	-6.09	117.96	121.00
1	AA	86	C	O5'-P-OP2	-6.09	100.22	105.70
1	AA	989	G	C6-N1-C2	-6.09	121.45	125.10
1	AA	1283	A	C4-C5-C6	6.09	120.04	117.00
2	AB	99	G	N1-C2-N2	-6.09	110.72	116.20
1	AA	290	G	N7-C8-N9	-6.08	110.06	113.10
1	AA	721	G	C5-C6-N1	6.08	114.54	111.50
34	BA	1524	C	C6-N1-C2	-6.08	117.87	120.30
1	CA	378	C	C6-N1-C2	6.08	122.73	120.30
1	AA	203	G	O4'-C1'-N9	6.08	113.07	108.20
1	AA	534	C	C6-N1-C2	6.08	122.73	120.30
1	AA	2586	G	C4-C5-C6	-6.08	115.15	118.80
22	AY	21	LYS	CD-CE-NZ	6.08	125.69	111.70
34	DA	800	G	N1-C6-O6	6.08	123.55	119.90
1	AA	82	G	C5-C6-O6	-6.08	124.95	128.60
1	AA	139	A	N3-C4-N9	-6.08	122.53	127.40
1	AA	445	G	C5-C6-O6	6.08	132.25	128.60
1	AA	1448	C	N3-C4-C5	6.08	124.33	121.90
1	AA	2397	C	C4-C5-C6	-6.08	114.36	117.40
1	AA	2551	C	C5-C4-N4	6.08	124.46	120.20
34	BA	543	C	C6-N1-C2	-6.08	117.87	120.30
1	AA	416	G	C8-N9-C4	6.08	108.83	106.40
1	AA	2488	A	C5-N7-C8	-6.08	100.86	103.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	AD	222	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	CA	154	G	C8-N9-C4	6.08	108.83	106.40
1	AA	1829	U	N3-C4-C5	6.08	118.25	114.60
1	AA	2244	U	C5-C4-O4	6.08	129.55	125.90
1	AA	2739	U	O5'-P-OP1	6.08	117.99	110.70
1	AA	2846	U	O5'-P-OP2	6.08	117.99	110.70
1	CA	19	C	C6-N1-C2	-6.08	117.87	120.30
1	CA	1623	G	N1-C6-O6	6.08	123.55	119.90
56	DW	6	G	N3-C4-C5	6.08	131.64	128.60
1	AA	737	G	N3-C2-N2	6.07	124.15	119.90
1	AA	846	G	C6-N1-C2	-6.07	121.46	125.10
1	AA	847	A	C4-C5-N7	-6.07	107.66	110.70
1	AA	1204	C	N1-C2-N3	6.07	123.45	119.20
1	AA	337	C	C5-C6-N1	-6.07	117.96	121.00
1	AA	2043	C	O5'-P-OP2	-6.07	100.23	105.70
34	BA	1482	G	N1-C6-O6	-6.07	116.26	119.90
1	CA	836	G	C2-N3-C4	6.07	114.94	111.90
1	CA	2394	C	C4-C5-C6	6.07	120.44	117.40
1	AA	907	U	N3-C4-C5	6.07	118.24	114.60
2	AB	107	G	N1-C6-O6	6.07	123.54	119.90
34	BA	504	C	C6-N1-C2	-6.07	117.87	120.30
1	AA	2896	G	O5'-P-OP2	6.07	117.98	110.70
1	CA	998	C	N1-C2-O2	6.07	122.54	118.90
1	CA	1367	A	N7-C8-N9	-6.07	110.77	113.80
1	AA	310	C	C2-N3-C4	-6.07	116.87	119.90
1	AA	1665	G	C8-N9-C4	6.07	108.83	106.40
1	AA	2372	A	C8-N9-C4	-6.07	103.37	105.80
1	CA	202	U	C5-C4-O4	-6.07	122.26	125.90
1	CA	1975	G	C4-C5-N7	6.07	113.23	110.80
1	AA	1028	C	N1-C2-O2	-6.07	115.26	118.90
1	AA	1312	G	C8-N9-C1'	6.07	134.88	127.00
1	AA	1754	G	C8-N9-C1'	-6.07	119.11	127.00
1	AA	2526	U	N3-C4-O4	-6.07	115.15	119.40
2	AB	63	G	O5'-P-OP2	-6.07	100.24	105.70
1	AA	2384	G	C5-C6-O6	-6.06	124.96	128.60
2	AB	38	C	N3-C4-N4	-6.06	113.75	118.00
1	CA	2832	U	C6-N1-C2	6.06	124.64	121.00
1	AA	640	A	N7-C8-N9	6.06	116.83	113.80
1	AA	1244	U	N3-C4-O4	-6.06	115.16	119.40
1	AA	1268	C	C4-C5-C6	6.06	120.43	117.40
1	AA	1373	C	N3-C2-O2	6.06	126.14	121.90
1	AA	2385	G	C5-C6-O6	-6.06	124.96	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2619	G	C5-C6-N1	6.06	114.53	111.50
1	AA	2620	G	C5-C6-N1	6.06	114.53	111.50
56	BW	17	C	C6-N1-C2	-6.06	117.88	120.30
1	CA	203	C	N3-C4-C5	6.06	124.33	121.90
1	AA	1037	C	C5-C4-N4	-6.06	115.96	120.20
1	AA	2054	G	C2-N3-C4	-6.06	108.87	111.90
1	AA	2794	A	C5-C6-N1	6.06	120.73	117.70
1	CA	2487	G	N3-C2-N2	6.06	124.14	119.90
1	AA	606	G	C2-N3-C4	6.06	114.93	111.90
1	AA	2641	A	N9-C1'-C2'	6.06	121.88	114.00
1	CA	2451	A	P-O3'-C3'	-6.06	112.43	119.70
1	AA	873	U	C5-C6-N1	-6.05	119.67	122.70
1	AA	1241	C	OP2-P-O3'	6.05	118.52	105.20
1	AA	2250	G	C2-N3-C4	6.05	114.93	111.90
1	CA	790	C	N3-C4-N4	6.05	122.24	118.00
1	CA	2576	G	C5-C6-O6	-6.05	124.97	128.60
1	AA	2414	C	C6-N1-C2	6.05	122.72	120.30
34	DA	919	A	N1-C6-N6	6.05	122.23	118.60
1	AA	715	G	N9-C4-C5	-6.05	102.98	105.40
1	AA	1303	C	C4-C5-C6	6.05	120.42	117.40
1	AA	1824	C	C2-N3-C4	-6.05	116.87	119.90
1	AA	35	G	C5-N7-C8	6.05	107.33	104.30
1	AA	1244	U	C5-C4-O4	6.05	129.53	125.90
1	AA	1690	G	O5'-P-OP2	-6.05	100.25	105.70
1	AA	1707	C	C2-N3-C4	-6.05	116.88	119.90
2	AB	5	C	N1-C2-O2	-6.05	115.27	118.90
34	BA	733	A	OP1-P-OP2	6.05	128.67	119.60
1	CA	2239	G	N1-C6-O6	-6.05	116.27	119.90
1	AA	50	G	N1-C2-N3	6.05	127.53	123.90
1	AA	2792	U	C6-N1-C2	-6.05	117.37	121.00
1	CA	1654	A	O5'-P-OP1	-6.05	100.26	105.70
1	AA	1188	A	O5'-P-OP1	-6.05	100.26	105.70
1	AA	2020	G	N9-C4-C5	-6.05	102.98	105.40
1	AA	2189	U	C2-N1-C1'	6.05	124.95	117.70
1	AA	2272	C	OP2-P-O3'	6.05	118.50	105.20
34	BA	578	C	N3-C2-O2	-6.05	117.67	121.90
31	C7	33	ARG	NE-CZ-NH1	-6.05	117.28	120.30
34	DA	1183	A	P-O3'-C3'	6.05	126.96	119.70
1	AA	633	G	C5-C6-O6	6.04	132.23	128.60
1	AA	2837	C	N3-C4-C5	6.04	124.32	121.90
1	AA	2571	C	N1-C2-O2	6.04	122.53	118.90
34	BA	581	G	N3-C4-N9	-6.04	122.37	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	528	A	C5-C6-N6	6.04	128.54	123.70
1	AA	489	G	OP1-P-O3'	6.04	118.49	105.20
1	AA	1665	G	C5-C6-O6	-6.04	124.97	128.60
1	CA	60	G	C6-C5-N7	-6.04	126.78	130.40
1	CA	1647	G	C5-C6-O6	-6.04	124.97	128.60
1	CA	2089	U	C4-C5-C6	6.04	123.33	119.70
1	CA	2570	G	N3-C2-N2	-6.04	115.67	119.90
1	AA	374	U	N3-C4-O4	-6.04	115.17	119.40
1	AA	555	G	C6-N1-C2	6.04	128.72	125.10
1	CA	1668	A	O5'-P-OP1	-6.04	100.26	105.70
1	CA	2576	G	C8-N9-C4	6.04	108.82	106.40
1	AA	990	A	N9-C1'-C2'	6.04	121.85	114.00
1	AA	2331	G	N9-C1'-C2'	6.04	121.85	114.00
34	DA	769	G	C5-C6-O6	-6.04	124.98	128.60
1	AA	231	G	C2-N3-C4	-6.04	108.88	111.90
34	BA	1286	A	C8-N9-C4	-6.04	103.39	105.80
1	CA	1329	U	C5-C6-N1	-6.04	119.68	122.70
1	AA	28	A	C6-N1-C2	6.04	122.22	118.60
1	AA	1707	C	N3-C4-C5	6.04	124.31	121.90
1	AA	2351	G	C5-C6-O6	6.04	132.22	128.60
1	AA	2359	C	N3-C2-O2	-6.04	117.67	121.90
1	AA	2461	U	C2-N3-C4	-6.04	123.38	127.00
1	AA	2600	G	C4-C5-N7	6.04	113.21	110.80
1	AA	2646	G	C5-C6-O6	-6.04	124.98	128.60
1	AA	2738	A	C8-N9-C4	6.04	108.21	105.80
1	AA	2753	A	C8-N9-C4	6.04	108.21	105.80
1	CA	191	A	OP1-P-O3'	-6.04	91.92	105.20
1	CA	2698	U	N3-C4-C5	6.04	118.22	114.60
2	CB	70	C	C5-C6-N1	6.04	124.02	121.00
1	AA	1001	G	C4-C5-C6	6.03	122.42	118.80
1	AA	1056	A	C4-C5-N7	-6.03	107.68	110.70
1	AA	1188	A	C8-N9-C1'	6.03	138.56	127.70
1	AA	2019	G	N3-C4-C5	-6.03	125.58	128.60
1	AA	2760	G	C5-C6-O6	-6.03	124.98	128.60
1	AA	2834	C	C5-C4-N4	-6.03	115.98	120.20
1	CA	2444	G	C5-C6-O6	6.03	132.22	128.60
1	AA	1067	A	C8-N9-C1'	6.03	138.56	127.70
1	AA	913	A	C4-N9-C1'	6.03	137.16	126.30
1	AA	1526	G	C4-C5-N7	6.03	113.21	110.80
1	AA	2059	G	N3-C2-N2	6.03	124.12	119.90
1	CA	1377	G	C6-C5-N7	-6.03	126.78	130.40
1	CA	2010	G	O5'-P-OP2	6.03	117.94	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	411	U	N3-C4-C5	6.03	118.22	114.60
1	AA	642	G	OP2-P-O3'	6.03	118.46	105.20
34	BA	1346	A	O4'-C1'-N9	6.03	113.02	108.20
1	CA	607	U	O5'-P-OP1	-6.03	100.27	105.70
1	CA	2486	G	C8-N9-C1'	-6.03	119.16	127.00
1	CA	2755	C	C5-C6-N1	6.03	124.01	121.00
34	DA	328	C	N3-C4-N4	6.03	122.22	118.00
1	AA	97	G	N1-C6-O6	-6.03	116.28	119.90
1	AA	738	C	C2-N3-C4	-6.03	116.89	119.90
1	AA	756	U	OP2-P-O3'	6.03	118.46	105.20
1	AA	841	G	N3-C2-N2	6.03	124.12	119.90
1	AA	894	U	C4-C5-C6	6.03	123.32	119.70
1	AA	2021	C	N3-C4-N4	-6.03	113.78	118.00
1	AA	2650	G	N3-C4-C5	-6.03	125.59	128.60
1	AA	2711	C	C6-N1-C2	6.03	122.71	120.30
34	BA	1201	A	P-O3'-C3'	6.03	126.93	119.70
1	AA	194	G	O5'-P-OP1	-6.02	100.28	105.70
1	AA	2498	G	C5-C6-O6	6.02	132.22	128.60
1	CA	1798	U	C2-N3-C4	-6.02	123.39	127.00
1	CA	1835	G	C4-N9-C1'	6.02	134.33	126.50
1	AA	1064	C	N3-C2-O2	-6.02	117.68	121.90
1	AA	2012	C	N3-C4-N4	6.02	122.22	118.00
1	AA	2070	G	N1-C6-O6	-6.02	116.29	119.90
1	AA	2895	C	N3-C4-C5	-6.02	119.49	121.90
1	AA	352	U	N3-C4-C5	6.02	118.21	114.60
1	CA	2038	G	N3-C2-N2	6.02	124.11	119.90
1	CA	2395	C	C2-N3-C4	-6.02	116.89	119.90
1	AA	1524	A	O5'-P-OP2	-6.02	100.28	105.70
1	AA	1487	G	N3-C2-N2	-6.02	115.69	119.90
56	BW	36	A	N3-C4-N9	-6.02	122.59	127.40
1	CA	1653	G	N7-C8-N9	6.02	116.11	113.10
1	AA	123	G	OP2-P-O3'	6.01	118.43	105.20
1	AA	859	C	C5-C6-N1	-6.01	117.99	121.00
1	AA	342	C	N1-C2-O2	-6.01	115.29	118.90
29	A5	20	ARG	NE-CZ-NH2	-6.01	117.29	120.30
56	BW	36	A	N3-C4-C5	6.01	131.01	126.80
1	AA	719	C	C6-N1-C2	-6.01	117.90	120.30
1	AA	1211	U	N1-C2-N3	6.01	118.51	114.90
1	CA	2607	G	N3-C2-N2	6.01	124.11	119.90
1	AA	585	U	N3-C4-O4	-6.01	115.19	119.40
1	AA	1817	A	N1-C2-N3	-6.01	126.30	129.30
1	CA	2675	A	OP1-P-OP2	-6.01	110.59	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1565	G	C5-C6-O6	6.01	132.20	128.60
1	AA	1725	G	C8-N9-C4	-6.01	104.00	106.40
1	CA	750	A	OP1-P-O3'	6.01	118.42	105.20
37	DD	157	LEU	CA-CB-CG	6.01	129.12	115.30
1	AA	207	A	N1-C6-N6	6.00	122.20	118.60
1	AA	1295	U	C2-N3-C4	-6.00	123.40	127.00
1	AA	1670	G	C5-C6-O6	6.00	132.20	128.60
1	CA	729	G	N3-C2-N2	6.00	124.10	119.90
1	AA	723	A	C6-N1-C2	6.00	122.20	118.60
1	AA	989	G	C4-C5-N7	6.00	113.20	110.80
1	CA	963	U	OP2-P-O3'	6.00	118.40	105.20
34	DA	913	A	P-O3'-C3'	6.00	126.90	119.70
1	AA	437	G	C2-N3-C4	-6.00	108.90	111.90
34	BA	652	U	N3-C2-O2	6.00	126.40	122.20
1	CA	37	C	N1-C2-O2	-6.00	115.30	118.90
1	CA	2071	A	O5'-P-OP1	-6.00	100.30	105.70
1	AA	444	C	O5'-P-OP1	-6.00	100.30	105.70
1	AA	1247	C	C6-N1-C2	6.00	122.70	120.30
1	AA	1984	C	C5-C6-N1	6.00	124.00	121.00
1	AA	2440	G	N3-C2-N2	6.00	124.10	119.90
1	CA	2612	C	O5'-P-OP2	-6.00	100.30	105.70
34	DA	583	A	N9-C4-C5	-6.00	103.40	105.80
39	DF	87	ARG	NE-CZ-NH1	6.00	123.30	120.30
1	AA	120	G	C2-N3-C4	6.00	114.90	111.90
1	AA	2037	A	N9-C4-C5	6.00	108.20	105.80
1	AA	2084	A	C4-C5-N7	-6.00	107.70	110.70
1	AA	2087	C	N3-C4-N4	-6.00	113.80	118.00
1	AA	2510	C	N1-C2-O2	-6.00	115.30	118.90
1	AA	2513	C	C2-N1-C1'	-6.00	112.20	118.80
34	BA	885	G	N1-C6-O6	6.00	123.50	119.90
1	CA	2381	C	O5'-P-OP1	6.00	117.89	110.70
1	AA	2060	G	N1-C2-N2	-5.99	110.81	116.20
1	AA	2484	G	C2-N3-C4	5.99	114.90	111.90
1	CA	2678	C	C5-C6-N1	-5.99	118.00	121.00
1	AA	17	G	N3-C2-N2	5.99	124.09	119.90
1	AA	325	G	N9-C4-C5	-5.99	103.00	105.40
1	AA	1734	G	C5-C6-O6	5.99	132.19	128.60
1	AA	2355	C	OP1-P-OP2	5.99	128.59	119.60
1	CA	2286	A	C5-N7-C8	-5.99	100.90	103.90
1	AA	114	C	N1-C2-O2	-5.99	115.31	118.90
1	AA	357	G	N1-C6-O6	-5.99	116.31	119.90
1	AA	2507	G	C5-C6-N1	-5.99	108.50	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	509	A	C4-C5-C6	5.99	120.00	117.00
1	CA	407	G	OP2-P-O3'	5.99	118.38	105.20
1	AA	886	U	N3-C4-C5	5.99	118.19	114.60
1	AA	998	A	O5'-P-OP2	5.99	117.89	110.70
34	BA	509	A	C6-C5-N7	-5.99	128.11	132.30
1	CA	312	G	O5'-P-OP1	-5.99	100.31	105.70
1	CA	2775	A	N1-C6-N6	5.99	122.19	118.60
34	DA	589	C	C6-N1-C2	-5.99	117.91	120.30
1	AA	233	A	C5-C6-N6	-5.99	118.91	123.70
1	AA	350	G	OP2-P-O3'	5.99	118.37	105.20
34	BA	718	G	C4-C5-N7	5.99	113.19	110.80
1	CA	1957	C	C6-N1-C2	-5.99	117.91	120.30
1	AA	1068	G	N3-C4-C5	5.99	131.59	128.60
1	AA	1274	G	N3-C4-N9	-5.99	122.41	126.00
1	AA	1310	G	O5'-P-OP2	-5.99	100.31	105.70
1	AA	1548	C	N3-C4-N4	5.99	122.19	118.00
34	BA	498	U	O5'-P-OP2	-5.99	100.31	105.70
1	CA	204	A	OP1-P-OP2	5.99	128.58	119.60
1	CA	2500	U	N3-C4-C5	5.99	118.19	114.60
1	AA	2331	G	O4'-C1'-N9	5.98	112.99	108.20
1	AA	2402	U	N3-C4-O4	-5.98	115.21	119.40
1	AA	29	U	N1-C2-N3	5.98	118.49	114.90
1	AA	405	C	C2-N3-C4	-5.98	116.91	119.90
57	BZ	88	VAL	CB-CA-C	-5.98	100.03	111.40
1	CA	1313	U	C6-N1-C2	-5.98	117.41	121.00
1	AA	751	G	C8-N9-C4	-5.98	104.01	106.40
1	AA	1744	G	OP1-P-O3'	5.98	118.36	105.20
1	AA	2778	A	O5'-P-OP1	-5.98	100.32	105.70
1	AA	2855	G	C2-N3-C4	5.98	114.89	111.90
1	CA	763	G	OP1-P-OP2	5.98	128.57	119.60
1	CA	1860	G	N7-C8-N9	5.98	116.09	113.10
1	AA	896	A	C5-C6-N1	-5.98	114.71	117.70
1	AA	1812	C	C6-N1-C1'	-5.98	113.62	120.80
1	AA	187	C	C5-C4-N4	-5.98	116.02	120.20
1	AA	2431	U	N3-C4-O4	-5.98	115.22	119.40
1	AA	2655	G	N1-C2-N3	5.98	127.49	123.90
1	CA	201	C	C2-N3-C4	-5.98	116.91	119.90
1	CA	964	C	N1-C2-O2	5.98	122.49	118.90
1	CA	2070	G	O5'-P-OP1	5.98	117.87	110.70
1	AA	34	C	C6-N1-C2	-5.98	117.91	120.30
1	AA	1847	G	N3-C4-N9	-5.97	122.42	126.00
1	AA	2776	G	OP2-P-O3'	5.97	118.34	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AB	85	G	N3-C4-N9	5.97	129.59	126.00
1	CA	940	G	N1-C6-O6	-5.97	116.31	119.90
1	CA	2436	G	O5'-P-OP1	-5.97	100.32	105.70
1	AA	785	G	C6-C5-N7	-5.97	126.82	130.40
1	AA	828	A	C5-N7-C8	5.97	106.89	103.90
1	AA	872	C	C2-N3-C4	-5.97	116.91	119.90
1	AA	1370	G	OP1-P-OP2	-5.97	110.64	119.60
1	AA	1477	U	N3-C4-O4	5.97	123.58	119.40
1	AA	1721	G	N1-C2-N3	5.97	127.48	123.90
1	CA	2079	U	C5-C6-N1	-5.97	119.71	122.70
1	CA	2721	A	O5'-P-OP1	-5.97	100.33	105.70
1	AA	991	G	N3-C4-C5	5.97	131.59	128.60
34	BA	1384	C	C5-C6-N1	5.97	123.99	121.00
1	AA	1846	A	OP1-P-OP2	5.97	128.55	119.60
1	AA	2098	U	C4-C5-C6	5.97	123.28	119.70
1	CA	1313	U	C5-C6-N1	5.97	125.69	122.70
1	CA	1675	C	O5'-P-OP2	5.97	117.86	110.70
1	AA	2399	U	OP1-P-O3'	-5.97	92.07	105.20
1	CA	376	C	C6-N1-C2	-5.97	117.91	120.30
34	DA	442	C	C5-C6-N1	5.97	123.98	121.00
1	AA	732	A	C5-N7-C8	-5.97	100.92	103.90
1	AA	980	C	N3-C4-C5	5.97	124.29	121.90
1	AA	1816	A	OP1-P-OP2	-5.97	110.65	119.60
1	AA	1847	G	N3-C2-N2	-5.97	115.72	119.90
1	AA	2370	G	C4-C5-N7	5.97	113.19	110.80
1	AA	2527	C	C6-N1-C2	5.97	122.69	120.30
1	AA	2551	C	N3-C4-N4	-5.97	113.82	118.00
1	AA	184	A	OP1-P-OP2	-5.96	110.65	119.60
1	AA	213	G	C5-C6-O6	5.96	132.18	128.60
1	AA	565	C	O5'-P-OP2	-5.96	100.33	105.70
1	AA	1705	C	C2-N3-C4	-5.96	116.92	119.90
1	AA	1859	G	C5-N7-C8	5.96	107.28	104.30
1	AA	2611	G	C2-N3-C4	5.96	114.88	111.90
1	CA	1397	U	C5-C6-N1	5.96	125.68	122.70
1	AA	854	U	C5-C4-O4	5.96	129.48	125.90
1	AA	1195	G	N9-C4-C5	5.96	107.78	105.40
1	AA	1252	C	C4-C5-C6	5.96	120.38	117.40
34	BA	731	G	O5'-P-OP2	-5.96	100.33	105.70
1	CA	822	U	C4-C5-C6	5.96	123.28	119.70
1	AA	1029	A	OP1-P-OP2	-5.96	110.66	119.60
1	AA	1846	A	C8-N9-C4	-5.96	103.42	105.80
1	AA	2372	A	O5'-P-OP1	5.96	117.85	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2455	C	C4-C5-C6	5.96	120.38	117.40
1	AA	2546	A	C5-C6-N1	5.96	120.68	117.70
1	AA	969	C	N1-C2-O2	-5.96	115.32	118.90
1	CA	2584	U	C6-N1-C2	5.96	124.58	121.00
1	AA	702	A	C8-N9-C4	-5.96	103.42	105.80
1	AA	757	G	C2-N3-C4	-5.96	108.92	111.90
1	AA	1022	C	N3-C4-N4	-5.96	113.83	118.00
1	AA	2528	G	C6-C5-N7	5.96	133.97	130.40
1	CA	2379	G	C6-C5-N7	-5.96	126.83	130.40
1	AA	244	A	C8-N9-C4	-5.96	103.42	105.80
1	AA	2849	G	C8-N9-C4	5.96	108.78	106.40
34	BA	687	A	P-O3'-C3'	5.96	126.85	119.70
1	AA	2015	U	C4-C5-C6	5.96	123.27	119.70
1	AA	2480	G	N3-C4-N9	5.96	129.57	126.00
1	CA	668	G	C4-C5-N7	5.96	113.18	110.80
1	CA	784	A	O5'-P-OP1	-5.96	100.34	105.70
1	CA	2242	G	C8-N9-C4	5.96	108.78	106.40
1	AA	502	G	N3-C4-N9	5.95	129.57	126.00
1	AA	797	A	N7-C8-N9	5.95	116.78	113.80
34	BA	1416	G	C4-C5-N7	-5.95	108.42	110.80
1	AA	126	C	C6-N1-C2	5.95	122.68	120.30
1	AA	989	G	OP1-P-OP2	5.95	128.53	119.60
1	AA	473	A	N1-C6-N6	-5.95	115.03	118.60
1	AA	530	A	C8-N9-C4	-5.95	103.42	105.80
1	AA	718	C	N3-C2-O2	-5.95	117.73	121.90
1	AA	1340	U	C4-C5-C6	5.95	123.27	119.70
1	AA	2102	G	N1-C2-N3	5.95	127.47	123.90
1	AA	2399	U	OP2-P-O3'	5.95	118.29	105.20
1	AA	2726	A	C4-C5-N7	-5.95	107.72	110.70
1	CA	1301	A	OP1-P-OP2	5.95	128.53	119.60
1	CA	1975	G	N9-C4-C5	-5.95	103.02	105.40
1	CA	2073	C	C5-C4-N4	-5.95	116.03	120.20
1	AA	50	G	C6-N1-C2	-5.95	121.53	125.10
1	AA	240	A	C5-N7-C8	5.95	106.87	103.90
34	BA	815	A	C5-N7-C8	5.95	106.88	103.90
1	CA	945	A	N3-C4-N9	-5.95	122.64	127.40
1	CA	1668	A	C5-C6-N6	-5.95	118.94	123.70
1	CA	2586	C	C2-N3-C4	-5.95	116.93	119.90
1	AA	1022	C	OP2-P-O3'	5.95	118.28	105.20
1	AA	199	C	OP2-P-O3'	5.95	118.28	105.20
1	AA	416	G	N7-C8-N9	-5.95	110.13	113.10
1	AA	540	A	N9-C4-C5	5.95	108.18	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1655	A	N1-C6-N6	5.95	122.17	118.60
1	CA	567	A	C5-N7-C8	-5.94	100.93	103.90
1	AA	705	C	C5-C6-N1	-5.94	118.03	121.00
1	AA	1518	A	N1-C2-N3	5.94	132.27	129.30
1	AA	1597	C	C5-C4-N4	-5.94	116.04	120.20
1	AA	2458	G	N9-C4-C5	5.94	107.78	105.40
1	CA	1154	G	C4-N9-C1'	5.94	134.22	126.50
1	CA	2393	A	C8-N9-C4	-5.94	103.42	105.80
34	DA	314	C	N1-C2-O2	-5.94	115.33	118.90
34	DA	1469	G	N3-C2-N2	-5.94	115.74	119.90
1	AA	431	C	N3-C4-C5	5.94	124.28	121.90
1	AA	460	C	OP2-P-O3'	5.94	118.27	105.20
1	AA	777	C	C2-N3-C4	-5.94	116.93	119.90
1	AA	1472	G	C2-N3-C4	5.94	114.87	111.90
1	AA	1724	A	C4-C5-C6	5.94	119.97	117.00
1	AA	2801	C	N1-C2-O2	-5.94	115.34	118.90
34	BA	797	C	O5'-P-OP2	5.94	117.83	110.70
1	CA	1359	A	N1-C2-N3	5.94	132.27	129.30
1	AA	985	G	C2-N3-C4	5.94	114.87	111.90
1	AA	1737	A	N1-C6-N6	-5.94	115.04	118.60
1	AA	2279	A	P-O3'-C3'	5.94	126.83	119.70
1	CA	513	A	C5-N7-C8	-5.94	100.93	103.90
1	AA	1376	C	C6-N1-C2	5.94	122.67	120.30
1	AA	1453	C	C2-N3-C4	-5.94	116.93	119.90
1	AA	1704	C	C6-N1-C2	5.94	122.67	120.30
1	AA	2757	G	C8-N9-C4	5.94	108.78	106.40
1	AA	2881	C	C5-C6-N1	-5.94	118.03	121.00
34	BA	652	U	N1-C2-N3	-5.94	111.34	114.90
1	AA	1003	U	O5'-P-OP2	-5.94	100.36	105.70
1	AA	1449	C	C5-C6-N1	-5.94	118.03	121.00
1	AA	1317	G	C6-N1-C2	-5.93	121.54	125.10
1	AA	2219	U	C5-C6-N1	-5.93	119.73	122.70
1	AA	2355	C	N1-C2-N3	5.93	123.35	119.20
2	AB	81	G	C2-N3-C4	-5.93	108.93	111.90
1	CA	2567	G	C8-N9-C4	5.93	108.77	106.40
1	AA	872	C	OP1-P-OP2	-5.93	110.70	119.60
1	CA	1615	C	N3-C4-C5	-5.93	119.53	121.90
34	DA	1406	U	C5-C6-N1	-5.93	119.73	122.70
1	AA	478	G	N7-C8-N9	-5.93	110.14	113.10
1	AA	838	C	C2-N1-C1'	-5.93	112.28	118.80
1	CA	1846	G	C4-C5-N7	-5.93	108.43	110.80
1	AA	491	G	C8-N9-C4	-5.93	104.03	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1988	A	N3-C4-N9	-5.93	122.66	127.40
1	CA	214	G	N7-C8-N9	-5.93	110.14	113.10
1	CA	1785	A	N1-C2-N3	5.93	132.26	129.30
1	CA	1946	U	O5'-P-OP2	-5.93	100.36	105.70
1	CA	2486	G	N3-C4-N9	5.93	129.56	126.00
34	DA	1464	G	O5'-P-OP1	-5.93	100.36	105.70
1	CA	2330	G	C6-C5-N7	-5.93	126.84	130.40
1	AA	2342	G	C5-C6-O6	-5.93	125.05	128.60
1	AA	2561	G	C5-N7-C8	5.93	107.26	104.30
34	BA	1067	A	P-O3'-C3'	5.93	126.81	119.70
1	CA	2509	G	N1-C6-O6	5.93	123.45	119.90
1	AA	1472	G	N3-C4-N9	5.92	129.56	126.00
1	AA	1725	G	N1-C2-N3	5.92	127.45	123.90
1	AA	2255	U	C4-C5-C6	5.92	123.25	119.70
1	AA	2296	C	C6-N1-C2	-5.92	117.93	120.30
1	AA	2403	G	C2-N3-C4	-5.92	108.94	111.90
1	AA	2448	G	N3-C4-N9	-5.92	122.44	126.00
2	AB	7	G	N3-C4-C5	5.92	131.56	128.60
34	BA	1529	G	C8-N9-C4	-5.92	104.03	106.40
1	CA	690	G	N3-C4-C5	-5.92	125.64	128.60
1	CA	1997	G	OP2-P-O3'	5.92	118.23	105.20
1	CA	1985	G	O5'-P-OP2	-5.92	100.37	105.70
1	AA	208	G	C5-C6-N1	5.92	114.46	111.50
1	AA	208	G	N3-C2-N2	5.92	124.05	119.90
1	AA	1248	G	N3-C4-N9	5.92	129.55	126.00
1	AA	1733	C	O5'-P-OP2	5.92	117.81	110.70
1	AA	1802	C	C5-C6-N1	-5.92	118.04	121.00
1	CA	1801	G	C8-N9-C4	5.92	108.77	106.40
1	CA	2494	G	C8-N9-C4	5.92	108.77	106.40
1	AA	1431	G	N3-C2-N2	-5.92	115.76	119.90
1	AA	1653	C	N3-C4-C5	-5.92	119.53	121.90
1	AA	2036	A	C5-N7-C8	5.92	106.86	103.90
1	AA	2046	G	C5-N7-C8	5.92	107.26	104.30
4	AD	54	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	AA	138	G	C8-N9-C4	-5.92	104.03	106.40
1	AA	837	C	N1-C2-O2	5.92	122.45	118.90
1	AA	1008	U	O5'-P-OP2	-5.92	100.37	105.70
1	AA	1270	C	C4-C5-C6	5.92	120.36	117.40
1	AA	1334	U	C5-C6-N1	-5.92	119.74	122.70
1	AA	1342	G	C6-C5-N7	5.92	133.95	130.40
1	AA	2073	A	N1-C6-N6	-5.92	115.05	118.60
1	CA	1770	G	N3-C2-N2	5.92	124.04	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	CB	78	A	N1-C2-N3	5.92	132.26	129.30
34	DA	283	C	N1-C2-O2	5.92	122.45	118.90
1	AA	1514	C	OP1-P-OP2	-5.92	110.72	119.60
1	AA	182	U	OP2-P-O3'	5.92	118.21	105.20
1	CA	474	G	C4-C5-N7	-5.92	108.43	110.80
1	CA	2295	C	N1-C2-O2	5.92	122.45	118.90
1	AA	502	G	N3-C2-N2	5.91	124.04	119.90
1	CA	1253	A	O4'-C1'-N9	-5.91	103.47	108.20
34	DA	577	G	C2-N3-C4	-5.91	108.94	111.90
1	AA	985	G	OP2-P-O3'	5.91	118.21	105.20
16	CS	67	ARG	NE-CZ-NH2	-5.91	117.34	120.30
1	AA	2667	G	O5'-P-OP2	-5.91	100.38	105.70
1	AA	989	G	O5'-P-OP2	-5.91	100.38	105.70
1	AA	2243	C	C4-C5-C6	5.91	120.35	117.40
1	CA	801	G	N9-C4-C5	5.91	107.76	105.40
1	AA	1518	A	C8-N9-C4	-5.91	103.44	105.80
1	AA	1787	G	N1-C2-N2	-5.91	110.88	116.20
1	AA	582	G	N7-C8-N9	5.91	116.05	113.10
1	AA	1003	U	OP1-P-OP2	5.91	128.46	119.60
1	AA	2599	A	C2-N3-C4	-5.91	107.65	110.60
8	AH	41	MET	CG-SD-CE	-5.91	90.75	100.20
1	AA	2076	A	C5-N7-C8	5.90	106.85	103.90
1	AA	211	A	O5'-P-OP1	-5.90	100.39	105.70
1	AA	829	A	C5-C6-N1	5.90	120.65	117.70
1	AA	1316	C	OP1-P-OP2	-5.90	110.75	119.60
1	AA	2580	C	C4-C5-C6	5.90	120.35	117.40
34	BA	894	G	C8-N9-C4	-5.90	104.04	106.40
1	CA	586	A	C8-N9-C4	5.90	108.16	105.80
1	CA	1376	C	N3-C4-N4	5.90	122.13	118.00
1	CA	2465	C	N3-C2-O2	-5.90	117.77	121.90
1	AA	1664	A	N1-C6-N6	-5.90	115.06	118.60
1	AA	1684	A	N3-C4-C5	5.90	130.93	126.80
1	AA	2068	G	N7-C8-N9	-5.90	110.15	113.10
1	AA	2600	G	N3-C4-C5	5.90	131.55	128.60
1	AA	2722	C	O5'-P-OP1	-5.90	100.39	105.70
34	BA	672	U	C5-C4-O4	5.90	129.44	125.90
1	CA	562	U	C5-C4-O4	5.90	129.44	125.90
1	CA	790	C	N3-C2-O2	5.90	126.03	121.90
34	DA	509	A	N7-C8-N9	5.90	116.75	113.80
1	AA	981	C	N3-C4-N4	5.90	122.13	118.00
1	AA	499	G	C5-C6-O6	5.90	132.14	128.60
1	AA	555	G	N9-C4-C5	5.90	107.76	105.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1287	A	C8-N9-C4	5.90	108.16	105.80
1	AA	2480	G	N3-C4-C5	-5.90	125.65	128.60
1	AA	2684	G	N3-C2-N2	-5.90	115.77	119.90
1	AA	2697	G	C6-C5-N7	5.90	133.94	130.40
34	BA	1069	C	C6-N1-C2	-5.90	117.94	120.30
34	DA	1079	G	C8-N9-C4	-5.90	104.04	106.40
1	AA	461	U	N1-C2-O2	-5.90	118.67	122.80
1	AA	96	C	N3-C2-O2	-5.89	117.77	121.90
1	AA	1852	A	N9-C4-C5	5.89	108.16	105.80
1	AA	2505	U	C4-C5-C6	5.89	123.24	119.70
2	AB	55	U	N1-C2-O2	-5.89	118.67	122.80
34	BA	766	A	C5-C6-N6	-5.89	118.98	123.70
1	CA	702	G	N1-C6-O6	-5.89	116.36	119.90
1	CA	2386	C	N1-C2-O2	-5.89	115.36	118.90
34	DA	275	G	N1-C6-O6	5.89	123.44	119.90
1	AA	21	A	N1-C6-N6	5.89	122.14	118.60
1	AA	341	G	C5-C6-O6	5.89	132.13	128.60
1	AA	557	A	C6-N1-C2	-5.89	115.06	118.60
1	AA	2852	G	OP1-P-O3'	5.89	118.17	105.20
1	CA	125	G	C4-C5-N7	5.89	113.16	110.80
1	CA	1423	G	C8-N9-C4	5.89	108.76	106.40
1	AA	585	U	O4'-C1'-N1	-5.89	103.49	108.20
1	CA	393	C	N1-C2-O2	-5.89	115.37	118.90
1	AA	60	G	C2-N3-C4	-5.89	108.95	111.90
1	AA	980	C	N1-C2-O2	-5.89	115.37	118.90
1	AA	1188	A	C5-C6-N6	5.89	128.41	123.70
1	AA	2655	G	C5-C6-N1	-5.89	108.56	111.50
1	AA	2745	G	N1-C2-N3	5.89	127.43	123.90
2	AB	65	C	O5'-P-OP2	5.89	117.77	110.70
1	CA	2224	G	N1-C6-O6	5.89	123.43	119.90
1	AA	178	G	OP1-P-OP2	5.89	128.43	119.60
1	AA	705	C	C2-N3-C4	-5.89	116.96	119.90
1	AA	2437	A	N1-C6-N6	5.89	122.13	118.60
1	AA	176	G	N7-C8-N9	5.89	116.04	113.10
1	AA	471	C	N3-C2-O2	5.89	126.02	121.90
1	AA	819	C	N3-C4-N4	5.89	122.12	118.00
1	AA	1858	C	N3-C2-O2	-5.89	117.78	121.90
1	CA	514	A	C5-C6-N1	-5.89	114.76	117.70
1	AA	578	U	N3-C4-C5	5.88	118.13	114.60
1	AA	721	G	N3-C2-N2	5.88	124.02	119.90
1	AA	1856	A	N1-C6-N6	-5.88	115.07	118.60
1	AA	2249	G	N1-C6-O6	-5.88	116.37	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2838	C	C5-C4-N4	-5.88	116.08	120.20
1	AA	1847	G	C4-C5-N7	-5.88	108.45	110.80
1	AA	2466	G	N1-C6-O6	-5.88	116.37	119.90
1	AA	2855	G	N3-C4-N9	5.88	129.53	126.00
34	BA	926	G	OP1-P-OP2	5.88	128.43	119.60
1	AA	1505	C	C4-C5-C6	5.88	120.34	117.40
1	AA	1720	U	OP1-P-OP2	-5.88	110.78	119.60
1	AA	1849	U	C6-N1-C2	5.88	124.53	121.00
1	AA	1922	A	C6-N1-C2	-5.88	115.07	118.60
1	AA	2361	G	C5-N7-C8	5.88	107.24	104.30
1	AA	2366	G	C4-N9-C1'	5.88	134.15	126.50
1	AA	2553	A	C2-N3-C4	-5.88	107.66	110.60
1	AA	2562	G	C4-C5-C6	5.88	122.33	118.80
34	BA	876	G	C5-C6-O6	5.88	132.13	128.60
1	AA	1858	C	N3-C4-N4	-5.88	113.88	118.00
1	AA	2570	C	OP1-P-O3'	5.88	118.14	105.20
1	AA	183	G	OP2-P-O3'	5.88	118.13	105.20
1	AA	343	C	N1-C2-O2	5.88	122.43	118.90
1	AA	786	G	OP2-P-O3'	-5.88	92.27	105.20
1	AA	1156	G	P-O3'-C3'	-5.88	112.65	119.70
1	AA	2658	C	C5-C6-N1	5.88	123.94	121.00
1	AA	2767	U	OP2-P-O3'	5.88	118.13	105.20
1	AA	2776	G	C4-C5-N7	5.88	113.15	110.80
1	CA	324	A	C2-N3-C4	-5.88	107.66	110.60
34	DA	777	A	O5'-P-OP1	5.88	117.75	110.70
34	DA	1483	A	C8-N9-C4	5.88	108.15	105.80
1	AA	355	A	OP1-P-OP2	5.88	128.41	119.60
1	AA	1046	A	C2-N3-C4	-5.88	107.66	110.60
1	AA	1423	G	C6-N1-C2	-5.88	121.57	125.10
1	CA	784	A	O4'-C1'-N9	5.88	112.90	108.20
1	CA	2041	U	C5-C6-N1	-5.88	119.76	122.70
1	CA	2598	A	OP1-P-OP2	-5.88	110.78	119.60
1	AA	989	G	N9-C4-C5	-5.88	103.05	105.40
1	AA	2468	C	N1-C2-O2	5.88	122.42	118.90
34	BA	1431	C	N1-C2-O2	-5.88	115.38	118.90
1	AA	832	G	N3-C2-N2	5.87	124.01	119.90
1	AA	2096	U	N1-C2-N3	5.87	118.42	114.90
34	BA	778	G	C5-C6-O6	5.87	132.12	128.60
1	CA	513	A	N1-C6-N6	5.87	122.12	118.60
1	CA	2784	C	O5'-P-OP2	-5.87	100.41	105.70
34	DA	699	C	C6-N1-C2	-5.87	117.95	120.30
1	AA	1030	A	P-O3'-C3'	5.87	126.75	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1278	G	O5'-P-OP2	5.87	117.75	110.70
1	AA	2372	A	O5'-P-OP2	-5.87	100.42	105.70
1	AA	2822	G	N9-C4-C5	5.87	107.75	105.40
1	CA	1362	C	C5-C6-N1	5.87	123.94	121.00
1	CA	2580	U	C6-N1-C1'	-5.87	112.98	121.20
34	DA	716	A	C8-N9-C4	-5.87	103.45	105.80
1	AA	585	U	N3-C2-O2	-5.87	118.09	122.20
1	AA	1385	G	O5'-P-OP2	5.87	117.74	110.70
1	AA	2660	C	C5-C4-N4	-5.87	116.09	120.20
1	AA	40	C	N3-C2-O2	5.87	126.01	121.90
1	AA	347	G	C8-N9-C4	-5.87	104.05	106.40
1	AA	364	A	N1-C2-N3	-5.87	126.36	129.30
1	AA	813	C	OP2-P-O3'	5.87	118.11	105.20
1	AA	879	G	C4-C5-N7	-5.87	108.45	110.80
1	AA	981	C	N3-C2-O2	5.87	126.01	121.90
1	AA	2112	G	C5-C6-N1	5.87	114.43	111.50
1	AA	2227	G	C8-N9-C1'	5.87	134.63	127.00
1	AA	2520	G	C5-C6-N1	5.87	114.43	111.50
34	DA	590	C	C6-N1-C2	-5.87	117.95	120.30
1	AA	798	A	O5'-P-OP1	-5.87	100.42	105.70
2	AB	98	G	N1-C6-O6	5.87	123.42	119.90
34	BA	1030(B)	C	N3-C2-O2	-5.87	117.79	121.90
1	CA	2044	C	O5'-P-OP2	-5.87	100.42	105.70
1	AA	290	G	N9-C4-C5	-5.87	103.05	105.40
1	AA	2097	U	C5-C6-N1	-5.87	119.77	122.70
1	AA	2236	G	OP1-P-OP2	-5.87	110.80	119.60
1	AA	2715	C	C6-N1-C2	-5.87	117.95	120.30
34	BA	731	G	N3-C2-N2	-5.87	115.80	119.90
1	CA	1367	A	C5-C6-N6	5.87	128.39	123.70
1	AA	369	A	C4-C5-N7	5.86	113.63	110.70
1	AA	829	A	OP1-P-OP2	5.86	128.40	119.60
1	AA	975	U	N3-C4-C5	-5.86	111.08	114.60
1	AA	1249	A	N9-C1'-C2'	5.86	121.62	114.00
1	AA	2416	C	C6-N1-C2	5.86	122.65	120.30
1	CA	1610	A	O5'-P-OP1	-5.86	100.42	105.70
1	AA	957	A	C5-C6-N1	5.86	120.63	117.70
1	AA	1801	G	OP1-P-OP2	5.86	128.39	119.60
1	AA	2053	A	N7-C8-N9	5.86	116.73	113.80
1	AA	2393	C	N1-C2-O2	-5.86	115.38	118.90
1	AA	2593	G	O4'-C1'-N9	5.86	112.89	108.20
2	AB	100	A	C5-N7-C8	-5.86	100.97	103.90
34	BA	20	U	OP2-P-O3'	5.86	118.09	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2035	G	C8-N9-C4	5.86	108.74	106.40
1	AA	805	C	N3-C4-C5	5.86	124.24	121.90
34	BA	523	A	C5-C6-N6	5.86	128.39	123.70
1	CA	676	A	C5-C6-N1	-5.86	114.77	117.70
1	CA	814	C	N1-C2-O2	-5.86	115.39	118.90
1	AA	2552	C	C6-N1-C2	5.86	122.64	120.30
1	AA	2572	C	N1-C2-N3	5.86	123.30	119.20
1	AA	2591	C	N1-C2-O2	-5.86	115.39	118.90
1	CA	702	G	N7-C8-N9	-5.86	110.17	113.10
1	AA	1297	C	C6-N1-C2	5.85	122.64	120.30
1	CA	1604	C	C5-C4-N4	-5.85	116.10	120.20
1	AA	32	C	N3-C4-N4	-5.85	113.90	118.00
1	AA	748	G	C8-N9-C4	5.85	108.74	106.40
1	AA	981	C	C5-C4-N4	-5.85	116.10	120.20
1	AA	1232	G	N3-C2-N2	5.85	124.00	119.90
1	AA	1235	G	C5-N7-C8	5.85	107.23	104.30
1	AA	1296	G	N3-C2-N2	5.85	124.00	119.90
1	AA	2279	A	N1-C2-N3	-5.85	126.37	129.30
1	AA	2335	G	C6-C5-N7	-5.85	126.89	130.40
2	AB	83	G	OP2-P-O3'	5.85	118.07	105.20
1	CA	1898	U	C5-C6-N1	-5.85	119.77	122.70
1	AA	2579	G	C5-C6-O6	5.85	132.11	128.60
34	BA	442	C	C5-C6-N1	5.85	123.93	121.00
1	CA	450	G	C8-N9-C4	-5.85	104.06	106.40
1	CA	1615	C	C6-N1-C2	-5.85	117.96	120.30
1	AA	917	A	N1-C6-N6	-5.85	115.09	118.60
1	AA	969	C	N3-C2-O2	5.85	126.00	121.90
1	AA	1155	C	C2-N3-C4	5.85	122.82	119.90
1	AA	2222	C	C6-N1-C2	5.85	122.64	120.30
1	AA	2303	U	C5-C4-O4	5.85	129.41	125.90
1	AA	2546	A	N1-C6-N6	5.85	122.11	118.60
1	AA	2556	G	N1-C2-N2	5.85	121.46	116.20
1	CA	2500	U	C2-N3-C4	-5.85	123.49	127.00
1	AA	199	C	C2-N1-C1'	-5.85	112.37	118.80
1	AA	778	C	N1-C2-O2	-5.85	115.39	118.90
1	AA	1472	G	N1-C6-O6	5.85	123.41	119.90
1	AA	1807	G	C5-C6-O6	-5.85	125.09	128.60
1	AA	722	A	C5-N7-C8	-5.84	100.98	103.90
1	CA	2229	C	C5-C6-N1	5.84	123.92	121.00
1	CA	2544	G	C4-C5-N7	5.84	113.14	110.80
34	DA	758	G	O5'-P-OP2	-5.84	100.44	105.70
1	AA	542	C	C2-N3-C4	-5.84	116.98	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1985	U	C2-N3-C4	5.84	130.51	127.00
1	AA	190	C	C2-N3-C4	-5.84	116.98	119.90
1	AA	1301	U	C6-N1-C2	-5.84	117.50	121.00
1	AA	1369	U	OP1-P-O3'	5.84	118.05	105.20
1	AA	1823	G	N7-C8-N9	-5.84	110.18	113.10
57	DZ	-20	LEU	CA-CB-CG	5.84	128.74	115.30
1	AA	1404	G	C5-C6-O6	5.84	132.10	128.60
1	AA	2347	A	N1-C6-N6	-5.84	115.10	118.60
1	AA	2749	G	C5-N7-C8	5.84	107.22	104.30
1	AA	245	A	N7-C8-N9	-5.84	110.88	113.80
1	AA	985	G	O5'-P-OP2	-5.84	100.45	105.70
1	AA	2712	C	C6-N1-C2	5.84	122.64	120.30
1	CA	587	C	C5-C6-N1	-5.84	118.08	121.00
1	CA	2497	A	C4-C5-C6	5.84	119.92	117.00
1	AA	122	G	C5-C6-O6	-5.84	125.10	128.60
1	AA	1083	G	O5'-P-OP1	5.84	117.70	110.70
1	AA	1229	G	C4-C5-N7	5.84	113.13	110.80
1	AA	2529	C	N3-C4-C5	5.84	124.23	121.90
34	DA	819	A	N1-C6-N6	5.84	122.10	118.60
1	AA	1539	C	C4-C5-C6	5.83	120.32	117.40
1	AA	2301	G	C4-C5-C6	-5.83	115.30	118.80
1	AA	40	C	C6-N1-C2	5.83	122.63	120.30
1	AA	643	C	N3-C2-O2	5.83	125.98	121.90
1	AA	1712	A	O5'-P-OP1	-5.83	100.45	105.70
1	AA	2073	A	C6-N1-C2	-5.83	115.10	118.60
1	AA	2458	G	C8-N9-C4	-5.83	104.07	106.40
1	AA	2846	U	C5-C4-O4	-5.83	122.40	125.90
34	BA	16	A	C5-N7-C8	5.83	106.82	103.90
34	BA	322	C	C6-N1-C2	5.83	122.63	120.30
34	BA	1397	C	C2-N3-C4	5.83	122.82	119.90
1	CA	1283	G	O5'-P-OP2	-5.83	100.45	105.70
1	CA	1314	C	C6-N1-C1'	-5.83	113.80	120.80
1	AA	372	G	C4-C5-N7	-5.83	108.47	110.80
1	AA	1007	G	C6-N1-C2	-5.83	121.60	125.10
1	AA	2448	G	N9-C4-C5	5.83	107.73	105.40
1	AA	1056	A	C2-N3-C4	5.83	113.52	110.60
1	AA	1249	A	C6-N1-C2	-5.83	115.10	118.60
1	AA	1322	A	N9-C4-C5	-5.83	103.47	105.80
1	AA	1830	G	N1-C6-O6	-5.83	116.40	119.90
1	AA	2261	U	C2-N1-C1'	5.83	124.69	117.70
1	AA	2554	A	N1-C2-N3	-5.83	126.39	129.30
1	CA	1190	G	N1-C6-O6	-5.83	116.40	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2422	A	C8-N9-C4	-5.83	103.47	105.80
1	CA	474	G	N9-C4-C5	5.83	107.73	105.40
1	AA	26	G	N3-C4-N9	5.83	129.50	126.00
1	AA	813	C	C6-N1-C2	5.83	122.63	120.30
1	AA	1652	G	N3-C2-N2	5.83	123.98	119.90
1	AA	2524	C	OP1-P-OP2	-5.83	110.86	119.60
1	AA	2691	A	C8-N9-C4	-5.83	103.47	105.80
1	CA	2385	C	N3-C4-C5	-5.83	119.57	121.90
1	AA	838	C	C6-N1-C2	5.82	122.63	120.30
1	AA	1076	G	N1-C6-O6	5.82	123.39	119.90
1	AA	2047	C	N3-C4-N4	-5.82	113.92	118.00
1	AA	2533	C	C6-N1-C2	5.82	122.63	120.30
1	AA	2720	G	C5-C6-N1	5.82	114.41	111.50
34	BA	902	G	C8-N9-C4	5.82	108.73	106.40
56	BW	45	U	N3-C2-O2	-5.82	118.12	122.20
1	CA	964	C	O5'-P-OP2	-5.82	100.46	105.70
1	CA	2381	C	O5'-P-OP2	-5.82	100.46	105.70
1	CA	2432	A	OP1-P-OP2	-5.82	110.86	119.60
1	AA	1013	G	N3-C2-N2	5.82	123.98	119.90
1	AA	1235	G	C4-C5-N7	-5.82	108.47	110.80
1	AA	2509	A	C6-N1-C2	-5.82	115.11	118.60
1	CA	2360	A	N1-C2-N3	5.82	132.21	129.30
1	CA	2540	C	C2-N3-C4	-5.82	116.99	119.90
1	AA	136	G	C5-C6-N1	5.82	114.41	111.50
1	AA	675	C	C5-C6-N1	-5.82	118.09	121.00
1	AA	745	C	C4-C5-C6	5.82	120.31	117.40
1	AA	1015	C	N3-C4-N4	5.82	122.07	118.00
1	AA	2105	G	N3-C4-N9	-5.82	122.51	126.00
1	AA	98	U	N3-C2-O2	-5.82	118.13	122.20
1	AA	114	C	OP2-P-O3'	5.82	118.00	105.20
1	AA	537	G	OP2-P-O3'	5.82	118.00	105.20
1	AA	1455	C	C6-N1-C2	-5.82	117.97	120.30
1	CA	137	C	C6-N1-C2	-5.82	117.97	120.30
1	CA	848	G	N3-C4-C5	-5.82	125.69	128.60
1	CA	2805	G	C8-N9-C1'	5.82	134.56	127.00
1	AA	356	A	C5-N7-C8	-5.82	100.99	103.90
1	AA	636	G	O5'-P-OP2	-5.82	100.47	105.70
1	AA	2459	G	O4'-C1'-N9	5.82	112.85	108.20
1	CA	740	U	C5-C4-O4	5.82	129.39	125.90
1	CA	2242	G	N3-C4-C5	5.82	131.51	128.60
1	CA	2432	A	C6-N1-C2	-5.82	115.11	118.60
1	AA	438	G	C6-N1-C2	-5.81	121.61	125.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1989	C	C2-N1-C1'	5.81	125.20	118.80
1	AA	2697	G	C5-C6-N1	5.81	114.41	111.50
34	BA	744	C	C2-N1-C1'	-5.81	112.41	118.80
1	AA	824	A	C4-C5-C6	5.81	119.91	117.00
1	AA	1292	A	N9-C4-C5	5.81	108.12	105.80
1	AA	1329	G	N9-C4-C5	-5.81	103.08	105.40
1	AA	2050	U	N3-C2-O2	-5.81	118.13	122.20
1	AA	2681	G	OP2-P-O3'	5.81	117.99	105.20
2	AB	108	U	OP1-P-OP2	5.81	128.32	119.60
34	DA	289	G	N9-C4-C5	-5.81	103.08	105.40
34	DA	352	C	C2-N1-C1'	5.81	125.19	118.80
34	DA	783	C	C6-N1-C2	5.81	122.62	120.30
1	AA	467	U	C5-C6-N1	-5.81	119.80	122.70
1	AA	1454	C	C4-C5-C6	5.81	120.31	117.40
34	BA	134	A	N1-C6-N6	5.81	122.08	118.60
1	CA	788	A	N9-C4-C5	-5.81	103.48	105.80
1	AA	556	C	N3-C4-N4	-5.81	113.94	118.00
1	AA	2867	G	N9-C4-C5	-5.81	103.08	105.40
34	BA	442	C	C6-N1-C2	-5.81	117.98	120.30
1	CA	1790	C	C5-C6-N1	5.81	123.90	121.00
1	AA	1334	U	C4-C5-C6	5.81	123.18	119.70
1	AA	1418	U	C2-N1-C1'	5.81	124.67	117.70
1	AA	2361	G	C4-C5-N7	-5.81	108.48	110.80
1	AA	2368	C	N3-C2-O2	5.81	125.96	121.90
1	AA	2656	G	OP1-P-OP2	-5.81	110.89	119.60
1	CA	908	C	N1-C2-O2	-5.81	115.42	118.90
1	AA	486	A	N9-C4-C5	-5.80	103.48	105.80
1	AA	2520	G	N9-C4-C5	5.80	107.72	105.40
1	CA	446	G	N1-C2-N2	5.80	121.42	116.20
1	CA	1914	C	N1-C2-O2	5.80	122.38	118.90
1	CA	2252	G	C2-N3-C4	-5.80	109.00	111.90
1	AA	1847	G	C4-N9-C1'	-5.80	118.96	126.50
1	AA	2781	C	C2-N3-C4	-5.80	117.00	119.90
34	DA	1466	C	C6-N1-C2	-5.80	117.98	120.30
1	AA	1009	C	N3-C2-O2	-5.80	117.84	121.90
1	AA	1194	A	N7-C8-N9	-5.80	110.90	113.80
1	AA	2471	A	N9-C4-C5	5.80	108.12	105.80
1	AA	2610	A	OP1-P-O3'	-5.80	92.44	105.20
1	AA	2773	C	C5-C6-N1	-5.80	118.10	121.00
34	BA	566	G	N3-C2-N2	-5.80	115.84	119.90
34	BA	890	G	C4-C5-N7	5.80	113.12	110.80
1	CA	2318	G	N7-C8-N9	-5.80	110.20	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1387	U	N1-C2-N3	-5.80	111.42	114.90
1	AA	1518	A	N7-C8-N9	5.80	116.70	113.80
1	AA	1628	G	C8-N9-C4	-5.80	104.08	106.40
1	AA	1694	G	N9-C4-C5	5.80	107.72	105.40
1	AA	2229	A	O4'-C1'-N9	5.80	112.84	108.20
1	AA	2290	A	C8-N9-C4	-5.80	103.48	105.80
1	AA	2297	C	O5'-P-OP1	5.80	117.66	110.70
34	BA	1497	G	N9-C4-C5	-5.80	103.08	105.40
1	AA	1525	G	O5'-P-OP2	-5.80	100.48	105.70
1	AA	2411	G	N1-C2-N3	5.80	127.38	123.90
14	AQ	14	ARG	NE-CZ-NH2	-5.80	117.40	120.30
56	DW	76	A	C6-C5-N7	-5.80	128.24	132.30
1	AA	184	A	OP1-P-O3'	-5.80	92.45	105.20
1	AA	200	A	C6-N1-C2	-5.80	115.12	118.60
1	AA	235	C	OP1-P-O3'	5.80	117.95	105.20
1	AA	2260	C	C4-C5-C6	5.80	120.30	117.40
1	AA	2351	G	N1-C2-N2	-5.80	110.98	116.20
34	BA	854	G	O5'-P-OP2	5.80	117.66	110.70
34	BA	1505	G	N9-C4-C5	5.80	107.72	105.40
1	AA	781	A	N1-C6-N6	5.79	122.08	118.60
1	AA	803	C	C5-C4-N4	-5.79	116.14	120.20
1	AA	129	G	C6-C5-N7	-5.79	126.92	130.40
1	AA	2266	C	O5'-P-OP1	5.79	117.65	110.70
25	A1	41	ARG	NE-CZ-NH2	-5.79	117.40	120.30
1	CA	1313	U	OP1-P-OP2	-5.79	110.91	119.60
1	CA	1760	A	C2-N3-C4	-5.79	107.70	110.60
1	CA	1905	C	C4-C5-C6	-5.79	114.50	117.40
1	CA	1965	C	N3-C4-C5	5.79	124.22	121.90
1	CA	2085	C	C2-N3-C4	-5.79	117.00	119.90
1	AA	420	C	C6-N1-C2	-5.79	117.98	120.30
1	AA	1542	A	OP2-P-O3'	5.79	117.94	105.20
1	AA	1737	A	C8-N9-C4	-5.79	103.48	105.80
34	BA	454	C	N1-C2-O2	5.79	122.37	118.90
34	BA	1510	U	C5-C6-N1	-5.79	119.81	122.70
1	CA	313	C	C6-N1-C2	5.79	122.62	120.30
1	CA	2042	A	N3-C4-N9	-5.79	122.77	127.40
1	AA	213	G	C4-C5-N7	-5.79	108.48	110.80
34	DA	1158	C	C2-N1-C1'	5.79	125.17	118.80
1	AA	1441	A	C5-C6-N6	-5.79	119.07	123.70
1	AA	1844	G	N9-C4-C5	-5.79	103.09	105.40
1	AA	2542	A	C6-C5-N7	-5.79	128.25	132.30
1	CA	51	G	N1-C6-O6	-5.79	116.43	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	194	G	N3-C4-N9	5.79	129.47	126.00
1	CA	2478	A	C8-N9-C4	5.79	108.11	105.80
1	AA	1435	G	OP1-P-OP2	-5.78	110.92	119.60
1	AA	1659	G	N9-C4-C5	-5.78	103.09	105.40
1	AA	1985	U	C2-N1-C1'	5.78	124.64	117.70
1	CA	113	G	N3-C4-C5	5.78	131.49	128.60
1	CA	2386	C	C2-N1-C1'	-5.78	112.44	118.80
1	AA	1305	G	C5-C6-O6	5.78	132.07	128.60
1	AA	2493	G	C6-C5-N7	-5.78	126.93	130.40
1	AA	1092	A	C8-N9-C4	-5.78	103.49	105.80
1	AA	1744	G	N7-C8-N9	-5.78	110.21	113.10
1	AA	2331	G	C4-C5-N7	5.78	113.11	110.80
2	AB	98	G	C5-C6-O6	-5.78	125.13	128.60
34	BA	809	G	N1-C6-O6	-5.78	116.43	119.90
1	CA	2258	C	C2-N3-C4	-5.78	117.01	119.90
34	DA	1414	U	C6-N1-C2	5.78	124.47	121.00
56	DW	73	A	C4-C5-N7	5.78	113.59	110.70
1	AA	2006	G	C4-C5-N7	-5.78	108.49	110.80
1	CA	48	G	C5-C6-N1	-5.78	108.61	111.50
1	CA	2829	C	C2-N3-C4	5.78	122.79	119.90
1	AA	630	U	C2-N1-C1'	-5.78	110.77	117.70
1	AA	761	U	C6-N1-C1'	-5.78	113.11	121.20
1	AA	1409	C	O5'-P-OP2	-5.78	100.50	105.70
1	AA	2608	U	N1-C2-O2	-5.78	118.75	122.80
2	AB	41	U	C2-N1-C1'	-5.78	110.77	117.70
2	AB	106	G	N9-C4-C5	5.78	107.71	105.40
1	CA	1331	A	N7-C8-N9	-5.78	110.91	113.80
1	CA	1759	A	N1-C2-N3	5.78	132.19	129.30
1	CA	1905	C	C5-C6-N1	5.78	123.89	121.00
1	AA	1001	G	C6-C5-N7	-5.78	126.94	130.40
1	AA	2673	G	C5-C6-O6	5.78	132.07	128.60
1	CA	451	C	C6-N1-C2	5.78	122.61	120.30
1	CA	2053	G	N1-C6-O6	5.78	123.36	119.90
1	CA	2358	G	O5'-P-OP2	-5.78	100.50	105.70
34	DA	352	C	C6-N1-C2	-5.78	117.99	120.30
1	AA	1007	G	OP1-P-O3'	5.77	117.90	105.20
1	AA	2526	U	N1-C2-O2	5.77	126.84	122.80
1	CA	124	G	C6-N1-C2	-5.77	121.64	125.10
34	DA	357	G	OP1-P-O3'	5.77	117.90	105.20
1	AA	485	U	C2-N3-C4	-5.77	123.54	127.00
1	AA	2004	C	C5-C6-N1	-5.77	118.11	121.00
1	AA	2289	G	N7-C8-N9	-5.77	110.21	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	174	C	C6-N1-C2	-5.77	117.99	120.30
1	CA	1901	A	N1-C6-N6	-5.77	115.14	118.60
1	CA	2584	U	N3-C4-C5	5.77	118.06	114.60
1	AA	2352	G	C8-N9-C4	-5.77	104.09	106.40
1	AA	2616	U	C2-N3-C4	-5.77	123.54	127.00
1	CA	1810	A	O5'-P-OP2	-5.77	100.51	105.70
1	AA	711	C	N3-C4-C5	5.77	124.21	121.90
1	AA	790	G	N1-C2-N2	-5.77	111.01	116.20
56	BW	74	C	C4-C5-C6	5.77	120.28	117.40
1	CA	2447	G	C4-N9-C1'	-5.77	119.00	126.50
1	CA	2893	G	N3-C4-C5	-5.77	125.72	128.60
1	AA	885	C	N1-C2-O2	5.77	122.36	118.90
1	AA	1049	G	N3-C2-N2	-5.77	115.86	119.90
1	AA	1188	A	C4-N9-C1'	-5.77	115.92	126.30
1	AA	2343	G	N7-C8-N9	-5.77	110.22	113.10
1	AA	2361	G	N7-C8-N9	-5.77	110.22	113.10
34	BA	784	C	C2-N3-C4	-5.77	117.02	119.90
1	CA	1199	U	N1-C2-N3	5.77	118.36	114.90
34	DA	241	C	N3-C4-C5	5.77	124.21	121.90
1	AA	479	C	C4-C5-C6	5.77	120.28	117.40
1	AA	959	U	N3-C4-O4	5.77	123.44	119.40
1	AA	2513	C	OP1-P-OP2	-5.77	110.95	119.60
1	AA	2779	G	N9-C4-C5	5.77	107.71	105.40
1	CA	814	C	N3-C2-O2	5.77	125.94	121.90
1	AA	2229	A	N7-C8-N9	5.76	116.68	113.80
34	BA	1339	A	O5'-P-OP2	-5.76	100.51	105.70
34	DA	656	C	N3-C2-O2	5.76	125.94	121.90
34	DA	898	G	N9-C4-C5	-5.76	103.09	105.40
1	AA	1242	G	C6-C5-N7	5.76	133.86	130.40
34	BA	815	A	C4-C5-N7	-5.76	107.82	110.70
1	CA	787	U	O5'-P-OP1	-5.76	100.51	105.70
1	AA	502	G	O5'-P-OP1	-5.76	100.51	105.70
1	AA	562	C	C5-C6-N1	-5.76	118.12	121.00
1	AA	2846	U	C2-N3-C4	-5.76	123.54	127.00
2	AB	51	G	OP1-P-OP2	-5.76	110.96	119.60
1	CA	527	C	O4'-C1'-N1	5.76	112.81	108.20
1	CA	2287	A	N3-C4-C5	5.76	130.83	126.80
1	AA	991	G	C6-C5-N7	5.76	133.86	130.40
1	AA	1078	A	C5-C6-N6	-5.76	119.09	123.70
1	AA	1211	U	N3-C2-O2	-5.76	118.17	122.20
34	BA	566	G	N9-C4-C5	5.76	107.70	105.40
34	BA	809	G	N9-C4-C5	5.76	107.70	105.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2250	G	OP1-P-OP2	5.76	128.24	119.60
1	CA	2526	G	N1-C6-O6	5.76	123.36	119.90
1	AA	904	C	OP1-P-OP2	5.76	128.24	119.60
1	AA	2621	U	OP2-P-O3'	5.76	117.87	105.20
1	AA	730	C	N3-C2-O2	-5.76	117.87	121.90
1	AA	910	A	N3-C4-C5	-5.76	122.77	126.80
1	AA	1210	G	N9-C4-C5	5.76	107.70	105.40
1	AA	1744	G	C5-C6-N1	5.76	114.38	111.50
1	AA	1977	U	N3-C4-C5	5.76	118.05	114.60
1	AA	2585	C	N3-C2-O2	5.76	125.93	121.90
1	AA	2776	G	C5-C6-N1	-5.76	108.62	111.50
34	DA	687	A	P-O3'-C3'	5.76	126.61	119.70
1	AA	589	U	C6-N1-C2	-5.75	117.55	121.00
1	AA	1078	A	C8-N9-C4	5.75	108.10	105.80
1	AA	1701	A	N7-C8-N9	-5.75	110.92	113.80
1	AA	2782	C	C5-C4-N4	-5.75	116.17	120.20
34	BA	1416	G	N9-C4-C5	5.75	107.70	105.40
1	CA	958	U	C6-N1-C2	-5.75	117.55	121.00
1	AA	1402	G	C8-N9-C4	5.75	108.70	106.40
1	AA	2285	A	C5-N7-C8	-5.75	101.02	103.90
1	AA	2518	U	N1-C2-O2	5.75	126.83	122.80
2	AB	56	G	N3-C4-N9	5.75	129.45	126.00
1	CA	668	G	OP2-P-O3'	5.75	117.85	105.20
1	AA	26	G	C6-N1-C2	-5.75	121.65	125.10
1	AA	150	C	C6-N1-C2	5.75	122.60	120.30
34	BA	1137	C	C6-N1-C2	-5.75	118.00	120.30
1	CA	330	A	C6-C5-N7	-5.75	128.28	132.30
1	AA	12	U	N1-C2-O2	5.75	126.82	122.80
1	AA	2088	C	OP1-P-O3'	5.75	117.85	105.20
1	AA	2471	A	N1-C6-N6	-5.75	115.15	118.60
34	BA	762	C	N1-C2-O2	5.75	122.35	118.90
1	CA	2286	A	C5-C6-N6	-5.75	119.10	123.70
1	AA	991	G	N7-C8-N9	-5.75	110.23	113.10
1	AA	991	G	C8-N9-C4	5.75	108.70	106.40
1	AA	1035	G	N7-C8-N9	-5.75	110.23	113.10
1	AA	2572	C	N1-C2-O2	-5.75	115.45	118.90
1	AA	2629	C	N1-C2-O2	-5.75	115.45	118.90
1	CA	2617	C	OP2-P-O3'	5.75	117.84	105.20
1	AA	803	C	N3-C4-C5	5.75	124.20	121.90
1	AA	1086	C	C6-N1-C2	5.75	122.60	120.30
1	AA	1825	U	OP2-P-O3'	5.75	117.84	105.20
1	AA	2456	G	N7-C8-N9	-5.75	110.23	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2752	U	OP1-P-OP2	-5.74	110.98	119.60
1	CA	272(E)	G	N3-C4-C5	5.74	131.47	128.60
1	CA	1955	U	N1-C2-N3	5.74	118.35	114.90
1	CA	2763	G	N1-C6-O6	-5.74	116.45	119.90
34	DA	890	G	O4'-C1'-N9	5.74	112.79	108.20
1	CA	845	G	C4-N9-C1'	5.74	133.96	126.50
1	CA	1619	G	OP1-P-O3'	5.74	117.83	105.20
1	CA	2698	U	N1-C2-O2	5.74	126.82	122.80
1	AA	181	C	OP1-P-OP2	5.74	128.21	119.60
1	AA	629	U	N3-C2-O2	-5.74	118.18	122.20
1	AA	652	A	N7-C8-N9	-5.74	110.93	113.80
1	AA	2651	A	C2-N3-C4	-5.74	107.73	110.60
2	AB	69	G	N9-C4-C5	-5.74	103.10	105.40
1	CA	1963	U	C5-C6-N1	5.74	125.57	122.70
1	CA	2852	G	O5'-P-OP1	-5.74	100.53	105.70
34	DA	1201	A	P-O3'-C3'	5.74	126.59	119.70
1	AA	210	A	C5'-C4'-O4'	-5.74	102.21	109.10
1	AA	908	A	C5-C6-N1	5.74	120.57	117.70
1	AA	992	G	C2-N3-C4	5.74	114.77	111.90
1	AA	1474	C	C6-N1-C1'	5.74	127.69	120.80
1	AA	2689	G	N9-C4-C5	-5.74	103.11	105.40
34	BA	816	A	O5'-P-OP2	-5.74	100.54	105.70
1	CA	674	G	N1-C2-N3	5.74	127.34	123.90
1	CA	2444	G	C8-N9-C1'	-5.74	119.54	127.00
1	CA	151	C	C5-C6-N1	-5.74	118.13	121.00
1	AA	106	U	OP2-P-O3'	5.74	117.82	105.20
1	AA	1705	C	C6-N1-C2	5.74	122.59	120.30
1	AA	1709	C	N3-C4-N4	-5.74	113.98	118.00
1	AA	2877	G	N3-C4-N9	-5.74	122.56	126.00
34	BA	652	U	C6-N1-C2	5.74	124.44	121.00
1	CA	139(A)	G	C8-N9-C1'	-5.74	119.54	127.00
1	AA	418	G	C6-C5-N7	-5.73	126.96	130.40
1	AA	990	A	C6-N1-C2	-5.73	115.16	118.60
34	DA	912	C	OP2-P-O3'	5.73	117.81	105.20
1	AA	1984	C	N1-C2-O2	5.73	122.34	118.90
1	AA	2556	G	C4-C5-N7	5.73	113.09	110.80
1	AA	2827	G	C6-N1-C2	-5.73	121.66	125.10
1	CA	994	C	N3-C4-C5	-5.73	119.61	121.90
1	CA	1962	C	OP1-P-OP2	-5.73	111.00	119.60
56	DW	74	C	C6-N1-C2	-5.73	118.01	120.30
1	AA	277	G	N3-C4-N9	-5.73	122.56	126.00
1	AA	471	C	C4-C5-C6	5.73	120.27	117.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	587	C	C2-N1-C1'	-5.73	112.50	118.80
1	AA	819	C	C5-C4-N4	-5.73	116.19	120.20
1	AA	956	A	N1-C6-N6	-5.73	115.16	118.60
1	AA	1885	A	C5-C6-N1	5.73	120.56	117.70
34	BA	809	G	C5-C6-O6	5.73	132.04	128.60
1	CA	1646	C	OP1-P-O3'	5.73	117.81	105.20
1	AA	69	G	C8-N9-C4	-5.73	104.11	106.40
1	AA	458	U	C5-C6-N1	-5.73	119.84	122.70
1	AA	1684	A	C2-N3-C4	-5.73	107.74	110.60
1	AA	1865	U	OP2-P-O3'	5.73	117.80	105.20
2	AB	6	C	N1-C2-O2	-5.73	115.46	118.90
1	CA	1944	U	O5'-P-OP2	-5.73	100.55	105.70
1	CA	2038	G	N1-C6-O6	-5.73	116.46	119.90
1	AA	916	G	O5'-P-OP2	-5.73	100.55	105.70
1	CA	1798	U	N3-C4-C5	5.73	118.03	114.60
1	AA	552	C	N1-C2-O2	-5.72	115.47	118.90
1	AA	789	G	C4-C5-N7	-5.72	108.51	110.80
1	AA	1700	G	C5-C6-N1	5.72	114.36	111.50
1	AA	1830	G	OP2-P-O3'	5.72	117.79	105.20
34	BA	1520	G	C5-C6-N1	5.72	114.36	111.50
56	BW	73	A	C4-C5-N7	5.72	113.56	110.70
1	CA	1979	C	C6-N1-C2	-5.72	118.01	120.30
34	DA	7	G	N3-C2-N2	-5.72	115.89	119.90
34	DA	50	A	N9-C4-C5	5.72	108.09	105.80
1	AA	597	C	C6-N1-C1'	5.72	127.67	120.80
1	AA	914	C	C6-N1-C2	5.72	122.59	120.30
1	AA	1274	G	C5-C6-N1	-5.72	108.64	111.50
1	AA	1678	A	C2-N3-C4	-5.72	107.74	110.60
1	AA	2689	G	N3-C2-N2	5.72	123.91	119.90
18	AU	53	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	CA	1227	G	N9-C4-C5	5.72	107.69	105.40
1	AA	1613	A	N1-C6-N6	-5.72	115.17	118.60
1	AA	1902	C	O5'-P-OP2	-5.72	100.55	105.70
1	CA	1992	G	C5-C6-O6	5.72	132.03	128.60
1	CA	2291	U	C5-C6-N1	-5.72	119.84	122.70
1	AA	600	G	OP2-P-O3'	5.72	117.78	105.20
1	AA	1375	U	O5'-P-OP1	5.72	117.56	110.70
1	AA	2022	G	O5'-P-OP2	-5.72	100.55	105.70
1	AA	2643	G	N9-C4-C5	5.72	107.69	105.40
1	AA	2736	C	N3-C4-N4	-5.72	114.00	118.00
2	AB	103	G	C4-C5-N7	5.72	113.09	110.80
1	AA	2544	G	C5-C6-O6	-5.72	125.17	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2563	U	N1-C2-O2	5.72	126.80	122.80
1	AA	290	G	N1-C6-O6	5.72	123.33	119.90
1	AA	2494	G	N3-C2-N2	-5.72	115.90	119.90
13	AP	50	ARG	NE-CZ-NH1	-5.72	117.44	120.30
1	CA	1369	G	N7-C8-N9	5.72	115.96	113.10
1	CA	2438	U	N3-C2-O2	-5.72	118.20	122.20
34	DA	1482	G	C4-N9-C1'	5.72	133.93	126.50
1	AA	363	U	N3-C4-O4	-5.71	115.40	119.40
1	AA	823	G	C8-N9-C4	5.71	108.69	106.40
2	AB	92	C	O5'-P-OP2	5.71	117.56	110.70
34	BA	1087	G	C8-N9-C4	-5.71	104.11	106.40
1	CA	2242	G	C6-N1-C2	5.71	128.53	125.10
34	DA	853	G	N1-C6-O6	5.71	123.33	119.90
1	AA	702	A	C5-C6-N6	5.71	128.27	123.70
1	AA	1211	U	O5'-P-OP2	-5.71	100.56	105.70
1	AA	1811	A	C5'-C4'-O4'	5.71	115.95	109.10
1	AA	468	G	OP1-P-OP2	-5.71	111.03	119.60
1	AA	761	U	C2-N1-C1'	5.71	124.55	117.70
1	AA	1015	C	C5-C4-N4	-5.71	116.20	120.20
1	AA	1457	C	C5-C6-N1	-5.71	118.14	121.00
1	AA	1843	A	C2-N3-C4	5.71	113.46	110.60
1	AA	2558	U	C5-C4-O4	5.71	129.33	125.90
1	AA	2626	A	OP1-P-OP2	-5.71	111.03	119.60
34	BA	546	G	N3-C4-C5	-5.71	125.74	128.60
1	CA	1637	A	N9-C4-C5	5.71	108.08	105.80
1	CA	1837	C	C6-N1-C2	-5.71	118.02	120.30
1	AA	2627	U	OP1-P-OP2	5.71	128.16	119.60
34	DA	879	C	N3-C4-N4	-5.71	114.00	118.00
1	AA	1301	U	OP1-P-OP2	5.71	128.16	119.60
1	AA	2275	C	N3-C4-C5	5.71	124.18	121.90
1	CA	2874	C	C5-C6-N1	5.71	123.85	121.00
34	DA	523	A	N3-C4-N9	-5.71	122.83	127.40
1	AA	589	U	OP1-P-O3'	5.71	117.75	105.20
1	AA	780	G	N3-C4-N9	5.71	129.42	126.00
1	AA	1282	G	O4'-C1'-N9	5.71	112.77	108.20
1	AA	1705	C	N3-C4-C5	5.71	124.18	121.90
34	BA	1402	C	C5-C4-N4	5.71	124.19	120.20
1	CA	51	G	N9-C4-C5	5.71	107.68	105.40
2	CB	47	C	C2-N3-C4	5.71	122.75	119.90
1	AA	2285	A	N1-C2-N3	-5.71	126.45	129.30
1	AA	2488	A	C5-C6-N6	-5.71	119.14	123.70
1	CA	2078	C	OP2-P-O3'	5.71	117.75	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	CB	101	G	C8-N9-C4	5.71	108.68	106.40
1	AA	15	G	N9-C1'-C2'	-5.70	105.72	112.00
1	AA	1000	C	C2-N3-C4	-5.70	117.05	119.90
1	AA	1195	G	C8-N9-C4	-5.70	104.12	106.40
1	AA	1312	G	N3-C2-N2	5.70	123.89	119.90
1	AA	2264	G	O4'-C1'-N9	5.70	112.76	108.20
34	BA	329	A	O4'-C1'-N9	-5.70	103.64	108.20
34	BA	675	A	OP1-P-O3'	5.70	117.75	105.20
34	BA	814	A	C8-N9-C4	5.70	108.08	105.80
34	BA	1436	U	O5'-P-OP1	5.70	117.54	110.70
1	CA	1352	U	O5'-P-OP1	5.70	117.54	110.70
1	AA	1369	U	N1-C2-N3	-5.70	111.48	114.90
1	AA	1853	G	N1-C6-O6	-5.70	116.48	119.90
1	AA	23	G	N3-C2-N2	-5.70	115.91	119.90
1	AA	1931	C	C5-C4-N4	-5.70	116.21	120.20
1	AA	2289	G	C5-N7-C8	5.70	107.15	104.30
1	AA	2892	A	O5'-P-OP2	-5.70	100.57	105.70
14	AQ	22	LYS	CD-CE-NZ	-5.70	98.59	111.70
1	CA	25	U	OP1-P-OP2	5.70	128.15	119.60
1	CA	446	G	C5-C6-O6	-5.70	125.18	128.60
1	CA	528	A	N1-C2-N3	5.70	132.15	129.30
1	AA	551	A	OP1-P-OP2	-5.70	111.05	119.60
1	AA	1201	A	C5-N7-C8	-5.70	101.05	103.90
1	AA	1859	G	C2-N3-C4	-5.70	109.05	111.90
1	AA	2343	G	C4-C5-N7	-5.70	108.52	110.80
1	AA	2632	C	C6-N1-C1'	5.70	127.64	120.80
1	AA	2674	A	OP1-P-O3'	5.70	117.74	105.20
1	AA	2841	G	C5-C6-N1	5.70	114.35	111.50
34	BA	730	G	O5'-P-OP2	-5.70	100.57	105.70
34	BA	1529	G	O4'-C1'-N9	5.70	112.76	108.20
1	CA	2070	G	N1-C2-N2	-5.70	111.07	116.20
1	AA	1263	C	OP1-P-OP2	5.70	128.15	119.60
1	CA	1141	U	C5-C6-N1	-5.70	119.85	122.70
1	CA	1571	A	C8-N9-C4	-5.70	103.52	105.80
1	CA	2572	A	N7-C8-N9	-5.70	110.95	113.80
1	AA	799	A	N1-C2-N3	5.70	132.15	129.30
1	AA	1092	A	N9-C4-C5	5.70	108.08	105.80
1	AA	1689	G	C4-C5-N7	-5.70	108.52	110.80
1	AA	2061	C	O5'-P-OP1	-5.70	100.57	105.70
1	AA	2097	U	OP2-P-O3'	5.70	117.73	105.20
56	BW	76	A	C6-C5-N7	-5.70	128.31	132.30
1	CA	1361	G	N1-C6-O6	-5.70	116.48	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1571	A	N9-C4-C5	5.70	108.08	105.80
1	AA	1431	G	N1-C2-N2	5.69	121.32	116.20
1	AA	2884	C	N1-C2-N3	5.69	123.19	119.20
1	AA	1653	C	C2-N3-C4	5.69	122.75	119.90
1	AA	2088	C	N3-C2-O2	-5.69	117.92	121.90
1	AA	2402	U	N3-C2-O2	-5.69	118.22	122.20
1	AA	59	G	C2-N3-C4	5.69	114.75	111.90
1	AA	1343	C	C4-C5-C6	5.69	120.25	117.40
1	AA	2272	C	C5-C6-N1	-5.69	118.16	121.00
1	CA	948	G	N1-C6-O6	-5.69	116.48	119.90
1	CA	2008	C	O5'-P-OP1	5.69	117.53	110.70
1	CA	2608	G	C8-N9-C4	5.69	108.68	106.40
1	AA	413	G	C5-C6-O6	-5.69	125.19	128.60
1	AA	575	G	C2-N3-C4	5.69	114.74	111.90
1	CA	772	C	C6-N1-C2	5.69	122.58	120.30
1	CA	2444	G	N1-C2-N2	-5.69	111.08	116.20
1	AA	723	A	OP1-P-OP2	5.69	128.13	119.60
1	AA	745	C	N3-C4-C5	-5.69	119.62	121.90
1	AA	1061	G	N1-C2-N3	5.69	127.31	123.90
1	AA	2539	C	OP2-P-O3'	5.69	117.71	105.20
1	CA	2382	G	C5-C6-O6	5.69	132.01	128.60
1	AA	25	U	N3-C4-O4	5.69	123.38	119.40
1	AA	514	G	N3-C2-N2	5.69	123.88	119.90
1	AA	1954	A	OP1-P-OP2	-5.69	111.07	119.60
1	AA	2064	A	N3-C4-C5	5.69	130.78	126.80
1	CA	1657	C	C5-C4-N4	5.69	124.18	120.20
1	CA	2893	G	N3-C4-N9	5.69	129.41	126.00
1	AA	744	C	C6-N1-C2	5.68	122.57	120.30
1	AA	781	A	C5-C6-N6	-5.68	119.15	123.70
1	AA	2771	A	C5-N7-C8	-5.68	101.06	103.90
1	CA	572	A	N9-C4-C5	-5.68	103.53	105.80
34	DA	739	C	N3-C4-C5	-5.68	119.63	121.90
1	AA	413	G	N1-C6-O6	5.68	123.31	119.90
1	CA	330	A	N3-C4-N9	-5.68	122.85	127.40
1	CA	614(B)	G	O4'-C1'-N9	5.68	112.75	108.20
1	AA	2066	C	C5-C4-N4	-5.68	116.22	120.20
1	CA	496	G	N1-C2-N3	5.68	127.31	123.90
1	CA	2250	G	O5'-P-OP2	-5.68	100.59	105.70
1	AA	126	C	OP1-P-OP2	5.68	128.12	119.60
1	AA	198	C	OP2-P-O3'	5.68	117.70	105.20
1	AA	829	A	C5-N7-C8	5.68	106.74	103.90
1	AA	1418	U	C4-C5-C6	5.68	123.11	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1735	U	C5-C4-O4	5.68	129.31	125.90
1	AA	2303	U	C4-C5-C6	5.68	123.11	119.70
1	AA	2369	U	N1-C2-N3	5.68	118.31	114.90
29	A5	16	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	CA	2582	G	C2-N3-C4	-5.68	109.06	111.90
1	AA	1897	C	C2-N1-C1'	5.68	125.05	118.80
1	AA	239	G	N1-C6-O6	5.68	123.31	119.90
1	AA	593	G	C6-C5-N7	-5.68	126.99	130.40
1	AA	1410	G	C5-C6-N1	5.68	114.34	111.50
1	AA	1895	U	N1-C2-O2	-5.68	118.83	122.80
1	AA	2078	G	N1-C2-N2	-5.68	111.09	116.20
34	BA	862	C	O5'-P-OP2	-5.68	100.59	105.70
57	BZ	-29	LEU	CA-CB-CG	5.68	128.36	115.30
1	CA	361	G	N1-C6-O6	5.68	123.31	119.90
1	CA	1288	U	N1-C2-O2	5.68	126.77	122.80
1	CA	2503	A	C5-C6-N1	5.68	120.54	117.70
1	CA	2608	G	O5'-P-OP1	5.68	117.51	110.70
1	AA	622	G	O5'-P-OP2	-5.67	100.59	105.70
1	AA	1346	U	C2-N1-C1'	5.67	124.51	117.70
34	BA	13	U	OP1-P-OP2	5.67	128.11	119.60
34	BA	771	G	C5-C6-N1	-5.67	108.66	111.50
1	CA	446	G	C8-N9-C4	5.67	108.67	106.40
1	CA	1206	G	N3-C2-N2	-5.67	115.93	119.90
1	AA	566	C	N3-C2-O2	-5.67	117.93	121.90
1	AA	733	G	N3-C4-N9	5.67	129.40	126.00
1	AA	1981	G	OP2-P-O3'	5.67	117.68	105.20
1	AA	2404	A	O4'-C1'-N9	-5.67	103.66	108.20
1	AA	192	C	C4-C5-C6	-5.67	114.56	117.40
1	AA	241	G	N7-C8-N9	5.67	115.94	113.10
1	AA	480	A	C5-C6-N1	-5.67	114.86	117.70
1	AA	998	A	C6-C5-N7	-5.67	128.33	132.30
1	AA	1427	G	C5-C6-N1	5.67	114.33	111.50
1	AA	1451	U	O5'-P-OP2	-5.67	100.59	105.70
1	AA	2041	A	N1-C6-N6	-5.67	115.20	118.60
2	AB	77	U	C5-C4-O4	-5.67	122.50	125.90
1	CA	686	G	N1-C6-O6	5.67	123.30	119.90
1	CA	2235	G	C4-C5-N7	-5.67	108.53	110.80
34	DA	835	U	O5'-P-OP1	-5.67	100.59	105.70
1	CA	1814	G	N1-C6-O6	5.67	123.30	119.90
1	CA	1857	G	N1-C6-O6	-5.67	116.50	119.90
1	CA	2504	U	OP1-P-O3'	5.67	117.67	105.20
1	AA	113	C	N1-C2-N3	-5.67	115.23	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	593	G	N3-C4-N9	5.67	129.40	126.00
1	AA	1187	U	O5'-P-OP1	-5.67	100.60	105.70
1	AA	2020	G	N1-C2-N2	-5.67	111.10	116.20
1	AA	2252	C	N1-C2-O2	-5.67	115.50	118.90
1	AA	2452	C	N3-C4-C5	5.67	124.17	121.90
1	AA	2871	G	C4-C5-N7	5.67	113.07	110.80
34	BA	1077	G	OP1-P-O3'	5.67	117.67	105.20
1	CA	2503	A	N1-C6-N6	5.67	122.00	118.60
1	CA	2570	G	N3-C4-N9	-5.67	122.60	126.00
1	CA	2596	U	N3-C4-O4	-5.67	115.43	119.40
34	DA	8	A	C8-N9-C4	5.67	108.07	105.80
34	DA	550	G	C8-N9-C4	5.67	108.67	106.40
1	AA	835	A	N1-C2-N3	-5.67	126.47	129.30
1	AA	1237	G	C8-N9-C4	5.67	108.67	106.40
1	AA	1340	U	N1-C2-O2	-5.67	118.83	122.80
1	AA	2597	U	N1-C2-N3	-5.67	111.50	114.90
1	CA	1626	G	N1-C6-O6	5.67	123.30	119.90
1	AA	1684	A	C6-N1-C2	5.67	122.00	118.60
1	CA	1558	A	N1-C2-N3	5.67	132.13	129.30
1	AA	820	U	C5-C4-O4	-5.66	122.50	125.90
1	AA	824	A	OP1-P-OP2	5.66	128.09	119.60
1	AA	1238	G	C2-N3-C4	5.66	114.73	111.90
1	AA	1977	U	C6-N1-C2	5.66	124.40	121.00
1	AA	2001	C	C2-N3-C4	-5.66	117.07	119.90
1	AA	2311	G	C8-N9-C4	-5.66	104.13	106.40
1	AA	2639	G	C4-C5-C6	-5.66	115.40	118.80
1	CA	804	A	O5'-P-OP2	5.66	117.50	110.70
1	AA	2651	A	OP1-P-O3'	5.66	117.66	105.20
1	CA	1142(A)	A	N1-C6-N6	5.66	122.00	118.60
1	AA	724	A	C4-C5-N7	-5.66	107.87	110.70
1	AA	887	C	N3-C4-C5	5.66	124.16	121.90
1	AA	1205	U	C2-N3-C4	-5.66	123.60	127.00
1	AA	2090	U	O5'-P-OP2	-5.66	100.61	105.70
1	CA	588	U	O5'-P-OP2	-5.66	100.61	105.70
1	CA	2590	A	N1-C6-N6	-5.66	115.20	118.60
34	DA	370	C	N3-C2-O2	-5.66	117.94	121.90
1	AA	2870	A	N1-C6-N6	-5.66	115.20	118.60
34	BA	402	G	C8-N9-C4	-5.66	104.14	106.40
1	CA	94(A)	G	N1-C6-O6	5.66	123.30	119.90
1	CA	1755	A	O5'-P-OP1	-5.66	100.61	105.70
1	AA	2238	C	C5-C6-N1	-5.66	118.17	121.00
1	AA	2633	A	N9-C4-C5	5.66	108.06	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
56	BW	75	C	C5-C6-N1	-5.66	118.17	121.00
1	CA	777	A	N7-C8-N9	5.66	116.63	113.80
1	CA	1975	G	C5-C6-O6	-5.66	125.21	128.60
1	AA	1390	G	C8-N9-C4	5.66	108.66	106.40
1	CA	2067	G	C5-C6-O6	5.66	131.99	128.60
1	AA	533	G	C4-C5-N7	-5.65	108.54	110.80
1	AA	1063	G	N9-C4-C5	5.65	107.66	105.40
1	AA	1394	G	C8-N9-C4	5.65	108.66	106.40
1	AA	2864	G	N9-C1'-C2'	-5.65	105.78	112.00
1	AA	714	U	N1-C2-N3	5.65	118.29	114.90
1	AA	1097	G	OP2-P-O3'	5.65	117.64	105.20
1	AA	1950	A	C8-N9-C4	5.65	108.06	105.80
1	AA	2004	C	N3-C4-C5	-5.65	119.64	121.90
1	AA	2014	G	C5-C6-N1	5.65	114.33	111.50
1	CA	777	A	C8-N9-C4	-5.65	103.54	105.80
1	CA	1624	G	N1-C6-O6	-5.65	116.51	119.90
1	AA	2463	A	C2-N3-C4	5.65	113.42	110.60
2	AB	103	G	C5-C6-N1	-5.65	108.67	111.50
56	BW	76	A	C5-N7-C8	-5.65	101.08	103.90
1	CA	1680	U	O5'-P-OP1	-5.65	100.61	105.70
1	CA	1814	G	C4-C5-N7	5.65	113.06	110.80
1	CA	2694	G	N7-C8-N9	5.65	115.92	113.10
1	AA	1845	G	C5-C6-O6	-5.65	125.21	128.60
1	AA	2515	A	N3-C4-N9	5.65	131.92	127.40
1	CA	945	A	N7-C8-N9	5.65	116.62	113.80
1	CA	2050	C	C5-C4-N4	-5.65	116.25	120.20
1	AA	625	G	N1-C2-N2	-5.65	111.12	116.20
1	AA	1042	A	C6-N1-C2	5.65	121.99	118.60
34	BA	896	C	C6-N1-C2	5.65	122.56	120.30
1	CA	1533	G	C4-N9-C1'	5.65	133.84	126.50
1	AA	1610	G	O5'-P-OP1	-5.65	100.62	105.70
1	AA	847	A	C2-N3-C4	5.64	113.42	110.60
1	AA	2020	G	C2-N3-C4	-5.64	109.08	111.90
1	AA	2074	G	N9-C4-C5	5.64	107.66	105.40
1	AA	2220	A	O5'-P-OP1	-5.64	100.62	105.70
1	AA	2450	U	C6-N1-C2	5.64	124.39	121.00
1	AA	210	A	P-O5'-C5'	-5.64	111.87	120.90
1	AA	1883	C	N3-C2-O2	-5.64	117.95	121.90
34	BA	890	G	N9-C4-C5	-5.64	103.14	105.40
1	CA	2875	C	C5-C6-N1	-5.64	118.18	121.00
34	DA	867	G	C8-N9-C4	-5.64	104.14	106.40
1	AA	1797	U	O5'-P-OP1	5.64	117.47	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1814	A	C5-C6-N1	-5.64	114.88	117.70
1	AA	2108	U	OP1-P-O3'	-5.64	92.79	105.20
1	CA	1673	U	C2-N1-C1'	-5.64	110.93	117.70
1	AA	1070	G	C6-N1-C2	-5.64	121.72	125.10
1	AA	1387	U	C2-N1-C1'	-5.64	110.93	117.70
1	AA	2511	C	C6-N1-C2	-5.64	118.04	120.30
1	AA	2554	A	N3-C4-C5	5.64	130.75	126.80
1	CA	2061	G	C5-C6-O6	-5.64	125.22	128.60
1	AA	202	A	OP2-P-O3'	5.64	117.61	105.20
1	AA	2039	U	N1-C2-N3	5.64	118.28	114.90
1	AA	2772	G	N3-C2-N2	-5.64	115.95	119.90
2	AB	92	C	C5-C4-N4	-5.64	116.25	120.20
34	BA	809	G	C4-C5-N7	-5.64	108.55	110.80
1	CA	1798	U	C5-C6-N1	-5.64	119.88	122.70
1	AA	574	G	OP2-P-O3'	5.64	117.60	105.20
1	AA	589	U	C5-C6-N1	5.64	125.52	122.70
1	AA	1427	G	N1-C2-N3	5.64	127.28	123.90
1	AA	1563	G	N3-C2-N2	-5.64	115.95	119.90
1	AA	2441	G	OP1-P-OP2	-5.64	111.14	119.60
2	AB	6	C	C6-N1-C2	5.64	122.55	120.30
34	BA	560	U	O5'-P-OP2	-5.64	100.63	105.70
1	CA	1379	A	N1-C6-N6	5.64	121.98	118.60
1	CA	2456	C	OP2-P-O3'	5.64	117.60	105.20
1	CA	2741	A	C8-N9-C4	5.64	108.06	105.80
1	AA	355	A	C5-C6-N1	5.63	120.52	117.70
1	AA	384	G	N9-C4-C5	-5.63	103.15	105.40
1	AA	1384	G	C5-C6-O6	5.63	131.98	128.60
1	AA	1518	A	C2-N3-C4	-5.63	107.78	110.60
1	AA	1609	A	C8-N9-C4	5.63	108.05	105.80
1	AA	2894	U	C2-N1-C1'	-5.63	110.94	117.70
2	AB	104	U	C2-N3-C4	-5.63	123.62	127.00
34	BA	1354	C	C6-N1-C2	-5.63	118.05	120.30
1	CA	2089	U	N1-C2-N3	5.63	118.28	114.90
34	DA	32	A	C2-N3-C4	5.63	113.42	110.60
1	AA	373	G	OP2-P-O3'	5.63	117.59	105.20
1	AA	1006	C	C5-C6-N1	-5.63	118.18	121.00
1	AA	1702	A	C4-C5-C6	-5.63	114.18	117.00
1	AA	2331	G	C8-N9-C1'	5.63	134.32	127.00
1	CA	450	G	C6-N1-C2	-5.63	121.72	125.10
1	CA	1339	G	C5-C6-O6	5.63	131.98	128.60
1	CA	2438	U	OP2-P-O3'	5.63	117.59	105.20
1	CA	2767	C	N3-C4-N4	5.63	121.94	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	171	A	C5-N7-C8	-5.63	101.08	103.90
1	AA	510	C	C5-C6-N1	5.63	123.82	121.00
1	AA	579	G	C5-C6-O6	5.63	131.98	128.60
1	AA	711	C	OP1-P-OP2	5.63	128.05	119.60
1	AA	807	G	O5'-P-OP2	5.63	117.46	110.70
1	AA	2366	G	O4'-C1'-N9	-5.63	103.69	108.20
1	CA	127	A	O5'-P-OP2	-5.63	100.63	105.70
1	CA	2503	A	N3-C4-N9	5.63	131.91	127.40
1	AA	878	G	C5-N7-C8	-5.63	101.48	104.30
1	AA	2272	C	C4-C5-C6	5.63	120.22	117.40
1	AA	2802	C	O4'-C1'-N1	5.63	112.70	108.20
1	AA	891	C	C6-N1-C2	5.63	122.55	120.30
1	AA	1849	U	C5-C4-O4	-5.63	122.52	125.90
1	AA	1859	G	N1-C2-N3	5.63	127.28	123.90
1	CA	1035	U	N1-C2-O2	5.63	126.74	122.80
1	CA	1047	G	N3-C4-C5	-5.63	125.78	128.60
1	CA	1646	C	C4-C5-C6	5.63	120.21	117.40
1	AA	900	G	C8-N9-C4	-5.63	104.15	106.40
1	AA	984	G	N9-C4-C5	-5.63	103.15	105.40
1	AA	1021	G	N3-C2-N2	5.63	123.84	119.90
1	CA	39	C	C5-C6-N1	-5.63	118.19	121.00
1	CA	50	U	N1-C2-N3	-5.63	111.52	114.90
2	AB	115	G	OP1-P-OP2	5.62	128.04	119.60
34	BA	574	A	C4-C5-N7	5.62	113.51	110.70
1	AA	244	A	C2-N3-C4	-5.62	107.79	110.60
34	DA	21	G	C4-C5-N7	5.62	113.05	110.80
34	DA	1515	C	C6-N1-C2	5.62	122.55	120.30
1	AA	28	A	N1-C6-N6	-5.62	115.23	118.60
1	AA	848	G	OP1-P-O3'	5.62	117.57	105.20
1	AA	1809	U	N3-C2-O2	5.62	126.14	122.20
1	CA	674	G	C2-N3-C4	-5.62	109.09	111.90
1	CA	1350	C	C2-N3-C4	-5.62	117.09	119.90
1	CA	1768	U	N3-C4-O4	-5.62	115.46	119.40
1	CA	2194	G	O5'-P-OP1	5.62	117.45	110.70
34	DA	1496	C	N1-C2-O2	-5.62	115.53	118.90
1	AA	129	G	N9-C4-C5	-5.62	103.15	105.40
1	AA	231	G	N1-C6-O6	-5.62	116.53	119.90
1	AA	1315	A	C6-N1-C2	5.62	121.97	118.60
1	AA	137	G	C4-N9-C1'	5.62	133.80	126.50
1	AA	777	C	C6-N1-C2	-5.62	118.05	120.30
1	AA	1003	U	C5-C6-N1	5.62	125.51	122.70
1	AA	1657	C	C4-C5-C6	5.62	120.21	117.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AB	116	G	N1-C2-N2	-5.62	111.14	116.20
34	BA	284	G	O5'-P-OP1	5.62	117.44	110.70
34	BA	766	A	C2-N3-C4	5.62	113.41	110.60
2	AB	109	C	O4'-C1'-N1	5.62	112.69	108.20
1	AA	2501	G	N3-C2-N2	-5.62	115.97	119.90
1	CA	1798	U	N3-C2-O2	-5.62	118.27	122.20
1	CA	2062	A	OP2-P-O3'	5.62	117.56	105.20
1	CA	2444	G	N1-C6-O6	-5.62	116.53	119.90
1	AA	223	C	N1-C2-O2	5.61	122.27	118.90
1	AA	400	U	C2-N3-C4	-5.61	123.63	127.00
1	AA	593	G	N3-C4-C5	-5.61	125.79	128.60
1	AA	708	C	N3-C4-C5	5.61	124.14	121.90
1	AA	1646	C	C6-N1-C2	5.61	122.55	120.30
1	AA	1752	G	N3-C2-N2	5.61	123.83	119.90
1	CA	643	A	C8-N9-C4	5.61	108.05	105.80
1	CA	2581	G	O5'-P-OP2	-5.61	100.65	105.70
1	AA	983	G	OP1-P-OP2	-5.61	111.18	119.60
1	AA	1830	G	O5'-P-OP1	5.61	117.44	110.70
1	AA	2655	G	N9-C4-C5	5.61	107.64	105.40
1	AA	2673	G	OP1-P-O3'	5.61	117.55	105.20
15	AR	113	LEU	CA-CB-CG	5.61	128.21	115.30
56	BW	45	U	C2-N1-C1'	5.61	124.44	117.70
1	AA	2005	C	C6-N1-C2	5.61	122.54	120.30
1	AA	2041	A	C4-C5-N7	-5.61	107.89	110.70
1	AA	2556	G	C6-N1-C2	-5.61	121.73	125.10
2	AB	38	C	OP2-P-O3'	5.61	117.54	105.20
34	BA	1442(A)	G	C5-C6-O6	-5.61	125.23	128.60
1	AA	1385	G	O5'-P-OP1	-5.61	100.65	105.70
1	AA	1594	C	N3-C2-O2	-5.61	117.97	121.90
1	AA	2048	C	OP1-P-O3'	-5.61	92.86	105.20
1	CA	224	G	O5'-P-OP2	-5.61	100.65	105.70
4	AD	258	LYS	CD-CE-NZ	-5.61	98.81	111.70
1	CA	52	A	N7-C8-N9	5.61	116.60	113.80
1	CA	1154	G	C8-N9-C1'	-5.61	119.71	127.00
1	CA	2793	G	C8-N9-C4	-5.61	104.16	106.40
1	AA	2041	A	N7-C8-N9	-5.60	111.00	113.80
1	CA	1896	G	C4-N9-C1'	-5.60	119.22	126.50
1	AA	41	C	N3-C4-N4	-5.60	114.08	118.00
1	AA	555	G	OP2-P-O3'	5.60	117.53	105.20
1	AA	1019	G	OP1-P-O3'	5.60	117.53	105.20
34	BA	188	C	C6-N1-C2	-5.60	118.06	120.30
1	CA	2067	G	C8-N9-C4	-5.60	104.16	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2870	C	C6-N1-C2	-5.60	118.06	120.30
1	AA	311	C	O5'-P-OP2	-5.60	100.66	105.70
1	AA	587	C	C6-N1-C1'	5.60	127.52	120.80
1	AA	1333	A	N1-C6-N6	-5.60	115.24	118.60
1	AA	2550	C	N1-C2-O2	-5.60	115.54	118.90
34	BA	296	U	C4-C5-C6	5.60	123.06	119.70
1	CA	679	C	OP1-P-OP2	-5.60	111.20	119.60
1	AA	490	U	C5-C6-N1	-5.60	119.90	122.70
1	AA	793	A	N9-C4-C5	5.60	108.04	105.80
1	AA	865	G	C6-C5-N7	5.60	133.76	130.40
1	AA	963	A	O4'-C1'-N9	-5.60	103.72	108.20
1	AA	1354	A	OP2-P-O3'	5.60	117.52	105.20
1	AA	1450	C	OP1-P-OP2	5.60	128.00	119.60
1	AA	2291	G	C5-N7-C8	5.60	107.10	104.30
13	AP	21	ARG	NE-CZ-NH1	-5.60	117.50	120.30
1	CA	1260	G	N7-C8-N9	-5.60	110.30	113.10
1	AA	202	A	O5'-P-OP2	-5.60	100.66	105.70
1	AA	1429	C	N3-C4-C5	5.60	124.14	121.90
1	AA	2045	G	C6-C5-N7	-5.60	127.04	130.40
1	AA	2464	C	N1-C2-N3	5.60	123.12	119.20
2	AB	41	U	N3-C4-O4	-5.60	115.48	119.40
1	AA	747	G	N7-C8-N9	-5.60	110.30	113.10
1	AA	1640	G	N1-C6-O6	5.60	123.26	119.90
1	AA	2632	C	C2-N3-C4	-5.60	117.10	119.90
1	CA	939	G	N1-C6-O6	5.60	123.26	119.90
1	CA	1657	C	N3-C4-N4	-5.60	114.08	118.00
1	AA	234	G	N3-C4-C5	-5.59	125.80	128.60
1	AA	437	G	C8-N9-C4	5.59	108.64	106.40
1	AA	2311	G	N9-C4-C5	5.59	107.64	105.40
1	CA	34	C	C2-N1-C1'	5.59	124.95	118.80
1	CA	318	C	N3-C4-N4	5.59	121.92	118.00
1	CA	480	A	N1-C6-N6	-5.59	115.24	118.60
1	CA	2373	G	C4-C5-C6	5.59	122.16	118.80
1	AA	1289	G	N1-C2-N2	-5.59	111.17	116.20
1	AA	2586	G	N1-C2-N3	-5.59	120.54	123.90
1	CA	516	C	N1-C2-O2	-5.59	115.54	118.90
1	CA	1647	G	C5-C6-N1	5.59	114.30	111.50
1	AA	630	U	C5-C6-N1	-5.59	119.90	122.70
1	AA	2532	C	N1-C2-N3	5.59	123.11	119.20
13	AP	55	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	CA	125	G	C5-N7-C8	-5.59	101.50	104.30
1	CA	576	U	O5'-P-OP2	-5.59	100.67	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	DA	372	C	C5-C4-N4	-5.59	116.28	120.20
34	DA	1149	C	C6-N1-C2	-5.59	118.06	120.30
1	AA	139	A	N1-C2-N3	5.59	132.09	129.30
1	AA	706	C	O5'-P-OP2	-5.59	100.67	105.70
1	AA	1457	C	C2-N3-C4	-5.59	117.11	119.90
1	AA	1632	A	C4-C5-C6	5.59	119.80	117.00
34	BA	254	G	O5'-P-OP1	-5.59	100.67	105.70
1	CA	1644	C	C2-N1-C1'	5.59	124.95	118.80
34	DA	874	G	N3-C4-N9	5.59	129.35	126.00
34	DA	1064	G	P-O3'-C3'	5.59	126.41	119.70
1	CA	1338	G	O5'-P-OP2	-5.59	100.67	105.70
34	DA	256	U	C5-C6-N1	-5.59	119.91	122.70
1	AA	889	G	N9-C4-C5	5.59	107.64	105.40
1	AA	1486	G	C5-N7-C8	5.59	107.09	104.30
1	AA	2645	G	N7-C8-N9	-5.59	110.31	113.10
1	AA	1533	G	C8-N9-C4	-5.58	104.17	106.40
1	AA	1974	A	C5-C6-N1	-5.58	114.91	117.70
34	BA	266	G	N7-C8-N9	5.58	115.89	113.10
15	CR	64	ARG	NE-CZ-NH2	-5.58	117.51	120.30
1	AA	1377	A	N9-C4-C5	5.58	108.03	105.80
1	AA	2019	G	C5'-C4'-O4'	5.58	115.80	109.10
2	AB	55	U	N1-C2-N3	5.58	118.25	114.90
1	AA	353	G	N3-C4-N9	5.58	129.35	126.00
1	AA	413	G	N3-C4-N9	5.58	129.35	126.00
1	AA	423	G	N1-C2-N2	5.58	121.22	116.20
1	AA	496	A	N7-C8-N9	5.58	116.59	113.80
1	AA	1821	C	C6-N1-C2	5.58	122.53	120.30
1	AA	2075	G	C5-C6-O6	5.58	131.95	128.60
1	AA	2436	C	OP1-P-OP2	5.58	127.97	119.60
34	BA	134	A	N9-C4-C5	-5.58	103.57	105.80
1	AA	962	G	N1-C2-N3	5.58	127.25	123.90
1	AA	2241	C	N3-C4-C5	5.58	124.13	121.90
1	CA	214	G	O4'-C1'-N9	5.58	112.66	108.20
1	CA	530	G	C8-N9-C4	-5.58	104.17	106.40
1	AA	889	G	OP2-P-O3'	5.58	117.47	105.20
1	CA	424	G	N1-C2-N2	5.58	121.22	116.20
1	CA	1309	G	C5-C6-O6	-5.58	125.25	128.60
1	AA	97	G	C5-C6-O6	5.58	131.95	128.60
1	AA	543	G	OP1-P-OP2	5.58	127.97	119.60
1	AA	1039	G	OP2-P-O3'	5.58	117.47	105.20
1	AA	1274	G	C8-N9-C4	5.58	108.63	106.40
1	AA	1482	G	C8-N9-C4	-5.58	104.17	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2075	G	O5'-P-OP1	-5.58	100.68	105.70
34	BA	741	G	O5'-P-OP2	-5.58	100.68	105.70
1	CA	948	G	N7-C8-N9	-5.58	110.31	113.10
1	CA	2596	U	C5-C6-N1	-5.58	119.91	122.70
34	DA	853	G	C8-N9-C4	-5.58	104.17	106.40
1	AA	340	C	N3-C4-N4	-5.57	114.10	118.00
1	AA	721	G	OP2-P-O3'	5.57	117.46	105.20
1	AA	895	G	N9-C4-C5	5.57	107.63	105.40
1	AA	1006	C	N3-C4-C5	5.57	124.13	121.90
1	AA	1653	C	OP1-P-O3'	5.57	117.46	105.20
1	AA	2039	U	N3-C2-O2	-5.57	118.30	122.20
1	CA	38	A	C8-N9-C4	-5.57	103.57	105.80
1	AA	1204	C	C4-C5-C6	5.57	120.19	117.40
1	AA	2613	C	C5-C4-N4	-5.57	116.30	120.20
1	CA	1318	C	N3-C4-C5	5.57	124.13	121.90
1	AA	554	A	N7-C8-N9	5.57	116.59	113.80
1	AA	731	G	C5-C6-N1	5.57	114.29	111.50
1	AA	864	C	OP1-P-OP2	-5.57	111.25	119.60
1	AA	922	G	N9-C4-C5	-5.57	103.17	105.40
1	AA	1239	A	C6-N1-C2	-5.57	115.26	118.60
1	AA	1702	A	N1-C6-N6	-5.57	115.26	118.60
1	AA	2410	U	C5-C4-O4	5.57	129.24	125.90
34	BA	312	C	OP2-P-O3'	5.57	117.46	105.20
34	BA	1523	G	C8-N9-C4	-5.57	104.17	106.40
1	CA	2415	G	N7-C8-N9	5.57	115.89	113.10
1	AA	845	G	N1-C6-O6	-5.57	116.56	119.90
1	AA	1041	C	OP1-P-O3'	5.57	117.45	105.20
1	AA	1684	A	C5-C6-N1	-5.57	114.92	117.70
1	CA	465	G	N3-C4-C5	-5.57	125.82	128.60
1	CA	1387	C	C6-N1-C2	-5.57	118.07	120.30
1	AA	226	C	C5-C6-N1	-5.57	118.22	121.00
1	AA	535	C	C5-C6-N1	-5.57	118.22	121.00
1	AA	780	G	N3-C4-C5	-5.57	125.82	128.60
1	AA	1690	G	C5-N7-C8	5.57	107.08	104.30
1	AA	1975	A	C4-C5-C6	-5.57	114.22	117.00
1	AA	2623	U	C2-N3-C4	-5.57	123.66	127.00
1	CA	1310	G	O4'-C1'-N9	5.57	112.65	108.20
1	AA	53	G	C4-C5-N7	-5.57	108.57	110.80
1	AA	593	G	C6-N1-C2	-5.57	121.76	125.10
1	AA	1689	G	N3-C4-C5	-5.57	125.82	128.60
1	AA	2561	G	C4-C5-N7	-5.57	108.57	110.80
2	AB	25	A	C8-N9-C4	5.57	108.03	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	809	G	C2-N3-C4	5.57	114.68	111.90
1	CA	2611	U	C2-N3-C4	-5.57	123.66	127.00
34	DA	718	G	N3-C2-N2	-5.57	116.00	119.90
1	AA	27	G	N1-C6-O6	-5.56	116.56	119.90
1	AA	345	G	C4-C5-N7	5.56	113.03	110.80
1	AA	1474	C	C2-N3-C4	-5.56	117.12	119.90
34	BA	236	G	OP1-P-O3'	5.56	117.44	105.20
1	CA	123	G	C6-N1-C2	-5.56	121.76	125.10
1	CA	387	U	N3-C2-O2	-5.56	118.31	122.20
1	CA	1219	G	N1-C6-O6	-5.56	116.56	119.90
1	AA	1922	A	N3-C4-C5	-5.56	122.91	126.80
1	AA	2454	C	OP1-P-OP2	-5.56	111.26	119.60
1	AA	2630	G	N7-C8-N9	5.56	115.88	113.10
1	AA	2645	G	C5-N7-C8	5.56	107.08	104.30
34	BA	897	C	OP1-P-O3'	5.56	117.44	105.20
1	CA	151	C	C6-N1-C2	5.56	122.53	120.30
1	AA	19	C	N3-C4-C5	5.56	124.12	121.90
1	AA	418	G	C2-N3-C4	-5.56	109.12	111.90
1	AA	883	G	N3-C4-N9	5.56	129.34	126.00
1	AA	2502	G	N1-C6-O6	-5.56	116.56	119.90
1	CA	1217	C	C6-N1-C2	-5.56	118.08	120.30
1	CA	1351	C	C5-C6-N1	-5.56	118.22	121.00
1	AA	337	C	C6-N1-C2	5.56	122.52	120.30
1	AA	1085	G	N7-C8-N9	-5.56	110.32	113.10
1	AA	1829	U	C2-N3-C4	-5.56	123.66	127.00
1	CA	1672	C	C2-N3-C4	-5.56	117.12	119.90
56	DW	40	C	P-O3'-C3'	5.56	126.37	119.70
1	AA	22	C	N3-C2-O2	-5.56	118.01	121.90
34	BA	804	U	N3-C4-C5	5.56	117.94	114.60
1	CA	1835	G	C6-C5-N7	-5.56	127.06	130.40
1	CA	2879	C	C4-C5-C6	5.56	120.18	117.40
34	DA	923	A	N7-C8-N9	5.56	116.58	113.80
1	AA	949	C	N1-C2-O2	-5.56	115.57	118.90
1	AA	1360	C	C4-C5-C6	5.56	120.18	117.40
1	CA	1788	C	N3-C4-C5	-5.56	119.68	121.90
1	CA	2010	G	O5'-P-OP1	-5.56	100.70	105.70
34	DA	289	G	C4-C5-N7	5.56	113.02	110.80
1	AA	426	G	N1-C2-N2	-5.55	111.20	116.20
1	AA	675	C	C2-N3-C4	-5.55	117.12	119.90
1	AA	710	G	N1-C2-N2	5.55	121.20	116.20
1	AA	757	G	C5-C6-O6	-5.55	125.27	128.60
1	AA	777	C	N1-C2-N3	5.55	123.09	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1270	C	N3-C4-C5	5.55	124.12	121.90
1	AA	1372	U	N3-C2-O2	-5.55	118.31	122.20
1	AA	1926	G	N9-C4-C5	5.55	107.62	105.40
34	BA	1523	G	C4-C5-N7	-5.55	108.58	110.80
1	CA	2069	G	N3-C2-N2	-5.55	116.01	119.90
1	AA	176	G	C2-N3-C4	-5.55	109.12	111.90
1	AA	317	U	OP1-P-OP2	5.55	127.93	119.60
1	AA	886	U	C4-C5-C6	-5.55	116.37	119.70
1	AA	1299	A	C5-N7-C8	-5.55	101.12	103.90
1	AA	2417	G	C6-C5-N7	-5.55	127.07	130.40
1	AA	501	U	N1-C2-N3	5.55	118.23	114.90
1	AA	2042	A	OP1-P-OP2	5.55	127.93	119.60
34	BA	558	G	N3-C4-N9	5.55	129.33	126.00
34	BA	733	A	O5'-P-OP1	-5.55	100.70	105.70
1	CA	265	A	C4-C5-N7	5.55	113.48	110.70
1	CA	1439	A	O5'-P-OP2	-5.55	100.70	105.70
1	CA	2849	U	OP2-P-O3'	5.55	117.41	105.20
34	DA	877	C	C6-N1-C2	5.55	122.52	120.30
1	AA	2290	A	N1-C6-N6	-5.55	115.27	118.60
1	AA	2497	G	N3-C4-N9	5.55	129.33	126.00
1	AA	2620	G	N1-C6-O6	5.55	123.23	119.90
1	AA	2645	G	C6-N1-C2	-5.55	121.77	125.10
1	CA	2374	C	C6-N1-C2	5.55	122.52	120.30
1	CA	2571	C	C4-C5-C6	5.55	120.17	117.40
1	AA	480	A	N9-C4-C5	5.55	108.02	105.80
57	BZ	87	HIS	N-CA-C	-5.55	96.02	111.00
1	CA	2315	G	N9-C4-C5	-5.55	103.18	105.40
1	AA	1462	G	OP1-P-OP2	5.55	127.92	119.60
1	AA	2114	U	N1-C2-N3	5.55	118.23	114.90
1	AA	2775	G	N1-C2-N2	-5.55	111.21	116.20
34	BA	283	C	N3-C2-O2	-5.55	118.02	121.90
56	BW	34	G	N1-C6-O6	5.55	123.23	119.90
1	AA	532	A	C8-N9-C4	-5.54	103.58	105.80
34	BA	1527	C	N1-C2-N3	5.54	123.08	119.20
1	CA	204	A	N9-C4-C5	5.54	108.02	105.80
1	CA	1941	C	N3-C4-C5	-5.54	119.68	121.90
1	AA	38	A	C6-C5-N7	5.54	136.18	132.30
1	AA	352	U	OP1-P-O3'	5.54	117.39	105.20
1	AA	743	G	N1-C6-O6	-5.54	116.57	119.90
1	AA	1248	G	N1-C2-N3	5.54	127.23	123.90
1	AA	2016	C	O5'-P-OP2	-5.54	100.71	105.70
1	AA	2479	C	C5-C6-N1	-5.54	118.23	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2598	C	OP1-P-OP2	-5.54	111.28	119.60
1	AA	2616	U	N3-C2-O2	-5.54	118.32	122.20
1	AA	2857	U	C2-N3-C4	-5.54	123.67	127.00
34	BA	1509	C	N3-C4-C5	5.54	124.12	121.90
1	AA	702	A	C5-C6-N1	-5.54	114.93	117.70
1	AA	1543	U	C6-N1-C1'	5.54	128.96	121.20
1	AA	2268	G	OP1-P-OP2	-5.54	111.29	119.60
1	AA	2279	A	C8-N9-C1'	-5.54	117.72	127.70
1	CA	2061	G	O5'-P-OP2	-5.54	100.71	105.70
1	AA	101	A	C5-C6-N6	-5.54	119.27	123.70
1	AA	999	G	OP1-P-OP2	-5.54	111.29	119.60
1	AA	2709	G	N3-C2-N2	5.54	123.78	119.90
1	CA	799	G	N3-C4-C5	-5.54	125.83	128.60
1	AA	1284	G	OP2-P-O3'	5.54	117.39	105.20
1	AA	1412	A	N1-C6-N6	-5.54	115.28	118.60
1	AA	1670	G	N3-C2-N2	5.54	123.78	119.90
1	AA	2081	A	N9-C4-C5	-5.54	103.58	105.80
1	AA	2703	C	N3-C4-N4	-5.54	114.12	118.00
34	BA	528	C	N3-C4-C5	5.54	124.11	121.90
1	CA	575	A	C2-N3-C4	5.54	113.37	110.60
1	CA	1609	A	C2-N3-C4	-5.54	107.83	110.60
1	CA	2388	A	O4'-C1'-N9	5.54	112.63	108.20
1	CA	2727	G	N1-C6-O6	-5.54	116.58	119.90
34	DA	705	U	C5-C6-N1	5.54	125.47	122.70
1	AA	1295	U	C5-C4-O4	-5.54	122.58	125.90
1	AA	2458	G	N1-C6-O6	-5.54	116.58	119.90
1	CA	2234	G	C5-N7-C8	-5.54	101.53	104.30
1	CA	2618	G	N7-C8-N9	-5.54	110.33	113.10
1	AA	69	G	O5'-P-OP1	5.54	117.34	110.70
1	AA	918	U	N3-C4-O4	5.54	123.27	119.40
1	AA	975	U	OP1-P-O3'	5.54	117.38	105.20
1	AA	1187	U	N3-C4-O4	5.54	123.28	119.40
1	AA	1661	C	N3-C4-N4	-5.54	114.12	118.00
1	AA	2088	C	N3-C4-N4	-5.54	114.13	118.00
1	AA	2832	G	C5-C6-O6	-5.54	125.28	128.60
34	BA	596	C	N1-C2-O2	5.54	122.22	118.90
1	CA	2452	C	OP2-P-O3'	5.54	117.38	105.20
1	AA	315	C	N3-C4-C5	5.53	124.11	121.90
1	AA	464	G	C6-N1-C2	-5.53	121.78	125.10
1	AA	1032	C	C2-N3-C4	-5.53	117.13	119.90
1	AA	1966	U	C2-N3-C4	-5.53	123.68	127.00
13	AP	61	ARG	NE-CZ-NH2	-5.53	117.53	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
21	AX	57	LEU	CA-CB-CG	5.53	128.03	115.30
1	CA	592	G	N3-C4-N9	5.53	129.32	126.00
1	CA	2057	A	N7-C8-N9	-5.53	111.03	113.80
1	CA	2558	C	C5-C4-N4	-5.53	116.33	120.20
34	DA	563	A	O4'-C1'-N9	5.53	112.63	108.20
1	AA	1690	G	N7-C8-N9	-5.53	110.33	113.10
1	AA	2647	C	O5'-P-OP1	5.53	117.34	110.70
1	AA	124	A	N7-C8-N9	-5.53	111.03	113.80
1	AA	2741	U	C4-C5-C6	5.53	123.02	119.70
1	CA	1428	C	C5-C4-N4	5.53	124.07	120.20
1	CA	1821	A	N1-C6-N6	-5.53	115.28	118.60
1	CA	2691	C	OP1-P-OP2	5.53	127.90	119.60
34	BA	1465	C	OP1-P-OP2	5.53	127.89	119.60
56	DW	15	G	C5-C6-O6	5.53	131.92	128.60
1	AA	345	G	C8-N9-C4	5.53	108.61	106.40
1	AA	515	G	N1-C2-N3	5.53	127.22	123.90
1	AA	590	A	C5-C6-N6	-5.53	119.28	123.70
1	AA	1078	A	C4-C5-N7	5.53	113.46	110.70
1	AA	1603	C	N3-C4-C5	-5.53	119.69	121.90
1	AA	2341	G	C5-N7-C8	5.53	107.06	104.30
1	CA	1377	G	N3-C4-N9	5.53	129.32	126.00
1	CA	1427	A	P-O3'-C3'	5.53	126.33	119.70
1	CA	2049	G	N3-C2-N2	5.53	123.77	119.90
1	AA	309	C	C2-N3-C4	-5.53	117.14	119.90
1	AA	1013	G	C2-N3-C4	5.53	114.66	111.90
1	AA	1738	C	N3-C4-N4	-5.53	114.13	118.00
34	BA	1471	G	O5'-P-OP2	-5.53	100.73	105.70
34	DA	873	A	N1-C6-N6	5.53	121.92	118.60
1	AA	16	G	C2-N3-C4	-5.52	109.14	111.90
1	AA	2883	A	N1-C2-N3	-5.52	126.54	129.30
1	CA	248	G	C8-N9-C4	-5.52	104.19	106.40
1	AA	1413	A	OP1-P-OP2	5.52	127.88	119.60
1	AA	2264	G	C2-N3-C4	-5.52	109.14	111.90
1	AA	2342	G	C4-C5-N7	5.52	113.01	110.80
1	AA	2676	G	C4-C5-N7	5.52	113.01	110.80
1	AA	2782	C	N1-C2-O2	-5.52	115.59	118.90
1	CA	34	C	C2-N3-C4	5.52	122.66	119.90
1	CA	384	U	C6-N1-C2	5.52	124.31	121.00
1	CA	1653	G	P-O3'-C3'	5.52	126.33	119.70
1	AA	465	G	C6-C5-N7	-5.52	127.09	130.40
1	AA	982	U	C6-N1-C2	5.52	124.31	121.00
1	AA	2785	C	C5-C6-N1	-5.52	118.24	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1303	C	O5'-P-OP2	5.52	117.32	110.70
1	AA	1693	C	N1-C2-O2	-5.52	115.59	118.90
1	CA	141	A	O4'-C1'-N9	5.52	112.62	108.20
1	CA	1441	G	N9-C4-C5	5.52	107.61	105.40
1	CA	2559	C	O5'-P-OP1	-5.52	100.73	105.70
1	CA	2689	U	C6-N1-C2	-5.52	117.69	121.00
1	CA	2895	U	C5-C6-N1	5.52	125.46	122.70
34	DA	36	C	C6-N1-C2	-5.52	118.09	120.30
1	AA	761	U	N1-C2-N3	-5.52	111.59	114.90
1	AA	1429	C	C6-N1-C2	5.52	122.51	120.30
1	AA	1668	G	OP2-P-O3'	5.52	117.34	105.20
2	AB	94	C	OP2-P-O3'	5.52	117.34	105.20
1	CA	220	G	N1-C6-O6	5.52	123.21	119.90
1	CA	1959	G	N3-C4-N9	-5.52	122.69	126.00
1	AA	57	G	C8-N9-C4	5.52	108.61	106.40
34	DA	266	G	N7-C8-N9	5.52	115.86	113.10
1	AA	1261	G	C6-C5-N7	-5.51	127.09	130.40
1	AA	2033	U	N1-C2-N3	5.51	118.21	114.90
1	AA	2774	G	OP2-P-O3'	5.51	117.33	105.20
1	AA	2803	A	N7-C8-N9	5.51	116.56	113.80
1	CA	491	G	N3-C4-N9	-5.51	122.69	126.00
1	CA	743	G	N3-C4-C5	-5.51	125.84	128.60
1	CA	2073	C	C2-N3-C4	-5.51	117.14	119.90
56	DW	17	C	C5-C6-N1	5.51	123.76	121.00
1	AA	28	A	C4-C5-C6	-5.51	114.24	117.00
1	AA	518	G	C6-C5-N7	-5.51	127.09	130.40
1	AA	2454	C	O5'-P-OP1	5.51	117.31	110.70
1	CA	2070	G	C5-N7-C8	5.51	107.06	104.30
1	AA	565	C	N1-C2-O2	-5.51	115.59	118.90
1	AA	734	C	N1-C2-N3	5.51	123.06	119.20
1	AA	907	U	OP1-P-O3'	5.51	117.32	105.20
1	AA	1234	A	C6-N1-C2	-5.51	115.29	118.60
1	AA	2839	C	OP2-P-O3'	5.51	117.32	105.20
34	BA	28	G	N3-C2-N2	-5.51	116.04	119.90
1	CA	535	C	N1-C2-O2	-5.51	115.59	118.90
1	AA	964	A	C2-N3-C4	-5.51	107.84	110.60
1	AA	1065	U	N1-C2-O2	-5.51	118.94	122.80
1	AA	1341	C	N3-C4-N4	-5.51	114.14	118.00
1	AA	1360	C	N3-C2-O2	-5.51	118.04	121.90
1	AA	1950	A	N1-C2-N3	-5.51	126.55	129.30
1	AA	2028	C	C2-N3-C4	-5.51	117.14	119.90
1	AA	2742	G	C6-N1-C2	5.51	128.41	125.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	172	C	N3-C4-N4	-5.51	114.14	118.00
1	AA	904	C	N1-C2-N3	5.51	123.06	119.20
1	AA	2362	C	C5-C6-N1	-5.51	118.25	121.00
1	CA	59	U	N1-C2-N3	5.51	118.20	114.90
1	CA	1846	G	N9-C4-C5	5.51	107.60	105.40
1	AA	508	A	N1-C2-N3	5.51	132.05	129.30
1	AA	2091	G	O5'-P-OP1	-5.51	100.74	105.70
1	AA	2240	G	N1-C6-O6	-5.51	116.60	119.90
56	BW	17	C	N3-C2-O2	-5.51	118.05	121.90
1	CA	416	C	OP1-P-OP2	-5.51	111.34	119.60
1	CA	2544	G	C6-C5-N7	-5.51	127.10	130.40
1	AA	1509	C	C2-N1-C1'	5.50	124.86	118.80
1	CA	25	U	N1-C2-O2	-5.50	118.95	122.80
1	AA	1398	U	OP1-P-OP2	-5.50	111.35	119.60
1	AA	1784	G	OP1-P-O3'	5.50	117.31	105.20
1	AA	2637	G	N3-C2-N2	-5.50	116.05	119.90
1	CA	659	C	C2-N1-C1'	-5.50	112.75	118.80
1	CA	1190	G	O5'-P-OP2	-5.50	100.75	105.70
1	CA	1597	A	O4'-C1'-N9	5.50	112.60	108.20
1	CA	1687	G	OP2-P-O3'	5.50	117.31	105.20
34	DA	353	A	N9-C4-C5	-5.50	103.60	105.80
1	AA	437	G	C5-C6-N1	-5.50	108.75	111.50
1	AA	790	G	C5-C6-O6	5.50	131.90	128.60
1	CA	571	A	OP2-P-O3'	5.50	117.30	105.20
1	AA	1256	U	OP1-P-OP2	-5.50	111.35	119.60
1	AA	2025	G	C5-C6-N1	5.50	114.25	111.50
1	AA	2065	C	C5-C6-N1	5.50	123.75	121.00
1	CA	2509	G	C2-N3-C4	-5.50	109.15	111.90
1	AA	702	A	N9-C4-C5	5.50	108.00	105.80
1	AA	2083	G	N7-C8-N9	5.50	115.85	113.10
1	AA	2399	U	C2-N3-C4	-5.50	123.70	127.00
1	AA	2401	G	OP1-P-O3'	5.50	117.30	105.20
1	AA	414	U	C5-C4-O4	-5.50	122.60	125.90
1	AA	1272	A	O5'-P-OP2	-5.50	100.75	105.70
1	AA	2243	C	C5-C6-N1	-5.50	118.25	121.00
2	AB	49	C	N3-C2-O2	5.50	125.75	121.90
2	AB	82	G	C5-N7-C8	5.50	107.05	104.30
1	AA	241	G	N3-C4-C5	-5.50	125.85	128.60
1	AA	2249	G	C8-N9-C4	5.50	108.60	106.40
1	CA	1812	A	N1-C6-N6	-5.50	115.30	118.60
1	CA	2850	A	O5'-P-OP1	5.50	117.29	110.70
1	AA	101	A	N1-C6-N6	5.49	121.90	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1548	C	C5-C4-N4	-5.49	116.36	120.20
1	AA	2108	U	C6-N1-C2	5.49	124.30	121.00
1	AA	2227	G	N3-C4-N9	-5.49	122.70	126.00
1	AA	2513	C	C5-C4-N4	5.49	124.05	120.20
1	AA	2791	A	OP2-P-O3'	5.49	117.28	105.20
1	CA	914	C	N3-C2-O2	-5.49	118.05	121.90
1	CA	1692	U	N1-C2-O2	-5.49	118.95	122.80
1	CA	2439	A	C5'-C4'-O4'	-5.49	102.51	109.10
1	AA	724	A	C4-C5-C6	5.49	119.75	117.00
1	AA	865	G	N1-C6-O6	-5.49	116.61	119.90
1	AA	1040	C	N1-C2-O2	-5.49	115.61	118.90
34	BA	291	C	C2-N1-C1'	-5.49	112.76	118.80
1	CA	1309	G	N1-C6-O6	5.49	123.19	119.90
1	CA	2552	U	N3-C4-O4	-5.49	115.56	119.40
34	DA	906	G	C4-C5-N7	5.49	113.00	110.80
1	AA	22	C	OP1-P-O3'	-5.49	93.12	105.20
1	AA	137	G	C6-C5-N7	-5.49	127.11	130.40
1	AA	2504	U	N3-C2-O2	-5.49	118.36	122.20
1	AA	2514	G	C5-N7-C8	5.49	107.05	104.30
1	AA	2639	G	N7-C8-N9	-5.49	110.36	113.10
1	AA	2803	A	C8-N9-C4	-5.49	103.60	105.80
34	BA	17	U	C5-C4-O4	5.49	129.19	125.90
34	BA	916	G	N3-C4-C5	-5.49	125.85	128.60
1	CA	1240	U	N1-C2-O2	-5.49	118.96	122.80
1	AA	385	G	C5-C6-O6	-5.49	125.31	128.60
1	AA	953	U	C5-C4-O4	-5.49	122.61	125.90
1	AA	1317	G	N1-C2-N2	-5.49	111.26	116.20
1	AA	1559	C	O5'-P-OP2	5.49	117.28	110.70
1	AA	1719	C	N3-C2-O2	-5.49	118.06	121.90
1	AA	2030	C	OP1-P-O3'	5.49	117.28	105.20
6	AF	89	VAL	C-N-CA	-5.49	107.98	121.70
34	BA	769	G	OP2-P-O3'	5.49	117.28	105.20
1	CA	128	C	O5'-P-OP2	-5.49	100.76	105.70
1	CA	2582	G	N1-C2-N3	5.49	127.19	123.90
1	CA	378	C	N1-C2-O2	5.49	122.19	118.90
34	DA	882	C	C4-C5-C6	5.49	120.14	117.40
1	AA	451	G	N3-C4-C5	5.49	131.34	128.60
1	AA	890	G	OP2-P-O3'	5.49	117.27	105.20
1	AA	1705	C	C5-C6-N1	-5.49	118.26	121.00
1	AA	1977	U	C2-N3-C4	-5.49	123.71	127.00
1	AA	1988	A	C2-N3-C4	-5.49	107.86	110.60
1	CA	2051	A	C6-N1-C2	-5.49	115.31	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	DA	619	U	C6-N1-C1'	5.48	128.88	121.20
1	AA	200	A	C2-N3-C4	-5.48	107.86	110.60
1	AA	998	A	C5-N7-C8	-5.48	101.16	103.90
1	AA	1931	C	N3-C4-N4	5.48	121.84	118.00
1	AA	2222	C	C5-C6-N1	-5.48	118.26	121.00
1	AA	2903	G	C5-N7-C8	5.48	107.04	104.30
34	BA	427	U	C6-N1-C2	-5.48	117.71	121.00
34	DA	353	A	C4-C5-N7	5.48	113.44	110.70
1	AA	98	U	O4'-C1'-N1	5.48	112.58	108.20
1	AA	496	A	N1-C6-N6	-5.48	115.31	118.60
1	AA	554	A	C8-N9-C4	-5.48	103.61	105.80
1	AA	2690	C	C6-N1-C2	-5.48	118.11	120.30
1	AA	2724	U	O4'-C1'-N1	5.48	112.58	108.20
34	BA	1413	A	OP1-P-O3'	5.48	117.26	105.20
1	CA	254	G	N3-C4-N9	-5.48	122.71	126.00
1	CA	948	G	C8-N9-C4	5.48	108.59	106.40
34	DA	50	A	OP1-P-OP2	5.48	127.82	119.60
1	AA	23	G	N7-C8-N9	-5.48	110.36	113.10
1	AA	1474	C	C2-N1-C1'	-5.48	112.77	118.80
1	AA	2554	A	C4-C5-C6	-5.48	114.26	117.00
1	CA	567	A	C5-C6-N1	5.48	120.44	117.70
1	CA	1408	C	C6-N1-C2	-5.48	118.11	120.30
1	AA	321	C	N3-C4-C5	-5.48	119.71	121.90
1	AA	566	C	N1-C2-O2	5.48	122.19	118.90
1	AA	798	A	N1-C2-N3	5.48	132.04	129.30
1	AA	1188	A	N7-C8-N9	5.48	116.54	113.80
1	AA	1819	C	C2-N3-C4	-5.48	117.16	119.90
1	CA	2607	G	C6-C5-N7	-5.48	127.11	130.40
1	CA	2719	G	C5-C6-O6	5.48	131.89	128.60
1	CA	2827	C	O5'-P-OP2	-5.48	100.77	105.70
1	AA	1312	G	C4-N9-C1'	-5.48	119.38	126.50
1	AA	1375	U	C2-N3-C4	-5.48	123.71	127.00
5	AE	47	VAL	CB-CA-C	-5.48	101.00	111.40
34	BA	801	U	N3-C4-C5	5.48	117.89	114.60
56	BW	11	C	O5'-P-OP2	-5.48	100.77	105.70
1	CA	1565	C	N3-C2-O2	5.48	125.73	121.90
1	AA	311	C	O5'-P-OP1	5.47	117.27	110.70
1	AA	547	G	OP2-P-O3'	5.47	117.24	105.20
1	AA	862	C	N3-C2-O2	-5.47	118.07	121.90
1	AA	2415	C	C6-N1-C2	-5.47	118.11	120.30
1	AA	2837	C	C6-N1-C2	5.47	122.49	120.30
27	A3	31	LEU	CB-CG-CD2	-5.47	101.69	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
56	BW	73	A	N9-C4-C5	-5.47	103.61	105.80
1	CA	527	C	N3-C4-N4	-5.47	114.17	118.00
1	CA	580	C	C5-C6-N1	-5.47	118.26	121.00
1	CA	1614	A	OP1-P-O3'	5.47	117.24	105.20
1	CA	2043	C	OP1-P-OP2	5.47	127.81	119.60
1	CA	2348	U	C5-C4-O4	5.47	129.19	125.90
34	DA	566	G	N1-C6-O6	-5.47	116.61	119.90
34	DA	1230	C	C6-N1-C2	-5.47	118.11	120.30
1	AA	867	A	C8-N9-C4	5.47	107.99	105.80
1	AA	868	A	C2-N3-C4	5.47	113.34	110.60
1	AA	1953	U	C6-N1-C2	-5.47	117.72	121.00
1	AA	2422	G	N7-C8-N9	-5.47	110.36	113.10
34	BA	901	A	O5'-P-OP2	5.47	117.27	110.70
1	CA	942	G	C5-C6-N1	5.47	114.24	111.50
1	AA	73	A	C6-N1-C2	-5.47	115.32	118.60
1	AA	1397	C	O5'-P-OP1	5.47	117.27	110.70
5	AE	144	ARG	NE-CZ-NH1	5.47	123.04	120.30
34	BA	1518	A	N1-C6-N6	-5.47	115.32	118.60
1	CA	37	C	C4-C5-C6	5.47	120.14	117.40
1	AA	31	C	N3-C2-O2	-5.47	118.07	121.90
1	AA	724	A	C5-C6-N6	5.47	128.08	123.70
1	AA	1610	G	OP2-P-O3'	5.47	117.23	105.20
1	AA	2620	G	N9-C1'-C2'	-5.47	105.98	112.00
1	AA	2650	G	N3-C4-N9	5.47	129.28	126.00
1	CA	1786	A	O4'-C1'-N9	5.47	112.58	108.20
1	CA	2439	A	O5'-P-OP1	5.47	117.26	110.70
1	CA	2841	C	O5'-P-OP2	-5.47	100.78	105.70
1	AA	757	G	OP2-P-O3'	5.47	117.23	105.20
1	AA	1851	U	C4-C5-C6	5.47	122.98	119.70
1	AA	2297	C	OP2-P-O3'	5.47	117.23	105.20
1	AA	2834	C	N1-C2-O2	-5.47	115.62	118.90
34	BA	339	C	N1-C2-O2	-5.47	115.62	118.90
1	CA	518	G	N3-C4-N9	-5.47	122.72	126.00
1	CA	1694	C	C6-N1-C2	5.47	122.49	120.30
1	CA	2456	C	OP1-P-O3'	-5.47	93.17	105.20
1	CA	2487	G	OP1-P-OP2	5.47	127.80	119.60
1	AA	790	G	N3-C2-N2	5.47	123.73	119.90
1	AA	894	U	N1-C2-O2	-5.47	118.97	122.80
1	AA	2718	G	C5-C6-O6	5.47	131.88	128.60
1	AA	2801	C	N1-C2-N3	5.47	123.03	119.20
1	CA	2444	G	C5-N7-C8	5.47	107.03	104.30
1	AA	580	U	C2-N3-C4	-5.46	123.72	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	725	C	C5-C4-N4	5.46	124.03	120.20
1	AA	1052	C	C6-N1-C1'	5.46	127.36	120.80
1	AA	1274	G	N1-C2-N3	5.46	127.18	123.90
1	AA	2404	A	N9-C4-C5	-5.46	103.61	105.80
1	AA	2609	G	N7-C8-N9	-5.46	110.37	113.10
1	AA	2886	G	N1-C6-O6	-5.46	116.62	119.90
1	CA	856	C	P-O3'-C3'	5.46	126.26	119.70
1	CA	2771	C	N1-C2-O2	-5.46	115.62	118.90
1	AA	1935	A	O4'-C1'-N9	-5.46	103.83	108.20
34	BA	731	G	N9-C4-C5	5.46	107.58	105.40
56	DW	56	C	C6-N1-C2	-5.46	118.11	120.30
1	AA	1057	G	N3-C4-N9	5.46	129.28	126.00
1	AA	2015	U	N3-C2-O2	-5.46	118.38	122.20
1	CA	702	G	C4-C5-N7	-5.46	108.62	110.80
1	AA	978	A	C5'-C4'-O4'	5.46	115.65	109.10
1	CA	2680	C	O5'-P-OP2	-5.46	100.79	105.70
1	AA	235	C	C4-C5-C6	5.46	120.13	117.40
1	AA	847	A	C6-N1-C2	-5.46	115.32	118.60
1	AA	1327	G	C8-N9-C4	-5.46	104.22	106.40
1	AA	2404	A	C2-N3-C4	-5.46	107.87	110.60
1	AA	2563	C	N1-C2-O2	-5.46	115.62	118.90
34	BA	46	G	N1-C6-O6	5.46	123.17	119.90
1	CA	1781	C	C6-N1-C1'	-5.46	114.25	120.80
34	DA	893	C	OP1-P-OP2	5.46	127.79	119.60
1	AA	400	U	N3-C4-C5	5.46	117.87	114.60
1	AA	768	C	N1-C2-O2	-5.46	115.63	118.90
1	AA	1444	C	C6-N1-C2	-5.46	118.12	120.30
2	AB	100	A	N3-C4-N9	-5.46	123.03	127.40
1	CA	766	C	C5-C6-N1	-5.46	118.27	121.00
1	AA	1735	U	C5-C6-N1	-5.46	119.97	122.70
1	CA	1656	C	N3-C2-O2	-5.46	118.08	121.90
1	CA	2232	U	C5-C4-O4	5.46	129.17	125.90
34	DA	729	A	N7-C8-N9	5.46	116.53	113.80
5	AE	182	LEU	CA-CB-CG	5.45	127.84	115.30
56	BW	1	G	N3-C4-C5	-5.45	125.87	128.60
1	CA	1372	U	N3-C4-O4	5.45	123.22	119.40
1	AA	992	G	N3-C4-C5	-5.45	125.87	128.60
1	AA	1230	C	C5-C6-N1	-5.45	118.28	121.00
34	BA	111	G	N1-C6-O6	5.45	123.17	119.90
1	CA	939	G	C5-C6-O6	-5.45	125.33	128.60
1	CA	1821	A	C5-N7-C8	5.45	106.62	103.90
1	CA	2234	G	C6-N1-C2	-5.45	121.83	125.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2420	C	N3-C4-C5	5.45	124.08	121.90
1	CA	2678	C	N1-C2-O2	-5.45	115.63	118.90
1	AA	972	A	C2-N3-C4	5.45	113.32	110.60
1	AA	1346	U	N3-C2-O2	-5.45	118.39	122.20
1	AA	1423	G	C8-N9-C4	-5.45	104.22	106.40
1	AA	2400	A	N3-C4-N9	-5.45	123.04	127.40
1	AA	2405	A	C2-N3-C4	-5.45	107.88	110.60
1	AA	2721	G	O5'-P-OP2	5.45	117.24	110.70
1	CA	787	U	O5'-P-OP2	5.45	117.24	110.70
1	CA	1323	U	C5-C6-N1	-5.45	119.98	122.70
1	CA	1823	G	C4-C5-N7	-5.45	108.62	110.80
1	AA	880	U	N3-C4-O4	5.45	123.21	119.40
1	AA	1387	U	N3-C2-O2	5.45	126.01	122.20
1	CA	324	A	C8-N9-C4	5.45	107.98	105.80
1	AA	724	A	C5-N7-C8	5.45	106.62	103.90
1	AA	1302	G	N7-C8-N9	-5.45	110.38	113.10
1	AA	2057	G	OP1-P-OP2	5.45	127.77	119.60
1	AA	624	C	C4-C5-C6	5.44	120.12	117.40
1	AA	2742	G	C8-N9-C4	5.44	108.58	106.40
34	BA	733	A	N1-C6-N6	5.44	121.87	118.60
1	AA	894	U	N3-C2-O2	-5.44	118.39	122.20
1	AA	2030	C	C6-N1-C2	-5.44	118.12	120.30
1	AA	2389	A	C8-N9-C4	5.44	107.98	105.80
1	CA	141	A	N1-C2-N3	5.44	132.02	129.30
1	CA	1032	A	C8-N9-C4	5.44	107.98	105.80
34	DA	572	A	N7-C8-N9	-5.44	111.08	113.80
34	DA	755	G	N1-C6-O6	5.44	123.17	119.90
34	DA	1473	A	OP2-P-O3'	5.44	117.17	105.20
1	AA	2558	U	N3-C4-O4	-5.44	115.59	119.40
1	CA	109	G	C6-C5-N7	5.44	133.66	130.40
1	CA	665	C	N3-C4-N4	5.44	121.81	118.00
1	CA	2571	C	N1-C2-N3	5.44	123.01	119.20
34	DA	97	G	O4'-C1'-N9	5.44	112.55	108.20
1	AA	1210	G	C4-C5-N7	-5.44	108.62	110.80
1	AA	1608	G	O5'-P-OP2	5.44	117.23	110.70
1	AA	1980	C	N3-C4-C5	-5.44	119.72	121.90
1	CA	1695	G	C5-C6-N1	-5.44	108.78	111.50
1	AA	518	G	N3-C4-N9	5.44	129.26	126.00
1	AA	1608	G	N3-C2-N2	-5.44	116.09	119.90
1	AA	1707	C	O5'-P-OP1	5.44	117.22	110.70
14	AQ	56	ARG	NE-CZ-NH1	5.44	123.02	120.30
34	BA	28	G	N1-C6-O6	5.44	123.16	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	CB	77	U	C6-N1-C2	-5.44	117.74	121.00
1	AA	2617	U	N3-C4-O4	-5.44	115.59	119.40
1	AA	2799	U	C5-C6-N1	-5.44	119.98	122.70
1	CA	2565	A	O5'-P-OP2	5.44	117.22	110.70
1	AA	122	G	N1-C6-O6	5.43	123.16	119.90
1	AA	518	G	O5'-P-OP1	5.43	117.22	110.70
1	AA	908	A	OP1-P-OP2	-5.43	111.45	119.60
1	AA	996	C	N1-C2-O2	5.43	122.16	118.90
1	AA	1597	C	N3-C4-N4	5.43	121.81	118.00
1	AA	1741	C	N1-C2-O2	-5.43	115.64	118.90
1	AA	2522	C	C4-C5-C6	5.43	120.12	117.40
1	AA	2550	C	N3-C4-C5	5.43	124.07	121.90
1	AA	2641	A	C4-C5-C6	5.43	119.72	117.00
1	CA	119	A	C2-N3-C4	-5.43	107.88	110.60
1	CA	2383	G	N1-C2-N2	-5.43	111.31	116.20
34	DA	283	C	N3-C2-O2	-5.43	118.09	121.90
34	DA	1158	C	N3-C2-O2	-5.43	118.10	121.90
1	AA	27	G	C4-C5-N7	-5.43	108.63	110.80
1	AA	357	G	N9-C4-C5	5.43	107.57	105.40
1	AA	1282	G	O5'-P-OP2	5.43	117.22	110.70
1	AA	2008	A	O5'-P-OP1	-5.43	100.81	105.70
1	AA	2635	G	C4-N9-C1'	5.43	133.56	126.50
34	BA	653	A	N1-C6-N6	5.43	121.86	118.60
1	CA	799	G	C2-N3-C4	5.43	114.62	111.90
1	CA	2023	G	C8-N9-C4	-5.43	104.23	106.40
34	DA	50	A	N7-C8-N9	5.43	116.52	113.80
1	AA	28	A	N1-C2-N3	-5.43	126.58	129.30
1	AA	38	A	C6-N1-C2	-5.43	115.34	118.60
1	AA	728	G	C5-N7-C8	5.43	107.02	104.30
1	AA	1237	G	C6-C5-N7	5.43	133.66	130.40
1	AA	2289	G	N1-C2-N3	5.43	127.16	123.90
1	CA	2244	U	O5'-P-OP2	-5.43	100.81	105.70
2	CB	47	C	N1-C2-N3	-5.43	115.40	119.20
1	AA	141	C	OP2-P-O3'	5.43	117.14	105.20
1	AA	353	G	C5-C6-N1	5.43	114.22	111.50
1	AA	546	G	OP1-P-OP2	-5.43	111.46	119.60
1	AA	673	G	N1-C2-N3	5.43	127.16	123.90
1	AA	901	G	C5-C6-N1	-5.43	108.78	111.50
1	AA	1015	C	C5-C6-N1	-5.43	118.28	121.00
1	AA	1612	C	OP1-P-OP2	5.43	127.75	119.60
34	BA	283	C	C6-N1-C2	-5.43	118.13	120.30
1	CA	933	A	N7-C8-N9	5.43	116.51	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2516	G	C6-C5-N7	-5.43	127.14	130.40
1	AA	1051	C	OP2-P-O3'	5.43	117.14	105.20
1	AA	1317	G	N3-C2-N2	5.43	123.70	119.90
1	AA	1463	C	OP1-P-OP2	-5.43	111.46	119.60
1	AA	12	U	C6-N1-C2	-5.43	117.74	121.00
1	AA	585	U	C2-N3-C4	-5.43	123.74	127.00
1	AA	2527	C	N3-C4-C5	5.43	124.07	121.90
1	CA	1656	C	OP2-P-O3'	5.43	117.14	105.20
1	CA	2555	U	OP1-P-O3'	-5.43	93.26	105.20
29	C5	16	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	AA	58	U	OP2-P-O3'	5.42	117.14	105.20
1	AA	528	A	O5'-P-OP2	-5.42	100.82	105.70
1	AA	1206	G	O5'-P-OP2	5.42	117.21	110.70
1	AA	2379	G	N3-C4-N9	5.42	129.25	126.00
1	AA	2636	G	C2-N3-C4	5.42	114.61	111.90
2	AB	11	C	N3-C2-O2	-5.42	118.10	121.90
11	AN	99	LEU	CB-CG-CD1	-5.42	101.78	111.00
34	BA	579	G	C5-C6-N1	-5.42	108.79	111.50
1	CA	515	A	C6-N1-C2	-5.42	115.34	118.60
1	AA	710	G	C5-C6-O6	-5.42	125.35	128.60
1	AA	1874	C	N3-C4-N4	-5.42	114.20	118.00
1	AA	2385	G	N1-C6-O6	5.42	123.15	119.90
1	CA	424	G	N3-C4-C5	5.42	131.31	128.60
1	CA	1622	G	C8-N9-C4	5.42	108.57	106.40
1	CA	1698	A	C4-C5-N7	5.42	113.41	110.70
34	DA	754	C	N3-C2-O2	-5.42	118.10	121.90
1	AA	2862	G	C5-C6-O6	5.42	131.85	128.60
1	CA	1768	U	C2-N1-C1'	-5.42	111.19	117.70
1	CA	2088	G	N9-C4-C5	5.42	107.57	105.40
1	CA	2540	C	OP1-P-OP2	5.42	127.73	119.60
1	CA	2588	G	C6-N1-C2	5.42	128.35	125.10
1	AA	2263	G	C5-N7-C8	5.42	107.01	104.30
1	AA	2379	G	C4-N9-C1'	5.42	133.54	126.50
1	CA	330	A	N7-C8-N9	5.42	116.51	113.80
1	CA	513	A	C6-C5-N7	-5.42	128.51	132.30
1	CA	1351	C	N3-C4-N4	-5.42	114.21	118.00
1	AA	243	G	N3-C2-N2	5.42	123.69	119.90
1	AA	879	G	OP1-P-O3'	5.42	117.12	105.20
1	AA	1082	G	C6-C5-N7	5.42	133.65	130.40
1	CA	567	A	C4-C5-N7	5.42	113.41	110.70
56	DW	59	U	N3-C2-O2	-5.42	118.41	122.20
1	AA	622	G	N7-C8-N9	-5.42	110.39	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	724	A	N1-C6-N6	-5.42	115.35	118.60
1	AA	1741	C	N3-C4-C5	5.42	124.07	121.90
1	AA	1742	G	N1-C2-N3	-5.42	120.65	123.90
1	CA	734	A	N1-C2-N3	5.42	132.01	129.30
1	CA	1334	G	C6-C5-N7	5.42	133.65	130.40
1	CA	1678	G	N1-C2-N3	5.42	127.15	123.90
1	CA	2230	G	C8-N9-C4	-5.42	104.23	106.40
1	AA	2249	G	N9-C4-C5	-5.42	103.23	105.40
1	AA	2475	C	C2-N3-C4	-5.42	117.19	119.90
34	BA	769	G	C5-C6-O6	-5.42	125.35	128.60
1	CA	1768	U	C6-N1-C1'	5.42	128.78	121.20
1	CA	2222	G	C5-C6-O6	5.42	131.85	128.60
1	AA	1850	A	C5-C6-N1	-5.41	114.99	117.70
34	BA	771	G	N3-C4-N9	-5.41	122.75	126.00
1	CA	34	C	C5-C6-N1	5.41	123.71	121.00
34	DA	1065	U	P-O3'-C3'	5.41	126.19	119.70
1	AA	756	U	C5-C6-N1	-5.41	119.99	122.70
1	AA	2026	G	C2-N3-C4	-5.41	109.19	111.90
1	AA	2449	U	OP1-P-OP2	5.41	127.72	119.60
1	AA	2518	U	OP1-P-OP2	-5.41	111.48	119.60
1	CA	328	U	C5-C6-N1	-5.41	119.99	122.70
1	CA	1659	U	C2-N3-C4	-5.41	123.75	127.00
1	AA	213	G	C8-N9-C4	-5.41	104.24	106.40
1	AA	240	A	N3-C4-N9	5.41	131.73	127.40
1	AA	1958	A	N1-C6-N6	5.41	121.85	118.60
1	AA	2574	U	C4-C5-C6	5.41	122.95	119.70
1	AA	2696	U	N1-C2-O2	-5.41	119.01	122.80
34	BA	715	A	N1-C6-N6	5.41	121.85	118.60
1	CA	2086	U	O5'-P-OP1	5.41	117.19	110.70
34	DA	108	G	C4-C5-N7	5.41	112.97	110.80
1	AA	132	C	OP2-P-O3'	5.41	117.10	105.20
1	AA	253	C	C6-N1-C2	5.41	122.46	120.30
1	AA	831	A	C8-N9-C4	5.41	107.96	105.80
1	AA	1423	G	N1-C2-N3	5.41	127.14	123.90
1	AA	1711	A	C5-C6-N1	-5.41	115.00	117.70
1	AA	1754	G	N9-C4-C5	-5.41	103.24	105.40
1	AA	2115	G	C8-N9-C1'	5.41	134.03	127.00
34	BA	1057	G	N3-C4-N9	-5.41	122.75	126.00
1	CA	24	G	C8-N9-C4	5.41	108.56	106.40
1	CA	2396	G	C5-C6-N1	5.41	114.20	111.50
1	CA	2739	U	C5-C6-N1	-5.41	120.00	122.70
34	DA	275	G	C5-C6-O6	-5.41	125.35	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	DA	665	A	OP1-P-OP2	5.41	127.71	119.60
34	BA	9	G	N3-C4-N9	5.41	129.24	126.00
1	CA	41	C	OP2-P-O3'	5.41	117.10	105.20
1	CA	1673	U	C2-N3-C4	-5.41	123.76	127.00
1	CA	2679	A	OP2-P-O3'	5.41	117.10	105.20
1	AA	76	C	C6-N1-C2	5.41	122.46	120.30
1	AA	1354	A	C5-C6-N1	-5.41	115.00	117.70
1	AA	1628	G	N3-C4-C5	-5.41	125.90	128.60
1	AA	1642	A	N7-C8-N9	-5.41	111.10	113.80
1	AA	2802	C	C5-C6-N1	-5.41	118.30	121.00
34	DA	39	G	O5'-P-OP2	-5.41	100.83	105.70
1	AA	729	G	C8-N9-C1'	-5.40	119.97	127.00
1	AA	1045	U	C5-C4-O4	-5.40	122.66	125.90
1	AA	1262	C	C5-C6-N1	-5.40	118.30	121.00
1	AA	1721	G	C4-C5-N7	5.40	112.96	110.80
4	AD	94	LEU	CB-CG-CD1	-5.40	101.81	111.00
34	BA	329	A	O5'-P-OP2	-5.40	100.84	105.70
1	CA	1760	A	OP1-P-OP2	5.40	127.71	119.60
1	CA	2820	A	C4-N9-C1'	5.40	136.03	126.30
1	AA	176	G	C6-C5-N7	-5.40	127.16	130.40
1	AA	1066	A	C4-C5-N7	5.40	113.40	110.70
1	AA	1264	G	C6-C5-N7	-5.40	127.16	130.40
1	AA	1611	C	OP2-P-O3'	5.40	117.09	105.20
1	AA	1962	U	P-O3'-C3'	5.40	126.18	119.70
1	CA	1558	A	O4'-C1'-N9	5.40	112.52	108.20
1	CA	2598	A	O5'-P-OP1	5.40	117.18	110.70
34	DA	922	G	C8-N9-C4	-5.40	104.24	106.40
1	AA	239	G	C4-C5-N7	5.40	112.96	110.80
1	AA	550	U	N3-C4-O4	-5.40	115.62	119.40
1	AA	617	U	C4-C5-C6	-5.40	116.46	119.70
1	AA	1493	C	N1-C2-O2	5.40	122.14	118.90
1	AA	1502	G	C5-C6-N1	5.40	114.20	111.50
1	AA	2062	C	C6-N1-C2	5.40	122.46	120.30
34	BA	1394	A	C8-N9-C4	5.40	107.96	105.80
1	CA	330	A	C6-N1-C2	5.40	121.84	118.60
1	CA	2259	G	N3-C4-C5	5.40	131.30	128.60
1	CA	2407	G	O5'-P-OP2	-5.40	100.84	105.70
34	DA	1482	G	C5-C6-N1	-5.40	108.80	111.50
1	AA	1022	C	N1-C2-O2	5.40	122.14	118.90
1	CA	127	A	C8-N9-C4	-5.40	103.64	105.80
1	CA	1381	G	C5-C6-N1	5.40	114.20	111.50
1	AA	649	C	N1-C2-N3	5.40	122.98	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	817	G	C5-N7-C8	-5.40	101.60	104.30
2	AB	6	C	C2-N3-C4	-5.40	117.20	119.90
1	CA	690	G	C5-C6-O6	-5.40	125.36	128.60
1	CA	1600	C	N1-C2-O2	-5.40	115.66	118.90
1	CA	2014	A	N1-C6-N6	5.40	121.84	118.60
1	AA	103	C	N3-C4-C5	5.40	124.06	121.90
1	AA	200	A	O5'-P-OP1	5.40	117.17	110.70
1	AA	327	U	N3-C4-C5	5.40	117.84	114.60
1	AA	1854	G	C5-C6-N1	5.40	114.20	111.50
1	AA	2595	G	O5'-P-OP2	-5.40	100.84	105.70
34	BA	757	U	C5-C6-N1	-5.40	120.00	122.70
1	CA	562	U	C6-N1-C2	-5.40	117.76	121.00
1	CA	1721	G	N3-C4-N9	5.40	129.24	126.00
1	AA	324	A	C5-C6-N6	-5.39	119.38	123.70
1	AA	743	G	N3-C2-N2	5.39	123.68	119.90
1	AA	761	U	N1-C2-O2	5.39	126.58	122.80
1	AA	957	A	C2-N3-C4	5.39	113.30	110.60
1	AA	1278	G	N3-C2-N2	-5.39	116.12	119.90
1	AA	1526	G	C5-C6-O6	-5.39	125.36	128.60
1	AA	1745	A	C8-N9-C1'	-5.39	117.99	127.70
1	AA	2698	G	OP1-P-OP2	5.39	127.69	119.60
1	CA	2073	C	C6-N1-C2	5.39	122.46	120.30
1	CA	2694	G	C8-N9-C4	-5.39	104.24	106.40
1	AA	120	G	N9-C4-C5	5.39	107.56	105.40
1	AA	593	G	OP2-P-O3'	5.39	117.06	105.20
1	AA	714	U	C4-C5-C6	5.39	122.94	119.70
1	AA	896	A	C4-C5-C6	5.39	119.70	117.00
1	AA	917	A	C2-N3-C4	5.39	113.30	110.60
1	AA	1061	G	N1-C2-N2	-5.39	111.35	116.20
23	AZ	77	ASP	CB-CG-OD1	5.39	123.15	118.30
34	BA	514	C	N3-C2-O2	5.39	125.67	121.90
34	BA	728	A	O5'-P-OP2	-5.39	100.85	105.70
34	BA	791	G	N3-C4-N9	5.39	129.24	126.00
1	CA	127	A	N9-C4-C5	5.39	107.96	105.80
1	CA	1352	U	O5'-P-OP2	-5.39	100.85	105.70
1	CA	1934	C	C5-C6-N1	-5.39	118.30	121.00
34	DA	748	C	P-O3'-C3'	5.39	126.17	119.70
1	AA	499	G	C2-N3-C4	-5.39	109.20	111.90
1	AA	904	C	C2-N3-C4	-5.39	117.20	119.90
1	AA	2375	C	OP1-P-OP2	5.39	127.69	119.60
1	CA	240	G	C5-C6-O6	-5.39	125.36	128.60
1	CA	811	U	O5'-P-OP1	-5.39	100.85	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2500	U	C6-N1-C2	5.39	124.23	121.00
1	CA	2681	C	C6-N1-C2	-5.39	118.14	120.30
1	AA	1314	A	N9-C4-C5	5.39	107.95	105.80
1	AA	1475	G	N3-C4-C5	-5.39	125.91	128.60
1	AA	1860	A	N7-C8-N9	-5.39	111.11	113.80
1	AA	1212	C	C5-C4-N4	-5.39	116.43	120.20
1	AA	2038	U	OP2-P-O3'	5.39	117.05	105.20
1	AA	2548	G	N3-C4-C5	-5.39	125.91	128.60
34	BA	611	A	O5'-P-OP2	-5.39	100.85	105.70
34	BA	1405	G	OP2-P-O3'	5.39	117.05	105.20
1	CA	2432	A	N1-C2-N3	5.39	131.99	129.30
1	AA	438	G	O5'-P-OP2	-5.38	100.86	105.70
1	AA	1245	C	C6-N1-C2	5.38	122.45	120.30
1	AA	2693	C	C5-C6-N1	-5.38	118.31	121.00
1	CA	2570	G	C5-C6-N1	-5.38	108.81	111.50
1	AA	2468	C	N1-C2-N3	-5.38	115.43	119.20
34	BA	1087	G	N7-C8-N9	5.38	115.79	113.10
1	AA	195	U	N1-C2-N3	5.38	118.13	114.90
1	AA	907	U	C2-N3-C4	-5.38	123.77	127.00
1	AA	1234	A	N1-C6-N6	5.38	121.83	118.60
1	AA	1720	U	N3-C4-C5	5.38	117.83	114.60
1	AA	1766	G	N7-C8-N9	5.38	115.79	113.10
1	AA	2516	U	C2-N3-C4	-5.38	123.77	127.00
1	AA	2645	G	N1-C6-O6	-5.38	116.67	119.90
1	AA	2705	A	C5-C6-N6	-5.38	119.40	123.70
2	AB	17	C	N3-C4-C5	-5.38	119.75	121.90
34	BA	399	G	C8-N9-C4	5.38	108.55	106.40
34	BA	899	C	N3-C2-O2	5.38	125.67	121.90
34	BA	1527	C	C5-C6-N1	-5.38	118.31	121.00
1	AA	1149	A	C2-N3-C4	-5.38	107.91	110.60
1	AA	1912	A	OP2-P-O3'	5.38	117.03	105.20
1	AA	1926	G	C2-N3-C4	5.38	114.59	111.90
1	CA	659	C	N1-C2-O2	-5.38	115.67	118.90
1	CA	1385	G	N3-C4-N9	-5.38	122.77	126.00
1	CA	1408	C	N1-C2-O2	-5.38	115.67	118.90
1	CA	2229	C	O5'-P-OP2	5.38	117.16	110.70
34	DA	618	C	N1-C2-O2	5.38	122.13	118.90
1	AA	254	A	N3-C4-N9	-5.38	123.10	127.40
1	AA	1423	G	C5-C6-O6	5.38	131.83	128.60
1	AA	1674	G	C2-N3-C4	5.38	114.59	111.90
1	AA	2446	A	N1-C2-N3	5.38	131.99	129.30
1	AA	2593	G	C8-N9-C4	-5.38	104.25	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	AD	257	LEU	CB-CG-CD1	-5.38	101.86	111.00
34	BA	1030(B)	C	C6-N1-C2	-5.38	118.15	120.30
1	CA	1660	C	O4'-C1'-N1	5.38	112.50	108.20
1	CA	2253	G	C5-C6-O6	-5.38	125.37	128.60
1	CA	2568	C	O5'-P-OP2	-5.38	100.86	105.70
1	AA	19	C	C5-C4-N4	-5.38	116.44	120.20
1	AA	177	G	C5-C6-O6	5.38	131.83	128.60
1	AA	2505	U	C6-N1-C2	-5.38	117.77	121.00
34	BA	900	A	C5-C6-N6	-5.38	119.40	123.70
1	CA	205	G	N9-C4-C5	-5.38	103.25	105.40
1	CA	2066	C	C5-C6-N1	-5.38	118.31	121.00
34	DA	115	G	P-O3'-C3'	5.38	126.15	119.70
1	AA	2019	G	O5'-P-OP1	5.38	117.15	110.70
1	AA	2397	C	C6-N1-C2	5.38	122.45	120.30
1	AA	483	A	C5-C6-N1	-5.37	115.01	117.70
1	AA	672	G	N3-C4-C5	-5.37	125.91	128.60
1	AA	1951	G	C5-C6-O6	5.37	131.82	128.60
1	AA	2622	C	C4-C5-C6	5.37	120.09	117.40
1	AA	2847	G	O5'-P-OP1	-5.37	100.86	105.70
1	CA	123	G	C5-C6-O6	-5.37	125.38	128.60
1	CA	450	G	N3-C4-C5	-5.37	125.91	128.60
1	CA	1265	A	C2-N3-C4	-5.37	107.91	110.60
1	CA	1552	G	C5-C6-O6	5.37	131.82	128.60
1	CA	2576	G	N3-C4-C5	-5.37	125.91	128.60
1	CA	2699	C	N3-C2-O2	-5.37	118.14	121.90
1	AA	16	G	N9-C4-C5	5.37	107.55	105.40
1	AA	1248	G	N1-C2-N2	-5.37	111.36	116.20
1	AA	1497	G	N1-C6-O6	-5.37	116.68	119.90
1	AA	2321	A	C8-N9-C4	-5.37	103.65	105.80
1	AA	2527	C	N3-C4-N4	5.37	121.76	118.00
34	BA	576	G	N1-C6-O6	5.37	123.12	119.90
1	CA	2588	G	N9-C4-C5	-5.37	103.25	105.40
1	AA	1282	G	C8-N9-C1'	-5.37	120.02	127.00
1	AA	2376	C	N1-C2-N3	-5.37	115.44	119.20
1	AA	599	U	N1-C2-O2	-5.37	119.04	122.80
1	AA	911	G	N3-C4-N9	5.37	129.22	126.00
1	AA	1357	G	C5-N7-C8	-5.37	101.62	104.30
1	AA	1974	A	C2-N3-C4	-5.37	107.92	110.60
34	BA	1402	C	N1-C2-O2	-5.37	115.68	118.90
1	CA	52	A	C2-N3-C4	-5.37	107.92	110.60
1	CA	513	A	N1-C2-N3	5.37	131.98	129.30
1	CA	2549	G	N3-C4-N9	5.37	129.22	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2691	C	C5-C6-N1	5.37	123.68	121.00
34	DA	577	G	OP2-P-O3'	5.37	117.01	105.20
1	AA	18	C	C6-N1-C2	5.37	122.45	120.30
1	AA	197	C	C6-N1-C2	5.37	122.45	120.30
1	AA	470	C	C5-C4-N4	5.37	123.96	120.20
1	AA	2057	G	O5'-P-OP2	-5.37	100.87	105.70
1	AA	2613	C	N3-C4-N4	5.37	121.76	118.00
34	BA	918	A	N9-C4-C5	5.37	107.95	105.80
1	CA	463	G	O5'-P-OP2	-5.37	100.87	105.70
1	CA	2570	G	C5-C6-O6	5.37	131.82	128.60
1	AA	854	U	N3-C4-O4	-5.37	115.64	119.40
1	AA	1009	C	O5'-P-OP2	-5.37	100.87	105.70
1	AA	1617	A	C2-N3-C4	-5.37	107.92	110.60
1	AA	1858	C	OP2-P-O3'	5.37	117.00	105.20
2	AB	98	G	O5'-P-OP2	-5.37	100.87	105.70
1	CA	64	A	O5'-P-OP1	-5.37	100.87	105.70
1	CA	298	G	N3-C4-N9	5.37	129.22	126.00
1	CA	1278	A	N1-C2-N3	5.37	131.98	129.30
1	AA	488	C	C6-N1-C1'	5.36	127.24	120.80
1	AA	1475	G	P-O3'-C3'	5.36	126.14	119.70
1	AA	1815	A	OP1-P-O3'	5.36	117.00	105.20
1	AA	2220	A	O5'-P-OP2	5.36	117.14	110.70
1	AA	2829	G	N1-C2-N3	5.36	127.12	123.90
34	BA	115	G	C2-N3-C4	5.36	114.58	111.90
34	BA	791	G	N3-C2-N2	5.36	123.65	119.90
34	BA	1484	C	N1-C2-O2	-5.36	115.68	118.90
1	CA	12	U	N3-C2-O2	-5.36	118.45	122.20
1	CA	687	C	C5-C6-N1	5.36	123.68	121.00
1	CA	749	C	O5'-P-OP2	5.36	117.14	110.70
1	AA	976	G	C4-C5-N7	-5.36	108.66	110.80
1	AA	2406	C	OP1-P-O3'	5.36	116.99	105.20
1	CA	312	G	C8-N9-C4	-5.36	104.26	106.40
1	CA	2707	G	C2-N3-C4	-5.36	109.22	111.90
34	DA	7	G	N3-C4-C5	5.36	131.28	128.60
1	AA	2236	G	N7-C8-N9	5.36	115.78	113.10
1	AA	2239	A	C4-C5-C6	5.36	119.68	117.00
34	DA	698	G	O5'-P-OP2	-5.36	100.88	105.70
1	AA	124	A	C5-N7-C8	5.36	106.58	103.90
1	AA	309	C	N3-C2-O2	-5.36	118.15	121.90
1	AA	990	A	C1'-O4'-C4'	-5.36	105.61	109.90
1	AA	1015	C	C6-N1-C1'	-5.36	114.37	120.80
1	AA	1371	G	C4-C5-N7	-5.36	108.66	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	504	C	C2-N1-C1'	5.36	124.69	118.80
34	BA	586	C	N1-C2-O2	-5.36	115.69	118.90
34	BA	824	C	OP2-P-O3'	5.36	116.99	105.20
1	CA	704	G	C5-C6-O6	-5.36	125.39	128.60
1	AA	345	G	N1-C6-O6	5.36	123.11	119.90
1	AA	845	G	N7-C8-N9	-5.36	110.42	113.10
1	AA	1737	A	OP1-P-OP2	-5.36	111.57	119.60
27	A3	56	VAL	CB-CA-C	-5.36	101.22	111.40
34	BA	768	A	C5-C6-N6	-5.36	119.42	123.70
1	CA	141	A	OP2-P-O3'	5.36	116.98	105.20
1	CA	592	G	N3-C2-N2	5.36	123.65	119.90
1	CA	662	G	C8-N9-C4	5.36	108.54	106.40
1	CA	1296	G	C8-N9-C4	-5.36	104.26	106.40
1	CA	1936	A	C8-N9-C4	5.36	107.94	105.80
34	DA	335	C	N1-C2-O2	5.36	122.11	118.90
1	AA	281	G	C8-N9-C4	5.35	108.54	106.40
1	AA	2858	G	N1-C2-N3	-5.35	120.69	123.90
34	BA	884	U	N3-C4-C5	-5.35	111.39	114.60
1	CA	829	A	O5'-P-OP1	-5.35	100.88	105.70
1	CA	2239	G	C5-C6-O6	5.35	131.81	128.60
1	AA	129	G	O5'-P-OP1	5.35	117.12	110.70
1	AA	510	C	C4-C5-C6	-5.35	114.72	117.40
1	AA	1363	A	C2-N3-C4	-5.35	107.92	110.60
1	AA	1684	A	N3-C4-N9	-5.35	123.12	127.40
1	AA	2854	G	C5-C6-N1	5.35	114.18	111.50
1	CA	1822	G	N3-C4-N9	-5.35	122.79	126.00
1	AA	2026	G	N1-C6-O6	5.35	123.11	119.90
1	AA	2829	G	C2-N3-C4	-5.35	109.22	111.90
31	A7	33	ARG	NE-CZ-NH1	5.35	122.98	120.30
34	BA	299	G	C6-C5-N7	-5.35	127.19	130.40
1	CA	2499	C	C5-C4-N4	-5.35	116.45	120.20
1	CA	2727	G	OP2-P-O3'	5.35	116.97	105.20
1	AA	211	A	N1-C6-N6	-5.35	115.39	118.60
1	AA	1017	G	N7-C8-N9	5.35	115.77	113.10
1	AA	2466	G	C5-C6-O6	5.35	131.81	128.60
1	AA	2484	G	N3-C4-N9	5.35	129.21	126.00
1	AA	2625	U	OP1-P-OP2	5.35	127.62	119.60
2	AB	37	C	OP1-P-O3'	-5.35	93.43	105.20
1	CA	561	G	N1-C6-O6	-5.35	116.69	119.90
1	CA	1968	G	OP2-P-O3'	5.35	116.97	105.20
1	AA	724	A	N9-C4-C5	5.35	107.94	105.80
1	AA	1975	A	N1-C6-N6	-5.35	115.39	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	429	U	O5'-P-OP1	-5.35	100.89	105.70
1	CA	668	G	N9-C4-C5	-5.35	103.26	105.40
1	CA	687	C	N3-C4-C5	5.35	124.04	121.90
1	CA	1558	A	C4-C5-N7	5.35	113.37	110.70
1	CA	2534	A	OP2-P-O3'	5.35	116.96	105.20
1	AA	192	C	N1-C2-O2	-5.35	115.69	118.90
1	AA	407	U	O5'-P-OP1	5.35	117.11	110.70
1	AA	413	G	C5-N7-C8	-5.35	101.63	104.30
1	CA	1675	C	C5-C6-N1	5.35	123.67	121.00
1	CA	1970	A	O4'-C1'-N9	-5.35	103.92	108.20
1	AA	1552	C	C6-N1-C2	-5.34	118.16	120.30
34	BA	1432	G	N1-C6-O6	-5.34	116.69	119.90
1	CA	389	G	C8-N9-C4	5.34	108.54	106.40
1	CA	2236	C	OP2-P-O3'	5.34	116.96	105.20
34	DA	353	A	C5-C6-N6	-5.34	119.42	123.70
1	AA	884	C	N1-C2-O2	-5.34	115.69	118.90
1	AA	2636	G	N3-C4-C5	-5.34	125.93	128.60
1	AA	2773	C	C6-N1-C2	5.34	122.44	120.30
34	BA	454	C	N3-C2-O2	-5.34	118.16	121.90
1	CA	1782	C	OP1-P-OP2	5.34	127.61	119.60
1	CA	2254	C	OP2-P-O3'	5.34	116.95	105.20
1	CA	2546	U	C5-C6-N1	-5.34	120.03	122.70
1	AA	1192	C	N3-C2-O2	5.34	125.64	121.90
1	AA	1476	C	C2-N3-C4	-5.34	117.23	119.90
1	AA	1962	U	OP1-P-O3'	5.34	116.95	105.20
47	BN	44	LEU	CA-CB-CG	5.34	127.58	115.30
56	BW	12	U	N3-C2-O2	-5.34	118.46	122.20
1	CA	1027	A	C8-N9-C4	5.34	107.94	105.80
1	CA	1609	A	N9-C4-C5	-5.34	103.66	105.80
1	CA	1905	C	N1-C2-N3	-5.34	115.46	119.20
34	DA	1482	G	C4-C5-C6	5.34	122.00	118.80
1	AA	256	C	C2-N3-C4	-5.34	117.23	119.90
1	AA	1186	U	C4-C5-C6	5.34	122.90	119.70
1	AA	1653	C	C6-N1-C1'	-5.34	114.39	120.80
1	CA	786	C	OP2-P-O3'	5.34	116.94	105.20
1	CA	2698	U	C4-C5-C6	-5.34	116.50	119.70
2	CB	78	A	C6-N1-C2	-5.34	115.40	118.60
34	DA	853	G	C6-C5-N7	-5.34	127.20	130.40
1	AA	744	C	O5'-P-OP2	-5.34	100.90	105.70
1	AA	1046	A	O5'-P-OP1	-5.34	100.90	105.70
1	AA	1225	C	C6-N1-C2	5.34	122.44	120.30
1	AA	1378	G	C2-N3-C4	5.34	114.57	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1703	C	OP1-P-OP2	-5.34	111.60	119.60
1	AA	2487	C	N1-C2-O2	5.34	122.10	118.90
1	AA	2506	G	C5-C6-N1	-5.34	108.83	111.50
34	BA	912	C	N3-C2-O2	5.34	125.64	121.90
34	BA	1529	G	N1-C6-O6	-5.34	116.70	119.90
1	CA	263	C	C5-C4-N4	-5.34	116.47	120.20
1	CA	424	G	N3-C4-N9	-5.34	122.80	126.00
1	AA	663	G	C5-C6-O6	5.33	131.80	128.60
1	AA	721	G	N7-C8-N9	5.33	115.77	113.10
1	AA	1449	C	C4-C5-C6	5.33	120.07	117.40
1	CA	942	G	N1-C6-O6	-5.33	116.70	119.90
1	CA	1785	A	O5'-P-OP1	-5.33	100.90	105.70
1	CA	2486	G	C4-N9-C1'	5.33	133.44	126.50
1	AA	471	C	C5-C6-N1	-5.33	118.33	121.00
1	AA	1786	A	OP1-P-OP2	5.33	127.60	119.60
1	AA	2705	A	C8-N9-C4	5.33	107.93	105.80
34	BA	123	C	N1-C2-O2	-5.33	115.70	118.90
56	BW	76	A	C4-C5-N7	5.33	113.37	110.70
1	CA	386	G	O4'-C1'-N9	5.33	112.47	108.20
1	AA	119	G	C6-N1-C2	-5.33	121.90	125.10
1	AA	586	G	C5-N7-C8	5.33	106.97	104.30
1	AA	1648	U	N1-C2-O2	5.33	126.53	122.80
1	AA	2742	G	C5-C6-O6	5.33	131.80	128.60
1	CA	1185	C	O5'-P-OP2	-5.33	100.90	105.70
1	CA	1763	G	N3-C4-C5	5.33	131.27	128.60
1	CA	1914	C	C2-N1-C1'	5.33	124.66	118.80
1	AA	623	G	N1-C6-O6	-5.33	116.70	119.90
1	AA	1023	G	C5-C6-N1	5.33	114.17	111.50
1	AA	2479	C	C6-N1-C2	5.33	122.43	120.30
9	AK	128	LEU	C-N-CA	5.33	144.38	122.00
34	BA	160	A	C8-N9-C4	-5.33	103.67	105.80
34	BA	1442	G	N1-C6-O6	5.33	123.10	119.90
1	CA	393	C	O5'-P-OP2	5.33	117.10	110.70
1	CA	2738	A	C8-N9-C4	5.33	107.93	105.80
1	AA	493	G	C5-C6-N1	5.33	114.17	111.50
1	AA	1922	A	N1-C6-N6	-5.33	115.40	118.60
1	AA	2289	G	C4-C5-N7	-5.33	108.67	110.80
1	AA	2600	G	C2-N3-C4	-5.33	109.23	111.90
17	AT	95	ARG	NE-CZ-NH2	5.33	122.96	120.30
34	BA	991	U	P-O3'-C3'	5.33	126.09	119.70
1	CA	571	A	C5-C6-N1	5.33	120.36	117.70
1	CA	2444	G	C4-N9-C1'	5.33	133.43	126.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	DA	926	G	N9-C4-C5	5.33	107.53	105.40
34	DA	1503	A	C8-N9-C4	5.33	107.93	105.80
1	AA	2638	C	C4-C5-C6	5.33	120.06	117.40
1	AA	310	C	C6-N1-C2	5.33	122.43	120.30
1	AA	348	A	OP1-P-O3'	5.33	116.92	105.20
1	AA	577	U	C5-C4-O4	-5.33	122.70	125.90
1	AA	1066	A	C6-C5-N7	-5.33	128.57	132.30
2	AB	85	G	C4-N9-C1'	5.33	133.42	126.50
34	BA	753	A	OP1-P-O3'	5.33	116.92	105.20
1	CA	681	G	C8-N9-C4	5.33	108.53	106.40
1	CA	1817	G	O5'-P-OP2	-5.33	100.91	105.70
1	CA	2286	A	C4-C5-C6	5.33	119.66	117.00
1	AA	1052	C	C5-C6-N1	-5.32	118.34	121.00
1	AA	1626	A	O5'-P-OP1	-5.32	100.91	105.70
1	AA	1792	C	N1-C2-O2	-5.32	115.71	118.90
1	AA	2411	G	N1-C6-O6	-5.32	116.71	119.90
1	AA	2412	G	N9-C4-C5	-5.32	103.27	105.40
1	CA	1274	A	C6-C5-N7	-5.32	128.57	132.30
1	CA	1648	C	C2-N1-C1'	-5.32	112.94	118.80
1	AA	1989	C	N3-C2-O2	-5.32	118.17	121.90
1	AA	2529	C	C5-C6-N1	-5.32	118.34	121.00
34	BA	1431	C	C2-N3-C4	-5.32	117.24	119.90
1	CA	277	C	C6-N1-C1'	-5.32	114.41	120.80
1	CA	777	A	C6-N1-C2	-5.32	115.41	118.60
1	CA	1769	G	N3-C2-N2	-5.32	116.17	119.90
1	AA	255	G	OP2-P-O3'	5.32	116.90	105.20
1	AA	748	G	OP1-P-OP2	5.32	127.58	119.60
1	AA	1331	G	C5-C6-N1	5.32	114.16	111.50
1	AA	1634	C	N3-C2-O2	5.32	125.62	121.90
1	AA	2622	C	C6-N1-C2	-5.32	118.17	120.30
1	AA	2882	G	OP1-P-OP2	5.32	127.58	119.60
2	AB	80	U	C6-N1-C2	5.32	124.19	121.00
2	AB	91	C	C5-C4-N4	-5.32	116.47	120.20
2	AB	102	A	OP2-P-O3'	5.32	116.91	105.20
34	BA	896	C	O5'-P-OP2	-5.32	100.91	105.70
34	BA	1524	C	C4-C5-C6	5.32	120.06	117.40
1	CA	1309	G	N3-C2-N2	-5.32	116.18	119.90
1	CA	2226	C	N1-C2-O2	5.32	122.09	118.90
1	CA	2674	G	N1-C6-O6	-5.32	116.71	119.90
1	AA	148	C	N3-C4-C5	5.32	124.03	121.90
1	AA	526	A	N1-C2-N3	5.32	131.96	129.30
1	AA	964	A	C6-N1-C2	5.32	121.79	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1691	C	C6-N1-C2	-5.32	118.17	120.30
1	AA	2383	G	N3-C4-C5	-5.32	125.94	128.60
1	AA	2493	G	C2-N3-C4	-5.32	109.24	111.90
1	CA	1428	C	N3-C4-N4	-5.32	114.28	118.00
1	AA	751	G	C2-N3-C4	5.32	114.56	111.90
1	AA	846	G	C4-C5-N7	-5.32	108.67	110.80
1	AA	1079	U	N3-C4-O4	-5.32	115.68	119.40
1	AA	1081	U	C5-C6-N1	-5.32	120.04	122.70
1	AA	2428	C	N3-C4-C5	-5.32	119.77	121.90
1	AA	2495	C	O5'-P-OP2	5.32	117.08	110.70
1	AA	2614	A	OP1-P-O3'	5.32	116.90	105.20
34	BA	295	C	C6-N1-C2	5.32	122.43	120.30
34	BA	1502	A	C8-N9-C4	-5.32	103.67	105.80
49	BP	51	VAL	CB-CA-C	-5.32	101.30	111.40
1	CA	915	C	N3-C4-C5	-5.32	119.77	121.90
2	CB	56	G	C8-N9-C4	-5.32	104.27	106.40
1	AA	559	U	N3-C4-O4	-5.32	115.68	119.40
1	AA	948	C	O5'-P-OP1	-5.32	100.92	105.70
1	AA	1617	A	N7-C8-N9	-5.32	111.14	113.80
1	AA	2293	C	C4-C5-C6	5.32	120.06	117.40
1	CA	2653	U	N3-C2-O2	-5.32	118.48	122.20
1	CA	2675	A	O5'-P-OP1	5.32	117.08	110.70
1	AA	1374	G	N3-C4-N9	5.31	129.19	126.00
1	AA	1650	C	N1-C2-N3	5.31	122.92	119.20
1	CA	1032	A	N7-C8-N9	-5.31	111.14	113.80
1	CA	1367	A	C2-N3-C4	-5.31	107.94	110.60
1	AA	411	U	C2-N3-C4	-5.31	123.81	127.00
1	AA	1256	U	C6-N1-C2	5.31	124.19	121.00
1	AA	2057	G	N7-C8-N9	-5.31	110.44	113.10
1	AA	2070	G	N3-C4-N9	5.31	129.19	126.00
1	AA	2601	A	C6-C5-N7	5.31	136.02	132.30
1	AA	2772	G	C4-C5-N7	-5.31	108.67	110.80
34	BA	1530	G	N3-C4-N9	-5.31	122.81	126.00
1	CA	150	C	N3-C4-N4	-5.31	114.28	118.00
1	CA	470	A	OP1-P-OP2	5.31	127.57	119.60
1	CA	2805	G	C4-N9-C1'	-5.31	119.59	126.50
34	DA	508	C	C2-N1-C1'	5.31	124.64	118.80
1	AA	567	C	OP1-P-OP2	5.31	127.56	119.60
1	AA	753	A	C5-N7-C8	-5.31	101.25	103.90
1	AA	1309	U	N3-C2-O2	-5.31	118.48	122.20
1	AA	1390	G	O4'-C1'-N9	5.31	112.45	108.20
1	AA	2241	C	N1-C2-O2	5.31	122.09	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AB	51	G	O5'-P-OP1	5.31	117.07	110.70
34	BA	1499	A	C8-N9-C4	5.31	107.92	105.80
1	CA	944	G	C5-C6-N1	-5.31	108.84	111.50
1	AA	534	C	O5'-P-OP2	-5.31	100.92	105.70
1	AA	895	G	C5-C6-O6	5.31	131.78	128.60
1	AA	1477	U	N1-C2-N3	-5.31	111.72	114.90
2	AB	38	C	C6-N1-C2	5.31	122.42	120.30
1	CA	596	G	C2-N3-C4	5.31	114.55	111.90
1	CA	795	C	C4-C5-C6	5.31	120.05	117.40
1	CA	2782	G	C6-C5-N7	-5.31	127.22	130.40
1	CA	2805	G	C4-C5-N7	-5.31	108.68	110.80
34	DA	673	G	N3-C4-C5	-5.31	125.95	128.60
1	AA	474	U	N1-C2-N3	5.30	118.08	114.90
1	AA	1379	C	C5-C4-N4	-5.30	116.49	120.20
1	AA	1491	A	OP2-P-O3'	5.30	116.87	105.20
1	AA	2484	G	N3-C4-C5	-5.30	125.95	128.60
2	AB	1	U	N1-C2-O2	5.30	126.51	122.80
34	BA	545	C	N1-C2-O2	5.30	122.08	118.90
1	CA	600	G	OP2-P-O3'	5.30	116.87	105.20
1	CA	1333	C	C5-C6-N1	5.30	123.65	121.00
1	CA	1408	C	N3-C4-C5	-5.30	119.78	121.90
1	CA	1968	G	C5-C6-O6	-5.30	125.42	128.60
1	CA	2233	U	N1-C2-O2	-5.30	119.09	122.80
1	CA	2501	C	N3-C4-C5	5.30	124.02	121.90
1	AA	178	G	C8-N9-C4	-5.30	104.28	106.40
1	AA	2084	A	OP2-P-O3'	5.30	116.86	105.20
1	AA	2509	A	O4'-C1'-N9	5.30	112.44	108.20
1	AA	2760	G	C6-N1-C2	-5.30	121.92	125.10
34	DA	530	G	N3-C4-N9	-5.30	122.82	126.00
34	DA	1383	C	C6-N1-C2	-5.30	118.18	120.30
1	AA	1565	G	N3-C4-N9	-5.30	122.82	126.00
1	AA	1752	G	C5-C6-O6	5.30	131.78	128.60
1	AA	2301	G	C5-N7-C8	-5.30	101.65	104.30
34	BA	523	A	N9-C4-C5	5.30	107.92	105.80
34	BA	1030(B)	C	C6-N1-C1'	-5.30	114.44	120.80
1	CA	271(A)	A	C8-N9-C4	5.30	107.92	105.80
1	AA	1239	A	C5-C6-N1	5.30	120.35	117.70
1	AA	1351	C	C2-N1-C1'	5.30	124.63	118.80
34	BA	818	G	N3-C4-C5	-5.30	125.95	128.60
1	CA	296	C	N3-C4-C5	-5.30	119.78	121.90
1	CA	1236	G	O5'-P-OP1	-5.30	100.93	105.70
1	CA	1271	G	N3-C4-N9	5.30	129.18	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	987	G	C4-C5-N7	-5.30	108.68	110.80
1	AA	1373	C	N1-C2-O2	-5.30	115.72	118.90
1	AA	1440	U	N3-C2-O2	-5.30	118.49	122.20
1	AA	2003	A	OP1-P-OP2	-5.30	111.65	119.60
1	CA	201	C	N3-C4-C5	5.30	124.02	121.90
1	CA	2056	G	OP2-P-O3'	5.30	116.85	105.20
1	CA	2379	G	C4-N9-C1'	5.30	133.39	126.50
34	DA	175	C	N1-C2-O2	5.30	122.08	118.90
1	AA	586	G	N9-C4-C5	5.29	107.52	105.40
1	AA	847	A	C5-C6-N6	5.29	127.94	123.70
1	AA	2240	G	N3-C2-N2	5.29	123.61	119.90
34	BA	122	G	C5-C6-O6	-5.29	125.42	128.60
1	CA	263	C	N3-C4-N4	5.29	121.71	118.00
1	CA	448	U	O5'-P-OP2	5.29	117.05	110.70
1	CA	2050	C	N3-C4-C5	5.29	124.02	121.90
1	CA	2320	A	C2-N3-C4	5.29	113.25	110.60
56	DW	73	A	O4'-C1'-N9	5.29	112.44	108.20
1	AA	649	C	C5-C6-N1	-5.29	118.35	121.00
1	AA	2282	G	O5'-P-OP2	5.29	117.05	110.70
1	AA	2618	C	N3-C4-C5	5.29	124.02	121.90
31	A7	9	ARG	NE-CZ-NH2	-5.29	117.65	120.30
1	CA	2501	C	C2-N1-C1'	-5.29	112.98	118.80
1	AA	1921	G	N9-C4-C5	5.29	107.52	105.40
1	AA	2620	G	OP1-P-O3'	5.29	116.84	105.20
2	AB	72	G	C4-C5-N7	5.29	112.92	110.80
1	CA	2581	G	N3-C2-N2	-5.29	116.19	119.90
1	AA	237	G	N1-C2-N3	-5.29	120.73	123.90
1	AA	2638	C	C2-N1-C1'	-5.29	112.98	118.80
1	CA	2429	G	N3-C4-N9	-5.29	122.83	126.00
1	AA	913	A	C8-N9-C4	-5.29	103.68	105.80
1	AA	1321	A	N1-C6-N6	-5.29	115.43	118.60
1	AA	1845	G	O5'-P-OP2	-5.29	100.94	105.70
1	AA	1847	G	C6-C5-N7	5.29	133.57	130.40
34	BA	347	G	N3-C4-C5	-5.29	125.96	128.60
34	BA	1505	G	N3-C4-N9	-5.29	122.83	126.00
1	CA	686	G	C5-C6-O6	-5.29	125.43	128.60
34	DA	26	A	N1-C2-N3	5.29	131.94	129.30
1	AA	1717	C	OP1-P-O3'	5.29	116.83	105.20
34	BA	339	C	N3-C2-O2	5.29	125.60	121.90
1	CA	1007	C	C5-C6-N1	-5.29	118.36	121.00
1	AA	649	C	O5'-P-OP2	5.29	117.04	110.70
1	AA	1879	A	C8-N9-C4	5.29	107.91	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	122	G	C8-N9-C4	5.29	108.51	106.40
1	CA	374	A	OP1-P-OP2	5.29	127.53	119.60
1	CA	1663	C	N1-C2-O2	5.29	122.07	118.90
1	CA	1695	G	N1-C6-O6	5.29	123.07	119.90
1	CA	1980	G	C5-C6-O6	-5.29	125.43	128.60
1	CA	2249	U	N3-C2-O2	-5.29	118.50	122.20
1	AA	128	C	OP2-P-O3'	5.28	116.82	105.20
1	AA	608	G	C2-N3-C4	5.28	114.54	111.90
1	AA	1611	C	C4-C5-C6	5.28	120.04	117.40
1	AA	1892	G	O5'-P-OP2	-5.28	100.94	105.70
1	AA	2403	G	OP1-P-OP2	5.28	127.52	119.60
1	AA	2537	G	C5-N7-C8	-5.28	101.66	104.30
1	AA	2833	A	C6-C5-N7	-5.28	128.60	132.30
1	AA	2879	G	C5-N7-C8	5.28	106.94	104.30
34	BA	804	U	C5-C6-N1	-5.28	120.06	122.70
1	CA	1211	U	N3-C2-O2	-5.28	118.50	122.20
1	CA	2038	G	N1-C2-N2	-5.28	111.44	116.20
34	DA	355	C	N3-C4-C5	-5.28	119.79	121.90
1	AA	1443	U	C2-N3-C4	-5.28	123.83	127.00
1	AA	422	U	O4'-C1'-N1	5.28	112.42	108.20
1	AA	621	G	N3-C2-N2	5.28	123.60	119.90
1	AA	1238	G	C5-C6-N1	5.28	114.14	111.50
1	AA	2089	G	N9-C4-C5	5.28	107.51	105.40
1	AA	2353	G	C6-C5-N7	5.28	133.57	130.40
1	AA	2376	C	C2-N1-C1'	-5.28	112.99	118.80
1	AA	2410	U	N1-C2-N3	5.28	118.07	114.90
34	BA	711	G	N3-C4-N9	5.28	129.17	126.00
1	CA	1760	A	N1-C2-N3	5.28	131.94	129.30
1	CA	1829	A	O5'-P-OP2	-5.28	100.95	105.70
1	CA	2446	G	N3-C2-N2	5.28	123.60	119.90
1	CA	12	U	C6-N1-C1'	-5.28	113.81	121.20
1	CA	265	A	C2-N3-C4	-5.28	107.96	110.60
1	CA	453	C	N3-C4-C5	5.28	124.01	121.90
1	CA	641	C	O5'-P-OP1	-5.28	100.95	105.70
1	CA	2521	C	N1-C2-O2	-5.28	115.73	118.90
1	AA	1403	U	N3-C4-C5	5.28	117.77	114.60
1	AA	1977	U	C5-C4-O4	-5.28	122.73	125.90
1	AA	2287	C	C2-N3-C4	-5.28	117.26	119.90
1	AA	2422	G	N9-C4-C5	-5.28	103.29	105.40
1	AA	2562	G	OP1-P-OP2	-5.28	111.68	119.60
1	AA	2890	C	C4-C5-C6	5.28	120.04	117.40
1	CA	2741	A	N7-C8-N9	-5.28	111.16	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	DA	774	G	OP2-P-O3'	5.28	116.81	105.20
1	AA	189	U	OP1-P-OP2	5.28	127.51	119.60
1	AA	893	C	O4'-C1'-N1	5.28	112.42	108.20
1	AA	1326	G	C5-C6-N1	5.28	114.14	111.50
1	AA	1844	G	N3-C4-N9	5.28	129.17	126.00
1	AA	1980	C	N3-C4-N4	5.28	121.69	118.00
1	AA	2503	U	C5-C4-O4	-5.28	122.73	125.90
1	AA	2631	C	N3-C2-O2	-5.28	118.21	121.90
1	AA	721	G	O5'-P-OP2	5.27	117.03	110.70
1	AA	748	G	OP2-P-O3'	5.27	116.80	105.20
1	AA	2456	G	C6-C5-N7	5.27	133.56	130.40
1	AA	2633	A	C5-C6-N6	5.27	127.92	123.70
1	CA	114	U	C6-N1-C1'	-5.27	113.82	121.20
34	DA	841	U	C5-C6-N1	5.27	125.34	122.70
1	AA	31	C	N1-C2-O2	5.27	122.06	118.90
1	AA	374	U	OP2-P-O3'	5.27	116.80	105.20
1	AA	1328	U	C4-C5-C6	5.27	122.86	119.70
1	AA	1816	A	N1-C6-N6	5.27	121.76	118.60
1	AA	1881	G	N1-C2-N3	5.27	127.06	123.90
1	AA	1972	G	OP2-P-O3'	5.27	116.80	105.20
1	AA	2528	G	OP2-P-O3'	5.27	116.80	105.20
32	A8	61	LEU	CB-CG-CD1	-5.27	102.04	111.00
1	CA	740	U	OP1-P-O3'	5.27	116.80	105.20
1	AA	500	G	C5-C6-O6	5.27	131.76	128.60
34	BA	767	A	N9-C4-C5	5.27	107.91	105.80
1	CA	486	C	OP2-P-O3'	5.27	116.80	105.20
1	AA	835	A	C8-N9-C4	5.27	107.91	105.80
1	AA	2343	G	C5-C6-N1	5.27	114.14	111.50
1	CA	574	C	O5'-P-OP1	-5.27	100.96	105.70
1	CA	2881	C	N3-C2-O2	-5.27	118.21	121.90
1	AA	1926	G	C6-C5-N7	5.27	133.56	130.40
1	AA	2656	G	C2-N3-C4	-5.27	109.27	111.90
2	AB	76	G	O5'-P-OP2	5.27	117.02	110.70
34	BA	574	A	N9-C4-C5	-5.27	103.69	105.80
1	CA	1239	G	C4-C5-N7	5.27	112.91	110.80
1	AA	136	G	C6-N1-C2	-5.27	121.94	125.10
1	AA	174	U	C2-N3-C4	-5.27	123.84	127.00
34	BA	821	G	OP1-P-OP2	-5.27	111.70	119.60
34	BA	855	G	C8-N9-C4	-5.27	104.29	106.40
1	CA	1244	G	C8-N9-C4	5.27	108.51	106.40
34	DA	758	G	C8-N9-C4	-5.27	104.29	106.40
1	AA	337	C	C2-N1-C1'	-5.26	113.01	118.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	658	A	N9-C4-C5	-5.26	103.69	105.80
1	AA	876	A	N1-C6-N6	5.26	121.76	118.60
1	AA	1329	G	C4-C5-N7	5.26	112.91	110.80
1	AA	1720	U	N3-C2-O2	5.26	125.89	122.20
1	AA	2403	G	C4-N9-C1'	-5.26	119.66	126.50
1	AA	2439	C	OP2-P-O3'	5.26	116.78	105.20
1	AA	2535	G	N1-C2-N3	5.26	127.06	123.90
1	AA	1206	G	C2-N3-C4	5.26	114.53	111.90
1	AA	1329	G	C5-C6-N1	-5.26	108.87	111.50
1	AA	2778	A	C5-N7-C8	5.26	106.53	103.90
2	AB	34	U	N3-C4-C5	-5.26	111.44	114.60
2	AB	49	C	C5-C4-N4	-5.26	116.52	120.20
34	BA	239	U	N1-C2-O2	5.26	126.48	122.80
1	CA	71	A	C2-N3-C4	5.26	113.23	110.60
1	CA	400	G	OP2-P-O3'	5.26	116.78	105.20
1	CA	1299	G	C6-N1-C2	5.26	128.26	125.10
1	CA	1903	G	N3-C4-C5	5.26	131.23	128.60
1	CA	2042	A	C5-C6-N6	5.26	127.91	123.70
34	DA	524	G	N3-C4-C5	5.26	131.23	128.60
34	DA	882	C	N1-C2-O2	5.26	122.06	118.90
1	AA	906	G	C4-N9-C1'	-5.26	119.66	126.50
1	AA	1199	C	OP2-P-O3'	5.26	116.77	105.20
1	AA	1528	U	N1-C2-O2	-5.26	119.12	122.80
1	AA	2219	U	C4-C5-C6	5.26	122.86	119.70
1	AA	2223	C	C2-N1-C1'	5.26	124.59	118.80
1	AA	2548	G	N1-C2-N3	5.26	127.06	123.90
1	AA	2627	U	N3-C4-O4	-5.26	115.72	119.40
1	AA	2745	G	C4-C5-N7	-5.26	108.70	110.80
34	BA	823	G	N3-C4-C5	5.26	131.23	128.60
1	AA	1189	A	C5-C6-N6	5.26	127.91	123.70
1	AA	1622	C	C2-N1-C1'	-5.26	113.02	118.80
1	AA	2835	C	C4-C5-C6	5.26	120.03	117.40
2	AB	8	U	C6-N1-C2	5.26	124.16	121.00
34	BA	916	G	N3-C4-N9	5.26	129.16	126.00
1	CA	1698	A	C8-N9-C1'	-5.26	118.23	127.70
1	AA	36	G	C5-N7-C8	5.26	106.93	104.30
1	AA	1253	C	C6-N1-C1'	-5.26	114.49	120.80
1	AA	2086	C	OP2-P-O3'	5.26	116.76	105.20
1	AA	2557	G	C5-C6-O6	-5.26	125.45	128.60
1	AA	2677	A	N9-C4-C5	-5.26	103.70	105.80
13	CP	21	ARG	NE-CZ-NH1	-5.26	117.67	120.30
1	AA	613	A	C8-N9-C4	-5.25	103.70	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1058	U	C2-N3-C4	-5.25	123.85	127.00
1	CA	223	A	C8-N9-C4	-5.25	103.70	105.80
34	DA	687	A	C8-N9-C4	-5.25	103.70	105.80
1	AA	922	G	C8-N9-C4	5.25	108.50	106.40
1	AA	1484	U	N1-C2-O2	-5.25	119.12	122.80
1	AA	2072	C	N1-C2-O2	-5.25	115.75	118.90
34	BA	328	C	OP1-P-O3'	5.25	116.76	105.20
56	BW	47	U	C5-C6-N1	5.25	125.33	122.70
1	CA	797	C	O5'-P-OP2	-5.25	100.97	105.70
1	CA	797	C	N1-C2-O2	-5.25	115.75	118.90
1	CA	2058	A	OP2-P-O3'	5.25	116.76	105.20
1	AA	361	C	OP2-P-O3'	5.25	116.75	105.20
1	AA	717	A	O4'-C1'-N9	-5.25	104.00	108.20
1	AA	954	C	O5'-P-OP2	-5.25	100.97	105.70
1	AA	1963	C	C2-N1-C1'	5.25	124.58	118.80
1	AA	2002	G	C2-N3-C4	-5.25	109.27	111.90
1	AA	2048	C	N3-C4-C5	-5.25	119.80	121.90
1	AA	2634	C	C5-C6-N1	-5.25	118.38	121.00
1	AA	2838	C	N3-C4-C5	5.25	124.00	121.90
15	AR	12	ARG	NE-CZ-NH1	-5.25	117.67	120.30
1	CA	1689	A	N1-C6-N6	-5.25	115.45	118.60
34	DA	798	G	C8-N9-C4	-5.25	104.30	106.40
1	AA	75	C	N3-C4-N4	5.25	121.67	118.00
1	AA	1410	G	C2-N3-C4	5.25	114.53	111.90
1	AA	2114	U	OP1-P-O3'	5.25	116.75	105.20
34	BA	1226	C	N3-C4-C5	-5.25	119.80	121.90
1	AA	1192	C	C5-C6-N1	-5.25	118.38	121.00
1	AA	1845	G	C5-C6-N1	5.25	114.12	111.50
1	AA	1967	G	C2-N3-C4	-5.25	109.28	111.90
1	CA	277	C	C6-N1-C2	-5.25	118.20	120.30
1	CA	1559	G	C4-N9-C1'	-5.25	119.68	126.50
34	DA	492	G	C8-N9-C4	-5.25	104.30	106.40
1	AA	87	G	C8-N9-C4	-5.25	104.30	106.40
1	AA	738	C	O5'-P-OP1	5.25	117.00	110.70
1	AA	1264	G	C2-N3-C4	-5.25	109.28	111.90
1	AA	1377	A	C4-C5-N7	-5.25	108.08	110.70
1	AA	1862	G	C8-N9-C4	-5.25	104.30	106.40
1	AA	1958	A	C5-C6-N6	-5.25	119.50	123.70
1	AA	2625	U	OP2-P-O3'	5.25	116.74	105.20
1	AA	2762	A	N3-C4-C5	-5.25	123.13	126.80
34	BA	519	C	C5-C6-N1	-5.25	118.38	121.00
51	BR	85	LEU	CA-CB-CG	5.25	127.37	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1313	U	N3-C4-O4	5.25	123.07	119.40
34	DA	863	U	C6-N1-C1'	5.25	128.54	121.20
1	AA	1442	U	O5'-P-OP1	-5.25	100.98	105.70
1	CA	1602	U	N1-C2-N3	5.25	118.05	114.90
1	AA	36	G	C6-C5-N7	5.24	133.55	130.40
1	AA	841	G	N3-C4-N9	5.24	129.15	126.00
1	AA	2556	G	O5'-P-OP2	5.24	116.99	110.70
34	BA	562	C	C5-C4-N4	-5.24	116.53	120.20
1	CA	1204	A	N7-C8-N9	5.24	116.42	113.80
34	DA	150	C	C6-N1-C2	-5.24	118.20	120.30
1	AA	1100	A	C6-C5-N7	-5.24	128.63	132.30
1	CA	195	A	O4'-C1'-N9	5.24	112.39	108.20
1	CA	1957	C	C5-C6-N1	5.24	123.62	121.00
1	AA	2298	A	N9-C1'-C2'	5.24	120.81	114.00
1	AA	2386	C	N3-C2-O2	5.24	125.57	121.90
1	AA	2471	A	C4-C5-N7	-5.24	108.08	110.70
2	AB	99	G	N3-C4-N9	5.24	129.14	126.00
34	BA	34	C	C6-N1-C2	5.24	122.40	120.30
34	BA	1468	A	N9-C4-C5	-5.24	103.70	105.80
1	CA	2734	A	C8-N9-C4	-5.24	103.70	105.80
1	AA	214	A	O5'-P-OP1	5.24	116.98	110.70
1	AA	1317	G	N1-C6-O6	5.24	123.04	119.90
1	AA	2426	G	N1-C6-O6	-5.24	116.76	119.90
34	BA	878	G	C4-C5-N7	5.24	112.90	110.80
1	CA	757	U	N3-C2-O2	-5.24	118.53	122.20
1	CA	1204	A	N1-C2-N3	5.24	131.92	129.30
1	AA	2011	G	N1-C6-O6	5.24	123.04	119.90
1	CA	1022	G	N1-C6-O6	-5.24	116.76	119.90
1	CA	2438	U	C2-N3-C4	-5.24	123.86	127.00
1	AA	548	C	C6-N1-C1'	5.24	127.08	120.80
1	AA	566	C	C2-N1-C1'	5.24	124.56	118.80
1	AA	1027	A	C5-N7-C8	5.24	106.52	103.90
1	AA	1231	G	C8-N9-C4	5.24	108.49	106.40
1	AA	1252	C	C5-C6-N1	-5.24	118.38	121.00
1	AA	1321	A	C4-C5-N7	-5.24	108.08	110.70
1	AA	1434	G	OP1-P-OP2	5.24	127.45	119.60
1	AA	1454	C	N3-C4-C5	-5.24	119.81	121.90
1	AA	2611	G	N3-C4-N9	5.24	129.14	126.00
1	AA	2659	U	C4-C5-C6	-5.24	116.56	119.70
34	BA	771	G	C6-N1-C2	5.24	128.24	125.10
1	CA	186	G	N1-C6-O6	5.24	123.04	119.90
1	CA	210	C	N3-C4-N4	-5.24	114.33	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2820	A	C8-N9-C1'	-5.24	118.28	127.70
34	DA	115	G	O5'-P-OP2	-5.24	100.99	105.70
34	DA	858	G	C6-C5-N7	-5.24	127.26	130.40
1	CA	1684	C	N1-C2-O2	-5.23	115.76	118.90
1	CA	2429	G	N9-C4-C5	5.23	107.49	105.40
34	DA	10	A	C8-N9-C4	-5.23	103.71	105.80
1	AA	279	G	N7-C8-N9	5.23	115.72	113.10
1	AA	2263	G	C6-N1-C2	-5.23	121.96	125.10
1	AA	2636	G	N3-C4-N9	5.23	129.14	126.00
21	AX	76	ARG	NE-CZ-NH2	-5.23	117.68	120.30
23	AZ	5	LEU	CA-CB-CG	5.23	127.33	115.30
34	BA	103	C	C6-N1-C2	-5.23	118.21	120.30
1	CA	923	C	C5-C6-N1	5.23	123.62	121.00
1	CA	2456	C	C2-N3-C4	-5.23	117.28	119.90
1	AA	455	A	C6-C5-N7	-5.23	128.64	132.30
1	AA	460	C	P-O3'-C3'	5.23	125.98	119.70
1	AA	1783	C	C2-N1-C1'	-5.23	113.05	118.80
1	AA	2267	G	C5-C6-O6	-5.23	125.46	128.60
1	AA	2455	C	N3-C2-O2	-5.23	118.24	121.90
1	AA	2478	C	N3-C4-C5	-5.23	119.81	121.90
1	AA	2528	G	C5-N7-C8	5.23	106.92	104.30
1	AA	2826	C	O5'-P-OP2	-5.23	100.99	105.70
34	BA	28	G	N1-C2-N2	5.23	120.91	116.20
34	BA	505	G	C8-N9-C4	-5.23	104.31	106.40
34	BA	611	A	C8-N9-C4	5.23	107.89	105.80
1	CA	513	A	N7-C8-N9	5.23	116.42	113.80
1	CA	1637	A	C8-N9-C4	-5.23	103.71	105.80
31	C7	33	ARG	NE-CZ-NH2	5.23	122.92	120.30
1	AA	1220	U	P-O3'-C3'	5.23	125.97	119.70
1	AA	1972	G	O5'-P-OP2	5.23	116.97	110.70
1	AA	2297	C	O5'-P-OP2	-5.23	100.99	105.70
1	AA	2374	G	C6-N1-C2	-5.23	121.96	125.10
34	BA	804	U	C2-N1-C1'	-5.23	111.43	117.70
1	CA	2568	C	O5'-P-OP1	5.23	116.97	110.70
1	CA	2879	C	N1-C2-O2	-5.23	115.76	118.90
34	DA	1077	G	C8-N9-C4	5.23	108.49	106.40
1	AA	549	U	C6-N1-C2	5.23	124.14	121.00
1	AA	1397	C	C6-N1-C2	-5.23	118.21	120.30
1	AA	1909	C	N3-C4-N4	5.23	121.66	118.00
1	AA	2737	C	N1-C2-O2	-5.23	115.76	118.90
1	CA	530	G	N3-C4-C5	5.23	131.21	128.60
1	CA	965	C	N1-C2-O2	5.23	122.04	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1238	G	C4-C5-N7	5.23	112.89	110.80
1	AA	326	C	C6-N1-C2	-5.23	118.21	120.30
1	AA	2100	C	O5'-P-OP1	-5.23	101.00	105.70
1	AA	1328	U	C6-N1-C2	5.22	124.14	121.00
1	AA	1604	C	C2-N3-C4	-5.22	117.29	119.90
1	AA	2359	C	C5-C6-N1	5.22	123.61	121.00
1	AA	2619	G	N3-C4-N9	5.22	129.13	126.00
1	AA	2852	G	N7-C8-N9	-5.22	110.49	113.10
34	BA	311	C	N3-C4-C5	5.22	123.99	121.90
56	BW	70	G	N3-C4-N9	-5.22	122.86	126.00
1	CA	183	C	C6-N1-C2	5.22	122.39	120.30
1	CA	1421	G	C8-N9-C4	-5.22	104.31	106.40
1	CA	1783	A	C5-C6-N6	5.22	127.88	123.70
1	CA	1901	A	C6-N1-C2	-5.22	115.47	118.60
1	CA	2522	U	OP2-P-O3'	5.22	116.69	105.20
1	AA	20	C	OP2-P-O3'	5.22	116.69	105.20
1	AA	474	U	C5-C6-N1	-5.22	120.09	122.70
34	BA	819	A	OP1-P-O3'	5.22	116.69	105.20
34	DA	861	G	N9-C4-C5	5.22	107.49	105.40
1	AA	773	G	N1-C6-O6	-5.22	116.77	119.90
1	AA	1266	C	OP1-P-OP2	-5.22	111.77	119.60
1	AA	1736	A	C5-C6-N1	5.22	120.31	117.70
34	BA	1064	G	N3-C2-N2	-5.22	116.25	119.90
1	CA	2274	A	C4-C5-N7	-5.22	108.09	110.70
1	CA	2401	U	O5'-P-OP1	-5.22	101.00	105.70
1	CA	2524	G	O5'-P-OP1	5.22	116.97	110.70
1	AA	16	G	N1-C2-N2	-5.22	111.50	116.20
1	AA	627	G	OP1-P-O3'	5.22	116.68	105.20
1	AA	2762	A	C4-C5-C6	5.22	119.61	117.00
1	AA	2786	C	C6-N1-C2	5.22	122.39	120.30
1	AA	2848	G	N7-C8-N9	-5.22	110.49	113.10
34	BA	330	C	N1-C2-O2	5.22	122.03	118.90
34	BA	756	C	N3-C2-O2	5.22	125.55	121.90
1	CA	1217	C	OP2-P-O3'	5.22	116.68	105.20
1	CA	1351	C	C2-N3-C4	-5.22	117.29	119.90
1	CA	2085	C	C4-C5-C6	5.22	120.01	117.40
1	AA	877	G	C2-N3-C4	-5.22	109.29	111.90
1	AA	1431	G	O5'-P-OP2	5.22	116.96	110.70
2	AB	5	C	N1-C2-N3	5.22	122.85	119.20
2	AB	92	C	N3-C4-N4	5.22	121.65	118.00
1	CA	374	A	O5'-P-OP1	-5.22	101.00	105.70
1	CA	1600	C	O5'-P-OP2	-5.22	101.00	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2487	G	O5'-P-OP2	-5.22	101.00	105.70
34	DA	718	G	N1-C2-N2	5.22	120.90	116.20
1	AA	131	C	N1-C2-O2	5.22	122.03	118.90
1	AA	991	G	N3-C4-N9	-5.22	122.87	126.00
34	BA	558	G	C6-C5-N7	-5.22	127.27	130.40
34	BA	1383	C	C5-C6-N1	5.22	123.61	121.00
1	CA	1293	C	C2-N3-C4	-5.22	117.29	119.90
1	CA	2489	G	C6-C5-N7	-5.22	127.27	130.40
2	CB	105	A	C8-N9-C4	5.22	107.89	105.80
1	AA	478	G	N9-C4-C5	5.21	107.49	105.40
1	AA	561	A	N7-C8-N9	-5.21	111.19	113.80
1	AA	1745	A	C6-N1-C2	-5.21	115.47	118.60
1	AA	2028	C	N1-C2-O2	-5.21	115.77	118.90
1	AA	2275	C	C2-N1-C1'	-5.21	113.06	118.80
1	AA	2554	A	N9-C4-C5	-5.21	103.71	105.80
34	BA	1036	G	C4-N9-C1'	5.21	133.28	126.50
1	CA	495	G	O5'-P-OP2	-5.21	101.01	105.70
1	CA	2265	U	C5-C6-N1	5.21	125.31	122.70
1	CA	2335	A	O4'-C1'-N9	5.21	112.37	108.20
1	CA	2569	G	C4-C5-N7	5.21	112.89	110.80
34	DA	175	C	N3-C2-O2	-5.21	118.25	121.90
34	DA	1513	A	N3-C4-C5	5.21	130.45	126.80
1	AA	471	C	C2-N3-C4	-5.21	117.29	119.90
1	AA	789	G	O5'-P-OP2	5.21	116.95	110.70
1	AA	862	C	C6-N1-C2	-5.21	118.22	120.30
34	BA	905	U	O5'-P-OP1	-5.21	101.01	105.70
1	CA	114	U	C2-N1-C1'	5.21	123.95	117.70
1	CA	328	U	OP1-P-O3'	5.21	116.67	105.20
1	CA	524	U	OP1-P-OP2	5.21	127.42	119.60
1	AA	750	U	C2-N1-C1'	5.21	123.95	117.70
1	AA	1455	C	OP2-P-O3'	5.21	116.67	105.20
1	AA	1766	G	N1-C6-O6	5.21	123.03	119.90
1	CA	1796	U	C2-N3-C4	-5.21	123.87	127.00
1	CA	2463	C	N3-C4-N4	-5.21	114.35	118.00
1	CA	2588	G	C5-C6-N1	-5.21	108.89	111.50
1	CA	2602	A	O4'-C1'-N9	5.21	112.37	108.20
1	CA	2655	G	C8-N9-C4	5.21	108.48	106.40
1	AA	308	U	C5-C6-N1	5.21	125.31	122.70
1	AA	556	C	C2-N3-C4	-5.21	117.30	119.90
1	AA	2559	U	N3-C4-C5	-5.21	111.47	114.60
34	BA	1404	C	N3-C4-C5	5.21	123.98	121.90
1	CA	2067	G	N1-C6-O6	-5.21	116.77	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	208	G	C2-N3-C4	5.21	114.50	111.90
1	AA	627	G	O5'-P-OP2	-5.21	101.01	105.70
1	AA	796	C	OP1-P-OP2	5.21	127.41	119.60
1	AA	914	C	C5-C6-N1	-5.21	118.40	121.00
1	AA	1233	U	OP2-P-O3'	5.21	116.66	105.20
1	AA	2423	A	N7-C8-N9	-5.21	111.20	113.80
1	AA	2632	C	C2-N1-C1'	-5.21	113.07	118.80
34	BA	302	G	OP2-P-O3'	5.21	116.66	105.20
56	BW	75	C	C2-N3-C4	-5.21	117.30	119.90
1	CA	450	G	C4-C5-N7	-5.21	108.72	110.80
1	CA	2340	G	N3-C4-N9	5.21	129.12	126.00
1	AA	177	G	C4-C5-N7	-5.21	108.72	110.80
1	AA	1648	U	N3-C4-C5	5.21	117.72	114.60
1	AA	2348	A	O4'-C1'-N9	-5.21	104.04	108.20
34	BA	768	A	C6-N1-C2	-5.21	115.48	118.60
34	BA	895	G	C6-C5-N7	5.21	133.52	130.40
1	CA	1951	U	N3-C2-O2	-5.21	118.56	122.20
1	CA	2016	U	C2-N3-C4	5.21	130.12	127.00
1	AA	720	C	OP2-P-O3'	5.21	116.65	105.20
1	AA	1231	G	N9-C4-C5	-5.21	103.32	105.40
1	AA	1473	A	N1-C2-N3	5.21	131.90	129.30
1	AA	2635	G	C8-N9-C4	-5.21	104.32	106.40
1	AA	786	G	N3-C4-N9	5.20	129.12	126.00
1	AA	817	G	C5-C6-N1	5.20	114.10	111.50
1	AA	1725	G	C4-C5-C6	5.20	121.92	118.80
1	AA	2834	C	C4-C5-C6	5.20	120.00	117.40
1	CA	125	G	O4'-C1'-N9	-5.20	104.04	108.20
1	CA	799	G	N3-C4-N9	5.20	129.12	126.00
1	CA	1368	G	OP2-P-O3'	5.20	116.65	105.20
1	CA	1651	G	N1-C6-O6	-5.20	116.78	119.90
1	CA	2406	U	O5'-P-OP1	-5.20	101.02	105.70
1	AA	560	C	N3-C4-C5	5.20	123.98	121.90
1	AA	1716	A	O4'-C1'-N9	5.20	112.36	108.20
34	BA	1502	A	O5'-P-OP2	-5.20	101.02	105.70
1	CA	2077	A	C5-C6-N6	-5.20	119.54	123.70
1	AA	775	G	C5-C6-O6	5.20	131.72	128.60
1	AA	915	U	O5'-P-OP1	-5.20	101.02	105.70
1	AA	1052	C	C6-N1-C2	5.20	122.38	120.30
1	AA	1251	G	N7-C8-N9	-5.20	110.50	113.10
1	AA	1253	C	C2-N1-C1'	5.20	124.52	118.80
1	AA	1619	A	O5'-P-OP2	-5.20	101.02	105.70
1	AA	1817	A	C6-N1-C2	5.20	121.72	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2097	U	C2-N3-C4	-5.20	123.88	127.00
1	AA	2611	G	C8-N9-C4	5.20	108.48	106.40
1	AA	2720	G	OP2-P-O3'	5.20	116.64	105.20
1	AA	2757	G	N1-C6-O6	5.20	123.02	119.90
1	AA	2833	A	C5-N7-C8	5.20	106.50	103.90
1	CA	252	G	C6-C5-N7	5.20	133.52	130.40
1	CA	1956	U	N3-C4-C5	5.20	117.72	114.60
1	CA	2342	C	C5-C6-N1	5.20	123.60	121.00
56	DW	43	C	C2-N1-C1'	5.20	124.52	118.80
1	AA	174	U	C5-C4-O4	-5.20	122.78	125.90
1	AA	608	G	OP1-P-O3'	5.20	116.64	105.20
1	AA	1296	G	C8-N9-C4	-5.20	104.32	106.40
1	AA	1962	U	C4-C5-C6	5.20	122.82	119.70
1	AA	2108	U	OP2-P-O3'	5.20	116.64	105.20
1	AA	2500	A	C5-C6-N6	5.20	127.86	123.70
1	CA	265	A	C8-N9-C4	-5.20	103.72	105.80
1	CA	1021	A	C5-N7-C8	-5.20	101.30	103.90
1	CA	1274	A	N1-C6-N6	5.20	121.72	118.60
1	CA	2253	G	N3-C2-N2	-5.20	116.26	119.90
56	DW	73	A	N7-C8-N9	5.20	116.40	113.80
1	AA	2050	U	N3-C4-O4	-5.20	115.76	119.40
1	AA	2063	U	OP2-P-O3'	5.20	116.63	105.20
1	AA	2626	A	N7-C8-N9	5.20	116.40	113.80
1	AA	2707	C	OP2-P-O3'	5.20	116.63	105.20
1	AA	2833	A	C6-N1-C2	-5.20	115.48	118.60
34	BA	275	G	C8-N9-C4	-5.20	104.32	106.40
1	CA	2049	G	N1-C2-N2	-5.20	111.52	116.20
1	CA	2883	A	O4'-C1'-N9	5.20	112.36	108.20
34	DA	758	G	N1-C6-O6	-5.20	116.78	119.90
1	AA	354	A	N1-C2-N3	5.20	131.90	129.30
1	AA	725	C	N1-C2-N3	5.20	122.84	119.20
1	AA	2074	G	OP1-P-O3'	5.20	116.63	105.20
1	AA	2353	G	N9-C4-C5	5.20	107.48	105.40
1	AA	2401	G	N3-C4-N9	-5.20	122.88	126.00
1	AA	2408	G	N1-C6-O6	-5.20	116.78	119.90
1	AA	2661	U	OP2-P-O3'	5.20	116.63	105.20
1	AA	2725	A	OP1-P-OP2	5.20	127.39	119.60
2	AB	102	A	C8-N9-C4	5.20	107.88	105.80
4	AD	222	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	CA	2045	C	N3-C4-C5	5.20	123.98	121.90
1	CA	2618	G	C5-N7-C8	5.20	106.90	104.30
1	AA	2520	G	C2-N3-C4	5.19	114.50	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1320	C	N3-C2-O2	5.19	125.54	121.90
5	CE	103	ASP	N-CA-C	-5.19	96.98	111.00
1	AA	1472	G	C6-C5-N7	-5.19	127.28	130.40
1	AA	1685	C	N1-C1'-C2'	-5.19	106.29	112.00
1	AA	2088	C	C5-C4-N4	5.19	123.83	120.20
1	AA	2189	U	N1-C2-O2	5.19	126.44	122.80
1	AA	2714	U	O4'-C1'-N1	-5.19	104.05	108.20
13	AP	55	ARG	NE-CZ-NH2	-5.19	117.70	120.30
23	AZ	77	ASP	CB-CG-OD2	-5.19	113.63	118.30
1	CA	300	A	C4-C5-C6	-5.19	114.40	117.00
1	CA	448	U	N1-C2-O2	5.19	126.43	122.80
1	CA	798	G	N9-C4-C5	5.19	107.48	105.40
1	CA	1598	C	OP1-P-O3'	5.19	116.62	105.20
1	CA	2767	C	C2-N1-C1'	5.19	124.51	118.80
1	AA	591	U	C2-N3-C4	-5.19	123.89	127.00
1	AA	657	A	OP1-P-O3'	5.19	116.62	105.20
1	AA	2869	G	C5-C6-O6	5.19	131.71	128.60
1	CA	2235	G	N3-C2-N2	-5.19	116.27	119.90
1	AA	524	U	N3-C2-O2	-5.19	118.57	122.20
1	AA	844	C	C4-C5-C6	5.19	120.00	117.40
1	AA	1376	C	N3-C4-N4	5.19	121.63	118.00
1	AA	2269	U	N1-C2-N3	5.19	118.01	114.90
1	AA	533	G	N9-C4-C5	5.19	107.47	105.40
1	AA	556	C	N3-C4-C5	5.19	123.97	121.90
1	AA	730	C	C4-C5-C6	5.19	119.99	117.40
1	AA	1484	U	N3-C4-O4	5.19	123.03	119.40
1	AA	1739	U	C5-C6-N1	-5.19	120.11	122.70
1	AA	2688	C	C6-N1-C2	5.19	122.38	120.30
1	AA	2858	G	C5-C6-N1	5.19	114.09	111.50
1	CA	1346	G	C5-C6-O6	5.19	131.71	128.60
1	CA	2255	G	C8-N9-C4	5.19	108.47	106.40
34	DA	303	A	N1-C6-N6	-5.19	115.49	118.60
34	DA	560	U	N3-C2-O2	-5.19	118.57	122.20
34	DA	731	G	OP2-P-O3'	5.19	116.61	105.20
34	DA	813	U	N3-C2-O2	5.19	125.83	122.20
1	AA	483	A	N1-C6-N6	5.19	121.71	118.60
1	AA	1193	C	C6-N1-C1'	5.19	127.02	120.80
1	AA	2627	U	C2-N3-C4	-5.19	123.89	127.00
13	AP	60	MET	CG-SD-CE	5.19	108.50	100.20
34	BA	923	A	O5'-P-OP1	-5.19	101.03	105.70
1	CA	705	A	O5'-P-OP2	-5.19	101.03	105.70
1	CA	795	C	C5-C6-N1	-5.19	118.41	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	798	A	C5-C6-N6	-5.18	119.55	123.70
1	AA	1264	G	O5'-P-OP1	5.18	116.92	110.70
1	AA	2484	G	C5-C6-N1	5.18	114.09	111.50
34	BA	1394	A	OP1-P-O3'	5.18	116.61	105.20
1	CA	2553	G	N1-C2-N2	-5.18	111.53	116.20
1	CA	2580	U	N3-C4-C5	5.18	117.71	114.60
1	AA	261	A	OP2-P-O3'	5.18	116.60	105.20
1	AA	473	A	N1-C2-N3	5.18	131.89	129.30
1	AA	1571	G	O5'-P-OP1	-5.18	101.03	105.70
1	AA	2849	G	N7-C8-N9	-5.18	110.51	113.10
34	BA	778	G	N1-C2-N2	-5.18	111.54	116.20
1	CA	350	U	N3-C2-O2	-5.18	118.57	122.20
1	CA	767	U	N1-C2-O2	5.18	126.43	122.80
1	CA	1905	C	C2-N3-C4	5.18	122.49	119.90
1	CA	2342	C	O5'-P-OP2	5.18	116.92	110.70
34	DA	716	A	N7-C8-N9	5.18	116.39	113.80
1	AA	1416	C	N1-C2-O2	-5.18	115.79	118.90
1	AA	1822	A	OP2-P-O3'	5.18	116.60	105.20
1	AA	712	C	N1-C2-O2	5.18	122.01	118.90
1	AA	835	A	C6-N1-C2	5.18	121.71	118.60
1	AA	1394	G	N9-C4-C5	-5.18	103.33	105.40
1	AA	2383	G	C6-N1-C2	-5.18	121.99	125.10
1	AA	2512	U	C2-N3-C4	-5.18	123.89	127.00
1	CA	1782	C	C2-N3-C4	5.18	122.49	119.90
1	CA	1999	C	N3-C4-C5	5.18	123.97	121.90
34	DA	665	A	O5'-P-OP2	-5.18	101.04	105.70
1	AA	46	C	C5-C4-N4	-5.18	116.58	120.20
1	AA	2374	G	C8-N9-C4	5.18	108.47	106.40
1	CA	2229	C	N3-C4-C5	-5.18	119.83	121.90
34	DA	7	G	C4-N9-C1'	-5.18	119.77	126.50
1	AA	952	G	N1-C6-O6	-5.18	116.79	119.90
1	AA	1080	G	N7-C8-N9	-5.18	110.51	113.10
34	BA	378	G	OP1-P-O3'	5.18	116.59	105.20
1	CA	1653	G	N1-C6-O6	-5.18	116.79	119.90
1	CA	2012	G	N7-C8-N9	5.18	115.69	113.10
1	CA	2444	G	N3-C2-N2	5.18	123.52	119.90
1	CA	2525	G	OP1-P-O3'	-5.18	93.81	105.20
34	DA	567	G	C8-N9-C4	-5.18	104.33	106.40
1	AA	100	G	N1-C6-O6	5.17	123.00	119.90
1	AA	183	G	C5-C6-O6	5.17	131.71	128.60
1	AA	546	G	C6-N1-C2	-5.17	122.00	125.10
1	AA	587	C	N3-C4-C5	-5.17	119.83	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	634	C	C2-N3-C4	-5.17	117.31	119.90
1	AA	640	A	C5-N7-C8	-5.17	101.31	103.90
1	AA	908	A	C4-C5-C6	-5.17	114.41	117.00
1	AA	982	U	N3-C2-O2	5.17	125.82	122.20
1	AA	1477	U	C2-N3-C4	5.17	130.10	127.00
1	AA	1664	A	O5'-P-OP2	5.17	116.91	110.70
1	AA	2387	G	OP1-P-OP2	5.17	127.36	119.60
1	AA	2454	C	N3-C4-N4	5.17	121.62	118.00
34	BA	882	C	C6-N1-C2	5.17	122.37	120.30
1	AA	2830	A	C5-N7-C8	-5.17	101.31	103.90
1	AA	2859	U	C2-N3-C4	-5.17	123.90	127.00
41	BH	112	LEU	CA-CB-CG	5.17	127.20	115.30
1	CA	946	G	C2-N3-C4	-5.17	109.31	111.90
1	CA	2893	G	C8-N9-C1'	-5.17	120.28	127.00
34	DA	105	G	N3-C4-N9	5.17	129.10	126.00
1	AA	776	G	OP2-P-O3'	5.17	116.58	105.20
1	AA	1241	C	N1-C2-O2	-5.17	115.80	118.90
1	AA	1416	C	N3-C4-N4	5.17	121.62	118.00
1	AA	1753	U	N1-C2-O2	5.17	126.42	122.80
1	AA	2256	U	C4-C5-C6	5.17	122.80	119.70
1	AA	2554	A	C8-N9-C4	5.17	107.87	105.80
1	AA	2559	U	C5-C6-N1	-5.17	120.11	122.70
1	AA	2832	G	N3-C4-N9	5.17	129.10	126.00
34	BA	879	C	OP2-P-O3'	5.17	116.58	105.20
1	CA	1441	G	C4-C5-N7	-5.17	108.73	110.80
34	DA	434	U	C6-N1-C2	-5.17	117.90	121.00
1	AA	554	A	C5'-C4'-C3'	-5.17	107.73	116.00
1	AA	655	G	O5'-P-OP2	-5.17	101.05	105.70
1	AA	1752	G	N1-C2-N2	-5.17	111.55	116.20
1	AA	167	G	N7-C8-N9	-5.17	110.52	113.10
1	AA	1501	U	N3-C2-O2	5.17	125.82	122.20
1	CA	741	G	N3-C4-N9	-5.17	122.90	126.00
1	CA	1460	A	O4'-C1'-N9	5.17	112.33	108.20
1	CA	2519	U	N3-C2-O2	5.17	125.82	122.20
1	CA	2606	C	OP2-P-O3'	5.17	116.57	105.20
1	AA	1678	A	C5-C6-N1	-5.17	115.12	117.70
1	AA	2235	G	OP2-P-O3'	5.17	116.57	105.20
1	AA	2594	G	C5-C6-N1	5.17	114.08	111.50
2	AB	38	C	C5-C6-N1	-5.17	118.42	121.00
34	BA	744	C	N3-C4-C5	5.17	123.97	121.90
1	CA	829	A	OP1-P-OP2	5.17	127.35	119.60
1	AA	449	A	OP1-P-OP2	-5.16	111.85	119.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	916	G	OP1-P-OP2	5.16	127.34	119.60
1	AA	2780	C	O5'-P-OP2	-5.16	101.05	105.70
2	AB	106	G	C2-N3-C4	5.16	114.48	111.90
4	AD	61	LEU	CA-CB-CG	5.16	127.18	115.30
13	AP	18	ARG	NE-CZ-NH1	5.16	122.88	120.30
34	BA	1519	A	C4-C5-C6	5.16	119.58	117.00
1	CA	764	A	O4'-C1'-N9	5.16	112.33	108.20
1	CA	1790	C	OP1-P-O3'	5.16	116.56	105.20
1	CA	2586	C	N1-C2-O2	5.16	122.00	118.90
1	CA	2618	G	C5-C6-N1	5.16	114.08	111.50
1	CA	2641	G	O5'-P-OP2	-5.16	101.05	105.70
1	CA	2673	G	C8-N9-C1'	-5.16	120.29	127.00
1	AA	188	A	OP2-P-O3'	5.16	116.56	105.20
1	AA	872	C	C4-C5-C6	5.16	119.98	117.40
1	AA	2000	A	C8-N9-C4	5.16	107.86	105.80
1	AA	911	G	O5'-P-OP1	5.16	116.89	110.70
1	AA	1071	G	C6-N1-C2	-5.16	122.00	125.10
1	AA	1423	G	C2-N3-C4	5.16	114.48	111.90
1	AA	2439	C	C5-C6-N1	-5.16	118.42	121.00
1	AA	2544	G	C5-N7-C8	-5.16	101.72	104.30
56	BY	74	C	C6-N1-C2	5.16	122.36	120.30
1	AA	150	C	C2-N3-C4	-5.16	117.32	119.90
1	AA	821	A	N7-C8-N9	5.16	116.38	113.80
1	AA	959	U	O5'-P-OP2	-5.16	101.06	105.70
1	AA	1017	G	OP1-P-O3'	5.16	116.55	105.20
1	AA	1273	G	C6-N1-C2	-5.16	122.00	125.10
1	AA	1502	G	N1-C6-O6	-5.16	116.81	119.90
1	AA	2013	U	N3-C4-O4	-5.16	115.79	119.40
1	AA	2487	C	N3-C2-O2	-5.16	118.29	121.90
1	CA	52	A	O5'-P-OP1	-5.16	101.06	105.70
1	CA	1721	G	C6-C5-N7	-5.16	127.31	130.40
1	CA	2050	C	C2-N3-C4	-5.16	117.32	119.90
34	DA	1528	U	C5-C6-N1	-5.16	120.12	122.70
1	AA	2011	G	C2-N3-C4	-5.16	109.32	111.90
1	AA	2050	U	N1-C2-O2	5.16	126.41	122.80
1	AA	2715	C	N1-C2-N3	5.16	122.81	119.20
56	BW	16	U	C2-N1-C1'	-5.16	111.51	117.70
1	AA	26	G	N9-C4-C5	-5.16	103.34	105.40
1	AA	478	G	N3-C2-N2	-5.16	116.29	119.90
1	AA	489	G	OP1-P-OP2	-5.16	111.87	119.60
1	AA	1422	C	N3-C2-O2	-5.16	118.29	121.90
1	AA	2011	G	N3-C4-C5	5.16	131.18	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2074	G	C6-C5-N7	5.16	133.49	130.40
1	AA	2113	U	C2-N3-C4	-5.16	123.91	127.00
1	CA	414	C	C4-C5-C6	5.16	119.98	117.40
1	CA	616	G	O5'-P-OP2	-5.16	101.06	105.70
1	CA	1645	G	N1-C2-N2	-5.16	111.56	116.20
1	CA	2393	A	N9-C4-C5	5.16	107.86	105.80
1	AA	325	G	C4-N9-C1'	-5.15	119.80	126.50
1	AA	874	U	OP1-P-OP2	5.15	127.33	119.60
1	AA	2462	A	C6-N1-C2	5.15	121.69	118.60
34	BA	784	C	N3-C4-C5	5.15	123.96	121.90
1	CA	315	G	O5'-P-OP1	5.15	116.89	110.70
1	AA	81	G	C4-C5-N7	5.15	112.86	110.80
1	AA	1067	A	C8-N9-C4	-5.15	103.74	105.80
1	CA	514	A	N1-C6-N6	5.15	121.69	118.60
1	CA	665	C	C6-N1-C2	-5.15	118.24	120.30
6	CF	74	ARG	NE-CZ-NH2	5.15	122.88	120.30
34	DA	7	G	C8-N9-C1'	5.15	133.70	127.00
57	DZ	216	LEU	CA-CB-CG	5.15	127.15	115.30
1	AA	277	G	N9-C4-C5	5.15	107.46	105.40
1	AA	1013	G	C5-C6-N1	5.15	114.08	111.50
1	AA	1266	C	C4-C5-C6	5.15	119.98	117.40
1	AA	1403	U	N3-C2-O2	-5.15	118.59	122.20
1	AA	2294	G	C5-N7-C8	-5.15	101.72	104.30
1	AA	2549	U	N3-C4-O4	-5.15	115.80	119.40
1	CA	483	A	O5'-P-OP1	-5.15	101.06	105.70
1	CA	2435	A	C5-N7-C8	-5.15	101.32	103.90
1	CA	2568	C	C6-N1-C2	5.15	122.36	120.30
1	AA	438	G	C8-N9-C4	-5.15	104.34	106.40
1	AA	2331	G	C4-C5-C6	-5.15	115.71	118.80
1	CA	1599	C	N3-C4-N4	-5.15	114.40	118.00
1	CA	2510	C	C4-C5-C6	5.15	119.97	117.40
1	AA	243	G	C6-N1-C2	-5.15	122.01	125.10
1	AA	435	G	OP2-P-O3'	5.15	116.53	105.20
1	AA	791	G	N1-C2-N2	-5.15	111.57	116.20
1	AA	855	G	N1-C2-N2	-5.15	111.57	116.20
1	AA	889	G	N3-C4-N9	-5.15	122.91	126.00
1	AA	1082	G	C4-N9-C1'	-5.15	119.81	126.50
1	AA	2464	C	C2-N3-C4	-5.15	117.33	119.90
1	AA	2620	G	C8-N9-C4	5.15	108.46	106.40
1	AA	2710	U	C5-C6-N1	-5.15	120.13	122.70
1	AA	2870	A	OP2-P-O3'	5.15	116.52	105.20
1	CA	578	A	C5-C6-N6	-5.15	119.58	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	2074	U	O5'-P-OP1	-5.15	101.07	105.70
34	DA	60	A	OP1-P-O3'	5.15	116.52	105.20
34	DA	355	C	C6-N1-C2	-5.15	118.24	120.30
1	AA	2434	A	N1-C6-N6	5.15	121.69	118.60
34	BA	644	G	O5'-P-OP2	-5.15	101.07	105.70
1	CA	702	G	O5'-P-OP2	-5.15	101.07	105.70
1	CA	1333	C	O5'-P-OP1	5.15	116.88	110.70
34	DA	882	C	C2-N1-C1'	5.15	124.46	118.80
1	AA	559	U	C5-C4-O4	5.14	128.99	125.90
1	AA	628	C	C6-N1-C2	5.14	122.36	120.30
1	AA	818	G	N9-C4-C5	5.14	107.46	105.40
1	AA	1020	C	C6-N1-C1'	5.14	126.97	120.80
1	AA	1290	G	C8-N9-C4	5.14	108.46	106.40
1	CA	296	C	C6-N1-C2	-5.14	118.24	120.30
1	CA	1296	G	N3-C2-N2	5.14	123.50	119.90
1	CA	1666	G	O5'-P-OP2	5.14	116.88	110.70
1	CA	2069	G	N9-C4-C5	5.14	107.46	105.40
34	DA	30	U	N1-C2-O2	-5.14	119.20	122.80
1	AA	577	U	OP1-P-OP2	-5.14	111.89	119.60
1	AA	847	A	N3-C4-C5	-5.14	123.20	126.80
1	AA	1070	G	C5-C6-O6	-5.14	125.52	128.60
1	AA	1237	G	C5-C6-O6	5.14	131.69	128.60
1	AA	1700	G	N1-C2-N2	-5.14	111.57	116.20
1	AA	1811	A	N9-C4-C5	5.14	107.86	105.80
1	AA	2764	G	O4'-C1'-N9	5.14	112.31	108.20
34	BA	397	A	N1-C2-N3	5.14	131.87	129.30
34	BA	881	G	O5'-P-OP1	5.14	116.87	110.70
34	BA	885	G	O5'-P-OP1	5.14	116.87	110.70
34	BA	1484	C	C5-C6-N1	-5.14	118.43	121.00
1	CA	787	U	C6-N1-C2	-5.14	117.91	121.00
1	CA	952	G	N1-C6-O6	-5.14	116.81	119.90
1	CA	964	C	OP2-P-O3'	5.14	116.51	105.20
1	CA	1779	U	O4'-C1'-N1	5.14	112.31	108.20
1	CA	2877	G	O5'-P-OP2	-5.14	101.07	105.70
34	DA	728	A	O5'-P-OP2	-5.14	101.07	105.70
34	DA	955	U	C5-C6-N1	5.14	125.27	122.70
1	AA	29	U	C5-C6-N1	-5.14	120.13	122.70
1	AA	478	G	C6-C5-N7	5.14	133.48	130.40
1	AA	2052	A	O5'-P-OP2	-5.14	101.07	105.70
1	AA	2081	A	N1-C2-N3	-5.14	126.73	129.30
1	AA	2257	U	O5'-P-OP1	5.14	116.87	110.70
56	BW	71	G	N3-C4-C5	5.14	131.17	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	114	C	C4-C5-C6	5.14	119.97	117.40
1	AA	1185	C	OP2-P-O3'	5.14	116.50	105.20
1	AA	2304	C	C4-C5-C6	5.14	119.97	117.40
1	CA	1323	U	C2-N3-C4	-5.14	123.92	127.00
1	AA	988	U	C5-C4-O4	-5.14	122.82	125.90
1	AA	1605	A	N7-C8-N9	5.14	116.37	113.80
1	AA	2471	A	C5-C6-N1	5.14	120.27	117.70
1	CA	2042	A	C2-N3-C4	-5.14	108.03	110.60
1	CA	2059	A	C8-N9-C4	5.14	107.86	105.80
1	CA	2885	C	N1-C2-O2	5.14	121.98	118.90
34	DA	221	C	N3-C2-O2	-5.14	118.30	121.90
1	AA	85	C	OP2-P-O3'	5.14	116.50	105.20
1	AA	449	A	C8-N9-C4	-5.14	103.75	105.80
1	AA	1610	G	C4-C5-N7	-5.14	108.75	110.80
1	AA	1664	A	C5-C6-N1	5.14	120.27	117.70
1	AA	1697	G	N1-C6-O6	-5.14	116.82	119.90
1	AA	2458	G	OP2-P-O3'	5.14	116.50	105.20
1	AA	2722	C	N1-C2-O2	5.14	121.98	118.90
34	BA	266	G	P-O3'-C3'	5.14	125.86	119.70
34	BA	809	G	C8-N9-C4	-5.14	104.34	106.40
34	BA	1402	C	C6-N1-C2	-5.14	118.25	120.30
1	CA	265	A	C6-C5-N7	-5.14	128.70	132.30
1	CA	2356	C	N3-C2-O2	5.14	125.50	121.90
1	CA	2673	G	C4-N9-C1'	5.14	133.18	126.50
1	CA	2677	G	C8-N9-C4	-5.14	104.35	106.40
1	AA	2437	A	C4-C5-N7	5.13	113.27	110.70
1	AA	2633	A	N3-C4-N9	-5.13	123.29	127.40
1	CA	468	G	C8-N9-C4	5.13	108.45	106.40
1	CA	508	G	N3-C4-N9	-5.13	122.92	126.00
1	CA	760	G	OP1-P-O3'	5.13	116.50	105.20
1	CA	1653	G	N9-C4-C5	5.13	107.45	105.40
1	CA	2255	G	O5'-P-OP2	-5.13	101.08	105.70
1	CA	2500	U	C2-N1-C1'	-5.13	111.54	117.70
1	CA	2851	A	N7-C8-N9	5.13	116.37	113.80
34	DA	678	U	N1-C2-O2	-5.13	119.21	122.80
38	DE	18	ARG	N-CA-C	-5.13	97.14	111.00
1	AA	1364	C	C5-C4-N4	-5.13	116.61	120.20
1	AA	2001	C	O5'-P-OP2	5.13	116.86	110.70
34	BA	1498	U	N3-C4-O4	-5.13	115.81	119.40
1	AA	235	C	C2-N3-C4	-5.13	117.33	119.90
1	AA	1364	C	OP2-P-O3'	5.13	116.49	105.20
1	AA	1403	U	N1-C2-O2	5.13	126.39	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1976	G	N3-C2-N2	5.13	123.49	119.90
1	AA	2056	U	C4-C5-C6	5.13	122.78	119.70
1	AA	2425	G	N3-C2-N2	-5.13	116.31	119.90
25	A1	46	LEU	CA-CB-CG	5.13	127.10	115.30
1	CA	761	A	OP1-P-O3'	5.13	116.49	105.20
1	CA	1767	C	C2-N1-C1'	-5.13	113.16	118.80
1	CA	2778	A	O5'-P-OP1	5.13	116.86	110.70
2	CB	10	C	C6-N1-C2	-5.13	118.25	120.30
1	AA	332	G	N1-C6-O6	-5.13	116.82	119.90
1	AA	2090	U	N3-C2-O2	5.13	125.79	122.20
1	CA	1752	C	N3-C4-C5	5.13	123.95	121.90
1	AA	351	G	C5-N7-C8	5.13	106.86	104.30
1	AA	1035	G	OP2-P-O3'	5.13	116.48	105.20
1	AA	1077	G	C5-C6-O6	-5.13	125.52	128.60
1	AA	1579	C	C5-C6-N1	5.13	123.56	121.00
1	AA	1628	G	N7-C8-N9	5.13	115.66	113.10
1	AA	1673	G	O5'-P-OP2	-5.13	101.08	105.70
1	AA	1698	G	C5-C6-O6	-5.13	125.52	128.60
34	BA	581	G	N3-C4-C5	5.13	131.16	128.60
1	CA	764	A	OP1-P-OP2	-5.13	111.91	119.60
1	CA	773	U	O5'-P-OP2	5.13	116.85	110.70
1	AA	1615	G	N3-C2-N2	5.13	123.49	119.90
1	AA	1656	A	O4'-C1'-N9	-5.13	104.10	108.20
1	AA	1789	G	OP2-P-O3'	5.13	116.48	105.20
1	AA	2641	A	C4-N9-C1'	5.13	135.53	126.30
1	AA	2691	A	N1-C6-N6	-5.13	115.52	118.60
1	AA	2780	C	N3-C4-N4	-5.13	114.41	118.00
1	AA	2835	C	C2-N3-C4	-5.13	117.34	119.90
1	CA	741	G	C5-C6-N1	-5.13	108.94	111.50
1	AA	249	G	N1-C6-O6	-5.12	116.83	119.90
1	AA	726	C	OP2-P-O3'	5.12	116.47	105.20
2	AB	67	G	C5-C6-O6	5.12	131.68	128.60
1	CA	1776	G	C8-N9-C1'	-5.12	120.34	127.00
1	CA	2816	C	O5'-P-OP1	-5.12	101.09	105.70
1	AA	241	G	O5'-P-OP1	-5.12	101.09	105.70
1	AA	824	A	N3-C4-C5	-5.12	123.21	126.80
1	AA	1242	G	C4-N9-C1'	-5.12	119.84	126.50
1	AA	1739	U	C2-N3-C4	-5.12	123.93	127.00
1	AA	1792	C	O5'-P-OP2	5.12	116.85	110.70
1	CA	271(W)	G	N1-C2-N3	5.12	126.97	123.90
1	CA	2276	G	N3-C2-N2	5.12	123.49	119.90
1	CA	2777	G	C2-N3-C4	-5.12	109.34	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	730	C	C6-N1-C2	5.12	122.35	120.30
1	AA	851	A	N3-C4-N9	-5.12	123.30	127.40
1	AA	1700	G	C2'-C3'-O3'	5.12	121.90	113.70
1	AA	1833	A	O4'-C1'-N9	-5.12	104.10	108.20
1	AA	2006	G	C5-C6-N1	-5.12	108.94	111.50
34	BA	554	C	N3-C2-O2	5.12	125.48	121.90
34	BA	807	A	C2-N3-C4	5.12	113.16	110.60
1	CA	704	G	C6-C5-N7	-5.12	127.33	130.40
1	CA	1100	C	C2-N1-C1'	5.12	124.44	118.80
1	CA	1309	G	C5-N7-C8	5.12	106.86	104.30
1	CA	2638	G	C6-C5-N7	5.12	133.47	130.40
1	AA	213	G	N1-C6-O6	-5.12	116.83	119.90
1	AA	308	U	C6-N1-C2	-5.12	117.93	121.00
1	AA	2305	C	C6-N1-C2	5.12	122.35	120.30
1	AA	2781	C	C5-C6-N1	-5.12	118.44	121.00
1	CA	1288	U	OP1-P-OP2	5.12	127.28	119.60
1	CA	1501	C	C5-C6-N1	5.12	123.56	121.00
1	CA	1984	G	O4'-C1'-N9	-5.12	104.10	108.20
1	AA	438	G	C2-N3-C4	5.12	114.46	111.90
1	AA	482	C	OP2-P-O3'	5.12	116.46	105.20
1	AA	793	A	C6-N1-C2	-5.12	115.53	118.60
1	AA	2488	A	N1-C6-N6	5.12	121.67	118.60
1	AA	2617	U	N1-C2-N3	5.12	117.97	114.90
2	AB	88	C	C2-N1-C1'	-5.12	113.17	118.80
1	CA	475	U	O5'-P-OP2	-5.12	101.09	105.70
1	CA	748	G	C4-C5-N7	-5.12	108.75	110.80
1	CA	1763	G	N9-C4-C5	-5.12	103.35	105.40
1	CA	1991	U	N1-C2-N3	5.12	117.97	114.90
1	CA	2269	A	N9-C4-C5	-5.12	103.75	105.80
1	CA	2540	C	N3-C4-N4	-5.12	114.42	118.00
1	AA	120	G	C8-N9-C4	-5.12	104.35	106.40
1	AA	332	G	N3-C4-N9	5.12	129.07	126.00
1	AA	1275	G	OP2-P-O3'	5.12	116.46	105.20
1	AA	2023	A	C5-C6-N6	5.12	127.79	123.70
1	AA	2530	A	C5-C6-N1	-5.12	115.14	117.70
1	AA	2607	G	N1-C6-O6	-5.12	116.83	119.90
34	BA	1507	A	N1-C6-N6	5.12	121.67	118.60
1	CA	2073	C	N1-C2-O2	-5.12	115.83	118.90
1	AA	1372	U	N1-C2-O2	5.12	126.38	122.80
1	AA	1650	C	N3-C4-C5	5.12	123.95	121.90
1	AA	2798	C	C2-N3-C4	-5.12	117.34	119.90
1	AA	2867	G	N3-C2-N2	5.12	123.48	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	503	C	O5'-P-OP1	-5.12	101.10	105.70
1	CA	848	G	C4-N9-C1'	5.12	133.15	126.50
1	CA	1451	C	C2-N1-C1'	-5.12	113.17	118.80
1	CA	1982	C	C4-C5-C6	-5.12	114.84	117.40
1	AA	724	A	N1-C2-N3	5.11	131.86	129.30
1	AA	887	C	C5-C6-N1	-5.11	118.44	121.00
1	AA	1487	G	OP1-P-OP2	-5.11	111.93	119.60
1	AA	1960	A	O4'-C1'-N9	5.11	112.29	108.20
34	BA	1530	G	C2-N3-C4	-5.11	109.34	111.90
1	CA	424	G	OP2-P-O3'	5.11	116.45	105.20
1	CA	1023	U	C5-C4-O4	5.11	128.97	125.90
1	CA	1345	C	O5'-P-OP1	-5.11	101.10	105.70
34	DA	1125	U	C5-C6-N1	5.11	125.26	122.70
34	DA	1469	G	N1-C2-N2	5.11	120.80	116.20
1	AA	337	C	C2-N3-C4	-5.11	117.34	119.90
1	AA	1685	C	C2-N3-C4	-5.11	117.34	119.90
1	AA	2597	U	N1-C2-O2	5.11	126.38	122.80
1	CA	1946	U	N3-C4-O4	5.11	122.98	119.40
1	CA	1992	G	P-O3'-C3'	5.11	125.83	119.70
1	CA	2510	C	N3-C4-C5	-5.11	119.86	121.90
34	DA	1462	G	C8-N9-C4	5.11	108.44	106.40
1	AA	403	C	C6-N1-C2	5.11	122.34	120.30
1	AA	542	C	C5-C6-N1	-5.11	118.44	121.00
1	AA	674	G	C8-N9-C4	-5.11	104.36	106.40
1	AA	723	A	C5-N7-C8	5.11	106.45	103.90
1	AA	1545	C	C5-C4-N4	-5.11	116.62	120.20
1	AA	2411	G	C5-C6-O6	5.11	131.67	128.60
34	BA	552	U	O5'-P-OP1	5.11	116.83	110.70
34	BA	782	A	N9-C4-C5	5.11	107.84	105.80
34	BA	1406	U	C5-C6-N1	-5.11	120.14	122.70
1	AA	591	U	C5-C6-N1	5.11	125.25	122.70
1	AA	1194	A	C4-C5-N7	-5.11	108.15	110.70
1	CA	1860	G	C8-N9-C4	-5.11	104.36	106.40
1	AA	439	A	OP2-P-O3'	5.11	116.44	105.20
1	AA	801	C	C2-N3-C4	-5.11	117.35	119.90
1	AA	893	C	N1-C2-O2	-5.11	115.84	118.90
11	AN	35	ARG	NE-CZ-NH2	5.11	122.85	120.30
1	CA	394	A	C4-C5-C6	-5.11	114.45	117.00
1	CA	569	U	N1-C2-O2	-5.11	119.22	122.80
1	CA	758	C	C5-C4-N4	-5.11	116.62	120.20
1	CA	1790	C	N3-C4-N4	5.11	121.58	118.00
1	CA	2080	G	OP2-P-O3'	5.11	116.44	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	348	A	C2-N3-C4	-5.11	108.05	110.60
1	AA	416	G	N1-C2-N3	5.11	126.96	123.90
1	AA	594	A	N1-C2-N3	-5.11	126.75	129.30
1	AA	622	G	N3-C4-N9	5.11	129.06	126.00
1	AA	1081	U	N3-C2-O2	-5.11	118.63	122.20
1	AA	1364	C	N3-C2-O2	5.11	125.47	121.90
1	AA	1698	G	N3-C4-C5	-5.11	126.05	128.60
1	AA	2289	G	OP2-P-O3'	5.11	116.43	105.20
1	AA	2450	U	OP2-P-O3'	5.11	116.43	105.20
1	CA	1993	U	C6-N1-C2	5.11	124.06	121.00
1	CA	2486	G	N3-C4-C5	-5.11	126.05	128.60
1	AA	263	C	C5-C6-N1	5.10	123.55	121.00
1	AA	1001	G	N1-C2-N3	5.10	126.96	123.90
1	AA	2607	G	C6-N1-C2	5.10	128.16	125.10
34	BA	301	G	N9-C4-C5	5.10	107.44	105.40
34	BA	1502	A	C5-C6-N6	-5.10	119.62	123.70
1	CA	704	G	N1-C6-O6	5.10	122.96	119.90
1	CA	1743	C	C6-N1-C2	-5.10	118.26	120.30
34	DA	105	G	N3-C4-C5	-5.10	126.05	128.60
34	DA	792	A	C2-N3-C4	5.10	113.15	110.60
1	AA	216	A	C4-C5-C6	5.10	119.55	117.00
1	AA	1037	C	OP1-P-O3'	5.10	116.42	105.20
1	AA	2515	A	N7-C8-N9	-5.10	111.25	113.80
1	AA	2768	C	O5'-P-OP1	-5.10	101.11	105.70
1	AA	2781	C	C6-N1-C2	5.10	122.34	120.30
1	AA	2836	A	C2-N3-C4	-5.10	108.05	110.60
2	AB	7	G	C5-C6-O6	-5.10	125.54	128.60
34	BA	1230	C	C6-N1-C2	-5.10	118.26	120.30
1	CA	2273	A	OP2-P-O3'	5.10	116.43	105.20
1	CA	2642	G	OP1-P-OP2	-5.10	111.94	119.60
1	CA	2878	U	OP1-P-OP2	-5.10	111.94	119.60
1	AA	416	G	C5-C6-O6	-5.10	125.54	128.60
1	AA	2399	U	N1-C2-N3	5.10	117.96	114.90
1	AA	474	U	P-O3'-C3'	-5.10	113.58	119.70
1	AA	1704	C	C5-C4-N4	-5.10	116.63	120.20
1	AA	2244	U	OP1-P-OP2	5.10	127.25	119.60
34	BA	553	A	C5-C6-N6	-5.10	119.62	123.70
1	CA	1775	U	N1-C2-N3	5.10	117.96	114.90
1	CA	2522	U	N1-C2-O2	5.10	126.37	122.80
1	CA	2625	G	N3-C4-N9	5.10	129.06	126.00
1	AA	250	G	N1-C2-N2	-5.10	111.61	116.20
1	AA	900	G	OP2-P-O3'	5.10	116.42	105.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1032	C	C2-N1-C1'	-5.10	113.19	118.80
1	AA	1211	U	C5-C4-O4	-5.10	122.84	125.90
1	AA	1276	C	O5'-P-OP2	-5.10	101.11	105.70
1	AA	1471	G	C5-C6-O6	-5.10	125.54	128.60
1	AA	1518	A	C4-C5-C6	5.10	119.55	117.00
1	AA	1685	C	C2-N1-C1'	-5.10	113.19	118.80
1	AA	1819	C	C4-C5-C6	5.10	119.95	117.40
1	AA	2634	C	O5'-P-OP2	-5.10	101.11	105.70
1	AA	2849	G	N1-C6-O6	-5.10	116.84	119.90
34	BA	125	U	OP2-P-O3'	5.10	116.42	105.20
34	BA	386	C	C6-N1-C2	-5.10	118.26	120.30
1	CA	535	C	C2-N3-C4	-5.10	117.35	119.90
34	DA	1482	G	C6-C5-N7	-5.10	127.34	130.40
1	AA	1746	G	O4'-C1'-N9	-5.10	104.12	108.20
1	AA	2641	A	C6-N1-C2	-5.10	115.54	118.60
2	AB	110	G	O5'-P-OP1	-5.10	101.11	105.70
1	CA	254	G	C5-C6-O6	5.10	131.66	128.60
1	AA	129	G	OP2-P-O3'	5.09	116.41	105.20
1	AA	1828	C	N3-C4-C5	5.09	123.94	121.90
1	AA	2072	C	C6-N1-C2	5.09	122.34	120.30
1	AA	2733	U	O5'-P-OP1	-5.09	101.11	105.70
34	BA	299	G	N1-C6-O6	5.09	122.96	119.90
34	BA	574	A	C5-C6-N6	-5.09	119.62	123.70
34	BA	774	G	O5'-P-OP1	-5.09	101.11	105.70
1	CA	197	A	OP1-P-OP2	-5.09	111.96	119.60
1	CA	383	U	O5'-P-OP1	-5.09	101.12	105.70
1	CA	2248	C	N1-C2-N3	5.09	122.77	119.20
1	CA	2587	A	N1-C6-N6	5.09	121.66	118.60
1	CA	2708	G	C8-N9-C1'	-5.09	120.38	127.00
34	DA	62	U	O5'-P-OP2	-5.09	101.11	105.70
1	AA	460	C	N1-C2-N3	5.09	122.77	119.20
1	AA	1741	C	OP1-P-O3'	5.09	116.40	105.20
1	AA	2491	G	C5-C6-O6	5.09	131.66	128.60
1	AA	2553	A	N1-C2-N3	5.09	131.85	129.30
1	AA	2837	C	N3-C2-O2	5.09	125.47	121.90
1	AA	2846	U	N3-C4-C5	5.09	117.66	114.60
1	CA	1791	A	C6-N1-C2	-5.09	115.54	118.60
1	CA	1837	C	O5'-P-OP2	5.09	116.81	110.70
1	AA	347	G	C5-C6-N1	5.09	114.05	111.50
1	AA	810	G	N3-C2-N2	5.09	123.46	119.90
1	AA	904	C	N1-C2-O2	-5.09	115.84	118.90
1	AA	2008	A	C4-C5-C6	5.09	119.55	117.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	724	G	OP1-P-O3'	5.09	116.40	105.20
1	CA	2678	C	C2-N3-C4	-5.09	117.35	119.90
1	AA	438	G	C4-C5-N7	-5.09	108.76	110.80
1	AA	652	A	C5-N7-C8	5.09	106.44	103.90
1	AA	733	G	N1-C2-N3	5.09	126.95	123.90
1	AA	1615	G	C5-C6-N1	5.09	114.05	111.50
1	AA	1971	G	C5-N7-C8	5.09	106.84	104.30
1	AA	2381	A	C2-N3-C4	5.09	113.14	110.60
2	AB	62	C	N1-C2-O2	-5.09	115.85	118.90
34	BA	366	C	N1-C2-O2	-5.09	115.85	118.90
34	BA	401	C	C6-N1-C2	-5.09	118.26	120.30
34	BA	720	C	N1-C2-O2	5.09	121.95	118.90
34	BA	914	A	OP1-P-O3'	5.09	116.40	105.20
1	CA	1142(A)	A	C4-C5-N7	5.09	113.25	110.70
1	CA	1648	C	C6-N1-C1'	5.09	126.91	120.80
1	CA	2897	U	C2-N1-C1'	5.09	123.81	117.70
34	DA	552	U	O5'-P-OP2	-5.09	101.12	105.70
1	AA	2104	A	OP2-P-O3'	5.09	116.39	105.20
1	AA	2350	G	N1-C6-O6	-5.09	116.85	119.90
1	AA	2397	C	N3-C4-N4	-5.09	114.44	118.00
1	CA	1975	G	C6-C5-N7	-5.09	127.35	130.40
34	DA	1205	U	C6-N1-C2	-5.09	117.95	121.00
1	AA	100	G	N3-C2-N2	5.09	123.46	119.90
1	AA	409	G	N1-C2-N2	-5.09	111.62	116.20
1	AA	801	C	O5'-P-OP2	-5.09	101.12	105.70
1	AA	1010	C	N3-C4-N4	-5.09	114.44	118.00
1	AA	1981	G	N3-C4-C5	-5.09	126.06	128.60
1	AA	2058	C	OP2-P-O3'	5.09	116.39	105.20
1	AA	2547	G	C2-N3-C4	-5.09	109.36	111.90
34	BA	194	C	C5-C6-N1	5.09	123.54	121.00
34	BA	1511	G	C6-N1-C2	-5.09	122.05	125.10
34	DA	175	C	C6-N1-C2	-5.09	118.27	120.30
1	AA	1353	A	C2-N3-C4	-5.08	108.06	110.60
1	CA	1195	G	C5-C6-O6	-5.08	125.55	128.60
1	CA	2483	C	N3-C4-C5	5.08	123.93	121.90
1	AA	515	G	C6-N1-C2	-5.08	122.05	125.10
1	AA	555	G	C4-N9-C1'	-5.08	119.89	126.50
1	AA	1299	A	C5-C6-N6	-5.08	119.63	123.70
1	AA	1397	C	N3-C4-C5	5.08	123.93	121.90
1	AA	1962	U	O5'-P-OP2	-5.08	101.12	105.70
1	AA	1975	A	C6-C5-N7	5.08	135.86	132.30
1	AA	2669	A	C8-N9-C4	-5.08	103.77	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	586	C	N1-C2-N3	5.08	122.76	119.20
34	BA	1527	C	N3-C2-O2	-5.08	118.34	121.90
1	CA	803	U	N1-C2-N3	5.08	117.95	114.90
1	CA	1378	A	N7-C8-N9	-5.08	111.26	113.80
1	CA	1644	C	C6-N1-C1'	-5.08	114.70	120.80
1	CA	2237	G	C5-N7-C8	5.08	106.84	104.30
34	DA	6	G	C5-C6-O6	-5.08	125.55	128.60
1	AA	711	C	C2-N1-C1'	-5.08	113.21	118.80
1	AA	855	G	N3-C2-N2	5.08	123.46	119.90
1	AA	1693	C	N3-C2-O2	5.08	125.46	121.90
1	AA	1790	A	C6-N1-C2	-5.08	115.55	118.60
1	AA	1873	G	N7-C8-N9	-5.08	110.56	113.10
34	BA	720	C	N3-C2-O2	-5.08	118.34	121.90
34	BA	1457	G	C8-N9-C4	5.08	108.43	106.40
1	CA	267	C	C6-N1-C2	5.08	122.33	120.30
1	CA	514	A	O5'-P-OP2	-5.08	101.13	105.70
1	CA	946	G	OP1-P-OP2	5.08	127.22	119.60
1	CA	1204	A	O4'-C1'-N9	5.08	112.27	108.20
1	CA	2769	C	C6-N1-C2	-5.08	118.27	120.30
34	DA	540	G	C8-N9-C4	-5.08	104.37	106.40
1	AA	137	G	N3-C4-N9	5.08	129.05	126.00
1	AA	1984	C	C4-C5-C6	-5.08	114.86	117.40
1	AA	2085	C	C5-C6-N1	-5.08	118.46	121.00
1	AA	2689	G	N1-C2-N2	-5.08	111.63	116.20
1	AA	2793	G	N3-C2-N2	-5.08	116.34	119.90
1	AA	2863	C	C6-N1-C2	5.08	122.33	120.30
1	AA	1341	C	C6-N1-C2	5.08	122.33	120.30
1	AA	1694	G	C5-N7-C8	5.08	106.84	104.30
1	AA	1912	A	N1-C6-N6	-5.08	115.55	118.60
1	AA	2448	G	OP1-P-O3'	5.08	116.37	105.20
34	BA	1487	G	O5'-P-OP2	-5.08	101.13	105.70
1	CA	60	G	C5-C6-O6	-5.08	125.55	128.60
1	CA	382	G	N3-C4-C5	-5.08	126.06	128.60
1	CA	733	G	O5'-P-OP2	-5.08	101.13	105.70
1	CA	1190	G	C5-C6-O6	5.08	131.65	128.60
1	CA	2599	G	O5'-P-OP1	-5.08	101.13	105.70
34	DA	1499	A	C2-N3-C4	-5.08	108.06	110.60
56	DW	5	G	C8-N9-C4	5.08	108.43	106.40
1	AA	1986	G	C6-C5-N7	-5.08	127.35	130.40
1	AA	2694	U	C2-N3-C4	-5.08	123.95	127.00
34	BA	875	C	OP1-P-O3'	5.08	116.37	105.20
1	CA	1227	G	C8-N9-C4	-5.08	104.37	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	769	A	OP1-P-OP2	5.08	127.21	119.60
1	AA	1657	C	C5-C6-N1	-5.08	118.46	121.00
1	AA	2437	A	N7-C8-N9	5.08	116.34	113.80
1	AA	2675	G	O5'-P-OP2	-5.08	101.13	105.70
1	AA	2818	U	N3-C2-O2	-5.08	118.65	122.20
15	AR	2	ARG	NE-CZ-NH1	-5.08	117.76	120.30
34	BA	804	U	OP2-P-O3'	5.08	116.37	105.20
1	CA	2508	G	N9-C1'-C2'	-5.08	106.42	112.00
1	AA	183	G	N3-C2-N2	-5.07	116.35	119.90
1	AA	995	G	C5-N7-C8	-5.07	101.76	104.30
1	AA	1020	C	N3-C4-C5	-5.07	119.87	121.90
1	AA	1259	A	O5'-P-OP2	-5.07	101.13	105.70
1	AA	2370	G	C5-N7-C8	-5.07	101.76	104.30
34	BA	33	A	C5-C6-N1	5.07	120.24	117.70
34	BA	594	G	C5-C6-O6	-5.07	125.56	128.60
1	CA	1778	U	C5-C6-N1	-5.07	120.16	122.70
1	AA	50	G	OP1-P-OP2	5.07	127.21	119.60
1	AA	289	G	C8-N9-C4	5.07	108.43	106.40
1	AA	541	C	O5'-P-OP1	5.07	116.79	110.70
1	AA	1766	G	C5-C6-O6	-5.07	125.56	128.60
34	BA	740	U	O5'-P-OP2	-5.07	101.14	105.70
1	CA	491	G	N9-C4-C5	5.07	107.43	105.40
1	CA	2581	G	N7-C8-N9	5.07	115.64	113.10
1	AA	199	C	O5'-P-OP1	5.07	116.78	110.70
1	AA	465	G	C5-C6-O6	-5.07	125.56	128.60
1	AA	726	C	C6-N1-C1'	-5.07	114.72	120.80
1	AA	1308	A	C8-N9-C4	5.07	107.83	105.80
34	BA	566	G	N1-C2-N2	5.07	120.76	116.20
1	CA	804	A	OP2-P-O3'	5.07	116.36	105.20
1	CA	1272	A	O4'-C1'-N9	5.07	112.26	108.20
1	CA	1333	C	N3-C4-C5	5.07	123.93	121.90
1	CA	1702	G	N3-C2-N2	-5.07	116.35	119.90
1	CA	1857	G	N1-C2-N2	-5.07	111.64	116.20
1	CA	2207	G	N7-C8-N9	5.07	115.64	113.10
1	CA	2698	U	O5'-P-OP2	-5.07	101.14	105.70
21	CX	57	LEU	CA-CB-CG	5.07	126.96	115.30
1	AA	627	G	C8-N9-C4	-5.07	104.37	106.40
1	AA	1306	G	OP2-P-O3'	5.07	116.35	105.20
2	AB	62	C	C5-C6-N1	-5.07	118.47	121.00
1	CA	137	C	O5'-P-OP1	-5.07	101.14	105.70
1	CA	2596	U	C2-N3-C4	-5.07	123.96	127.00
34	DA	1232	U	C5-C6-N1	5.07	125.23	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	324	A	N9-C4-C5	-5.07	103.77	105.80
1	AA	1782	C	OP2-P-O3'	5.07	116.35	105.20
1	AA	2014	G	N9-C4-C5	5.07	107.43	105.40
1	AA	2750	G	C8-N9-C4	-5.07	104.37	106.40
34	BA	1507	A	C5-C6-N6	-5.07	119.65	123.70
1	CA	365	C	N3-C2-O2	-5.07	118.35	121.90
1	CA	1947	C	C2-N3-C4	-5.07	117.37	119.90
1	AA	785	G	C5-N7-C8	-5.07	101.77	104.30
1	AA	1194	A	C5-C6-N6	5.07	127.75	123.70
1	AA	1632	A	C2-N3-C4	-5.07	108.07	110.60
1	AA	1858	C	C4-C5-C6	-5.07	114.87	117.40
1	AA	1884	A	C5-N7-C8	-5.07	101.37	103.90
1	AA	2222	C	C4-C5-C6	5.07	119.93	117.40
1	AA	2293	C	N3-C2-O2	5.07	125.45	121.90
1	AA	2782	C	C6-N1-C2	5.07	122.33	120.30
2	AB	23	G	C8-N9-C4	-5.07	104.37	106.40
1	CA	420	C	N3-C4-N4	-5.07	114.45	118.00
1	CA	589	C	C5-C4-N4	5.07	123.75	120.20
1	AA	1401	G	OP2-P-O3'	5.06	116.34	105.20
5	AE	146	THR	C-N-CD	-5.06	109.46	120.60
34	BA	730	G	N3-C2-N2	5.06	123.44	119.90
1	CA	793	A	OP1-P-O3'	5.06	116.34	105.20
1	AA	1064	C	O4'-C1'-N1	5.06	112.25	108.20
1	AA	1351	C	O5'-P-OP2	-5.06	101.14	105.70
1	AA	1428	G	C8-N9-C4	5.06	108.42	106.40
1	AA	2090	U	N1-C2-O2	-5.06	119.26	122.80
1	AA	2252	C	C6-N1-C2	5.06	122.33	120.30
1	CA	592	G	N1-C2-N2	-5.06	111.64	116.20
1	CA	614	U	C5-C4-O4	5.06	128.94	125.90
1	CA	1321	A	C5-C6-N6	-5.06	119.65	123.70
1	CA	1675	C	C6-N1-C2	-5.06	118.28	120.30
1	CA	1920	C	O5'-P-OP2	-5.06	101.14	105.70
1	CA	2469	A	OP1-P-OP2	5.06	127.19	119.60
1	CA	2616	C	O5'-P-OP1	-5.06	101.14	105.70
1	CA	2805	G	C5-C6-O6	5.06	131.64	128.60
57	DZ	92	ILE	N-CA-C	-5.06	97.33	111.00
1	AA	474	U	C5'-C4'-O4'	5.06	115.17	109.10
1	AA	582	G	N1-C6-O6	-5.06	116.86	119.90
34	BA	698	G	C5-N7-C8	-5.06	101.77	104.30
1	CA	440	G	C8-N9-C4	5.06	108.42	106.40
1	CA	2531	A	C8-N9-C4	5.06	107.83	105.80
2	CB	115	G	C8-N9-C4	5.06	108.42	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
56	DW	76	A	C2-N3-C4	-5.06	108.07	110.60
1	AA	52	A	N1-C2-N3	5.06	131.83	129.30
1	AA	103	C	OP1-P-O3'	-5.06	94.07	105.20
1	AA	455	A	C8-N9-C4	-5.06	103.78	105.80
1	AA	1463	C	O5'-P-OP2	5.06	116.77	110.70
1	AA	2735	G	C8-N9-C4	5.06	108.42	106.40
2	AB	97	G	OP2-P-O3'	5.06	116.33	105.20
34	BA	1498	U	N1-C2-N3	5.06	117.94	114.90
1	CA	113	G	C5-C6-O6	-5.06	125.56	128.60
1	CA	587	C	C2-N3-C4	-5.06	117.37	119.90
1	CA	748	G	C6-C5-N7	5.06	133.44	130.40
34	DA	583	A	N1-C6-N6	5.06	121.64	118.60
1	AA	805	C	O5'-P-OP1	5.06	116.77	110.70
1	AA	1288	A	OP2-P-O3'	5.06	116.32	105.20
1	AA	2361	G	C6-N1-C2	-5.06	122.07	125.10
1	AA	2450	U	C5-C4-O4	-5.06	122.86	125.90
34	BA	1384	C	N3-C4-C5	-5.06	119.88	121.90
1	CA	458	G	C6-C5-N7	-5.06	127.37	130.40
1	CA	992	C	OP1-P-O3'	5.06	116.33	105.20
1	CA	1315	C	O5'-P-OP1	-5.06	101.15	105.70
1	CA	1333	C	C4-C5-C6	-5.06	114.87	117.40
34	DA	619	U	N1-C2-O2	-5.06	119.26	122.80
1	AA	72	A	C5-C6-N1	5.06	120.23	117.70
1	AA	1648	U	N3-C2-O2	-5.06	118.66	122.20
1	AA	2400	A	O4'-C1'-N9	5.06	112.25	108.20
1	AA	2738	A	N7-C8-N9	-5.06	111.27	113.80
1	AA	2895	C	N1-C2-O2	-5.06	115.87	118.90
31	A7	42	LEU	CB-CG-CD1	-5.06	102.40	111.00
1	CA	2555	U	OP2-P-O3'	5.06	116.32	105.20
1	AA	1076	G	C5-C6-N1	5.05	114.03	111.50
1	AA	2280	A	N1-C6-N6	5.05	121.63	118.60
1	AA	2837	C	C4-C5-C6	-5.05	114.87	117.40
34	BA	193	C	N3-C2-O2	-5.05	118.36	121.90
34	BA	204	U	N1-C2-O2	5.05	126.34	122.80
1	CA	387	U	OP1-P-O3'	5.05	116.32	105.20
1	CA	1654	A	C5-C6-N1	-5.05	115.17	117.70
1	CA	2061	G	OP1-P-OP2	5.05	127.18	119.60
34	DA	574	A	C8-N9-C4	5.05	107.82	105.80
56	DW	6	G	N7-C8-N9	-5.05	110.57	113.10
1	AA	582	G	C5-C6-O6	5.05	131.63	128.60
1	AA	952	G	OP2-P-O3'	5.05	116.32	105.20
1	AA	2720	G	N3-C2-N2	5.05	123.44	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	791	G	C4-C5-N7	5.05	112.82	110.80
1	AA	806	G	N3-C2-N2	-5.05	116.36	119.90
1	AA	1208	G	C8-N9-C4	5.05	108.42	106.40
1	AA	1642	A	OP1-P-O3'	-5.05	94.09	105.20
1	AA	1789	G	O5'-P-OP2	5.05	116.76	110.70
1	AA	2355	C	C2-N3-C4	-5.05	117.38	119.90
1	AA	2470	G	OP1-P-OP2	5.05	127.18	119.60
1	CA	424	G	N1-C2-N3	-5.05	120.87	123.90
1	CA	1604	C	C6-N1-C2	5.05	122.32	120.30
1	CA	2033	A	N1-C6-N6	5.05	121.63	118.60
1	CA	2548	G	C4-C5-N7	5.05	112.82	110.80
1	AA	209	G	C5-C6-O6	5.05	131.63	128.60
1	AA	1077	G	OP2-P-O3'	5.05	116.31	105.20
1	AA	1080	G	C8-N9-C4	5.05	108.42	106.40
1	AA	1264	G	C8-N9-C4	5.05	108.42	106.40
1	AA	1721	G	C4-N9-C1'	5.05	133.06	126.50
1	AA	1989	C	OP1-P-OP2	5.05	127.17	119.60
1	AA	980	C	N1-C2-N3	5.05	122.73	119.20
1	AA	2377	G	O5'-P-OP1	5.05	116.76	110.70
1	AA	2529	C	O5'-P-OP2	-5.05	101.16	105.70
1	CA	2006	C	OP1-P-OP2	5.05	127.17	119.60
1	AA	96	C	OP1-P-OP2	5.05	127.17	119.60
1	AA	747	G	OP1-P-O3'	-5.05	94.10	105.20
1	AA	1843	A	N1-C6-N6	-5.05	115.57	118.60
1	AA	1851	U	C2-N3-C4	-5.05	123.97	127.00
1	AA	2721	G	C5-C6-N1	-5.05	108.98	111.50
34	BA	266	G	O4'-C1'-N9	-5.05	104.16	108.20
34	BA	303	A	N7-C8-N9	-5.05	111.28	113.80
1	CA	738	G	O5'-P-OP1	5.05	116.75	110.70
1	CA	1333	C	C2-N1-C1'	5.05	124.35	118.80
1	CA	1397	U	N1-C2-N3	5.05	117.93	114.90
1	CA	2023	G	N9-C4-C5	5.05	107.42	105.40
1	AA	348	A	N7-C8-N9	-5.04	111.28	113.80
1	AA	648	G	C2-N3-C4	5.04	114.42	111.90
1	AA	1814	A	C5-C6-N6	5.04	127.74	123.70
34	BA	1466	C	OP1-P-O3'	5.04	116.30	105.20
1	AA	569	G	C8-N9-C4	-5.04	104.38	106.40
1	AA	856	G	OP2-P-O3'	5.04	116.30	105.20
1	AA	1264	G	N1-C2-N2	-5.04	111.66	116.20
1	AA	1291	G	OP1-P-O3'	5.04	116.30	105.20
1	AA	2633	A	N1-C2-N3	-5.04	126.78	129.30
1	AA	2757	G	C5-C6-O6	-5.04	125.57	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
34	BA	558	G	C4-C5-N7	5.04	112.82	110.80
34	BA	566	G	C8-N9-C4	-5.04	104.38	106.40
34	BA	841	U	C2-N1-C1'	5.04	123.75	117.70
34	BA	899	C	C2-N3-C4	-5.04	117.38	119.90
1	CA	410	G	C6-C5-N7	5.04	133.43	130.40
1	CA	446	G	N7-C8-N9	-5.04	110.58	113.10
1	CA	817	C	N3-C4-C5	-5.04	119.88	121.90
1	CA	1721	G	N3-C4-C5	-5.04	126.08	128.60
1	CA	1845	G	N9-C4-C5	5.04	107.42	105.40
1	AA	402	C	O5'-P-OP2	-5.04	101.16	105.70
1	AA	612	C	OP1-P-OP2	5.04	127.16	119.60
1	AA	884	C	N3-C4-N4	-5.04	114.47	118.00
1	AA	1498	C	C6-N1-C2	5.04	122.32	120.30
34	BA	864	A	OP1-P-OP2	-5.04	112.04	119.60
34	BA	1425	U	OP2-P-O3'	5.04	116.29	105.20
1	CA	512	G	O5'-P-OP2	-5.04	101.16	105.70
1	CA	848	G	N3-C4-N9	5.04	129.03	126.00
1	CA	1391	U	O5'-P-OP1	-5.04	101.16	105.70
1	CA	1930	G	C4-C5-N7	-5.04	108.78	110.80
1	CA	2056	G	O4'-C1'-N9	-5.04	104.17	108.20
1	AA	749	G	OP1-P-OP2	5.04	127.16	119.60
1	AA	1356	G	OP2-P-O3'	5.04	116.29	105.20
1	AA	2611	G	N9-C4-C5	-5.04	103.38	105.40
1	CA	1898	U	N3-C4-O4	-5.04	115.87	119.40
1	AA	139	A	O5'-P-OP2	5.04	116.75	110.70
1	AA	407	U	OP2-P-O3'	5.04	116.28	105.20
1	AA	566	C	N3-C4-C5	5.04	123.92	121.90
1	AA	624	C	N1-C2-O2	5.04	121.92	118.90
1	AA	1365	G	C8-N9-C4	-5.04	104.38	106.40
1	AA	1908	C	OP1-P-O3'	5.04	116.28	105.20
1	AA	2048	C	O4'-C1'-N1	5.04	112.23	108.20
1	AA	2444	A	C4-C5-N7	5.04	113.22	110.70
1	AA	2466	G	C8-N9-C4	-5.04	104.39	106.40
1	AA	2575	U	O5'-P-OP2	-5.04	101.17	105.70
1	AA	2599	A	N1-C6-N6	5.04	121.62	118.60
1	AA	2612	A	C6-N1-C2	5.04	121.62	118.60
1	AA	2828	G	O5'-P-OP2	5.04	116.75	110.70
34	BA	694	A	N1-C6-N6	5.04	121.62	118.60
34	BA	777	A	C8-N9-C4	-5.04	103.78	105.80
34	BA	985	C	C6-N1-C2	-5.04	118.28	120.30
1	CA	217	G	C8-N9-C1'	-5.04	120.45	127.00
1	CA	1897	G	C8-N9-C4	5.04	108.42	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	CA	1999	C	C2-N3-C4	-5.04	117.38	119.90
1	AA	477	C	C2-N3-C4	-5.04	117.38	119.90
1	AA	816	G	N3-C4-N9	5.04	129.02	126.00
1	AA	1265	A	C8-N9-C4	-5.04	103.78	105.80
1	AA	1690	G	N3-C4-N9	5.04	129.02	126.00
1	AA	1860	A	C4-C5-N7	-5.04	108.18	110.70
1	CA	430	G	C5-C6-O6	-5.04	125.58	128.60
1	CA	665	C	N3-C4-C5	-5.04	119.89	121.90
34	DA	314	C	C5-C4-N4	-5.04	116.67	120.20
34	DA	766	A	N1-C2-N3	-5.04	126.78	129.30
34	DA	926	G	C8-N9-C4	-5.04	104.39	106.40
1	AA	544	U	C2-N3-C4	-5.04	123.98	127.00
1	AA	977	G	C5-C6-N1	5.04	114.02	111.50
1	AA	2361	G	N1-C6-O6	-5.04	116.88	119.90
1	AA	2497	G	N7-C8-N9	-5.04	110.58	113.10
34	BA	226	G	C8-N9-C4	5.04	108.41	106.40
1	CA	202	U	OP1-P-OP2	5.04	127.15	119.60
1	CA	797	C	N3-C4-C5	-5.04	119.89	121.90
34	DA	331	G	C5-C6-O6	-5.04	125.58	128.60
1	AA	28	A	C5-C6-N6	5.03	127.73	123.70
1	AA	216	A	O5'-P-OP1	-5.03	101.17	105.70
1	AA	762	G	N1-C6-O6	5.03	122.92	119.90
1	AA	866	A	O5'-P-OP2	-5.03	101.17	105.70
1	AA	1046	A	C6-N1-C2	5.03	121.62	118.60
1	AA	1065	U	C5-C4-O4	-5.03	122.88	125.90
1	AA	2363	G	N3-C4-N9	5.03	129.02	126.00
1	AA	2460	A	C8-N9-C4	-5.03	103.79	105.80
1	AA	2593	G	O5'-P-OP2	-5.03	101.17	105.70
2	AB	56	G	C4-N9-C1'	5.03	133.04	126.50
34	BA	266	G	C5-C6-O6	-5.03	125.58	128.60
1	CA	1233	C	C6-N1-C2	-5.03	118.29	120.30
1	CA	1295	C	OP2-P-O3'	5.03	116.27	105.20
34	DA	352	C	O5'-P-OP1	-5.03	101.17	105.70
1	AA	2605	U	OP2-P-O3'	5.03	116.27	105.20
1	CA	25	U	C2-N1-C1'	-5.03	111.66	117.70
1	AA	98	U	N1-C2-O2	5.03	126.32	122.80
1	AA	1307	C	N3-C4-C5	5.03	123.91	121.90
1	CA	210	C	OP2-P-O3'	5.03	116.27	105.20
34	DA	1508	G	O5'-P-OP2	-5.03	101.17	105.70
1	AA	1079	U	C5-C6-N1	-5.03	120.19	122.70
1	AA	1148	C	C2-N1-C1'	5.03	124.33	118.80
1	AA	1655	A	C6-C5-N7	-5.03	128.78	132.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	1804	A	P-O3'-C3'	-5.03	113.67	119.70
1	CA	1984	G	C4-N9-C1'	5.03	133.04	126.50
1	CA	2851	A	C5-N7-C8	-5.03	101.39	103.90
34	DA	912	C	C2-N1-C1'	-5.03	113.27	118.80
1	AA	789	G	OP1-P-OP2	-5.03	112.06	119.60
1	AA	906	G	C5-C6-O6	-5.03	125.58	128.60
1	AA	1155	C	N3-C4-N4	5.03	121.52	118.00
1	AA	1413	A	O5'-P-OP2	-5.03	101.17	105.70
1	AA	1652	G	OP1-P-O3'	5.03	116.26	105.20
1	AA	1721	G	O4'-C1'-N9	-5.03	104.18	108.20
1	AA	2047	C	N3-C4-C5	5.03	123.91	121.90
1	AA	2409	G	OP2-P-O3'	5.03	116.26	105.20
1	AA	2528	G	C8-N9-C4	5.03	108.41	106.40
1	AA	2586	G	C6-C5-N7	5.03	133.42	130.40
2	AB	80	U	N3-C4-C5	5.03	117.62	114.60
34	BA	968	A	C8-N9-C4	5.03	107.81	105.80
34	BA	1509	C	OP1-P-OP2	5.03	127.14	119.60
1	CA	474	G	N3-C4-C5	-5.03	126.09	128.60
1	AA	30	G	OP1-P-O3'	5.03	116.26	105.20
1	AA	555	G	C6-C5-N7	5.03	133.42	130.40
1	AA	633	G	N1-C2-N2	-5.03	111.68	116.20
1	AA	912	C	C4-C5-C6	5.03	119.91	117.40
1	AA	1098	C	C3'-C2'-C1'	-5.03	97.48	101.50
1	AA	1641	G	N3-C4-C5	-5.03	126.09	128.60
1	AA	2727	G	C2-N3-C4	5.03	114.41	111.90
1	CA	195	A	N9-C4-C5	5.03	107.81	105.80
1	CA	562	U	O5'-P-OP2	-5.03	101.18	105.70
1	CA	1142(A)	A	C5-C6-N1	-5.03	115.19	117.70
1	AA	415	G	C4-C5-N7	5.02	112.81	110.80
1	AA	743	G	C4-C5-C6	-5.02	115.79	118.80
1	AA	1659	G	OP1-P-OP2	-5.02	112.06	119.60
34	BA	1485	U	N3-C2-O2	-5.02	118.68	122.20
1	CA	1992	G	C6-C5-N7	5.02	133.41	130.40
1	AA	556	C	N3-C2-O2	-5.02	118.39	121.90
1	AA	1276	C	C2-N3-C4	-5.02	117.39	119.90
1	AA	1399	A	C8-N9-C4	5.02	107.81	105.80
1	AA	1648	U	O5'-P-OP1	-5.02	101.18	105.70
1	AA	1698	G	N3-C4-N9	5.02	129.01	126.00
1	AA	2069	U	C5-C4-O4	-5.02	122.89	125.90
1	AA	2498	G	N1-C6-O6	-5.02	116.89	119.90
1	AA	2510	C	O5'-P-OP1	5.02	116.73	110.70
1	AA	2783	G	N7-C8-N9	-5.02	110.59	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2839	C	N1-C2-O2	-5.02	115.89	118.90
2	AB	78	A	N1-C6-N6	5.02	121.61	118.60
34	BA	757	U	O5'-P-OP1	5.02	116.73	110.70
1	CA	1276	A	N9-C4-C5	-5.02	103.79	105.80
1	CA	1533	G	N3-C4-N9	5.02	129.01	126.00
1	CA	1757	U	C2-N1-C1'	-5.02	111.67	117.70
1	AA	1646	C	N3-C4-C5	5.02	123.91	121.90
1	AA	2657	G	C8-N9-C1'	5.02	133.53	127.00
1	AA	2759	U	C5-C4-O4	5.02	128.91	125.90
34	BA	364	A	N9-C4-C5	5.02	107.81	105.80
34	BA	672	U	N3-C2-O2	-5.02	118.69	122.20
1	CA	1665	A	C6-N1-C2	-5.02	115.59	118.60
1	CA	2557	G	OP2-P-O3'	5.02	116.25	105.20
1	AA	151	C	C6-N1-C2	5.02	122.31	120.30
1	AA	1393	G	N9-C4-C5	5.02	107.41	105.40
1	AA	1482	G	OP1-P-O3'	5.02	116.25	105.20
1	AA	2450	U	N3-C4-C5	5.02	117.61	114.60
34	BA	579	G	C6-C5-N7	-5.02	127.39	130.40
34	BA	585	G	N3-C4-C5	5.02	131.11	128.60
1	CA	1204	A	N9-C4-C5	-5.02	103.79	105.80
1	CA	1695	G	N3-C2-N2	-5.02	116.39	119.90
34	DA	286	G	N7-C8-N9	-5.02	110.59	113.10
1	AA	642	G	N1-C6-O6	-5.02	116.89	119.90
1	AA	1595	C	OP1-P-OP2	-5.02	112.07	119.60
1	AA	1627	A	C6-C5-N7	-5.02	128.79	132.30
1	AA	2017	U	C5-C4-O4	-5.02	122.89	125.90
1	AA	2256	U	C5-C4-O4	5.02	128.91	125.90
1	AA	2647	C	O5'-P-OP2	-5.02	101.19	105.70
1	AA	2737	C	N3-C4-N4	5.02	121.51	118.00
1	CA	1385	G	O4'-C1'-N9	5.02	112.21	108.20
1	CA	2038	G	OP2-P-O3'	5.02	116.24	105.20
1	AA	598	A	C5-C6-N6	-5.02	119.69	123.70
1	AA	1665	G	N9-C4-C5	-5.02	103.39	105.40
1	AA	2702	C	OP2-P-O3'	5.02	116.23	105.20
34	BA	1405	G	C5-C6-O6	5.02	131.61	128.60
1	CA	826	U	OP1-P-O3'	5.02	116.23	105.20
1	CA	2569	G	N1-C6-O6	5.02	122.91	119.90
1	CA	2582	G	C6-C5-N7	-5.02	127.39	130.40
1	AA	454	U	N3-C2-O2	5.01	125.71	122.20
1	AA	612	C	N1-C2-O2	-5.01	115.89	118.90
1	AA	1198	C	OP1-P-OP2	-5.01	112.08	119.60
1	AA	1211	U	C4-C5-C6	5.01	122.71	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	2047	C	C6-N1-C2	5.01	122.31	120.30
1	AA	2673	G	N1-C6-O6	-5.01	116.89	119.90
34	BA	560	U	C3'-C2'-C1'	5.01	105.51	101.50
34	BA	810	C	OP1-P-O3'	5.01	116.23	105.20
45	BL	77	LEU	CB-CG-CD1	-5.01	102.47	111.00
1	CA	1215	G	OP2-P-O3'	5.01	116.23	105.20
1	CA	1637	A	C5-C6-N1	-5.01	115.19	117.70
1	CA	1936	A	O4'-C1'-N9	5.01	112.21	108.20
2	CB	10	C	N3-C2-O2	-5.01	118.39	121.90
1	AA	249	G	N9-C4-C5	5.01	107.41	105.40
1	AA	316	C	C5-C4-N4	-5.01	116.69	120.20
1	AA	1873	G	OP2-P-O3'	5.01	116.23	105.20
1	AA	2761	A	N1-C2-N3	-5.01	126.79	129.30
2	AB	99	G	O5'-P-OP2	-5.01	101.19	105.70
34	BA	1467	G	N1-C2-N2	-5.01	111.69	116.20
34	BA	1505	G	N1-C6-O6	-5.01	116.89	119.90
1	CA	1808	U	C6-N1-C2	5.01	124.01	121.00
1	AA	499	G	N1-C2-N2	-5.01	111.69	116.20
1	AA	621	G	N1-C6-O6	-5.01	116.89	119.90
1	AA	1100	A	C4-C5-N7	5.01	113.21	110.70
1	AA	1149	A	N3-C4-C5	5.01	130.31	126.80
1	AA	2635	G	N3-C4-C5	-5.01	126.09	128.60
1	AA	2700	U	C5-C4-O4	-5.01	122.89	125.90
34	BA	174	C	N1-C2-O2	5.01	121.91	118.90
34	BA	1383	C	C2-N3-C4	5.01	122.41	119.90
1	CA	1423	G	N9-C4-C5	-5.01	103.39	105.40
34	DA	692	U	N3-C4-C5	-5.01	111.59	114.60
34	DA	765	G	C5-C6-O6	5.01	131.61	128.60
1	AA	478	G	C4-C5-N7	-5.01	108.80	110.80
1	AA	2307	C	C5-C6-N1	5.01	123.50	121.00
1	AA	2637	G	C4-C5-N7	-5.01	108.80	110.80
1	CA	762	U	C6-N1-C1'	-5.01	114.19	121.20
1	CA	2083	G	N1-C6-O6	5.01	122.91	119.90
1	CA	2487	G	N9-C4-C5	-5.01	103.40	105.40
1	CA	2727	G	P-O3'-C3'	5.01	125.71	119.70
34	DA	906	G	N9-C4-C5	-5.01	103.40	105.40
1	AA	1683	C	P-O3'-C3'	-5.01	113.69	119.70
2	AB	91	C	C6-N1-C2	5.01	122.30	120.30
1	CA	1567	A	N1-C6-N6	-5.01	115.59	118.60
1	CA	2080	G	O5'-P-OP1	5.01	116.71	110.70
34	DA	550	G	N3-C4-C5	5.01	131.10	128.60
1	AA	309	C	O5'-P-OP2	-5.01	101.19	105.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	676	G	C5-C6-O6	-5.01	125.60	128.60
1	AA	858	U	C2-N3-C4	-5.01	124.00	127.00
1	AA	1289	G	N3-C2-N2	5.01	123.41	119.90
1	AA	1958	A	N9-C4-C5	-5.01	103.80	105.80
1	AA	2290	A	OP2-P-O3'	5.01	116.21	105.20
2	AB	101	G	N9-C4-C5	-5.01	103.40	105.40
1	CA	1350	C	N1-C2-N3	5.01	122.70	119.20
1	CA	1666	G	OP1-P-OP2	-5.01	112.09	119.60
1	CA	1830	C	C6-N1-C2	-5.01	118.30	120.30
1	CA	2059	A	N7-C8-N9	-5.01	111.30	113.80
34	DA	503	C	C5-C6-N1	5.01	123.50	121.00
1	AA	1948	U	O5'-P-OP1	-5.00	101.20	105.70
1	CA	1380	G	O5'-P-OP2	-5.00	101.19	105.70
1	CA	1804	C	OP1-P-O3'	5.00	116.21	105.20
1	CA	1888	G	N3-C4-C5	-5.00	126.10	128.60
1	AA	129	G	N1-C2-N2	-5.00	111.70	116.20
1	AA	493	G	N3-C4-C5	-5.00	126.10	128.60
1	AA	1411	A	C8-N9-C4	5.00	107.80	105.80
1	AA	2870	A	C5-C6-N6	5.00	127.70	123.70
34	BA	801	U	N3-C2-O2	-5.00	118.70	122.20
57	BZ	640	ALA	N-CA-C	5.00	124.51	111.00
1	CA	1698	A	C5-C6-N1	-5.00	115.20	117.70
1	CA	1975	G	N1-C6-O6	5.00	122.90	119.90
1	CA	2057	A	C8-N9-C4	5.00	107.80	105.80
1	CA	2242	G	C5-C6-N1	-5.00	109.00	111.50
2	CB	55	U	O5'-P-OP1	-5.00	101.20	105.70
1	AA	789	G	N9-C4-C5	5.00	107.40	105.40
1	AA	2453	C	N3-C2-O2	-5.00	118.40	121.90
1	AA	2696	U	C5-C6-N1	-5.00	120.20	122.70
1	AA	2838	C	OP2-P-O3'	5.00	116.20	105.20
34	DA	619	U	C5-C4-O4	5.00	128.90	125.90

There are no chirality outliers.

All (15) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
24	A0	11	ARG	Peptide
25	A1	2	SER	Peptide
28	A4	52	THR	Peptide
19	AV	54	GLY	Peptide
21	AX	93	GLU	Peptide
23	AZ	176	PRO	Peptide

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Mol	Chain	Res	Type	Group
53	BT	9	ASN	Peptide
57	BZ	404	VAL	Peptide
1	CA	512	G	Sidechain
17	CT	27	THR	Peptide
21	CX	93	GLU	Peptide
53	DT	9	ASN	Peptide
57	DZ	159	ALA	Peptide
57	DZ	87	HIS	Peptide
57	DZ	91	THR	Peptide

5.2 Too-close contacts [i](#)

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	AA	61426	0	30937	933	0
1	CA	61337	0	30928	1107	0
2	AB	2573	0	1306	39	0
2	CB	2573	0	1306	49	0
3	AC	1063	0	1089	162	0
3	CC	1063	0	1091	203	0
4	AD	2136	0	2218	72	0
4	CD	2142	0	2229	72	0
5	AE	1559	0	1618	46	0
5	CE	1559	0	1618	92	0
6	AF	1584	0	1625	58	0
6	CF	1580	0	1619	68	0
7	AG	1425	0	1443	69	0
7	CG	1424	0	1434	59	0
8	AH	1330	0	1407	40	0
8	CH	1330	0	1407	51	0
9	AK	641	0	309	13	0
9	CK	641	0	309	13	0
10	AL	1025	0	1066	54	0
10	CL	1025	0	1066	50	0
11	AN	1117	0	1184	32	0
11	CN	1117	0	1184	45	0
12	AO	933	0	996	32	0
12	CO	933	0	996	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	AP	1139	0	1223	48	0
13	CP	1135	0	1212	46	0
14	AQ	1122	0	1179	35	0
14	CQ	1122	0	1179	54	0
15	AR	968	0	1033	26	0
15	CR	968	0	1033	44	0
16	AS	877	0	938	30	0
16	CS	870	0	923	43	0
17	AT	1091	0	1151	35	0
17	CT	1083	0	1136	55	0
18	AU	959	0	1019	30	0
18	CU	959	0	1018	34	0
19	AV	771	0	829	24	0
19	CV	771	0	830	16	0
20	AW	886	0	940	26	0
20	CW	886	0	940	31	0
21	AX	750	0	814	27	0
21	CX	750	0	814	27	0
22	AY	806	0	881	23	0
22	CY	806	0	881	39	0
23	AZ	1451	0	1457	56	0
23	CZ	1451	0	1457	66	0
24	A0	653	0	674	29	0
24	C0	653	0	674	22	0
25	A1	755	0	826	24	0
25	C1	755	0	826	23	0
26	A2	588	0	643	13	0
26	C2	588	0	643	19	0
27	A3	469	0	518	13	0
27	C3	464	0	514	11	0
28	A4	558	0	547	24	0
28	C4	532	0	505	14	0
29	A5	455	0	465	18	0
29	C5	455	0	465	20	0
30	A6	453	0	473	18	0
30	C6	449	0	469	13	0
31	A7	418	0	467	17	0
31	C7	418	0	467	17	0
32	A8	517	0	582	22	0
32	C8	517	0	582	28	0
33	A9	307	0	335	10	0
33	C9	307	0	335	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
34	BA	32141	0	16224	675	0
34	DA	32268	0	16287	690	0
35	BB	1846	0	1867	102	0
35	DB	1825	0	1828	82	0
36	BC	1552	0	1546	59	0
36	DC	1544	0	1524	63	0
37	BD	1659	0	1678	93	0
37	DD	1678	0	1720	90	0
38	BE	1129	0	1185	65	0
38	DE	1133	0	1191	64	0
39	BF	812	0	804	27	0
39	DF	820	0	814	22	0
40	BG	1231	0	1238	42	0
40	DG	1235	0	1249	32	0
41	BH	1088	0	1126	62	0
41	DH	1088	0	1126	36	0
42	BI	986	0	995	41	0
42	DI	978	0	966	40	0
43	BJ	709	0	650	37	0
43	DJ	714	0	672	47	0
44	BK	833	0	836	31	0
44	DK	833	0	836	25	0
45	BL	930	0	980	37	0
45	DL	930	0	980	44	0
46	BM	923	0	970	29	0
46	DM	950	0	988	46	0
47	BN	492	0	529	29	0
47	DN	492	0	531	20	0
48	BO	728	0	760	29	0
48	DO	728	0	760	18	0
49	BP	681	0	697	51	0
49	DP	677	0	686	33	0
50	BQ	823	0	891	26	0
50	DQ	823	0	891	35	0
51	BR	555	0	618	22	0
51	DR	555	0	618	25	0
52	BS	661	0	675	34	0
52	DS	646	0	644	30	0
53	BT	728	0	798	35	0
53	DT	731	0	807	24	0
54	BU	199	0	208	5	0
54	DU	199	0	208	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
55	BV	148	0	76	5	0
55	DV	123	0	66	5	0
56	BW	1631	0	839	22	0
56	BY	1581	0	805	24	0
56	DW	1631	0	839	45	0
56	DY	1561	0	796	40	0
57	BZ	5690	0	5783	272	0
57	DZ	5690	0	5783	322	0
58	A0	3	0	0	0	0
58	A2	2	0	0	0	0
58	A4	1	0	0	0	0
58	A5	1	0	0	0	0
58	A6	1	0	0	0	0
58	A7	3	0	0	0	0
58	A8	2	0	0	0	0
58	A9	1	0	0	0	0
58	AA	835	0	0	0	0
58	AB	23	0	0	0	0
58	AD	10	0	0	0	0
58	AE	4	0	0	0	0
58	AF	5	0	0	0	0
58	AG	2	0	0	0	0
58	AH	2	0	0	0	0
58	AN	3	0	0	0	0
58	AO	1	0	0	0	0
58	AP	2	0	0	0	0
58	AQ	3	0	0	0	0
58	AR	1	0	0	0	0
58	AU	3	0	0	0	0
58	AV	3	0	0	0	0
58	AW	4	0	0	0	0
58	AX	1	0	0	0	0
58	AY	1	0	0	0	0
58	AZ	2	0	0	0	0
58	BA	211	0	0	0	0
58	BB	1	0	0	0	0
58	BD	1	0	0	0	0
58	BE	1	0	0	0	0
58	BF	1	0	0	0	0
58	BK	1	0	0	0	0
58	BL	2	0	0	0	0
58	BM	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
58	BN	1	0	0	0	0
58	BT	1	0	0	0	0
58	BV	1	0	0	0	0
58	BW	3	0	0	0	0
58	BZ	2	0	0	0	0
58	C3	1	0	0	0	0
58	C5	1	0	0	0	0
58	C7	1	0	0	0	0
58	C8	1	0	0	0	0
58	CA	664	0	0	0	0
58	CB	13	0	0	0	0
58	CD	3	0	0	0	0
58	CE	6	0	0	0	0
58	CF	5	0	0	0	0
58	CG	1	0	0	0	0
58	CN	1	0	0	0	0
58	CO	2	0	0	0	0
58	CP	3	0	0	0	0
58	CQ	5	0	0	0	0
58	CR	1	0	0	0	0
58	CU	1	0	0	0	0
58	CV	2	0	0	0	0
58	CW	1	0	0	0	0
58	CY	1	0	0	0	0
58	DA	168	0	0	0	0
58	DD	1	0	0	0	0
58	DE	2	0	0	0	0
58	DF	1	0	0	0	0
58	DJ	1	0	0	0	0
58	DK	1	0	0	0	0
58	DT	1	0	0	0	0
58	DW	3	0	0	0	0
58	DZ	2	0	0	0	0
59	A4	1	0	0	0	0
59	A5	1	0	0	0	0
59	A6	1	0	0	0	0
59	A9	1	0	0	0	0
59	AY	1	0	0	0	0
59	BN	1	0	0	0	0
59	C4	1	0	0	0	0
59	C5	1	0	0	0	0
59	C6	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
59	C9	1	0	0	0	0
59	CY	1	0	0	0	0
59	DN	1	0	0	0	0
60	BD	8	0	0	1	0
60	DD	8	0	0	2	0
61	BZ	37	0	47	11	0
61	DZ	37	0	47	17	0
62	BZ	28	0	12	5	0
62	DZ	28	0	12	9	0
63	A0	7	0	0	1	0
63	A1	3	0	0	0	0
63	A3	1	0	0	0	0
63	A5	2	0	0	0	0
63	A6	1	0	0	0	0
63	A7	3	0	0	2	0
63	A8	11	0	0	2	0
63	AA	1408	0	0	60	0
63	AB	36	0	0	1	0
63	AD	15	0	0	1	0
63	AE	19	0	0	5	0
63	AF	8	0	0	1	0
63	AG	3	0	0	1	0
63	AH	1	0	0	0	0
63	AN	2	0	0	1	0
63	AO	3	0	0	0	0
63	AP	15	0	0	0	0
63	AQ	3	0	0	0	0
63	AR	3	0	0	0	0
63	AS	1	0	0	0	0
63	AT	2	0	0	0	0
63	AU	6	0	0	0	0
63	AW	1	0	0	0	0
63	AX	2	0	0	0	0
63	AZ	1	0	0	0	0
63	BA	205	0	0	13	0
63	BD	3	0	0	0	0
63	BE	3	0	0	0	0
63	BJ	1	0	0	0	0
63	BL	2	0	0	0	0
63	BM	1	0	0	0	0
63	BO	1	0	0	0	0
63	BV	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
63	BW	1	0	0	0	0
63	BZ	3	0	0	0	0
63	C0	5	0	0	0	0
63	C1	3	0	0	0	0
63	C3	1	0	0	0	0
63	C5	1	0	0	0	0
63	C7	3	0	0	0	0
63	C8	3	0	0	1	0
63	CA	981	0	0	65	0
63	CB	9	0	0	0	0
63	CD	15	0	0	0	0
63	CE	9	0	0	1	0
63	CF	6	0	0	0	0
63	CP	13	0	0	3	0
63	CQ	1	0	0	0	0
63	CT	3	0	0	0	0
63	CU	4	0	0	1	0
63	CV	1	0	0	0	0
63	CW	1	0	0	0	0
63	CX	1	0	0	0	0
63	CY	1	0	0	0	0
63	DA	153	0	0	11	0
63	DE	2	0	0	0	0
63	DH	1	0	0	1	0
63	DJ	1	0	0	0	0
63	DK	2	0	0	0	0
63	DL	1	0	0	0	0
63	DP	1	0	0	0	0
63	DT	1	0	0	0	0
63	DY	1	0	0	0	0
63	DZ	2	0	0	0	0
All	All	310279	0	209988	7291	0

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 15.

All (7291) 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)
1:AA:1891:G:C5'	3:AC:206:LYS:CG	1.80	1.55
1:AA:1891:G:H5''	3:AC:206:LYS:CG	1.26	1.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1891:G:C5'	3:AC:206:LYS:HG3	1.40	1.40
1:CA:2128:C:H5''	3:CC:219:MET:CE	1.55	1.37
1:AA:2143:G:N2	3:AC:169:THR:OG1	1.57	1.36
1:CA:2176:A:H4'	3:CC:45:HIS:CD2	1.62	1.32
1:AA:1891:G:H5'	3:AC:206:LYS:CD	1.61	1.30
1:CA:2121:G:O2'	3:CC:168:LYS:HB3	1.17	1.26
1:CA:2121:G:N2	3:CC:169:THR:OG1	1.63	1.26
1:CA:2177:C:H1'	3:CC:171:ALA:CB	1.64	1.25
1:CA:2177:C:H4'	3:CC:46:ALA:O	1.40	1.20
57:DZ:87:HIS:NE2	61:DZ:703:FUA:H283	1.55	1.20
1:CA:2176:A:O2'	3:CC:45:HIS:ND1	1.74	1.19
1:CA:1859:A:O2'	3:CC:206:LYS:HE3	1.42	1.18
1:CA:2121:G:C1'	3:CC:168:LYS:HD3	1.75	1.16
1:AA:2143:G:O2'	3:AC:168:LYS:HD3	1.43	1.16
1:CA:2121:G:H1'	3:CC:168:LYS:CD	1.74	1.16
1:AA:2143:G:O2'	3:AC:168:LYS:HB3	1.48	1.14
46:DM:123:ALA:HB3	57:DZ:573:HIS:HB2	1.31	1.12
1:CA:2177:C:H1'	3:CC:171:ALA:HB1	1.32	1.11
1:AA:1891:G:H5'	3:AC:206:LYS:HD2	1.12	1.10
1:AA:1891:G:C5'	3:AC:206:LYS:CD	2.21	1.10
1:AA:2143:G:C2'	3:AC:168:LYS:HD3	1.79	1.09
1:AA:1891:G:H5''	3:AC:206:LYS:HG2	1.25	1.09
61:BZ:703:FUA:H5	61:BZ:703:FUA:H202	1.29	1.08
1:CA:2121:G:O2'	3:CC:168:LYS:CB	2.02	1.07
1:CA:2176:A:O2'	3:CC:45:HIS:CG	2.06	1.07
61:DZ:703:FUA:H5	61:DZ:703:FUA:H202	1.29	1.06
1:AA:1891:G:C4'	3:AC:206:LYS:HG3	1.84	1.06
1:AA:1891:G:C5'	3:AC:206:LYS:HD2	1.85	1.05
17:CT:54:ARG:HA	17:CT:59:THR:HB	1.33	1.05
1:CA:1859:A:H2'	3:CC:206:LYS:CD	1.85	1.05
1:CA:1859:A:H2'	3:CC:206:LYS:HD2	1.09	1.05
1:CA:2176:A:O2'	3:CC:45:HIS:CE1	2.09	1.04
1:CA:1861:G:P	3:CC:206:LYS:HA	1.98	1.04
1:CA:1798:U:H5'	4:CD:259:THR:HG22	1.35	1.03
1:CA:2121:G:C1'	3:CC:168:LYS:CD	2.34	1.02
1:CA:2177:C:H1'	3:CC:171:ALA:HB2	1.37	1.02
57:DZ:160:ARG:NH1	57:DZ:256:THR:OG1	1.93	1.01
1:CA:1859:A:C2'	3:CC:206:LYS:HD2	1.92	0.99
35:DB:185:ILE:HG22	35:DB:199:TYR:HB2	1.40	0.99
1:AA:2331:G:H22	16:AS:3:ARG:HG2	1.26	0.98
1:CA:2128:C:C5'	3:CC:219:MET:CE	2.42	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2124:G:H4'	3:CC:175:PRO:HG3	1.44	0.98
1:CA:2128:C:C5'	3:CC:219:MET:HE3	1.94	0.98
21:AX:31:HIS:HD2	21:AX:33:LYS:H	1.09	0.97
4:CD:60:ARG:HD3	4:CD:86:PRO:HB2	1.45	0.97
1:CA:2128:C:H5''	3:CC:219:MET:HE3	0.98	0.97
1:AA:2143:G:O2'	3:AC:168:LYS:CD	2.13	0.97
1:CA:1860:G:H5'	3:CC:206:LYS:C	1.60	0.96
1:AA:9:U:H3	1:AA:2641:A:H2	1.06	0.96
1:CA:2176:A:C4'	3:CC:45:HIS:CD2	2.48	0.96
57:BZ:644:ARG:HB2	57:BZ:644:ARG:HH11	1.29	0.96
1:CA:2132:U:H1'	3:CC:6:LYS:HB3	1.48	0.95
1:AA:1219:A:H1'	1:AA:1220:U:H5''	1.46	0.95
4:AD:148:GLU:HB2	4:AD:151:LYS:HD2	1.49	0.95
1:CA:2121:G:H1'	3:CC:168:LYS:CE	1.97	0.95
35:BB:16:HIS:HB2	35:BB:204:ASN:HB3	1.48	0.94
1:AA:1891:G:H5'	3:AC:206:LYS:CG	1.77	0.93
57:DZ:191:ASP:OD1	57:DZ:267:LYS:NZ	2.02	0.93
34:DA:396:G:OP1	57:DZ:349:LYS:NZ	2.01	0.93
1:AA:1736:A:H62	1:AA:1745:A:H2	1.17	0.92
1:AA:2459:G:OP2	63:AA:4488:HOH:O	1.87	0.92
1:AA:1154:U:HO2'	1:AA:1155:C:H6	0.97	0.92
1:CA:2121:G:HO2'	3:CC:168:LYS:HB3	1.31	0.92
1:CA:2206:G:H3'	1:CA:2207:G:C8	2.05	0.92
31:A7:24:THR:HG22	31:A7:27:GLY:H	1.32	0.92
1:AA:2143:G:O2'	3:AC:168:LYS:CB	2.17	0.92
1:CA:1861:G:OP2	3:CC:206:LYS:HA	1.69	0.92
1:CA:1782:C:OP1	63:CA:4469:HOH:O	1.88	0.91
1:AA:1829:U:H5'	4:AD:259:THR:HG22	1.51	0.91
14:CQ:110:THR:HG23	14:CQ:113:GLN:HB2	1.52	0.91
1:CA:1019:U:HO2'	1:CA:1021:A:H2	0.98	0.91
53:DT:57:ARG:HH22	53:DT:100:ILE:HD12	1.33	0.91
57:BZ:210:ARG:HB2	57:BZ:210:ARG:HH11	1.35	0.91
1:CA:1271:G:OP2	63:CA:4162:HOH:O	1.89	0.91
1:CA:1689:A:H62	1:CA:1698:A:H2	1.15	0.91
7:AG:41:GLN:NE2	7:AG:154:GLY:O	2.03	0.90
1:AA:1090:G:O2'	1:AA:1157:A:N6	2.04	0.90
57:BZ:13:ARG:HH12	57:BZ:247:ARG:HH12	0.96	0.90
1:CA:2177:C:O2	3:CC:171:ALA:CB	2.19	0.90
34:BA:559:A:OP1	38:BE:126:ARG:NH2	2.04	0.90
34:BA:160:A:N6	34:BA:345:C:OP2	2.02	0.90
1:AA:1891:G:H4'	3:AC:206:LYS:CD	2.01	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:169:GLY:H	57:DZ:170:ARG:NH1	1.69	0.90
37:DD:13:ARG:NH1	37:DD:38:TYR:O	2.05	0.89
37:BD:41:GLY:O	37:BD:43:HIS:N	2.05	0.89
1:CA:1021:A:H62	1:CA:1141:U:H3	1.18	0.89
20:AW:14:PRO:HG2	20:AW:78:GLU:HG2	1.53	0.89
37:BD:104:VAL:HG11	37:BD:146:ILE:HD13	1.55	0.89
1:AA:1100:A:H62	1:AA:1151:U:H3	1.20	0.89
3:CC:31:LYS:NZ	3:CC:181:PHE:O	2.06	0.89
1:CA:1332:G:OP1	63:CA:4126:HOH:O	1.89	0.89
1:CA:2788:C:OP1	5:CE:61:ARG:NH2	2.06	0.89
39:DF:87:ARG:HH11	39:DF:87:ARG:HG3	1.36	0.89
34:DA:1129:C:H42	34:DA:1143:G:H1	1.17	0.88
22:AY:102:CYS:SG	22:AY:103:GLY:N	2.44	0.88
34:DA:922:G:H4'	38:DE:20:GLN:HA	1.53	0.88
1:CA:198:C:OP2	63:CA:4246:HOH:O	1.89	0.88
1:CA:2177:C:O2	3:CC:171:ALA:HB3	1.74	0.88
1:AA:535:C:OP1	63:AA:4769:HOH:O	1.90	0.88
57:BZ:114:VAL:HG23	57:BZ:152:THR:HB	1.56	0.87
1:CA:2046:G:H5'	29:C5:19:ARG:HA	1.55	0.87
1:AA:553:A:C8	1:AA:553:A:H3'	2.09	0.87
57:DZ:88:VAL:O	57:DZ:90:PHE:N	2.08	0.87
16:AS:52:SER:HB2	16:AS:55:ALA:H	1.38	0.87
56:DY:7:A:H61	56:DY:66:U:H3	1.23	0.87
3:AC:31:LYS:NZ	3:AC:181:PHE:O	2.06	0.86
57:BZ:373:ASP:OD2	57:BZ:374:LEU:N	2.07	0.86
1:AA:1065:U:HO2'	1:AA:1067:A:H2	1.21	0.86
24:C0:11:ARG:O	24:C0:14:ARG:NH2	2.09	0.86
37:BD:167:GLY:H	37:BD:168:ARG:HH12	1.20	0.86
1:CA:2177:C:C1'	3:CC:171:ALA:HB2	2.04	0.86
1:CA:2121:G:C1'	3:CC:168:LYS:CE	2.53	0.86
1:CA:2124:G:H4'	3:CC:175:PRO:CG	2.06	0.86
34:DA:1502:A:H2	34:DA:1505:G:H1	1.19	0.86
3:AC:52:PRO:HG2	3:AC:53:ARG:HD3	1.57	0.85
34:BA:560:U:H5'	34:BA:566:G:N2	1.91	0.85
57:DZ:169:GLY:HA3	57:DZ:174:PHE:HA	1.58	0.85
34:DA:959:A:O2'	34:DA:984:C:O2'	1.93	0.85
34:DA:1320:C:N3	52:DS:36:ARG:NH2	2.24	0.85
1:AA:778:C:OP2	63:AA:4735:HOH:O	1.95	0.85
34:BA:406:G:H21	37:BD:119:GLN:HE22	1.22	0.85
57:BZ:132:ARG:HH11	57:BZ:160:ARG:NH1	1.75	0.85
1:CA:1268:A:OP1	63:CA:3952:HOH:O	1.94	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1891:G:C5'	3:AC:206:LYS:HG2	1.84	0.85
34:BA:343:U:H3	34:BA:347:G:H22	1.21	0.85
1:CA:1315:C:OP2	63:CA:4126:HOH:O	1.94	0.85
57:BZ:225:GLU:HA	57:BZ:228:MET:HB3	1.59	0.84
1:AA:874:U:OP1	63:AA:4781:HOH:O	1.93	0.84
1:AA:2143:G:H1'	3:AC:168:LYS:HG2	1.57	0.84
28:C4:40:HIS:HB3	28:C4:43:TYR:HB2	1.58	0.84
1:CA:2121:G:H1'	3:CC:168:LYS:HE2	1.59	0.84
1:CA:2206:G:H3'	1:CA:2207:G:H8	1.42	0.84
34:BA:1125:U:H4'	43:BJ:5:ARG:HH22	1.42	0.84
3:CC:52:PRO:HG2	3:CC:53:ARG:HD3	1.58	0.84
34:BA:1356:G:H2'	34:BA:1357:A:C8	2.13	0.84
63:AE:415:HOH:O	15:AR:3:HIS:NE2	2.10	0.84
56:DW:40:C:H4'	56:DY:36:A:H5'	1.59	0.83
34:BA:78:G:H22	34:BA:92:C:H42	1.27	0.83
34:DA:355:C:OP2	63:DA:3334:HOH:O	1.96	0.83
1:AA:1541:A:OP2	63:AA:4073:HOH:O	1.96	0.83
1:AA:553:A:H2'	1:AA:554:A:H5'	1.60	0.83
21:CX:43:VAL:HG11	21:CX:81:VAL:HG21	1.60	0.83
6:AF:50:SER:HB2	6:AF:94:PRO:HD3	1.61	0.83
21:CX:60:ARG:HH22	31:C7:47:ARG:HH22	1.26	0.83
17:CT:39:ARG:NH2	34:DA:345:C:OP2	2.11	0.83
14:AQ:14:ARG:HG2	14:AQ:41:TRP:HH2	1.43	0.83
1:CA:1022:G:H22	1:CA:1142(A):A:H2	1.24	0.82
20:CW:34:ASN:OD1	20:CW:37:ARG:NH2	2.11	0.82
8:AH:41:MET:HE1	8:AH:65:HIS:HA	1.62	0.82
34:BA:812:C:N3	63:BA:5133:HOH:O	2.10	0.82
55:BV:15:A:HO2'	55:BV:16:U:H5	0.89	0.82
1:CA:1204:A:H2	1:CA:1241:A:H62	1.26	0.82
1:AA:2801:C:OP1	5:AE:61:ARG:NH2	2.11	0.82
1:AA:427:G:N7	63:AA:4918:HOH:O	2.11	0.82
1:AA:656:A:OP1	13:AP:65:ARG:NH1	2.12	0.82
2:CB:66:A:H61	2:CB:109:C:H5'	1.43	0.82
43:BJ:35:SER:HB3	43:BJ:73:ASP:HB2	1.61	0.82
57:BZ:13:ARG:NH1	57:BZ:247:ARG:HH12	1.76	0.82
1:CA:397:G:N7	63:CA:4557:HOH:O	2.11	0.82
34:BA:456:C:H42	34:BA:475:G:H1	1.25	0.82
57:DZ:553:GLY:H	57:DZ:557:GLY:HA2	1.42	0.82
34:BA:262:A:H2'	34:BA:263:A:C8	2.15	0.82
1:CA:218:A:OP2	63:CA:3751:HOH:O	1.96	0.82
1:CA:2836:U:H2'	1:CA:2837:G:C8	2.13	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:1279:A:O2'	34:DA:1282:C:N4	2.13	0.82
36:DC:180:ALA:HB1	36:DC:182:ILE:HG13	1.62	0.82
57:BZ:357:ARG:NH1	57:BZ:373:ASP:OD1	2.12	0.82
1:CA:1658:C:OP1	63:CA:4668:HOH:O	1.95	0.82
34:DA:266:G:H5''	34:DA:268:C:H41	1.45	0.82
1:AA:1891:G:C4'	3:AC:206:LYS:CD	2.57	0.82
1:CA:1861:G:OP1	3:CC:205:ALA:O	1.97	0.82
6:AF:53:THR:HG22	6:AF:55:GLY:H	1.45	0.82
34:BA:1158:C:H5	34:BA:1181:G:H1	1.28	0.82
1:CA:2121:G:H1'	3:CC:168:LYS:CG	2.10	0.82
57:BZ:13:ARG:NH1	57:BZ:280:LEU:O	2.11	0.81
34:DA:742:G:OP2	48:DO:35:ARG:NH2	2.13	0.81
34:BA:392:G:H2'	34:BA:393:A:H8	1.45	0.81
1:CA:2287:A:H62	1:CA:2344:U:H3	1.27	0.81
1:CA:1859:A:O2'	3:CC:206:LYS:CE	2.27	0.81
1:AA:1016:C:OP2	63:AA:5197:HOH:O	1.97	0.81
3:AC:54:ARG:NH2	3:AC:56:ASP:HB3	1.95	0.81
34:BA:1399:C:H4'	34:BA:1400:C:H5''	1.63	0.81
34:BA:538:G:H5''	45:BL:114:LYS:HB2	1.60	0.81
57:BZ:93:GLU:OE2	57:BZ:96:ARG:NH2	2.13	0.81
34:DA:998:G:H1	34:DA:1043:C:H42	1.29	0.81
1:AA:878:G:OP1	63:AA:4750:HOH:O	1.98	0.81
1:AA:2143:G:O2'	3:AC:168:LYS:CG	2.29	0.81
1:CA:2448:A:N1	63:CA:4239:HOH:O	2.13	0.81
3:CC:54:ARG:NH2	3:CC:56:ASP:HB3	1.95	0.81
1:CA:855:G:O2'	24:C0:27:GLU:OE2	1.98	0.81
38:DE:126:ARG:HA	38:DE:131:ILE:HD11	1.62	0.81
46:DM:25:ILE:HG13	46:DM:29:ARG:HG2	1.62	0.81
10:CL:84:LEU:HD21	10:CL:96:VAL:HB	1.63	0.81
35:DB:69:LEU:HB3	35:DB:162:ILE:HG22	1.61	0.81
1:CA:601:C:OP1	6:CF:108:LYS:NZ	2.14	0.80
37:DD:13:ARG:HB2	37:DD:40:PRO:HD3	1.61	0.80
42:BI:17:VAL:HG21	42:BI:81:ILE:HG22	1.63	0.80
49:DP:53:VAL:HG13	49:DP:79:VAL:HG13	1.63	0.80
34:BA:1502:A:H2	34:BA:1505:G:H1	1.29	0.80
57:BZ:92:ILE:HG21	57:BZ:437:THR:HG21	1.64	0.80
14:CQ:34:LEU:HD11	14:CQ:129:THR:HB	1.61	0.80
25:C1:3:LYS:HB2	25:C1:61:ARG:NH1	1.97	0.80
1:AA:927:G:N2	1:AA:944:C:N3	2.29	0.80
56:BW:26:A:H61	56:BW:44:G:H1	1.29	0.80
41:DH:64:LYS:HG2	41:DH:79:VAL:HG21	1.63	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BD:178:VAL:O	37:BD:180:GLY:N	2.14	0.80
34:BA:1375:A:H4'	40:BG:29:LYS:HE2	1.64	0.80
1:CA:1082:U:H5'	10:CL:117:THR:HA	1.63	0.80
34:BA:674:G:N2	34:BA:717:C:O2	2.14	0.80
1:CA:2176:A:H4'	3:CC:45:HIS:HD2	1.40	0.80
34:DA:838:G:H1	34:DA:848:C:H42	1.30	0.80
35:DB:178:ARG:NH1	35:DB:196:LEU:O	2.15	0.80
1:CA:1557:C:OP2	1:CA:1558:A:O2'	2.00	0.79
5:CE:106:GLY:HA3	5:CE:189:PRO:HB2	1.63	0.79
9:CK:73:GLY:O	9:CK:75:GLN:N	2.13	0.79
6:AF:191:ARG:HG2	6:AF:191:ARG:HH11	1.46	0.79
56:BW:28:G:H1	56:BW:42:C:H42	1.29	0.79
35:DB:16:HIS:HB2	35:DB:204:ASN:HB3	1.64	0.79
1:AA:1891:G:C4'	3:AC:206:LYS:CG	2.53	0.79
1:CA:2121:G:O4'	3:CC:168:LYS:HD3	1.82	0.79
5:AE:105:THR:OG1	5:AE:199:ARG:NH2	2.16	0.79
17:AT:65:LYS:HE2	17:AT:67:SER:HB2	1.64	0.79
44:BK:79:SER:HA	44:BK:104:GLN:HB2	1.62	0.79
1:AA:553:A:H3'	1:AA:553:A:H8	1.48	0.79
14:AQ:109:VAL:HG13	14:AQ:113:GLN:HB3	1.63	0.79
38:BE:147:ASP:OD1	38:BE:147:ASP:N	2.16	0.79
1:AA:1151:U:H2'	1:AA:1152:G:H8	1.47	0.79
20:AW:18:ARG:NH1	20:AW:76:VAL:O	2.16	0.79
1:AA:1891:G:H5''	3:AC:206:LYS:HG3	0.83	0.79
15:CR:33:ARG:NH2	29:C5:57:VAL:O	2.16	0.79
1:CA:568:U:O4	63:CA:3804:HOH:O	1.99	0.79
13:AP:100:LEU:HD12	13:AP:112:LEU:HD11	1.65	0.79
1:AA:2299:A:H62	1:AA:2356:U:H3	1.30	0.78
38:BE:100:VAL:O	38:BE:107:ARG:NH2	2.16	0.78
37:DD:64:LEU:HD13	37:DD:198:VAL:HG11	1.65	0.78
1:AA:2328:C:H2'	1:AA:2329:C:H6	1.48	0.78
5:AE:111:ARG:HG3	5:AE:160:TYR:CD2	2.18	0.78
34:BA:1305:G:H22	34:BA:1331:G:H1'	1.48	0.78
35:BB:111:ARG:HH11	35:BB:111:ARG:HG2	1.49	0.78
29:C5:45:VAL:HG11	29:C5:58:LEU:HD13	1.65	0.78
34:BA:452:A:OP1	49:BP:43:LYS:NZ	2.15	0.78
34:DA:115:G:OP1	63:DA:3268:HOH:O	2.00	0.78
1:AA:2143:G:H1'	3:AC:168:LYS:CG	2.05	0.78
1:AA:554:A:N6	1:AA:2063:U:O2	2.17	0.78
1:AA:1201:A:OP1	18:AU:55:ARG:HD3	1.83	0.78
1:CA:2137:C:H42	1:CA:2154:G:H1	1.30	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:CL:6:ALA:HB3	10:CL:30:HIS:HE1	1.48	0.78
14:CQ:34:LEU:HB2	14:CQ:118:LEU:HD22	1.66	0.78
57:DZ:-66:MET:N	57:DZ:-46:VAL:O	2.15	0.78
1:AA:1378:G:OP1	63:AA:4578:HOH:O	2.01	0.78
34:BA:1353:G:N2	34:BA:1369:C:O2	2.14	0.78
14:CQ:18:LYS:O	14:CQ:98:LYS:NZ	2.16	0.78
36:BC:70:VAL:HG22	36:BC:72:LYS:H	1.47	0.78
34:DA:656:C:O2'	48:DO:28:GLN:NE2	2.13	0.78
1:AA:1749:G:N7	63:AA:4933:HOH:O	2.17	0.78
1:CA:641:C:O2'	1:CA:2350:C:OP1	2.01	0.78
15:CR:97:VAL:HG22	15:CR:114:VAL:HG13	1.65	0.78
20:CW:18:ARG:NH1	20:CW:76:VAL:O	2.17	0.78
57:DZ:87:HIS:NE2	61:DZ:703:FUA:C28	2.42	0.78
1:AA:11:G:H2'	1:AA:12:U:H5''	1.66	0.78
57:DZ:182:ARG:O	57:DZ:184:LYS:N	2.17	0.78
1:AA:1312:G:O5'	20:AW:15:ARG:NH2	2.17	0.78
34:BA:972:C:OP2	43:BJ:57:LYS:NZ	2.17	0.78
57:BZ:132:ARG:HH11	57:BZ:160:ARG:HH12	1.32	0.78
25:C1:3:LYS:HB2	25:C1:61:ARG:HH12	1.47	0.78
34:DA:1086:U:H3	34:DA:1099:G:H22	1.29	0.78
34:BA:514:C:O2	34:BA:538:G:N2	2.17	0.78
1:CA:2124:G:C4'	3:CC:175:PRO:HG3	2.14	0.78
21:CX:35:THR:HG22	21:CX:38:GLU:H	1.48	0.78
57:DZ:169:GLY:H	57:DZ:170:ARG:HH12	1.28	0.78
1:AA:1151:U:H2'	1:AA:1152:G:C8	2.19	0.77
1:AA:1501:U:OP1	15:AR:77:ARG:NH1	2.16	0.77
38:BE:33:VAL:HG13	38:BE:112:LEU:HD12	1.66	0.77
57:BZ:119:GLU:OE1	57:BZ:156:ARG:NH1	2.17	0.77
57:BZ:330:VAL:HG12	57:BZ:371:ALA:HA	1.63	0.77
1:CA:1970:A:OP1	63:CA:3911:HOH:O	2.02	0.77
1:CA:2022:U:OP1	63:CA:4132:HOH:O	2.01	0.77
34:DA:1004:A:H62	34:DA:1037:C:H2'	1.48	0.77
38:BE:48:ALA:H	38:BE:54:ALA:HB2	1.48	0.77
1:CA:761:A:OP1	63:CA:4232:HOH:O	2.03	0.77
2:AB:105:A:OP1	23:AZ:72:ARG:NH1	2.18	0.77
1:AA:839:G:H5''	1:AA:840:A:H5'	1.66	0.77
21:AX:31:HIS:CD2	21:AX:33:LYS:H	2.00	0.77
34:BA:382:A:H2'	34:BA:383:A:H8	1.48	0.77
34:BA:128:G:O2'	50:BQ:3:LYS:NZ	2.17	0.77
56:DW:39:PSU:O3'	56:DY:35:A:O2'	2.02	0.77
34:BA:978:A:OP2	34:BA:1363:C:N4	2.17	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:24:ASP:O	3:CC:28:ARG:HG3	1.85	0.77
1:AA:2198:A:O2'	3:AC:45:HIS:CD2	2.38	0.77
34:BA:1086:U:H3	34:BA:1099:G:H22	1.29	0.77
34:BA:656:C:O2'	48:BO:28:GLN:NE2	2.15	0.77
1:CA:2178:C:O2	3:CC:169:THR:HG21	1.85	0.77
34:DA:878:G:H5'	41:DH:89:PRO:HG2	1.66	0.77
49:DP:43:LYS:HG2	49:DP:48:TRP:CD2	2.19	0.77
57:BZ:87:HIS:O	57:BZ:89:ASP:N	2.17	0.77
1:CA:2124:G:O3'	3:CC:175:PRO:CG	2.33	0.77
57:BZ:148:LEU:O	57:BZ:152:THR:OG1	2.02	0.77
3:AC:20:VAL:O	3:AC:21:TYR:HB2	1.83	0.77
3:AC:24:ASP:O	3:AC:28:ARG:HG3	1.85	0.77
49:BP:53:VAL:HG13	49:BP:79:VAL:HG13	1.66	0.77
4:CD:69:ARG:NH2	4:CD:128:GLY:O	2.18	0.77
6:AF:53:THR:CG2	6:AF:55:GLY:H	1.97	0.77
14:AQ:56:ARG:HG3	14:AQ:56:ARG:HH11	1.50	0.77
34:BA:1118:C:H1'	34:BA:1179:A:C4	2.19	0.77
57:BZ:428:LEU:HD13	57:BZ:440:VAL:HG21	1.66	0.77
34:DA:358:U:OP1	57:DZ:381:LYS:NZ	2.15	0.77
34:BA:392:G:H2'	34:BA:393:A:C8	2.19	0.76
1:CA:2808:U:O2	1:CA:2892:A:N6	2.18	0.76
57:DZ:170:ARG:HH11	57:DZ:170:ARG:N	1.82	0.76
3:AC:27:ALA:O	3:AC:30:VAL:HG22	1.85	0.76
35:BB:47:THR:HA	35:BB:202:PRO:HG2	1.67	0.76
55:BV:15:A:O2'	55:BV:16:U:H5	1.66	0.76
7:AG:41:GLN:HG3	7:AG:60:LEU:HD21	1.67	0.76
39:BF:97:PHE:HD1	51:BR:31:LEU:HD21	1.50	0.76
1:CA:2285:C:OP2	30:C6:6:ARG:NH1	2.18	0.76
1:CA:2070:G:OP2	63:CA:4419:HOH:O	2.03	0.76
4:CD:276:LYS:H	4:CD:276:LYS:HD3	1.48	0.76
38:DE:122:GLU:O	38:DE:126:ARG:NH1	2.18	0.76
16:AS:58:LEU:HD22	16:AS:59:LYS:HG3	1.67	0.76
13:CP:44:GLY:O	63:CP:302:HOH:O	2.04	0.76
23:AZ:52:SER:OG	23:AZ:53:ILE:N	2.17	0.76
3:CC:20:VAL:O	3:CC:21:TYR:HB2	1.83	0.76
3:CC:27:ALA:O	3:CC:30:VAL:HG22	1.85	0.76
1:CA:1278:A:OP1	15:CR:36:THR:HG22	1.86	0.76
57:DZ:74:TRP:CE2	57:DZ:273:LEU:HB3	2.19	0.76
34:BA:1274:G:N2	34:BA:1275:A:N7	2.33	0.76
19:AV:29:PRO:HA	19:AV:61:VAL:HG22	1.66	0.76
1:CA:1627:G:OP1	63:CA:4466:HOH:O	2.04	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:1302:U:OP2	46:DM:21:TYR:OH	2.04	0.76
57:BZ:422:GLU:HA	57:BZ:425:SER:HB2	1.68	0.76
36:BC:58:GLU:HB3	43:BJ:92:THR:HG21	1.66	0.76
33:C9:25:VAL:HB	33:C9:34:GLN:HB2	1.68	0.76
57:DZ:526:VAL:HB	57:DZ:566:THR:HA	1.68	0.76
34:BA:941:G:O6	34:BA:1342:C:N4	2.17	0.76
1:CA:2849:U:O4	17:CT:23:ARG:NH2	2.18	0.76
1:CA:1859:A:HO2'	3:CC:206:LYS:HE3	1.51	0.75
1:CA:1817:G:OP1	4:CD:88:ARG:NH2	2.20	0.75
1:CA:1021:A:OP2	11:CN:65:LYS:NZ	2.19	0.75
1:AA:2772:G:OP2	63:AA:4053:HOH:O	2.03	0.75
36:BC:19:GLU:HB3	36:BC:40:ARG:HH22	1.51	0.75
1:CA:1653:G:OP1	1:CA:2822:G:N2	2.19	0.75
35:DB:189:ASP:N	35:DB:189:ASP:OD1	2.19	0.75
23:AZ:72:ARG:NH2	23:AZ:97:GLU:O	2.19	0.75
1:CA:2121:G:C2'	3:CC:168:LYS:HD3	2.17	0.75
16:CS:27:SER:HA	16:CS:88:ASP:HB3	1.67	0.75
23:CZ:144:LEU:HD11	23:CZ:150:LEU:HD23	1.67	0.75
1:AA:1249:A:N6	63:AA:4783:HOH:O	2.15	0.75
1:CA:54:G:N7	63:CA:4438:HOH:O	2.18	0.75
34:DA:1318:A:H1'	52:DS:37:ARG:HD3	1.67	0.75
57:DZ:33:LEU:HD12	57:DZ:360:ALA:HB2	1.68	0.75
1:CA:300:A:OP1	22:CY:86:ARG:NH2	2.19	0.75
43:BJ:49:VAL:HG23	47:BN:41:ARG:HB2	1.68	0.75
34:BA:972:C:O2'	43:BJ:55:LYS:O	2.05	0.75
57:BZ:404:VAL:HG13	57:BZ:405:PRO:HD3	1.67	0.75
1:AA:1831:C:OP1	4:AD:260:ARG:NH2	2.19	0.75
61:BZ:703:FUA:H202	61:BZ:703:FUA:C5	2.15	0.75
1:AA:1084:C:H42	1:AA:1163:G:H1	1.31	0.75
34:BA:1182:G:H4'	34:BA:1183:A:H5'	1.67	0.75
34:BA:401:C:OP2	37:BD:73:ARG:NH1	2.20	0.75
17:CT:16:ARG:HD2	17:CT:19:LEU:HD11	1.69	0.75
34:DA:493:G:O6	63:DA:3343:HOH:O	2.02	0.75
57:BZ:13:ARG:HH12	57:BZ:247:ARG:NH1	1.81	0.74
38:DE:33:VAL:HG21	38:DE:109:ILE:HA	1.67	0.74
57:DZ:13:ARG:NE	57:DZ:280:LEU:O	2.18	0.74
8:AH:98:LEU:HD22	8:AH:125:VAL:HG23	1.70	0.74
1:CA:825:C:OP1	63:CA:4609:HOH:O	2.05	0.74
6:CF:129:PHE:HB2	6:CF:132:VAL:HG21	1.68	0.74
7:CG:122:PRO:HB3	7:CG:170:ARG:HH21	1.50	0.74
41:BH:113:SER:HB2	41:BH:134:ILE:HD11	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BO:26:GLU:OE2	48:BO:77:ARG:NE	2.16	0.74
37:DD:8:VAL:HG22	37:DD:21:LEU:HD13	1.69	0.74
46:DM:15:VAL:HG13	46:DM:45:VAL:HG22	1.69	0.74
16:AS:84:GLN:HA	16:AS:111:GLU:HB2	1.70	0.74
44:BK:86:GLY:N	44:BK:112:THR:OG1	2.16	0.74
14:CQ:85:LYS:HG2	24:C0:7:LEU:HB3	1.70	0.74
56:DW:39:PSU:O2'	56:DY:35:A:H1'	1.87	0.74
13:CP:89:ALA:O	13:CP:121:LYS:NZ	2.18	0.74
3:CC:48:LEU:HB3	3:CC:50:ILE:HD12	1.70	0.74
35:DB:163:PHE:HD1	35:DB:185:ILE:HG13	1.51	0.74
1:AA:1093:G:H21	1:AA:1157:A:H2	1.35	0.74
1:AA:1891:G:C4'	3:AC:206:LYS:HD2	2.18	0.74
34:BA:200:G:H1	34:BA:217:C:H42	1.35	0.74
6:CF:164:ARG:HD2	6:CF:175:THR:HG23	1.70	0.74
1:AA:1154:U:O2'	1:AA:1155:C:H5''	1.88	0.74
34:BA:382:A:H2'	34:BA:383:A:C8	2.22	0.74
1:CA:2612:C:OP2	29:C5:2:ALA:N	2.20	0.74
12:CO:2:ILE:HB	12:CO:33:ALA:HB3	1.70	0.74
1:AA:1199:C:OP2	63:AA:4586:HOH:O	2.06	0.74
1:AA:2317:A:H5''	7:AG:134:GLY:HA3	1.69	0.74
37:DD:135:LEU:O	37:DD:137:SER:N	2.19	0.74
1:AA:1094:A:OP2	1:AA:1155:C:N4	2.21	0.74
19:AV:40:LEU:HB2	19:AV:46:VAL:HG13	1.70	0.74
53:BT:10:LEU:HB3	53:BT:12:ALA:H	1.52	0.74
61:DZ:703:FUA:H202	61:DZ:703:FUA:C5	2.15	0.74
61:DZ:703:FUA:H5	61:DZ:703:FUA:C20	2.12	0.74
1:AA:1001:G:OP2	14:AQ:14:ARG:NH2	2.21	0.73
34:BA:1197:G:OP2	63:BA:5175:HOH:O	2.05	0.73
57:BZ:132:ARG:HD3	57:BZ:160:ARG:HH12	1.51	0.73
1:CA:2121:G:O4'	3:CC:168:LYS:NZ	2.20	0.73
15:CR:70:LEU:O	15:CR:72:ASP:N	2.19	0.73
34:DA:337:C:H2'	34:DA:338:A:C8	2.22	0.73
15:CR:85:PRO:O	15:CR:87:TYR:N	2.22	0.73
35:DB:16:HIS:CD2	35:DB:17:PHE:H	2.06	0.73
2:AB:13:A:N1	2:AB:69:G:O2'	2.19	0.73
34:BA:972:C:OP1	63:BA:5237:HOH:O	2.07	0.73
1:CA:11:G:H2'	1:CA:12:U:H5''	1.70	0.73
47:BN:27:CYS:SG	47:BN:28:GLY:N	2.61	0.73
7:CG:43:LEU:HD11	7:CG:153:ARG:HG2	1.70	0.73
29:C5:16:ARG:HG2	29:C5:16:ARG:HH11	1.53	0.73
1:CA:2128:C:C5'	3:CC:219:MET:HE1	2.14	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DF:2:ARG:NE	39:DF:69:GLU:HG2	2.02	0.73
57:DZ:69:VAL:HG21	57:DZ:374:LEU:HD23	1.71	0.73
1:AA:2601:A:OP2	63:AA:4561:HOH:O	2.06	0.73
13:AP:38:GLN:HG2	13:AP:45:LEU:HD23	1.70	0.73
34:BA:711:G:OP1	39:BF:54:LYS:NZ	2.17	0.73
1:CA:2107:C:H42	1:CA:2182:G:H1	1.36	0.73
49:DP:21:VAL:HG13	49:DP:34:GLU:HB3	1.71	0.73
4:AD:8:PRO:HB3	4:AD:14:ARG:HB2	1.70	0.73
1:AA:2143:G:HO2'	3:AC:168:LYS:HB3	1.52	0.73
50:BQ:66:SER:O	50:BQ:70:ARG:NH1	2.22	0.73
53:BT:10:LEU:HD23	53:BT:12:ALA:HB2	1.71	0.73
57:BZ:210:ARG:HH11	57:BZ:210:ARG:CB	2.01	0.73
30:C6:6:ARG:NH1	30:C6:26:ASN:HB2	2.04	0.73
13:CP:100:LEU:HD12	13:CP:112:LEU:HD11	1.71	0.73
34:BA:820:U:H4'	34:BA:821:G:OP2	1.89	0.73
6:CF:101:LEU:O	6:CF:106:ARG:NH1	2.21	0.73
7:CG:113:ARG:NH1	7:CG:139:LEU:O	2.21	0.73
34:DA:1347:G:N2	34:DA:1373:G:H2'	2.04	0.73
1:AA:965:G:N2	1:AA:2281:A:OP2	2.22	0.73
22:AY:92:ASN:HB2	22:AY:94:LYS:H	1.51	0.73
34:BA:560:U:H5'	34:BA:566:G:H22	1.51	0.73
34:BA:564:C:O2'	41:BH:91:ARG:NH2	2.20	0.73
39:BF:97:PHE:HB2	51:BR:32:ARG:HD2	1.68	0.73
34:DA:664:G:H22	34:DA:741:G:H1	1.37	0.73
1:AA:2199:C:O2	3:AC:173:HIS:CE1	2.42	0.72
5:AE:110:GLY:O	63:AE:415:HOH:O	2.06	0.72
11:AN:75:TYR:CE2	11:AN:77:GLY:HA2	2.23	0.72
4:CD:206:LEU:HD22	4:CD:211:ARG:HG2	1.70	0.72
13:CP:88:LEU:HD11	13:CP:114:ILE:HD12	1.69	0.72
6:AF:135:LYS:HB2	6:AF:138:GLU:HG3	1.70	0.72
37:BD:101:LEU:O	37:BD:103:ASN:N	2.22	0.72
41:BH:6:ILE:HG22	41:BH:10:LEU:HD21	1.71	0.72
1:CA:1418:G:N7	63:CA:4067:HOH:O	2.20	0.72
1:CA:2178:C:OP1	3:CC:47:LYS:HG2	1.88	0.72
1:CA:854:G:O6	63:CA:4561:HOH:O	2.07	0.72
2:CB:86:G:N2	2:CB:91:C:O2	2.18	0.72
35:BB:179:LYS:HA	41:BH:72:PRO:HG3	1.70	0.72
41:DH:86:ILE:HG21	41:DH:133:LEU:HD13	1.69	0.72
37:BD:167:GLY:H	37:BD:168:ARG:NH1	1.87	0.72
34:BA:664:G:H22	34:BA:741:G:H1	1.37	0.72
57:DZ:103:GLY:H	57:DZ:130:VAL:HG23	1.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2200:C:H4'	3:AC:47:LYS:NZ	2.05	0.72
10:AL:17:ALA:HB3	10:AL:38:VAL:HG13	1.70	0.72
35:DB:119:GLU:OE2	35:DB:153:ARG:NH1	2.22	0.72
1:AA:2297:C:OP2	30:A6:6:ARG:NH1	2.23	0.72
3:AC:48:LEU:HB3	3:AC:50:ILE:HD12	1.70	0.72
1:CA:1798:U:OP2	4:CD:274:ARG:NH2	2.23	0.72
1:CA:981:A:OP1	63:CA:4070:HOH:O	2.08	0.72
34:DA:1304:G:N2	34:DA:1332:A:OP2	2.21	0.72
35:DB:178:ARG:HH21	41:DH:74:PRO:HB3	1.55	0.72
37:BD:176:LEU:HG	37:BD:178:VAL:HG22	1.71	0.72
1:CA:2121:G:C1'	3:CC:168:LYS:HE2	2.16	0.72
1:CA:2439:A:C8	1:CA:2439:A:H5''	2.24	0.72
1:CA:2660:A:N6	57:DZ:661:SER:OG	2.22	0.72
1:AA:1154:U:O2'	1:AA:1155:C:H6	1.72	0.72
8:CH:3:ARG:HH22	8:CH:5:GLY:H	1.37	0.72
14:CQ:57:HIS:CD2	14:CQ:117:ALA:HB2	2.25	0.72
37:BD:13:ARG:NH1	37:BD:38:TYR:O	2.23	0.72
34:BA:395:C:O3'	57:BZ:349:LYS:NZ	2.22	0.72
1:CA:740:U:OP2	63:CA:4170:HOH:O	2.08	0.72
34:DA:1344:C:H4'	42:DI:120:ARG:HB3	1.72	0.72
34:DA:677:U:H3	34:DA:713:G:H22	1.38	0.72
1:AA:2199:C:O2	3:AC:173:HIS:HE1	1.71	0.71
38:BE:152:ARG:HB3	41:BH:43:GLY:HA3	1.72	0.71
57:DZ:132:ARG:N	57:DZ:132:ARG:HD2	2.05	0.71
57:DZ:127:LYS:NZ	57:DZ:404:VAL:HG11	2.05	0.71
1:AA:2457:G:OP1	6:AF:74:ARG:NH2	2.23	0.71
1:CA:1268:A:OP1	63:CA:3954:HOH:O	2.08	0.71
1:CA:1310:G:OP2	31:C7:9:ARG:NH1	2.22	0.71
1:CA:2445:G:OP1	6:CF:74:ARG:NH2	2.23	0.71
8:CH:149:ARG:NH1	8:CH:167:GLU:OE2	2.23	0.71
57:DZ:146:LEU:HD12	57:DZ:167:PRO:HD3	1.70	0.71
47:BN:3:ARG:O	47:BN:6:LEU:N	2.19	0.71
16:CS:83:LYS:HG3	16:CS:84:GLN:HG3	1.72	0.71
23:CZ:29:TYR:HB3	23:CZ:34:ASN:HD22	1.54	0.71
56:DY:19:G:N2	56:DY:56:C:N3	2.37	0.71
30:A6:14:THR:HB	30:A6:48:VAL:O	1.90	0.71
1:AA:1466:U:O2'	1:AA:1467:G:OP1	2.09	0.71
1:AA:1480:A:H61	1:AA:1605:A:N6	1.88	0.71
1:AA:2154:U:C5	3:AC:6:LYS:HB2	2.25	0.71
12:AO:64:ARG:NH1	12:AO:81:ASP:OD1	2.18	0.71
17:AT:54:ARG:HA	17:AT:59:THR:HB	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
47:BN:3:ARG:O	47:BN:5:ALA:N	2.24	0.71
52:BS:63:THR:OG1	52:BS:65:ASN:ND2	2.24	0.71
57:DZ:92:ILE:HD13	57:DZ:93:GLU:H	1.55	0.71
34:DA:1106:G:H5''	36:DC:172:ARG:HG2	1.72	0.71
57:DZ:71:THR:HG22	57:DZ:80:ASN:OD1	1.90	0.71
3:CC:51:ASP:HB3	3:CC:57:GLN:OE1	1.91	0.71
11:CN:22:THR:HB	11:CN:25:ARG:HG3	1.71	0.71
45:DL:75:HIS:CD2	45:DL:77:LEU:H	2.08	0.71
13:AP:62:LEU:O	32:A8:13:ARG:HD3	1.91	0.71
1:AA:2108:U:N3	1:AA:2245:U:O4	2.18	0.71
41:BH:86:ILE:HG21	41:BH:133:LEU:HD13	1.72	0.71
1:CA:1971:A:OP2	4:CD:242:ARG:NH2	2.24	0.71
46:DM:123:ALA:HB3	57:DZ:573:HIS:CB	2.16	0.71
2:CB:56:G:H5'	7:CG:27:ASN:ND2	2.05	0.71
7:CG:27:ASN:HB3	7:CG:30:GLU:HG3	1.72	0.71
12:CO:104:ARG:HH22	17:CT:43:GLN:HE22	1.36	0.71
14:CQ:63:LYS:HG2	23:CZ:178:GLU:HG2	1.73	0.71
34:BA:782:A:OP1	63:BA:5146:HOH:O	2.08	0.71
35:BB:187:LEU:HA	35:BB:201:ILE:HB	1.73	0.71
1:CA:1386:C:H2'	1:CA:1387:C:H6	1.54	0.71
1:CA:1434:A:H61	1:CA:1558:A:H62	1.39	0.71
1:CA:2176:A:H4'	3:CC:45:HIS:NE2	2.06	0.71
2:CB:62:C:H2'	2:CB:63:G:H8	1.55	0.71
1:CA:2132:U:C1'	3:CC:6:LYS:HB3	2.21	0.71
35:BB:69:LEU:HB3	35:BB:162:ILE:HG22	1.72	0.71
1:CA:1064:C:H4'	10:CL:89:HIS:HA	1.72	0.71
19:CV:6:LYS:HB2	19:CV:38:LEU:HD21	1.71	0.71
34:DA:920:U:H2'	34:DA:921:U:C6	2.25	0.71
6:AF:75:HIS:ND1	63:AF:406:HOH:O	2.23	0.70
57:BZ:169:GLY:O	57:BZ:173:THR:OG1	2.09	0.70
1:CA:1647:G:OP1	63:CA:4162:HOH:O	2.09	0.70
1:CA:2839:G:H5'	15:CR:46:GLY:HA2	1.73	0.70
19:CV:40:LEU:HB2	19:CV:46:VAL:HG13	1.73	0.70
37:DD:132:ARG:HB3	37:DD:132:ARG:HH21	1.56	0.70
42:DI:9:ARG:H	42:DI:79:LEU:HD23	1.55	0.70
1:AA:181:C:H2'	1:AA:182:U:H5'	1.71	0.70
36:BC:130:VAL:HG11	36:BC:153:VAL:HG11	1.73	0.70
44:BK:98:LEU:O	44:BK:101:SER:OG	2.07	0.70
37:DD:119:GLN:HG3	37:DD:123:HIS:CD2	2.26	0.70
4:AD:95:LEU:HD11	4:AD:105:ILE:HD13	1.73	0.70
6:AF:185:ASP:HA	6:AF:188:ARG:HD3	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AK:73:GLY:O	9:AK:75:GLN:N	2.19	0.70
57:BZ:-53:ASP:H	57:BZ:-50:GLN:NE2	1.89	0.70
1:CA:1309:G:H4'	31:C7:7:PRO:HB2	1.74	0.70
1:CA:9:U:H3	1:CA:2629:A:H2	1.39	0.70
1:CA:2121:G:O2'	3:CC:168:LYS:HD3	1.91	0.70
10:CL:76:TYR:HD1	10:CL:79:ARG:HH21	1.39	0.70
11:CN:14:VAL:HG11	11:CN:138:LEU:HD12	1.71	0.70
34:DA:1075:C:OP1	35:DB:179:LYS:NZ	2.24	0.70
35:DB:189:ASP:O	35:DB:192:SER:OG	2.09	0.70
34:DA:1226:C:O2'	46:DM:111:LYS:NZ	2.24	0.70
1:AA:455:A:H8	1:AA:455:A:OP2	1.74	0.70
56:BY:60:U:H5''	56:BY:61:C:H5	1.56	0.70
61:BZ:703:FUA:H122	61:BZ:703:FUA:H231	1.73	0.70
1:CA:2121:G:H21	3:CC:169:THR:HG1	1.38	0.70
12:CO:2:ILE:HD12	12:CO:6:THR:HG21	1.72	0.70
34:BA:501:C:H2'	34:BA:502:G:C8	2.27	0.70
41:DH:120:THR:H	41:DH:123:GLU:HB3	1.56	0.70
47:DN:27:CYS:SG	47:DN:28:GLY:N	2.64	0.70
6:AF:9:ILE:HG21	6:AF:125:LEU:HD22	1.72	0.70
38:BE:36:ASP:OD1	38:BE:38:GLN:N	2.20	0.70
1:CA:2631:G:O2'	1:CA:2810:A:N1	2.24	0.70
35:DB:162:ILE:HD11	35:DB:184:VAL:HG22	1.73	0.70
61:DZ:703:FUA:H122	61:DZ:703:FUA:H231	1.73	0.70
3:AC:51:ASP:HB3	3:AC:57:GLN:OE1	1.91	0.70
3:CC:15:VAL:O	3:CC:16:ASP:HB3	1.92	0.70
19:CV:25:LEU:N	19:CV:92:THR:OG1	2.21	0.70
57:DZ:110:SER:OG	57:DZ:137:ASN:O	2.08	0.70
44:BK:86:GLY:H	44:BK:112:THR:HG1	1.37	0.70
30:C6:8:LYS:HD3	32:C8:34:TRP:CD2	2.27	0.70
1:CA:2124:G:O3'	3:CC:175:PRO:HG3	1.91	0.70
8:CH:86:GLU:HB2	8:CH:165:ALA:HB2	1.74	0.70
34:DA:1111:A:N1	36:DC:177:THR:OG1	2.22	0.70
34:DA:1240:U:OP2	40:DG:116:ALA:N	2.23	0.70
38:DE:50:GLU:HB2	38:DE:53:LEU:HD13	1.73	0.70
1:CA:323:G:HO2'	1:CA:1205:U:H3	1.40	0.69
57:DZ:117:GLN:NE2	57:DZ:120:THR:OG1	2.24	0.69
57:DZ:92:ILE:O	57:DZ:94:VAL:N	2.25	0.69
1:AA:1740:U:O2'	4:AD:14:ARG:NH2	2.24	0.69
37:BD:23:GLY:N	37:BD:26:CYS:SG	2.65	0.69
14:CQ:57:HIS:HD2	14:CQ:117:ALA:HB2	1.56	0.69
15:AR:67:LEU:HD13	15:AR:76:VAL:HG21	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:BZ:264:LEU:HB2	62:BZ:704:GDP:C6	2.26	0.69
1:CA:1349:A:OP1	63:CA:4335:HOH:O	2.10	0.69
8:CH:3:ARG:NH2	8:CH:5:GLY:H	1.91	0.69
10:AL:72:PRO:O	10:AL:111:LYS:NZ	2.24	0.69
1:AA:2827:G:OP1	15:AR:99:LYS:NZ	2.25	0.69
34:DA:23:C:OP2	34:DA:561:U:N3	2.21	0.69
41:DH:51:VAL:HG11	41:DH:60:ARG:HH11	1.56	0.69
46:DM:108:ARG:HD2	46:DM:114:ARG:HD2	1.74	0.69
57:DZ:619:ASP:HB3	57:DZ:662:LYS:HD2	1.74	0.69
30:A6:13:CYS:SG	30:A6:47:THR:HG21	2.32	0.69
44:BK:73:MET:HG2	44:BK:103:LEU:HD21	1.73	0.69
47:BN:3:ARG:HH21	47:BN:3:ARG:HB3	1.56	0.69
57:BZ:428:LEU:HA	57:BZ:431:LEU:HB2	1.74	0.69
57:BZ:552:SER:O	57:BZ:591:LYS:NZ	2.25	0.69
1:CA:2150:U:H2'	1:CA:2151:G:H8	1.58	0.69
3:CC:46:ALA:HB3	3:CC:172:ILE:HG22	1.75	0.69
1:CA:1842:G:O2'	4:CD:253:GLN:NE2	2.26	0.69
1:AA:2118:U:H2'	1:AA:2119:C:C6	2.26	0.69
28:A4:61:ARG:HH21	52:BS:42:PRO:HD2	1.55	0.69
57:BZ:485:GLU:HG2	57:BZ:554:PRO:HD2	1.74	0.69
1:CA:323:G:O2'	1:CA:1205:U:N3	2.26	0.69
57:DZ:485:GLU:HG2	57:DZ:555:LEU:HD12	1.72	0.69
40:BG:111:ARG:NH2	40:BG:126:ASP:OD2	2.26	0.69
1:CA:2022:U:OP1	63:CA:4131:HOH:O	2.10	0.69
1:AA:1189:A:OP1	11:AN:25:ARG:NH2	2.25	0.69
3:AC:183:PRO:HG2	3:AC:184:GLU:OE2	1.92	0.69
3:AC:57:GLN:O	3:AC:57:GLN:HG3	1.93	0.69
8:AH:3:ARG:HH12	8:AH:65:HIS:HB3	1.58	0.69
34:BA:36:C:O2'	45:BL:117:ARG:NH2	2.26	0.69
1:CA:1253:A:N7	63:CA:3852:HOH:O	2.26	0.69
1:CA:1860:G:C5'	3:CC:207:GLY:N	2.56	0.69
1:CA:2836:U:H2'	1:CA:2837:G:H8	1.58	0.69
3:CC:57:GLN:HG3	3:CC:57:GLN:O	1.93	0.69
6:CF:110:LEU:HD11	6:CF:181:LEU:HD23	1.73	0.69
2:CB:104:U:O3'	23:CZ:72:ARG:NH1	2.25	0.69
34:DA:352:C:OP2	63:DA:3237:HOH:O	2.11	0.69
1:AA:1056:A:OP2	63:AA:4596:HOH:O	2.11	0.69
1:AA:555:G:O4'	1:AA:555:G:N3	2.18	0.69
15:AR:55:ALA:HB2	15:AR:79:LEU:HD13	1.74	0.69
61:BZ:703:FUA:H5	61:BZ:703:FUA:C20	2.12	0.69
1:CA:1837:C:O2'	1:CA:1927:A:N3	2.22	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CE:33:VAL:HG21	5:CE:36:ARG:HE	1.58	0.69
34:DA:188:C:H2'	34:DA:189:G:H8	1.57	0.69
34:DA:838:G:N2	34:DA:848:C:N3	2.37	0.69
1:AA:9:U:N3	1:AA:2641:A:H2	1.88	0.69
1:CA:1651:G:OP1	15:CR:40:LYS:NZ	2.25	0.69
1:CA:2059:A:OP2	63:CA:3916:HOH:O	2.11	0.69
1:CA:646:A:H2'	1:CA:647:G:O4'	1.93	0.69
1:AA:2255:U:OP1	63:AA:3945:HOH:O	2.10	0.68
1:CA:816:C:OP2	63:CA:4594:HOH:O	2.10	0.68
1:CA:2177:C:H5'	3:CC:45:HIS:HB3	1.75	0.68
4:AD:108:PRO:HD2	4:AD:111:LEU:HG	1.75	0.68
7:AG:83:ARG:N	7:AG:86:MET:SD	2.60	0.68
27:C3:39:ASP:OD2	27:C3:44:ARG:NH1	2.27	0.68
3:CC:183:PRO:HG2	3:CC:184:GLU:OE2	1.92	0.68
14:CQ:65:PHE:HB2	14:CQ:105:GLU:HB2	1.74	0.68
34:DA:1279:A:O2'	34:DA:1281:U:OP2	2.10	0.68
1:AA:121:G:OP2	63:AA:3922:HOH:O	2.11	0.68
1:AA:2564:U:OP2	63:AA:5097:HOH:O	2.11	0.68
1:AA:2825:C:H5'	29:A5:29:THR:HG21	1.74	0.68
3:AC:55:SER:O	3:AC:57:GLN:N	2.22	0.68
1:CA:1588:C:H2'	1:CA:1589:C:H6	1.58	0.68
1:CA:1860:G:H5'	3:CC:207:GLY:N	2.08	0.68
36:DC:122:GLU:O	36:DC:126:ARG:NH1	2.26	0.68
48:DO:16:ALA:HB1	48:DO:21:ASP:HB3	1.75	0.68
1:AA:2764:G:H4'	8:AH:4:ILE:HD11	1.76	0.68
23:AZ:45:ASP:OD1	23:AZ:49:ARG:HD2	1.92	0.68
1:CA:2286:A:H4'	1:CA:2287:A:O4'	1.94	0.68
1:CA:528:A:O2'	1:CA:529:A:H5''	1.94	0.68
1:CA:991:C:OP2	63:CA:4149:HOH:O	2.12	0.68
3:CC:25:GLU:HA	3:CC:28:ARG:HD2	1.74	0.68
16:CS:34:HIS:O	16:CS:97:ARG:NH2	2.26	0.68
34:DA:624:C:H2'	34:DA:625:G:H8	1.58	0.68
1:AA:243:G:O6	32:A8:5:LYS:HG2	1.94	0.68
34:DA:1129:C:N3	34:DA:1143:G:N2	2.33	0.68
34:DA:457:C:H2'	34:DA:458:C:H6	1.59	0.68
34:DA:522:C:H41	45:DL:53:ARG:HH22	1.39	0.68
51:DR:32:ARG:HA	51:DR:69:THR:HG21	1.76	0.68
3:AC:25:GLU:HA	3:AC:28:ARG:HD2	1.74	0.68
35:BB:21:ARG:HH21	35:BB:21:ARG:H	1.40	0.68
1:CA:1153:C:OP1	18:CU:92:ARG:NH1	2.20	0.68
3:AC:46:ALA:HB3	3:AC:172:ILE:HG22	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:AY:23:ARG:HG2	22:AY:42:VAL:HG22	1.74	0.68
34:BA:1238:A:OP2	63:BA:5228:HOH:O	2.12	0.68
38:BE:50:GLU:HB2	38:BE:53:LEU:HD13	1.76	0.68
49:BP:58:TYR:O	49:BP:61:SER:OG	2.11	0.68
57:BZ:99:ARG:O	57:BZ:101:LEU:N	2.27	0.68
1:CA:1336:A:OP2	21:CX:64:LYS:NZ	2.25	0.68
36:DC:40:ARG:NH2	36:DC:55:VAL:O	2.26	0.68
3:AC:15:VAL:O	3:AC:16:ASP:HB3	1.92	0.68
34:BA:259:G:H2'	34:BA:260:G:H8	1.59	0.68
40:BG:113:GLU:HB2	40:BG:119:ARG:HG2	1.76	0.68
53:BT:26:ASN:ND2	53:BT:71:THR:OG1	2.23	0.68
1:CA:1405:U:H2'	1:CA:1406:U:C6	2.29	0.68
3:CC:30:VAL:HG23	3:CC:31:LYS:N	2.09	0.68
3:CC:41:THR:O	3:CC:42:VAL:HB	1.94	0.68
34:DA:736:C:H2'	34:DA:737:A:C8	2.27	0.68
1:AA:1249:A:H2	1:AA:1287:A:H62	1.42	0.68
3:AC:30:VAL:HG23	3:AC:31:LYS:H	1.58	0.68
3:AC:46:ALA:HB3	3:AC:172:ILE:CG2	2.23	0.68
9:AK:70:GLU:O	9:AK:72:ASP:N	2.27	0.68
57:BZ:388:THR:HG23	57:BZ:399:LEU:HD22	1.74	0.68
25:C1:25:LYS:O	25:C1:27:GLU:N	2.27	0.68
1:CA:1784:A:O2'	63:CA:4517:HOH:O	2.11	0.68
23:CZ:45:ASP:OD1	23:CZ:49:ARG:NH1	2.27	0.68
41:DH:10:LEU:HD22	41:DH:83:ILE:HD11	1.76	0.68
57:DZ:285:ASP:N	57:DZ:285:ASP:OD2	2.25	0.68
57:DZ:533:VAL:HG12	57:DZ:534:ILE:HG13	1.76	0.68
1:AA:2623:U:H6	1:AA:2623:U:H5'	1.58	0.67
34:BA:406:G:N2	37:BD:119:GLN:HE22	1.92	0.67
7:CG:64:THR:HB	7:CG:94:LEU:HD21	1.76	0.67
23:CZ:183:LEU:O	23:CZ:185:GLU:N	2.28	0.67
23:CZ:5:LEU:HD23	23:CZ:47:VAL:HG21	1.76	0.67
1:AA:2771:A:OP2	63:AA:4053:HOH:O	2.12	0.67
35:BB:21:ARG:HB3	35:BB:39:ILE:HG12	1.75	0.67
57:BZ:316:ILE:HD12	57:BZ:326:THR:HG23	1.76	0.67
57:BZ:99:ARG:NH1	57:BZ:401:SER:O	2.27	0.67
34:DA:1335:C:O2	63:DA:3259:HOH:O	2.09	0.67
34:DA:1352:C:H2'	34:DA:1353:G:C8	2.29	0.67
1:AA:2775:G:OP2	63:AA:5117:HOH:O	2.12	0.67
23:AZ:31:ARG:NH1	23:AZ:94:GLU:OE1	2.27	0.67
34:BA:1369:C:H2'	34:BA:1370:G:C8	2.29	0.67
34:BA:67:C:H2'	34:BA:68:G:C8	2.29	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
51:BR:59:SER:H	51:BR:62:GLU:HG3	1.59	0.67
57:BZ:573:HIS:CD2	57:BZ:576:ASP:H	2.12	0.67
57:BZ:616:TYR:CE2	57:BZ:664:GLN:HG3	2.30	0.67
1:CA:2357:U:OP1	24:C0:20:ARG:NH1	2.28	0.67
5:CE:175:VAL:HG23	5:CE:177:PRO:HD3	1.77	0.67
34:DA:512:U:H2'	34:DA:513:C:C6	2.30	0.67
34:DA:512:U:H2'	34:DA:513:C:H6	1.58	0.67
34:DA:891:U:OP2	63:DA:3261:HOH:O	2.11	0.67
1:AA:2703:C:OP2	63:AA:4913:HOH:O	2.12	0.67
3:AC:30:VAL:HG23	3:AC:31:LYS:N	2.09	0.67
34:BA:738:C:H2'	34:BA:739:C:H6	1.58	0.67
1:CA:2684:U:H1'	12:CO:70:LYS:HD2	1.76	0.67
1:CA:2120:G:H21	3:CC:168:LYS:HE2	1.59	0.67
45:DL:80:HIS:ND1	57:DZ:425:SER:OG	2.26	0.67
52:DS:15:LEU:HD22	52:DS:33:THR:HB	1.76	0.67
57:DZ:-65:LYS:HB3	57:DZ:-28:ALA:HB3	1.75	0.67
38:BE:78:HIS:CD2	38:BE:142:LEU:HD23	2.29	0.67
1:CA:1269:A:OP2	63:CA:4473:HOH:O	2.11	0.67
1:CA:994:C:OP1	18:CU:53:ARG:NH2	2.28	0.67
34:DA:392:G:H2'	34:DA:393:A:H8	1.60	0.67
34:DA:542:G:OP1	37:DD:10:ARG:NH1	2.26	0.67
37:DD:153:ARG:O	37:DD:155:LEU:N	2.28	0.67
50:BQ:55:ASP:O	50:BQ:57:VAL:HG13	1.95	0.67
57:BZ:369:LEU:HD21	57:BZ:375:GLY:HA3	1.76	0.67
1:CA:330:A:HO2'	1:CA:331:A:H8	1.43	0.67
3:CC:30:VAL:HG23	3:CC:31:LYS:H	1.58	0.67
1:AA:1360:C:OP1	63:AA:4578:HOH:O	2.11	0.67
57:BZ:276:VAL:HG13	57:BZ:280:LEU:HD12	1.75	0.67
57:BZ:84:THR:HB	57:BZ:94:VAL:HG22	1.75	0.67
1:CA:2150:U:H2'	1:CA:2151:G:C8	2.30	0.67
22:CY:19:LYS:NZ	22:CY:20:TYR:OH	2.27	0.67
36:DC:47:LEU:HG	36:DC:68:VAL:HG11	1.76	0.67
1:AA:1133:G:N2	1:AA:1148:C:O2	2.25	0.67
1:AA:1846:A:OP1	1:AA:1846:A:H8	1.78	0.67
1:AA:2442:A:OP2	63:AA:4781:HOH:O	2.13	0.67
1:AA:553:A:C8	1:AA:553:A:C3'	2.76	0.67
41:BH:101:PRO:HG3	41:BH:133:LEU:HD11	1.76	0.67
45:BL:89:ARG:HA	45:BL:97:ARG:HA	1.77	0.67
57:BZ:275:ALA:HA	57:BZ:278:ASP:HB2	1.77	0.67
1:CA:1859:A:C2'	3:CC:206:LYS:CD	2.64	0.67
3:CC:46:ALA:HB3	3:CC:172:ILE:CG2	2.23	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DE:7:GLU:OE1	38:DE:37:ARG:NH2	2.20	0.67
49:DP:52:ASP:O	49:DP:54:GLU:N	2.28	0.67
53:DT:50:GLU:H	53:DT:99:LEU:HD12	1.60	0.67
56:DW:41:C:P	56:DY:36:A:OP1	2.53	0.67
38:BE:140:ARG:HG3	38:BE:140:ARG:O	1.93	0.67
1:CA:2693:A:H2'	1:CA:2694:G:H8	1.60	0.67
1:CA:572:A:N7	63:CA:4663:HOH:O	2.27	0.67
1:CA:668:G:H5'	1:CA:669:G:OP2	1.94	0.67
3:CC:55:SER:O	3:CC:57:GLN:N	2.22	0.67
57:DZ:120:THR:HG22	57:DZ:123:ARG:HH22	1.59	0.67
5:AE:40:GLU:CD	5:AE:40:GLU:H	1.97	0.67
34:BA:1137:C:H4'	34:BA:1138:G:C2	2.30	0.67
37:BD:157:LEU:O	37:BD:161:ASN:ND2	2.28	0.67
41:BH:51:VAL:HG12	41:BH:52:ASP:H	1.59	0.67
6:CF:101:LEU:HD12	6:CF:102:PRO:HD2	1.77	0.67
14:CQ:65:PHE:N	14:CQ:105:GLU:O	2.24	0.67
34:DA:352:C:O2'	34:DA:354:G:OP1	2.12	0.67
57:DZ:138:LYS:HG2	62:DZ:704:GDP:C6	2.29	0.67
1:AA:2562:G:OP1	63:AA:5041:HOH:O	2.13	0.66
1:AA:929:G:N2	1:AA:941:U:O2	2.28	0.66
3:CC:42:VAL:HG13	3:CC:43:GLU:N	2.10	0.66
7:CG:170:ARG:NH1	7:CG:174:GLU:OE2	2.27	0.66
34:DA:619:U:N3	37:DD:134:ASP:OD1	2.27	0.66
46:DM:25:ILE:HG23	46:DM:29:ARG:HB3	1.75	0.66
1:AA:1716:A:OP2	63:AA:5041:HOH:O	2.13	0.66
3:AC:41:THR:O	3:AC:42:VAL:HB	1.94	0.66
34:BA:1399:C:C2	34:BA:1502:A:N6	2.64	0.66
34:BA:403:C:H2'	34:BA:404:U:H6	1.61	0.66
1:CA:2450:A:O2'	56:DW:76:A:N6	2.28	0.66
1:AA:1336:C:H2'	1:AA:1337:C:H6	1.58	0.66
1:AA:2807:C:H42	1:AA:2813:G:H1	1.43	0.66
45:BL:59:ARG:HE	57:BZ:422:GLU:HG2	1.59	0.66
1:CA:2177:C:H5'	3:CC:45:HIS:CB	2.25	0.66
15:CR:74:LYS:HG2	15:CR:77:ARG:HH21	1.60	0.66
35:BB:16:HIS:O	35:BB:18:GLY:N	2.29	0.66
34:BA:881:G:P	45:BL:12:ARG:HH22	2.19	0.66
57:BZ:546:ILE:HG23	57:BZ:590:ILE:HG13	1.78	0.66
1:CA:2121:G:H1'	3:CC:168:LYS:HG2	1.76	0.66
3:CC:31:LYS:NZ	3:CC:180:SER:O	2.28	0.66
4:CD:242:ARG:HD3	4:CD:242:ARG:N	2.10	0.66
7:CG:13:GLU:O	7:CG:15:VAL:N	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:DG:133:GLY:HA2	40:DG:136:LYS:HB2	1.75	0.66
29:A5:16:ARG:HG2	29:A5:16:ARG:HH11	1.60	0.66
1:AA:664:U:H2'	1:AA:665:C:C6	2.30	0.66
39:BF:55:ASP:OD2	39:BF:86:ARG:NH1	2.26	0.66
43:BJ:50:ILE:HD11	43:BJ:57:LYS:HD3	1.77	0.66
61:BZ:703:FUA:H201	61:BZ:703:FUA:O1	1.95	0.66
1:CA:1919:A:N1	34:DA:1495:U:O2'	2.23	0.66
1:CA:2177:C:C1'	3:CC:171:ALA:CB	2.56	0.66
19:CV:25:LEU:H	19:CV:92:THR:HG1	1.41	0.66
3:AC:176:VAL:HG11	3:AC:190:ILE:HD13	1.76	0.66
17:AT:24:PRO:HA	17:AT:49:VAL:HG22	1.76	0.66
34:BA:17:U:H2'	34:BA:18:C:C6	2.31	0.66
1:CA:2025:C:H2'	1:CA:2026:C:C6	2.30	0.66
1:CA:2121:G:O4'	3:CC:168:LYS:CE	2.44	0.66
3:CC:176:VAL:HG11	3:CC:190:ILE:HD13	1.76	0.66
43:DJ:61:GLU:OE2	47:DN:49:HIS:NE2	2.27	0.66
1:AA:1525:G:O2'	1:AA:1605:A:H2	1.77	0.66
1:CA:195:A:N7	63:CA:4246:HOH:O	2.27	0.66
4:CD:68:LYS:HD2	4:CD:70:TRP:CZ2	2.31	0.66
34:DA:1311:G:N2	34:DA:1326:C:O2	2.27	0.66
1:CA:2422:A:O4'	56:DY:76:A:N6	2.29	0.66
4:AD:69:ARG:NH2	4:AD:128:GLY:O	2.26	0.66
34:BA:1304:G:OP2	63:BA:5206:HOH:O	2.14	0.66
57:BZ:603:GLU:HG2	57:BZ:679:VAL:HG12	1.77	0.66
1:CA:2632:A:HO2'	1:CA:2811:G:HO2'	1.27	0.66
5:CE:36:ARG:NH2	5:CE:88:GLY:O	2.29	0.66
34:DA:1279:A:OP2	43:DJ:9:ARG:NH1	2.29	0.66
1:AA:2146:G:N2	1:AA:2196:C:N3	2.34	0.66
3:AC:65:LEU:HB3	3:AC:189:ASN:ND2	2.11	0.66
38:BE:102:ALA:H	38:BE:107:ARG:NH2	1.94	0.66
1:CA:2695:C:H2'	1:CA:2696:U:H6	1.61	0.66
1:CA:858:U:O2	1:CA:2268:A:H2'	1.96	0.66
14:CQ:63:LYS:HA	23:CZ:178:GLU:HG3	1.77	0.66
34:DA:999:C:H42	34:DA:1042:G:H1	1.42	0.66
5:CE:36:ARG:HG2	5:CE:47:VAL:HG12	1.78	0.66
34:DA:975:A:H4'	34:DA:976:G:H5''	1.78	0.66
3:AC:42:VAL:HG13	3:AC:43:GLU:N	2.10	0.65
3:AC:63:VAL:O	3:AC:161:ARG:HA	1.96	0.65
34:BA:1047:G:O2'	34:BA:1215:G:O2'	2.13	0.65
1:CA:400:G:N7	63:CA:4346:HOH:O	2.28	0.65
10:CL:88:ALA:O	10:CL:90:LYS:N	2.29	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:CT:117:ASP:OD2	17:CT:120:ARG:NE	2.26	0.65
34:DA:222:U:H2'	34:DA:223:U:C6	2.31	0.65
57:DZ:13:ARG:HH21	57:DZ:247:ARG:NH1	1.94	0.65
1:AA:1100:A:N6	1:AA:1151:U:H3	1.92	0.65
36:BC:111:LEU:HB3	36:BC:204:LEU:HD21	1.77	0.65
48:BO:39:LEU:HD13	48:BO:56:LEU:HB2	1.78	0.65
56:BY:53:G:H1	56:BY:61:C:H42	1.42	0.65
20:CW:88:ARG:NH1	20:CW:94:ASP:OD2	2.29	0.65
57:DZ:601:ILE:HD12	57:DZ:684:GLN:HB2	1.78	0.65
34:BA:1329:A:H5''	46:BM:26:GLY:H	1.61	0.65
42:BI:45:ALA:HA	42:BI:48:GLU:HB2	1.77	0.65
1:CA:1019:U:H3	1:CA:1142(A):A:H62	1.42	0.65
1:CA:2583:G:OP2	63:CA:3973:HOH:O	2.14	0.65
12:CO:75:SER:HB2	17:CT:75:ILE:O	1.97	0.65
34:DA:229:U:O2'	49:DP:23:ASP:OD2	2.14	0.65
1:AA:2255:U:H2'	1:AA:2256:U:C6	2.32	0.65
3:AC:31:LYS:NZ	3:AC:180:SER:O	2.28	0.65
1:AA:2574:U:H1'	12:AO:23:ARG:HH11	1.60	0.65
3:CC:65:LEU:HB3	3:CC:189:ASN:ND2	2.11	0.65
45:DL:24:VAL:HG12	45:DL:98:TYR:CE1	2.32	0.65
34:DA:1312:G:H5'	52:DS:5:LEU:HD11	1.78	0.65
1:AA:1825:U:H2'	1:AA:1826:C:H6	1.60	0.65
1:AA:1891:G:H4'	3:AC:206:LYS:CG	2.26	0.65
1:AA:542:C:OP1	29:A5:16:ARG:NH2	2.29	0.65
3:AC:206:LYS:NZ	3:AC:206:LYS:HB3	2.12	0.65
16:AS:15:ARG:O	16:AS:19:LYS:HG2	1.97	0.65
36:BC:134:ILE:HG23	36:BC:151:VAL:HB	1.79	0.65
1:CA:2572:A:N7	5:CE:145:LYS:HB2	2.12	0.65
17:CT:24:PRO:HG3	17:CT:52:ILE:HG13	1.77	0.65
34:DA:812:C:N3	63:DA:3216:HOH:O	2.30	0.65
46:DM:124:PRO:HB3	57:DZ:500:GLN:HG3	1.78	0.65
61:DZ:703:FUA:H201	61:DZ:703:FUA:O1	1.95	0.65
34:BA:1445:C:O2'	34:BA:1447:A:N6	2.29	0.65
34:BA:259:G:H2'	34:BA:260:G:C8	2.31	0.65
34:BA:542:G:OP1	37:BD:10:ARG:NH2	2.26	0.65
41:BH:34:GLU:OE1	41:BH:37:ARG:NH2	2.30	0.65
57:BZ:498:ILE:HB	57:BZ:507:TYR:HD1	1.61	0.65
2:CB:105:A:P	23:CZ:72:ARG:HH12	2.18	0.65
43:DJ:30:SER:O	43:DJ:81:THR:OG1	2.10	0.65
50:DQ:95:TYR:HA	50:DQ:98:LEU:HD12	1.77	0.65
1:AA:139:A:H8	1:AA:1454:C:HO2'	1.42	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AH:149:ARG:NH1	8:AH:167:GLU:OE2	2.29	0.65
43:BJ:17:ASP:OD1	43:BJ:70:ARG:NH1	2.30	0.65
1:CA:614(B):G:H2'	6:CF:44:ARG:HH11	1.61	0.65
11:CN:58:ASP:OD1	11:CN:125:GLY:N	2.24	0.65
43:DJ:6:ILE:HG12	43:DJ:98:ILE:HG12	1.78	0.65
1:AA:1067:A:H3'	1:AA:1067:A:C8	2.32	0.65
1:AA:1117:G:O2'	1:AA:1135:G:OP2	2.13	0.65
4:AD:77:ALA:O	4:AD:116:GLN:HA	1.97	0.65
42:BI:40:LEU:O	42:BI:42:ARG:N	2.29	0.65
57:BZ:120:THR:HG22	57:BZ:123:ARG:NH1	2.12	0.65
1:CA:1316:U:H2'	1:CA:1317:A:C8	2.31	0.65
1:CA:2316:C:H2'	1:CA:2317:C:C6	2.30	0.65
3:CC:63:VAL:O	3:CC:161:ARG:HA	1.96	0.65
11:CN:15:LEU:HD23	11:CN:53:VAL:HB	1.79	0.65
34:DA:881:G:P	45:DL:12:ARG:HH22	2.20	0.65
37:DD:129:ASN:ND2	37:DD:145:GLU:H	1.95	0.65
57:DZ:90:PHE:CE1	61:DZ:703:FUA:H122	2.31	0.65
1:AA:2227:G:H5'	1:AA:2228:G:N7	2.11	0.65
1:AA:638:U:H6	1:AA:638:U:O5'	1.79	0.65
14:AQ:21:THR:HG21	14:AQ:101:ARG:HB2	1.78	0.65
34:BA:757:U:H2'	34:BA:758:G:O4'	1.97	0.65
41:BH:81:HIS:N	41:BH:138:TRP:O	2.30	0.65
3:CC:206:LYS:NZ	3:CC:206:LYS:HB3	2.12	0.65
16:CS:63:THR:HG23	16:CS:64:GLU:H	1.62	0.65
1:AA:426:G:OP2	63:AA:4920:HOH:O	2.14	0.65
8:AH:159:GLU:HG3	8:AH:169:VAL:HG11	1.79	0.65
34:BA:1015:A:N3	34:BA:1218:C:O2'	2.30	0.65
1:CA:2176:A:O2'	3:CC:45:HIS:CD2	2.50	0.65
7:CG:28:VAL:O	7:CG:31:VAL:HG12	1.97	0.65
34:DA:1227:A:OP1	52:DS:80:TYR:OH	2.07	0.65
34:DA:235:C:H5'	50:DQ:70:ARG:HG2	1.79	0.65
35:DB:74:LYS:NZ	35:DB:206:ASP:OD1	2.30	0.65
42:DI:53:VAL:O	42:DI:55:ALA:N	2.31	0.65
57:DZ:181:LEU:HD12	57:DZ:216:LEU:HD21	1.78	0.65
57:DZ:505:GLY:HA2	57:DZ:576:ASP:HB2	1.79	0.65
57:DZ:631:ILE:HD12	57:DZ:645:ALA:HB2	1.79	0.65
24:A0:32:ARG:H	24:A0:35:ASN:ND2	1.96	0.64
10:AL:13:PRO:HA	10:AL:52:ILE:HG12	1.79	0.64
34:BA:164:U:H2'	34:BA:165:C:C6	2.32	0.64
1:CA:1803:A:O2'	4:CD:259:THR:HG21	1.96	0.64
1:CA:831:G:O2'	13:CP:38:GLN:NE2	2.30	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:DC:47:LEU:HB3	36:DC:52:LEU:HB3	1.79	0.64
1:AA:1356:G:OP2	31:A7:9:ARG:NH1	2.29	0.64
7:AG:173:LEU:HA	7:AG:176:LEU:HD12	1.80	0.64
33:C9:14:CYS:HA	33:C9:27:CYS:HB2	1.80	0.64
49:DP:43:LYS:HG2	49:DP:48:TRP:CG	2.31	0.64
56:DY:50:U:H3	56:DY:64:A:H61	1.45	0.64
1:AA:1829:U:OP2	4:AD:274:ARG:NH2	2.30	0.64
1:AA:553:A:C2'	1:AA:554:A:H5'	2.26	0.64
30:C6:9:LEU:HA	30:C6:54:ILE:HB	1.79	0.64
1:CA:2268:A:OP1	63:CA:4086:HOH:O	2.14	0.64
23:CZ:93:ASP:HB2	23:CZ:131:ARG:HH22	1.62	0.64
34:DA:1207:G:H2'	34:DA:1208:C:C6	2.32	0.64
35:DB:95:GLN:HG3	35:DB:148:TYR:HA	1.78	0.64
1:AA:1848:G:H2'	1:AA:1849:U:H5'	1.79	0.64
34:BA:108:G:C6	53:BT:15:ARG:HD2	2.33	0.64
1:CA:2124:G:O3'	3:CC:175:PRO:HG2	1.98	0.64
1:CA:527:C:OP1	63:CA:4143:HOH:O	2.15	0.64
10:CL:79:ARG:NH1	10:CL:85:GLU:O	2.31	0.64
11:CN:62:VAL:HG21	11:CN:87:LEU:HD11	1.79	0.64
34:DA:1097:C:O2'	34:DA:1169:A:N3	2.30	0.64
34:DA:179:A:H2'	34:DA:180:U:C6	2.32	0.64
37:DD:171:GLY:O	37:DD:174:LEU:N	2.30	0.64
50:DQ:32:TYR:O	50:DQ:34:LYS:N	2.30	0.64
57:DZ:127:LYS:HZ3	57:DZ:404:VAL:HG11	1.62	0.64
57:DZ:518:PRO:O	57:DZ:521:SER:OG	2.12	0.64
34:BA:1014:A:H2'	34:BA:1015:A:C8	2.33	0.64
34:BA:1442:G:O2'	34:BA:1442(A):G:OP1	2.16	0.64
34:BA:973:G:H3'	34:BA:974:A:H5''	1.80	0.64
48:BO:17:ARG:HG3	48:BO:17:ARG:HH11	1.62	0.64
25:C1:8:SER:HB3	25:C1:66:HIS:CD2	2.32	0.64
5:CE:97:LYS:N	5:CE:100:GLU:OE1	2.19	0.64
41:DH:104:ARG:NH2	63:DH:4001:HOH:O	2.22	0.64
57:DZ:532:GLY:O	57:DZ:534:ILE:N	2.31	0.64
1:AA:2510:C:OP2	63:AA:4620:HOH:O	2.14	0.64
3:AC:69:LEU:O	3:AC:178:LYS:HG3	1.97	0.64
8:AH:54:ARG:HD3	8:AH:65:HIS:ND1	2.12	0.64
11:AN:74:ARG:HH12	11:AN:85:ILE:CD1	2.10	0.64
17:AT:23:ARG:HG3	17:AT:120:ARG:CZ	2.28	0.64
30:C6:43:CYS:O	30:C6:45:LYS:N	2.30	0.64
1:CA:1466:G:O2'	1:CA:1546:C:O2'	2.13	0.64
34:DA:978:A:O2'	34:DA:1322:C:N3	2.29	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DI:46:ALA:HA	42:DI:78:LYS:HB2	1.79	0.64
49:DP:23:ASP:OD1	49:DP:25:ARG:HD3	1.97	0.64
1:AA:2860:A:OP2	1:AA:2876:U:H5	1.79	0.64
34:BA:1077:G:N2	34:BA:1080:A:OP2	2.30	0.64
57:BZ:99:ARG:C	57:BZ:101:LEU:H	2.00	0.64
1:CA:1237:A:OP1	63:CA:4442:HOH:O	2.15	0.64
1:CA:1588:C:H2'	1:CA:1589:C:C6	2.32	0.64
1:CA:848:G:H2'	1:CA:849:A:C8	2.31	0.64
8:CH:18:GLU:HB3	8:CH:25:LYS:HB2	1.77	0.64
1:CA:2820:A:C4	15:CR:4:LEU:HD21	2.32	0.64
16:CS:35:ILE:HD12	16:CS:101:LEU:HD12	1.78	0.64
34:DA:1435:G:H2'	34:DA:1436:U:C6	2.33	0.64
1:AA:1233:U:H4'	19:AV:79:VAL:HG22	1.78	0.64
1:AA:1475:G:H2'	1:AA:1476:C:C6	2.32	0.64
1:AA:1634:C:H2'	1:AA:1635:C:H6	1.62	0.64
1:AA:2052:A:H5''	1:AA:2053:A:OP1	1.98	0.64
1:AA:483:A:OP1	63:AA:5245:HOH:O	2.15	0.64
3:AC:44:VAL:CG2	3:AC:176:VAL:HG21	2.28	0.64
23:AZ:145:GLU:N	23:AZ:148:ASP:OD2	2.27	0.64
35:BB:195:ASP:O	41:BH:68:ARG:NH2	2.31	0.64
56:BY:8:4SU:H4'	56:BY:48:C:H4'	1.79	0.64
1:CA:2456:C:N4	63:CA:4095:HOH:O	2.30	0.64
8:CH:38:SER:HB3	8:CH:41:MET:HG2	1.80	0.64
13:CP:38:GLN:HG2	13:CP:45:LEU:H	1.62	0.64
34:DA:922:G:H2'	34:DA:923:A:C8	2.33	0.64
1:AA:2434:A:O4'	56:BY:76:A:N6	2.31	0.64
34:BA:390:C:H2'	34:BA:391:G:H8	1.63	0.64
1:CA:2218:U:O4'	25:C1:52:ARG:NH2	2.31	0.64
10:CL:30:HIS:HA	10:CL:59:ILE:HD12	1.80	0.64
1:AA:1675:U:O4	63:AA:4127:HOH:O	2.12	0.64
1:AA:2832:G:OP2	63:AE:415:HOH:O	2.15	0.64
34:BA:435:C:H2'	34:BA:436:C:H6	1.62	0.64
37:BD:155:LEU:HD13	37:BD:158:ILE:HD11	1.79	0.64
1:CA:2096:U:H2'	1:CA:2097:C:C6	2.33	0.64
4:CD:71:ASP:CG	4:CD:103:ARG:HH22	2.01	0.64
1:CA:997:G:OP1	18:CU:92:ARG:HG2	1.97	0.64
34:DA:517:G:N2	34:DA:533:A:OP2	2.31	0.64
37:DD:165:MET:SD	37:DD:168:ARG:NH1	2.60	0.64
37:DD:57:ARG:NH2	37:DD:205:GLU:OE2	2.26	0.64
11:AN:46:VAL:HG23	11:AN:48:MET:HG2	1.80	0.63
34:BA:626:U:H2'	34:BA:627:G:H8	1.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2064:C:OP2	63:CA:4245:HOH:O	2.14	0.63
1:CA:2168:G:H2'	1:CA:2169:A:H8	1.62	0.63
3:CC:68:GLY:N	3:CC:189:ASN:HD21	1.96	0.63
3:CC:29:LEU:O	3:CC:32:GLU:N	2.31	0.63
3:CC:69:LEU:O	3:CC:178:LYS:HG3	1.97	0.63
34:DA:673:G:H2'	34:DA:674:G:C8	2.33	0.63
39:DF:23:LYS:NZ	39:DF:42:GLU:OE2	2.31	0.63
57:DZ:179:ASP:OD2	57:DZ:182:ARG:HD2	1.97	0.63
57:DZ:168:ILE:HG23	57:DZ:205:TYR:HE2	1.63	0.63
57:DZ:517:LEU:HG	57:DZ:518:PRO:HD2	1.80	0.63
1:AA:303:C:H42	1:AA:385:G:H1	1.46	0.63
3:AC:29:LEU:O	3:AC:32:GLU:N	2.32	0.63
37:BD:57:ARG:NH2	37:BD:205:GLU:OE2	2.31	0.63
30:C6:13:CYS:SG	30:C6:47:THR:HG21	2.39	0.63
1:CA:2867:G:OP2	17:CT:119:LYS:NZ	2.32	0.63
34:DA:1143:G:H2'	34:DA:1144:G:H8	1.63	0.63
34:DA:403:C:H2'	34:DA:404:U:H6	1.64	0.63
44:DK:79:SER:HA	44:DK:104:GLN:HB3	1.79	0.63
46:DM:33:ALA:HA	46:DM:59:TYR:HE2	1.63	0.63
1:AA:671:A:H2'	1:AA:672:G:O4'	1.99	0.63
3:AC:68:GLY:N	3:AC:189:ASN:HD21	1.96	0.63
6:AF:32:LEU:HB3	6:AF:112:MET:HE1	1.80	0.63
34:BA:1063:C:H2'	34:BA:1064:G:C8	2.32	0.63
46:BM:90:LEU:HD23	46:BM:93:ARG:HD2	1.79	0.63
1:AA:1935:A:OP1	57:BZ:499:ARG:NH2	2.31	0.63
16:CS:84:GLN:HA	16:CS:111:GLU:HB2	1.80	0.63
17:CT:16:ARG:HH11	17:CT:19:LEU:HD21	1.61	0.63
34:DA:976:G:OP2	34:DA:1358:U:O2'	2.15	0.63
56:DW:40:C:H1'	56:DY:35:A:H4'	1.79	0.63
57:DZ:230:LYS:HB3	57:DZ:235:GLU:HB3	1.78	0.63
1:AA:2367:C:H1'	24:A0:39:ARG:HH21	1.62	0.63
3:AC:7:ARG:O	3:AC:11:LEU:HD23	1.99	0.63
53:BT:56:MET:CE	53:BT:85:MET:HA	2.29	0.63
1:CA:2037:G:O6	63:CA:4152:HOH:O	2.11	0.63
3:CC:44:VAL:CG2	3:CC:176:VAL:HG21	2.28	0.63
5:CE:143:ASN:HD22	5:CE:147:PRO:HD3	1.63	0.63
5:CE:73:GLU:OE2	5:CE:73:GLU:N	2.31	0.63
34:DA:148:G:H2'	34:DA:149:A:H8	1.62	0.63
28:A4:53:GLU:HB3	28:A4:54:GLY:HA2	1.80	0.63
1:AA:1296:G:OP2	13:AP:21:ARG:NH1	2.31	0.63
1:AA:2326:C:H2'	1:AA:2327:G:H8	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
22:AY:8:LYS:HD3	22:AY:97:ARG:NH1	2.14	0.63
48:BO:16:ALA:HB1	48:BO:21:ASP:HB3	1.79	0.63
1:CA:2646:C:OP2	1:CA:2732:G:O2'	2.15	0.63
7:AG:143:GLU:O	28:A4:28:LYS:NZ	2.30	0.63
1:AA:1701:A:OP1	15:AR:1:MET:N	2.25	0.63
57:BZ:264:LEU:HB2	62:BZ:704:GDP:C5	2.33	0.63
8:CH:113:VAL:HG11	8:CH:151:ILE:HG21	1.81	0.63
18:CU:49:HIS:HA	18:CU:52:ARG:HB3	1.80	0.63
34:DA:811:C:O2'	34:DA:901:A:N1	2.30	0.63
35:DB:16:HIS:CG	35:DB:17:PHE:H	2.16	0.63
34:DA:404:U:H5'	37:DD:122:ARG:HE	1.61	0.63
57:DZ:497:PHE:CE2	57:DZ:506:GLN:HG3	2.32	0.63
1:AA:354:A:H2	1:AA:1255:A:HO2'	1.46	0.63
34:BA:921:U:O2	38:BE:19:MET:HB2	1.99	0.63
57:BZ:629:GLY:HA3	57:BZ:647:VAL:HG12	1.79	0.63
57:BZ:-9:LEU:O	57:BZ:-6:ARG:N	2.20	0.63
1:CA:214:G:O2'	1:CA:216:A:O2'	2.08	0.63
23:CZ:104:PHE:HA	23:CZ:139:VAL:HG22	1.80	0.63
34:DA:932:C:H2'	34:DA:933:G:C8	2.33	0.63
30:A6:35:GLU:OE2	30:A6:50:ARG:NH1	2.31	0.63
2:AB:66:A:H61	2:AB:108:U:H2'	1.64	0.63
9:AK:28:ASN:O	9:AK:30:GLN:N	2.31	0.63
57:BZ:239:GLU:O	57:BZ:241:GLU:N	2.32	0.63
57:BZ:508:GLY:HA3	57:BZ:581:ALA:O	1.98	0.63
1:CA:830:G:H4'	1:CA:831:G:OP2	1.97	0.63
3:CC:7:ARG:O	3:CC:11:LEU:HD23	1.99	0.63
15:CR:57:ARG:O	15:CR:59:ASP:N	2.32	0.63
36:DC:113:ALA:HB2	36:DC:202:ILE:HG13	1.79	0.63
57:DZ:497:PHE:O	57:DZ:507:TYR:HA	1.99	0.63
1:AA:1003:U:OP2	14:AQ:14:ARG:HD3	1.99	0.63
1:AA:2152:U:H4'	1:AA:2155:G:H4'	1.80	0.63
8:AH:88:LEU:HD12	8:AH:130:ARG:HG2	1.79	0.63
10:AL:100:THR:HA	10:AL:139:VAL:HB	1.81	0.63
35:BB:55:PHE:CD1	35:BB:58:ILE:HD12	2.34	0.63
40:BG:31:MET:SD	40:BG:34:GLY:HA2	2.39	0.63
1:CA:1859:A:C2'	3:CC:206:LYS:CE	2.77	0.63
1:CA:486:C:O2'	20:CW:60:ASN:OD1	2.12	0.63
34:DA:1456:G:O3'	53:DT:39:LYS:NZ	2.31	0.63
56:DW:55:PSU:O2'	56:DW:57:G:N7	2.20	0.63
57:DZ:215:LYS:HA	57:DZ:218:GLU:HB3	1.81	0.63
57:DZ:524:GLU:HB3	57:DZ:564:LYS:HA	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2331:G:H22	16:AS:3:ARG:CG	2.06	0.62
1:AA:1324:A:OP1	15:AR:36:THR:HG22	1.98	0.62
1:CA:1705:G:O6	63:CA:4615:HOH:O	2.13	0.62
1:CA:340:A:H2'	1:CA:341:G:O4'	1.99	0.62
7:CG:80:PHE:O	7:CG:82:LEU:N	2.31	0.62
10:CL:71:THR:O	10:CL:111:LYS:NZ	2.32	0.62
11:CN:42:TRP:CH2	11:CN:44:PRO:HB3	2.34	0.62
14:CQ:38:GLU:HG3	14:CQ:127:ILE:HG22	1.80	0.62
34:DA:437:U:O2'	37:DD:125:HIS:HE1	1.81	0.62
37:DD:23:GLY:N	37:DD:26:CYS:SG	2.61	0.62
54:DU:3:LYS:HB3	54:DU:14:TRP:CD1	2.33	0.62
1:AA:482:C:H4'	63:AA:5245:HOH:O	1.99	0.62
10:AL:45:THR:O	10:AL:45:THR:OG1	2.16	0.62
20:AW:4:LYS:HB2	20:AW:106:ILE:HG12	1.82	0.62
34:BA:1240:U:OP2	40:BG:116:ALA:N	2.26	0.62
57:BZ:356:LEU:HD12	57:BZ:365:GLU:HA	1.79	0.62
57:BZ:549:ALA:HB1	57:BZ:591:LYS:HE2	1.81	0.62
1:CA:1913:A:H4'	1:CA:1914:C:H5'	1.80	0.62
57:DZ:332:SER:HA	57:DZ:371:ALA:HB2	1.79	0.62
1:AA:714:U:O2	32:A8:2:PRO:HD2	1.98	0.62
1:AA:831:A:OP2	63:AA:4559:HOH:O	2.15	0.62
2:AB:66:A:N6	2:AB:108:U:H2'	2.15	0.62
36:BC:35:GLU:OE2	36:BC:59:ARG:NH2	2.31	0.62
34:BA:437:U:O2'	37:BD:125:HIS:HE1	1.81	0.62
1:CA:922:U:H2'	1:CA:923:C:C6	2.33	0.62
1:CA:1859:A:C2'	3:CC:206:LYS:HE3	2.29	0.62
3:CC:6:LYS:HG3	3:CC:7:ARG:N	2.14	0.62
8:CH:45:VAL:HG22	8:CH:50:VAL:HG22	1.81	0.62
1:CA:328:U:H4'	22:CY:68:HIS:CD2	2.34	0.62
34:DA:1228:C:OP1	46:DM:115:LYS:N	2.29	0.62
34:DA:473:G:H2'	34:DA:474:G:H8	1.62	0.62
4:AD:223:GLY:HA3	4:AD:231:HIS:CE1	2.34	0.62
6:AF:24:LEU:HB3	6:AF:115:ALA:HB2	1.79	0.62
20:AW:14:PRO:CG	20:AW:78:GLU:HG2	2.29	0.62
34:BA:565:U:OP2	34:BA:566:G:O2'	2.14	0.62
35:BB:55:PHE:HD1	35:BB:58:ILE:HD12	1.65	0.62
1:CA:1815:A:OP2	4:CD:54:ARG:NH2	2.32	0.62
35:DB:13:ALA:N	35:DB:14:GLY:HA3	2.14	0.62
47:DN:27:CYS:SG	47:DN:29:ARG:N	2.73	0.62
1:AA:2175:G:H2'	1:AA:2176:G:H8	1.63	0.62
4:AD:145:VAL:HG12	4:AD:146:GLU:O	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:AU:66:ASN:O	18:AU:70:ARG:HG3	2.00	0.62
34:BA:1238:A:O5'	34:BA:1336:C:N4	2.33	0.62
37:BD:57:ARG:HG2	37:BD:202:LEU:HD22	1.81	0.62
34:BA:1316:G:H4'	47:BN:18:VAL:HG13	1.81	0.62
1:CA:813:U:H2'	1:CA:814:C:C6	2.34	0.62
34:DA:560:U:O2'	34:DA:561:U:OP2	2.18	0.62
57:DZ:-9:LEU:O	57:DZ:-6:ARG:N	2.32	0.62
28:A4:10:VAL:HG21	28:A4:29:PRO:HG3	1.81	0.62
1:AA:116:A:C8	1:AA:117:A:C8	2.87	0.62
1:AA:1627:A:H8	1:AA:1627:A:OP2	1.82	0.62
3:AC:53:ARG:HD3	3:AC:53:ARG:H	1.65	0.62
34:BA:499:A:O2'	34:BA:546:G:N2	2.32	0.62
50:BQ:6:LEU:HD23	50:BQ:23:VAL:HG11	1.80	0.62
28:C4:44:THR:O	28:C4:46:GLN:N	2.31	0.62
3:CC:68:GLY:H	3:CC:189:ASN:HD21	1.47	0.62
1:AA:553:A:H2	1:AA:2065:C:H5'	1.63	0.62
1:AA:552:C:C5	1:AA:2792:U:H2'	2.35	0.62
3:AC:6:LYS:HG3	3:AC:7:ARG:N	2.14	0.62
36:BC:116:VAL:HG21	36:BC:202:ILE:HD11	1.82	0.62
57:BZ:432:ALA:HA	57:BZ:438:PHE:HE1	1.65	0.62
1:CA:2124:G:C3'	3:CC:175:PRO:HG3	2.30	0.62
1:CA:307:G:N1	1:CA:310:A:OP2	2.29	0.62
6:CF:118:ALA:HB2	6:CF:123:LEU:HD23	1.81	0.62
9:CK:104:ILE:HA	9:CK:109:SER:HA	1.82	0.62
1:CA:2562:U:H1'	12:CO:23:ARG:HH11	1.63	0.62
34:DA:1292:U:OP2	40:DG:41:ARG:NH2	2.32	0.62
34:DA:1339:A:O3'	56:DY:35:A:OP1	2.16	0.62
57:DZ:87:HIS:CD2	61:DZ:703:FUA:H283	2.31	0.62
1:AA:1891:G:H4'	3:AC:206:LYS:HD3	1.79	0.62
34:BA:501:C:H2'	34:BA:502:G:H8	1.64	0.62
35:BB:166:ASP:HB3	35:BB:169:LYS:HB2	1.81	0.62
34:DA:1347:G:H22	34:DA:1373:G:H2'	1.62	0.62
34:DA:9:G:OP2	38:DE:121:LYS:NZ	2.31	0.62
36:DC:11:ARG:HB3	36:DC:15:THR:HB	1.82	0.62
1:AA:215:G:H21	1:AA:217:A:H62	1.48	0.62
40:BG:36:LYS:HA	40:BG:39:ALA:HB3	1.81	0.62
61:BZ:703:FUA:O2	61:BZ:703:FUA:H211	2.00	0.62
1:CA:1671:U:HO2'	1:CA:1673:U:H5	1.48	0.62
5:CE:47:VAL:HG11	5:CE:86:PRO:HD2	1.81	0.62
51:DR:58:LEU:HD12	51:DR:62:GLU:HB3	1.82	0.62
6:AF:148:LEU:HD13	6:AF:154:VAL:HG21	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:1497:G:O6	63:BA:5225:HOH:O	2.15	0.62
38:BE:31:LEU:HD23	38:BE:45:PHE:HB2	1.81	0.62
1:CA:1877:A:H5'	1:CA:1878:G:OP2	2.00	0.62
61:DZ:703:FUA:O2	61:DZ:703:FUA:H211	2.00	0.62
57:DZ:138:LYS:HG2	62:DZ:704:GDP:C5	2.34	0.62
1:AA:1825:U:H2'	1:AA:1826:C:C6	2.34	0.61
14:AQ:135:ASP:HB3	14:AQ:137:TYR:H	1.64	0.61
54:BU:3:LYS:HB3	54:BU:14:TRP:CG	2.35	0.61
1:CA:2632:A:O2'	1:CA:2811:G:O2'	2.05	0.61
1:CA:2892:A:H2'	1:CA:2893:G:H8	1.65	0.61
3:CC:53:ARG:HD3	3:CC:53:ARG:H	1.65	0.61
1:CA:1665:A:H1'	12:CO:1:MET:HG3	1.82	0.61
12:CO:4:PRO:O	12:CO:5:GLN:HB2	1.99	0.61
34:DA:1016:A:HO2'	34:DA:1217:C:HO2'	1.44	0.61
34:DA:1468:A:H2'	34:DA:1469:G:O4'	1.99	0.61
45:DL:113:ARG:HG2	45:DL:117:ARG:HG2	1.80	0.61
34:DA:255:G:H1'	50:DQ:16:GLN:NE2	2.15	0.61
3:AC:68:GLY:H	3:AC:189:ASN:HD21	1.47	0.61
7:AG:180:PHE:O	7:AG:182:LYS:N	2.33	0.61
57:BZ:210:ARG:HB2	57:BZ:210:ARG:NH1	2.12	0.61
35:DB:55:PHE:HA	35:DB:58:ILE:HG13	1.81	0.61
38:DE:77:PRO:HD2	38:DE:142:LEU:HD22	1.82	0.61
57:DZ:328:ILE:HD12	57:DZ:377:VAL:HG12	1.81	0.61
57:DZ:363:ARG:CG	57:DZ:363:ARG:HH11	2.13	0.61
20:AW:13:SER:HB3	20:AW:16:LYS:HD2	1.82	0.61
34:BA:1318:A:H2'	34:BA:1319:A:H5''	1.82	0.61
34:BA:345:C:H4'	34:BA:346:G:C2	2.35	0.61
25:C1:83:GLU:OE1	25:C1:83:GLU:N	2.33	0.61
2:CB:76:G:H2'	2:CB:77:U:O4'	2.00	0.61
6:CF:20:LEU:HD13	6:CF:21:ALA:H	1.65	0.61
34:DA:1059:C:OP2	36:DC:199:LYS:NZ	2.33	0.61
34:DA:353:A:H5'	34:DA:353:A:H8	1.66	0.61
1:AA:868:A:H2'	1:AA:991:G:H5''	1.82	0.61
6:AF:191:ARG:HG2	6:AF:191:ARG:NH1	2.11	0.61
34:BA:1013:G:N2	34:BA:1016:A:OP2	2.32	0.61
42:BI:110:GLU:OE2	42:BI:113:LYS:NZ	2.33	0.61
46:BM:84:ILE:HB	52:BS:74:PHE:HE1	1.65	0.61
6:CF:120:GLU:HB2	6:CF:122:LYS:HG2	1.81	0.61
36:DC:111:LEU:HD22	36:DC:146:ALA:HB2	1.83	0.61
37:DD:187:ARG:NH2	37:DD:193:ASP:OD2	2.34	0.61
34:DA:559:A:OP1	38:DE:126:ARG:NH2	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:DE:57:LYS:HD3	38:DE:61:TYR:HE2	1.64	0.61
57:DZ:177:ILE:HG13	57:DZ:188:TYR:HE2	1.64	0.61
57:DZ:223:PHE:CZ	57:DZ:249:GLY:HA3	2.35	0.61
30:A6:25:LYS:HE3	30:A6:30:THR:O	2.00	0.61
1:AA:1935:A:H4'	1:AA:1936:C:H5''	1.81	0.61
1:AA:1935:A:H4'	1:AA:1936:C:C5'	2.30	0.61
34:BA:1347:G:H5''	42:BI:107:ARG:HB3	1.81	0.61
34:BA:277:C:P	50:BQ:41:LYS:HZ1	2.23	0.61
55:BV:15:A:O2'	55:BV:16:U:C5	2.39	0.61
34:DA:1494:G:H5''	57:DZ:506:GLN:HE22	1.65	0.61
34:BA:347:G:H2'	34:BA:348:G:O4'	1.99	0.61
34:BA:407:G:H5''	37:BD:115:ARG:HB3	1.83	0.61
45:BL:34:ARG:HG2	45:BL:35:GLY:N	2.16	0.61
1:CA:2631:G:N3	1:CA:2810:A:H2	1.98	0.61
10:CL:64:SER:OG	10:CL:65:PHE:N	2.33	0.61
57:DZ:264:LEU:HD12	62:DZ:704:GDP:C4	2.35	0.61
5:AE:116:VAL:HG13	5:AE:122:PHE:HB2	1.83	0.61
7:AG:124:SER:HB2	7:AG:131:TYR:CE1	2.36	0.61
7:AG:140:ILE:HD12	7:AG:140:ILE:H	1.66	0.61
12:AO:64:ARG:HD3	17:AT:70:VAL:HG11	1.83	0.61
18:AU:76:TYR:CE1	18:AU:80:ILE:HG13	2.36	0.61
34:BA:1372:U:OP1	42:BI:72:GLY:N	2.33	0.61
49:BP:67:THR:HB	49:BP:70:ALA:HB2	1.82	0.61
57:DZ:74:TRP:NE1	57:DZ:273:LEU:HB3	2.16	0.61
1:AA:2023:A:H2'	1:AA:2024:G:C8	2.36	0.61
20:AW:14:PRO:HG2	20:AW:78:GLU:CG	2.30	0.61
34:BA:308:C:H2'	34:BA:309:G:H8	1.66	0.61
50:BQ:86:GLU:O	50:BQ:88:TYR:N	2.33	0.61
51:BR:32:ARG:HA	51:BR:69:THR:HG21	1.83	0.61
61:BZ:703:FUA:H12	61:BZ:703:FUA:O1	2.00	0.61
1:CA:1803:A:H4'	4:CD:259:THR:HG23	1.83	0.61
3:CC:11:LEU:HD12	3:CC:33:LEU:HA	1.82	0.61
4:CD:238:GLY:O	4:CD:239:ARG:HB2	1.98	0.61
50:DQ:10:VAL:HG13	50:DQ:19:VAL:HB	1.83	0.61
56:DW:47:U:O2'	56:DW:48:C:OP1	2.18	0.61
1:CA:1095:A:N6	57:DZ:614:GLU:OE1	2.33	0.61
1:AA:2658:C:H2'	1:AA:2659:U:O4'	2.01	0.61
45:BL:33:ARG:HD3	45:BL:62:SER:HB3	1.83	0.61
57:BZ:512:ILE:HD12	57:BZ:589:ALA:HB1	1.82	0.61
1:CA:338:G:N7	63:CA:3740:HOH:O	2.31	0.61
1:CA:2121:G:O2'	3:CC:168:LYS:CG	2.49	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:CG:126:ASP:HB3	7:CG:128:ARG:H	1.66	0.61
17:CT:83:ILE:HD13	17:CT:86:ILE:HD11	1.83	0.61
49:DP:53:VAL:HG22	49:DP:79:VAL:HG22	1.82	0.61
25:A1:49:VAL:HG21	25:A1:67:ILE:HG23	1.81	0.61
21:AX:65:ARG:HB2	21:AX:70:LEU:HD22	1.83	0.61
22:AY:92:ASN:H	22:AY:92:ASN:HD22	1.49	0.61
34:BA:1227:A:O3'	46:BM:115:LYS:HD2	2.01	0.61
35:BB:178:ARG:NH1	35:BB:196:LEU:O	2.34	0.61
1:CA:775:G:N2	1:CA:794:G:H5'	2.15	0.61
34:DA:407:G:H5''	37:DD:115:ARG:HD2	1.81	0.61
42:DI:3:GLN:OE1	42:DI:20:ARG:NH2	2.30	0.61
56:DW:9:A:O2'	56:DW:10:G:N7	2.34	0.61
57:DZ:169:GLY:N	57:DZ:170:ARG:HH12	1.99	0.61
1:AA:2451:A:C8	1:AA:2451:A:H5'	2.36	0.60
1:AA:2695:C:O2	12:AO:70:LYS:NZ	2.29	0.60
1:AA:591:U:OP1	13:AP:36:LYS:HE3	2.01	0.60
2:AB:45:A:OP2	7:AG:96:ARG:NH2	2.30	0.60
1:AA:1125:C:O2'	10:AL:132:ARG:NH1	2.34	0.60
17:AT:60:THR:HG22	17:AT:77:PRO:HA	1.83	0.60
34:BA:1055:A:H2'	36:BC:156:ARG:HD2	1.83	0.60
40:BG:69:VAL:HG22	40:BG:135:VAL:HG22	1.82	0.60
40:BG:92:SER:O	40:BG:95:ARG:N	2.32	0.60
51:BR:59:SER:H	51:BR:62:GLU:CG	2.14	0.60
1:CA:2168:G:H2'	1:CA:2169:A:C8	2.36	0.60
34:DA:972:C:OP2	43:DJ:57:LYS:NZ	2.24	0.60
57:DZ:32:ILE:O	57:DZ:36:THR:OG1	2.19	0.60
1:AA:1848:G:OP1	4:AD:88:ARG:NH2	2.33	0.60
1:AA:2328:C:H2'	1:AA:2329:C:C6	2.34	0.60
1:AA:2349:G:OP1	63:AA:4052:HOH:O	2.16	0.60
1:AA:2859:U:OP2	17:AT:95:ARG:NH1	2.34	0.60
16:AS:10:ARG:O	16:AS:14:VAL:HG13	2.01	0.60
18:AU:25:TRP:O	18:AU:28:ARG:HB2	2.01	0.60
34:BA:1221:G:H4'	52:BS:77:THR:HG21	1.83	0.60
57:BZ:99:ARG:NH1	57:BZ:312:LEU:HD11	2.16	0.60
1:CA:1693:U:O2'	4:CD:14:ARG:NH2	2.33	0.60
5:CE:35:GLN:OE1	5:CE:66:HIS:HE1	1.84	0.60
5:CE:68:ALA:O	5:CE:70:ALA:N	2.34	0.60
8:CH:144:VAL:O	8:CH:148:ILE:HG12	2.01	0.60
11:CN:21:LYS:NZ	11:CN:140:VAL:OXT	2.31	0.60
36:DC:157:ILE:HD12	36:DC:164:ARG:HB3	1.82	0.60
10:AL:30:HIS:HD2	10:AL:65:PHE:HB2	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BE:152:ARG:HA	41:BH:64:LYS:NZ	2.17	0.60
1:CA:1364:G:N7	25:C1:3:LYS:HE2	2.16	0.60
1:CA:1912:A:O3'	57:DZ:499:ARG:NH2	2.34	0.60
1:CA:2113:U:H2'	1:CA:2114:A:H8	1.64	0.60
1:CA:2298:A:H2'	1:CA:2299:G:O4'	2.01	0.60
1:CA:77:C:OP1	26:C2:59:ARG:HD3	2.00	0.60
1:CA:2177:C:C2'	3:CC:171:ALA:HB2	2.31	0.60
34:DA:620:C:C2	37:DD:135:LEU:HG	2.36	0.60
35:DB:219:VAL:HA	35:DB:222:ILE:HG12	1.83	0.60
57:DZ:637:ARG:C	57:DZ:639:ASN:H	2.05	0.60
7:AG:43:LEU:HD11	7:AG:153:ARG:HG2	1.83	0.60
7:AG:77:ILE:HG22	7:AG:80:PHE:H	1.65	0.60
34:BA:509:A:H3'	34:BA:509:A:C8	2.36	0.60
1:CA:2162:G:OP1	1:CA:2172:U:O2'	2.17	0.60
1:CA:2364:C:H2'	1:CA:2365:G:O4'	2.01	0.60
23:CZ:132:ASN:O	23:CZ:134:PRO:HD3	2.02	0.60
34:DA:343:U:O2'	34:DA:344:A:H2'	2.02	0.60
34:DA:982:U:H5"	47:DN:6:LEU:HD21	1.81	0.60
11:AN:75:TYR:CZ	11:AN:77:GLY:HA2	2.37	0.60
16:AS:27:SER:HA	16:AS:88:ASP:HB3	1.84	0.60
41:BH:51:VAL:HG21	41:BH:60:ARG:HD2	1.83	0.60
1:CA:1709:U:H2'	1:CA:1710:C:C6	2.37	0.60
37:DD:36:ARG:HG3	37:DD:38:TYR:CE2	2.37	0.60
61:DZ:703:FUA:H12	61:DZ:703:FUA:O1	2.00	0.60
1:AA:2402:U:P	32:A8:35:GLN:HE22	2.25	0.60
1:AA:2331:G:N1	16:AS:3:ARG:HA	2.16	0.60
34:BA:167:G:H2'	34:BA:168:G:C8	2.36	0.60
34:BA:1372:U:H5"	42:BI:71:SER:HB3	1.81	0.60
53:BT:77:ALA:O	53:BT:81:LYS:HG3	2.00	0.60
57:BZ:350:GLU:OE1	57:BZ:381:LYS:N	2.17	0.60
1:CA:1683:C:H2'	1:CA:1684:C:C6	2.36	0.60
16:CS:14:VAL:O	16:CS:18:ILE:HG12	2.01	0.60
23:CZ:69:THR:HG22	23:CZ:90:VAL:HG22	1.82	0.60
56:DW:40:C:O2'	56:DY:36:A:OP1	2.19	0.60
1:AA:1405:A:N1	1:AA:1418:U:C4	2.69	0.60
1:AA:469:A:H1'	1:AA:1246:C:O4'	2.02	0.60
46:BM:16:ASP:N	46:BM:16:ASP:OD1	2.31	0.60
24:C0:10:THR:HG22	24:C0:12:ASN:H	1.66	0.60
1:CA:889:C:O2'	1:CA:890:A:O4'	2.19	0.60
3:CC:41:THR:HG22	3:CC:42:VAL:N	2.17	0.60
24:A0:11:ARG:O	24:A0:14:ARG:NH2	2.26	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:214:TYR:CE2	3:AC:224:ARG:HG2	2.37	0.60
34:BA:959:A:HO2'	34:BA:984:C:HO2'	1.36	0.60
42:BI:23:ASN:HD22	42:BI:25:LYS:HG2	1.66	0.60
1:CA:1420:U:O2'	1:CA:1421:G:OP1	2.19	0.60
1:CA:2317:C:N4	1:CA:2318:G:O6	2.35	0.60
1:CA:748:G:O6	20:CW:90:ARG:NH1	2.35	0.60
5:CE:16:ARG:NH1	5:CE:171:GLU:OE2	2.31	0.60
6:CF:129:PHE:O	6:CF:132:VAL:HG22	2.02	0.60
34:DA:1347:G:C8	42:DI:107:ARG:HB3	2.36	0.60
35:DB:166:ASP:OD2	35:DB:169:LYS:N	2.34	0.60
36:DC:65:ALA:HA	36:DC:100:ALA:HB3	1.83	0.60
1:AA:2157:A:N6	1:AA:2178:G:O2'	2.23	0.60
1:AA:2511:C:N3	63:AA:4228:HOH:O	2.31	0.60
34:BA:1030(C):G:H2'	34:BA:1030(D):A:H8	1.65	0.60
25:C1:50:ARG:HG2	25:C1:59:THR:HB	1.83	0.60
53:DT:18:GLN:O	53:DT:22:ARG:HG3	2.01	0.60
18:AU:76:TYR:CZ	18:AU:80:ILE:HG13	2.37	0.60
45:BL:53:ARG:HG3	45:BL:93:LEU:HD21	1.84	0.60
3:CC:194:ILE:HD11	3:CC:227:PRO:CB	2.32	0.60
1:CA:2307:G:N1	7:CG:43:LEU:O	2.34	0.60
34:DA:187:C:OP1	53:DT:82:SER:OG	2.18	0.60
28:A4:58:ARG:O	28:A4:60:GLN:N	2.35	0.59
1:AA:818:G:OP1	31:A7:10:ARG:NH1	2.35	0.59
1:AA:1065:U:O2'	1:AA:1067:A:H2	1.83	0.59
1:AA:2473:C:H2'	1:AA:2474:U:C6	2.37	0.59
3:AC:194:ILE:HD11	3:AC:227:PRO:CB	2.32	0.59
34:BA:321:A:N7	34:BA:328:C:O2'	2.29	0.59
34:BA:828:A:H2'	34:BA:829:G:O4'	2.02	0.59
34:BA:508:C:OP1	37:BD:209:ARG:NH2	2.34	0.59
34:BA:279:A:N6	50:BQ:98:LEU:O	2.35	0.59
57:BZ:74:TRP:CD1	57:BZ:273:LEU:HB3	2.37	0.59
57:BZ:639:ASN:HA	57:BZ:640:ALA:O	2.02	0.59
1:CA:774:A:HO2'	1:CA:775:G:H8	1.49	0.59
38:DE:11:ILE:HG22	38:DE:31:LEU:HB3	1.84	0.59
49:DP:55:ARG:O	49:DP:58:TYR:N	2.35	0.59
57:DZ:510:VAL:HG21	57:DZ:542:VAL:HG21	1.84	0.59
57:DZ:606:MET:HG3	57:DZ:649:LEU:HD23	1.84	0.59
24:A0:6:GLY:O	56:BW:1:G:O2'	2.19	0.59
1:AA:1219:A:H4'	1:AA:1220:U:OP1	2.01	0.59
1:AA:2230:U:O4'	25:A1:52:ARG:NH2	2.34	0.59
3:AC:11:LEU:HD12	3:AC:33:LEU:HA	1.82	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1359:A:H2'	1:CA:1360:A:H5'	1.83	0.59
6:CF:36:VAL:HG11	6:CF:183:VAL:HG13	1.83	0.59
43:DJ:52:GLY:O	47:DN:41:ARG:NH2	2.33	0.59
57:DZ:630:GLN:O	57:DZ:646:PHE:N	2.34	0.59
1:AA:483:A:H5''	63:AA:5245:HOH:O	2.03	0.59
5:AE:120:TRP:CD2	5:AE:155:LYS:HG2	2.37	0.59
6:AF:191:ARG:HH11	6:AF:191:ARG:CG	2.14	0.59
7:AG:130:ASN:HB3	7:AG:160:VAL:HA	1.84	0.59
10:AL:99:ILE:HG23	10:AL:103:GLN:HB3	1.84	0.59
13:AP:101:VAL:HG23	13:AP:106:LEU:HB3	1.84	0.59
34:BA:671:G:H5'	39:BF:77:ARG:HH22	1.66	0.59
56:BY:6:G:O6	56:BY:7:A:N6	2.36	0.59
57:BZ:512:ILE:HB	57:BZ:566:THR:O	2.02	0.59
4:CD:146:GLU:HB2	4:CD:189:CYS:HB3	1.84	0.59
10:CL:125:ARG:HA	10:CL:128:ALA:HB3	1.84	0.59
17:CT:50:ILE:HA	17:CT:99:LEU:HB2	1.84	0.59
34:DA:493:G:N2	34:DA:494:U:O4	2.36	0.59
38:DE:88:LYS:HB3	38:DE:123:LEU:HB2	1.84	0.59
49:DP:28:ARG:NH1	49:DP:29:ASP:OD1	2.36	0.59
1:AA:1410:G:C8	25:A1:3:LYS:HE2	2.37	0.59
1:AA:794:U:O2	1:AA:2036:A:H1'	2.02	0.59
1:AA:2658:C:O5'	1:AA:2658:C:H6	1.85	0.59
10:AL:51:ALA:HB1	10:AL:72:PRO:HB3	1.85	0.59
34:BA:539:A:OP2	45:BL:115:LYS:NZ	2.35	0.59
57:BZ:210:ARG:CG	57:BZ:210:ARG:HH11	2.16	0.59
13:CP:63:PRO:HG2	32:C8:25:MET:HB2	1.83	0.59
1:CA:1859:A:N6	1:CA:1883:G:O2'	2.35	0.59
1:CA:2128:C:N3	1:CA:2160:G:N2	2.40	0.59
1:CA:81:G:O6	63:CA:4119:HOH:O	2.17	0.59
6:CF:129:PHE:CD2	6:CF:163:VAL:HG21	2.37	0.59
1:CA:528:A:OP2	11:CN:114:ARG:NH1	2.35	0.59
1:CA:72:U:OP2	21:CX:1:MET:N	2.35	0.59
34:DA:1133:G:H2'	34:DA:1134:G:H8	1.65	0.59
37:DD:26:CYS:HA	60:DD:501:SF4:S3	2.43	0.59
34:DA:737:A:H1'	39:DF:73:ASN:OD1	2.03	0.59
42:DI:9:ARG:HG2	42:DI:14:VAL:HG12	1.84	0.59
49:DP:5:ARG:HH12	49:DP:28:ARG:HA	1.67	0.59
53:DT:33:ILE:O	53:DT:37:SER:OG	2.11	0.59
1:AA:1935:A:H8	1:AA:1935:A:H5'	1.67	0.59
16:AS:14:VAL:O	16:AS:18:ILE:HG12	2.02	0.59
34:BA:435:C:H2'	34:BA:436:C:C6	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BB:204:ASN:OD1	35:BB:205:ASP:N	2.35	0.59
41:BH:4:ASP:OD2	41:BH:85:ARG:NH1	2.33	0.59
34:BA:1525:G:P	44:BK:120:ARG:HH22	2.25	0.59
44:BK:84:VAL:HG21	44:BK:95:ILE:HD11	1.85	0.59
57:BZ:338:GLY:O	57:BZ:351:ARG:NH2	2.34	0.59
57:BZ:402:ILE:HG23	57:BZ:403:GLU:H	1.67	0.59
57:BZ:421:GLN:O	57:BZ:424:LEU:HB3	2.01	0.59
28:C4:41:PRO:HA	28:C4:44:THR:HG22	1.84	0.59
1:CA:1292:U:H2'	1:CA:1293:C:C6	2.38	0.59
1:CA:1466:G:HO2'	1:CA:1546:C:HO2'	1.50	0.59
7:CG:15:VAL:HA	7:CG:175:LEU:HD23	1.83	0.59
10:CL:100:THR:HA	10:CL:139:VAL:HB	1.84	0.59
34:DA:1457:G:H5''	53:DT:35:THR:HG21	1.84	0.59
1:AA:1793:A:H2'	63:AA:5101:HOH:O	2.02	0.59
1:AA:2298:A:H4'	1:AA:2299:A:O4'	2.01	0.59
7:AG:146:TYR:O	7:AG:149:VAL:HG12	2.03	0.59
20:AW:62:HIS:O	20:AW:64:MET:HG3	2.03	0.59
34:BA:1356:G:H2'	34:BA:1357:A:H8	1.63	0.59
37:BD:18:LYS:HG2	37:BD:33:MET:HG2	1.84	0.59
26:C2:19:VAL:HA	26:C2:22:GLU:HG3	1.85	0.59
34:DA:1342:C:O2'	42:DI:124:GLN:HG2	2.02	0.59
1:AA:2331:G:N2	16:AS:3:ARG:HG2	2.09	0.59
18:AU:28:ARG:HG2	18:AU:38:THR:OG1	2.02	0.59
36:BC:40:ARG:HG2	36:BC:55:VAL:HG11	1.85	0.59
14:CQ:27:VAL:O	14:CQ:29:PHE:N	2.35	0.59
16:CS:50:SER:OG	16:CS:50:SER:O	2.17	0.59
1:CA:25:U:H5''	20:CW:80:PRO:HD3	1.84	0.59
35:DB:127:ILE:HG12	35:DB:128:GLU:H	1.65	0.59
42:DI:16:ARG:HB2	42:DI:64:THR:HG23	1.84	0.59
33:A9:15:LYS:HE2	33:A9:17:ILE:HD11	1.85	0.59
1:AA:2362:C:OP2	63:AA:3980:HOH:O	2.17	0.59
1:AA:41:C:H42	1:AA:464:G:H1	1.49	0.59
3:AC:214:TYR:CZ	3:AC:224:ARG:HG2	2.37	0.59
34:BA:222:U:H2'	34:BA:223:U:C6	2.37	0.59
34:BA:43:C:H42	34:BA:399:G:H1	1.50	0.59
34:BA:688:G:O2'	34:BA:704:A:N1	2.27	0.59
43:BJ:45:ARG:HH11	47:BN:36:PHE:HE1	1.49	0.59
49:BP:69:THR:HA	49:BP:72:ARG:HB2	1.84	0.59
49:BP:74:LEU:O	49:BP:79:VAL:HG23	2.03	0.59
53:BT:56:MET:HE2	53:BT:85:MET:HG2	1.84	0.59
57:BZ:601:ILE:O	57:BZ:679:VAL:HG13	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2121:G:O4'	3:CC:168:LYS:CD	2.47	0.59
23:CZ:156:LYS:HD2	23:CZ:158:PRO:HD3	1.85	0.59
34:DA:1363(A):A:H1'	34:DA:1365:G:N7	2.17	0.59
34:DA:390:C:O3'	49:DP:28:ARG:NH2	2.36	0.59
56:DY:33:U:H2'	56:DY:35:A:OP2	2.02	0.59
6:AF:178:PRO:HB3	6:AF:198:ALA:HB1	1.83	0.59
7:AG:129:GLY:O	7:AG:161:THR:HG22	2.03	0.59
34:BA:147:G:N2	34:BA:175:C:O2	2.31	0.59
37:BD:105:VAL:HG13	37:BD:110:PHE:HB2	1.85	0.59
34:BA:412:A:H8	37:BD:35:ARG:HH21	1.50	0.59
1:CA:1044:G:H21	1:CA:1111:A:H2	1.51	0.59
1:CA:1055:G:H3'	1:CA:1056:G:H8	1.66	0.59
1:CA:1429:G:H2'	1:CA:1430:C:C6	2.37	0.59
3:CC:214:TYR:CZ	3:CC:224:ARG:HG2	2.37	0.59
9:CK:27:VAL:HA	9:CK:113:GLN:HA	1.85	0.59
17:CT:6:LEU:O	17:CT:10:VAL:HG23	2.03	0.59
12:CO:104:ARG:HH22	17:CT:43:GLN:NE2	2.00	0.59
34:DA:624:C:H2'	34:DA:625:G:C8	2.38	0.59
34:DA:981:U:H5'	47:DN:21:TYR:CE2	2.38	0.59
34:BA:447:G:H2'	34:BA:485:G:N2	2.18	0.59
56:BY:67:C:H2'	56:BY:68:C:C6	2.38	0.59
17:CT:16:ARG:NH1	17:CT:19:LEU:HD21	2.18	0.59
34:DA:1143:G:H2'	34:DA:1144:G:C8	2.38	0.59
43:DJ:6:ILE:HB	43:DJ:72:VAL:HG23	1.84	0.59
57:DZ:413:ILE:HB	57:DZ:476:VAL:HG12	1.84	0.59
4:AD:148:GLU:O	4:AD:151:LYS:HB2	2.03	0.58
29:C5:20:ARG:O	29:C5:23:HIS:HB2	2.02	0.58
1:CA:1359:A:H2	1:CA:1372:U:O4	1.86	0.58
15:CR:33:ARG:HE	15:CR:113:LEU:HD22	1.67	0.58
34:DA:1348:U:H4'	42:DI:120:ARG:HD3	1.85	0.58
34:DA:978:A:OP2	34:DA:1363:C:N4	2.36	0.58
34:DA:35:G:O2'	45:DL:118:SER:O	2.13	0.58
57:DZ:114:VAL:HG21	57:DZ:156:ARG:HB2	1.85	0.58
57:DZ:264:LEU:HB2	62:DZ:704:GDP:C5	2.38	0.58
1:AA:2250:G:N3	1:AA:2250:G:H2'	2.18	0.58
1:AA:595:A:H5''	1:AA:596:G:OP2	2.02	0.58
1:AA:692:C:H2'	1:AA:693:G:C8	2.38	0.58
14:AQ:38:GLU:HA	14:AQ:99:PRO:HG3	1.85	0.58
34:BA:186:C:H2'	34:BA:187:C:H6	1.68	0.58
35:BB:170:GLU:HB3	35:BB:173:ALA:HB3	1.83	0.58
35:BB:158:LEU:HG	35:BB:182:ILE:HD11	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BC:11:ARG:HH21	36:BC:180:ALA:HB3	1.67	0.58
36:BC:6:HIS:HB2	47:BN:49:HIS:HD2	1.67	0.58
38:BE:152:ARG:HA	41:BH:64:LYS:HZ2	1.68	0.58
44:BK:44:SER:OG	44:BK:47:VAL:HG23	2.03	0.58
1:CA:946:G:OP1	63:CA:4242:HOH:O	2.16	0.58
34:DA:537:G:H2'	34:DA:538:G:C8	2.38	0.58
34:DA:426:G:OP1	37:DD:38:TYR:OH	2.21	0.58
41:DH:44:PHE:HB3	41:DH:80:ILE:HD11	1.85	0.58
27:A3:3:ARG:HH11	27:A3:60:GLU:CD	2.07	0.58
6:AF:161:GLU:OE1	6:AF:165:ARG:NH1	2.36	0.58
6:AF:53:THR:HG22	6:AF:55:GLY:N	2.17	0.58
23:AZ:183:LEU:O	23:AZ:185:GLU:N	2.37	0.58
34:BA:1095:U:OP1	34:BA:1108:G:N1	2.33	0.58
34:BA:1305:G:N2	34:BA:1331:G:H1'	2.15	0.58
1:CA:1006:C:O2'	11:CN:106:MET:HB3	2.03	0.58
1:CA:284:U:H2'	1:CA:285:C:C6	2.38	0.58
1:CA:2861:G:H2'	1:CA:2862:G:H8	1.67	0.58
1:CA:829:A:N7	1:CA:2247:A:O2'	2.37	0.58
2:CB:66:A:N6	2:CB:109:C:H5''	2.16	0.58
3:CC:214:TYR:CE2	3:CC:224:ARG:HG2	2.37	0.58
56:DW:36:A:N6	56:DW:37:MIA:H152	2.18	0.58
56:DY:5:G:H1	56:DY:68:C:H42	1.48	0.58
57:DZ:225:GLU:HA	57:DZ:228:MET:HB3	1.84	0.58
1:AA:1338:U:H2'	1:AA:1339:C:C6	2.38	0.58
13:AP:52:GLU:HG3	13:AP:57:THR:HG22	1.86	0.58
1:AA:1700:G:H3'	15:AR:2:ARG:HD3	1.84	0.58
34:BA:1226:C:P	46:BM:91:ARG:HH22	2.26	0.58
34:BA:1239:A:H4'	34:BA:1240:U:H5''	1.86	0.58
34:BA:102:G:O2'	34:BA:151:A:N3	2.32	0.58
34:BA:358:U:OP1	57:BZ:381:LYS:NZ	2.37	0.58
43:BJ:38:ILE:HD11	43:BJ:71:LEU:HD23	1.85	0.58
57:BZ:-4:ALA:HA	57:BZ:-1:GLU:HB3	1.85	0.58
57:BZ:637:ARG:C	57:BZ:639:ASN:H	2.06	0.58
2:CB:6:C:H2'	2:CB:7:G:O4'	2.04	0.58
34:DA:976:G:H5'	34:DA:1358:U:O2'	2.03	0.58
57:DZ:35:TYR:CE2	57:DZ:269:VAL:HB	2.38	0.58
1:AA:116:A:H5'	1:AA:117:A:H8	1.68	0.58
1:AA:992:G:OP2	63:AA:4762:HOH:O	2.17	0.58
3:AC:41:THR:HG22	3:AC:42:VAL:N	2.17	0.58
34:BA:243:A:H4'	34:BA:244:U:O5'	2.02	0.58
34:BA:942:G:H21	42:BI:124:GLN:NE2	2.00	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:BC:11:ARG:NH2	36:BC:177:THR:O	2.36	0.58
38:BE:57:LYS:HG2	38:BE:61:TYR:HE2	1.67	0.58
33:C9:18:ARG:NH1	33:C9:21:GLY:HA2	2.18	0.58
1:CA:787:U:OP2	63:CA:3874:HOH:O	2.17	0.58
1:CA:2121:G:C4'	3:CC:168:LYS:HD3	2.34	0.58
11:CN:58:ASP:N	11:CN:58:ASP:OD1	2.35	0.58
34:DA:670:G:H21	39:DF:73:ASN:HD21	1.51	0.58
27:A3:3:ARG:NH1	27:A3:60:GLU:OE1	2.37	0.58
34:BA:1125:U:H4'	43:BJ:5:ARG:NH2	2.17	0.58
34:BA:1376:U:OP1	40:BG:98:SER:OG	2.21	0.58
14:CQ:52:VAL:HG22	23:CZ:183:LEU:HD11	1.86	0.58
34:DA:1189:C:OP1	43:DJ:51:ARG:NH2	2.36	0.58
34:DA:1346:A:OP1	42:DI:120:ARG:NH2	2.28	0.58
34:DA:60:A:N1	34:DA:107:G:O2'	2.26	0.58
34:DA:1190:G:OP1	36:DC:5:ILE:N	2.36	0.58
57:DZ:99:ARG:HD2	57:DZ:289:ILE:HD11	1.86	0.58
1:AA:1452:U:H2'	1:AA:1453:C:C6	2.39	0.58
1:AA:1891:G:H4'	3:AC:206:LYS:HG3	1.80	0.58
34:BA:1030(C):G:N7	34:BA:1031:G:N2	2.52	0.58
41:BH:64:LYS:HG2	41:BH:79:VAL:HG21	1.85	0.58
1:CA:1840:G:OP2	63:CA:4373:HOH:O	2.17	0.58
17:CT:23:ARG:HG3	17:CT:120:ARG:NH1	2.19	0.58
34:DA:678:U:H2'	34:DA:679:C:C6	2.37	0.58
28:A4:3:GLU:O	28:A4:5:ILE:N	2.36	0.58
1:AA:409:G:OP2	63:AA:4963:HOH:O	2.16	0.58
1:AA:662:A:H8	13:AP:117:GLU:HG3	1.68	0.58
10:AL:53:VAL:HG12	10:AL:69:THR:HB	1.85	0.58
12:AO:18:LYS:HB2	12:AO:45:GLU:HB3	1.85	0.58
34:BA:452:A:H4'	49:BP:72:ARG:NH1	2.19	0.58
35:BB:215:LEU:O	35:BB:219:VAL:HG23	2.04	0.58
34:DA:1071:C:H2'	34:DA:1072:G:H8	1.69	0.58
34:DA:1412:C:H2'	34:DA:1413:A:C8	2.38	0.58
34:DA:456:C:H42	34:DA:475:G:H1	1.50	0.58
42:DI:13:ALA:HB2	42:DI:68:GLY:HA3	1.86	0.58
49:DP:5:ARG:HB3	49:DP:67:THR:HG23	1.86	0.58
57:DZ:15:ILE:HA	57:DZ:103:GLY:O	2.04	0.58
57:DZ:175:SER:O	57:DZ:188:TYR:N	2.34	0.58
57:DZ:316:ILE:HG12	57:DZ:385:THR:HG22	1.86	0.58
28:A4:16:CYS:SG	28:A4:17:GLY:N	2.75	0.58
10:AL:50:ASP:OD1	10:AL:50:ASP:N	2.37	0.58
35:BB:15:VAL:HB	35:BB:209:ARG:HD3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:492:A:H2'	1:CA:493:G:O4'	2.03	0.58
12:CO:35:VAL:HG13	12:CO:65:THR:HG23	1.85	0.58
34:DA:337:C:H2'	34:DA:338:A:H8	1.67	0.58
34:DA:433:C:H2'	34:DA:434:U:H6	1.67	0.58
1:AA:1273:G:OP1	18:AU:13:LYS:HE3	2.03	0.58
1:AA:2856:G:H2'	1:AA:2857:U:O4'	2.04	0.58
6:AF:34:TRP:CE2	13:AP:8:PRO:HD3	2.38	0.58
17:AT:16:ARG:NH2	17:AT:83:ILE:O	2.33	0.58
34:BA:42:G:OP1	63:BA:5286:HOH:O	2.16	0.58
57:BZ:71:THR:HG22	57:BZ:80:ASN:OD1	2.04	0.58
1:CA:2864:G:O2'	1:CA:2865:U:H5'	2.04	0.58
1:CA:307:G:H21	1:CA:330:A:H62	1.52	0.58
43:DJ:34:VAL:HG12	43:DJ:74:ILE:HG12	1.86	0.58
49:DP:74:LEU:O	49:DP:79:VAL:HG23	2.04	0.58
1:AA:1652:G:H5''	1:AA:1653:C:OP1	2.03	0.57
1:AA:2274:U:OP2	24:A0:19:LYS:NZ	2.37	0.57
34:BA:434:U:H2'	34:BA:435:C:O4'	2.04	0.57
34:BA:687:A:N3	34:BA:688:G:H1'	2.19	0.57
34:BA:955:U:H2'	34:BA:956:U:H6	1.69	0.57
35:BB:18:GLY:O	35:BB:19:HIS:HB3	2.03	0.57
49:BP:49:LEU:HD12	49:BP:50:LYS:N	2.19	0.57
57:BZ:318:ALA:O	57:BZ:319:ASP:HB2	2.04	0.57
1:CA:1240:U:O4	63:CA:4307:HOH:O	2.13	0.57
1:CA:1902:C:H5'	4:CD:246:PRO:HD3	1.85	0.57
7:CG:107:LEU:HD21	7:CG:178:PHE:CE2	2.39	0.57
10:CL:101:TRP:HE1	10:CL:140:GLY:HA3	1.69	0.57
42:DI:112:LYS:NZ	42:DI:113:LYS:O	2.35	0.57
57:DZ:519:ARG:NH1	57:DZ:678:GLU:H	2.02	0.57
1:AA:2303:U:H2'	1:AA:2304:C:C6	2.38	0.57
34:BA:222:U:H2'	34:BA:223:U:H6	1.68	0.57
34:BA:661:G:H1	34:BA:744:C:H42	1.52	0.57
34:BA:814:A:H2'	34:BA:816:A:H5''	1.86	0.57
37:BD:117:ALA:O	37:BD:121:VAL:HG23	2.04	0.57
1:CA:1263:U:H1'	29:C5:10:LYS:HG3	1.86	0.57
32:C8:6:THR:HG22	32:C8:64:TYR:HD2	1.68	0.57
1:CA:2387:U:OP1	24:C0:55:ARG:NH1	2.32	0.57
34:DA:1224:G:OP1	63:DA:3269:HOH:O	2.17	0.57
36:DC:78:GLY:HA3	36:DC:83:ARG:H	1.68	0.57
1:AA:1055:A:OP2	63:AA:4596:HOH:O	2.17	0.57
1:AA:1285:G:H2'	1:AA:1286:U:O4'	2.04	0.57
8:AH:7:LEU:O	8:AH:69:ARG:HD3	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:BB:100:GLY:N	35:BB:176:GLU:OE2	2.34	0.57
35:BB:178:ARG:HH22	41:BH:68:ARG:NH1	2.02	0.57
34:BA:1014:A:H4'	52:BS:14:HIS:CE1	2.39	0.57
1:CA:2345:G:N3	1:CA:2381:C:H2'	2.19	0.57
1:CA:530:G:N1	63:CA:4131:HOH:O	2.22	0.57
34:DA:1289:A:H3'	34:DA:1290:G:H8	1.69	0.57
34:DA:765:G:H5''	34:DA:766:A:OP1	2.04	0.57
36:DC:44:GLU:HG3	36:DC:52:LEU:HD11	1.86	0.57
46:DM:65:LYS:HB2	46:DM:69:GLU:HB2	1.85	0.57
43:DJ:63:PHE:HE2	47:DN:45:ARG:HA	1.69	0.57
51:DR:25:THR:O	51:DR:25:THR:OG1	2.23	0.57
54:DU:9:ARG:O	54:DU:13:ILE:HG13	2.04	0.57
57:DZ:402:ILE:HG23	57:DZ:403:GLU:H	1.69	0.57
57:DZ:494:GLU:HB2	57:DZ:511:LYS:HE2	1.86	0.57
32:A8:39:LYS:O	32:A8:43:GLN:HB2	2.05	0.57
34:BA:1070:U:H2'	34:BA:1071:C:H6	1.69	0.57
34:BA:458:C:N4	34:BA:474:G:O6	2.37	0.57
34:BA:920:U:H2'	34:BA:921:U:C6	2.39	0.57
50:BQ:45:HIS:HB3	50:BQ:72:ARG:HB3	1.86	0.57
57:BZ:227:ILE:HA	57:BZ:230:LYS:HG3	1.86	0.57
1:CA:2751:G:N2	8:CH:2:SER:OG	2.38	0.57
12:CO:48:PRO:HB3	34:DA:1422:G:H5'	1.86	0.57
1:CA:1339:G:H5''	21:CX:16:LYS:HD2	1.85	0.57
1:AA:2144:U:O2'	3:AC:167:ASP:HB3	2.03	0.57
22:AY:54:LYS:HA	22:AY:56:PRO:HD3	1.86	0.57
34:BA:375:U:C2	34:BA:376:G:C8	2.93	0.57
35:BB:219:VAL:HA	35:BB:222:ILE:HD12	1.86	0.57
53:BT:56:MET:HE2	53:BT:85:MET:HA	1.86	0.57
1:CA:2661:G:O6	8:CH:175:LYS:NZ	2.37	0.57
1:CA:674:G:H1'	6:CF:74:ARG:HD3	1.86	0.57
17:CT:82:LEU:HD12	17:CT:82:LEU:H	1.68	0.57
1:AA:2289:G:P	24:A0:10:THR:HG21	2.45	0.57
1:AA:1451:U:H2'	1:AA:1452:U:C6	2.40	0.57
1:AA:2054:G:O2'	5:AE:145:LYS:HE3	2.05	0.57
1:AA:2245:U:H2'	1:AA:2246:G:C8	2.40	0.57
42:BI:21:PRO:HA	42:BI:59:PHE:HA	1.87	0.57
43:BJ:8:LEU:HB2	43:BJ:70:ARG:HB2	1.85	0.57
1:CA:2539:C:H4'	33:C9:3:VAL:HG21	1.85	0.57
1:CA:2320:A:H2'	1:CA:2320:A:N3	2.18	0.57
1:CA:526:A:N1	1:CA:2625:G:O2'	2.29	0.57
7:CG:120:LEU:HB3	7:CG:131:TYR:OH	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:798:G:OP1	44:DK:122:LYS:NZ	2.37	0.57
57:DZ:169:GLY:N	57:DZ:170:ARG:NH1	2.48	0.57
1:AA:2357:G:OP2	30:A6:38:LYS:HD3	2.04	0.57
1:AA:801:C:H2'	1:AA:802:C:C6	2.39	0.57
5:AE:4:ILE:HD11	5:AE:29:GLY:HA2	1.86	0.57
12:AO:16:ALA:HB2	12:AO:52:VAL:CG2	2.35	0.57
16:AS:39:ILE:HB	16:AS:49:VAL:HG12	1.86	0.57
34:BA:413:G:H1'	34:BA:428:G:H21	1.69	0.57
34:BA:67:C:H2'	34:BA:68:G:H8	1.69	0.57
43:BJ:38:ILE:HG13	43:BJ:71:LEU:HB3	1.86	0.57
1:CA:1154:G:O5'	1:CA:1154:G:H8	1.87	0.57
1:CA:2177:C:O2	3:CC:171:ALA:HB2	2.04	0.57
3:CC:6:LYS:HG3	3:CC:7:ARG:H	1.69	0.57
5:CE:181:LEU:HD21	17:CT:6:LEU:HD12	1.86	0.57
34:DA:1151:A:O2'	34:DA:1152:A:O5'	2.21	0.57
57:DZ:363:ARG:HG2	57:DZ:363:ARG:HH11	1.68	0.57
1:AA:2559:U:H2'	1:AA:2560:G:C8	2.40	0.57
6:AF:28:ILE:HD13	6:AF:119:ARG:HH21	1.69	0.57
7:AG:11:TYR:HA	7:AG:15:VAL:HB	1.86	0.57
42:BI:93:ARG:HB2	42:BI:93:ARG:HH11	1.70	0.57
57:BZ:526:VAL:HG23	57:BZ:566:THR:HA	1.87	0.57
26:C2:64:LEU:HD11	26:C2:68:ARG:HH21	1.69	0.57
1:CA:1386:C:H2'	1:CA:1387:C:C6	2.39	0.57
1:CA:1530:C:H42	1:CA:1539:G:H1	1.51	0.57
1:CA:2267:A:H5'	1:CA:2268:A:H5'	1.86	0.57
1:CA:2365:G:O6	32:C8:39:LYS:HE3	2.04	0.57
1:CA:731:C:OP1	63:CA:4294:HOH:O	2.18	0.57
34:DA:1288:A:N1	34:DA:1371:G:H1'	2.20	0.57
34:DA:1356:G:H2'	34:DA:1357:A:C8	2.39	0.57
34:DA:17:U:H2'	34:DA:18:C:C6	2.40	0.57
37:DD:103:ASN:OD1	37:DD:114:ARG:NE	2.36	0.57
39:DF:82:ARG:HB2	39:DF:85:VAL:HG23	1.86	0.57
48:DO:26:GLU:OE2	48:DO:77:ARG:NE	2.20	0.57
57:DZ:177:ILE:HG13	57:DZ:188:TYR:CE2	2.39	0.57
1:AA:1529:G:O6	1:AA:1553:A:N6	2.38	0.57
1:AA:927:G:OP2	1:AA:927:G:H8	1.88	0.57
34:BA:189(C):C:H2'	34:BA:189(D):C:O4'	2.05	0.57
47:BN:4:LYS:HA	47:BN:7:ILE:HG23	1.87	0.57
49:BP:49:LEU:HD11	49:BP:51:VAL:HG23	1.86	0.57
52:BS:3:ARG:NH1	52:BS:10:PHE:HB2	2.20	0.57
4:CD:13:ARG:HD2	4:CD:16:MET:HE3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:CL:8:VAL:HG21	10:CL:26:ALA:HB1	1.87	0.57
34:DA:157:G:H2'	34:DA:158:G:H8	1.70	0.57
34:DA:646:U:H2'	34:DA:647:C:C6	2.39	0.57
42:DI:77:ILE:O	42:DI:81:ILE:HG22	2.05	0.57
43:DJ:43:ARG:HB2	43:DJ:67:THR:HG23	1.85	0.57
57:DZ:210:ARG:HH11	57:DZ:210:ARG:HG3	1.70	0.57
57:DZ:326:THR:HB	57:DZ:377:VAL:HG13	1.86	0.57
1:AA:847:A:OP1	1:AA:847:A:H8	1.88	0.57
1:AA:630:U:OP1	6:AF:102:PRO:HA	2.05	0.57
7:AG:41:GLN:CG	7:AG:60:LEU:HD21	2.35	0.57
23:AZ:19:ARG:NH1	23:AZ:84:GLU:O	2.37	0.57
34:BA:408:A:OP1	37:BD:113:SER:OG	2.18	0.57
37:BD:110:PHE:CE2	37:BD:148:VAL:HG23	2.40	0.57
38:BE:36:ASP:OD1	38:BE:39:GLY:N	2.34	0.57
41:BH:77:GLU:HG3	41:BH:78:GLN:N	2.19	0.57
56:BW:44:G:O2'	56:BW:45:U:H5'	2.05	0.57
57:BZ:-10:ARG:HB2	57:BZ:-10:ARG:HH11	1.70	0.57
1:CA:184:C:H2'	1:CA:185:U:C6	2.39	0.57
1:CA:1956:U:H2'	1:CA:1957:C:H5'	1.86	0.57
1:CA:2291:U:H2'	1:CA:2292:C:C6	2.40	0.57
15:CR:18:LEU:HD22	15:CR:22:ARG:HD2	1.87	0.57
34:DA:1022:G:H2'	34:DA:1023:G:H8	1.69	0.57
34:DA:1305:G:O2'	34:DA:1331:G:N2	2.37	0.57
41:DH:8:ASP:OD2	41:DH:12:ARG:NH1	2.38	0.57
45:DL:76:ASN:ND2	45:DL:106:ASP:O	2.37	0.57
52:DS:33:THR:OG1	52:DS:35:SER:O	2.23	0.57
14:AQ:56:ARG:HG3	14:AQ:56:ARG:NH1	2.19	0.56
34:BA:1530:G:H4'	34:BA:1530:G:OP1	2.05	0.56
34:BA:633:G:H2'	34:BA:634:C:C6	2.40	0.56
36:BC:53:ALA:HB2	36:BC:115:LEU:HD13	1.86	0.56
1:CA:1270:C:H5''	1:CA:1271:G:O5'	2.05	0.56
1:CA:2554:U:H2'	1:CA:2555:U:C6	2.39	0.56
7:CG:117:PHE:CE1	7:CG:119:GLY:HA2	2.40	0.56
1:CA:64:A:O3'	21:CX:71:GLY:HA3	2.05	0.56
34:DA:1412:C:H2'	34:DA:1413:A:H8	1.69	0.56
50:DQ:76:LEU:HD12	50:DQ:77:VAL:H	1.69	0.56
57:DZ:326:THR:HG21	57:DZ:380:LEU:HD23	1.87	0.56
57:DZ:555:LEU:HD11	57:DZ:599:PRO:HB2	1.87	0.56
33:A9:2:LYS:HE2	33:A9:31:LYS:O	2.05	0.56
1:AA:1084:C:N4	1:AA:1163:G:H1	2.03	0.56
3:AC:6:LYS:HG3	3:AC:7:ARG:H	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:AL:99:ILE:O	10:AL:139:VAL:N	2.36	0.56
12:AO:64:ARG:HG2	12:AO:79:PHE:CG	2.40	0.56
13:AP:81:GLN:OE1	13:AP:106:LEU:HD23	2.05	0.56
38:BE:78:HIS:HD2	38:BE:142:LEU:HD23	1.69	0.56
42:BI:9:ARG:H	42:BI:79:LEU:HD23	1.69	0.56
48:BO:69:TYR:HA	48:BO:72:ARG:HH11	1.70	0.56
6:CF:167:ALA:HB1	6:CF:173:VAL:HG11	1.87	0.56
10:CL:6:ALA:HB3	10:CL:30:HIS:CE1	2.36	0.56
34:DA:1220:G:N2	52:DS:54:GLY:O	2.38	0.56
35:DB:58:ILE:HA	35:DB:61:LEU:HB3	1.86	0.56
1:AA:1487:G:H2'	1:AA:1488:G:H8	1.70	0.56
1:AA:2379:G:N7	63:AA:5202:HOH:O	2.33	0.56
1:AA:2736:C:OP2	5:AE:109:LYS:HE2	2.06	0.56
7:AG:16:ARG:O	7:AG:20:ILE:HG13	2.06	0.56
34:BA:394:G:H2'	34:BA:395:C:H6	1.70	0.56
35:BB:212:GLN:NE2	35:BB:234:PRO:O	2.38	0.56
40:BG:138:LYS:NZ	40:BG:142:GLU:OE2	2.34	0.56
49:BP:6:LEU:HB3	49:BP:17:TYR:CD1	2.40	0.56
53:BT:9:ASN:O	53:BT:10:LEU:HB2	2.05	0.56
32:C8:34:TRP:CD2	32:C8:35:GLN:HG2	2.40	0.56
1:CA:1657:C:H2'	1:CA:1658:C:C6	2.41	0.56
1:CA:2692:C:H2'	1:CA:2693:A:H8	1.70	0.56
6:CF:157:VAL:HG11	6:CF:181:LEU:HD12	1.87	0.56
22:CY:77:PRO:HD3	22:CY:106:LEU:HD23	1.86	0.56
34:DA:392:G:H2'	34:DA:393:A:C8	2.39	0.56
34:DA:685:G:C2	34:DA:686:U:C4	2.93	0.56
57:DZ:15:ILE:HB	57:DZ:104:ALA:HA	1.87	0.56
57:DZ:222:ASP:N	57:DZ:222:ASP:OD2	2.35	0.56
1:AA:1532:A:H2'	1:AA:1533:G:H8	1.70	0.56
1:AA:1250:U:C2	6:AF:171:PRO:HB3	2.41	0.56
36:BC:18:TRP:CD1	47:BN:54:PRO:HA	2.41	0.56
49:BP:18:ARG:O	49:BP:20:VAL:HB	2.06	0.56
3:CC:49:GLY:N	3:CC:209:PHE:O	2.39	0.56
34:DA:174:C:H2'	34:DA:175:C:H6	1.70	0.56
35:DB:51:LEU:HD11	35:DB:201:ILE:HG23	1.88	0.56
35:DB:93:VAL:HG21	35:DB:97:TRP:CD1	2.40	0.56
37:DD:3:ARG:HG2	37:DD:118:ARG:NE	2.20	0.56
57:DZ:373:ASP:OD2	57:DZ:374:LEU:N	2.33	0.56
57:DZ:546:ILE:HG23	57:DZ:590:ILE:HG13	1.86	0.56
1:AA:815:G:O2'	1:AA:1425:A:N1	2.35	0.56
1:AA:2326:C:H2'	1:AA:2327:G:C8	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:650:G:N7	13:AP:107:LYS:NZ	2.54	0.56
1:AA:2851:C:OP1	15:AR:53:HIS:NE2	2.39	0.56
21:AX:63:LYS:O	21:AX:64:LYS:HD3	2.04	0.56
1:AA:2672:A:N6	57:BZ:661:SER:OG	2.38	0.56
1:CA:1899:G:O2'	1:CA:1900:A:OP2	2.18	0.56
1:CA:528:A:C2	1:CA:2043:C:H4'	2.39	0.56
2:CB:100:A:H3'	2:CB:101:G:H8	1.71	0.56
1:CA:1022:G:N7	11:CN:66:LYS:HE2	2.20	0.56
6:CF:34:TRP:CZ2	13:CP:8:PRO:HG3	2.40	0.56
1:CA:566:U:O4	19:CV:78:LYS:HE2	2.05	0.56
34:DA:519:C:H2'	34:DA:520:A:O4'	2.06	0.56
56:DY:55:PSU:N3	56:DY:58:A:N7	2.47	0.56
1:AA:1405:A:C2	1:AA:1418:U:O4	2.58	0.56
1:AA:2340:A:H2'	1:AA:2341:G:C8	2.40	0.56
1:AA:2574:U:H1'	12:AO:23:ARG:HD3	1.87	0.56
1:AA:2807:C:N4	1:AA:2813:G:H1	2.03	0.56
11:AN:120:LEU:HD22	11:AN:122:VAL:HG23	1.88	0.56
34:BA:507:C:OP2	34:BA:508:C:O2'	2.17	0.56
34:BA:559:A:H4'	34:BA:560:U:H3'	1.87	0.56
37:BD:23:GLY:HA3	37:BD:112:VAL:HG12	1.88	0.56
42:BI:5:TYR:OH	42:BI:7:THR:OG1	2.21	0.56
57:BZ:257:PRO:HB2	57:BZ:259:PHE:HE1	1.69	0.56
33:C9:9:ARG:HG2	33:C9:14:CYS:HB2	1.87	0.56
1:CA:2343:C:O2'	1:CA:2373:G:O2'	2.16	0.56
1:CA:2693:A:H2'	1:CA:2694:G:C8	2.39	0.56
5:CE:135:HIS:H	5:CE:135:HIS:CD2	2.21	0.56
34:DA:1150:U:O4	34:DA:1151:A:N6	2.39	0.56
34:DA:691:G:H2'	34:DA:692:U:C6	2.41	0.56
35:DB:104:ASN:HB3	35:DB:108:ILE:HD11	1.88	0.56
38:DE:139:LEU:O	38:DE:141:GLN:N	2.37	0.56
57:DZ:13:ARG:HH21	57:DZ:247:ARG:HH12	1.51	0.56
57:DZ:32:ILE:HG23	57:DZ:273:LEU:HD21	1.88	0.56
1:AA:597:C:H4'	1:AA:598:A:O5'	2.05	0.56
1:AA:613:A:OP1	6:AF:95:ARG:NH1	2.38	0.56
48:BO:74:ASP:CG	48:BO:77:ARG:HG3	2.26	0.56
57:BZ:494:GLU:HG2	57:BZ:511:LYS:HG2	1.88	0.56
1:CA:1090:U:H2'	1:CA:1091:G:C8	2.40	0.56
1:CA:271(E):U:H2'	1:CA:271(F):C:C6	2.40	0.56
1:CA:898:C:H2'	1:CA:899:A:O4'	2.05	0.56
1:CA:987:G:O2'	1:CA:1000:A:N3	2.34	0.56
23:CZ:138:GLU:H	23:CZ:156:LYS:HZ1	1.54	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:CZ:160:GLY:HA2	23:CZ:161:VAL:HG12	1.88	0.56
34:DA:652:U:O4	34:DA:752:G:O2'	2.16	0.56
37:DD:118:ARG:O	37:DD:121:VAL:N	2.38	0.56
1:AA:2661:U:H2'	1:AA:2662:U:C6	2.41	0.56
13:AP:94:GLU:HG3	13:AP:124:LYS:HB3	1.88	0.56
34:BA:1305:G:H5''	54:BU:4:GLY:HA3	1.87	0.56
34:BA:1388:C:H2'	34:BA:1389:C:C6	2.41	0.56
34:BA:1530:G:H2'	34:BA:1531:A:O4'	2.04	0.56
34:BA:328:C:H4'	34:BA:329:A:H5''	1.88	0.56
1:CA:747:U:O2	1:CA:2014:A:H1'	2.06	0.56
1:CA:477:A:H2'	1:CA:478:A:C8	2.40	0.56
1:CA:869:G:N1	1:CA:909:A:C6	2.74	0.56
1:CA:2124:G:H4'	3:CC:175:PRO:CD	2.35	0.56
3:CC:42:VAL:O	3:CC:216:THR:O	2.24	0.56
4:CD:75:ILE:HD12	4:CD:75:ILE:H	1.70	0.56
34:DA:1237:C:H2'	34:DA:1336:C:C5	2.41	0.56
1:CA:2572:A:C8	5:CE:144:ARG:HD2	2.40	0.56
8:CH:159:GLU:HG3	8:CH:169:VAL:HG11	1.87	0.56
13:CP:139:LYS:O	13:CP:141:ALA:N	2.38	0.56
34:DA:537:G:H2'	34:DA:538:G:H8	1.71	0.56
36:DC:71:ALA:HB2	36:DC:106:VAL:HB	1.87	0.56
37:DD:10:ARG:HA	37:DD:13:ARG:HG3	1.86	0.56
37:DD:150:GLU:HA	37:DD:153:ARG:HE	1.71	0.56
1:AA:1900:G:H2'	1:AA:1901:C:C6	2.41	0.56
1:AA:2262:G:OP1	14:AQ:85:LYS:NZ	2.22	0.56
1:AA:2798:C:OP1	5:AE:41:LYS:HE3	2.05	0.56
1:AA:533:G:N2	20:AW:80:PRO:HG2	2.21	0.56
1:AA:1856:A:OP1	4:AD:249:PRO:HD3	2.06	0.56
5:AE:36:ARG:HH11	5:AE:85:ASN:ND2	2.03	0.56
7:AG:15:VAL:HG22	7:AG:175:LEU:HB3	1.88	0.56
34:BA:1401:G:C2	34:BA:1402:C:H1'	2.41	0.56
34:BA:1496:C:OP2	57:BZ:504:ARG:NH2	2.38	0.56
45:BL:71:PRO:O	45:BL:102:ARG:NH1	2.33	0.56
13:CP:59:LEU:O	32:C8:13:ARG:HD2	2.06	0.56
1:CA:1266:G:O2'	1:CA:2012:G:O6	2.20	0.56
1:CA:2641:G:P	11:CN:74:ARG:HH21	2.29	0.56
13:CP:94:GLU:HG3	13:CP:124:LYS:HD3	1.86	0.56
23:CZ:63:ASP:O	23:CZ:65:GLN:N	2.34	0.56
34:DA:1022:G:H2'	34:DA:1023:G:C8	2.41	0.56
24:A0:27:GLU:HG3	24:A0:68:GLU:HA	1.86	0.56
1:AA:418:G:H1'	1:AA:438:G:O4'	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:165:THR:OG1	7:AG:168:GLU:HG3	2.06	0.56
45:BL:34:ARG:HG3	45:BL:105:TYR:CE2	2.41	0.56
49:BP:6:LEU:HB3	49:BP:17:TYR:HD1	1.71	0.56
49:BP:3:LYS:O	49:BP:21:VAL:HA	2.06	0.56
50:BQ:86:GLU:C	50:BQ:88:TYR:H	2.09	0.56
1:CA:127:A:H5''	1:CA:128:C:C6	2.41	0.56
1:CA:1657:C:H2'	1:CA:1658:C:H6	1.71	0.56
1:CA:2379:G:H5''	1:CA:2379:G:H8	1.71	0.56
1:CA:321:G:OP2	6:CF:135:LYS:HG3	2.05	0.56
1:CA:796:C:H2'	1:CA:797:C:C6	2.41	0.56
1:CA:829:A:N7	1:CA:2248:C:H5'	2.21	0.56
7:CG:41:GLN:NE2	7:CG:154:GLY:O	2.27	0.56
34:DA:97:G:O2'	34:DA:98:G:OP2	2.19	0.56
34:DA:407:G:O2'	37:DD:116:GLN:HG3	2.06	0.56
57:DZ:176:GLY:HA2	57:DZ:187:THR:HA	1.88	0.56
57:DZ:484:ARG:HD2	57:DZ:676:TYR:CE1	2.40	0.56
1:AA:1114:G:O2'	1:AA:1142:A:O2'	2.24	0.55
1:AA:2571:C:O2'	1:AA:2572:C:H5'	2.07	0.55
1:AA:2679:C:H2'	1:AA:2680:G:O4'	2.05	0.55
22:AY:8:LYS:HD3	22:AY:97:ARG:HH12	1.70	0.55
23:AZ:110:GLY:N	23:AZ:144:LEU:O	2.39	0.55
23:AZ:150:LEU:O	23:AZ:171:ILE:HG13	2.07	0.55
34:BA:1053:G:O2'	34:BA:1199:U:OP2	2.14	0.55
34:BA:1326:C:H2'	34:BA:1327:C:C6	2.40	0.55
34:BA:353:A:H5'	34:BA:353:A:H8	1.71	0.55
34:BA:742:G:OP2	48:BO:35:ARG:NH2	2.38	0.55
35:BB:19:HIS:HE1	35:BB:189:ASP:HB3	1.71	0.55
35:BB:218:ALA:O	35:BB:222:ILE:HG13	2.07	0.55
36:BC:3:ASN:OD1	36:BC:3:ASN:N	2.37	0.55
1:CA:1778:U:H2'	1:CA:1784:A:N6	2.21	0.55
1:CA:205:G:OP2	63:CA:4260:HOH:O	2.18	0.55
1:CA:2474:C:H5''	1:CA:2475:C:OP2	2.06	0.55
15:CR:85:PRO:C	15:CR:87:TYR:H	2.09	0.55
34:DA:1318:A:OP1	52:DS:3:ARG:NH1	2.38	0.55
50:DQ:63:ARG:HG2	50:DQ:64:PRO:HD2	1.88	0.55
57:DZ:132:ARG:HD3	57:DZ:160:ARG:CZ	2.36	0.55
57:DZ:573:HIS:CE1	57:DZ:575:VAL:HB	2.41	0.55
1:AA:1099:C:H2'	1:AA:1100:A:H5''	1.87	0.55
1:AA:2153:G:OP1	3:AC:6:LYS:CD	2.54	0.55
1:AA:2331:G:H1	16:AS:3:ARG:HA	1.71	0.55
34:BA:1177:G:H2'	34:BA:1178:G:O4'	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:617:G:C6	34:BA:618:C:C5	2.94	0.55
52:BS:33:THR:OG1	52:BS:34:TRP:N	2.39	0.55
31:C7:19:ARG:HG2	31:C7:19:ARG:HH11	1.71	0.55
8:CH:46:GLU:HB2	8:CH:49:VAL:HG12	1.87	0.55
35:DB:171:ALA:O	35:DB:175:ARG:N	2.37	0.55
36:DC:8:ILE:HD13	36:DC:184:TYR:HD2	1.72	0.55
37:DD:64:LEU:HD11	37:DD:94:LEU:HD21	1.89	0.55
39:DF:35:ALA:HA	39:DF:67:MET:HB3	1.87	0.55
1:AA:1211:U:H2'	1:AA:1212:C:C6	2.40	0.55
1:AA:2585:C:H3'	63:AA:4115:HOH:O	2.05	0.55
21:AX:57:LEU:CD1	21:AX:78:LYS:HB2	2.36	0.55
34:BA:1005:A:H1'	34:BA:1036:G:H22	1.71	0.55
34:BA:736:C:H2'	34:BA:737:A:C8	2.40	0.55
57:BZ:177:ILE:HD12	57:BZ:188:TYR:CE2	2.41	0.55
57:BZ:498:ILE:HB	57:BZ:507:TYR:CD1	2.41	0.55
1:CA:1796:U:H2'	1:CA:1797:C:C6	2.42	0.55
1:CA:479:A:N3	1:CA:481:G:H5''	2.21	0.55
1:CA:2176:A:C4'	3:CC:45:HIS:NE2	2.67	0.55
13:CP:62:LEU:HD11	32:C8:50:LEU:HD11	1.87	0.55
17:CT:16:ARG:NH1	17:CT:18:ASP:OD2	2.40	0.55
34:DA:173:U:H5''	34:DA:197:A:O4'	2.07	0.55
43:DJ:13:HIS:HB3	43:DJ:68:HIS:CE1	2.41	0.55
55:DV:13:A:O2'	55:DV:14:A:H5''	2.06	0.55
57:DZ:264:LEU:HD12	62:DZ:704:GDP:N3	2.20	0.55
57:DZ:512:ILE:HG13	57:DZ:514:VAL:HG23	1.88	0.55
32:A8:23:VAL:HG11	32:A8:47:LYS:HD3	1.89	0.55
1:AA:1417:G:H2'	1:AA:1418:U:H5	1.70	0.55
1:AA:1634:C:H2'	1:AA:1635:C:C6	2.41	0.55
8:AH:3:ARG:HG2	8:AH:6:ARG:HG2	1.88	0.55
34:BA:200:G:H1	34:BA:217:C:N4	2.04	0.55
49:BP:4:ILE:HA	49:BP:20:VAL:O	2.07	0.55
57:BZ:416:LYS:HE2	57:BZ:475:ASN:HD22	1.70	0.55
1:CA:2059:A:O2'	6:CF:69:HIS:HD2	1.89	0.55
7:CG:126:ASP:HB2	7:CG:130:ASN:H	1.71	0.55
7:CG:29:TRP:O	7:CG:33:ARG:NH1	2.38	0.55
34:DA:339:C:H2'	34:DA:340:U:C6	2.41	0.55
36:DC:45:LYS:HG3	36:DC:46:GLU:HG2	1.87	0.55
1:AA:1558:G:H2'	1:AA:1559:C:O4'	2.07	0.55
1:AA:935:C:N3	1:AA:936:C:N4	2.54	0.55
34:BA:598:U:H4'	41:BH:94:TYR:CD2	2.42	0.55
35:BB:73:THR:OG1	35:BB:170:GLU:OE2	2.20	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BD:110:PHE:HE2	37:BD:148:VAL:HG23	1.71	0.55
39:BF:61:LEU:HD12	39:BF:63:TYR:OH	2.07	0.55
51:BR:39:VAL:O	51:BR:42:ARG:HB2	2.06	0.55
57:BZ:227:ILE:HG12	57:BZ:237:PRO:HG3	1.89	0.55
14:CQ:109:VAL:HG13	14:CQ:113:GLN:HB3	1.88	0.55
19:CV:37:VAL:HG12	19:CV:39:LEU:H	1.71	0.55
34:DA:1409:C:H2'	34:DA:1410:G:H8	1.71	0.55
57:DZ:25:LYS:NZ	62:DZ:704:GDP:O2B	2.27	0.55
1:AA:1735:U:O2	1:AA:1747:A:H5'	2.05	0.55
1:AA:554:A:H62	1:AA:2063:U:H3	1.55	0.55
22:AY:92:ASN:CB	22:AY:94:LYS:H	2.19	0.55
34:BA:1000:U:H2'	34:BA:1001:A:H8	1.71	0.55
34:BA:1227:A:OP2	46:BM:111:LYS:HD2	2.07	0.55
12:AO:49:ARG:NH2	34:BA:1423:G:OP1	2.30	0.55
57:BZ:309:LEU:HA	57:BZ:333:GLY:HA3	1.87	0.55
7:CG:136:ARG:HH11	7:CG:137:GLU:H	1.54	0.55
35:DB:16:HIS:HB2	35:DB:204:ASN:CB	2.36	0.55
57:DZ:38:ARG:NH1	57:DZ:270:GLN:NE2	2.55	0.55
57:DZ:-37:LEU:HD23	57:DZ:-32:LEU:HD12	1.89	0.55
1:AA:1653:C:H4'	1:AA:1654:A:O5'	2.06	0.55
1:AA:561:A:H2'	1:AA:562:C:C6	2.41	0.55
34:BA:1225:A:H2'	34:BA:1226:C:C5	2.42	0.55
34:BA:489:C:OP1	37:BD:132:ARG:NH2	2.39	0.55
34:BA:127:G:OP1	34:BA:635:G:H1'	2.06	0.55
35:BB:174:VAL:O	35:BB:178:ARG:HB2	2.06	0.55
36:BC:19:GLU:OE1	36:BC:40:ARG:NH2	2.40	0.55
42:BI:23:ASN:ND2	42:BI:25:LYS:HG2	2.22	0.55
56:BW:7:A:H5''	56:BW:7:A:H8	1.70	0.55
1:CA:1494:A:H2'	1:CA:1495:A:C8	2.42	0.55
1:CA:1510:G:H2'	1:CA:1511:C:O4'	2.07	0.55
1:CA:500:G:N2	1:CA:502:A:H3'	2.21	0.55
3:CC:52:PRO:HB2	3:CC:168:LYS:O	2.06	0.55
6:CF:153:SER:OG	6:CF:190:GLU:N	2.38	0.55
10:CL:21:PRO:HD2	57:DZ:641:GLN:NE2	2.21	0.55
11:CN:19:GLU:HG3	11:CN:59:LYS:HB3	1.89	0.55
15:CR:38:VAL:HB	15:CR:39:PRO:HD3	1.89	0.55
15:CR:72:ASP:OD2	15:CR:75:LEU:HB2	2.06	0.55
34:DA:1260:C:O5'	34:DA:1284:C:H4'	2.07	0.55
35:DB:95:GLN:HB2	35:DB:148:TYR:HD1	1.72	0.55
40:DG:113:GLU:HB2	40:DG:119:ARG:HG2	1.87	0.55
57:DZ:132:ARG:H	57:DZ:132:ARG:HD2	1.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:264:LEU:HB2	62:DZ:704:GDP:C6	2.42	0.55
1:AA:2140:U:C4	1:AA:2171:G:H1'	2.42	0.55
35:BB:17:PHE:HA	35:BB:44:LEU:HD11	1.88	0.55
37:BD:134:ASP:O	37:BD:136:PRO:HD3	2.06	0.55
57:BZ:644:ARG:CB	57:BZ:644:ARG:HH11	2.10	0.55
1:CA:2022:U:OP2	29:C5:15:ARG:NH2	2.40	0.55
1:CA:30:G:H2'	1:CA:31:C:C6	2.41	0.55
2:CB:56:G:H5'	7:CG:27:ASN:HD22	1.70	0.55
34:DA:363:A:OP2	45:DL:34:ARG:NH2	2.40	0.55
34:DA:622:A:C8	34:DA:623:C:C6	2.94	0.55
37:DD:15:GLU:OE1	37:DD:63:LYS:HG3	2.06	0.55
38:DE:57:LYS:HD3	38:DE:61:TYR:CE2	2.41	0.55
43:DJ:63:PHE:HE1	47:DN:58:LYS:HG2	1.71	0.55
49:DP:4:ILE:O	49:DP:66:PRO:HA	2.06	0.55
34:DA:263:A:OP1	53:DT:79:ARG:NH1	2.40	0.55
1:AA:1072:U:H4'	1:AA:1073:A:OP1	2.07	0.55
1:AA:142:G:H1'	21:AX:37:THR:HG21	1.88	0.55
1:AA:278:G:H2'	1:AA:279:G:H5''	1.89	0.55
34:BA:1435:G:H2'	34:BA:1436:U:C6	2.41	0.55
51:BR:47:THR:HG23	51:BR:49:LYS:HG3	1.89	0.55
57:BZ:127:LYS:HB3	57:BZ:128:TYR:CD2	2.41	0.55
57:BZ:329:ARG:HB2	57:BZ:374:LEU:HG	1.89	0.55
26:C2:35:LEU:HD12	26:C2:53:LEU:HD12	1.88	0.55
33:C9:25:VAL:HG11	33:C9:34:GLN:NE2	2.21	0.55
1:CA:375:C:H2'	1:CA:376:C:C6	2.42	0.55
1:CA:528:A:N1	1:CA:2042:A:H2'	2.22	0.55
7:CG:146:TYR:O	7:CG:149:VAL:HG12	2.07	0.55
34:DA:1141:C:H2'	34:DA:1142:G:H8	1.71	0.55
34:DA:1305:G:N2	34:DA:1331:G:H1'	2.21	0.55
34:DA:481:G:H21	34:DA:482:A:N6	2.05	0.55
13:AP:63:PRO:HG2	32:A8:25:MET:HB2	1.89	0.55
1:AA:1525:G:HO2'	1:AA:1605:A:H2	1.53	0.55
3:AC:42:VAL:O	3:AC:216:THR:O	2.24	0.55
23:AZ:115:GLY:HA2	23:AZ:177:PRO:HB3	1.89	0.55
57:BZ:445:GLU:OE1	57:BZ:484:ARG:NH1	2.39	0.55
26:C2:2:LYS:O	26:C2:5:GLU:N	2.40	0.55
1:CA:2386:C:H2'	1:CA:2387:U:C6	2.42	0.55
1:CA:754:C:H2'	1:CA:755:C:C6	2.42	0.55
3:CC:44:VAL:HG23	3:CC:176:VAL:HG21	1.89	0.55
1:CA:2132:U:C1'	3:CC:6:LYS:CB	2.86	0.55
34:DA:275:G:H5'	50:DQ:14:LYS:HB3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DI:23:ASN:ND2	42:DI:23:ASN:H	2.05	0.55
44:DK:123:LYS:O	44:DK:126:ARG:HG3	2.06	0.55
1:AA:1553:A:O2'	1:AA:1554:A:O5'	2.25	0.54
1:AA:354:A:HO2'	1:AA:355:A:H8	1.53	0.54
3:AC:194:ILE:HD11	3:AC:227:PRO:HB3	1.89	0.54
3:AC:54:ARG:CZ	3:AC:56:ASP:HB3	2.37	0.54
16:AS:82:ILE:HD12	16:AS:82:ILE:H	1.71	0.54
17:AT:119:LYS:HB2	34:BA:1442(A):G:N2	2.22	0.54
57:BZ:12:LEU:HD12	57:BZ:78:ARG:HB3	1.90	0.54
28:C4:59:PHE:HA	28:C4:61:ARG:N	2.23	0.54
1:CA:2102:U:H2'	1:CA:2103:C:C6	2.42	0.54
3:CC:6:LYS:HA	3:CC:9:ARG:HH11	1.73	0.54
34:DA:1289:A:N1	34:DA:1371:G:O2'	2.32	0.54
34:DA:67:C:H2'	34:DA:68:G:C8	2.41	0.54
34:DA:93:G:H2'	34:DA:96:U:O4'	2.06	0.54
38:DE:8:GLU:HG2	38:DE:34:VAL:HG23	1.89	0.54
38:DE:36:ASP:OD2	38:DE:40:ARG:HB2	2.06	0.54
40:DG:93:PRO:HA	40:DG:96:GLN:HB2	1.88	0.54
43:DJ:49:VAL:HG23	47:DN:41:ARG:HB2	1.88	0.54
51:DR:70:ILE:O	51:DR:74:ARG:HG3	2.07	0.54
53:DT:53:LEU:HA	53:DT:56:MET:HG2	1.89	0.54
27:A3:59:VAL:O	27:A3:60:GLU:HG2	2.07	0.54
1:AA:1218:G:O2'	1:AA:1219:A:O4'	2.21	0.54
1:AA:1305:G:O2'	1:AA:1306:G:H5'	2.08	0.54
1:AA:1848:G:C2'	1:AA:1849:U:H5'	2.37	0.54
5:AE:128:SER:OG	5:AE:129:HIS:N	2.38	0.54
1:AA:2324:U:H5'	7:AG:88:ILE:HD11	1.89	0.54
23:AZ:45:ASP:O	23:AZ:49:ARG:HG3	2.07	0.54
35:BB:17:PHE:HB2	35:BB:44:LEU:HD21	1.89	0.54
42:BI:4:TYR:CE1	42:BI:88:TYR:HA	2.42	0.54
45:BL:31:PRO:HB2	45:BL:32:PHE:CD2	2.42	0.54
57:BZ:179:ASP:OD2	57:BZ:182:ARG:HD2	2.07	0.54
30:C6:21:TYR:CE1	30:C6:38:LYS:HG2	2.42	0.54
1:CA:1404:C:O2'	1:CA:1405:U:H5'	2.06	0.54
1:CA:749:C:O2	1:CA:1618:A:H2'	2.07	0.54
1:CA:639:U:H2'	1:CA:640:C:C6	2.41	0.54
5:CE:60:ASN:OD1	5:CE:62:PRO:HD2	2.07	0.54
11:CN:128:HIS:CE1	11:CN:135:PRO:HG2	2.43	0.54
13:CP:91:PHE:O	13:CP:121:LYS:NZ	2.39	0.54
23:CZ:92:SER:O	23:CZ:130:PRO:HG2	2.07	0.54
34:DA:1504:G:OP1	34:DA:1507:A:H4'	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:435:C:H2'	34:DA:436:C:C6	2.42	0.54
38:DE:53:LEU:H	38:DE:53:LEU:HD12	1.71	0.54
1:AA:553:A:C2	1:AA:2064:A:H2'	2.42	0.54
13:AP:83:VAL:HG13	13:AP:112:LEU:HD21	1.89	0.54
34:BA:100:C:H2'	34:BA:101:A:O4'	2.06	0.54
34:BA:161:A:H2'	34:BA:162:A:C8	2.42	0.54
35:BB:20:GLU:HA	35:BB:21:ARG:NH2	2.21	0.54
39:BF:70:ASP:HB2	39:BF:71:ARG:HG2	1.89	0.54
46:BM:11:ARG:HA	46:BM:45:VAL:HB	1.89	0.54
36:BC:6:HIS:HB2	47:BN:49:HIS:CD2	2.41	0.54
1:CA:1969:A:O2'	1:CA:1972:A:N3	2.31	0.54
1:CA:631:A:OP1	13:CP:65:ARG:NH1	2.39	0.54
1:CA:98:G:OP1	26:C2:3:LEU:N	2.31	0.54
14:CQ:37:LEU:HB2	14:CQ:128:LYS:HB2	1.89	0.54
34:DA:137:C:H2'	34:DA:138:G:C8	2.42	0.54
34:DA:45:U:O5'	34:DA:45:U:H6	1.90	0.54
34:DA:499:A:H4'	34:DA:500:G:OP1	2.08	0.54
20:AW:19:LEU:HB3	29:A5:25:LEU:HD11	1.90	0.54
1:AA:1093:G:N2	1:AA:1157:A:H2	2.04	0.54
3:AC:49:GLY:N	3:AC:209:PHE:O	2.39	0.54
37:BD:104:VAL:O	37:BD:107:ARG:N	2.40	0.54
42:BI:86:VAL:O	42:BI:90:PRO:HG3	2.07	0.54
34:BA:1216:G:OP1	47:BN:2:ALA:N	2.40	0.54
57:BZ:443:HIS:ND1	57:BZ:446:THR:HG22	2.22	0.54
1:CA:1914:C:H5''	1:CA:1915:U:OP2	2.08	0.54
1:CA:2389:G:H5''	1:CA:2390:U:O4'	2.06	0.54
1:CA:2394:C:OP1	32:C8:30:ARG:NH1	2.40	0.54
1:CA:221:A:O2'	1:CA:266:G:N7	2.32	0.54
1:CA:45:C:OP2	1:CA:215:G:H5'	2.08	0.54
1:CA:2121:G:O2'	3:CC:168:LYS:CD	2.54	0.54
10:CL:19:PRO:HB3	10:CL:34:ILE:HD12	1.89	0.54
16:CS:23:ARG:HH21	16:CS:84:GLN:HB3	1.71	0.54
23:CZ:63:ASP:C	23:CZ:65:GLN:H	2.10	0.54
41:DH:25:ASP:OD1	41:DH:25:ASP:N	2.40	0.54
50:DQ:56:VAL:HB	50:DQ:78:GLU:HB3	1.88	0.54
1:AA:1286:U:H3'	63:AA:4783:HOH:O	2.06	0.54
3:AC:52:PRO:HB2	3:AC:168:LYS:O	2.07	0.54
19:AV:52:VAL:HG22	19:AV:55:ALA:HB3	1.90	0.54
34:BA:1095:U:H2'	34:BA:1096:C:O4'	2.07	0.54
34:BA:167:G:H2'	34:BA:168:G:H8	1.72	0.54
35:BB:204:ASN:OD1	35:BB:206:ASP:N	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
37:BD:31:CYS:SG	37:BD:33:MET:N	2.80	0.54
39:BF:97:PHE:CB	51:BR:32:ARG:HD2	2.37	0.54
52:BS:40:ILE:HB	52:BS:67:VAL:HA	1.89	0.54
57:BZ:406:GLU:HG2	57:BZ:439:ARG:HH22	1.73	0.54
25:C1:23:LYS:HB3	25:C1:29:GLY:HA3	1.90	0.54
5:CE:2:LYS:HG3	5:CE:200:GLU:HB2	1.90	0.54
34:DA:909:A:H2'	34:DA:910:C:O4'	2.07	0.54
36:DC:137:ALA:HA	36:DC:140:ARG:HD3	1.90	0.54
53:DT:16:HIS:O	53:DT:19:SER:OG	2.16	0.54
3:AC:52:PRO:HG2	3:AC:53:ARG:H	1.73	0.54
6:AF:150:GLY:HA2	6:AF:172:TRP:CD2	2.43	0.54
34:BA:677:U:H3	34:BA:713:G:H22	1.56	0.54
38:BE:67:VAL:HG12	38:BE:69:VAL:HG12	1.88	0.54
38:BE:93:PRO:HG2	41:BH:105:ARG:HE	1.72	0.54
57:BZ:329:ARG:HD3	57:BZ:331:TYR:CZ	2.42	0.54
26:C2:10:LEU:HD21	26:C2:59:ARG:HD2	1.90	0.54
31:C7:34:ARG:NH1	31:C7:41:ARG:O	2.41	0.54
1:CA:1843:C:H5'	4:CD:253:GLN:NE2	2.22	0.54
1:CA:1970:A:H4'	1:CA:1971:A:OP1	2.08	0.54
1:CA:299:A:N3	1:CA:319:C:O2'	2.33	0.54
10:CL:75:SER:HA	10:CL:78:ILE:HG22	1.88	0.54
14:CQ:38:GLU:OE2	14:CQ:128:LYS:HG2	2.08	0.54
12:CO:49:ARG:NH1	34:DA:1423:G:OP1	2.41	0.54
32:A8:42:ARG:HD2	63:A8:6307:HOH:O	2.08	0.54
1:AA:611:U:H2'	1:AA:612:C:C6	2.43	0.54
23:AZ:28:MET:HE3	23:AZ:59:LEU:HD12	1.89	0.54
34:BA:1068:G:H8	34:BA:1068:G:OP2	1.90	0.54
34:BA:193:C:H2'	34:BA:194:C:C6	2.42	0.54
34:BA:373:A:C2	34:BA:374:A:C8	2.94	0.54
34:BA:975:A:N1	43:BJ:48:THR:HB	2.22	0.54
37:BD:149:ALA:HB3	37:BD:152:SER:HB2	1.89	0.54
57:BZ:466:LEU:O	57:BZ:470:PHE:HB2	2.07	0.54
1:CA:1297:C:H2'	1:CA:1298:C:H6	1.72	0.54
1:CA:1297:C:OP1	1:CA:2710:C:H4'	2.08	0.54
1:CA:2726:U:O2'	1:CA:2727:G:H5'	2.08	0.54
1:CA:861:A:C2	1:CA:917:A:C4	2.96	0.54
3:CC:30:VAL:HG23	3:CC:31:LYS:HG2	1.89	0.54
3:CC:52:PRO:HG2	3:CC:53:ARG:H	1.73	0.54
1:CA:7:G:H4'	11:CN:13:TRP:HH2	1.72	0.54
21:CX:57:LEU:HD13	21:CX:78:LYS:HB3	1.89	0.54
34:DA:416:G:C5	34:DA:417:C:C4	2.96	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DB:47:THR:O	35:DB:51:LEU:HD22	2.07	0.54
57:DZ:35:TYR:HE2	57:DZ:269:VAL:HB	1.73	0.54
26:A2:2:LYS:O	26:A2:6:VAL:HG23	2.08	0.54
1:AA:1633:A:H2'	1:AA:1634:C:C6	2.42	0.54
4:AD:8:PRO:CB	4:AD:14:ARG:HB2	2.38	0.54
14:AQ:16:ARG:C	14:AQ:17:LEU:HD23	2.27	0.54
34:BA:1086:U:H2'	34:BA:1087:G:O4'	2.07	0.54
34:BA:1338:G:H2'	34:BA:1339:A:C8	2.42	0.54
34:BA:435:C:C2	34:BA:436:C:C5	2.96	0.54
34:BA:487:A:H2'	34:BA:488:C:O4'	2.08	0.54
34:BA:643:C:H5'	41:BH:31:PHE:CD1	2.43	0.54
46:BM:3:ARG:NH2	46:BM:11:ARG:HH12	2.04	0.54
34:BA:134:A:H61	49:BP:25:ARG:HH12	1.56	0.54
1:CA:1786:A:H1'	1:CA:1938:A:N6	2.22	0.54
1:CA:2590:A:O2'	1:CA:2591:C:H5'	2.07	0.54
5:CE:2:LYS:HB2	5:CE:95:ILE:HD12	1.90	0.54
8:CH:154:PRO:HA	8:CH:161:GLY:HA3	1.90	0.54
14:CQ:70:PRO:HA	14:CQ:95:ALA:HB2	1.90	0.54
15:CR:13:HIS:O	15:CR:15:SER:N	2.41	0.54
50:DQ:40:LYS:HD3	50:DQ:42:TYR:CZ	2.42	0.54
34:DA:333:G:H4'	53:DT:16:HIS:CE1	2.42	0.54
56:DW:40:C:H4'	56:DY:35:A:O3'	2.08	0.54
1:CA:2432:A:H4'	56:DY:76:A:P	2.47	0.54
25:A1:80:LEU:HB3	25:A1:82:LEU:HG	1.90	0.54
34:BA:1070:U:H2'	34:BA:1071:C:C6	2.43	0.54
34:BA:685:G:C2	34:BA:686:U:C4	2.96	0.54
34:BA:973:G:OP1	43:BJ:57:LYS:HE3	2.08	0.54
35:BB:12:GLU:HA	35:BB:213:LEU:HD11	1.89	0.54
36:BC:155:GLY:HA3	36:BC:196:LEU:HD12	1.88	0.54
1:CA:1708:C:H2'	1:CA:1709:U:H6	1.72	0.54
1:CA:601:C:O2'	1:CA:605:C:H5''	2.08	0.54
1:CA:192:C:O2'	1:CA:802:A:N3	2.37	0.54
5:CE:59:VAL:HG12	5:CE:64:LYS:HE2	1.90	0.54
20:CW:29:LEU:O	20:CW:33:ARG:HG3	2.08	0.54
34:DA:1207:G:H2'	34:DA:1208:C:H6	1.71	0.54
34:DA:1281:U:P	34:DA:1282:C:H41	2.30	0.54
34:DA:177:C:H2'	34:DA:178:C:H6	1.73	0.54
35:DB:25:ASN:O	35:DB:27:LYS:N	2.40	0.54
37:DD:25:ARG:HG2	37:DD:25:ARG:O	2.08	0.54
40:DG:26:PHE:O	40:DG:30:ILE:HG13	2.08	0.54
1:AA:1476:C:H2'	1:AA:1477:U:C6	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:AG:123:ASN:O	63:AG:301:HOH:O	2.18	0.54
8:AH:4:ILE:O	8:AH:69:ARG:HG2	2.07	0.54
34:BA:630:G:O2'	34:BA:631:G:H5'	2.08	0.54
39:BF:10:LEU:HB2	39:BF:59:TYR:HB3	1.90	0.54
56:BY:53:G:H1	56:BY:61:C:N4	2.06	0.54
57:BZ:503:GLY:O	57:BZ:505:GLY:N	2.41	0.54
57:BZ:484:ARG:HG2	57:BZ:561:VAL:HG22	1.90	0.54
1:CA:1991:U:H2'	1:CA:1992:G:H5''	1.89	0.54
1:CA:375:C:H2'	1:CA:376:C:H6	1.73	0.54
2:CB:62:C:H2'	2:CB:63:G:C8	2.41	0.54
3:CC:171:ALA:HB1	3:CC:173:HIS:CE1	2.43	0.54
5:CE:119:ARG:HD2	5:CE:120:TRP:CE2	2.43	0.54
5:CE:101:ARG:NH1	5:CE:169:ASN:O	2.40	0.54
34:DA:1004:A:H8	34:DA:1005:A:H4'	1.73	0.54
41:DH:116:LYS:HD2	41:DH:129:VAL:HG11	1.90	0.54
57:DZ:514:VAL:HG22	57:DZ:565:VAL:HA	1.90	0.54
57:DZ:620:VAL:O	57:DZ:624:LEU:HB2	2.08	0.54
24:A0:43:THR:HG23	24:A0:43:THR:O	2.08	0.53
1:AA:559:U:H2'	1:AA:560:C:C6	2.43	0.53
4:AD:242:ARG:HD3	4:AD:242:ARG:N	2.23	0.53
6:AF:20:LEU:HD22	6:AF:21:ALA:O	2.08	0.53
7:AG:131:TYR:O	7:AG:159:VAL:HG12	2.08	0.53
11:AN:65:LYS:NZ	11:AN:65:LYS:HB2	2.23	0.53
23:AZ:63:ASP:OD1	23:AZ:65:GLN:HB2	2.07	0.53
34:BA:502:G:C2	34:BA:503:C:C2	2.96	0.53
50:BQ:88:TYR:HD2	50:BQ:89:LEU:HD23	1.71	0.53
52:BS:50:ALA:HB1	52:BS:57:HIS:HB3	1.90	0.53
57:BZ:655:TYR:CE2	57:BZ:659:LEU:HG	2.42	0.53
38:DE:80:ILE:HD13	41:DH:104:ARG:HH21	1.72	0.53
48:DO:70:LEU:HD23	48:DO:78:TYR:HA	1.88	0.53
57:DZ:250:THR:HG21	57:DZ:279:TYR:O	2.08	0.53
25:A1:6:GLU:HB2	25:A1:61:ARG:O	2.08	0.53
1:AA:1336:C:H2'	1:AA:1337:C:C6	2.41	0.53
1:AA:1540:A:H2'	1:AA:1541:A:C8	2.43	0.53
1:AA:1604:C:H5''	1:AA:1605:A:OP2	2.08	0.53
1:AA:2096:U:H2'	1:AA:2097:U:C6	2.43	0.53
1:AA:254:A:N6	1:AA:454:U:O2'	2.41	0.53
3:AC:171:ALA:HB1	3:AC:173:HIS:CE1	2.44	0.53
34:BA:345:C:O5'	34:BA:345:C:H6	1.91	0.53
34:BA:300:A:H1'	34:BA:565:U:O2	2.08	0.53
39:BF:37:VAL:HG12	39:BF:38:GLU:H	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BG:115:ARG:HH11	40:BG:118:VAL:HG21	1.73	0.53
1:CA:2355:C:H1'	24:C0:39:ARG:HH21	1.73	0.53
32:C8:36:LYS:HB2	32:C8:41:ILE:HD11	1.90	0.53
1:CA:2454:G:H1'	63:CA:3887:HOH:O	2.07	0.53
1:CA:2880:C:O3'	15:CR:90:ARG:NH1	2.37	0.53
1:CA:2892:A:H2'	1:CA:2893:G:C8	2.43	0.53
3:CC:194:ILE:HD11	3:CC:227:PRO:HB3	1.89	0.53
17:CT:26:ASP:O	17:CT:49:VAL:HG12	2.09	0.53
34:DA:1073:U:OP1	38:DE:57:LYS:HE3	2.08	0.53
34:DA:160:A:H1'	34:DA:344:A:C5	2.43	0.53
1:AA:1136:U:O2	1:AA:1148:C:H1'	2.08	0.53
3:AC:42:VAL:HG13	3:AC:43:GLU:H	1.73	0.53
34:BA:1241:G:H2'	34:BA:1242:C:C6	2.43	0.53
34:BA:196:A:OP1	53:BT:68:LYS:NZ	2.40	0.53
34:BA:626:U:H2'	34:BA:627:G:C8	2.43	0.53
39:BF:94:GLN:OE1	51:BR:32:ARG:NH2	2.42	0.53
50:BQ:59:ILE:HG22	50:BQ:73:VAL:HA	1.90	0.53
46:BM:118:ALA:HB1	56:BW:28:G:H4'	1.89	0.53
57:BZ:289:ILE:HD11	57:BZ:331:TYR:CG	2.43	0.53
57:BZ:490:PRO:HG3	57:BZ:515:GLU:HB3	1.90	0.53
1:CA:1558:A:H4'	1:CA:1559:G:O5'	2.08	0.53
1:CA:2315:G:H2'	1:CA:2316:C:C6	2.43	0.53
1:CA:2408:U:OP2	63:CA:3857:HOH:O	2.19	0.53
8:CH:80:SER:OG	8:CH:81:GLU:N	2.41	0.53
20:CW:6:ILE:HD13	20:CW:104:THR:HG23	1.90	0.53
34:DA:827:U:H5''	34:DA:828:A:OP2	2.08	0.53
37:DD:3:ARG:NH1	37:DD:5:ILE:HG13	2.24	0.53
57:DZ:490:PRO:HG3	57:DZ:516:PRO:HD2	1.89	0.53
1:AA:1093:G:N2	1:AA:1156:G:O2'	2.41	0.53
1:AA:2198:A:H2'	1:AA:2199:C:C6	2.42	0.53
1:AA:2209:G:O2'	1:AA:2210:C:OP1	2.24	0.53
1:AA:2576:A:OP1	1:AA:2660:C:O2'	2.23	0.53
34:BA:1510:U:H2'	34:BA:1511:G:C8	2.42	0.53
34:BA:396:G:P	57:BZ:349:LYS:HZ1	2.32	0.53
35:BB:45:GLN:O	35:BB:49:GLU:HB2	2.08	0.53
57:BZ:269:VAL:O	57:BZ:272:LEU:HB3	2.08	0.53
1:CA:1792:G:O2'	1:CA:1830:C:OP1	2.26	0.53
5:CE:176:ILE:HB	5:CE:181:LEU:HB2	1.91	0.53
11:CN:120:LEU:HD22	11:CN:122:VAL:HG23	1.90	0.53
23:CZ:8:TYR:HB2	23:CZ:38:TYR:CZ	2.44	0.53
34:DA:1142:G:H2'	34:DA:1143:G:O4'	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:269:C:H2'	34:DA:270:A:C8	2.43	0.53
34:DA:828:A:H2'	34:DA:829:G:O4'	2.08	0.53
40:DG:50:ILE:HD11	40:DG:58:PRO:HB3	1.89	0.53
41:DH:26:VAL:HG23	41:DH:27:PRO:O	2.09	0.53
39:DF:99:ALA:O	51:DR:28:GLU:HA	2.08	0.53
1:AA:354:A:H2	1:AA:1255:A:O2'	1.92	0.53
13:AP:39:LYS:HG3	13:AP:45:LEU:HD11	1.90	0.53
1:AA:325:G:OP2	22:AY:84:ARG:NH2	2.42	0.53
54:BU:3:LYS:HA	54:BU:11:GLY:HA2	1.88	0.53
1:CA:1070:A:H2'	1:CA:1097:U:OP1	2.08	0.53
1:CA:1860:G:OP2	1:CA:1860:G:H8	1.92	0.53
1:CA:2837:G:O2'	1:CA:2838:G:H5'	2.09	0.53
1:CA:2850:A:H2'	1:CA:2851:A:C8	2.43	0.53
2:CB:54:G:H2'	2:CB:55:U:H6	1.74	0.53
4:CD:108:PRO:HD2	4:CD:111:LEU:HG	1.89	0.53
11:CN:42:TRP:HD1	11:CN:48:MET:HE1	1.74	0.53
1:CA:329:G:OP2	22:CY:71:LYS:HD2	2.08	0.53
34:DA:448:A:C4	34:DA:487:A:C2	2.96	0.53
34:DA:642:A:N3	41:DH:113:SER:OG	2.37	0.53
34:DA:737:A:H2'	34:DA:738:C:C6	2.44	0.53
34:DA:953:G:H5'	34:DA:965:A:H61	1.73	0.53
40:DG:30:ILE:HD13	40:DG:120:ILE:HD13	1.91	0.53
1:AA:1587:U:H2'	1:AA:1588:G:O4'	2.09	0.53
1:AA:1604:C:OP2	1:AA:1605:A:O2'	2.24	0.53
1:AA:721:G:H1'	6:AF:74:ARG:HD3	1.90	0.53
3:AC:30:VAL:HG23	3:AC:31:LYS:HG2	1.89	0.53
1:AA:2325:C:H4'	7:AG:91:ARG:HG3	1.89	0.53
1:AA:1941:A:O3'	34:BA:1517:G:H1'	2.09	0.53
48:BO:24:SER:O	48:BO:28:GLN:N	2.34	0.53
53:BT:87:LYS:O	53:BT:91:LEU:HG	2.09	0.53
1:CA:1769:G:O2'	1:CA:1958:C:OP1	2.20	0.53
1:CA:583:G:OP2	18:CU:10:ARG:HD2	2.09	0.53
5:CE:21:VAL:HG23	5:CE:185:LYS:HD2	1.90	0.53
8:CH:107:VAL:HG11	8:CH:162:ILE:HD11	1.90	0.53
13:CP:99:LEU:HA	13:CP:102:ARG:HB2	1.90	0.53
18:CU:92:ARG:HA	18:CU:95:LEU:HB2	1.88	0.53
34:DA:1129:C:H2'	34:DA:1139:G:N7	2.24	0.53
34:DA:1304:G:C6	34:DA:1305:G:N1	2.77	0.53
35:DB:12:GLU:C	35:DB:14:GLY:HA3	2.28	0.53
38:DE:131:ILE:O	38:DE:135:THR:OG1	2.26	0.53
39:DF:78:GLU:C	39:DF:80:ARG:H	2.11	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1133:G:C2	1:AA:1149:A:C2	2.97	0.53
1:AA:2294:G:H4'	1:AA:2401:G:O2'	2.09	0.53
7:AG:111:LEU:HD22	7:AG:120:LEU:HD21	1.89	0.53
34:BA:1207:G:H2'	34:BA:1208:C:C6	2.44	0.53
49:BP:75:ARG:O	49:BP:78:GLY:N	2.31	0.53
57:BZ:330:VAL:CG1	57:BZ:371:ALA:HA	2.37	0.53
1:CA:1059:G:H5''	1:CA:1060:U:H2'	1.91	0.53
1:CA:2531:A:H5''	8:CH:157:TYR:CE2	2.44	0.53
1:CA:301:G:OP2	22:CY:84:ARG:NH2	2.42	0.53
1:CA:579:G:H2'	1:CA:580:C:C6	2.44	0.53
3:CC:51:ASP:OD2	3:CC:54:ARG:HB2	2.09	0.53
3:CC:54:ARG:CZ	3:CC:56:ASP:HB3	2.37	0.53
7:CG:19:LEU:HG	7:CG:175:LEU:HD22	1.91	0.53
34:DA:107:G:H2'	34:DA:108:G:O4'	2.08	0.53
37:DD:121:VAL:HG22	37:DD:126:ILE:HG13	1.91	0.53
57:DZ:526:VAL:HG11	57:DZ:566:THR:HG23	1.91	0.53
1:AA:1014:U:O3'	27:A3:14:GLY:HA2	2.08	0.53
1:AA:1100:A:H2'	1:AA:1101:G:O4'	2.09	0.53
1:AA:2044:U:O2'	1:AA:2629:C:H5'	2.09	0.53
1:AA:2150:C:H4'	3:AC:219:MET:HE3	1.91	0.53
3:AC:6:LYS:HA	3:AC:9:ARG:HH11	1.72	0.53
8:AH:11:VAL:HG13	8:AH:15:VAL:HG22	1.91	0.53
11:AN:12:ARG:HH21	11:AN:138:LEU:HD21	1.73	0.53
34:BA:1278:U:H5'	34:BA:1279:A:H5'	1.91	0.53
34:BA:1431:C:H2'	34:BA:1432:G:O4'	2.09	0.53
34:BA:148:G:H2'	34:BA:149:A:C8	2.43	0.53
34:BA:1516:G:N1	34:BA:1519:A:OP2	2.42	0.53
39:BF:19:LEU:HD11	39:BF:59:TYR:CE2	2.44	0.53
57:BZ:608:VAL:HG21	57:BZ:647:VAL:HG22	1.91	0.53
25:C1:25:LYS:C	25:C1:27:GLU:H	2.12	0.53
1:CA:2689:U:P	1:CA:2719:G:H22	2.32	0.53
1:CA:2869:G:H2'	1:CA:2870:C:O4'	2.09	0.53
2:CB:116:G:OP2	2:CB:116:G:H8	1.92	0.53
4:CD:142:VAL:HG12	4:CD:163:ALA:HB3	1.91	0.53
8:CH:157:TYR:HE1	8:CH:172:LYS:HG2	1.74	0.53
1:CA:910:A:H62	14:CQ:12:GLN:HA	1.73	0.53
16:CS:65:VAL:O	16:CS:68:GLN:HB2	2.09	0.53
17:CT:60:THR:HG22	17:CT:77:PRO:HA	1.91	0.53
34:DA:786:G:H2'	34:DA:787:A:O4'	2.09	0.53
37:DD:129:ASN:HD21	37:DD:144:ASP:HA	1.74	0.53
48:DO:5:LYS:HD2	48:DO:5:LYS:H	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:609:GLU:HG2	57:DZ:644:ARG:HG2	1.91	0.53
1:AA:1325:G:H4'	15:AR:31:HIS:CD2	2.44	0.53
1:AA:1358:U:H4'	1:AA:1359:U:O5'	2.09	0.53
1:AA:1571:G:H2'	1:AA:1572:G:O4'	2.09	0.53
1:AA:1686:U:O2'	1:AA:1687:C:H5'	2.08	0.53
7:AG:178:PHE:O	7:AG:180:PHE:HD2	1.91	0.53
38:BE:93:PRO:HG2	41:BH:105:ARG:NE	2.24	0.53
34:BA:1292:U:OP2	40:BG:41:ARG:NH2	2.41	0.53
43:BJ:17:ASP:O	43:BJ:21:GLN:HB2	2.09	0.53
44:BK:99:GLN:HG2	44:BK:105:VAL:HG11	1.91	0.53
52:BS:36:ARG:NH1	52:BS:52:TYR:O	2.37	0.53
57:BZ:273:LEU:O	57:BZ:276:VAL:N	2.42	0.53
57:BZ:490:PRO:HB3	57:BZ:515:GLU:HG2	1.91	0.53
27:C3:18:ASP:OD1	27:C3:18:ASP:N	2.42	0.53
1:CA:2120:G:H21	3:CC:168:LYS:CE	2.20	0.53
2:CB:66:A:N6	2:CB:108:U:H3'	2.24	0.53
4:CD:3:VAL:HG13	4:CD:17:THR:HB	1.90	0.53
1:CA:323:G:C8	6:CF:171:PRO:HG3	2.44	0.53
34:DA:612:C:O2	34:DA:629:G:N2	2.41	0.53
29:A5:16:ARG:NH1	29:A5:17:ASP:OD1	2.39	0.53
1:AA:1188:A:C4	1:AA:1190:G:C8	2.97	0.53
1:AA:1452:U:H2'	1:AA:1453:C:H6	1.74	0.53
2:AB:89:G:H2'	2:AB:90:A:C8	2.43	0.53
3:AC:44:VAL:HG23	3:AC:176:VAL:HG21	1.89	0.53
17:AT:108:ARG:HH12	17:AT:112:ARG:HE	1.57	0.53
34:BA:368:U:N3	57:BZ:354:ARG:NH1	2.50	0.53
34:BA:375:U:N3	34:BA:376:G:N7	2.57	0.53
34:BA:38:G:H22	34:BA:397:A:H5''	1.73	0.53
35:BB:21:ARG:HA	35:BB:39:ILE:HG23	1.90	0.53
43:BJ:45:ARG:HG2	43:BJ:47:PHE:CZ	2.44	0.53
1:CA:1046:A:H3'	1:CA:1047:G:H5'	1.91	0.53
1:CA:1049:C:H3'	1:CA:1050:A:H8	1.73	0.53
1:CA:1639:U:H2'	1:CA:1640:C:H5''	1.91	0.53
1:CA:2887:U:H2'	1:CA:2888:C:H6	1.74	0.53
1:CA:443:A:H1'	1:CA:1201:C:O4'	2.08	0.53
1:CA:2682:U:H5'	5:CE:11:MET:O	2.08	0.53
11:CN:14:VAL:HG12	11:CN:15:LEU:H	1.73	0.53
16:CS:66:ALA:O	16:CS:69:VAL:N	2.42	0.53
20:CW:20:VAL:O	20:CW:23:LEU:HB2	2.09	0.53
34:DA:253:U:H2'	34:DA:254:G:C8	2.44	0.53
34:DA:544:G:OP1	37:DD:59:ARG:NH2	2.31	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:1179:A:H4'	42:DI:103:THR:HA	1.90	0.53
47:DN:7:ILE:HG22	47:DN:23:ARG:HD2	1.90	0.53
57:DZ:486:THR:OG1	57:DZ:487:ILE:N	2.42	0.53
31:A7:33:ARG:NH2	63:A7:201:HOH:O	2.41	0.52
1:AA:1405:A:N6	1:AA:1418:U:H3	2.07	0.52
3:AC:65:LEU:HD22	3:AC:189:ASN:HB3	1.91	0.52
23:AZ:23:LYS:HD2	23:AZ:40:ASP:HA	1.90	0.52
34:BA:251:G:H4'	34:BA:252:U:O5'	2.09	0.52
34:BA:731:G:H5'	34:BA:766:A:H4'	1.91	0.52
35:BB:51:LEU:HD22	35:BB:55:PHE:CE2	2.44	0.52
35:BB:77:ALA:HB2	35:BB:211:ILE:HD13	1.90	0.52
45:BL:34:ARG:HG2	45:BL:35:GLY:H	1.74	0.52
46:BM:20:THR:HA	46:BM:25:ILE:O	2.08	0.52
49:BP:18:ARG:NH1	49:BP:32:TYR:OH	2.42	0.52
57:BZ:182:ARG:O	57:BZ:184:LYS:N	2.42	0.52
57:BZ:342:TYR:N	57:BZ:390:VAL:O	2.38	0.52
1:CA:1068:G:H21	1:CA:1096:A:H5'	1.74	0.52
1:CA:2572:A:N7	5:CE:144:ARG:HD2	2.25	0.52
3:CC:48:LEU:CB	3:CC:50:ILE:HD12	2.38	0.52
5:CE:70:ALA:O	5:CE:72:VAL:N	2.38	0.52
6:CF:25:PRO:HD2	6:CF:115:ALA:HB2	1.91	0.52
17:CT:99:LEU:HD22	17:CT:101:PHE:HE1	1.74	0.52
17:CT:66:VAL:HA	17:CT:71:GLY:HA2	1.91	0.52
23:CZ:85:HIS:HE1	23:CZ:87:ASP:OD2	1.92	0.52
34:DA:938:A:H2'	34:DA:939:G:O4'	2.09	0.52
37:DD:60:GLU:OE1	37:DD:199:ASN:N	2.40	0.52
38:DE:53:LEU:O	38:DE:57:LYS:HB2	2.08	0.52
57:DZ:612:THR:OG1	57:DZ:613:PRO:O	2.26	0.52
1:AA:2584:A:N7	5:AE:144:ARG:HD2	2.24	0.52
1:AA:2880:C:H2'	1:AA:2881:C:O4'	2.08	0.52
1:AA:934:A:OP1	1:AA:935:C:N4	2.42	0.52
2:AB:24:G:H3'	63:AB:3101:HOH:O	2.08	0.52
7:AG:110:ALA:HB1	7:AG:140:ILE:HG22	1.91	0.52
34:BA:1243:C:N3	34:BA:1295:G:N2	2.57	0.52
35:BB:111:ARG:NH1	35:BB:111:ARG:HG2	2.20	0.52
50:BQ:88:TYR:CD2	50:BQ:89:LEU:HD23	2.44	0.52
57:BZ:133:ILE:HG22	57:BZ:257:PRO:HB2	1.90	0.52
1:CA:1639:U:C2'	1:CA:1640:C:H5''	2.39	0.52
1:CA:1651:G:N2	1:CA:2007:C:C2	2.77	0.52
1:CA:2148:G:H2'	1:CA:2149:G:C8	2.43	0.52
1:CA:2458:G:O2'	1:CA:2460:U:O4	2.21	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:272:G:H4'	1:CA:272(A):U:H5''	1.90	0.52
14:CQ:37:LEU:HD12	14:CQ:128:LYS:HB3	1.91	0.52
20:CW:76:VAL:HG22	20:CW:103:ILE:HG23	1.92	0.52
34:DA:1219:U:OP1	47:DN:19:ARG:NH1	2.41	0.52
34:DA:1347:G:O2'	34:DA:1373:G:O6	2.27	0.52
34:DA:148:G:H2'	34:DA:149:A:C8	2.43	0.52
36:DC:182:ILE:HG12	36:DC:203:PHE:HD1	1.74	0.52
57:DZ:546:ILE:HD12	57:DZ:546:ILE:H	1.74	0.52
57:DZ:534:ILE:HD11	57:DZ:570:GLY:HA3	1.91	0.52
1:AA:1698:G:OP1	15:AR:40:LYS:HE3	2.08	0.52
1:AA:346:A:H4'	1:AA:347:G:OP2	2.09	0.52
3:AC:48:LEU:CB	3:AC:50:ILE:HD12	2.38	0.52
3:AC:64:SER:HA	3:AC:161:ARG:H	1.74	0.52
13:AP:120:ALA:HB1	13:AP:138:LEU:HD12	1.92	0.52
13:AP:132:LYS:O	13:AP:136:GLU:HG3	2.08	0.52
34:BA:659:U:C2	34:BA:660:G:C8	2.98	0.52
41:BH:2:LEU:HD11	41:BH:8:ASP:HB2	1.91	0.52
45:BL:24:VAL:HB	45:BL:27:LEU:HD22	1.92	0.52
56:BW:35:A:H4'	57:BZ:575:VAL:HG21	1.92	0.52
1:CA:2712:U:H2'	1:CA:2714:G:H5''	1.90	0.52
1:CA:664:C:H2'	1:CA:665:C:H6	1.74	0.52
13:CP:97:PRO:HG3	13:CP:112:LEU:HD12	1.91	0.52
34:DA:1189:C:O5'	34:DA:1189:C:H6	1.92	0.52
34:DA:1057:G:N2	34:DA:1204:A:H1'	2.24	0.52
38:DE:127:ASN:O	38:DE:131:ILE:HG12	2.10	0.52
53:DT:59:ALA:O	53:DT:62:LEU:N	2.42	0.52
1:AA:1222:A:N3	1:AA:1222:A:H2'	2.24	0.52
1:AA:1636:U:H2'	1:AA:1637:G:C8	2.45	0.52
1:AA:2187:G:H1	1:AA:2194:U:H5	1.57	0.52
1:AA:238:C:O2'	13:AP:64:LYS:HE3	2.09	0.52
1:AA:2413:U:OP1	30:A6:18:ARG:NH2	2.42	0.52
14:AQ:27:VAL:N	14:AQ:138:ASP:OD1	2.39	0.52
1:AA:1199:C:OP1	18:AU:92:ARG:NH1	2.43	0.52
34:BA:1131:G:OP1	42:BI:20:ARG:NH2	2.41	0.52
34:BA:310:G:OP2	49:BP:27:LYS:NZ	2.38	0.52
34:BA:397:A:N6	34:BA:548:G:C5	2.78	0.52
34:BA:922:G:C6	34:BA:923:A:C6	2.98	0.52
53:BT:82:SER:O	53:BT:86:ARG:HG3	2.09	0.52
57:BZ:166:LEU:HB3	57:BZ:178:ILE:HD12	1.92	0.52
57:BZ:273:LEU:HA	57:BZ:276:VAL:HG23	1.92	0.52
57:BZ:323:GLY:O	57:BZ:325:LEU:N	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:C1:67:ILE:N	25:C1:68:PRO:HD2	2.24	0.52
1:CA:1001:A:H2'	1:CA:1002:G:O4'	2.08	0.52
1:CA:2166:G:H3'	1:CA:2167:U:H5''	1.91	0.52
1:CA:2648:C:H2'	1:CA:2649:U:C6	2.44	0.52
1:CA:394:A:O2'	1:CA:395:U:H5'	2.09	0.52
34:DA:1053:G:O5'	34:DA:1054:C:H5'	2.10	0.52
34:DA:411:A:H62	34:DA:413:G:H21	1.55	0.52
34:DA:933:G:O6	40:DG:3:ARG:NH2	2.40	0.52
35:DB:93:VAL:HG21	35:DB:97:TRP:HD1	1.74	0.52
37:DD:3:ARG:HE	37:DD:118:ARG:HD3	1.75	0.52
44:DK:48:ILE:O	44:DK:50:TYR:N	2.42	0.52
56:DW:53:G:N2	56:DW:61:C:O2	2.36	0.52
57:DZ:363:ARG:HG2	57:DZ:363:ARG:NH1	2.24	0.52
7:AG:66:GLN:HG3	28:A4:1:MET:CE	2.40	0.52
1:AA:1106:U:O4	10:AL:130:SER:OG	2.11	0.52
1:AA:2211:U:H2'	1:AA:2212:G:C8	2.44	0.52
1:AA:70:A:N7	21:AX:31:HIS:HE1	2.06	0.52
1:AA:1857:G:H4'	4:AD:242:ARG:CZ	2.39	0.52
7:AG:27:ASN:HB3	7:AG:30:GLU:HG3	1.92	0.52
7:AG:47:LYS:O	7:AG:51:ARG:HG2	2.09	0.52
10:AL:59:ILE:HD11	10:AL:63:ARG:HA	1.90	0.52
1:AA:1296:G:N7	13:AP:18:ARG:NH2	2.57	0.52
34:BA:402:G:C6	34:BA:403:C:C4	2.98	0.52
34:BA:714:G:H2'	34:BA:715:A:C8	2.45	0.52
34:BA:814:A:N7	34:BA:816:A:C4	2.77	0.52
57:BZ:443:HIS:ND1	57:BZ:445:GLU:O	2.39	0.52
57:BZ:627:ARG:HH22	57:BZ:658:ASP:CG	2.12	0.52
27:C3:30:ARG:HB3	27:C3:33:GLN:HB2	1.92	0.52
1:CA:108:U:H2'	1:CA:109:G:H8	1.75	0.52
1:CA:2028:U:H2'	1:CA:2029:G:O4'	2.10	0.52
1:CA:2756:U:H1'	1:CA:2757:A:H5''	1.91	0.52
1:CA:635:C:O2'	1:CA:639:U:OP1	2.27	0.52
8:CH:24:VAL:HG13	8:CH:37:VAL:HG21	1.91	0.52
1:CA:1058:G:O2'	10:CL:114:ASP:O	2.26	0.52
15:CR:51:LEU:HD23	15:CR:66:VAL:HG22	1.91	0.52
16:CS:69:VAL:O	16:CS:72:ALA:HB3	2.10	0.52
22:CY:86:ARG:HB2	22:CY:98:VAL:HG23	1.91	0.52
41:DH:20:TYR:HD2	41:DH:65:TYR:CE2	2.27	0.52
44:DK:59:TYR:CE1	44:DK:63:LEU:HD21	2.44	0.52
1:AA:1475:G:H2'	1:AA:1476:C:H6	1.74	0.52
1:AA:2143:G:N2	3:AC:169:THR:CB	2.68	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:27:G:N7	63:AA:4832:HOH:O	2.34	0.52
1:AA:616:G:O4'	32:A8:4:MET:HE2	2.09	0.52
1:AA:2200:C:OP1	3:AC:47:LYS:HG2	2.09	0.52
13:AP:52:GLU:HB2	13:AP:55:ARG:HD2	1.91	0.52
23:AZ:30:ASN:ND2	23:AZ:90:VAL:HB	2.23	0.52
37:BD:55:ALA:O	37:BD:58:LEU:N	2.43	0.52
43:BJ:5:ARG:NH2	43:BJ:73:ASP:OD2	2.43	0.52
46:BM:15:VAL:O	46:BM:19:LEU:HD13	2.09	0.52
28:C4:36:CYS:SG	28:C4:37:SER:N	2.83	0.52
1:CA:1401:G:C6	1:CA:1402:C:C4	2.98	0.52
1:CA:2682:U:OP2	63:CA:3790:HOH:O	2.19	0.52
1:CA:2727:G:O3'	12:CO:70:LYS:NZ	2.38	0.52
1:CA:740:U:H2'	1:CA:741:G:C8	2.45	0.52
1:CA:996:A:C2	1:CA:997:G:C8	2.97	0.52
4:CD:121:PRO:HB3	4:CD:135:PHE:CE2	2.45	0.52
6:CF:53:THR:HG22	6:CF:56:GLU:HG3	1.92	0.52
6:CF:53:THR:HB	6:CF:56:GLU:OE2	2.09	0.52
10:CL:99:ILE:O	10:CL:139:VAL:N	2.41	0.52
21:CX:43:VAL:HG21	21:CX:81:VAL:HG11	1.91	0.52
34:DA:1095:U:H5'	34:DA:1109:C:O2	2.10	0.52
34:DA:1224:G:O2'	34:DA:1322:C:OP1	2.26	0.52
35:DB:100:GLY:O	35:DB:104:ASN:N	2.39	0.52
38:DE:143:ARG:HD2	41:DH:77:GLU:OE2	2.10	0.52
36:DC:22:TRP:HA	43:DJ:93:GLY:HA2	1.91	0.52
57:DZ:528:ALA:O	57:DZ:568:TYR:HA	2.10	0.52
6:AF:34:TRP:CH2	13:AP:8:PRO:HB3	2.45	0.52
13:AP:121:LYS:O	13:AP:123:LEU:N	2.43	0.52
34:BA:1063:C:H3'	34:BA:1064:G:H2'	1.92	0.52
34:BA:175:C:H2'	34:BA:176:C:C6	2.44	0.52
34:BA:44:G:H2'	34:BA:45:U:O4'	2.09	0.52
34:BA:509:A:H3'	34:BA:509:A:H8	1.73	0.52
38:BE:43:LEU:HD21	38:BE:132:ALA:HB1	1.92	0.52
52:BS:40:ILE:HG12	52:BS:71:LEU:HD12	1.91	0.52
1:CA:1653:G:H3'	15:CR:2:ARG:HD3	1.92	0.52
1:CA:2120:G:N2	3:CC:168:LYS:HE2	2.23	0.52
1:CA:2683:C:H4'	5:CE:13:ARG:NH2	2.24	0.52
13:CP:84:ASN:CG	13:CP:117:GLU:HB2	2.29	0.52
34:DA:1161:C:H2'	34:DA:1162:C:H6	1.75	0.52
44:DK:98:LEU:O	44:DK:101:SER:OG	2.19	0.52
49:DP:51:VAL:O	49:DP:53:VAL:HG23	2.10	0.52
33:A9:15:LYS:HG2	33:A9:17:ILE:HD13	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1851:U:C2	4:AD:202:LYS:HG2	2.45	0.52
3:AC:51:ASP:OD2	3:AC:54:ARG:HB2	2.09	0.52
4:AD:147:LEU:HD22	4:AD:155:LEU:HD11	1.91	0.52
15:AR:97:VAL:HG22	15:AR:114:VAL:HG13	1.92	0.52
23:AZ:110:GLY:O	23:AZ:113:ALA:HB3	2.10	0.52
34:BA:153:C:H42	34:BA:169:C:N4	2.06	0.52
34:BA:368:U:C4	57:BZ:354:ARG:NH1	2.77	0.52
34:BA:735:C:H2'	34:BA:736:C:H6	1.75	0.52
35:BB:108:ILE:O	35:BB:111:ARG:HB2	2.10	0.52
35:BB:62:ALA:HB1	35:BB:226:ARG:HD3	1.92	0.52
37:BD:108:LEU:HB3	37:BD:110:PHE:CE1	2.45	0.52
38:BE:100:VAL:HG22	38:BE:118:ILE:HG22	1.92	0.52
57:BZ:89:ASP:OD1	57:BZ:457:LEU:HB2	2.09	0.52
1:CA:2653:U:O2'	8:CH:110:SER:HB2	2.10	0.52
11:CN:112:LEU:O	11:CN:115:ARG:N	2.42	0.52
16:CS:105:ALA:O	16:CS:110:LEU:HB2	2.10	0.52
21:CX:53:LYS:HB3	21:CX:82:GLN:HB3	1.90	0.52
23:CZ:125:LEU:HB3	23:CZ:165:VAL:HG13	1.90	0.52
34:DA:149:A:H2'	34:DA:150:C:C6	2.45	0.52
34:DA:253:U:H2'	34:DA:254:G:H8	1.74	0.52
34:DA:255:G:H1'	50:DQ:16:GLN:HE21	1.75	0.52
34:DA:977:A:H2'	34:DA:978:A:H5''	1.92	0.52
35:DB:178:ARG:O	41:DH:71:GLY:HA2	2.08	0.52
34:DA:522:C:H41	45:DL:53:ARG:NH2	2.05	0.52
56:DW:73:A:H5''	56:DW:74:C:H5'	1.91	0.52
1:AA:1525:G:H2'	1:AA:1526:G:H8	1.75	0.52
1:AA:2124:U:H2'	1:AA:2125:C:C6	2.44	0.52
1:AA:692:C:H2'	1:AA:693:G:H8	1.75	0.52
3:AC:29:LEU:O	3:AC:30:VAL:C	2.48	0.52
10:AL:134:MET:HG3	10:AL:136:VAL:HG12	1.91	0.52
1:AA:1068:G:N7	11:AN:66:LYS:HE2	2.24	0.52
12:AO:100:GLY:O	12:AO:119:PRO:HD2	2.10	0.52
41:BH:87:SER:HA	41:BH:93:VAL:HG23	1.92	0.52
57:BZ:217:VAL:HG22	57:BZ:242:LEU:HD21	1.92	0.52
1:CA:2327:A:H2'	1:CA:2328:A:C8	2.44	0.52
1:CA:2600:A:N6	63:CA:3928:HOH:O	2.42	0.52
1:CA:570:G:H2'	1:CA:2030:A:C5	2.44	0.52
2:CB:15:A:H1'	2:CB:110:G:C5	2.45	0.52
3:CC:67:HIS:CG	3:CC:185:LYS:HD2	2.45	0.52
4:CD:148:GLU:OE1	4:CD:151:LYS:NZ	2.33	0.52
7:CG:37:VAL:HG23	7:CG:99:MET:HG3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:CQ:122:GLY:HA2	14:CQ:125:LEU:HD12	1.92	0.52
18:CU:65:ILE:O	18:CU:69:CYS:HB2	2.09	0.52
20:CW:14:PRO:HG2	20:CW:78:GLU:HG2	1.91	0.52
34:DA:229:U:H5''	49:DP:33:ILE:HD13	1.92	0.52
43:DJ:55:LYS:O	43:DJ:57:LYS:N	2.42	0.52
1:AA:1846:A:O3'	63:AA:4736:HOH:O	2.19	0.52
1:AA:2891:C:H2'	1:AA:2892:A:O4'	2.10	0.52
4:AD:133:LEU:HB3	4:AD:173:VAL:HG11	1.92	0.52
4:AD:18:VAL:HG12	4:AD:211:ARG:NH2	2.25	0.52
4:AD:206:LEU:HD22	4:AD:211:ARG:HG2	1.91	0.52
17:AT:15:VAL:HG13	17:AT:79:HIS:CE1	2.46	0.52
38:BE:109:ILE:HD12	38:BE:135:THR:HB	1.91	0.52
34:BA:559:A:P	38:BE:126:ARG:HH22	2.31	0.52
44:BK:115:PRO:HB2	44:BK:118:GLY:H	1.75	0.52
49:BP:54:GLU:HG3	49:BP:55:ARG:N	2.24	0.52
57:BZ:539:ILE:HA	57:BZ:542:VAL:HG12	1.91	0.52
1:CA:2788:C:O2'	1:CA:2809:A:N3	2.39	0.52
1:CA:374:A:H2'	1:CA:375:C:H5'	1.92	0.52
1:CA:527:C:H4'	1:CA:528:A:O5'	2.10	0.52
1:CA:857:C:H2'	1:CA:858:U:C6	2.44	0.52
3:CC:65:LEU:HD22	3:CC:189:ASN:HB3	1.91	0.52
1:CA:2177:C:O2'	3:CC:47:LYS:HD3	2.09	0.52
13:CP:81:GLN:NE2	13:CP:105:LEU:O	2.42	0.52
20:CW:12:ILE:HD13	20:CW:17:VAL:HG13	1.92	0.52
1:CA:1337:G:OP2	21:CX:73:ARG:NH2	2.42	0.52
34:DA:297:G:N2	34:DA:300:A:OP2	2.42	0.52
34:DA:433:C:H2'	34:DA:434:U:C6	2.44	0.52
34:DA:967:C:H6	34:DA:967:C:O5'	1.93	0.52
38:DE:107:ARG:HG2	38:DE:108:ALA:N	2.25	0.52
46:DM:37:THR:HG21	46:DM:56:LEU:HA	1.92	0.52
48:DO:64:ARG:O	48:DO:68:ARG:N	2.43	0.52
1:AA:1067:A:H8	1:AA:1067:A:H3'	1.72	0.51
1:AA:1218:G:OP2	1:AA:1218:G:H2'	2.10	0.51
1:AA:1817:A:H1'	1:AA:1960:A:N6	2.25	0.51
1:AA:2705:A:H2'	1:AA:2706:G:H8	1.75	0.51
3:AC:67:HIS:CG	3:AC:185:LYS:HD2	2.45	0.51
11:AN:51:PHE:CZ	11:AN:119:ARG:HG2	2.45	0.51
1:AA:2022:G:OP1	15:AR:5:LYS:NZ	2.43	0.51
34:BA:15:G:C4	34:BA:16:A:C8	2.98	0.51
35:BB:145:LEU:O	35:BB:149:LEU:HB2	2.09	0.51
36:BC:186:PHE:CG	36:BC:187:ALA:N	2.78	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:1191:A:H5''	36:BC:4:LYS:NZ	2.25	0.51
40:BG:27:ILE:HD12	40:BG:40:ALA:HA	1.92	0.51
49:BP:71:ARG:HA	49:BP:74:LEU:HB2	1.90	0.51
53:BT:56:MET:O	53:BT:60:GLU:HB2	2.10	0.51
57:BZ:424:LEU:O	57:BZ:428:LEU:HG	2.10	0.51
1:CA:1032:A:H1'	33:C9:23:VAL:HG21	1.92	0.51
2:CB:15:A:H5'	2:CB:16:G:C8	2.45	0.51
3:CC:6:LYS:N	3:CC:9:ARG:NH1	2.58	0.51
4:CD:132:PRO:HD3	4:CD:190:TYR:CZ	2.45	0.51
37:DD:158:ILE:HG22	37:DD:162:LEU:HD12	1.91	0.51
53:DT:9:ASN:O	53:DT:10:LEU:HB2	2.08	0.51
26:A2:35:LEU:HD12	26:A2:53:LEU:HD12	1.91	0.51
1:AA:2803:A:H5''	1:AA:2804:C:H5''	1.92	0.51
3:AC:218:THR:HG22	3:AC:219:MET:SD	2.50	0.51
5:AE:93:VAL:N	63:AE:417:HOH:O	2.35	0.51
13:AP:52:GLU:OE1	13:AP:55:ARG:NH1	2.33	0.51
17:AT:116:ALA:HB1	17:AT:121:ILE:HD11	1.92	0.51
34:BA:1286:A:H2'	34:BA:1287:A:H4'	1.92	0.51
34:BA:316:G:OP2	34:BA:351:G:O2'	2.28	0.51
40:BG:103:TRP:HA	40:BG:106:GLN:HB2	1.92	0.51
24:C0:40:GLN:HE21	24:C0:59:LEU:HG	1.76	0.51
1:CA:1046:A:H3'	1:CA:1047:G:C5'	2.39	0.51
3:CC:64:SER:HA	3:CC:161:ARG:H	1.74	0.51
3:CC:44:VAL:HG21	3:CC:176:VAL:HG21	1.92	0.51
3:CC:50:ILE:HD13	3:CC:50:ILE:H	1.76	0.51
21:CX:47:PHE:O	21:CX:49:VAL:HG13	2.10	0.51
23:CZ:100:VAL:O	23:CZ:124:ILE:N	2.43	0.51
2:CB:106:G:H5'	23:CZ:31:ARG:HG2	1.93	0.51
34:DA:992:U:H3	34:DA:1044:A:H62	1.58	0.51
34:DA:1402:C:H2'	34:DA:1403:C:O4'	2.10	0.51
44:DK:59:TYR:CZ	44:DK:63:LEU:HD21	2.46	0.51
57:DZ:276:VAL:O	57:DZ:280:LEU:HB2	2.09	0.51
1:AA:1386:U:H4'	1:AA:1387:U:OP2	2.10	0.51
1:AA:2146:G:H1	1:AA:2196:C:H42	1.58	0.51
1:AA:905:U:O2	1:AA:2280:A:H2'	2.09	0.51
3:AC:68:GLY:H	3:AC:189:ASN:ND2	2.09	0.51
11:AN:74:ARG:HH11	11:AN:74:ARG:HG3	1.74	0.51
34:BA:406:G:H21	37:BD:119:GLN:NE2	2.02	0.51
57:BZ:181:LEU:HD23	57:BZ:182:ARG:HG3	1.93	0.51
57:BZ:72:CYS:SG	57:BZ:79:ILE:HB	2.51	0.51
57:BZ:88:VAL:HG13	57:BZ:117:GLN:NE2	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:C4:8:LYS:O	28:C4:27:THR:HA	2.10	0.51
1:CA:196:A:H62	13:CP:38:GLN:HE22	1.56	0.51
1:CA:2336:A:H61	24:C0:43:THR:HG22	1.75	0.51
1:CA:68:G:N2	1:CA:74:A:OP2	2.43	0.51
1:CA:918:A:C5	1:CA:919:G:H1'	2.45	0.51
3:CC:29:LEU:O	3:CC:30:VAL:C	2.48	0.51
6:CF:126:VAL:HG21	6:CF:129:PHE:CE1	2.45	0.51
34:DA:1148:U:H2'	34:DA:1149:C:O4'	2.10	0.51
34:DA:1281:U:OP2	34:DA:1282:C:N4	2.38	0.51
34:DA:406:G:H1	34:DA:436:C:H42	1.58	0.51
34:DA:837:G:H1	34:DA:849:C:H42	1.58	0.51
34:DA:401:C:OP2	37:DD:73:ARG:NH2	2.42	0.51
46:DM:33:ALA:HA	46:DM:59:TYR:CE2	2.45	0.51
56:DW:40:C:H2'	56:DW:41:C:C6	2.45	0.51
57:DZ:150:ILE:HA	57:DZ:153:MET:HB3	1.92	0.51
1:AA:1940:A:O2'	1:AA:1942:C:N4	2.43	0.51
1:AA:2802:C:O2'	1:AA:2803:A:O4'	2.28	0.51
13:AP:8:PRO:HB2	13:AP:12:ALA:HB3	1.92	0.51
35:BB:201:ILE:HG21	35:BB:214:ILE:HG21	1.93	0.51
35:BB:54:THR:HG23	35:BB:199:TYR:HB3	1.91	0.51
37:BD:79:PHE:HB2	37:BD:93:PHE:CZ	2.45	0.51
57:BZ:180:VAL:O	57:BZ:213:HIS:HD2	1.93	0.51
57:BZ:-63:ILE:HG12	57:BZ:-49:VAL:HG23	1.93	0.51
1:CA:2016:U:H2'	1:CA:2017:U:H6	1.76	0.51
1:CA:2124:G:O6	1:CA:2174:C:N4	2.38	0.51
1:CA:794:G:H2'	1:CA:795:C:C6	2.46	0.51
8:CH:15:VAL:HG23	8:CH:28:GLY:HA3	1.93	0.51
17:CT:27:THR:HB	17:CT:90:GLN:HB3	1.91	0.51
34:DA:1112:C:O2	36:DC:179:ARG:HG2	2.10	0.51
34:DA:289:G:OP2	63:DA:3239:HOH:O	2.19	0.51
34:DA:391:G:C6	34:DA:392:G:C5	2.99	0.51
34:DA:539:A:OP2	45:DL:115:LYS:NZ	2.42	0.51
39:DF:62:TRP:CD1	51:DR:35:ARG:HD3	2.45	0.51
45:DL:119:LYS:O	45:DL:121:GLY:N	2.44	0.51
57:DZ:201:ILE:HG21	57:DZ:206:LEU:HD13	1.93	0.51
1:AA:2473:C:H2'	1:AA:2474:U:H6	1.75	0.51
1:AA:397:G:H8	1:AA:397:G:OP2	1.94	0.51
1:AA:861:C:O2'	1:AA:862:C:H5'	2.11	0.51
10:AL:125:ARG:O	10:AL:129:GLY:N	2.43	0.51
34:BA:1251:A:H2'	34:BA:1252:A:C8	2.45	0.51
34:BA:131:C:H2'	34:BA:132:C:C6	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:232:G:H2'	34:BA:233:C:H6	1.76	0.51
34:BA:975:A:N6	34:BA:1367:C:O4'	2.44	0.51
47:BN:51:GLY:O	47:BN:53:LEU:N	2.43	0.51
53:BT:97:ALA:N	53:BT:98:PRO:HD3	2.26	0.51
1:CA:1866:C:H2'	1:CA:1876:A:O4'	2.11	0.51
1:CA:20:C:OP1	18:CU:22:LYS:NZ	2.30	0.51
1:CA:2335:A:O2'	1:CA:2336:A:OP2	2.20	0.51
1:CA:804:A:H5''	1:CA:805:G:OP1	2.11	0.51
2:CB:73:A:C4	2:CB:105:A:C2	2.99	0.51
3:CC:42:VAL:CG1	3:CC:43:GLU:N	2.73	0.51
11:CN:14:VAL:HG12	11:CN:15:LEU:N	2.25	0.51
12:CO:31:LYS:HB3	12:CO:32:TYR:CE2	2.46	0.51
13:CP:139:LYS:C	13:CP:141:ALA:H	2.14	0.51
34:DA:1014:A:H4'	52:DS:14:HIS:CE1	2.46	0.51
34:DA:1144:G:N2	34:DA:1146:A:H62	2.08	0.51
34:DA:1513:A:H2'	34:DA:1514:C:C6	2.45	0.51
35:DB:192:SER:O	35:DB:194:PRO:HD3	2.11	0.51
35:DB:88:ALA:HB2	35:DB:219:VAL:HG13	1.93	0.51
48:DO:24:SER:O	48:DO:28:GLN:HG3	2.10	0.51
56:DY:51:U:H3	56:DY:63:G:H1	1.58	0.51
57:DZ:103:GLY:N	57:DZ:130:VAL:HG23	2.24	0.51
28:A4:59:PHE:CD2	52:BS:42:PRO:HB3	2.45	0.51
1:AA:1104:G:N2	1:AA:1127:U:H1'	2.26	0.51
1:AA:142:G:H2'	1:AA:143:C:C6	2.46	0.51
1:AA:2769:U:H1'	1:AA:2770:A:H5''	1.91	0.51
3:AC:44:VAL:HG21	3:AC:176:VAL:HG21	1.92	0.51
3:AC:54:ARG:HD2	3:AC:55:SER:H	1.76	0.51
3:AC:54:ARG:HH22	3:AC:56:ASP:HB3	1.76	0.51
34:BA:1072:G:C5	34:BA:1073:U:C4	2.98	0.51
34:BA:243:A:C2	34:BA:246:A:C8	2.99	0.51
34:BA:401:C:O2'	34:BA:621:A:N3	2.39	0.51
51:BR:58:LEU:HB3	51:BR:62:GLU:HG3	1.91	0.51
52:BS:44:MET:O	52:BS:47:HIS:HB2	2.11	0.51
1:CA:1093:G:H21	1:CA:1098:A:H62	1.58	0.51
1:CA:1110:G:H2'	1:CA:1110:G:N3	2.25	0.51
1:CA:2175:C:H2'	1:CA:2176:A:O4'	2.11	0.51
1:CA:2543:G:H2'	1:CA:2544:G:C8	2.46	0.51
1:CA:556:G:H2'	1:CA:557:U:C6	2.46	0.51
1:CA:846:C:H4'	1:CA:847:U:O5'	2.10	0.51
3:CC:57:GLN:HB2	3:CC:202:PRO:HG2	1.93	0.51
1:CA:673:C:H5''	6:CF:81:PRO:HD2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CH:87:LEU:HD11	8:CH:148:ILE:HB	1.92	0.51
12:CO:7:TYR:HE1	12:CO:20:MET:HE3	1.75	0.51
13:CP:54:GLY:O	63:CP:309:HOH:O	2.19	0.51
21:CX:57:LEU:CD1	21:CX:78:LYS:HB3	2.40	0.51
34:DA:1145:C:H4'	34:DA:1146:A:H5'	1.93	0.51
34:DA:565:U:OP2	34:DA:566:G:O2'	2.13	0.51
43:DJ:11:PHE:CE1	43:DJ:67:THR:HB	2.46	0.51
53:DT:57:ARG:HH12	53:DT:101:GLY:H	1.58	0.51
1:AA:85:C:H4'	1:AA:102:U:H1'	1.93	0.51
1:AA:1321:A:N3	1:AA:1322:A:H1'	2.25	0.51
1:AA:956:A:N1	1:AA:2289:G:H1'	2.25	0.51
1:AA:2901:A:N7	1:AA:2902:G:C6	2.79	0.51
1:AA:417:A:H4'	1:AA:418:G:H5'	1.92	0.51
3:AC:57:GLN:HB2	3:AC:202:PRO:HG2	1.93	0.51
6:AF:7:TYR:O	6:AF:21:ALA:HA	2.10	0.51
8:AH:7:LEU:HD23	8:AH:69:ARG:NH1	2.26	0.51
34:BA:1307:U:H2'	34:BA:1308:U:C6	2.45	0.51
34:BA:28:G:O2'	34:BA:296:U:OP1	2.27	0.51
42:BI:121:ARG:NH1	42:BI:122:ALA:O	2.40	0.51
56:BY:19:G:H4'	56:BY:20:U:OP2	2.09	0.51
57:BZ:289:ILE:HD11	57:BZ:331:TYR:CD1	2.46	0.51
57:BZ:490:PRO:HG3	57:BZ:516:PRO:HD2	1.92	0.51
1:CA:144:C:H5'	21:CX:2:LYS:HE2	1.92	0.51
1:CA:2025:C:H2'	1:CA:2026:C:H6	1.71	0.51
1:CA:2537:U:H2'	1:CA:2538:C:C6	2.46	0.51
1:CA:271(O):C:H2'	1:CA:271(P):C:C6	2.46	0.51
1:CA:958:U:O2	2:CB:90:A:O2'	2.23	0.51
1:CA:2124:G:C4'	3:CC:175:PRO:CG	2.82	0.51
3:CC:65:LEU:HB3	3:CC:189:ASN:HD22	1.75	0.51
5:CE:12:THR:HG21	17:CT:11:GLU:OE2	2.11	0.51
5:CE:169:ASN:HD22	5:CE:203:LYS:HB2	1.76	0.51
6:CF:184:TYR:CE2	6:CF:188:ARG:HD2	2.45	0.51
22:CY:11:ASP:OD2	22:CY:97:ARG:NH2	2.43	0.51
34:DA:436:C:H2'	34:DA:437:U:H6	1.75	0.51
36:DC:54:ARG:HB3	36:DC:54:ARG:HH11	1.76	0.51
44:DK:79:SER:HB2	44:DK:106:LYS:HE3	1.92	0.51
1:AA:2559:U:H2'	1:AA:2560:G:H8	1.76	0.51
1:AA:2556:G:H1'	1:AA:2658:C:H4'	1.93	0.51
1:AA:552:C:O2	1:AA:552:C:O4'	2.27	0.51
1:AA:864:C:H4'	1:AA:977:G:C5	2.46	0.51
3:AC:65:LEU:HB3	3:AC:189:ASN:HD22	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:AH:98:LEU:HD12	8:AH:102:ALA:O	2.11	0.51
1:AA:1186:U:H6	11:AN:63:THR:HG1	1.57	0.51
1:AA:956:A:H62	14:AQ:12:GLN:HA	1.74	0.51
34:BA:153:C:H42	34:BA:169:C:H42	1.58	0.51
34:BA:441:A:H8	34:BA:441:A:OP2	1.94	0.51
34:BA:771:G:H2'	34:BA:772:U:C6	2.46	0.51
38:BE:78:HIS:NE2	38:BE:142:LEU:HA	2.26	0.51
3:CC:218:THR:HG22	3:CC:219:MET:SD	2.50	0.51
6:CF:108:LYS:O	6:CF:112:MET:HG3	2.11	0.51
7:CG:49:ASP:C	7:CG:51:ARG:H	2.13	0.51
13:CP:97:PRO:HD3	13:CP:126:VAL:O	2.10	0.51
1:CA:2469:A:O3'	14:CQ:56:ARG:NH2	2.43	0.51
17:CT:115:ARG:H	17:CT:115:ARG:HD2	1.75	0.51
23:CZ:29:TYR:HB3	23:CZ:34:ASN:ND2	2.25	0.51
23:CZ:39:VAL:HG21	23:CZ:44:PHE:HB2	1.92	0.51
34:DA:1029:C:H2'	34:DA:1030:C:H5''	1.92	0.51
34:DA:1372:U:H5''	42:DI:71:SER:HB3	1.93	0.51
37:DD:196:LEU:O	37:DD:198:VAL:N	2.41	0.51
57:DZ:12:LEU:O	57:DZ:283:PRO:HD3	2.10	0.51
57:DZ:20:HIS:O	57:DZ:25:LYS:NZ	2.29	0.51
57:DZ:138:LYS:HA	62:DZ:704:GDP:N1	2.25	0.51
1:AA:1704:C:H4'	5:AE:133:LYS:HB3	1.92	0.51
1:AA:2304:C:P	16:AS:17:ARG:HH12	2.34	0.51
1:AA:7:G:H2'	1:AA:8:A:C8	2.45	0.51
4:AD:221:VAL:HG22	4:AD:226:MET:CE	2.41	0.51
1:AA:515:G:N7	20:AW:49:LYS:NZ	2.57	0.51
23:AZ:48:PHE:CE2	23:AZ:52:SER:HA	2.46	0.51
34:BA:551:U:H2'	34:BA:552:U:C6	2.46	0.51
36:BC:134:ILE:O	36:BC:138:VAL:HG23	2.11	0.51
37:BD:13:ARG:HB3	37:BD:13:ARG:HH11	1.76	0.51
38:BE:88:LYS:HE2	38:BE:123:LEU:HD12	1.92	0.51
46:BM:15:VAL:HA	46:BM:18:ALA:HB3	1.92	0.51
52:BS:16:LEU:HD12	52:BS:19:VAL:HB	1.92	0.51
57:BZ:240:GLU:H	57:BZ:240:GLU:CD	2.14	0.51
57:BZ:639:ASN:N	57:BZ:640:ALA:HB3	2.26	0.51
57:BZ:647:VAL:HG21	57:BZ:652:MET:SD	2.51	0.51
26:C2:28:LYS:HE3	26:C2:56:GLN:OE1	2.11	0.51
1:CA:1049:C:H3'	1:CA:1050:A:C8	2.46	0.51
1:CA:2112:G:N7	1:CA:2113:U:H1'	2.26	0.51
1:CA:2854:G:H2'	1:CA:2855:C:C6	2.45	0.51
1:CA:1861:G:OP1	3:CC:205:ALA:C	2.48	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:CR:37:THR:OG1	15:CR:40:LYS:HB2	2.10	0.51
18:CU:65:ILE:HD11	18:CU:95:LEU:HB3	1.92	0.51
34:DA:1003:G:H2'	34:DA:1004:A:O4'	2.11	0.51
34:DA:1003:G:N2	34:DA:1025:U:O4	2.43	0.51
34:DA:428:G:C5	34:DA:430:A:C6	2.99	0.51
34:DA:460:G:O6	34:DA:470:C:H5''	2.11	0.51
34:DA:503:C:OP2	45:DL:116:SER:HB3	2.10	0.51
57:DZ:548:GLU:O	57:DZ:551:GLN:HG2	2.09	0.51
57:DZ:556:ILE:HG13	57:DZ:558:PHE:HD2	1.75	0.51
57:DZ:409:ILE:HD11	57:DZ:654:GLY:HA2	1.92	0.51
1:AA:2343:G:H4'	24:A0:43:THR:H	1.76	0.51
1:AA:843:C:H2'	1:AA:844:C:C6	2.45	0.51
1:AA:990:A:C4	1:AA:2460:A:C2	2.99	0.51
9:AK:73:GLY:C	9:AK:75:GLN:H	2.11	0.51
10:AL:30:HIS:CG	10:AL:59:ILE:HB	2.46	0.51
18:AU:90:VAL:HB	18:AU:95:LEU:HD13	1.92	0.51
40:BG:111:ARG:NH1	40:BG:113:GLU:OE2	2.43	0.51
56:BY:58:A:C2	56:BY:60:U:H2'	2.46	0.51
57:BZ:114:VAL:CG2	57:BZ:152:THR:HB	2.36	0.51
1:CA:2395:C:O2'	25:C1:30:VAL:HG22	2.11	0.51
1:CA:1041:C:H42	1:CA:1114:G:H1	1.57	0.51
1:CA:2349:G:H3'	1:CA:2350:C:H5''	1.93	0.51
1:CA:2850:A:H2'	1:CA:2851:A:H8	1.76	0.51
1:CA:580:C:H2'	1:CA:581:C:C6	2.46	0.51
3:CC:68:GLY:H	3:CC:189:ASN:ND2	2.09	0.51
7:CG:37:VAL:O	7:CG:94:LEU:N	2.35	0.51
11:CN:102:ALA:O	11:CN:106:MET:HG3	2.11	0.51
12:CO:18:LYS:HB2	12:CO:45:GLU:HB3	1.92	0.51
1:CA:994:C:H1'	19:CV:10:LYS:HE3	1.92	0.51
34:DA:1016:A:O2'	34:DA:1217:C:O2'	2.18	0.51
34:DA:1040:U:N3	34:DA:1041:A:N7	2.58	0.51
34:DA:1118:C:H1'	34:DA:1179:A:C4	2.46	0.51
34:DA:502:G:C6	34:DA:503:C:N3	2.79	0.51
35:DB:163:PHE:CD1	35:DB:185:ILE:HG13	2.41	0.51
41:DH:11:THR:O	41:DH:15:ASN:ND2	2.39	0.51
57:DZ:415:PRO:HB3	57:DZ:424:LEU:HD23	1.92	0.51
24:A0:70:GLN:OE1	24:A0:80:HIS:NE2	2.37	0.50
25:A1:53:VAL:HG21	25:A1:94:LEU:HD11	1.93	0.50
16:AS:10:ARG:HG2	16:AS:91:PRO:HA	1.93	0.50
34:BA:345:C:H4'	34:BA:346:G:C4	2.47	0.50
34:BA:390:C:H2'	34:BA:391:G:C8	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:448:A:C4	34:BA:487:A:C2	2.99	0.50
34:BA:626:U:C2	34:BA:627:G:C8	2.99	0.50
35:BB:105:PHE:O	35:BB:107:THR:N	2.45	0.50
41:BH:10:LEU:HD11	41:BH:85:ARG:HG2	1.93	0.50
1:CA:2788:C:N4	1:CA:2789:C:H41	2.09	0.50
1:CA:2823:A:OP1	5:CE:159:HIS:NE2	2.32	0.50
1:CA:247:G:H4'	1:CA:386:G:C5	2.46	0.50
34:DA:1103:C:H2'	34:DA:1104:G:O4'	2.11	0.50
34:DA:137:C:H2'	34:DA:138:G:H8	1.76	0.50
34:DA:475:G:H2'	34:DA:476:G:H8	1.76	0.50
34:DA:987:G:N2	34:DA:1219:U:O2	2.44	0.50
40:DG:118:VAL:HG13	40:DG:122:HIS:NE2	2.26	0.50
41:DH:6:ILE:O	41:DH:10:LEU:HG	2.11	0.50
1:AA:1128:U:H5''	1:AA:1129:U:OP2	2.12	0.50
1:AA:552:C:H4'	1:AA:553:A:O5'	2.11	0.50
3:AC:6:LYS:N	3:AC:9:ARG:NH1	2.58	0.50
34:BA:1234:C:H1'	34:BA:1364:U:O2	2.11	0.50
38:BE:40:ARG:NH2	38:BE:68:GLU:HA	2.26	0.50
34:BA:1381:U:O4'	40:BG:79:ARG:NE	2.43	0.50
43:BJ:47:PHE:HB2	43:BJ:63:PHE:HB2	1.93	0.50
34:BA:376:G:O3'	49:BP:5:ARG:HD2	2.10	0.50
56:BY:5:G:H1	56:BY:68:C:H42	1.59	0.50
57:BZ:-6:ARG:C	57:BZ:-4:ALA:H	2.14	0.50
1:CA:2207:G:H3'	1:CA:2208:A:H5''	1.93	0.50
1:CA:2420:C:P	32:C8:33:ASN:H	2.34	0.50
5:CE:49:LEU:HD22	5:CE:81:ILE:HG13	1.93	0.50
13:CP:86:LYS:HD3	13:CP:117:GLU:HB3	1.93	0.50
38:DE:90:VAL:HG23	38:DE:121:LYS:HB3	1.92	0.50
39:DF:36:ARG:NH1	39:DF:38:GLU:OE2	2.44	0.50
52:DS:27:GLU:HG2	52:DS:47:HIS:CD2	2.46	0.50
1:AA:1160:G:H2'	1:AA:1161:G:C8	2.46	0.50
1:AA:1993:A:OP2	4:AD:242:ARG:NH2	2.44	0.50
9:AK:26:LEU:HA	9:AK:84:GLU:HA	1.94	0.50
9:AK:4:LYS:N	9:AK:5:ARG:HA	2.27	0.50
13:AP:113:LYS:HG3	13:AP:129:ALA:HB3	1.93	0.50
22:AY:86:ARG:HB2	22:AY:98:VAL:HG23	1.93	0.50
34:BA:27:G:H2'	34:BA:28:G:C8	2.46	0.50
34:BA:302:G:N3	34:BA:556:C:H4'	2.27	0.50
41:BH:39:LEU:HB3	41:BH:45:ILE:HG12	1.93	0.50
34:BA:1525:G:OP1	44:BK:120:ARG:NH2	2.44	0.50
57:BZ:616:TYR:HE2	57:BZ:664:GLN:HG3	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1274:A:N3	1:CA:1297:C:H1'	2.26	0.50
1:CA:1359:A:C2	1:CA:1372:U:O4	2.65	0.50
1:CA:1745(A):C:H5'	1:CA:1746:G:OP2	2.11	0.50
1:CA:2112:G:C5	1:CA:2113:U:H1'	2.47	0.50
3:CC:42:VAL:HG13	3:CC:43:GLU:H	1.73	0.50
1:CA:1971:A:C4	4:CD:241:PRO:HD3	2.46	0.50
6:CF:132:VAL:HG21	6:CF:163:VAL:HG22	1.92	0.50
7:CG:46:ALA:O	7:CG:51:ARG:HA	2.10	0.50
8:CH:90:LYS:HD3	8:CH:159:GLU:HG2	1.94	0.50
16:CS:15:ARG:O	16:CS:19:LYS:HG2	2.12	0.50
34:DA:1256:A:H61	34:DA:1278:U:H1'	1.76	0.50
34:DA:1320:C:C2	52:DS:72:GLY:HA3	2.46	0.50
34:DA:792:A:H4'	34:DA:793:U:O5'	2.12	0.50
46:DM:124:PRO:HD2	57:DZ:507:TYR:HB2	1.92	0.50
26:A2:46:GLN:HB2	26:A2:49:LYS:HD2	1.94	0.50
1:AA:1221:G:H1'	1:AA:1222:A:H5'	1.94	0.50
1:AA:1537:G:C5	1:AA:1546:G:N2	2.80	0.50
1:AA:1636:U:O2'	1:AA:1637:G:H5'	2.12	0.50
1:AA:2221:A:H3'	1:AA:2222:C:H6	1.77	0.50
1:AA:2896:G:O2'	29:A5:32:PRO:HD2	2.11	0.50
1:AA:441:C:H2'	1:AA:442:A:C8	2.47	0.50
3:AC:50:ILE:HD13	3:AC:50:ILE:H	1.76	0.50
14:AQ:10:ARG:HB2	14:AQ:10:ARG:CZ	2.41	0.50
34:BA:1148:U:H2'	34:BA:1149:C:O4'	2.11	0.50
44:BK:99:GLN:HG2	44:BK:105:VAL:HG21	1.92	0.50
45:BL:32:PHE:HB3	45:BL:84:LEU:HD11	1.92	0.50
45:BL:84:LEU:HD23	45:BL:105:TYR:HE2	1.76	0.50
55:BV:15:A:C2'	55:BV:16:U:H5	2.24	0.50
25:C1:95:LEU:O	25:C1:98:LEU:HB2	2.12	0.50
31:C7:24:THR:O	31:C7:28:ARG:HG3	2.11	0.50
1:CA:1472:A:N6	1:CA:1519:G:H1'	2.27	0.50
1:CA:754:C:H2'	1:CA:755:C:H6	1.76	0.50
11:CN:24:GLY:O	11:CN:28:THR:HG23	2.12	0.50
34:DA:447:G:H2'	34:DA:485:G:N2	2.25	0.50
34:DA:665:A:H2'	34:DA:732:C:O2	2.11	0.50
34:DA:707:C:H2'	34:DA:708:C:H6	1.76	0.50
34:DA:836:G:OP1	51:DR:61:LYS:NZ	2.43	0.50
34:DA:90:U:H2'	34:DA:91:C:C6	2.46	0.50
34:DA:986:A:O2'	52:DS:55:LYS:O	2.30	0.50
34:DA:1316:G:N7	52:DS:7:LYS:NZ	2.60	0.50
57:DZ:221:ALA:HB1	57:DZ:228:MET:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:A1:34:THR:HG22	25:A1:36:GLY:H	1.75	0.50
30:A6:50:ARG:HG3	30:A6:51:GLU:N	2.26	0.50
1:AA:1106:U:H4'	1:AA:1107:U:H5'	1.93	0.50
1:AA:2155:G:O2'	1:AA:2178:G:N2	2.44	0.50
1:AA:1834:A:O2'	4:AD:259:THR:HG21	2.11	0.50
10:AL:55:VAL:HG13	10:AL:57:ILE:HD11	1.93	0.50
11:AN:5:VAL:HG23	11:AN:6:PRO:HD2	1.94	0.50
23:AZ:70:LEU:HD11	23:AZ:98:MET:CE	2.42	0.50
34:BA:1466:C:H2'	34:BA:1467:G:O4'	2.11	0.50
34:BA:389:A:N6	34:BA:390:C:O2	2.44	0.50
35:BB:88:ALA:HB2	35:BB:219:VAL:HG13	1.93	0.50
37:BD:42:GLN:O	37:BD:42:GLN:HG3	2.11	0.50
40:BG:26:PHE:O	40:BG:30:ILE:HG13	2.11	0.50
45:BL:88:GLY:O	45:BL:99:HIS:CD2	2.64	0.50
1:CA:1364:G:C8	25:C1:3:LYS:HE2	2.47	0.50
26:C2:16:LEU:HD13	26:C2:20:GLU:HB3	1.92	0.50
1:CA:2394:C:P	32:C8:30:ARG:HH11	2.34	0.50
1:CA:1059:G:OP2	1:CA:1060:U:H3'	2.11	0.50
1:CA:2636:U:H1'	1:CA:2783:G:N2	2.26	0.50
1:CA:487:C:H1'	20:CW:53:SER:HA	1.93	0.50
1:CA:503:A:H4'	1:CA:504:U:H5''	1.93	0.50
1:CA:614(B):G:H2'	6:CF:44:ARG:NH1	2.26	0.50
1:CA:2177:C:O3'	3:CC:47:LYS:HB2	2.10	0.50
34:DA:976:G:N2	34:DA:1362:C:H2'	2.27	0.50
34:DA:144:G:H2'	34:DA:145:G:H8	1.76	0.50
34:DA:189(B):C:H2'	34:DA:189(C):C:C6	2.47	0.50
34:DA:685:G:N1	34:DA:686:U:O4	2.44	0.50
35:DB:150:SER:O	35:DB:153:ARG:HG2	2.11	0.50
38:DE:101:ILE:HG13	38:DE:119:LEU:HD23	1.93	0.50
56:DY:25:C:H2'	56:DY:26:A:H8	1.77	0.50
1:AA:1410:G:OP2	25:A1:3:LYS:HD2	2.12	0.50
28:A4:47:GLN:HG2	28:A4:49:PHE:H	1.76	0.50
1:AA:1739:U:O2'	1:AA:1740:U:H2'	2.11	0.50
1:AA:1827:U:H2'	1:AA:1828:C:C6	2.46	0.50
1:AA:641:G:OP2	6:AF:43:LYS:NZ	2.44	0.50
1:AA:997:G:OP1	14:AQ:16:ARG:NH2	2.45	0.50
4:AD:112:GLN:O	4:AD:115:GLN:HB3	2.11	0.50
4:AD:89:SER:HB2	4:AD:159:ALA:H	1.76	0.50
10:AL:88:ALA:O	10:AL:90:LYS:N	2.45	0.50
12:AO:16:ALA:HB2	12:AO:52:VAL:HG21	1.94	0.50
18:AU:102:GLU:HA	18:AU:104:GLN:HE22	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:1030(D):A:H62	34:BA:1031:G:H21	1.58	0.50
34:BA:142:G:H2'	34:BA:143:A:H8	1.77	0.50
38:BE:6:PHE:CE2	38:BE:36:ASP:HB3	2.47	0.50
38:BE:41:VAL:O	38:BE:67:VAL:N	2.44	0.50
40:BG:93:PRO:HA	40:BG:96:GLN:HB2	1.94	0.50
41:BH:40:ALA:O	41:BH:42:GLU:N	2.45	0.50
39:BF:96:PRO:HB3	51:BR:30:ASP:OD2	2.12	0.50
57:BZ:225:GLU:HA	57:BZ:228:MET:CB	2.38	0.50
1:CA:265:A:C8	1:CA:266:G:H1'	2.47	0.50
1:CA:912:C:H2'	1:CA:913:U:C6	2.46	0.50
3:CC:191:ARG:O	3:CC:195:ARG:HG2	2.11	0.50
5:CE:96:PHE:O	5:CE:175:VAL:HG11	2.11	0.50
34:DA:377:G:OP1	49:DP:3:LYS:HD2	2.12	0.50
45:DL:69:TYR:HB3	45:DL:99:HIS:ND1	2.27	0.50
47:DN:40:CYS:O	47:DN:43:CYS:N	2.45	0.50
49:DP:6:LEU:HD23	49:DP:17:TYR:CD1	2.46	0.50
56:DY:12:U:H3	56:DY:23:A:H61	1.59	0.50
57:DZ:35:TYR:OH	57:DZ:266:ASN:HB3	2.11	0.50
1:AA:1772:C:H6	1:AA:1772:C:O5'	1.94	0.50
1:AA:2119:C:H2'	1:AA:2120:U:O4'	2.11	0.50
3:AC:191:ARG:O	3:AC:194:ILE:HG22	2.12	0.50
3:AC:42:VAL:CG1	3:AC:43:GLU:N	2.73	0.50
4:AD:221:VAL:HG22	4:AD:226:MET:HE3	1.94	0.50
7:AG:61:ALA:O	28:A4:7:PRO:HG2	2.12	0.50
9:AK:69:PRO:C	9:AK:71:LEU:H	2.15	0.50
43:BJ:13:HIS:HA	43:BJ:16:LEU:HB3	1.92	0.50
43:BJ:16:LEU:HD21	43:BJ:70:ARG:HG2	1.94	0.50
49:BP:43:LYS:HG2	49:BP:48:TRP:CE2	2.47	0.50
57:BZ:-38:TYR:O	57:BZ:-35:PRO:HD2	2.12	0.50
1:CA:2360:A:H2'	1:CA:2361:A:O4'	2.11	0.50
1:CA:469:G:C2'	1:CA:470:A:H5''	2.42	0.50
2:CB:78:A:C2	2:CB:100:A:C4	3.00	0.50
4:CD:228:PRO:HD3	4:CD:235:GLY:HA3	1.93	0.50
4:CD:89:SER:HB2	4:CD:159:ALA:HB2	1.94	0.50
7:CG:137:GLU:HB3	7:CG:140:ILE:HD13	1.94	0.50
9:CK:74:LEU:O	9:CK:76:GLY:N	2.45	0.50
12:CO:26:LYS:O	12:CO:30:ALA:HB2	2.11	0.50
20:CW:60:ASN:N	20:CW:60:ASN:HD22	2.09	0.50
34:DA:1240:U:H5'	34:DA:1241:G:C8	2.47	0.50
34:DA:509:A:H8	34:DA:509:A:H3'	1.77	0.50
34:DA:520:A:N1	34:DA:536:C:H1'	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
41:DH:44:PHE:HB3	41:DH:80:ILE:CD1	2.41	0.50
34:DA:1268:A:O2'	54:DU:19:GLY:HA2	2.12	0.50
57:DZ:132:ARG:HD3	57:DZ:160:ARG:NH1	2.27	0.50
57:DZ:498:ILE:HG22	57:DZ:507:TYR:CE2	2.47	0.50
29:A5:35:GLU:HG3	29:A5:51:TYR:CD2	2.46	0.50
1:AA:1157:A:H8	1:AA:1158:G:H1'	1.75	0.50
1:AA:2221:A:H5''	1:AA:2222:C:OP2	2.12	0.50
1:AA:2317:A:N1	7:AG:154:GLY:N	2.54	0.50
1:AA:2737:C:OP1	5:AE:118:LYS:NZ	2.45	0.50
17:AT:16:ARG:HH12	17:AT:19:LEU:HD21	1.76	0.50
34:BA:1320:C:H2'	34:BA:1321:C:O4'	2.12	0.50
34:BA:437:U:H2'	34:BA:438:G:H5'	1.92	0.50
34:BA:937:A:OP2	63:BA:5214:HOH:O	2.18	0.50
37:BD:64:LEU:HD11	37:BD:97:LEU:HD13	1.92	0.50
46:BM:24:GLY:O	46:BM:29:ARG:NH1	2.43	0.50
56:BY:26:A:H61	56:BY:44:G:H1	1.59	0.50
30:C6:10:LEU:HD23	30:C6:22:ALA:HB2	1.94	0.50
1:CA:1680:U:O2	1:CA:1763:G:H3'	2.11	0.50
1:CA:1860:G:OP1	3:CC:207:GLY:CA	2.60	0.50
1:CA:2203:U:H2'	1:CA:2205:C:H6	1.77	0.50
1:CA:2729:G:C6	1:CA:2730:C:C4	3.00	0.50
1:CA:34:C:O2	1:CA:34:C:H2'	2.12	0.50
3:CC:20:VAL:O	3:CC:21:TYR:CB	2.58	0.50
11:CN:38:HIS:NE2	11:CN:50:ASP:OD2	2.38	0.50
34:DA:1123:A:H4'	43:DJ:36:GLY:HA3	1.94	0.50
34:DA:418:C:H2'	34:DA:419:C:C6	2.47	0.50
46:DM:17:VAL:O	46:DM:20:THR:OG1	2.22	0.50
57:DZ:412:ALA:HB2	57:DZ:479:PRO:HB3	1.92	0.50
1:AA:2144:U:H1'	3:AC:167:ASP:HB2	1.92	0.50
1:AA:326:C:H2'	1:AA:327:U:H6	1.77	0.50
1:AA:779:C:H2'	1:AA:780:G:O4'	2.12	0.50
1:AA:860:U:H2'	1:AA:861:C:C6	2.47	0.50
1:AA:1232:G:H5''	19:AV:81:TYR:CE1	2.47	0.50
34:BA:1106:G:C6	34:BA:1107:C:C4	2.99	0.50
38:BE:146:ALA:O	38:BE:149:GLU:N	2.45	0.50
44:BK:84:VAL:HG11	44:BK:91:ARG:HH11	1.77	0.50
48:BO:3:ILE:HG21	48:BO:34:LEU:HD21	1.93	0.50
57:BZ:247:ARG:O	57:BZ:251:ILE:HG13	2.12	0.50
1:CA:1268:A:C2	1:CA:2013:A:C4	3.00	0.50
23:CZ:139:VAL:HG23	23:CZ:141:VAL:HG13	1.94	0.50
34:DA:1132:C:H2'	34:DA:1133:G:C8	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
29:A5:48:GLU:O	29:A5:60:VAL:HG11	2.12	0.49
32:A8:34:TRP:CG	32:A8:35:GLN:N	2.80	0.49
1:AA:2219:U:C6	1:AA:2236:G:C6	2.99	0.49
1:AA:2287:C:H6	1:AA:2287:C:H5'	1.77	0.49
3:AC:183:PRO:C	3:AC:185:LYS:H	2.16	0.49
4:AD:181:GLU:OE2	4:AD:270:ILE:HG12	2.11	0.49
1:AA:642:G:OP2	6:AF:106:ARG:NH2	2.45	0.49
12:AO:71:ARG:NH2	12:AO:105:GLU:OE1	2.44	0.49
2:AB:48:A:H4'	16:AS:95:HIS:HD2	1.77	0.49
34:BA:107:G:H2'	34:BA:108:G:O4'	2.12	0.49
34:BA:520:A:N1	34:BA:536:C:H1'	2.26	0.49
34:BA:622:A:N7	34:BA:623:C:C2	2.80	0.49
41:BH:20:TYR:HA	41:BH:65:TYR:OH	2.11	0.49
57:BZ:180:VAL:O	57:BZ:213:HIS:CD2	2.65	0.49
57:BZ:612:THR:OG1	57:BZ:616:TYR:HB2	2.12	0.49
1:CA:1340:U:OP1	21:CX:16:LYS:NZ	2.43	0.49
1:CA:860:U:C2	1:CA:2268:A:C8	3.00	0.49
1:CA:272(C):G:H2'	1:CA:272(D):G:O4'	2.12	0.49
6:CF:29:ASN:H	6:CF:112:MET:HE2	1.76	0.49
16:CS:48:LEU:HD23	16:CS:82:ILE:HD11	1.93	0.49
1:CA:489:G:N7	20:CW:49:LYS:NZ	2.60	0.49
34:DA:474:G:H2'	34:DA:475:G:H8	1.76	0.49
34:DA:561:U:O2'	34:DA:562:C:OP2	2.29	0.49
36:DC:179:ARG:NH1	36:DC:206:GLU:OE1	2.44	0.49
38:DE:92:LYS:HB3	38:DE:119:LEU:HB2	1.94	0.49
39:DF:87:ARG:NH1	39:DF:87:ARG:HG3	2.13	0.49
44:DK:65:ALA:HB1	44:DK:98:LEU:HD23	1.94	0.49
34:DA:1325:C:H4'	54:DU:17:THR:HG21	1.93	0.49
1:AA:2398:C:H2'	1:AA:2399:U:C6	2.47	0.49
7:AG:68:PRO:HG2	7:AG:90:LEU:HD22	1.93	0.49
1:AA:612:C:P	13:AP:16:ARG:HH22	2.34	0.49
1:AA:26:G:OP1	20:AW:80:PRO:HB3	2.12	0.49
1:AA:142:G:H1'	21:AX:37:THR:CG2	2.42	0.49
34:BA:7:G:H5'	34:BA:298:A:O4'	2.11	0.49
34:BA:453:A:O2'	49:BP:68:ASP:O	2.31	0.49
37:BD:196:LEU:O	37:BD:198:VAL:N	2.39	0.49
38:BE:28:PHE:CD1	38:BE:28:PHE:N	2.80	0.49
48:BO:88:ARG:HB3	48:BO:88:ARG:CZ	2.41	0.49
53:BT:16:HIS:O	53:BT:19:SER:OG	2.27	0.49
57:BZ:674:ASP:HB3	57:BZ:675:HIS:HD2	1.77	0.49
1:CA:1721:G:H8	1:CA:1741:A:H62	1.59	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:236:C:H2'	1:CA:237:C:C6	2.47	0.49
1:CA:58:G:O2'	1:CA:73:A:N1	2.43	0.49
3:CC:54:ARG:HD2	3:CC:55:SER:H	1.76	0.49
1:CA:2823:A:P	5:CE:159:HIS:HE2	2.35	0.49
34:DA:1273:G:H3'	34:DA:1274:G:C8	2.47	0.49
34:DA:1443:G:N2	34:DA:1460:A:H1'	2.26	0.49
34:DA:1399:C:C2	34:DA:1502:A:N6	2.80	0.49
34:DA:509:A:H3'	34:DA:509:A:C8	2.48	0.49
35:DB:174:VAL:O	35:DB:178:ARG:HB2	2.12	0.49
37:DD:72:GLU:O	37:DD:76:ARG:HB3	2.12	0.49
34:DA:1374:A:H4'	40:DG:28:ASN:ND2	2.27	0.49
38:DE:93:PRO:O	41:DH:105:ARG:NH2	2.45	0.49
41:DH:36:LEU:HA	41:DH:39:LEU:HD23	1.95	0.49
13:AP:63:PRO:HD3	32:A8:27:THR:HG22	1.93	0.49
1:AA:553:A:OP2	11:AN:114:ARG:NH1	2.45	0.49
1:AA:801:C:H2'	1:AA:802:C:H6	1.77	0.49
14:AQ:11:LYS:HE2	14:AQ:88:GLY:O	2.12	0.49
18:AU:81:HIS:CE1	18:AU:85:LYS:HD2	2.47	0.49
23:AZ:156:LYS:O	23:AZ:157:LEU:HB2	2.10	0.49
34:BA:1318:A:O2'	52:BS:37:ARG:HB3	2.13	0.49
17:AT:41:ARG:NH2	34:BA:346:G:OP1	2.43	0.49
35:BB:170:GLU:O	35:BB:174:VAL:HG23	2.12	0.49
36:BC:43:LEU:HB3	36:BC:47:LEU:HD12	1.93	0.49
38:BE:121:LYS:HG3	38:BE:122:GLU:N	2.26	0.49
38:BE:139:LEU:O	38:BE:141:GLN:N	2.46	0.49
38:BE:40:ARG:HB3	38:BE:66:MET:CE	2.42	0.49
56:BY:7:A:O2'	56:BY:49:C:H5'	2.13	0.49
1:CA:1533:G:H2'	1:CA:1536:C:H42	1.76	0.49
1:CA:1956:U:C2'	1:CA:1957:C:H5'	2.41	0.49
1:CA:300:A:H2'	1:CA:334:C:H1'	1.95	0.49
1:CA:866:A:C6	1:CA:914:C:C5	3.00	0.49
6:CF:57:VAL:HG13	6:CF:59:TYR:CD2	2.47	0.49
8:CH:163:TYR:CE2	8:CH:169:VAL:HG22	2.47	0.49
9:CK:49:ALA:N	9:CK:90:ALA:HB1	2.26	0.49
18:CU:28:ARG:HD3	18:CU:38:THR:OG1	2.11	0.49
21:CX:11:PRO:HB3	21:CX:92:LEU:HD21	1.93	0.49
34:DA:130:A:C8	50:DQ:63:ARG:HG3	2.46	0.49
34:DA:201:C:H42	34:DA:216:G:H1	1.60	0.49
34:DA:536:C:H6	34:DA:536:C:O5'	1.96	0.49
34:DA:620:C:H2'	34:DA:621:A:O4'	2.13	0.49
34:DA:866:C:C4	34:DA:867:G:H1'	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:982:U:O2	34:DA:1222:G:N1	2.37	0.49
37:DD:19:LEU:O	37:DD:21:LEU:N	2.45	0.49
37:DD:38:TYR:CE1	37:DD:45:GLN:HG3	2.47	0.49
38:DE:43:LEU:HD12	38:DE:44:GLY:N	2.28	0.49
40:DG:32:ARG:O	40:DG:34:GLY:N	2.45	0.49
43:DJ:50:ILE:HA	43:DJ:60:ARG:HD3	1.95	0.49
57:DZ:170:ARG:N	57:DZ:170:ARG:NH1	2.57	0.49
61:DZ:703:FUA:O1	61:DZ:703:FUA:C1	2.60	0.49
28:A4:24:THR:OG1	28:A4:25:TYR:N	2.46	0.49
1:AA:116:A:H3'	1:AA:117:A:H5''	1.95	0.49
1:AA:1402:G:H2'	1:AA:1403:U:C6	2.48	0.49
1:AA:1785:C:N3	1:AA:2729:U:O2'	2.44	0.49
2:AB:73:A:C4	2:AB:105:A:C2	3.01	0.49
4:AD:102:LYS:C	4:AD:103:ARG:HG2	2.28	0.49
23:AZ:184:ALA:O	23:AZ:185:GLU:HB3	2.13	0.49
23:AZ:30:ASN:HD22	23:AZ:90:VAL:HB	1.77	0.49
34:BA:931:C:H42	34:BA:1386:G:H1	1.61	0.49
34:BA:192:U:H2'	34:BA:193:C:H6	1.77	0.49
36:BC:195:VAL:O	36:BC:196:LEU:HD13	2.12	0.49
36:BC:43:LEU:HD22	36:BC:47:LEU:HD11	1.93	0.49
40:BG:61:VAL:O	40:BG:65:ALA:N	2.45	0.49
51:BR:44:LEU:HD21	51:BR:70:ILE:HG21	1.94	0.49
57:BZ:166:LEU:HD21	57:BZ:208:GLN:HG2	1.95	0.49
1:CA:2577:A:O4'	29:C5:3:LYS:HB2	2.13	0.49
1:CA:1652:A:OP1	15:CR:8:ARG:NH1	2.45	0.49
1:CA:1688:U:H1'	1:CA:1701:A:C6	2.48	0.49
1:CA:2512:C:H2'	1:CA:2513:G:O4'	2.12	0.49
1:CA:2557:G:H2'	1:CA:2558:C:C6	2.48	0.49
1:CA:828:U:H4'	1:CA:831:G:N1	2.27	0.49
7:CG:124:SER:HB2	7:CG:131:TYR:CE1	2.46	0.49
12:CO:23:ARG:HG3	12:CO:24:VAL:N	2.27	0.49
34:DA:1063:C:OP2	34:DA:1064:G:O2'	2.15	0.49
34:DA:188:C:H2'	34:DA:189:G:C8	2.44	0.49
34:DA:538:G:H5''	45:DL:114:LYS:HB2	1.93	0.49
35:DB:51:LEU:O	35:DB:55:PHE:N	2.35	0.49
36:DC:186:PHE:CE2	36:DC:188:LEU:HB2	2.47	0.49
38:DE:43:LEU:HD21	38:DE:132:ALA:HB1	1.94	0.49
34:DA:1347:G:H8	42:DI:107:ARG:HB3	1.77	0.49
1:AA:2697:G:H5'	12:AO:68:GLU:OE1	2.12	0.49
1:AA:868:A:C2'	1:AA:991:G:H5''	2.42	0.49
3:AC:191:ARG:O	3:AC:195:ARG:HG2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:20:ASP:N	4:AD:20:ASP:OD1	2.45	0.49
37:BD:188:LEU:H	37:BD:188:LEU:HD23	1.77	0.49
40:BG:38:LEU:HA	40:BG:41:ARG:HB2	1.92	0.49
45:BL:27:LEU:HD13	45:BL:98:TYR:CE2	2.47	0.49
57:BZ:264:LEU:HD12	62:BZ:704:GDP:N3	2.27	0.49
57:BZ:467:LYS:O	57:BZ:469:GLU:N	2.38	0.49
1:CA:1063:G:H4'	10:CL:134:MET:HB3	1.94	0.49
1:CA:1607:C:H4'	1:CA:1608:A:O5'	2.12	0.49
1:CA:1805:U:O2	4:CD:50:THR:HB	2.11	0.49
1:CA:2123:G:H21	3:CC:45:HIS:HE1	1.60	0.49
1:CA:2646:C:H6	1:CA:2646:C:O5'	1.95	0.49
1:CA:2661:G:H2'	1:CA:2662:A:O4'	2.12	0.49
10:CL:22:PRO:HA	10:CL:25:PRO:HD2	1.95	0.49
15:CR:96:ARG:HD2	15:CR:115:GLU:OE1	2.13	0.49
17:CT:62:THR:OG1	17:CT:75:ILE:HG12	2.12	0.49
22:CY:74:PRO:O	22:CY:82:PRO:HA	2.12	0.49
22:CY:90:LEU:HB2	22:CY:92:ASN:HB3	1.94	0.49
34:DA:1114:C:H42	34:DA:1186:G:H1	1.60	0.49
34:DA:1376:U:H2'	34:DA:1377:A:C8	2.47	0.49
34:DA:859:A:H2'	34:DA:860:A:O4'	2.12	0.49
36:DC:43:LEU:O	36:DC:47:LEU:N	2.24	0.49
34:DA:1235:U:H5''	54:DU:3:LYS:HB2	1.94	0.49
57:DZ:127:LYS:NZ	57:DZ:404:VAL:HG21	2.28	0.49
1:AA:2149:G:H21	1:AA:2195:A:H1'	1.77	0.49
1:AA:842:C:H2'	1:AA:843:C:C6	2.48	0.49
1:AA:1891:G:C3'	3:AC:206:LYS:HG3	2.43	0.49
13:AP:89:ALA:HA	13:AP:121:LYS:HE2	1.94	0.49
14:AQ:38:GLU:HB2	14:AQ:39:PRO:HD2	1.95	0.49
34:BA:794:A:OP2	63:BA:5146:HOH:O	2.20	0.49
35:BB:178:ARG:NH2	41:BH:74:PRO:HB3	2.28	0.49
35:BB:19:HIS:HE1	35:BB:189:ASP:CB	2.26	0.49
47:BN:32:SER:HB3	47:BN:41:ARG:HG2	1.95	0.49
56:BW:45:U:H2'	56:BW:45:U:OP2	2.13	0.49
1:CA:1922:G:H2'	1:CA:1923:U:O4'	2.13	0.49
1:CA:2840:C:H5''	15:CR:53:HIS:ND1	2.28	0.49
5:CE:52:LEU:O	5:CE:75:VAL:HA	2.11	0.49
9:CK:26:LEU:O	9:CK:114:GLY:N	2.43	0.49
10:CL:112:MET:HG3	10:CL:113:PRO:HD3	1.94	0.49
22:CY:38:ILE:HD11	22:CY:66:PRO:HG3	1.93	0.49
22:CY:49:VAL:HG21	22:CY:61:ILE:HG23	1.93	0.49
34:DA:1058:G:H2'	34:DA:1059:C:C6	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:1202:G:O4'	47:DN:29:ARG:NH1	2.46	0.49
34:DA:1286:A:C8	34:DA:1287:A:H4'	2.47	0.49
34:DA:430:A:OP1	37:DD:9:CYS:HB2	2.12	0.49
34:DA:491:G:H2'	34:DA:492:G:H8	1.78	0.49
34:DA:623:C:C4	34:DA:624:C:C5	3.01	0.49
42:DI:75:ASP:HA	42:DI:78:LYS:HE3	1.94	0.49
56:DW:37:MIA:H3'	56:DW:38:A:H8	1.77	0.49
57:DZ:133:ILE:HA	57:DZ:257:PRO:HD2	1.94	0.49
57:DZ:247:ARG:NH2	57:DZ:285:ASP:OD1	2.46	0.49
57:DZ:590:ILE:O	57:DZ:594:VAL:HG23	2.12	0.49
57:DZ:639:ASN:N	57:DZ:640:ALA:HB3	2.27	0.49
1:AA:1112:U:H2'	1:AA:1113:A:C8	2.47	0.49
1:AA:2207:C:H2'	1:AA:2208:G:H8	1.77	0.49
1:AA:894:U:H5	1:AA:978:A:H62	1.60	0.49
8:AH:125:VAL:HG12	8:AH:127:GLU:O	2.13	0.49
21:AX:35:THR:HG22	21:AX:38:GLU:CB	2.42	0.49
34:BA:1266:G:N2	34:BA:1270:C:N3	2.61	0.49
34:BA:232:G:H2'	34:BA:233:C:C6	2.48	0.49
38:BE:139:LEU:C	38:BE:141:GLN:H	2.15	0.49
40:BG:16:LEU:HD12	42:BI:45:ALA:HB2	1.94	0.49
2:CB:61:G:C6	2:CB:62:C:C4	3.00	0.49
3:CC:191:ARG:O	3:CC:194:ILE:HG22	2.12	0.49
5:CE:167:VAL:HG21	63:CE:3102:HOH:O	2.12	0.49
34:DA:1258:G:H2'	34:DA:1259:C:C6	2.47	0.49
34:DA:1392:G:O2'	34:DA:1393:U:H5'	2.13	0.49
34:DA:1460:A:H2'	34:DA:1461:G:O4'	2.13	0.49
34:DA:757:U:H2'	34:DA:758:G:O4'	2.12	0.49
35:DB:162:ILE:O	35:DB:185:ILE:HG12	2.12	0.49
37:DD:112:VAL:HG22	37:DD:116:GLN:OE1	2.12	0.49
37:DD:59:ARG:O	37:DD:61:LYS:N	2.45	0.49
42:DI:7:THR:O	42:DI:83:ARG:NH1	2.43	0.49
1:AA:1097:G:C6	1:AA:1098:C:C4	3.00	0.49
1:AA:672:G:H2'	1:AA:673:G:O4'	2.13	0.49
14:AQ:109:VAL:HG13	14:AQ:113:GLN:CB	2.39	0.49
19:AV:98:GLU:CD	19:AV:100:ARG:HH11	2.15	0.49
34:BA:1126:U:HO2'	34:BA:1127:G:H8	1.61	0.49
34:BA:1152:A:H4'	43:BJ:13:HIS:ND1	2.27	0.49
34:BA:924:C:O2'	34:BA:1502:A:N6	2.46	0.49
34:BA:975:A:C8	34:BA:1357:A:H2	2.31	0.49
35:BB:104:ASN:OD1	35:BB:107:THR:OG1	2.23	0.49
35:BB:42:ILE:HD12	35:BB:203:GLY:HA2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
40:BG:23:VAL:O	40:BG:27:ILE:HG12	2.12	0.49
56:BW:36:A:H1'	57:BZ:503:GLY:H	1.78	0.49
57:BZ:401:SER:OG	57:BZ:402:ILE:N	2.46	0.49
27:C3:8:LEU:HD12	27:C3:30:ARG:O	2.13	0.49
32:C8:63:PRO:HG2	32:C8:64:TYR:CE2	2.47	0.49
1:CA:1021:A:H8	1:CA:1022:G:H5''	1.78	0.49
1:CA:1472:A:H2'	1:CA:1473:G:O4'	2.12	0.49
1:CA:244:A:C2	1:CA:255:A:C4	3.01	0.49
2:CB:14:U:O3'	2:CB:108:U:O2'	2.30	0.49
12:CO:20:MET:HE3	12:CO:44:LYS:HE3	1.95	0.49
35:DB:133:LYS:C	35:DB:135:GLN:H	2.16	0.49
35:DB:78:GLN:HA	35:DB:94:ASN:ND2	2.27	0.49
44:DK:15:ALA:HB1	44:DK:78:GLN:HG3	1.95	0.49
45:DL:119:LYS:C	45:DL:121:GLY:H	2.15	0.49
1:AA:2377:G:O6	32:A8:39:LYS:HE3	2.12	0.49
1:AA:2286:A:O2'	1:AA:2288:G:OP1	2.22	0.49
8:AH:56:SER:OG	8:AH:57:ASP:N	2.45	0.49
1:AA:1134:A:N6	10:AL:133:SER:OG	2.46	0.49
34:BA:1279:A:H5''	34:BA:1280:A:OP1	2.13	0.49
34:BA:748:C:H4'	34:BA:749:C:O5'	2.13	0.49
34:BA:791:G:C6	34:BA:792:A:N7	2.81	0.49
35:BB:16:HIS:CG	35:BB:17:PHE:N	2.80	0.49
50:BQ:32:TYR:O	50:BQ:34:LYS:N	2.41	0.49
57:BZ:132:ARG:N	57:BZ:132:ARG:HD2	2.27	0.49
32:C8:63:PRO:HG2	32:C8:64:TYR:CD2	2.48	0.49
1:CA:2177:C:H5'	3:CC:45:HIS:HB2	1.95	0.49
1:CA:657:U:H2'	1:CA:658:C:C6	2.47	0.49
3:CC:183:PRO:C	3:CC:185:LYS:H	2.16	0.49
7:CG:64:THR:HG21	7:CG:92:VAL:HG11	1.94	0.49
15:CR:61:HIS:O	15:CR:65:LEU:HD22	2.13	0.49
34:DA:1367:C:H4'	43:DJ:48:THR:HG21	1.94	0.49
34:DA:1072:G:H21	35:DB:107:THR:HG21	1.77	0.49
20:AW:83:LYS:O	20:AW:84:ARG:HD3	2.13	0.49
34:BA:1346:A:N6	34:BA:1375:A:OP2	2.44	0.49
34:BA:1478:C:H2'	34:BA:1479:C:C6	2.48	0.49
38:BE:72:GLN:O	38:BE:75:THR:HG22	2.13	0.49
49:BP:67:THR:HG22	49:BP:69:THR:HG23	1.95	0.49
57:BZ:15:ILE:HA	57:BZ:103:GLY:O	2.13	0.49
57:BZ:600:VAL:HG21	57:BZ:678:GLU:HG3	1.94	0.49
1:CA:2420:C:OP2	32:C8:33:ASN:HB2	2.13	0.49
1:CA:1061:U:H4'	1:CA:1070:A:H1'	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1109:C:H5'	1:CA:1110:G:OP2	2.13	0.49
1:CA:141:A:H8	1:CA:1408:C:HO2'	1.60	0.49
1:CA:1420:U:HO2'	1:CA:1421:G:P	2.35	0.49
1:CA:729:G:OP2	4:CD:13:ARG:HD3	2.13	0.49
5:CE:108:SER:O	5:CE:162:ALA:HA	2.13	0.49
17:CT:93:ARG:HG2	17:CT:93:ARG:HH11	1.78	0.49
17:CT:99:LEU:HD22	17:CT:101:PHE:CE1	2.47	0.49
35:DB:178:ARG:HH22	41:DH:68:ARG:HH22	1.61	0.49
46:DM:10:PRO:HG2	46:DM:21:TYR:CD1	2.48	0.49
34:DA:1494:G:O3'	57:DZ:499:ARG:NH1	2.46	0.49
24:A0:24:LYS:O	24:A0:25:ARG:HD3	2.12	0.48
24:A0:73:GLY:O	24:A0:75:LEU:N	2.46	0.48
1:AA:139:A:H8	1:AA:1454:C:O2'	1.96	0.48
1:AA:1532:A:H2'	1:AA:1533:G:C8	2.46	0.48
1:AA:2701:U:P	1:AA:2732:G:H22	2.35	0.48
1:AA:520:G:N3	20:AW:61:ASN:ND2	2.61	0.48
7:AG:126:ASP:HB3	7:AG:128:ARG:H	1.78	0.48
34:BA:1310:G:H5'	46:BM:77:ASN:ND2	2.28	0.48
34:BA:516:U:C4	34:BA:517:G:C6	3.01	0.48
34:BA:939:G:OP1	40:BG:95:ARG:NH2	2.41	0.48
35:BB:32:ILE:HD13	35:BB:40:HIS:CG	2.48	0.48
41:BH:73:ASP:OD1	41:BH:75:ARG:HD3	2.13	0.48
34:BA:1123:A:O2'	43:BJ:37:PRO:O	2.28	0.48
48:BO:56:LEU:O	48:BO:60:VAL:HG23	2.12	0.48
39:BF:99:ALA:HB1	51:BR:23:LYS:HE3	1.95	0.48
28:A4:59:PHE:CE1	52:BS:45:VAL:HG21	2.48	0.48
53:BT:63:ILE:HD12	53:BT:81:LYS:HG2	1.95	0.48
57:BZ:554:PRO:HG3	57:BZ:594:VAL:O	2.12	0.48
29:C5:33:CYS:O	29:C5:37:LYS:N	2.46	0.48
1:CA:2630:G:H2'	1:CA:2631:G:C8	2.48	0.48
1:CA:2695:C:H2'	1:CA:2696:U:C6	2.46	0.48
1:CA:2748:A:C2	8:CH:63:SER:HB2	2.48	0.48
1:CA:2787:C:H1'	5:CE:62:PRO:HG3	1.95	0.48
1:CA:1812:A:O2'	4:CD:45:ASN:N	2.45	0.48
6:CF:129:PHE:HB2	6:CF:132:VAL:CG2	2.41	0.48
7:CG:136:ARG:HD2	7:CG:137:GLU:N	2.28	0.48
18:CU:79:PHE:HE1	18:CU:83:LEU:HD13	1.77	0.48
23:CZ:154:ASP:N	23:CZ:154:ASP:OD2	2.46	0.48
39:DF:42:GLU:OE1	39:DF:59:TYR:OH	2.19	0.48
45:DL:34:ARG:O	45:DL:61:THR:HG23	2.13	0.48
50:DQ:12:SER:HB3	50:DQ:20:THR:OG1	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:-6:ARG:O	57:DZ:-2:ALA:HB3	2.13	0.48
57:DZ:456:GLU:C	57:DZ:458:HIS:H	2.17	0.48
57:DZ:-55:LEU:HD22	57:DZ:-48:VAL:HG21	1.95	0.48
1:AA:1204:C:H4'	27:A3:32:GLN:HB2	1.95	0.48
1:AA:1216:G:N2	1:AA:1225:C:C2	2.81	0.48
1:AA:2517:G:H2'	1:AA:2588:G:O6	2.12	0.48
1:AA:2846:U:H2'	1:AA:2847:G:C8	2.48	0.48
10:AL:112:MET:HG2	10:AL:113:PRO:HD3	1.95	0.48
35:BB:56:ARG:HB2	35:BB:56:ARG:HE	1.37	0.48
36:BC:138:VAL:HG22	36:BC:151:VAL:HG23	1.95	0.48
46:BM:87:TYR:O	46:BM:91:ARG:HG2	2.13	0.48
57:BZ:-38:TYR:O	57:BZ:-34:ARG:HG2	2.13	0.48
57:BZ:461:ILE:HD12	57:BZ:462:ILE:H	1.78	0.48
57:BZ:534:ILE:HD11	57:BZ:570:GLY:HA3	1.96	0.48
1:CA:1005:C:H2'	1:CA:1006:C:C6	2.48	0.48
1:CA:1773:A:H2'	1:CA:1774:C:O4'	2.13	0.48
1:CA:2303:G:O2'	7:CG:132:ASN:HB2	2.13	0.48
1:CA:7:G:H4'	11:CN:13:TRP:CH2	2.48	0.48
3:CC:42:VAL:HA	3:CC:216:THR:O	2.13	0.48
5:CE:167:VAL:HG22	5:CE:170:LEU:HD11	1.94	0.48
6:CF:132:VAL:O	6:CF:132:VAL:HG23	2.13	0.48
11:CN:38:HIS:O	18:CU:67:ALA:HB1	2.13	0.48
16:CS:66:ALA:HA	16:CS:69:VAL:HG12	1.94	0.48
17:CT:53:ARG:NH1	17:CT:58:ASN:O	2.45	0.48
22:CY:16:ALA:HB2	22:CY:73:ARG:HG3	1.96	0.48
22:CY:46:LYS:HG2	22:CY:60:PHE:CD2	2.48	0.48
34:DA:1087:G:N2	34:DA:1099:G:H1'	2.28	0.48
34:DA:942:G:C2	34:DA:1342:C:C2	3.01	0.48
34:DA:1531:A:N7	34:DA:1532:U:C4	2.81	0.48
34:DA:513:C:H2'	34:DA:514:C:H6	1.78	0.48
34:DA:913:A:OP2	45:DL:91:LYS:NZ	2.45	0.48
35:DB:91:PRO:HG3	35:DB:155:LEU:HG	1.96	0.48
38:DE:76:ILE:HB	38:DE:77:PRO:HD2	1.93	0.48
43:DJ:38:ILE:HG12	43:DJ:71:LEU:O	2.13	0.48
48:DO:33:THR:HG21	48:DO:85:LEU:HD22	1.94	0.48
56:DY:62:C:H2'	56:DY:63:G:C8	2.48	0.48
1:AA:1105:G:H1'	10:AL:126:MET:HE3	1.94	0.48
1:AA:628:C:H2'	1:AA:629:U:O4'	2.13	0.48
5:AE:143:ASN:HD22	5:AE:147:PRO:CD	2.25	0.48
5:AE:68:ALA:C	5:AE:70:ALA:H	2.17	0.48
6:AF:157:VAL:HB	6:AF:194:MET:HG2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AO:100:GLY:H	12:AO:119:PRO:HG2	1.78	0.48
18:AU:69:CYS:HB3	18:AU:74:LEU:HD13	1.94	0.48
23:AZ:74:VAL:HG13	23:AZ:86:VAL:HG13	1.95	0.48
34:BA:1277:C:O2'	34:BA:1279:A:H1'	2.13	0.48
34:BA:1260:C:O5'	34:BA:1284:C:H4'	2.13	0.48
34:BA:475:G:H2'	34:BA:476:G:C8	2.48	0.48
34:BA:48:C:H5''	34:BA:365:U:O4	2.13	0.48
39:BF:60:PHE:CE2	51:BR:78:LEU:HD21	2.48	0.48
40:BG:107:ALA:HB3	40:BG:134:ALA:HB2	1.94	0.48
34:BA:580:U:H5''	48:BO:58:MET:HG2	1.95	0.48
57:BZ:116:PRO:O	57:BZ:118:SER:N	2.43	0.48
57:BZ:406:GLU:HG2	57:BZ:439:ARG:NH2	2.28	0.48
1:CA:1614:A:H8	1:CA:1614:A:P	2.35	0.48
1:CA:2097:C:H2'	1:CA:2098:U:C6	2.48	0.48
1:CA:2113:U:H2'	1:CA:2114:A:C8	2.47	0.48
1:CA:2267:A:H5''	1:CA:2268:A:C5'	2.43	0.48
1:CA:197:A:N6	1:CA:2430:A:H2'	2.27	0.48
1:CA:468:G:N7	31:C7:39:ARG:NH2	2.50	0.48
1:CA:55:G:N3	1:CA:127:A:H2	2.11	0.48
1:CA:741:G:O2'	1:CA:742:G:H5'	2.14	0.48
3:CC:184:GLU:O	3:CC:188:ASP:OD2	2.31	0.48
7:CG:41:GLN:HB3	7:CG:43:LEU:HD22	1.95	0.48
8:CH:140:LYS:O	8:CH:144:VAL:HG23	2.13	0.48
10:CL:77:LEU:HD11	10:CL:111:LYS:HD2	1.96	0.48
23:CZ:54:HIS:HA	23:CZ:98:MET:HE1	1.95	0.48
34:DA:309:G:H1'	34:DA:608:A:C2	2.47	0.48
28:A4:14:ILE:HG23	28:A4:31:ILE:HG22	1.94	0.48
1:AA:2141:A:O2'	1:AA:2142:G:H5'	2.13	0.48
1:AA:2518:U:O2'	1:AA:2519:C:O5'	2.30	0.48
23:AZ:111:VAL:HG12	23:AZ:112:ARG:H	1.79	0.48
34:BA:1274:G:N2	34:BA:1275:A:H62	2.11	0.48
34:BA:27:G:H2'	34:BA:28:G:H8	1.78	0.48
42:BI:111:ARG:NH2	43:BJ:62:HIS:HD2	2.11	0.48
42:BI:83:ARG:HA	42:BI:86:VAL:HG22	1.95	0.48
44:BK:31:THR:HA	44:BK:42:TRP:HA	1.94	0.48
34:BA:1202:G:O4'	47:BN:29:ARG:NH1	2.45	0.48
51:BR:59:SER:OG	51:BR:62:GLU:HG2	2.13	0.48
53:BT:99:LEU:HA	53:BT:100:ILE:O	2.12	0.48
57:BZ:132:ARG:H	57:BZ:132:ARG:HD2	1.78	0.48
57:BZ:127:LYS:NZ	57:BZ:404:VAL:HG11	2.29	0.48
57:BZ:493:VAL:HG23	57:BZ:512:ILE:HD11	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2814:C:O2'	29:C5:29:THR:HG21	2.13	0.48
1:CA:1540:U:O2'	1:CA:1541:G:H5'	2.14	0.48
1:CA:1589:C:H2'	1:CA:1590:U:C6	2.48	0.48
1:CA:2306:C:H3'	1:CA:2307:G:H8	1.79	0.48
1:CA:2529:G:H5''	1:CA:2530:A:H5''	1.95	0.48
1:CA:271(A):A:N7	1:CA:271(W):G:N2	2.58	0.48
1:CA:969:U:H2'	1:CA:970:C:C6	2.49	0.48
34:DA:33:A:N3	45:DL:32:PHE:HE2	2.12	0.48
35:DB:44:LEU:HD22	35:DB:44:LEU:H	1.78	0.48
37:DD:119:GLN:HG3	37:DD:123:HIS:HD2	1.78	0.48
34:DA:401:C:P	37:DD:73:ARG:HE	2.35	0.48
37:DD:76:ARG:O	37:DD:80:GLU:HG2	2.13	0.48
38:DE:110:LEU:HD13	38:DE:118:ILE:HD13	1.95	0.48
38:DE:68:GLU:HG2	38:DE:69:VAL:N	2.28	0.48
40:DG:44:TYR:HA	40:DG:47:CYS:SG	2.53	0.48
57:DZ:503:GLY:C	57:DZ:505:GLY:H	2.17	0.48
57:DZ:495:GLY:HA3	57:DZ:589:ALA:HB2	1.95	0.48
28:A4:36:CYS:SG	28:A4:37:SER:N	2.86	0.48
1:AA:2118:U:H2'	1:AA:2119:C:H6	1.77	0.48
1:AA:2504:U:H2'	1:AA:2505:U:C6	2.49	0.48
1:AA:2627:U:H2'	1:AA:2628:C:H6	1.78	0.48
1:AA:929:G:H1	1:AA:940:C:H42	1.61	0.48
1:AA:2200:C:H4'	3:AC:47:LYS:CE	2.43	0.48
5:AE:59:VAL:O	5:AE:64:LYS:HE2	2.13	0.48
21:AX:84:ALA:HB3	21:AX:87:GLN:NE2	2.29	0.48
34:BA:659:U:N3	34:BA:660:G:N7	2.61	0.48
38:BE:6:PHE:HE2	38:BE:36:ASP:HB3	1.79	0.48
40:BG:45:ASP:O	40:BG:49:ILE:HG13	2.12	0.48
41:BH:121:ASP:OD1	41:BH:125:ARG:NH2	2.47	0.48
44:BK:87:THR:OG1	44:BK:87:THR:O	2.28	0.48
26:C2:22:GLU:OE2	26:C2:68:ARG:NH2	2.45	0.48
1:CA:1448:G:N2	1:CA:1463:C:O2	2.43	0.48
1:CA:235:U:C4	1:CA:236:C:N4	2.81	0.48
1:CA:2686:G:C2	1:CA:2724:C:O2	2.66	0.48
1:CA:2704:C:H2'	1:CA:2705:A:O4'	2.13	0.48
1:CA:897:C:H3'	1:CA:898:C:C6	2.49	0.48
4:CD:221:VAL:HG22	4:CD:226:MET:CE	2.44	0.48
1:CA:1063:G:H21	10:CL:91:PRO:HG2	1.78	0.48
14:CQ:54:MET:HG2	14:CQ:117:ALA:O	2.14	0.48
19:CV:32:THR:HG22	19:CV:60:GLU:HB2	1.95	0.48
22:CY:45:VAL:HG23	22:CY:63:LYS:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DB:42:ILE:HG21	35:DB:202:PRO:O	2.13	0.48
37:DD:149:ALA:O	37:DD:152:SER:N	2.45	0.48
37:DD:156:GLU:O	37:DD:160:GLN:HG3	2.13	0.48
41:DH:29:SER:HB3	41:DH:32:LYS:HB2	1.96	0.48
57:DZ:119:GLU:O	57:DZ:122:TRP:N	2.46	0.48
57:DZ:187:THR:HB	57:DZ:199:ILE:HD11	1.96	0.48
1:AA:1684:A:H5'	1:AA:1791:A:O2'	2.13	0.48
1:AA:2705:A:H2'	1:AA:2706:G:C8	2.49	0.48
2:AB:1:U:O2	2:AB:1:U:H2'	2.13	0.48
4:AD:245:PRO:HB3	4:AD:253:GLN:NE2	2.29	0.48
17:AT:16:ARG:NH1	17:AT:19:LEU:HD21	2.29	0.48
1:AA:2857:U:OP1	17:AT:98:LYS:HD3	2.13	0.48
18:AU:76:TYR:HH	18:AU:92:ARG:HH11	1.61	0.48
23:AZ:128:VAL:HG22	23:AZ:129:SER:O	2.14	0.48
34:BA:1236:A:H2'	34:BA:1237:C:C6	2.48	0.48
34:BA:1511:G:H2'	34:BA:1512:U:O4'	2.14	0.48
34:BA:186:C:H2'	34:BA:187:C:C6	2.46	0.48
34:BA:603:U:H2'	34:BA:604:G:C8	2.48	0.48
35:BB:105:PHE:C	35:BB:107:THR:H	2.16	0.48
37:BD:63:LYS:HE3	37:BD:197:PRO:O	2.13	0.48
57:BZ:248:LYS:O	57:BZ:252:ASP:N	2.42	0.48
1:CA:1165:U:H2'	1:CA:1166:C:C6	2.48	0.48
1:CA:1316:U:H2'	1:CA:1317:A:H8	1.74	0.48
1:CA:1683:C:H2'	1:CA:1684:C:H6	1.77	0.48
1:CA:2279:G:N7	24:C0:14:ARG:NH1	2.62	0.48
1:CA:2622:C:H5'	5:CE:159:HIS:ND1	2.28	0.48
2:CB:19:G:H2'	2:CB:20:C:O4'	2.14	0.48
6:CF:187:VAL:HG13	13:CP:1:MET:O	2.13	0.48
23:CZ:108:PRO:HB3	23:CZ:144:LEU:HB2	1.96	0.48
23:CZ:111:VAL:O	23:CZ:112:ARG:HB2	2.13	0.48
34:DA:1161:C:H2'	34:DA:1162:C:C6	2.47	0.48
34:DA:426:G:C6	34:DA:427:U:C4	3.01	0.48
34:DA:7:G:O2'	38:DE:120:THR:O	2.32	0.48
36:DC:113:ALA:HB2	36:DC:183:ASP:HB3	1.96	0.48
36:DC:149:ALA:HA	36:DC:201:TYR:O	2.13	0.48
49:DP:29:ASP:N	49:DP:29:ASP:OD1	2.44	0.48
57:DZ:635:GLU:HG3	57:DZ:636:PRO:HD2	1.96	0.48
30:A6:25:LYS:NZ	30:A6:51:GLU:OE1	2.46	0.48
1:AA:1359:U:H2'	1:AA:1656:A:C2	2.49	0.48
1:AA:1402:G:H2'	1:AA:1403:U:H6	1.78	0.48
1:AA:2679:C:H1'	8:AH:109:PHE:CD1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2735:G:H2'	1:AA:2736:C:C6	2.49	0.48
1:AA:2798:C:H2'	1:AA:2799:U:O4'	2.13	0.48
1:AA:768:C:H2'	1:AA:769:A:C8	2.49	0.48
4:AD:67:PHE:HB3	4:AD:153:ALA:H	1.77	0.48
8:AH:3:ARG:HG2	8:AH:6:ARG:HD2	1.94	0.48
34:BA:1092:A:C6	34:BA:1093:A:C6	3.02	0.48
34:BA:1238:A:N3	34:BA:1241:G:O2'	2.40	0.48
34:BA:6:G:C4	38:BE:119:LEU:HD11	2.48	0.48
42:BI:17:VAL:HG23	42:BI:63:ILE:HG12	1.96	0.48
43:BJ:11:PHE:HE1	43:BJ:67:THR:HG22	1.78	0.48
51:BR:40:LEU:HD22	51:BR:70:ILE:HG12	1.95	0.48
57:BZ:138:LYS:HA	62:BZ:704:GDP:N1	2.28	0.48
1:CA:1810:A:H2'	1:CA:1811:G:O4'	2.13	0.48
1:CA:1861:G:P	3:CC:206:LYS:CA	2.86	0.48
1:CA:2250:G:H5''	1:CA:2250:G:N3	2.28	0.48
1:CA:910:A:N1	1:CA:2277:G:H1'	2.28	0.48
3:CC:17:PRO:HG2	3:CC:18:ASN:H	1.79	0.48
5:CE:4:ILE:HD11	5:CE:29:GLY:HA2	1.96	0.48
5:CE:98:PRO:HD3	5:CE:175:VAL:HG13	1.96	0.48
10:CL:13:PRO:HA	10:CL:52:ILE:HG23	1.95	0.48
1:CA:955:C:OP1	14:CQ:87:LYS:HE2	2.13	0.48
34:DA:107:G:C2	34:DA:108:G:H1'	2.49	0.48
34:DA:243:A:C2	34:DA:245:C:C2	3.02	0.48
35:DB:174:VAL:HG13	35:DB:184:VAL:HG11	1.93	0.48
42:DI:114:TYR:HE1	43:DJ:60:ARG:H	1.59	0.48
34:DA:707:C:H4'	44:DK:20:TYR:CD2	2.49	0.48
46:DM:19:LEU:O	46:DM:22:ILE:HG12	2.13	0.48
46:DM:66:LEU:O	46:DM:68:GLY:N	2.47	0.48
56:DW:45:U:O2'	56:DW:46:7MG:H82	2.14	0.48
57:DZ:9:LEU:O	57:DZ:283:PRO:HD2	2.14	0.48
25:A1:8:SER:HB3	25:A1:66:HIS:CD2	2.49	0.48
1:AA:118:U:H5'	1:AA:118:U:H6	1.78	0.48
3:AC:42:VAL:HA	3:AC:216:THR:O	2.13	0.48
13:AP:94:GLU:OE2	13:AP:124:LYS:HD3	2.14	0.48
20:AW:5:ALA:C	20:AW:6:ILE:HG13	2.33	0.48
35:BB:192:SER:O	35:BB:194:PRO:HD3	2.14	0.48
43:BJ:11:PHE:CE1	43:BJ:67:THR:HG22	2.49	0.48
43:BJ:44:VAL:HG22	43:BJ:66:ARG:HE	1.78	0.48
34:BA:375:U:H4'	49:BP:17:TYR:CE1	2.48	0.48
56:BW:19:G:H4'	56:BW:20:U:OP2	2.14	0.48
56:BY:57:G:H2'	56:BY:58:A:H5'	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:BZ:13:ARG:HH21	57:BZ:77:HIS:CE1	2.32	0.48
57:BZ:238:THR:HG23	57:BZ:241:GLU:HG2	1.94	0.48
1:CA:118:A:H1'	1:CA:178:G:O4'	2.13	0.48
1:CA:1809:A:H2'	1:CA:1810:A:C8	2.48	0.48
1:CA:899:A:O2'	1:CA:900:A:H5'	2.13	0.48
19:CV:76:LYS:O	19:CV:79:VAL:HG12	2.13	0.48
22:CY:102:CYS:SG	22:CY:104:GLY:N	2.80	0.48
23:CZ:73:GLN:H	23:CZ:87:ASP:HB2	1.79	0.48
34:DA:1024:G:H2'	34:DA:1025:U:H5''	1.95	0.48
34:DA:1179:A:H2'	34:DA:1180:A:O4'	2.13	0.48
34:DA:535:A:H4'	34:DA:536:C:OP2	2.14	0.48
37:DD:59:ARG:NE	37:DD:59:ARG:HA	2.29	0.48
43:DJ:11:PHE:HE1	43:DJ:67:THR:HB	1.78	0.48
50:DQ:62:SER:HB3	50:DQ:72:ARG:HD3	1.94	0.48
34:DA:1317:C:O2	52:DS:37:ARG:NH1	2.46	0.48
46:DM:124:PRO:HB3	56:DW:36:A:H4'	1.96	0.48
56:DY:69:G:H2'	56:DY:70:G:O4'	2.14	0.48
57:DZ:346:LYS:HZ1	57:DZ:384:ILE:CD1	2.27	0.48
57:DZ:-66:MET:HB2	57:DZ:-65:LYS:H	1.47	0.48
24:A0:18:ALA:HB3	24:A0:20:ARG:HH21	1.78	0.48
1:AA:1293:A:OP1	6:AF:95:ARG:NH2	2.47	0.48
1:AA:1495:G:H1'	1:AA:1574:A:N1	2.27	0.48
1:AA:185:A:H2'	1:AA:185:A:N3	2.29	0.48
1:AA:2122:G:C2	1:AA:2212:G:C2	3.02	0.48
1:AA:217:A:H3'	1:AA:218:A:C5'	2.44	0.48
1:AA:2673:G:H5''	57:BZ:21:ILE:HG13	1.95	0.48
1:AA:733:G:H4'	1:AA:734:C:OP2	2.12	0.48
1:AA:2144:U:H1'	3:AC:167:ASP:CB	2.44	0.48
5:AE:52:LEU:O	5:AE:75:VAL:HA	2.14	0.48
6:AF:178:PRO:C	6:AF:180:GLY:H	2.17	0.48
10:AL:12:LEU:HD11	10:AL:23:VAL:HG22	1.96	0.48
14:AQ:14:ARG:HG2	14:AQ:41:TRP:CH2	2.35	0.48
15:AR:44:LEU:HD22	15:AR:48:VAL:HG23	1.95	0.48
16:AS:89:ARG:HD2	16:AS:92:TYR:O	2.13	0.48
18:AU:74:LEU:HD12	18:AU:74:LEU:H	1.78	0.48
21:AX:60:ARG:NH2	31:A7:47:ARG:HH21	2.12	0.48
23:AZ:111:VAL:C	23:AZ:113:ALA:H	2.17	0.48
34:BA:1323:G:H2'	34:BA:1324:A:C8	2.48	0.48
34:BA:557:G:C6	34:BA:558:G:C6	3.01	0.48
27:C3:30:ARG:NH2	27:C3:33:GLN:HE21	2.12	0.48
1:CA:103:A:H5'	26:C2:3:LEU:HD11	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1268:A:H2'	1:CA:1269:A:O5'	2.13	0.48
1:CA:1452:A:O2'	1:CA:1453:U:H2'	2.14	0.48
1:CA:2271:G:C6	1:CA:2272:U:C4	3.01	0.48
1:CA:2852:G:C2	1:CA:2853:C:O2	2.67	0.48
1:CA:862:G:H2'	1:CA:863:A:O4'	2.14	0.48
6:CF:21:ALA:HB3	6:CF:22:ALA:HA	1.96	0.48
9:CK:40:LEU:HA	9:CK:43:ALA:HB3	1.96	0.48
22:CY:83:THR:HG21	22:CY:99:CYS:HA	1.96	0.48
34:DA:151:A:N6	34:DA:171:A:N7	2.62	0.48
34:DA:382:A:H2'	34:DA:383:A:C8	2.48	0.48
34:DA:544:G:C2	34:DA:545:C:C2	3.01	0.48
34:DA:741:G:H2'	34:DA:742:G:O4'	2.13	0.48
34:DA:1060:C:C4	36:DC:2:GLY:HA3	2.49	0.48
36:DC:57:ILE:HG13	36:DC:66:VAL:HG22	1.96	0.48
39:DF:87:ARG:HH11	39:DF:87:ARG:CG	2.17	0.48
42:DI:28:VAL:HG22	42:DI:63:ILE:HB	1.96	0.48
43:DJ:47:PHE:N	43:DJ:63:PHE:O	2.41	0.48
44:DK:95:ILE:O	44:DK:98:LEU:N	2.47	0.48
52:DS:3:ARG:HH11	52:DS:7:LYS:HE2	1.79	0.48
57:DZ:15:ILE:HG21	57:DZ:280:LEU:HD13	1.94	0.48
27:A3:8:LEU:HD13	27:A3:31:LEU:HD23	1.95	0.48
33:A9:7:VAL:HG12	33:A9:34:GLN:HB3	1.94	0.48
1:AA:2509:A:H5''	63:AA:3911:HOH:O	2.13	0.48
1:AA:941:U:H5'	1:AA:942:A:P	2.54	0.48
5:AE:29:GLY:O	5:AE:51:PHE:HE1	1.97	0.48
13:AP:59:LEU:HD11	32:A8:10:ALA:HB2	1.96	0.48
22:AY:92:ASN:HB2	22:AY:94:LYS:N	2.25	0.48
34:BA:1183:A:O2'	34:BA:1184:G:OP1	2.29	0.48
34:BA:343:U:H3	34:BA:347:G:N2	2.01	0.48
34:BA:520:A:C2	34:BA:536:C:H1'	2.49	0.48
34:BA:617:G:O6	34:BA:623:C:N4	2.40	0.48
34:BA:662:G:H2'	34:BA:663:A:C8	2.49	0.48
34:BA:9:G:C2	34:BA:26:A:N1	2.82	0.48
41:BH:82:HIS:O	41:BH:137:VAL:HA	2.14	0.48
49:BP:69:THR:O	49:BP:69:THR:OG1	2.32	0.48
57:BZ:485:GLU:O	57:BZ:486:THR:HB	2.14	0.48
7:CG:112:PRO:HG3	28:C4:43:TYR:CE2	2.49	0.48
29:C5:36:CYS:O	29:C5:38:ALA:N	2.47	0.48
32:C8:30:ARG:O	63:C8:6001:HOH:O	2.19	0.48
1:CA:1651:G:C2	1:CA:2007:C:C2	3.02	0.48
1:CA:1899:G:H2'	1:CA:1899:G:N3	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2839:G:H5'	15:CR:46:GLY:CA	2.44	0.48
6:CF:53:THR:HG23	6:CF:55:GLY:H	1.79	0.48
21:CX:9:LEU:HB2	21:CX:29:TRP:O	2.14	0.48
1:CA:480:A:OP2	22:CY:47:LYS:HE2	2.13	0.48
34:DA:57:G:C2	34:DA:58:C:C2	3.02	0.48
43:DJ:13:HIS:HB3	43:DJ:68:HIS:HE1	1.77	0.48
47:DN:22:THR:HB	47:DN:33:VAL:HB	1.95	0.48
49:DP:5:ARG:NH1	49:DP:28:ARG:HA	2.28	0.48
51:DR:38:GLU:HA	51:DR:41:LYS:NZ	2.29	0.48
52:DS:41:VAL:O	52:DS:43:GLU:N	2.47	0.48
57:DZ:20:HIS:CG	57:DZ:21:ILE:N	2.82	0.48
1:AA:2289:G:OP2	24:A0:10:THR:HG21	2.14	0.47
1:AA:2283:G:OP1	24:A0:18:ALA:HB1	2.13	0.47
1:AA:1540:A:OP2	63:AA:4072:HOH:O	2.20	0.47
1:AA:555:G:C5	1:AA:2044:U:H5''	2.49	0.47
1:AA:215:G:N2	1:AA:217:A:H62	2.12	0.47
1:AA:645:G:N3	1:AA:645:G:H2'	2.28	0.47
1:AA:821:A:HO2'	1:AA:822:G:H8	1.61	0.47
3:AC:184:GLU:O	3:AC:188:ASP:OD2	2.31	0.47
3:AC:68:GLY:N	3:AC:189:ASN:ND2	2.62	0.47
1:AA:1834:A:H4'	4:AD:259:THR:HG23	1.95	0.47
10:AL:56:GLU:O	10:AL:67:PHE:HA	2.14	0.47
11:AN:58:ASP:N	11:AN:58:ASP:OD1	2.36	0.47
34:BA:1034:G:H3'	34:BA:1035:A:C8	2.48	0.47
34:BA:1155:G:H2'	34:BA:1156:G:O4'	2.13	0.47
34:BA:1311:G:H2'	34:BA:1312:G:O4'	2.13	0.47
34:BA:67:C:O2'	34:BA:171:A:H1'	2.14	0.47
34:BA:177:C:H2'	34:BA:178:C:H6	1.79	0.47
34:BA:604:G:C2	34:BA:635:G:C5	3.02	0.47
38:BE:39:GLY:O	38:BE:69:VAL:HG13	2.14	0.47
53:BT:63:ILE:HD13	53:BT:80:ARG:HB2	1.96	0.47
53:BT:71:THR:O	53:BT:72:LEU:HD23	2.14	0.47
57:BZ:655:TYR:O	57:BZ:657:THR:N	2.47	0.47
14:CQ:81:VAL:HB	24:C0:7:LEU:HD11	1.95	0.47
30:C6:3:SER:OG	30:C6:5:VAL:HG13	2.14	0.47
1:CA:1469:A:H2'	1:CA:1470:G:C8	2.49	0.47
1:CA:1488:G:H5'	1:CA:1489:U:OP2	2.13	0.47
1:CA:2124:G:H4'	3:CC:175:PRO:HD3	1.96	0.47
1:CA:2309:A:N6	1:CA:2310:A:N1	2.62	0.47
1:CA:2611:U:C4	29:C5:3:LYS:HG2	2.49	0.47
1:CA:2626:C:H2'	1:CA:2627:G:O4'	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2845:G:H2'	1:CA:2846:G:C8	2.49	0.47
1:CA:2861:G:H2'	1:CA:2862:G:C8	2.49	0.47
14:CQ:75:THR:HA	14:CQ:89:ASN:O	2.14	0.47
16:CS:67:ARG:HH11	16:CS:67:ARG:HG2	1.79	0.47
17:CT:14:TYR:N	17:CT:14:TYR:CD2	2.82	0.47
34:DA:1096:C:H2'	34:DA:1097:C:H6	1.79	0.47
34:DA:171:A:H2'	34:DA:172:A:C8	2.49	0.47
34:DA:502:G:N2	34:DA:544:G:C4	2.82	0.47
35:DB:69:LEU:HD12	35:DB:70:PHE:N	2.28	0.47
36:DC:32:LEU:HD12	36:DC:59:ARG:NH1	2.30	0.47
45:DL:34:ARG:HG2	45:DL:35:GLY:N	2.29	0.47
51:DR:40:LEU:HD22	51:DR:70:ILE:HG12	1.96	0.47
56:DW:51:U:H3	56:DW:63:G:H1	1.60	0.47
25:A1:3:LYS:HG2	25:A1:4:VAL:N	2.29	0.47
1:AA:27:G:C4	1:AA:537:G:N2	2.82	0.47
3:AC:30:VAL:CG2	3:AC:31:LYS:H	2.27	0.47
8:AH:174:GLY:O	8:AH:175:LYS:HB2	2.15	0.47
14:AQ:138:ASP:OD2	23:AZ:81:ARG:NH1	2.46	0.47
34:BA:269:C:H2'	34:BA:270:A:C8	2.49	0.47
34:BA:445:G:C6	34:BA:446:G:C5	3.03	0.47
38:BE:132:ALA:O	38:BE:134:ALA:N	2.47	0.47
38:BE:74:GLY:HA3	38:BE:116:THR:HG22	1.95	0.47
1:CA:1434:A:H61	1:CA:1558:A:N6	2.10	0.47
1:CA:1577:C:H2'	1:CA:1578:U:C1'	2.43	0.47
1:CA:1932:A:H2'	1:CA:1933:G:O4'	2.14	0.47
1:CA:2531:A:H5''	8:CH:157:TYR:CZ	2.49	0.47
1:CA:25:U:H5'	20:CW:79:GLY:HA2	1.95	0.47
11:CN:56:ASN:N	11:CN:125:GLY:O	2.37	0.47
11:CN:14:VAL:CG1	11:CN:138:LEU:HB2	2.43	0.47
15:CR:33:ARG:NH2	29:C5:57:VAL:HG12	2.29	0.47
16:CS:67:ARG:CG	16:CS:67:ARG:HH11	2.27	0.47
18:CU:79:PHE:CZ	18:CU:83:LEU:HD22	2.49	0.47
34:DA:1217:C:H2'	34:DA:1218:C:C6	2.48	0.47
34:DA:1304:G:C5	34:DA:1305:G:C6	3.02	0.47
34:DA:266:G:H5'	34:DA:266:G:C8	2.49	0.47
34:DA:922:G:C2	34:DA:923:A:C4	3.01	0.47
56:DW:23:A:H2'	56:DW:24:G:C8	2.49	0.47
1:AA:1220:U:O3'	1:AA:1221:G:H4'	2.14	0.47
1:AA:217:A:H8	1:AA:218:A:H5'	1.78	0.47
7:AG:126:ASP:HB2	7:AG:130:ASN:O	2.13	0.47
7:AG:77:ILE:HB	7:AG:82:LEU:HD12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:AV:72:VAL:HG11	19:AV:85:LYS:HD2	1.96	0.47
22:AY:20:TYR:CE1	22:AY:43:ASN:HA	2.50	0.47
34:BA:1142:G:H2'	34:BA:1143:G:O4'	2.14	0.47
34:BA:391:G:C6	34:BA:392:G:C5	3.02	0.47
34:BA:785:G:N2	34:BA:798:G:C4	2.82	0.47
39:BF:44:GLY:O	39:BF:60:PHE:N	2.42	0.47
53:BT:44:ALA:HB1	53:BT:91:LEU:HB2	1.95	0.47
56:BW:19:G:C4	56:BW:57:G:N2	2.82	0.47
56:BY:58:A:O2'	56:BY:60:U:H5	1.98	0.47
57:BZ:445:GLU:O	57:BZ:447:GLY:N	2.42	0.47
57:BZ:99:ARG:C	57:BZ:101:LEU:N	2.67	0.47
1:CA:1204:A:H2	1:CA:1241:A:N6	2.04	0.47
1:CA:1469:A:H2'	1:CA:1470:G:H8	1.79	0.47
1:CA:1685:C:H2'	1:CA:1686:C:C6	2.49	0.47
1:CA:1912:A:O2'	34:DA:1494:G:O2'	2.32	0.47
1:CA:1939:U:OP1	1:CA:2604:U:O2'	2.25	0.47
1:CA:2334:G:H5'	16:CS:9:ARG:HG2	1.95	0.47
1:CA:2611:U:H6	1:CA:2611:U:H5'	1.79	0.47
1:CA:706:A:H2'	1:CA:707:G:O4'	2.14	0.47
1:CA:892:G:H2'	1:CA:893:C:O4'	2.14	0.47
8:CH:157:TYR:CE1	8:CH:172:LYS:HG2	2.49	0.47
18:CU:79:PHE:CE1	18:CU:83:LEU:HD13	2.49	0.47
23:CZ:101:PRO:O	23:CZ:102:LEU:HD12	2.14	0.47
34:DA:1070:U:H2'	34:DA:1071:C:H6	1.79	0.47
34:DA:303:A:HO2'	34:DA:555:C:HO2'	1.56	0.47
34:DA:434:U:H2'	34:DA:435:C:C6	2.49	0.47
37:DD:108:LEU:HD21	37:DD:183:GLY:HA3	1.96	0.47
37:DD:59:ARG:O	37:DD:62:GLN:N	2.47	0.47
61:DZ:703:FUA:O1	61:DZ:703:FUA:C20	2.62	0.47
57:DZ:-9:LEU:O	57:DZ:-7:GLU:N	2.47	0.47
30:A6:8:LYS:HD3	32:A8:34:TRP:CD2	2.49	0.47
1:AA:1447:G:H2'	1:AA:1448:C:O4'	2.14	0.47
1:AA:1451:U:H2'	1:AA:1452:U:H6	1.79	0.47
1:AA:2053:A:C6	1:AA:2510:C:H1'	2.49	0.47
1:AA:2059:G:O6	63:AA:4616:HOH:O	2.18	0.47
1:AA:2586:G:H2'	1:AA:2587:C:C6	2.50	0.47
1:AA:265:U:H2'	1:AA:266:C:C6	2.50	0.47
1:AA:310:C:H2'	1:AA:311:C:H6	1.79	0.47
4:AD:261:LYS:HZ1	4:AD:263:ARG:NH2	2.12	0.47
7:AG:33:ARG:O	7:AG:162:THR:HG23	2.14	0.47
20:AW:9:TYR:HA	20:AW:100:THR:HG23	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:AZ:70:LEU:HD11	23:AZ:98:MET:HE3	1.96	0.47
34:BA:1037:C:H2'	34:BA:1038:C:C6	2.49	0.47
34:BA:1053:G:C4	34:BA:1199:U:C5	3.03	0.47
34:BA:1187:G:H2'	34:BA:1188:A:C8	2.49	0.47
34:BA:33:A:H2'	34:BA:34:C:C6	2.49	0.47
34:BA:652:U:O4	34:BA:752:G:O2'	2.26	0.47
34:BA:707:C:H2'	34:BA:708:C:C6	2.48	0.47
34:BA:899:C:H2'	34:BA:900:A:C8	2.49	0.47
37:BD:111:ALA:HB2	37:BD:120:LEU:HD12	1.95	0.47
38:BE:139:LEU:HA	38:BE:142:LEU:HD12	1.96	0.47
48:BO:62:GLN:O	48:BO:65:ARG:N	2.47	0.47
53:BT:44:ALA:HB3	53:BT:91:LEU:HD12	1.95	0.47
57:BZ:357:ARG:HH11	57:BZ:373:ASP:CG	2.17	0.47
57:BZ:417:THR:HA	57:BZ:418:LYS:CB	2.45	0.47
1:CA:1075:C:H5'	1:CA:1076:C:OP2	2.14	0.47
1:CA:2103:C:H1'	1:CA:2187:G:N2	2.29	0.47
1:CA:237:C:H2'	1:CA:238:C:H6	1.79	0.47
1:CA:495:G:O2'	20:CW:61:ASN:ND2	2.47	0.47
5:CE:163:GLU:HG2	5:CE:164:ARG:N	2.30	0.47
8:CH:118:PRO:HG2	8:CH:121:ILE:HB	1.96	0.47
14:CQ:63:LYS:HA	23:CZ:178:GLU:CG	2.41	0.47
16:CS:66:ALA:O	16:CS:69:VAL:HG12	2.14	0.47
17:CT:99:LEU:O	17:CT:101:PHE:N	2.47	0.47
18:CU:66:ASN:HD21	18:CU:70:ARG:NH2	2.12	0.47
34:DA:303:A:H2'	34:DA:304:U:O4'	2.14	0.47
34:DA:438:G:O2'	34:DA:494:U:O4	2.33	0.47
45:DL:84:LEU:HB2	45:DL:105:TYR:CE2	2.50	0.47
45:DL:24:VAL:HG12	45:DL:98:TYR:HE1	1.79	0.47
57:DZ:26:THR:O	57:DZ:30:GLU:HB2	2.14	0.47
1:AA:1099:C:C2	1:AA:1153:G:C2	3.02	0.47
1:AA:2653:G:H5''	1:AA:2653:G:H8	1.80	0.47
1:AA:2864:G:H2'	1:AA:2865:C:C6	2.49	0.47
1:AA:311:C:H2'	1:AA:312:C:C6	2.50	0.47
1:AA:509:A:H2'	1:AA:510:C:H5'	1.95	0.47
7:AG:125:PHE:CE1	7:AG:131:TYR:HB2	2.49	0.47
17:AT:9:LEU:HD23	17:AT:9:LEU:C	2.34	0.47
34:BA:1295:G:O2'	46:BM:14:ARG:NH1	2.47	0.47
34:BA:503:C:H2'	34:BA:504:C:H6	1.78	0.47
34:BA:959:A:O2'	34:BA:984:C:O2'	2.16	0.47
35:BB:141:GLU:O	35:BB:145:LEU:HB2	2.13	0.47
37:BD:53:ASP:O	37:BD:57:ARG:HG3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BE:77:PRO:HD2	38:BE:142:LEU:HD22	1.96	0.47
49:BP:34:GLU:OE2	49:BP:59:TRP:NE1	2.36	0.47
57:BZ:128:TYR:N	57:BZ:128:TYR:CD2	2.81	0.47
27:C3:8:LEU:HD23	27:C3:54:VAL:HG23	1.97	0.47
1:CA:1115:G:H2'	1:CA:1116:C:C6	2.50	0.47
1:CA:1813:G:H2'	1:CA:1814:G:H5'	1.97	0.47
1:CA:2497:A:H5''	63:CA:3708:HOH:O	2.14	0.47
1:CA:2725:A:O2'	1:CA:2726:U:OP2	2.30	0.47
3:CC:68:GLY:N	3:CC:189:ASN:ND2	2.62	0.47
11:CN:53:VAL:HA	11:CN:121:LYS:O	2.13	0.47
18:CU:83:LEU:HG	18:CU:88:ILE:HB	1.96	0.47
20:CW:16:LYS:O	20:CW:19:LEU:HB2	2.14	0.47
34:DA:256:U:OP1	50:DQ:17:LYS:NZ	2.39	0.47
34:DA:380:G:C2	34:DA:384:G:C6	3.02	0.47
34:DA:599:C:H5''	41:DH:95:VAL:O	2.14	0.47
34:DA:883:C:C2'	34:DA:884:U:H5'	2.44	0.47
36:DC:78:GLY:O	36:DC:80:GLY:N	2.45	0.47
44:DK:104:GLN:HG2	44:DK:106:LYS:HG2	1.95	0.47
48:DO:46:HIS:O	48:DO:48:LYS:N	2.48	0.47
57:DZ:272:LEU:HD12	57:DZ:272:LEU:HA	1.46	0.47
1:AA:2207:C:H2'	1:AA:2208:G:C8	2.50	0.47
1:AA:2796:G:H2'	1:AA:2797:C:C6	2.49	0.47
1:AA:455:A:C8	1:AA:455:A:OP2	2.62	0.47
4:AD:130:ALA:C	4:AD:131:LEU:HD12	2.34	0.47
5:AE:9:VAL:HB	17:AT:3:ARG:HG2	1.95	0.47
7:AG:131:TYR:HB3	7:AG:159:VAL:HG13	1.97	0.47
7:AG:66:GLN:HG3	28:A4:1:MET:HE1	1.96	0.47
13:AP:128:HIS:CE1	13:AP:148:LEU:HD21	2.49	0.47
14:AQ:21:THR:CG2	14:AQ:101:ARG:HB2	2.44	0.47
14:AQ:61:GLY:HA3	23:AZ:178:GLU:O	2.15	0.47
16:AS:66:ALA:O	16:AS:69:VAL:HG13	2.14	0.47
34:BA:1392:G:H21	34:BA:1502:A:H8	1.61	0.47
12:AO:48:PRO:HB3	34:BA:1422:G:H5'	1.97	0.47
34:BA:284:G:H2'	34:BA:285:G:C8	2.50	0.47
35:BB:200:ILE:HB	35:BB:202:PRO:HD3	1.96	0.47
35:BB:229:VAL:HG12	35:BB:230:VAL:H	1.79	0.47
36:BC:12:LEU:HD23	36:BC:16:ARG:HB3	1.97	0.47
37:BD:121:VAL:O	37:BD:134:ASP:HA	2.15	0.47
34:BA:923:A:OP1	38:BE:21:ALA:HB2	2.14	0.47
41:BH:6:ILE:O	41:BH:9:MET:N	2.48	0.47
45:BL:111:LYS:O	45:BL:112:ASP:HB2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
48:BO:62:GLN:NE2	48:BO:65:ARG:HH12	2.13	0.47
49:BP:57:ARG:NE	49:BP:79:VAL:O	2.35	0.47
57:BZ:587:SER:OG	57:BZ:588:MET:N	2.48	0.47
1:CA:330:A:H2	1:CA:1210:A:H2'	1.79	0.47
1:CA:2097:C:H2'	1:CA:2098:U:H6	1.78	0.47
1:CA:2203:U:H2'	1:CA:2205:C:C6	2.49	0.47
1:CA:2306:C:H3'	1:CA:2307:G:H2'	1.96	0.47
1:CA:2328:A:H2'	1:CA:2329:G:C8	2.50	0.47
1:CA:2542:A:H4'	1:CA:2543:G:C8	2.49	0.47
3:CC:6:LYS:HA	3:CC:9:ARG:NH1	2.30	0.47
8:CH:17:VAL:HG11	8:CH:50:VAL:HG21	1.95	0.47
34:DA:1286:A:H8	34:DA:1287:A:H4'	1.79	0.47
34:DA:503:C:H2'	34:DA:504:C:H6	1.79	0.47
34:DA:693:G:H2'	34:DA:694:A:C8	2.50	0.47
35:DB:200:ILE:O	35:DB:202:PRO:HD3	2.14	0.47
34:DA:1307:U:OP1	46:DM:101:GLN:NE2	2.47	0.47
46:DM:3:ARG:N	46:DM:7:VAL:O	2.48	0.47
57:DZ:152:THR:C	57:DZ:154:GLN:H	2.18	0.47
57:DZ:505:GLY:HA2	57:DZ:576:ASP:CB	2.45	0.47
57:DZ:527:ASN:O	57:DZ:529:ILE:N	2.48	0.47
1:AA:1310:G:OP1	29:A5:19:ARG:NH2	2.35	0.47
1:AA:919:A:H2'	1:AA:920:G:O4'	2.14	0.47
1:AA:895:G:N9	1:AA:978:A:H8	2.13	0.47
2:AB:17:C:H2'	2:AB:18:G:O4'	2.15	0.47
3:AC:180:SER:O	3:AC:181:PHE:O	2.33	0.47
16:AS:93:LYS:HG2	16:AS:95:HIS:HB2	1.97	0.47
34:BA:475:G:H2'	34:BA:476:G:H8	1.79	0.47
34:BA:567:G:O6	45:BL:15:ARG:NH1	2.40	0.47
43:BJ:64:GLU:HB3	47:BN:59:ALA:HB2	1.95	0.47
57:BZ:249:GLY:HA2	57:BZ:252:ASP:HB2	1.97	0.47
29:C5:16:ARG:HG2	29:C5:16:ARG:NH1	2.22	0.47
1:CA:1366:A:H2'	1:CA:1367:A:O4'	2.15	0.47
1:CA:2006:C:OP2	63:CA:4474:HOH:O	2.20	0.47
1:CA:2093:G:H1	1:CA:2196:C:H42	1.63	0.47
1:CA:580:C:H2'	1:CA:581:C:H6	1.78	0.47
1:CA:687:C:H2'	1:CA:688:U:O4'	2.14	0.47
5:CE:12:THR:HG22	17:CT:58:ASN:HD21	1.79	0.47
5:CE:15:PHE:CD2	17:CT:81:PRO:HD2	2.50	0.47
10:CL:93:ARG:HB3	10:CL:93:ARG:HE	1.34	0.47
23:CZ:157:LEU:HD11	23:CZ:163:LEU:HB2	1.97	0.47
34:DA:177:C:H2'	34:DA:178:C:C6	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:324:G:N2	34:DA:327:A:C8	2.82	0.47
38:DE:40:ARG:HB3	38:DE:66:MET:CE	2.45	0.47
45:DL:110:VAL:HG23	45:DL:120:TYR:O	2.15	0.47
50:DQ:62:SER:CB	50:DQ:72:ARG:HD3	2.45	0.47
56:DW:44:G:O2'	56:DW:45:U:H5'	2.14	0.47
56:DY:7:A:N6	56:DY:66:U:H3	2.02	0.47
57:DZ:213:HIS:C	57:DZ:213:HIS:HD1	2.18	0.47
57:DZ:79:ILE:HD13	57:DZ:276:VAL:HG11	1.95	0.47
1:AA:1692:G:H5''	1:AA:1693:C:H5'	1.96	0.47
1:AA:1888:G:C2'	1:AA:1889:G:H5'	2.44	0.47
1:AA:2105:G:O5'	1:AA:2105:G:H8	1.98	0.47
1:AA:2760:G:OP1	8:AH:138:LYS:NZ	2.39	0.47
1:AA:2859:U:O4	17:AT:23:ARG:NH2	2.33	0.47
1:AA:2859:U:N3	1:AA:2877:G:O4'	2.48	0.47
1:AA:417:A:C6	13:AP:71:VAL:HG11	2.50	0.47
2:AB:32:C:C2	2:AB:51:G:N2	2.83	0.47
4:AD:69:ARG:C	4:AD:71:ASP:H	2.18	0.47
7:AG:77:ILE:N	7:AG:82:LEU:O	2.32	0.47
22:AY:55:TYR:N	22:AY:55:TYR:CD1	2.82	0.47
34:BA:1151:A:O2'	34:BA:1152:A:H8	1.98	0.47
34:BA:1318:A:H5''	52:BS:3:ARG:NH2	2.30	0.47
34:BA:1318:A:OP1	52:BS:3:ARG:NH2	2.43	0.47
34:BA:255:G:H2'	34:BA:256:U:C6	2.50	0.47
34:BA:376:G:C2	34:BA:389:A:C2	3.03	0.47
34:BA:653:A:OP1	41:BH:56:LYS:NZ	2.48	0.47
34:BA:741:G:H2'	34:BA:742:G:O4'	2.15	0.47
34:BA:964:A:N3	34:BA:969:A:O2'	2.35	0.47
38:BE:102:ALA:H	38:BE:107:ARG:HH21	1.62	0.47
41:BH:116:LYS:O	41:BH:119:LEU:HD21	2.15	0.47
45:BL:75:HIS:ND1	45:BL:77:LEU:HB2	2.30	0.47
51:BR:31:LEU:HD23	51:BR:31:LEU:H	1.78	0.47
56:BW:17:C:H5'	56:BW:18:G:OP1	2.14	0.47
57:BZ:138:LYS:HG2	62:BZ:704:GDP:C2	2.49	0.47
57:BZ:590:ILE:HA	57:BZ:590:ILE:HD13	1.76	0.47
1:CA:1832:C:H2'	1:CA:1833:U:O4'	2.14	0.47
1:CA:1860:G:H5''	3:CC:207:GLY:N	2.27	0.47
1:CA:2684:U:C4	1:CA:2685:G:N7	2.82	0.47
1:CA:652(E):G:O5'	1:CA:652(E):G:H8	1.98	0.47
1:CA:794:G:H2'	1:CA:795:C:H6	1.80	0.47
1:CA:952:G:C6	1:CA:966:G:C6	3.03	0.47
4:CD:126:GLN:O	4:CD:193:VAL:HG22	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:CR:9:LYS:HA	15:CR:17:ARG:HE	1.80	0.47
20:CW:70:TYR:O	20:CW:107:LEU:HA	2.15	0.47
21:CX:84:ALA:HB3	21:CX:87:GLN:OE1	2.14	0.47
35:DB:63:MET:HG2	35:DB:225:ALA:HB1	1.97	0.47
37:DD:135:LEU:C	37:DD:137:SER:H	2.13	0.47
37:DD:31:CYS:SG	37:DD:33:MET:N	2.87	0.47
39:DF:33:TYR:HD1	39:DF:75:LEU:HD23	1.79	0.47
39:DF:78:GLU:O	39:DF:80:ARG:N	2.48	0.47
40:DG:111:ARG:HD2	40:DG:123:GLU:HB2	1.95	0.47
44:DK:14:VAL:HG12	44:DK:15:ALA:H	1.80	0.47
50:DQ:32:TYR:C	50:DQ:34:LYS:H	2.17	0.47
57:DZ:97:SER:HA	57:DZ:100:VAL:HG12	1.96	0.47
57:DZ:168:ILE:HG23	57:DZ:205:TYR:CE2	2.48	0.47
57:DZ:213:HIS:ND1	57:DZ:213:HIS:O	2.39	0.47
57:DZ:-58:LEU:HD13	57:DZ:-58:LEU:HA	1.65	0.47
29:A5:35:GLU:HG3	29:A5:51:TYR:CG	2.50	0.47
1:AA:1158:G:H2'	1:AA:1159:U:O4'	2.14	0.47
1:AA:1314:A:C2	1:AA:2035:A:C4	3.03	0.47
1:AA:1825:U:O2'	1:AA:1826:C:H5'	2.14	0.47
1:AA:2583:C:H5''	1:AA:2584:A:H5''	1.96	0.47
3:AC:223:VAL:HG23	3:AC:223:VAL:O	2.15	0.47
11:AN:14:VAL:HG12	11:AN:52:VAL:HA	1.96	0.47
34:BA:111:G:O5'	34:BA:111:G:H8	1.97	0.47
34:BA:437:U:O2'	37:BD:125:HIS:CE1	2.67	0.47
34:BA:8:A:N7	37:BD:208:SER:OG	2.44	0.47
36:BC:66:VAL:HG12	36:BC:68:VAL:HG23	1.97	0.47
39:BF:10:LEU:HD21	39:BF:26:ILE:HD11	1.97	0.47
41:BH:56:LYS:HB2	41:BH:58:TYR:HE1	1.78	0.47
49:BP:74:LEU:HG	49:BP:79:VAL:HG21	1.97	0.47
57:BZ:177:ILE:HD12	57:BZ:188:TYR:HE2	1.79	0.47
26:C2:10:LEU:O	26:C2:14:ARG:HB2	2.15	0.47
32:C8:20:GLY:O	32:C8:59:LYS:HE2	2.15	0.47
1:CA:819:A:C4	1:CA:1189:A:C2	3.03	0.47
1:CA:300:A:P	22:CY:86:ARG:NH2	2.88	0.47
1:CA:453:C:OP1	63:CA:4580:HOH:O	2.20	0.47
1:CA:729:G:OP2	4:CD:13:ARG:NH1	2.32	0.47
1:CA:774:A:O2'	1:CA:775:G:H8	1.97	0.47
2:CB:20:C:H42	2:CB:63:G:H1	1.63	0.47
3:CC:15:VAL:O	3:CC:21:TYR:OH	2.29	0.47
1:CA:1843:C:H5'	4:CD:253:GLN:HE21	1.79	0.47
16:CS:53:SER:O	16:CS:57:LYS:N	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:232:G:H1'	34:DA:262:A:N1	2.30	0.47
38:DE:152:ARG:HG2	41:DH:43:GLY:HA3	1.96	0.47
42:DI:8:GLY:HA3	42:DI:76:ALA:O	2.15	0.47
48:DO:41:GLU:O	48:DO:44:LYS:HB2	2.15	0.47
57:DZ:162:VAL:H	57:DZ:258:VAL:HG23	1.78	0.47
61:DZ:703:FUA:H16	61:DZ:703:FUA:H322	1.74	0.47
1:AA:1026:A:N3	1:AA:2059:G:O2'	2.39	0.47
1:AA:2184:G:H4'	1:AA:2194:U:H2'	1.96	0.47
1:AA:2137:G:H21	1:AA:2193:A:H61	1.62	0.47
1:AA:2650:G:P	5:AE:82:ARG:NH2	2.88	0.47
1:AA:821:A:N3	1:AA:821:A:H2'	2.30	0.47
3:AC:42:VAL:CG1	3:AC:43:GLU:H	2.27	0.47
3:AC:46:ALA:O	3:AC:47:LYS:HB2	2.15	0.47
1:AA:2679:C:H1'	8:AH:109:PHE:HD1	1.79	0.47
22:AY:1:MET:HB2	22:AY:2:ARG:H	1.59	0.47
23:AZ:180:VAL:O	23:AZ:183:LEU:HB2	2.15	0.47
34:BA:1157:A:C6	34:BA:1180:A:C6	3.03	0.47
34:BA:363:A:N7	45:BL:30:ALA:HB1	2.29	0.47
34:BA:977:A:H1'	34:BA:982:U:O4	2.15	0.47
36:BC:22:TRP:CE2	47:BN:54:PRO:HG3	2.49	0.47
37:BD:114:ARG:HA	37:BD:117:ALA:HB3	1.96	0.47
45:BL:26:ALA:HB3	45:BL:98:TYR:OH	2.14	0.47
51:BR:73:ALA:HB3	51:BR:79:LEU:HD12	1.96	0.47
57:BZ:120:THR:HG22	57:BZ:123:ARG:HH12	1.78	0.47
57:BZ:389:LEU:HA	57:BZ:389:LEU:HD12	1.61	0.47
1:CA:1024:G:HO2'	1:CA:1144:G:HO2'	1.58	0.47
1:CA:211:A:H2'	1:CA:212:G:O4'	2.14	0.47
1:CA:2711:A:H1'	63:CA:4654:HOH:O	2.14	0.47
1:CA:325:G:O2'	1:CA:326:G:H5'	2.14	0.47
3:CC:180:SER:O	3:CC:181:PHE:O	2.33	0.47
3:CC:42:VAL:CG1	3:CC:43:GLU:H	2.28	0.47
4:CD:137:PRO:O	4:CD:140:THR:OG1	2.29	0.47
4:CD:85:ASP:OD2	4:CD:88:ARG:NH1	2.46	0.47
5:CE:181:LEU:HD12	5:CE:181:LEU:HA	1.67	0.47
6:CF:21:ALA:CB	6:CF:22:ALA:HA	2.45	0.47
9:CK:70:GLU:O	9:CK:72:ASP:N	2.48	0.47
19:CV:55:ALA:HA	19:CV:100:ARG:O	2.15	0.47
23:CZ:28:MET:HA	23:CZ:88:PHE:O	2.15	0.47
23:CZ:8:TYR:HB2	23:CZ:38:TYR:CE2	2.49	0.47
34:DA:1014:A:N3	34:DA:1219:U:O2'	2.38	0.47
34:DA:181:G:N1	34:DA:195:A:C8	2.83	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:457:C:H2'	34:DA:458:C:C6	2.43	0.47
34:DA:814:A:H2'	34:DA:816:A:H5''	1.97	0.47
43:DJ:78:ASN:O	43:DJ:80:LYS:N	2.47	0.47
50:DQ:74:LEU:HD23	50:DQ:74:LEU:HA	1.72	0.47
56:DW:41:C:H2'	56:DW:42:C:O4'	2.15	0.47
57:DZ:21:ILE:CD1	57:DZ:87:HIS:HA	2.44	0.47
30:A6:40:CYS:SG	30:A6:42:TRP:HB2	2.54	0.47
1:AA:1073:A:C2	1:AA:2500:A:H5'	2.50	0.47
1:AA:1841:A:C2'	1:AA:1842:G:H5'	2.45	0.47
1:AA:1911:A:N1	1:AA:2246:G:H1'	2.29	0.47
1:AA:2021:C:H4'	1:AA:2736:C:O2	2.14	0.47
1:AA:2175:G:H2'	1:AA:2176:G:C8	2.48	0.47
1:AA:2439:C:H5''	1:AA:2440:G:OP1	2.14	0.47
1:AA:2199:C:O2'	3:AC:46:ALA:O	2.31	0.47
5:AE:47:VAL:HG22	5:AE:84:PHE:O	2.15	0.47
6:AF:18:ARG:NH2	6:AF:127:GLU:OE1	2.48	0.47
10:AL:73:PRO:O	10:AL:77:LEU:HG	2.15	0.47
11:AN:134:ARG:N	11:AN:135:PRO:HD3	2.30	0.47
18:AU:104:GLN:H	18:AU:104:GLN:CD	2.17	0.47
23:AZ:150:LEU:HD12	23:AZ:150:LEU:HA	1.75	0.47
34:BA:1112:C:N3	36:BC:178:LEU:N	2.51	0.47
34:BA:1119:C:OP1	42:BI:83:ARG:NH1	2.42	0.47
34:BA:216:G:H2'	34:BA:217:C:C6	2.50	0.47
35:BB:83:MET:HB2	35:BB:84:GLU:OE2	2.14	0.47
34:BA:826:C:H5'	41:BH:12:ARG:NH1	2.30	0.47
41:BH:4:ASP:OD1	41:BH:7:ALA:N	2.48	0.47
56:BW:61:C:H2'	56:BW:62:C:H6	1.79	0.47
1:CA:2112:G:H2'	1:CA:2113:U:H4'	1.97	0.47
1:CA:2340:G:H2'	1:CA:2341:G:H8	1.80	0.47
1:CA:2849:U:H4'	1:CA:2868:A:C2	2.49	0.47
1:CA:530:G:C5	1:CA:2022:U:H5''	2.50	0.47
1:CA:868:U:C4	1:CA:869:G:N7	2.83	0.47
2:CB:22:U:H3	2:CB:61:G:H1	1.61	0.47
14:CQ:76:LYS:HG2	14:CQ:77:LYS:O	2.15	0.47
23:CZ:101:PRO:HA	23:CZ:123:ASP:HA	1.96	0.47
34:DA:1312:G:N7	52:DS:2:PRO:HD2	2.30	0.47
34:DA:1442(A):G:H3'	34:DA:1442(B):A:H8	1.80	0.47
34:DA:300:A:H1'	34:DA:565:U:O2	2.15	0.47
34:DA:60:A:H4'	34:DA:61:G:O5'	2.15	0.47
38:DE:82:VAL:HG11	38:DE:137:GLU:HB3	1.97	0.47
41:DH:63:LEU:HD23	41:DH:65:TYR:OH	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:DI:71:SER:HA	42:DI:74:ILE:HD12	1.95	0.47
48:DO:26:GLU:H	48:DO:26:GLU:HG2	1.43	0.47
52:DS:33:THR:HG21	52:DS:71:LEU:HD21	1.97	0.47
52:DS:40:ILE:HB	52:DS:67:VAL:O	2.15	0.47
56:DY:19:G:H1	56:DY:56:C:H42	1.62	0.47
57:DZ:-3:GLU:HG3	57:DZ:0:ARG:HH11	1.80	0.47
57:DZ:630:GLN:O	57:DZ:645:ALA:HB1	2.15	0.47
10:CL:21:PRO:HD2	57:DZ:641:GLN:HE22	1.79	0.47
1:AA:1038:C:O2'	1:AA:1039:G:H5'	2.14	0.46
1:AA:2871:G:O2'	1:AA:2872:G:H5'	2.15	0.46
2:AB:94:C:H2'	2:AB:95:C:H6	1.79	0.46
3:AC:17:PRO:HG2	3:AC:18:ASN:H	1.79	0.46
3:AC:6:LYS:HA	3:AC:9:ARG:NH1	2.30	0.46
8:AH:41:MET:CE	8:AH:65:HIS:HA	2.40	0.46
8:AH:3:ARG:HG2	8:AH:6:ARG:CG	2.45	0.46
9:AK:53:VAL:O	9:AK:85:ASP:HA	2.15	0.46
13:AP:39:LYS:HG3	13:AP:45:LEU:CD1	2.44	0.46
19:AV:21:ARG:HG3	19:AV:91:TYR:CD1	2.50	0.46
34:BA:1411:C:H2'	34:BA:1412:C:C6	2.50	0.46
34:BA:406:G:N3	37:BD:119:GLN:NE2	2.63	0.46
34:BA:913:A:H4'	34:BA:914:A:O5'	2.15	0.46
35:BB:36:ARG:O	35:BB:38:GLY:N	2.38	0.46
48:BO:18:PHE:HD1	48:BO:20:GLY:H	1.62	0.46
52:BS:40:ILE:O	52:BS:67:VAL:HG13	2.15	0.46
1:CA:284:U:H2'	1:CA:285:C:H6	1.79	0.46
1:CA:442:G:H21	6:CF:48:THR:HB	1.79	0.46
1:CA:684:G:OP1	31:C7:16:HIS:ND1	2.46	0.46
4:CD:68:LYS:HD2	4:CD:70:TRP:CH2	2.49	0.46
7:CG:173:LEU:HD22	7:CG:178:PHE:CE1	2.50	0.46
7:CG:43:LEU:HB3	7:CG:44:GLY:H	1.58	0.46
7:CG:76:SER:CB	7:CG:84:LYS:H	2.28	0.46
16:CS:29:PHE:HD1	16:CS:30:ARG:N	2.13	0.46
34:DA:1107:C:C4	34:DA:1108:G:C8	3.04	0.46
34:DA:983:A:N1	34:DA:1222:G:N2	2.64	0.46
34:DA:1409:C:H2'	34:DA:1410:G:C8	2.50	0.46
34:DA:1478:C:H2'	34:DA:1479:C:H6	1.80	0.46
34:DA:17:U:H5''	38:DE:14:ARG:NH1	2.30	0.46
34:DA:20:U:H2'	34:DA:21:G:O4'	2.15	0.46
34:DA:622:A:H2'	34:DA:623:C:H5'	1.96	0.46
34:DA:767:A:H2'	34:DA:768:A:O4'	2.15	0.46
34:DA:986:A:N3	52:DS:52:TYR:OH	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
35:DB:71:VAL:HG23	35:DB:164:VAL:HG13	1.98	0.46
36:DC:54:ARG:HB3	36:DC:54:ARG:NH1	2.30	0.46
37:DD:22:LYS:O	37:DD:113:SER:HB3	2.15	0.46
38:DE:18:ARG:HH21	38:DE:25:ARG:CG	2.27	0.46
38:DE:78:HIS:HA	41:DH:105:ARG:HG3	1.97	0.46
52:DS:41:VAL:O	52:DS:44:MET:N	2.42	0.46
53:DT:79:ARG:HD2	53:DT:83:ARG:NH1	2.30	0.46
56:DW:19:G:N2	56:DW:57:G:H1'	2.31	0.46
56:DY:30:G:H2'	56:DY:31:A:H8	1.79	0.46
56:DY:9:A:H5'	56:DY:46:7MG:H1'	1.98	0.46
57:DZ:181:LEU:HD12	57:DZ:216:LEU:HD11	1.95	0.46
57:DZ:92:ILE:HG21	57:DZ:437:THR:HG21	1.97	0.46
57:DZ:539:ILE:HA	57:DZ:542:VAL:HG12	1.98	0.46
1:AA:2399:U:OP1	24:A0:55:ARG:NH2	2.48	0.46
1:AA:1470:G:H2'	1:AA:1471:G:O4'	2.15	0.46
1:AA:1938:A:H2'	1:AA:1939:U:O4'	2.16	0.46
1:AA:2760:G:C2	1:AA:2769:U:C5	3.03	0.46
1:AA:495:G:O6	31:A7:37:LYS:HE2	2.16	0.46
7:AG:60:LEU:HA	7:AG:63:ILE:HD12	1.97	0.46
21:AX:53:LYS:HB3	21:AX:82:GLN:HB3	1.97	0.46
22:AY:9:LYS:HE3	22:AY:28:LYS:O	2.15	0.46
34:BA:944:G:O2'	34:BA:1339:A:N6	2.44	0.46
34:BA:1456:G:H1'	53:BT:39:LYS:NZ	2.30	0.46
34:BA:414:A:C5	34:BA:431:A:C2	3.04	0.46
34:BA:753:A:H4'	34:BA:754:C:H5''	1.97	0.46
53:BT:57:ARG:HH12	53:BT:100:ILE:HG13	1.80	0.46
56:BW:7:A:H5''	56:BW:7:A:C8	2.50	0.46
1:CA:1021:A:C8	1:CA:1021:A:H3'	2.50	0.46
1:CA:1220:A:OP2	18:CU:19:LYS:NZ	2.43	0.46
1:CA:729:G:C4	1:CA:1775:U:O2	2.68	0.46
1:CA:2103:C:H1'	1:CA:2187:G:H22	1.79	0.46
1:CA:225:A:N6	1:CA:226:G:C2	2.84	0.46
1:CA:736:C:H2'	1:CA:737:C:C6	2.50	0.46
1:CA:912:C:H2'	1:CA:913:U:H6	1.79	0.46
1:CA:2277:G:H5''	14:CQ:87:LYS:HB3	1.97	0.46
34:DA:1241:G:H2'	34:DA:1242:C:C6	2.50	0.46
34:DA:1326:C:H5''	54:DU:18:TYR:O	2.15	0.46
35:DB:121:LEU:O	35:DB:127:ILE:HB	2.16	0.46
35:DB:145:LEU:O	35:DB:149:LEU:HB2	2.14	0.46
40:DG:78:ARG:HG2	40:DG:79:ARG:H	1.80	0.46
34:DA:1126:U:H3	43:DJ:40:LEU:HD11	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:DM:60:VAL:HG23	46:DM:64:TRP:CE3	2.50	0.46
50:DQ:24:GLU:HB3	50:DQ:39:SER:HB3	1.97	0.46
52:DS:15:LEU:HG	52:DS:18:LYS:HD3	1.96	0.46
55:DV:17:U:C2	56:DW:36:A:C2	3.03	0.46
57:DZ:2:LYS:O	57:DZ:6:GLU:HB2	2.15	0.46
57:DZ:637:ARG:O	57:DZ:639:ASN:N	2.47	0.46
1:AA:1501:U:O2'	1:AA:1502:G:N7	2.48	0.46
1:AA:1480:A:N6	1:AA:1605:A:N6	2.61	0.46
1:AA:2074:G:H4'	5:AE:143:ASN:O	2.15	0.46
1:AA:2149:G:H2'	1:AA:2150:C:C6	2.51	0.46
1:AA:2178:G:H2'	1:AA:2179:G:C4	2.51	0.46
1:AA:2729:U:O2'	1:AA:2730:G:H5'	2.15	0.46
1:AA:941:U:H5'	1:AA:942:A:OP2	2.16	0.46
2:AB:43:C:OP1	28:A4:6:HIS:NE2	2.42	0.46
4:AD:70:TRP:HB3	4:AD:190:TYR:CZ	2.50	0.46
18:AU:65:ILE:HD11	18:AU:95:LEU:HB3	1.97	0.46
34:BA:1427:U:H2'	34:BA:1428:A:C8	2.50	0.46
34:BA:1429:C:H2'	34:BA:1430:C:C6	2.50	0.46
34:BA:354:G:H2'	34:BA:355:C:H5'	1.98	0.46
34:BA:431:A:H2'	34:BA:432:A:O4'	2.15	0.46
34:BA:763:G:H2'	34:BA:764:C:H6	1.79	0.46
37:BD:101:LEU:O	37:BD:104:VAL:N	2.48	0.46
37:BD:8:VAL:HG22	37:BD:21:LEU:HD13	1.98	0.46
38:BE:28:PHE:CD2	38:BE:51:VAL:HG22	2.51	0.46
34:BA:684:A:O2'	44:BK:39:PRO:O	2.33	0.46
44:BK:43:SER:OG	44:BK:44:SER:N	2.49	0.46
46:BM:80:ARG:NH2	52:BS:69:HIS:HE1	2.13	0.46
56:BY:28:G:H2'	56:BY:29:G:H8	1.80	0.46
1:CA:1225:G:C2	1:CA:1226:A:C2	3.03	0.46
1:CA:1762:A:N1	63:CA:4249:HOH:O	2.36	0.46
1:CA:1842:G:O3'	4:CD:253:GLN:NE2	2.43	0.46
1:CA:2577:A:O2'	29:C5:2:ALA:HB1	2.15	0.46
4:CD:89:SER:HB2	4:CD:159:ALA:CB	2.45	0.46
5:CE:13:ARG:O	17:CT:57:PHE:HE2	1.98	0.46
34:DA:920:U:H2'	34:DA:921:U:H6	1.79	0.46
36:DC:186:PHE:HE2	36:DC:188:LEU:HB2	1.80	0.46
37:DD:83:SER:HA	37:DD:89:THR:OG1	2.15	0.46
42:DI:47:LEU:HB3	42:DI:50:LEU:HD12	1.96	0.46
45:DL:78:GLN:HG3	45:DL:79:GLU:O	2.16	0.46
51:DR:33:ASP:CG	51:DR:36:ASN:HB2	2.34	0.46
57:DZ:329:ARG:HD3	57:DZ:331:TYR:CZ	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:499:ARG:HB2	57:DZ:506:GLN:HB3	1.97	0.46
57:DZ:639:ASN:HA	57:DZ:640:ALA:O	2.15	0.46
57:DZ:94:VAL:HG21	57:DZ:121:VAL:HG23	1.96	0.46
57:DZ:-9:LEU:O	57:DZ:-6:ARG:HB2	2.16	0.46
25:A1:64:ALA:HA	25:A1:67:ILE:HG13	1.97	0.46
1:AA:1764:G:C2'	1:AA:1765:U:H5'	2.45	0.46
1:AA:2149:G:N2	1:AA:2195:A:H1'	2.30	0.46
1:AA:2198:A:O2'	3:AC:45:HIS:CG	2.68	0.46
1:AA:2594:G:C2	1:AA:2595:G:C8	3.04	0.46
11:AN:91:LEU:HA	11:AN:91:LEU:HD23	1.81	0.46
20:AW:71:VAL:HA	20:AW:107:LEU:HD12	1.97	0.46
34:BA:1226:C:O2'	46:BM:111:LYS:NZ	2.48	0.46
34:BA:1346:A:N1	34:BA:1374:A:H5''	2.30	0.46
34:BA:1412:C:H2'	34:BA:1413:A:C8	2.51	0.46
34:BA:453:A:C6	34:BA:454:C:C4	3.04	0.46
34:BA:590:C:H2'	34:BA:591:U:H6	1.79	0.46
35:BB:60:ASP:OD2	35:BB:64:ARG:NH2	2.48	0.46
38:BE:91:LEU:HD12	38:BE:120:THR:HG22	1.97	0.46
42:BI:100:GLY:O	42:BI:103:THR:HG22	2.15	0.46
44:BK:20:TYR:O	44:BK:30:VAL:HA	2.15	0.46
44:BK:70:LYS:HA	44:BK:73:MET:HE3	1.96	0.46
57:BZ:169:GLY:HA3	57:BZ:174:PHE:HA	1.97	0.46
1:CA:336:C:OP1	22:CY:84:ARG:HG2	2.15	0.46
1:CA:99:U:O4	22:CY:8:LYS:NZ	2.28	0.46
3:CC:30:VAL:CG2	3:CC:31:LYS:N	2.78	0.46
4:CD:124:PRO:HD2	4:CD:129:ASN:ND2	2.30	0.46
5:CE:46:ALA:HB2	5:CE:82:ARG:HA	1.97	0.46
8:CH:124:GLU:HB2	8:CH:132:ARG:HB3	1.97	0.46
8:CH:35:VAL:HA	8:CH:36:PRO:HD2	1.60	0.46
10:CL:10:LEU:N	10:CL:55:VAL:O	2.48	0.46
13:CP:96:THR:OG1	13:CP:98:GLU:HG2	2.15	0.46
14:CQ:104:PHE:O	14:CQ:105:GLU:HG3	2.14	0.46
34:DA:396:G:H2'	34:DA:397:A:H5''	1.97	0.46
34:DA:743:U:H2'	34:DA:744:C:C6	2.50	0.46
34:DA:782:A:O3'	34:DA:1515:C:H4'	2.16	0.46
34:DA:834:C:H2'	34:DA:835:U:H6	1.80	0.46
35:DB:127:ILE:C	35:DB:129:GLU:H	2.18	0.46
35:DB:29:ALA:HA	35:DB:32:ILE:HD12	1.97	0.46
46:DM:3:ARG:NH1	46:DM:4:ILE:HB	2.31	0.46
50:DQ:66:SER:OG	50:DQ:69:LYS:HB2	2.15	0.46
52:DS:3:ARG:NH2	52:DS:10:PHE:HB2	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:435:ASP:O	57:DZ:438:PHE:HD1	1.98	0.46
27:A3:32:GLN:HA	27:A3:32:GLN:NE2	2.29	0.46
7:AG:105:LYS:HG2	28:A4:24:THR:HG21	1.96	0.46
1:AA:1120:G:H2'	1:AA:1121:C:C6	2.51	0.46
1:AA:1153:G:H2'	1:AA:1154:U:O2	2.15	0.46
1:AA:1466:U:HO2'	1:AA:1467:G:P	2.36	0.46
1:AA:1817:A:H8	63:AA:5177:HOH:O	1.98	0.46
1:AA:1934:A:N1	34:BA:1407:C:O2'	2.40	0.46
8:AH:3:ARG:HG2	8:AH:6:ARG:CD	2.45	0.46
10:AL:95:LYS:HG2	10:AL:137:GLU:HB3	1.97	0.46
18:AU:108:GLU:OE2	18:AU:112:ARG:NH1	2.44	0.46
18:AU:24:TYR:HB3	18:AU:28:ARG:HB3	1.97	0.46
1:AA:1387:U:O2	21:AX:80:ILE:HD12	2.15	0.46
34:BA:669:U:C2	34:BA:670:G:C8	3.03	0.46
35:BB:64:ARG:HE	35:BB:64:ARG:HB2	1.29	0.46
38:BE:37:ARG:HH12	38:BE:111:GLU:HG2	1.81	0.46
38:BE:144:THR:OG1	38:BE:147:ASP:OD1	2.27	0.46
42:BI:9:ARG:HG2	42:BI:14:VAL:HG22	1.97	0.46
33:C9:22:ARG:HB2	33:C9:24:TYR:CE1	2.50	0.46
1:CA:1633:G:C5	1:CA:1635:G:C5	3.03	0.46
1:CA:2024:G:O2'	1:CA:2025:C:H5'	2.16	0.46
1:CA:20:C:H2'	1:CA:21:A:H8	1.81	0.46
1:CA:910:A:C5	14:CQ:13:GLN:HG3	2.50	0.46
3:CC:176:VAL:HG12	3:CC:176:VAL:O	2.15	0.46
3:CC:20:VAL:O	3:CC:224:ARG:O	2.33	0.46
12:CO:111:PHE:O	12:CO:115:VAL:HG23	2.16	0.46
2:CB:114:C:H4'	16:CS:46:VAL:HG22	1.97	0.46
34:DA:1039:C:H2'	34:DA:1040:U:O4'	2.15	0.46
34:DA:1239:A:H62	34:DA:1299:A:N6	2.13	0.46
12:CO:97:ARG:NH1	34:DA:339:C:OP2	2.44	0.46
34:DA:954:G:H21	34:DA:1227:A:H62	1.63	0.46
37:DD:117:ALA:HA	37:DD:120:LEU:HB2	1.97	0.46
50:DQ:85:VAL:O	50:DQ:89:LEU:HB2	2.16	0.46
57:DZ:106:VAL:HG23	57:DZ:132:ARG:HB2	1.97	0.46
57:DZ:36:THR:HG22	57:DZ:74:TRP:HA	1.97	0.46
26:A2:22:GLU:OE2	26:A2:68:ARG:NH2	2.49	0.46
21:AX:10:ALA:HA	26:A2:37:PHE:CE1	2.51	0.46
29:A5:16:ARG:HG2	29:A5:17:ASP:OD1	2.15	0.46
1:AA:1068:G:C5	1:AA:1185:C:C4	3.03	0.46
1:AA:1324:A:OP1	15:AR:36:THR:CG2	2.63	0.46
1:AA:1370:G:C4	1:AA:1374:G:O6	2.68	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1884:A:O2'	1:AA:1885:A:H5'	2.15	0.46
1:AA:1932:G:C2	1:AA:1943:G:C4	3.04	0.46
1:AA:606:G:OP2	18:AU:10:ARG:HD2	2.15	0.46
34:BA:1323:G:H4'	34:BA:1363:C:N3	2.30	0.46
34:BA:965:A:C2	34:BA:969:A:C2	3.04	0.46
35:BB:37:ASN:OD1	35:BB:37:ASN:N	2.48	0.46
35:BB:78:GLN:O	35:BB:94:ASN:ND2	2.37	0.46
49:BP:71:ARG:O	49:BP:75:ARG:N	2.35	0.46
1:CA:1432:C:H2'	1:CA:1433:U:O4'	2.16	0.46
1:CA:2887:U:H2'	1:CA:2888:C:C6	2.50	0.46
3:CC:223:VAL:HG23	3:CC:223:VAL:O	2.15	0.46
4:CD:221:VAL:HG22	4:CD:226:MET:HE3	1.96	0.46
7:CG:107:LEU:HD11	7:CG:178:PHE:CE1	2.51	0.46
8:CH:13:LYS:HA	8:CH:14:GLY:HA2	1.61	0.46
14:CQ:29:PHE:HB3	14:CQ:65:PHE:CD2	2.50	0.46
14:CQ:72:LYS:HA	14:CQ:73:PRO:HD3	1.84	0.46
22:CY:31:LEU:HA	22:CY:31:LEU:HD23	1.79	0.46
34:DA:1078:U:H1'	38:DE:130:ASN:HD21	1.81	0.46
34:DA:1086:U:H3	34:DA:1099:G:N2	2.06	0.46
34:DA:113:G:N3	34:DA:353:A:O2'	2.44	0.46
34:DA:1307:U:H2'	34:DA:1308:U:C6	2.51	0.46
17:CT:41:ARG:CZ	34:DA:345:C:H3'	2.45	0.46
34:DA:530:G:O2'	34:DA:531:U:OP1	2.27	0.46
45:DL:35:GLY:HA2	45:DL:60:LEU:HA	1.97	0.46
50:DQ:6:LEU:HB3	50:DQ:23:VAL:HG21	1.97	0.46
57:DZ:619:ASP:HB3	57:DZ:662:LYS:CD	2.44	0.46
1:AA:1248:G:OP2	1:AA:1249:A:O2'	2.28	0.46
1:AA:2078:G:C2	1:AA:2079:A:C8	3.04	0.46
1:AA:2087:C:H2'	1:AA:2088:C:C6	2.51	0.46
1:AA:2326:C:C2	1:AA:2327:G:C8	3.03	0.46
1:AA:733:G:H1	31:A7:16:HIS:CD2	2.34	0.46
1:AA:753:A:H2'	1:AA:754:G:O4'	2.16	0.46
5:AE:120:TRP:CE2	5:AE:155:LYS:HG2	2.51	0.46
7:AG:18:GLU:O	7:AG:21:ARG:HB2	2.16	0.46
9:AK:74:LEU:O	9:AK:76:GLY:N	2.39	0.46
17:AT:105:LEU:HA	17:AT:105:LEU:HD23	1.62	0.46
34:BA:1437:C:H2'	34:BA:1438:G:C8	2.50	0.46
34:BA:502:G:C6	34:BA:503:C:C4	3.03	0.46
34:BA:812:C:OP1	34:BA:903:G:H1'	2.15	0.46
34:BA:1191:A:H5''	36:BC:4:LYS:HZ2	1.80	0.46
57:BZ:166:LEU:HA	57:BZ:166:LEU:HD12	1.71	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:BZ:309:LEU:CA	57:BZ:333:GLY:HA3	2.45	0.46
61:BZ:703:FUA:C20	61:BZ:703:FUA:O1	2.62	0.46
57:BZ:75:LYS:HD2	57:BZ:75:LYS:HA	1.64	0.46
33:C9:17:ILE:HG23	33:C9:24:TYR:HB2	1.98	0.46
1:CA:1097:U:O2	10:CL:22:PRO:HG3	2.16	0.46
1:CA:55:G:N3	1:CA:127:A:C2	2.84	0.46
1:CA:2094:G:C2	1:CA:2196:C:C2	3.03	0.46
1:CA:2272:U:H5'	1:CA:2273:A:OP1	2.16	0.46
1:CA:2563:U:O2	1:CA:2565:A:H8	1.98	0.46
1:CA:909:A:C6	1:CA:912:C:C2	3.03	0.46
6:CF:123:LEU:HD12	6:CF:124:LEU:N	2.31	0.46
1:CA:2690:C:OP2	15:CR:14:SER:HB2	2.15	0.46
34:DA:109:A:H2'	34:DA:326:G:N2	2.31	0.46
37:DD:5:ILE:O	37:DD:5:ILE:HG23	2.15	0.46
38:DE:151:LEU:HD11	41:DH:77:GLU:OE1	2.16	0.46
34:DA:935:A:H61	40:DG:3:ARG:HG3	1.80	0.46
43:DJ:35:SER:HB3	43:DJ:73:ASP:HB2	1.96	0.46
34:DA:986:A:H1'	52:DS:55:LYS:HA	1.97	0.46
57:DZ:199:ILE:HB	57:DZ:200:PRO:HD2	1.97	0.46
27:A3:26:LEU:O	27:A3:35:ARG:NE	2.45	0.46
1:AA:1160:G:H2'	1:AA:1161:G:H8	1.81	0.46
1:AA:1321:A:H4'	1:AA:1322:A:OP1	2.15	0.46
1:AA:1487:G:H2'	1:AA:1488:G:C8	2.50	0.46
1:AA:1566:U:H2'	1:AA:1567:G:O4'	2.16	0.46
1:AA:2162:C:O2	1:AA:2162:C:H2'	2.15	0.46
1:AA:2183:C:O2'	1:AA:2195:A:H4'	2.15	0.46
1:AA:2116:G:C2	1:AA:2218:C:C2	3.04	0.46
1:AA:2319:G:H4'	1:AA:2320:G:O5'	2.16	0.46
1:AA:311:C:H2'	1:AA:312:C:H6	1.81	0.46
1:AA:509:A:C2'	1:AA:510:C:H5'	2.46	0.46
1:AA:750:U:H2'	1:AA:751:G:O4'	2.15	0.46
2:AB:111:G:H2'	2:AB:112:U:H6	1.81	0.46
3:AC:176:VAL:O	3:AC:176:VAL:HG12	2.15	0.46
4:AD:3:VAL:HG13	4:AD:17:THR:HB	1.98	0.46
1:AA:624:C:OP1	6:AF:108:LYS:HE3	2.15	0.46
34:BA:1001:A:H2'	34:BA:1001(A):G:C8	2.51	0.46
34:BA:976:G:C8	34:BA:1362:C:N4	2.84	0.46
37:BD:22:LYS:HG3	60:BD:501:SF4:S4	2.56	0.46
42:BI:9:ARG:HD2	42:BI:104:ARG:NH1	2.31	0.46
46:BM:8:GLU:OE2	46:BM:11:ARG:NH2	2.48	0.46
29:C5:35:GLU:HG2	29:C5:51:TYR:CD2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1265:A:O4'	1:CA:1267:U:C6	2.69	0.46
1:CA:1449:A:H5'	1:CA:1450:G:OP2	2.16	0.46
1:CA:1682:G:H2'	1:CA:1683:C:C6	2.51	0.46
1:CA:2439:A:H5''	1:CA:2439:A:H8	1.75	0.46
3:CC:30:VAL:CG2	3:CC:31:LYS:H	2.27	0.46
7:CG:135:LEU:HD23	7:CG:155:MET:SD	2.56	0.46
7:CG:143:GLU:HA	28:C4:31:ILE:HD13	1.98	0.46
15:CR:67:LEU:HD13	15:CR:76:VAL:HG21	1.98	0.46
16:CS:103:GLU:O	16:CS:107:GLU:HG3	2.15	0.46
20:CW:54:ALA:HB1	20:CW:107:LEU:HD22	1.96	0.46
34:DA:345:C:H4'	34:DA:346:G:O4'	2.15	0.46
34:DA:416:G:C6	34:DA:417:C:C4	3.03	0.46
34:DA:622:A:C8	34:DA:623:C:C5	3.04	0.46
34:DA:730:G:C5	34:DA:731:G:H1'	2.50	0.46
40:DG:46:ALA:HB1	40:DG:121:ALA:HB2	1.98	0.46
1:AA:2356:U:OP1	30:A6:37:ARG:HD3	2.15	0.46
1:AA:2843:G:H4'	1:AA:2844:G:OP2	2.15	0.46
1:AA:343:C:H2'	1:AA:344:A:O4'	2.15	0.46
7:AG:79:ASN:OD1	7:AG:79:ASN:N	2.32	0.46
13:AP:29:LYS:HG2	13:AP:30:THR:HG23	1.98	0.46
34:BA:1036:G:H2'	34:BA:1036:G:N3	2.31	0.46
34:BA:1298:C:C5	40:BG:114:ARG:HD2	2.50	0.46
34:BA:43:C:N4	34:BA:399:G:H1	2.14	0.46
34:BA:572:A:H5'	34:BA:573:A:OP2	2.15	0.46
34:BA:865:A:N3	34:BA:918:A:O2'	2.35	0.46
48:BO:61:GLY:O	48:BO:64:ARG:HB3	2.16	0.46
50:BQ:45:HIS:O	50:BQ:73:VAL:HG23	2.16	0.46
57:BZ:141:LYS:O	57:BZ:171:GLU:HB3	2.16	0.46
57:BZ:97:SER:O	57:BZ:101:LEU:HD12	2.16	0.46
14:CQ:81:VAL:HB	24:C0:7:LEU:HD21	1.98	0.46
25:C1:21:ARG:HD3	25:C1:35:THR:HG21	1.96	0.46
1:CA:2625:G:H2'	1:CA:2626:C:O4'	2.16	0.46
1:CA:288:C:O2'	1:CA:289:A:H5'	2.16	0.46
1:CA:637:A:H4'	1:CA:638:G:O5'	2.16	0.46
1:CA:2784:C:H1'	5:CE:37:ARG:NH1	2.30	0.46
13:CP:135:LEU:HA	13:CP:135:LEU:HD23	1.86	0.46
23:CZ:7:ALA:HB2	23:CZ:59:LEU:HD22	1.97	0.46
23:CZ:59:LEU:O	23:CZ:61:LEU:HD22	2.16	0.46
34:DA:298:A:O5'	34:DA:298:A:H8	1.98	0.46
34:DA:687:A:H4'	34:DA:688:G:O5'	2.16	0.46
34:DA:717:C:H4'	44:DK:117:ASN:HD22	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:73:G:C6	34:DA:97:G:C6	3.03	0.46
35:DB:7:VAL:HG12	35:DB:8:LYS:HG2	1.98	0.46
39:DF:2:ARG:CZ	39:DF:69:GLU:HG2	2.46	0.46
56:DY:65:G:H2'	56:DY:66:U:C6	2.51	0.46
57:DZ:151:ARG:O	57:DZ:154:GLN:HG2	2.15	0.46
26:A2:51:ARG:HD3	26:A2:55:ARG:NH1	2.30	0.46
1:AA:2047:C:H2'	1:AA:2048:C:C6	2.51	0.46
1:AA:553:A:H2	1:AA:2065:C:C5'	2.28	0.46
1:AA:596:G:O2'	1:AA:597:C:H3'	2.15	0.46
1:AA:645:G:H5'	1:AA:645:G:N3	2.30	0.46
1:AA:732:A:H1'	1:AA:735:U:O4	2.16	0.46
5:AE:18:ASP:HB3	17:AT:82:LEU:HD11	1.97	0.46
7:AG:75:LYS:HE3	7:AG:77:ILE:HD11	1.98	0.46
20:AW:2:GLU:OE2	20:AW:72:LYS:NZ	2.31	0.46
34:BA:1203:C:H2'	34:BA:1204:A:H8	1.81	0.46
34:BA:1401:G:N2	34:BA:1402:C:H1'	2.31	0.46
34:BA:27:G:C5	34:BA:557:G:C2	3.04	0.46
34:BA:403:C:H2'	34:BA:404:U:C6	2.47	0.46
34:BA:581:G:O2'	34:BA:582:U:H5'	2.15	0.46
35:BB:69:LEU:HD12	35:BB:70:PHE:N	2.31	0.46
36:BC:181:ASN:HB3	36:BC:204:LEU:HB2	1.97	0.46
37:BD:101:LEU:HG	37:BD:121:VAL:HG11	1.98	0.46
48:BO:9:GLN:HA	48:BO:12:ILE:HD12	1.98	0.46
54:BU:8:THR:O	54:BU:12:LYS:HB2	2.17	0.46
32:C8:34:TRP:CE2	32:C8:35:GLN:HG2	2.50	0.46
1:CA:1711:C:H2'	1:CA:1712:C:H6	1.80	0.46
1:CA:1882:C:H2'	1:CA:1883:G:O4'	2.16	0.46
1:CA:2169:A:H2'	1:CA:2170:A:C8	2.51	0.46
1:CA:2176:A:C3'	3:CC:45:HIS:CD2	3.00	0.46
1:CA:271(S):G:C6	1:CA:271(T):C:C4	3.03	0.46
2:CB:59:A:H2'	2:CB:60:C:O4'	2.16	0.46
3:CC:46:ALA:O	3:CC:47:LYS:HB2	2.15	0.46
10:CL:44:ALA:O	10:CL:47:ASN:HB3	2.16	0.46
23:CZ:19:ARG:HA	23:CZ:23:LYS:O	2.16	0.46
23:CZ:53:ILE:HD12	23:CZ:99:TYR:HB2	1.98	0.46
34:DA:302:G:N3	34:DA:556:C:H4'	2.30	0.46
35:DB:116:GLU:HA	35:DB:119:GLU:HB2	1.98	0.46
39:DF:33:TYR:CD1	39:DF:75:LEU:HD23	2.51	0.46
50:DQ:52:LYS:HE3	50:DQ:52:LYS:HB2	1.74	0.46
57:DZ:188:TYR:CD1	57:DZ:196:ILE:HD13	2.51	0.46
57:DZ:488:THR:O	57:DZ:516:PRO:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
28:A4:16:CYS:HA	28:A4:33:VAL:HB	1.98	0.45
1:AA:464:G:H8	1:AA:464:G:O5'	1.98	0.45
3:AC:20:VAL:O	3:AC:224:ARG:O	2.34	0.45
3:AC:211:ARG:HH11	3:AC:211:ARG:HG2	1.81	0.45
13:AP:81:GLN:OE1	13:AP:106:LEU:HA	2.16	0.45
2:AB:27:C:H5''	16:AS:54:LEU:HD13	1.98	0.45
17:AT:61:PHE:CE1	17:AT:78:LEU:HD23	2.51	0.45
34:BA:232:G:H1'	34:BA:262:A:N1	2.30	0.45
35:BB:213:LEU:HD23	35:BB:213:LEU:O	2.15	0.45
37:BD:107:ARG:HH22	37:BD:194:LEU:HD11	1.81	0.45
42:BI:118:LYS:HG3	42:BI:121:ARG:HB3	1.98	0.45
56:BW:14:A:C2	56:BW:15:G:H1'	2.51	0.45
57:BZ:162:VAL:O	57:BZ:164:MET:HG2	2.16	0.45
1:CA:1562:A:H2'	1:CA:1563:G:C8	2.51	0.45
11:CN:84:LYS:C	11:CN:85:ILE:HG13	2.35	0.45
34:DA:1124:G:O2'	34:DA:1145:C:C4	2.69	0.45
34:DA:201:C:H42	34:DA:216:G:H22	1.63	0.45
34:DA:350:G:O2'	34:DA:351:G:H5'	2.16	0.45
34:DA:407:G:C6	34:DA:408:A:C6	3.04	0.45
34:DA:926:G:H5''	34:DA:927:G:O5'	2.16	0.45
35:DB:101:MET:HA	35:DB:108:ILE:HG13	1.98	0.45
35:DB:28:PHE:CD1	35:DB:190:THR:HG22	2.51	0.45
40:DG:59:LEU:HG	40:DG:63:LYS:HE2	1.97	0.45
44:DK:22:HIS:HB3	44:DK:29:ILE:HB	1.98	0.45
51:DR:56:THR:HB	51:DR:58:LEU:HD23	1.96	0.45
56:DW:65:G:H2'	56:DW:66:U:C6	2.51	0.45
56:DY:9:A:H8	56:DY:11:C:H41	1.64	0.45
1:AA:1157:A:H2'	1:AA:1158:G:O4'	2.16	0.45
1:AA:1773:C:H2'	1:AA:1774:C:C6	2.51	0.45
1:AA:2500:A:H2'	1:AA:2501:G:O4'	2.16	0.45
1:AA:26:G:C6	1:AA:27:G:N1	2.85	0.45
1:AA:886:U:H1'	1:AA:1236:G:H1'	1.98	0.45
1:AA:950:C:H2'	1:AA:951:U:C6	2.51	0.45
2:AB:63:G:H2'	2:AB:64:C:C6	2.52	0.45
3:AC:48:LEU:HD23	3:AC:59:VAL:HG21	1.98	0.45
7:AG:126:ASP:CG	7:AG:130:ASN:HD22	2.19	0.45
7:AG:137:GLU:C	7:AG:140:ILE:HD13	2.36	0.45
14:AQ:2:LEU:HG	14:AQ:69:PHE:CD2	2.51	0.45
15:AR:65:LEU:HD12	15:AR:65:LEU:HA	1.73	0.45
16:AS:39:ILE:HB	16:AS:49:VAL:CG1	2.46	0.45
22:AY:99:CYS:SG	22:AY:101:LYS:HB2	2.56	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:1241:G:H1	34:BA:1296:C:H42	1.63	0.45
34:BA:642:A:H2'	34:BA:643:C:C6	2.51	0.45
34:BA:781:A:C5	34:BA:802:A:C2	3.04	0.45
37:BD:178:VAL:HG12	37:BD:179:GLU:H	1.81	0.45
38:BE:37:ARG:NH1	38:BE:111:GLU:HG2	2.31	0.45
42:BI:23:ASN:N	42:BI:60:ASP:OD1	2.49	0.45
45:BL:88:GLY:O	45:BL:99:HIS:HD2	1.99	0.45
31:C7:19:ARG:HG2	31:C7:19:ARG:NH1	2.31	0.45
1:CA:271(H):G:O2'	1:CA:271(I):G:H8	1.99	0.45
2:CB:94:C:H2'	2:CB:95:C:H6	1.81	0.45
1:CA:2177:C:O2'	3:CC:171:ALA:HB2	2.16	0.45
6:CF:53:THR:CG2	6:CF:55:GLY:H	2.28	0.45
7:CG:106:LEU:O	7:CG:110:ALA:HB3	2.16	0.45
1:CA:1054:A:O2'	9:CK:30:GLN:O	2.35	0.45
10:CL:134:MET:HG3	10:CL:136:VAL:HG12	1.98	0.45
17:CT:23:ARG:HG3	17:CT:120:ARG:CZ	2.46	0.45
18:CU:17:ILE:HA	18:CU:17:ILE:HD13	1.87	0.45
34:DA:1410:G:H2'	34:DA:1411:C:C6	2.51	0.45
34:DA:45:U:H2'	34:DA:46:G:C8	2.51	0.45
34:DA:688:G:H2'	34:DA:689:C:H6	1.81	0.45
35:DB:16:HIS:CG	35:DB:17:PHE:N	2.84	0.45
37:DD:79:PHE:CE1	37:DD:204:ILE:HD13	2.51	0.45
34:DA:718:G:H5'	44:DK:117:ASN:HB2	1.98	0.45
50:DQ:65:ILE:HD11	50:DQ:72:ARG:HD2	1.98	0.45
57:DZ:184:LYS:NZ	57:DZ:184:LYS:HB2	2.32	0.45
57:DZ:182:ARG:NH2	57:DZ:278:ASP:OD2	2.49	0.45
57:DZ:411:VAL:HB	57:DZ:459:LEU:HD13	1.99	0.45
57:DZ:7:ASN:O	57:DZ:11:ARG:NH2	2.45	0.45
57:DZ:21:ILE:HD12	57:DZ:87:HIS:HA	1.98	0.45
1:AA:1114:G:N2	1:AA:1142:A:H5'	2.31	0.45
1:AA:1136:U:C2	1:AA:1148:C:H1'	2.51	0.45
1:AA:1632:A:O5'	1:AA:1632:A:H8	1.99	0.45
1:AA:180:A:H2'	1:AA:181:C:C6	2.51	0.45
6:AF:184:TYR:CE2	6:AF:188:ARG:HD2	2.50	0.45
10:AL:20:ALA:HB1	57:BZ:636:PRO:HG2	1.97	0.45
34:BA:1064:G:H4'	34:BA:1065:U:OP1	2.14	0.45
34:BA:1437:C:H2'	34:BA:1438:G:H8	1.82	0.45
34:BA:510:A:N3	34:BA:543:C:H1'	2.31	0.45
34:BA:938:A:C6	34:BA:939:G:C5	3.05	0.45
36:BC:26:LYS:HG3	36:BC:26:LYS:H	1.59	0.45
41:BH:40:ALA:C	41:BH:42:GLU:H	2.19	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
46:BM:40:ASN:HA	46:BM:41:PRO:HD3	1.81	0.45
48:BO:15:PHE:CZ	48:BO:84:LYS:HD2	2.51	0.45
34:BA:958:A:N6	52:BS:77:THR:O	2.48	0.45
57:BZ:388:THR:OG1	57:BZ:399:LEU:HB2	2.16	0.45
57:BZ:39:ILE:HD12	57:BZ:40:HIS:H	1.81	0.45
57:BZ:505:GLY:HA2	57:BZ:576:ASP:OD2	2.16	0.45
1:CA:1068:G:O2'	1:CA:1096:A:O2'	2.26	0.45
1:CA:188:G:H2'	1:CA:189:G:H5'	1.98	0.45
1:CA:1996:C:H4'	1:CA:1997:G:OP1	2.17	0.45
1:CA:901:A:H5''	1:CA:902:C:OP2	2.17	0.45
1:CA:942:G:O2'	1:CA:943:U:H5'	2.16	0.45
3:CC:37:LYS:O	3:CC:38:PHE:HB3	2.17	0.45
4:CD:101:GLU:HG3	4:CD:102:LYS:N	2.31	0.45
5:CE:120:TRP:CE3	5:CE:155:LYS:HD3	2.52	0.45
6:CF:137:LYS:HB3	6:CF:137:LYS:HE2	1.55	0.45
8:CH:3:ARG:CZ	8:CH:5:GLY:H	2.30	0.45
10:CL:107:ILE:HG21	10:CL:127:ILE:HG21	1.98	0.45
10:CL:59:ILE:HG12	10:CL:60:TYR:N	2.31	0.45
14:CQ:135:ASP:HB2	14:CQ:137:TYR:HB2	1.98	0.45
22:CY:2:ARG:CZ	22:CY:4:LYS:HD2	2.46	0.45
34:DA:1353:G:H2'	34:DA:1354:C:H6	1.82	0.45
34:DA:1428:A:H2'	34:DA:1429:C:O4'	2.15	0.45
34:DA:245:C:O2	34:DA:283:C:N3	2.48	0.45
34:DA:49:U:H3	34:DA:362:G:H1'	1.80	0.45
34:DA:617:G:C6	34:DA:618:C:C5	3.04	0.45
36:DC:18:TRP:HE3	36:DC:18:TRP:H	1.63	0.45
36:DC:179:ARG:HD2	36:DC:206:GLU:HB2	1.97	0.45
45:DL:75:HIS:CD2	45:DL:77:LEU:N	2.82	0.45
46:DM:96:LEU:C	46:DM:110:ARG:HG2	2.37	0.45
34:DA:110:C:O2'	49:DP:25:ARG:O	2.33	0.45
53:DT:55:ILE:O	53:DT:58:LYS:N	2.49	0.45
57:DZ:169:GLY:C	57:DZ:170:ARG:HH11	2.19	0.45
57:DZ:309:LEU:HD12	57:DZ:310:ALA:H	1.82	0.45
1:AA:1096:A:H2'	1:AA:1097:G:C8	2.51	0.45
1:AA:1102:G:H5''	1:AA:1103:A:O4'	2.16	0.45
1:AA:1699:A:C2'	1:AA:1700:G:H5'	2.47	0.45
1:AA:354:A:O2'	1:AA:355:A:H8	1.99	0.45
1:AA:372:G:H2'	1:AA:373:G:O4'	2.17	0.45
1:AA:646:A:OP2	13:AP:108:LYS:NZ	2.48	0.45
3:AC:30:VAL:CG2	3:AC:31:LYS:N	2.78	0.45
3:AC:37:LYS:O	3:AC:38:PHE:HB3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AC:7:ARG:HH22	3:AC:219:MET:HB2	1.82	0.45
6:AF:110:LEU:HD23	6:AF:110:LEU:HA	1.80	0.45
12:AO:68:GLU:HB3	12:AO:78:ARG:HB2	1.99	0.45
23:AZ:163:LEU:HD12	23:AZ:163:LEU:HA	1.57	0.45
34:BA:165:C:H2'	34:BA:166:G:C8	2.52	0.45
34:BA:509:A:C8	34:BA:509:A:C3'	2.99	0.45
34:BA:540:G:H2'	34:BA:541:G:O4'	2.16	0.45
34:BA:948:C:O2'	34:BA:949:A:H5'	2.16	0.45
35:BB:12:GLU:C	35:BB:14:GLY:H	2.20	0.45
35:BB:42:ILE:HG21	35:BB:202:PRO:O	2.16	0.45
38:BE:51:VAL:O	38:BE:55:VAL:HG23	2.17	0.45
39:BF:4:TYR:CD1	39:BF:92:LYS:HA	2.52	0.45
44:BK:21:ILE:HD13	44:BK:94:ALA:HB3	1.98	0.45
49:BP:29:ASP:N	49:BP:29:ASP:OD1	2.49	0.45
50:BQ:31:LEU:HD23	50:BQ:32:TYR:CZ	2.52	0.45
57:BZ:507:TYR:HB3	57:BZ:577:SER:HB2	1.98	0.45
24:C0:27:GLU:HG3	24:C0:68:GLU:HA	1.98	0.45
1:CA:1203:G:OP2	1:CA:1204:A:O2'	2.28	0.45
1:CA:154(A):C:H42	1:CA:171:G:H1	1.65	0.45
1:CA:1668:A:C8	1:CA:1674:G:C6	3.05	0.45
1:CA:2250:G:C8	1:CA:2496:C:H5''	2.51	0.45
1:CA:2353:G:H2'	1:CA:2354:G:O4'	2.15	0.45
1:CA:2478:A:H1'	1:CA:2528:U:O2'	2.16	0.45
1:CA:2853:C:H2'	1:CA:2854:G:H8	1.81	0.45
1:CA:2880:C:O2'	15:CR:90:ARG:NH1	2.50	0.45
3:CC:54:ARG:HH22	3:CC:56:ASP:HB3	1.75	0.45
1:CA:694:U:OP1	4:CD:59:LYS:NZ	2.50	0.45
10:CL:90:LYS:HE3	10:CL:90:LYS:HB3	1.83	0.45
1:CA:2276:G:H5'	14:CQ:86:GLY:HA2	1.97	0.45
21:CX:26:TYR:HD1	21:CX:92:LEU:HD12	1.80	0.45
23:CZ:117:LEU:HD12	23:CZ:174:VAL:HG22	1.98	0.45
34:DA:1141:C:H2'	34:DA:1142:G:C8	2.51	0.45
34:DA:303:A:O2'	34:DA:555:C:O2'	2.30	0.45
34:DA:401:C:H1'	34:DA:622:A:H1'	1.98	0.45
40:DG:92:SER:O	40:DG:95:ARG:N	2.48	0.45
42:DI:8:GLY:O	42:DI:15:ALA:N	2.39	0.45
48:DO:74:ASP:OD2	48:DO:77:ARG:HG3	2.16	0.45
54:DU:10:ARG:HA	54:DU:13:ILE:HD12	1.99	0.45
57:DZ:314:PHE:CE2	57:DZ:327:PHE:HB3	2.50	0.45
24:A0:72:ARG:HB2	24:A0:75:LEU:HB2	1.97	0.45
1:AA:1066:A:N1	1:AA:1186:U:O2'	2.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1529:G:C6	1:AA:1553:A:C6	3.04	0.45
1:AA:2045:G:H5'	1:AA:2629:C:H4'	1.98	0.45
1:AA:37:C:H2'	1:AA:38:A:C8	2.51	0.45
1:AA:626:A:O4'	1:AA:702:A:N6	2.50	0.45
1:AA:820:U:H2'	1:AA:821:A:H5'	1.98	0.45
5:AE:29:GLY:HA3	63:AE:417:HOH:O	2.16	0.45
7:AG:161:THR:HG23	7:AG:163:ALA:H	1.81	0.45
11:AN:68:GLU:HG2	11:AN:88:GLU:OE2	2.16	0.45
12:AO:117:LEU:HA	12:AO:117:LEU:HD23	1.52	0.45
22:AY:79:CYS:SG	22:AY:81:LYS:HG3	2.57	0.45
34:BA:597:G:C4	34:BA:644:G:C2	3.04	0.45
40:BG:14:PRO:HB3	40:BG:19:GLY:C	2.37	0.45
40:BG:21:VAL:HG23	40:BG:22:LEU:N	2.31	0.45
43:BJ:30:SER:HB2	43:BJ:81:THR:HG21	1.98	0.45
57:BZ:112:GLN:HG2	57:BZ:115:GLU:CD	2.37	0.45
57:BZ:-23:LEU:HA	57:BZ:-20:LEU:HB2	1.99	0.45
30:C6:14:THR:OG1	30:C6:48:VAL:O	2.20	0.45
31:C7:5:TRP:CD1	31:C7:7:PRO:HD3	2.51	0.45
1:CA:1069:A:N7	1:CA:1073:A:N6	2.64	0.45
1:CA:1206:G:C2	1:CA:1207:C:C2	3.05	0.45
1:CA:1378:A:O2'	1:CA:1380:G:N7	2.35	0.45
1:CA:1537:G:H2'	1:CA:1538:G:H8	1.80	0.45
1:CA:2128:C:H5'	1:CA:2173:A:N3	2.32	0.45
1:CA:957:A:C2	1:CA:2459:A:H5'	2.50	0.45
1:CA:2687:U:H2'	1:CA:2688:U:O4'	2.16	0.45
1:CA:271(D):G:H2'	1:CA:271(E):U:C6	2.51	0.45
5:CE:4:ILE:HD13	5:CE:28:ALA:HB1	1.97	0.45
5:CE:77:ILE:HD11	5:CE:79:ARG:HH12	1.81	0.45
7:CG:180:PHE:O	7:CG:182:LYS:N	2.49	0.45
10:CL:72:PRO:HA	10:CL:73:PRO:HD3	1.83	0.45
13:CP:59:LEU:HD11	32:C8:10:ALA:HA	1.99	0.45
14:CQ:84:GLY:O	14:CQ:85:LYS:HB2	2.15	0.45
15:CR:62:ALA:HA	15:CR:65:LEU:HD23	1.98	0.45
16:CS:3:ARG:HE	16:CS:4:LEU:H	1.64	0.45
22:CY:49:VAL:HG11	22:CY:55:TYR:CD2	2.51	0.45
23:CZ:144:LEU:HD22	23:CZ:148:ASP:HB3	1.98	0.45
34:DA:1060:C:H4'	43:DJ:51:ARG:HB3	1.97	0.45
34:DA:1292:U:H2'	34:DA:1293:G:H8	1.82	0.45
34:DA:376:G:H5''	49:DP:5:ARG:HB2	1.99	0.45
34:DA:489:C:N4	34:DA:490:G:O6	2.50	0.45
34:DA:589:C:H2'	34:DA:590:C:H6	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:893:C:C4	34:DA:894:G:N7	2.84	0.45
34:DA:9:G:H2'	34:DA:10:A:C8	2.52	0.45
34:DA:9:G:H2'	34:DA:10:A:H8	1.80	0.45
42:DI:36:TYR:HD2	42:DI:37:PHE:CE2	2.35	0.45
43:DJ:44:VAL:HG13	43:DJ:66:ARG:HG2	1.98	0.45
45:DL:69:TYR:CD2	45:DL:99:HIS:HE1	2.35	0.45
49:DP:23:ASP:OD2	49:DP:25:ARG:NH1	2.49	0.45
28:C4:61:ARG:HE	52:DS:42:PRO:HG3	1.81	0.45
57:DZ:606:MET:N	57:DZ:647:VAL:O	2.46	0.45
1:AA:1186:U:OP1	11:AN:25:ARG:NH1	2.45	0.45
1:AA:11:G:C2'	1:AA:12:U:H5''	2.42	0.45
1:AA:1480:A:C2	1:AA:1481:G:C4	3.04	0.45
1:AA:1857:G:H2'	1:AA:1858:C:H6	1.82	0.45
1:AA:1911:A:H1'	1:AA:2109:G:O4'	2.16	0.45
1:AA:2190:G:O6	1:AA:2193:A:H2'	2.16	0.45
1:AA:329:U:H2'	1:AA:330:U:C6	2.51	0.45
12:AO:47:ILE:HB	12:AO:48:PRO:HD2	1.98	0.45
17:AT:23:ARG:HG3	17:AT:120:ARG:NH1	2.31	0.45
19:AV:43:GLU:H	19:AV:43:GLU:HG2	1.45	0.45
34:BA:294:U:OP1	34:BA:610:G:O2'	2.27	0.45
35:BB:69:LEU:HD12	35:BB:70:PHE:H	1.81	0.45
36:BC:125:GLU:HA	36:BC:191:THR:HG22	1.98	0.45
34:BA:528:C:H41	45:BL:49:ASN:HD21	1.62	0.45
46:BM:4:ILE:HA	46:BM:5:ALA:HA	1.72	0.45
54:BU:5:ASP:O	54:BU:11:GLY:HA3	2.17	0.45
57:BZ:287:PRO:HA	57:BZ:288:PRO:HD3	1.76	0.45
57:BZ:432:ALA:HA	57:BZ:438:PHE:CE1	2.49	0.45
61:BZ:703:FUA:C1	61:BZ:703:FUA:O1	2.60	0.45
1:CA:1666:G:O2'	1:CA:1667:G:H5'	2.16	0.45
1:CA:2582:G:OP2	1:CA:2582:G:H3'	2.16	0.45
1:CA:478:A:N1	1:CA:500:G:H4'	2.31	0.45
1:CA:587:C:OP2	13:CP:21:ARG:NH2	2.49	0.45
1:CA:632:A:H2'	1:CA:633:A:C8	2.51	0.45
3:CC:22:THR:HG23	3:CC:25:GLU:OE1	2.17	0.45
4:CD:260:ARG:O	4:CD:260:ARG:HG3	2.17	0.45
6:CF:36:VAL:HG11	6:CF:183:VAL:CG1	2.46	0.45
12:CO:68:GLU:HG2	12:CO:68:GLU:O	2.17	0.45
19:CV:16:PRO:HD3	19:CV:99:ILE:HD11	1.98	0.45
22:CY:31:LEU:HA	22:CY:32:PRO:HD3	1.71	0.45
22:CY:76:CYS:SG	22:CY:99:CYS:HB2	2.56	0.45
34:DA:1064:G:H21	34:DA:1190:G:H2'	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:25:C:C5	34:DA:558:G:N2	2.84	0.45
37:DD:99:SER:O	37:DD:140:VAL:HG23	2.16	0.45
38:DE:72:GLN:O	38:DE:75:THR:HG22	2.17	0.45
56:DY:23:A:N6	56:DY:24:G:O6	2.50	0.45
32:A8:62:LEU:HB3	32:A8:65:GLU:HG2	1.98	0.45
1:AA:1775:C:H5'	1:AA:1776:G:OP2	2.17	0.45
1:AA:2221:A:OP2	1:AA:2222:C:H5	1.99	0.45
3:AC:31:LYS:HG2	3:AC:31:LYS:H	1.57	0.45
21:AX:12:VAL:HG22	21:AX:29:TRP:CE2	2.52	0.45
22:AY:76:CYS:SG	22:AY:78:ALA:HB3	2.56	0.45
34:BA:1295:G:O5'	34:BA:1295:G:H8	1.98	0.45
34:BA:1376:U:H2'	34:BA:1377:A:C8	2.52	0.45
34:BA:456:C:N4	34:BA:475:G:H1	2.03	0.45
34:BA:839:U:H4'	34:BA:840:C:OP2	2.16	0.45
35:BB:91:PRO:HG2	35:BB:155:LEU:HD13	1.97	0.45
35:BB:166:ASP:HA	35:BB:167:PRO:HD3	1.64	0.45
37:BD:61:LYS:HG3	37:BD:203:VAL:HG22	1.98	0.45
39:BF:44:GLY:HA2	39:BF:59:TYR:CE1	2.51	0.45
41:BH:31:PHE:CE2	41:BH:35:ILE:HD11	2.51	0.45
57:BZ:-32:LEU:HA	57:BZ:-32:LEU:HD23	1.71	0.45
24:C0:68:GLU:OE2	24:C0:82:ARG:NH1	2.49	0.45
32:C8:6:THR:CG2	32:C8:64:TYR:HD2	2.29	0.45
1:CA:1557:C:H5''	1:CA:1558:A:OP2	2.17	0.45
1:CA:1796:U:H2'	1:CA:1797:C:H6	1.80	0.45
1:CA:1651:G:C2	1:CA:2007:C:N3	2.85	0.45
1:CA:212:G:H2'	1:CA:213:A:O4'	2.17	0.45
1:CA:2460:U:H2'	1:CA:2461:C:O4'	2.17	0.45
1:CA:359:A:H2'	1:CA:360:G:O4'	2.16	0.45
1:CA:518:G:H2'	1:CA:519:U:C6	2.52	0.45
1:CA:840:C:H2'	1:CA:841:A:C8	2.52	0.45
7:CG:2:PRO:HB2	7:CG:3:LEU:H	1.57	0.45
9:CK:49:ALA:H	9:CK:90:ALA:HB1	1.82	0.45
11:CN:4:TYR:O	18:CU:64:ARG:NH2	2.39	0.45
2:CB:27:C:H5''	16:CS:54:LEU:HD11	1.98	0.45
22:CY:41:GLY:N	22:CY:64:GLU:OE2	2.40	0.45
34:DA:1292:U:H2'	34:DA:1293:G:C8	2.52	0.45
34:DA:933:G:C2	34:DA:1385:G:C2	3.04	0.45
37:DD:59:ARG:HE	37:DD:59:ARG:HA	1.80	0.45
53:DT:56:MET:HE1	53:DT:85:MET:HG2	1.98	0.45
56:DW:23:A:H2'	56:DW:24:G:H8	1.81	0.45
34:DA:1340:A:P	56:DY:35:A:OP1	2.75	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:608:VAL:O	57:DZ:644:ARG:HA	2.17	0.45
1:AA:1097:G:H5'	1:AA:1098:C:OP2	2.16	0.45
1:AA:1112:U:O2	1:AA:1114:G:C8	2.70	0.45
1:AA:1193:C:O2'	1:AA:1194:A:H5'	2.16	0.45
1:AA:1481:G:H21	1:AA:1525:G:H5'	1.82	0.45
1:AA:197:C:H2'	1:AA:198:C:C6	2.52	0.45
1:AA:2036:A:H2'	1:AA:2037:A:C8	2.52	0.45
1:AA:233:A:C2	1:AA:244:A:C4	3.05	0.45
1:AA:346:A:H5'	1:AA:364:A:H1'	1.98	0.45
7:AG:125:PHE:HB3	7:AG:166:ASP:OD1	2.16	0.45
2:AB:57:A:H1'	7:AG:29:TRP:HB2	1.98	0.45
7:AG:43:LEU:HB2	7:AG:89:GLY:HA2	1.99	0.45
8:AH:71:LEU:HA	8:AH:71:LEU:HD12	1.82	0.45
34:BA:1101:A:H4'	34:BA:1102:A:O5'	2.17	0.45
34:BA:110:C:H2'	34:BA:111:G:O4'	2.17	0.45
34:BA:1361:G:H2'	34:BA:1362:C:O4'	2.17	0.45
34:BA:1418:A:C2	34:BA:1483:A:C2	3.05	0.45
36:BC:44:GLU:HA	36:BC:52:LEU:HD12	1.99	0.45
37:BD:61:LYS:HD3	37:BD:206:PHE:CE1	2.52	0.45
41:BH:12:ARG:NH2	41:BH:27:PRO:HD3	2.31	0.45
41:BH:51:VAL:HG21	41:BH:60:ARG:CD	2.47	0.45
56:BY:48:C:H2'	56:BY:48:C:OP1	2.17	0.45
57:BZ:160:ARG:HB2	57:BZ:255:ILE:HA	1.99	0.45
1:CA:768:G:O2'	1:CA:1379:A:N1	2.45	0.45
1:CA:185:U:H4'	1:CA:218:A:H4'	1.99	0.45
1:CA:2582:G:C2	1:CA:2583:G:C8	3.05	0.45
1:CA:2702:U:H4'	1:CA:2703:C:OP1	2.17	0.45
1:CA:2761:G:C2	1:CA:2762:G:C8	3.05	0.45
1:CA:2841:C:C2	1:CA:2877:G:N2	2.84	0.45
1:CA:874:G:N2	1:CA:904:C:C2	2.85	0.45
2:CB:22:U:O2	2:CB:61:G:N2	2.35	0.45
2:CB:79:C:H2'	2:CB:80:U:O4'	2.17	0.45
4:CD:10:THR:HG23	4:CD:13:ARG:HG2	1.99	0.45
4:CD:5:LYS:HG2	4:CD:17:THR:HG22	1.98	0.45
4:CD:228:PRO:HD3	4:CD:235:GLY:CA	2.47	0.45
12:CO:118:ALA:HA	12:CO:119:PRO:HD3	1.71	0.45
14:CQ:32:TYR:HE2	14:CQ:111:GLU:HA	1.82	0.45
34:DA:1006:C:H2'	34:DA:1007:C:O4'	2.16	0.45
34:DA:403:C:H2'	34:DA:404:U:C6	2.49	0.45
34:DA:834:C:H2'	34:DA:835:U:C6	2.51	0.45
35:DB:108:ILE:HG12	35:DB:108:ILE:H	1.54	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:DC:136:GLN:O	36:DC:140:ARG:N	2.43	0.45
40:DG:38:LEU:O	40:DG:42:ILE:HG13	2.16	0.45
44:DK:102:GLY:O	44:DK:103:LEU:HD22	2.17	0.45
45:DL:6:THR:HG23	45:DL:9:GLN:CD	2.36	0.45
49:DP:5:ARG:CZ	49:DP:22:THR:HG21	2.47	0.45
57:DZ:236:GLU:HA	57:DZ:237:PRO:HD3	1.79	0.45
57:DZ:357:ARG:HH12	57:DZ:373:ASP:CG	2.20	0.45
57:DZ:390:VAL:HG23	57:DZ:391:GLY:O	2.17	0.45
46:DM:123:ALA:HB2	57:DZ:507:TYR:CD1	2.51	0.45
1:AA:1116:A:O2'	1:AA:1117:G:OP1	2.32	0.45
1:AA:1401:G:C6	1:AA:1402:G:C5	3.05	0.45
1:AA:1556:A:H3'	1:AA:1557:A:H8	1.82	0.45
1:AA:1935:A:C8	1:AA:1935:A:H5'	2.48	0.45
1:AA:2287:C:O2	14:AQ:85:LYS:HG3	2.17	0.45
1:AA:2589:A:O4'	29:A5:3:LYS:HB2	2.17	0.45
1:AA:809:U:H4'	1:AA:810:G:O5'	2.17	0.45
3:AC:16:ASP:OD2	3:AC:19:LYS:HB2	2.17	0.45
10:AL:17:ALA:HB1	10:AL:38:VAL:HG22	1.99	0.45
10:AL:6:ALA:H	10:AL:59:ILE:HG22	1.82	0.45
1:AA:2874:G:OP1	17:AT:119:LYS:HD2	2.17	0.45
19:AV:39:LEU:HD23	19:AV:40:LEU:N	2.32	0.45
21:AX:41:ASN:O	21:AX:45:THR:HG23	2.17	0.45
23:AZ:150:LEU:HB3	23:AZ:171:ILE:HD11	1.99	0.45
34:BA:189(F):U:C4	50:BQ:72:ARG:NH1	2.85	0.45
34:BA:340:U:H2'	34:BA:341:C:C6	2.52	0.45
34:BA:512:U:H2'	34:BA:513:C:C6	2.52	0.45
34:BA:514:C:O2	34:BA:538:G:C2	2.69	0.45
34:BA:545:C:OP2	37:BD:65:ARG:NH2	2.50	0.45
34:BA:559:A:H2'	34:BA:559:A:N3	2.31	0.45
34:BA:738:C:C2	34:BA:739:C:C5	3.04	0.45
34:BA:781:A:H5'	34:BA:782:A:OP2	2.17	0.45
35:BB:28:PHE:CD1	35:BB:190:THR:HA	2.52	0.45
39:BF:41:GLU:O	39:BF:43:LEU:N	2.50	0.45
42:BI:99:LEU:HB3	42:BI:101:PHE:CE1	2.52	0.45
47:BN:34:TYR:N	47:BN:39:LEU:O	2.48	0.45
49:BP:58:TYR:O	49:BP:62:VAL:HG22	2.16	0.45
52:BS:31:ILE:HB	52:BS:49:ILE:HG23	1.97	0.45
29:C5:20:ARG:HA	29:C5:23:HIS:ND1	2.32	0.45
1:CA:667:U:O2	32:C8:2:PRO:HD2	2.17	0.45
1:CA:687:C:C2	1:CA:788:A:H5'	2.52	0.45
1:CA:904:C:H2'	1:CA:905:U:C6	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CC:48:LEU:HD23	3:CC:59:VAL:HG21	1.98	0.45
5:CE:201:THR:HG23	5:CE:203:LYS:H	1.81	0.45
5:CE:68:ALA:C	5:CE:70:ALA:N	2.70	0.45
6:CF:178:PRO:HB3	6:CF:198:ALA:HB1	1.98	0.45
7:CG:114:ILE:HB	7:CG:117:PHE:HD2	1.82	0.45
17:CT:24:PRO:HA	17:CT:49:VAL:HG13	1.98	0.45
22:CY:29:GLU:HB3	22:CY:38:ILE:HG13	1.99	0.45
34:DA:437:U:H5'	37:DD:155:LEU:HD21	1.98	0.45
34:DA:475:G:H2'	34:DA:476:G:C8	2.52	0.45
34:DA:707:C:H2'	34:DA:708:C:C6	2.51	0.45
34:DA:778:G:C6	34:DA:779:C:C4	3.05	0.45
37:DD:127:THR:HG23	37:DD:147:ALA:HB3	1.99	0.45
37:DD:79:PHE:HE1	37:DD:204:ILE:HD13	1.81	0.45
34:DA:921:U:O2	38:DE:19:MET:HB2	2.17	0.45
36:DC:58:GLU:HB3	43:DJ:92:THR:HG21	1.99	0.45
44:DK:34:ASP:HB2	44:DK:35:PRO:HD2	1.98	0.45
46:DM:123:ALA:HB1	57:DZ:507:TYR:HB3	1.98	0.45
57:DZ:-62:LEU:HD12	57:DZ:-62:LEU:N	2.32	0.45
1:AA:116:A:H3'	1:AA:117:A:C5'	2.45	0.45
1:AA:1250:U:H4'	1:AA:1251:G:OP2	2.17	0.45
1:AA:1921:G:H2'	1:AA:1921:G:N3	2.32	0.45
2:AB:46:A:C5	2:AB:47:C:C5	3.05	0.45
3:AC:179:ALA:O	3:AC:180:SER:O	2.35	0.45
3:AC:203:GLU:N	3:AC:203:GLU:CD	2.70	0.45
6:AF:135:LYS:HB2	6:AF:138:GLU:CG	2.43	0.45
7:AG:67:LYS:NZ	7:AG:68:PRO:O	2.50	0.45
7:AG:76:SER:N	7:AG:84:LYS:HB2	2.32	0.45
13:AP:138:LEU:HD12	13:AP:138:LEU:HA	1.77	0.45
15:AR:36:THR:HB	15:AR:37:THR:H	1.51	0.45
34:BA:1157:A:C5	34:BA:1181:G:C6	3.05	0.45
34:BA:1343:G:H2'	34:BA:1344:C:C6	2.51	0.45
34:BA:375:U:C4	34:BA:376:G:N7	2.85	0.45
34:BA:986:A:H1'	52:BS:54:GLY:O	2.17	0.45
35:BB:187:LEU:HD11	35:BB:204:ASN:O	2.17	0.45
36:BC:22:TRP:CH2	36:BC:32:LEU:HB2	2.52	0.45
37:BD:105:VAL:HG21	37:BD:126:ILE:HD12	1.98	0.45
39:BF:19:LEU:HD11	39:BF:59:TYR:CD2	2.52	0.45
40:BG:92:SER:HA	40:BG:93:PRO:HD2	1.66	0.45
41:BH:10:LEU:HD23	41:BH:10:LEU:H	1.81	0.45
57:BZ:312:LEU:O	57:BZ:328:ILE:HA	2.17	0.45
57:BZ:624:LEU:HA	57:BZ:624:LEU:HD12	1.69	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:C1:94:LEU:O	25:C1:97:LEU:HB2	2.16	0.45
1:CA:1512:U:H2'	1:CA:1513:C:C6	2.52	0.45
1:CA:184:C:H2'	1:CA:185:U:H6	1.81	0.45
1:CA:1911:U:H2'	1:CA:1918:A:N1	2.31	0.45
1:CA:2439:A:C5'	1:CA:2439:A:C8	2.97	0.45
1:CA:422:A:H2'	1:CA:423:A:C8	2.51	0.45
1:CA:671:C:H2'	1:CA:672:C:C6	2.52	0.45
1:CA:96:G:OP1	26:C2:46:GLN:NE2	2.50	0.45
2:CB:72:G:O2'	2:CB:105:A:N6	2.45	0.45
14:CQ:109:VAL:HG22	14:CQ:113:GLN:OE1	2.16	0.45
22:CY:86:ARG:HG3	22:CY:100:ALA:HB2	1.99	0.45
34:DA:1327:C:H2'	34:DA:1328:C:C6	2.52	0.45
34:DA:1466:C:H2'	34:DA:1467:G:O4'	2.17	0.45
34:DA:293:G:C5	34:DA:294:U:C5	3.05	0.45
34:DA:57:G:N2	34:DA:58:C:O2	2.49	0.45
34:DA:596:C:H2'	34:DA:597:G:H8	1.82	0.45
34:DA:61:G:H2'	34:DA:62:U:O4'	2.17	0.45
34:DA:931:C:H1'	34:DA:1387:G:N2	2.32	0.45
34:DA:967:C:H2'	34:DA:968:A:N7	2.31	0.45
46:DM:65:LYS:NZ	46:DM:73:GLU:OE2	2.50	0.45
50:DQ:81:ARG:HH21	50:DQ:84:LEU:HD21	1.82	0.45
53:DT:43:LEU:HD13	53:DT:51:GLU:HB3	1.99	0.45
57:DZ:20:HIS:HB2	57:DZ:118:SER:CB	2.47	0.45
57:DZ:223:PHE:HD2	57:DZ:245:ALA:O	2.00	0.45
57:DZ:138:LYS:HA	62:DZ:704:GDP:C6	2.51	0.45
26:A2:8:LYS:HD3	26:A2:8:LYS:HA	1.48	0.44
27:A3:31:LEU:HD23	27:A3:31:LEU:HA	1.44	0.44
1:AA:2154:U:C6	3:AC:6:LYS:CB	3.00	0.44
1:AA:2339:A:H2'	1:AA:2340:A:C8	2.53	0.44
7:AG:7:LEU:HD12	7:AG:104:GLU:HA	2.00	0.44
10:AL:74:ALA:O	10:AL:78:ILE:HG22	2.17	0.44
34:BA:300:A:H2'	34:BA:301:G:O4'	2.16	0.44
34:BA:345:C:H4'	34:BA:346:G:C6	2.52	0.44
34:BA:42:G:H5''	63:BA:5286:HOH:O	2.16	0.44
34:BA:738:C:H2'	34:BA:739:C:C6	2.44	0.44
35:BB:12:GLU:O	35:BB:14:GLY:N	2.50	0.44
35:BB:189:ASP:OD1	35:BB:189:ASP:N	2.46	0.44
45:BL:45:PRO:HB3	45:BL:92:ASP:HB3	1.99	0.44
50:BQ:45:HIS:CE1	50:BQ:47:PRO:HG3	2.51	0.44
55:BV:16:U:O5'	55:BV:16:U:C6	2.70	0.44
56:BW:75:C:H2'	56:BW:76:A:C2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:BZ:114:VAL:O	57:BZ:118:SER:HB2	2.16	0.44
57:BZ:485:GLU:HG3	57:BZ:555:LEU:HD12	1.98	0.44
1:CA:1378:A:OP1	31:C7:10:ARG:NH2	2.50	0.44
1:CA:2557:G:H2'	1:CA:2558:C:H6	1.82	0.44
1:CA:2721:A:H2'	1:CA:2722:G:O4'	2.16	0.44
1:CA:300:A:P	22:CY:86:ARG:HH21	2.40	0.44
1:CA:465:G:OP1	31:C7:12:ARG:NH2	2.46	0.44
1:CA:814:C:H2'	1:CA:815:C:H6	1.83	0.44
3:CC:211:ARG:HH11	3:CC:211:ARG:HG2	1.81	0.44
6:CF:185:ASP:HA	6:CF:188:ARG:HD3	1.98	0.44
8:CH:98:LEU:HD22	8:CH:125:VAL:HG23	1.99	0.44
10:CL:16:LYS:HB2	10:CL:16:LYS:HE3	1.71	0.44
1:CA:1096:A:H2	10:CL:22:PRO:HD3	1.82	0.44
11:CN:38:HIS:CE1	11:CN:39:ARG:HG3	2.52	0.44
22:CY:95:LYS:HE3	22:CY:95:LYS:HB3	1.69	0.44
34:DA:114:U:O2'	34:DA:115:G:H5'	2.17	0.44
34:DA:1258:G:O2'	34:DA:1259:C:H5'	2.16	0.44
34:DA:646:U:H2'	34:DA:647:C:H6	1.82	0.44
34:DA:736:C:H2'	34:DA:737:A:H8	1.78	0.44
34:DA:885:G:O2'	34:DA:914:A:N1	2.48	0.44
37:DD:63:LYS:NZ	37:DD:197:PRO:O	2.41	0.44
38:DE:72:GLN:C	38:DE:73:ASN:HD22	2.20	0.44
57:DZ:18:ALA:HB1	57:DZ:121:VAL:HG21	1.99	0.44
57:DZ:232:LEU:HD13	57:DZ:232:LEU:HA	1.86	0.44
57:DZ:327:PHE:CE1	57:DZ:376:ALA:HB2	2.52	0.44
1:AA:1347:A:C8	1:AA:1349:G:C8	3.05	0.44
1:AA:1636:U:H2'	1:AA:1637:G:H8	1.82	0.44
7:AG:140:ILE:CD1	7:AG:140:ILE:H	2.30	0.44
1:AA:2417:G:OP1	13:AP:77:ARG:NH2	2.50	0.44
16:AS:24:LEU:HA	16:AS:24:LEU:HD23	1.79	0.44
18:AU:108:GLU:O	18:AU:112:ARG:HG2	2.17	0.44
34:BA:15:G:H2'	34:BA:16:A:H8	1.82	0.44
34:BA:256:U:H2'	34:BA:257:G:C8	2.53	0.44
34:BA:262:A:C6	34:BA:263:A:C6	3.05	0.44
34:BA:541:G:C6	34:BA:542:G:C5	3.05	0.44
38:BE:144:THR:OG1	38:BE:146:ALA:HB3	2.18	0.44
41:BH:5:PRO:O	41:BH:8:ASP:HB3	2.17	0.44
43:BJ:70:ARG:HD3	43:BJ:70:ARG:HA	1.80	0.44
57:BZ:-7:GLU:O	57:BZ:-6:ARG:NH1	2.49	0.44
1:CA:1365:A:OP1	25:C1:41:ARG:NH1	2.50	0.44
28:C4:59:PHE:HA	28:C4:61:ARG:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1480:G:C6	1:CA:1481:U:C4	3.05	0.44
1:CA:1489:U:H5'	1:CA:1490:A:OP1	2.18	0.44
1:CA:2075:U:OP2	1:CA:2238:G:O2'	2.29	0.44
1:CA:2846:G:H2'	1:CA:2847:U:O4'	2.17	0.44
1:CA:2850:A:C6	1:CA:2851:A:C6	3.06	0.44
1:CA:793:A:OP2	1:CA:2071:A:O2'	2.32	0.44
3:CC:203:GLU:CD	3:CC:203:GLU:N	2.70	0.44
3:CC:31:LYS:HG2	3:CC:31:LYS:H	1.57	0.44
3:CC:7:ARG:HH22	3:CC:219:MET:HB2	1.81	0.44
5:CE:68:ALA:C	5:CE:70:ALA:H	2.20	0.44
7:CG:168:GLU:O	7:CG:171:ALA:HB3	2.17	0.44
8:CH:56:SER:HB3	8:CH:58:GLU:HG2	2.00	0.44
34:DA:1040:U:C2	34:DA:1041:A:C8	3.05	0.44
34:DA:1203:C:OP1	47:DN:3:ARG:HG3	2.17	0.44
34:DA:130:A:N6	34:DA:234:C:O4'	2.51	0.44
34:DA:432:A:N7	34:DA:433:C:C4	2.85	0.44
38:DE:18:ARG:HH21	38:DE:25:ARG:HG2	1.80	0.44
49:DP:9:PHE:N	49:DP:16:HIS:O	2.50	0.44
46:DM:124:PRO:O	57:DZ:576:ASP:HB2	2.17	0.44
57:DZ:607:ARG:HH22	57:DZ:672:PHE:HD2	1.65	0.44
57:DZ:78:ARG:NH1	57:DZ:357:ARG:NH2	2.66	0.44
25:A1:50:ARG:HG2	25:A1:59:THR:HB	1.99	0.44
28:A4:40:HIS:HA	28:A4:41:PRO:HD2	1.52	0.44
31:A7:24:THR:HG23	63:A7:203:HOH:O	2.17	0.44
1:AA:2146:G:H2'	1:AA:2147:G:O4'	2.18	0.44
1:AA:2240:G:OP1	4:AD:261:LYS:HE2	2.16	0.44
1:AA:2430:A:H2'	1:AA:2431:U:C6	2.52	0.44
1:AA:895:G:C4	1:AA:978:A:H8	2.35	0.44
2:AB:63:G:H2'	2:AB:64:C:H6	1.81	0.44
3:AC:206:LYS:HB3	3:AC:206:LYS:HZ3	1.81	0.44
3:AC:39:ASP:O	3:AC:178:LYS:HE3	2.17	0.44
5:AE:51:PHE:O	5:AE:75:VAL:HG13	2.18	0.44
9:AK:88:ALA:C	9:AK:90:ALA:H	2.20	0.44
10:AL:22:PRO:O	10:AL:27:LEU:HD13	2.17	0.44
12:AO:118:ALA:HA	12:AO:119:PRO:HD3	1.87	0.44
5:AE:9:VAL:HG23	17:AT:3:ARG:HG2	1.98	0.44
19:AV:69:LYS:HG3	19:AV:70:ILE:N	2.32	0.44
21:AX:57:LEU:HD13	21:AX:78:LYS:HB2	2.00	0.44
14:AQ:62:GLY:H	23:AZ:178:GLU:HB2	1.83	0.44
34:BA:115:G:H4'	34:BA:116:A:O5'	2.17	0.44
34:BA:38:G:C2	34:BA:397:A:C2	3.06	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:49:U:O4	34:BA:365:U:H5	2.01	0.44
34:BA:625:G:H2'	34:BA:626:U:H6	1.83	0.44
36:BC:6:HIS:ND1	47:BN:49:HIS:HB3	2.33	0.44
37:BD:13:ARG:NH1	37:BD:13:ARG:HB3	2.32	0.44
38:BE:110:LEU:HD13	38:BE:118:ILE:HG21	1.99	0.44
48:BO:8:LYS:O	48:BO:12:ILE:HG13	2.17	0.44
53:BT:29:LYS:O	53:BT:33:ILE:HG13	2.17	0.44
57:BZ:125:ALA:HB1	57:BZ:132:ARG:HH21	1.83	0.44
57:BZ:637:ARG:C	57:BZ:639:ASN:N	2.70	0.44
1:CA:2432:A:N1	25:C1:35:THR:HG22	2.32	0.44
33:C9:2:LYS:HD3	33:C9:4:ARG:NH2	2.32	0.44
1:CA:1045:A:O4'	1:CA:1047:G:H8	2.00	0.44
1:CA:1912:A:C8	1:CA:1918:A:C2	3.05	0.44
1:CA:2218:U:O2	25:C1:52:ARG:NE	2.48	0.44
1:CA:2306:C:H3'	1:CA:2307:G:C8	2.51	0.44
1:CA:2857:G:C2	1:CA:2861:G:C6	3.05	0.44
3:CC:55:SER:C	3:CC:57:GLN:N	2.71	0.44
5:CE:55:ASN:HA	5:CE:56:PRO:HD3	1.88	0.44
6:CF:119:ARG:HB3	6:CF:119:ARG:CZ	2.45	0.44
9:CK:4:LYS:HA	9:CK:5:ARG:HA	1.78	0.44
13:CP:27:HIS:NE2	63:CP:311:HOH:O	2.35	0.44
34:DA:124:G:H4'	34:DA:291:C:O2'	2.17	0.44
34:DA:587:G:C2	34:DA:755:G:C5	3.06	0.44
34:DA:654:G:H2'	34:DA:655:A:O4'	2.17	0.44
35:DB:135:GLN:O	35:DB:139:LYS:HB2	2.17	0.44
43:DJ:78:ASN:C	43:DJ:80:LYS:H	2.19	0.44
45:DL:45:PRO:HB2	45:DL:92:ASP:HB3	1.98	0.44
49:DP:60:LEU:HD13	49:DP:60:LEU:HA	1.80	0.44
53:DT:36:LEU:HD13	53:DT:36:LEU:HA	1.78	0.44
57:DZ:223:PHE:CE2	57:DZ:249:GLY:HA3	2.52	0.44
57:DZ:90:PHE:CD1	61:DZ:703:FUA:H121	2.53	0.44
1:AA:861:C:H4'	1:AA:1270:C:O2	2.18	0.44
1:AA:2747:A:H2'	1:AA:2748:G:O4'	2.17	0.44
3:AC:24:ASP:C	3:AC:24:ASP:OD1	2.55	0.44
3:AC:22:THR:HG23	3:AC:25:GLU:OE1	2.17	0.44
8:AH:88:LEU:HD22	8:AH:165:ALA:HA	1.98	0.44
9:AK:103:GLY:HA2	9:AK:110:GLY:HA3	1.98	0.44
10:AL:108:ALA:O	10:AL:111:LYS:N	2.47	0.44
16:AS:25:ARG:HD3	16:AS:42:ASP:OD2	2.18	0.44
17:AT:14:TYR:HB2	17:AT:57:PHE:CE1	2.53	0.44
19:AV:72:VAL:HG13	19:AV:85:LYS:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:AX:60:ARG:HH22	31:A7:47:ARG:HH21	1.65	0.44
34:BA:1066:C:C2'	34:BA:1067:A:H5'	2.47	0.44
34:BA:1101:A:OP2	35:BB:96:ARG:HD3	2.17	0.44
34:BA:1280:A:H5'	43:BJ:40:LEU:HD22	1.99	0.44
34:BA:1468:A:H5''	34:BA:1469:G:OP2	2.17	0.44
37:BD:166:LYS:N	37:BD:168:ARG:HH12	2.16	0.44
38:BE:89:ILE:HD12	38:BE:90:VAL:H	1.82	0.44
44:BK:48:ILE:O	44:BK:50:TYR:N	2.41	0.44
50:BQ:81:ARG:HA	50:BQ:81:ARG:HD2	1.60	0.44
52:BS:15:LEU:O	52:BS:19:VAL:HG23	2.17	0.44
53:BT:63:ILE:HG22	53:BT:77:ALA:HB1	1.99	0.44
57:BZ:340:TYR:CD1	57:BZ:349:LYS:HD3	2.52	0.44
32:C8:33:ASN:HA	32:C8:36:LYS:HG3	2.00	0.44
1:CA:191:A:C2	1:CA:192:C:C2	3.06	0.44
1:CA:2166:G:H3'	1:CA:2167:U:C5'	2.47	0.44
1:CA:2415:G:C6	1:CA:2416:C:C4	3.05	0.44
1:CA:2418:A:OP2	32:C8:29:LYS:NZ	2.30	0.44
1:CA:469:G:H2'	1:CA:470:A:H5''	2.00	0.44
1:CA:638:G:H2'	1:CA:639:U:O4'	2.17	0.44
1:CA:921:G:C5	1:CA:922:U:C4	3.06	0.44
1:CA:2749:A:H1'	8:CH:63:SER:HB3	1.99	0.44
11:CN:23:LEU:HA	11:CN:60:ILE:HD11	2.00	0.44
12:CO:87:ILE:HD12	12:CO:91:LEU:HA	1.99	0.44
1:CA:2493:U:O2'	14:CQ:80:GLU:OE1	2.20	0.44
17:CT:26:ASP:OD1	17:CT:120:ARG:NH2	2.45	0.44
20:CW:12:ILE:HG13	20:CW:42:ARG:HH11	1.82	0.44
34:DA:1118:C:H1'	34:DA:1179:A:C5	2.52	0.44
34:DA:1165:C:H2'	34:DA:1166:G:O4'	2.16	0.44
34:DA:426:G:OP1	37:DD:36:ARG:HD2	2.17	0.44
34:DA:965:A:H5'	34:DA:969:A:O4'	2.16	0.44
38:DE:33:VAL:HG13	38:DE:112:LEU:HD12	1.98	0.44
40:DG:50:ILE:HG22	40:DG:125:MET:HG3	1.99	0.44
34:DA:952:U:O4	46:DM:104:ARG:HD3	2.17	0.44
51:DR:47:THR:OG1	51:DR:47:THR:O	2.34	0.44
57:DZ:515:GLU:HG3	57:DZ:564:LYS:HB3	2.00	0.44
24:A0:48:GLY:HA3	24:A0:80:HIS:ND1	2.33	0.44
1:AA:1400:A:H2'	1:AA:1401:G:O4'	2.17	0.44
1:AA:1617:A:H2'	1:AA:1618:A:C8	2.53	0.44
1:AA:596:G:N1	1:AA:2053:A:OP2	2.35	0.44
1:AA:704:U:H2'	1:AA:705:C:C6	2.52	0.44
6:AF:41:LEU:HD23	6:AF:41:LEU:HA	1.78	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AN:63:THR:O	11:AN:66:LYS:HG3	2.18	0.44
1:AA:2856:G:OP2	17:AT:54:ARG:HB2	2.18	0.44
23:AZ:156:LYS:HG2	23:AZ:157:LEU:N	2.33	0.44
34:BA:108:G:O6	53:BT:15:ARG:HD2	2.17	0.44
34:BA:1165:C:H2'	34:BA:1166:G:O4'	2.17	0.44
34:BA:606:G:N2	34:BA:631:G:N7	2.65	0.44
35:BB:164:VAL:HB	35:BB:186:ALA:HB2	1.99	0.44
44:BK:95:ILE:O	44:BK:99:GLN:HG3	2.18	0.44
49:BP:67:THR:C	49:BP:69:THR:H	2.21	0.44
56:BY:28:G:H2'	56:BY:29:G:C8	2.52	0.44
57:BZ:100:VAL:HG12	57:BZ:100:VAL:O	2.17	0.44
57:BZ:573:HIS:HB3	57:BZ:577:SER:H	1.82	0.44
32:C8:8:LYS:HB3	32:C8:12:LYS:HE3	2.00	0.44
1:CA:1239:G:H2'	1:CA:1240:U:O4'	2.17	0.44
1:CA:2056:G:N2	29:C5:5:PRO:HA	2.33	0.44
1:CA:2658:C:O3'	8:CH:158:HIS:HE1	2.00	0.44
1:CA:2830:G:O2'	1:CA:2883:A:N1	2.43	0.44
1:CA:652(C):G:H5''	1:CA:652(D):C:OP2	2.17	0.44
1:CA:65:C:H2'	1:CA:66:C:C6	2.52	0.44
1:CA:864:G:N2	1:CA:913:U:C2	2.86	0.44
2:CB:100:A:H3'	2:CB:101:G:C8	2.52	0.44
3:CC:179:ALA:O	3:CC:180:SER:O	2.35	0.44
10:CL:12:LEU:HD11	10:CL:23:VAL:HG21	1.98	0.44
12:CO:92:GLU:HA	12:CO:93:PRO:HD2	1.79	0.44
22:CY:9:LYS:HA	22:CY:10:GLY:HA2	1.57	0.44
34:DA:839:U:H5''	34:DA:840:C:H5	1.82	0.44
37:DD:117:ALA:O	37:DD:121:VAL:HG23	2.18	0.44
42:DI:18:PHE:O	42:DI:61:ALA:HA	2.17	0.44
49:DP:23:ASP:CG	49:DP:25:ARG:HH11	2.21	0.44
56:DW:54:5MU:H73	56:DW:55:PSU:C2	2.53	0.44
56:DY:55:PSU:HN1	56:DY:57:G:H5'	1.82	0.44
57:DZ:137:ASN:OD1	57:DZ:263:ALA:N	2.38	0.44
57:DZ:-55:LEU:HD23	57:DZ:-55:LEU:HA	1.82	0.44
57:DZ:8:ASP:O	57:DZ:10:LYS:N	2.50	0.44
33:A9:3:VAL:C	33:A9:4:ARG:HG3	2.37	0.44
1:AA:1135:G:OP2	1:AA:1135:G:H2'	2.17	0.44
1:AA:1358:U:C2	1:AA:1649:A:C2	3.05	0.44
1:AA:1405:A:H2	1:AA:1418:U:O4	2.01	0.44
1:AA:2357:G:N3	1:AA:2393:C:H2'	2.33	0.44
1:AA:354:A:H2	1:AA:1255:A:C2'	2.31	0.44
2:AB:111:G:H2'	2:AB:112:U:C6	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AB:15:A:OP2	2:AB:69:G:N2	2.45	0.44
2:AB:32:C:C2	2:AB:51:G:C2	3.05	0.44
3:AC:55:SER:C	3:AC:57:GLN:N	2.71	0.44
4:AD:20:ASP:OD1	4:AD:21:PHE:N	2.51	0.44
14:AQ:109:VAL:CG1	14:AQ:113:GLN:HB3	2.43	0.44
19:AV:39:LEU:C	19:AV:39:LEU:HD23	2.38	0.44
34:BA:1388:C:H2'	34:BA:1389:C:H6	1.82	0.44
34:BA:47:C:C6	34:BA:365:U:H2'	2.53	0.44
34:BA:654:G:C4	34:BA:753:A:C6	3.04	0.44
34:BA:909:A:H2'	34:BA:910:C:O4'	2.16	0.44
37:BD:194:LEU:HD12	37:BD:195:ALA:N	2.33	0.44
40:BG:137:LYS:HG3	40:BG:137:LYS:O	2.17	0.44
41:BH:100:ILE:HA	41:BH:101:PRO:HD2	1.77	0.44
41:BH:51:VAL:HG11	41:BH:60:ARG:HH11	1.83	0.44
45:BL:28:LYS:HG3	45:BL:62:SER:HB2	2.00	0.44
50:BQ:29:HIS:HA	50:BQ:30:PRO:HD2	1.74	0.44
57:BZ:556:ILE:HG13	57:BZ:558:PHE:HD2	1.82	0.44
57:BZ:-9:LEU:O	57:BZ:-6:ARG:HB2	2.17	0.44
1:CA:1031:G:H21	33:C9:36:GLN:HE22	1.65	0.44
1:CA:2056:G:H2'	1:CA:2056:G:N3	2.33	0.44
1:CA:2318:G:N3	1:CA:2318:G:O2'	2.40	0.44
1:CA:584:C:OP2	18:CU:6:THR:OG1	2.21	0.44
3:CC:16:ASP:OD2	3:CC:19:LYS:HB2	2.17	0.44
1:CA:2591:C:OP1	4:CD:239:ARG:HD2	2.18	0.44
4:CD:77:ALA:HA	4:CD:97:TYR:HA	1.99	0.44
5:CE:22:PRO:O	5:CE:185:LYS:HA	2.17	0.44
8:CH:7:LEU:HA	8:CH:8:PRO:HD3	1.82	0.44
17:CT:11:GLU:O	17:CT:15:VAL:HG23	2.18	0.44
22:CY:68:HIS:HB3	22:CY:71:LYS:HG3	1.99	0.44
23:CZ:98:MET:HB2	23:CZ:98:MET:HE2	1.73	0.44
34:DA:1079:G:C6	34:DA:1080:A:N6	2.85	0.44
34:DA:1226:C:OP2	46:DM:91:ARG:NH1	2.36	0.44
34:DA:409:G:H1	34:DA:433:C:N4	2.15	0.44
34:DA:566:G:H4'	34:DA:567:G:OP1	2.18	0.44
34:DA:1104:G:H4'	35:DB:111:ARG:NH1	2.32	0.44
36:DC:130:VAL:HG11	36:DC:153:VAL:HG11	2.00	0.44
43:DJ:40:LEU:HB2	43:DJ:69:ASN:HB3	2.00	0.44
57:DZ:515:GLU:OE2	57:DZ:564:LYS:HD3	2.18	0.44
30:A6:30:THR:HG22	30:A6:30:THR:O	2.18	0.44
1:AA:174:U:H4'	1:AA:207:A:H4'	2.00	0.44
1:AA:644:G:O6	6:AF:103:LYS:HE3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1828:C:H4'	4:AD:257:LEU:O	2.18	0.44
6:AF:106:ARG:H	6:AF:106:ARG:HG2	1.60	0.44
8:AH:167:GLU:HA	8:AH:168:PRO:HD3	1.74	0.44
10:AL:29:GLN:O	10:AL:59:ILE:HD12	2.17	0.44
12:AO:111:PHE:O	12:AO:115:VAL:HG23	2.18	0.44
34:BA:1189:C:H5''	34:BA:1190:G:OP2	2.18	0.44
17:AT:41:ARG:NH1	34:BA:346:G:OP1	2.47	0.44
37:BD:4:TYR:O	37:BD:5:ILE:HG22	2.17	0.44
37:BD:78:LEU:HA	37:BD:78:LEU:HD23	1.77	0.44
40:BG:113:GLU:H	40:BG:113:GLU:HG2	1.39	0.44
53:BT:60:GLU:O	53:BT:63:ILE:HB	2.17	0.44
56:BW:18:G:H4'	56:BW:60:U:C6	2.53	0.44
57:BZ:292:THR:HA	57:BZ:298:VAL:HA	2.00	0.44
57:BZ:349:LYS:HG2	57:BZ:350:GLU:N	2.33	0.44
34:BA:1494:G:O3'	57:BZ:499:ARG:NH1	2.50	0.44
57:BZ:11:ARG:O	57:BZ:77:HIS:HA	2.17	0.44
24:C0:82:ARG:HB2	24:C0:82:ARG:HH11	1.83	0.44
33:C9:17:ILE:HG21	33:C9:19:ARG:HH21	1.83	0.44
1:CA:1071:G:H1'	1:CA:1089:G:C8	2.53	0.44
1:CA:1443:G:N2	1:CA:1549:C:C2	2.86	0.44
1:CA:1300:U:C2	1:CA:1626:G:C6	3.06	0.44
1:CA:1820:U:H4'	1:CA:1821:A:OP2	2.17	0.44
1:CA:1860:G:P	3:CC:206:LYS:HE2	2.52	0.44
1:CA:487:C:C2'	1:CA:488:G:H5'	2.48	0.44
2:CB:7:G:H4'	16:CS:29:PHE:CD2	2.53	0.44
5:CE:102:VAL:HB	5:CE:103:ASP:H	1.58	0.44
7:CG:36:LYS:HD3	7:CG:95:ARG:CZ	2.47	0.44
1:CA:1063:G:N3	10:CL:91:PRO:HG2	2.33	0.44
14:CQ:58:PHE:O	14:CQ:59:ARG:HB2	2.18	0.44
1:CA:2723:C:H4'	15:CR:1:MET:HG3	2.00	0.44
1:CA:2318:G:N2	16:CS:3:ARG:HH12	2.15	0.44
17:CT:109:GLU:HG2	17:CT:112:ARG:HH22	1.81	0.44
21:CX:26:TYR:HB3	21:CX:92:LEU:HD12	1.98	0.44
21:CX:61:GLY:HA3	21:CX:73:ARG:O	2.18	0.44
22:CY:77:PRO:CD	22:CY:106:LEU:HD23	2.47	0.44
23:CZ:166:SER:HA	23:CZ:167:PRO:HD3	1.81	0.44
23:CZ:89:PHE:CE1	23:CZ:96:VAL:HG21	2.52	0.44
34:DA:1091:U:C2	34:DA:1095:U:C4	3.06	0.44
34:DA:1128:C:H1'	34:DA:1148:U:H3	1.83	0.44
34:DA:1153:C:H2'	34:DA:1154:G:H5''	1.99	0.44
34:DA:1355:G:H2'	34:DA:1356:G:O4'	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:32:A:C2	34:DA:33:A:C4	3.05	0.44
34:DA:546:G:P	37:DD:72:GLU:HB3	2.57	0.44
35:DB:104:ASN:O	35:DB:108:ILE:HG12	2.18	0.44
38:DE:30:ALA:O	38:DE:45:PHE:HD1	2.01	0.44
40:DG:105:VAL:O	40:DG:108:ALA:HB3	2.17	0.44
41:DH:20:TYR:HD2	41:DH:65:TYR:CD2	2.35	0.44
45:DL:105:TYR:O	45:DL:107:ALA:N	2.51	0.44
57:DZ:453:GLY:HA3	57:DZ:459:LEU:HG	2.00	0.44
57:DZ:540:PRO:O	57:DZ:543:GLN:HB3	2.18	0.44
10:CL:21:PRO:CG	57:DZ:614:GLU:HB2	2.48	0.44
29:A5:16:ARG:HD2	29:A5:20:ARG:NH1	2.33	0.44
1:AA:1036:A:H5''	1:AA:1037:C:P	2.58	0.44
1:AA:1157:A:O2'	1:AA:1158:G:H4'	2.18	0.44
1:AA:1496:A:H5'	1:AA:1497:G:OP2	2.17	0.44
1:AA:2282:G:H2'	1:AA:2283:G:O4'	2.17	0.44
1:AA:237:G:OP1	63:AA:4921:HOH:O	2.21	0.44
1:AA:2784:C:H2'	1:AA:2785:C:C6	2.53	0.44
1:AA:282:G:H2'	1:AA:283:G:O4'	2.18	0.44
4:AD:16:MET:HG3	4:AD:206:LEU:O	2.18	0.44
9:AK:25:PHE:O	9:AK:84:GLU:HA	2.17	0.44
10:AL:103:GLN:O	10:AL:106:GLU:HG3	2.18	0.44
15:AR:70:LEU:HD23	15:AR:70:LEU:HA	1.66	0.44
21:AX:5:TYR:CZ	26:A2:30:ARG:HB2	2.53	0.44
23:AZ:31:ARG:HG3	23:AZ:31:ARG:H	1.28	0.44
34:BA:1027:C:H5''	34:BA:1028:C:OP2	2.18	0.44
34:BA:299:G:C6	34:BA:300:A:C6	3.06	0.44
34:BA:453:A:H62	34:BA:479:C:H42	1.66	0.44
34:BA:612:C:O2	34:BA:629:G:N2	2.51	0.44
34:BA:673:G:H2'	34:BA:674:G:C8	2.53	0.44
34:BA:960:U:H1'	34:BA:1223:C:H5'	1.99	0.44
36:BC:122:GLU:O	36:BC:126:ARG:NH1	2.44	0.44
36:BC:174:PRO:HD2	36:BC:182:ILE:HD11	2.00	0.44
37:BD:59:ARG:NE	37:BD:59:ARG:HA	2.33	0.44
53:BT:14:LYS:O	53:BT:18:GLN:HG3	2.18	0.44
56:BW:17:C:O2	56:BW:17:C:H2'	2.17	0.44
57:BZ:88:VAL:HG13	57:BZ:117:GLN:HE22	1.83	0.44
57:BZ:191:ASP:HA	57:BZ:265:LYS:O	2.17	0.44
1:CA:1035:U:H2'	1:CA:1036:G:C8	2.53	0.44
1:CA:1355:G:P	4:CD:38:LYS:HE2	2.58	0.44
1:CA:2016:U:H2'	1:CA:2017:U:C6	2.52	0.44
1:CA:2643:G:N2	1:CA:2772:C:C2	2.86	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2886:G:H2'	1:CA:2887:U:H6	1.83	0.44
1:CA:889:C:O2'	1:CA:890:A:O5'	2.34	0.44
3:CC:24:ASP:OD1	3:CC:24:ASP:C	2.55	0.44
7:CG:36:LYS:HD3	7:CG:95:ARG:NH1	2.33	0.44
11:CN:85:ILE:HG21	11:CN:90:MET:HE2	1.99	0.44
16:CS:63:THR:HG23	16:CS:64:GLU:N	2.30	0.44
23:CZ:136:PHE:HA	23:CZ:136:PHE:HD2	1.78	0.44
34:DA:196:A:N3	34:DA:222:U:H1'	2.33	0.44
34:DA:402:G:O2'	34:DA:620:C:N3	2.49	0.44
34:DA:904:C:H2'	34:DA:905:U:O4'	2.17	0.44
35:DB:15:VAL:HG23	35:DB:16:HIS:ND1	2.33	0.44
36:DC:22:TRP:HZ3	36:DC:24:ALA:HB2	1.83	0.44
37:DD:110:PHE:CD1	37:DD:110:PHE:N	2.86	0.44
40:DG:47:CYS:HB3	40:DG:58:PRO:HB3	2.00	0.44
45:DL:59:ARG:NH1	45:DL:65:GLU:OE1	2.51	0.44
57:DZ:149:VAL:O	57:DZ:153:MET:HB2	2.18	0.44
57:DZ:74:TRP:HE1	57:DZ:274:ASP:N	2.15	0.44
27:A3:18:ASP:N	27:A3:18:ASP:OD1	2.42	0.44
31:A7:31:LEU:HD22	31:A7:42:LEU:HB3	1.99	0.44
1:AA:1140:U:H2'	1:AA:1141:A:H3'	2.00	0.44
1:AA:1740:U:H1'	4:AD:14:ARG:NH2	2.33	0.44
1:AA:2225:U:O2'	1:AA:2226:C:H5'	2.17	0.44
1:AA:263:C:H6	1:AA:263:C:H5''	1.83	0.44
4:AD:108:PRO:HG3	4:AD:143:HIS:CE1	2.53	0.44
4:AD:253:GLN:HE21	4:AD:253:GLN:HB3	1.63	0.44
4:AD:71:ASP:OD2	4:AD:103:ARG:NH2	2.47	0.44
6:AF:120:GLU:HG3	6:AF:122:LYS:HG2	2.00	0.44
14:AQ:45:GLN:N	14:AQ:45:GLN:OE1	2.46	0.44
18:AU:74:LEU:HD12	18:AU:74:LEU:N	2.33	0.44
23:AZ:102:LEU:HD11	23:AZ:124:ILE:HB	2.00	0.44
34:BA:1234:C:H2'	34:BA:1235:U:C6	2.53	0.44
34:BA:1340:A:H2'	34:BA:1341:U:O4'	2.18	0.44
34:BA:662:G:O2'	34:BA:836:G:H5'	2.17	0.44
34:BA:967:C:O5'	34:BA:967:C:H6	2.01	0.44
36:BC:178:LEU:HA	36:BC:178:LEU:HD13	1.83	0.44
38:BE:69:VAL:HA	38:BE:70:PRO:HD2	1.83	0.44
49:BP:40:ASP:HA	49:BP:41:PRO:HD2	1.66	0.44
52:BS:32:LYS:HA	52:BS:50:ALA:HB3	1.99	0.44
57:BZ:350:GLU:HG2	57:BZ:380:LEU:HA	2.00	0.44
57:BZ:435:ASP:OD2	57:BZ:437:THR:OG1	2.34	0.44
33:C9:23:VAL:HB	33:C9:36:GLN:NE2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2063:C:O2	1:CA:2450:A:N1	2.50	0.44
1:CA:908:C:OP2	14:CQ:22:LYS:HD3	2.18	0.44
3:CC:194:ILE:HD11	3:CC:227:PRO:HB2	1.99	0.44
4:CD:61:LEU:O	4:CD:63:ARG:NH1	2.51	0.44
5:CE:116:VAL:HG13	5:CE:122:PHE:HB2	2.00	0.44
5:CE:11:MET:HG2	5:CE:24:THR:HG22	1.99	0.44
11:CN:34:LEU:HD21	11:CN:120:LEU:HG	2.00	0.44
15:CR:55:ALA:HA	15:CR:80:PHE:CZ	2.53	0.44
18:CU:97:ASP:OD1	18:CU:101:ARG:NH1	2.51	0.44
22:CY:87:LYS:HD2	22:CY:95:LYS:HD2	1.99	0.44
34:DA:1089:G:C6	34:DA:1090:U:C4	3.06	0.44
34:DA:1142:G:H3'	34:DA:1143:G:C8	2.52	0.44
36:DC:104:GLN:HE21	36:DC:104:GLN:HB3	1.62	0.44
38:DE:51:VAL:O	38:DE:55:VAL:HG23	2.18	0.44
44:DK:72:ALA:HB1	44:DK:77:MET:HB2	1.98	0.44
44:DK:20:TYR:CD1	44:DK:83:ILE:HB	2.53	0.44
45:DL:69:TYR:CD2	45:DL:99:HIS:CE1	3.06	0.44
46:DM:3:ARG:O	46:DM:57:ARG:NH2	2.47	0.44
47:DN:2:ALA:O	47:DN:6:LEU:HD22	2.18	0.44
51:DR:26:LEU:HD23	51:DR:29:PHE:CE1	2.52	0.44
57:DZ:8:ASP:C	57:DZ:10:LYS:H	2.21	0.44
57:DZ:363:ARG:CG	57:DZ:363:ARG:NH1	2.78	0.44
57:DZ:74:TRP:NE1	57:DZ:273:LEU:C	2.72	0.44
32:A8:62:LEU:HB3	32:A8:65:GLU:CG	2.48	0.43
1:AA:1067:A:C3'	1:AA:1067:A:C8	2.98	0.43
1:AA:1074:A:N3	1:AA:2498:G:O2'	2.37	0.43
1:AA:1154:U:H1'	1:AA:1155:C:OP1	2.18	0.43
1:AA:2191:A:N3	1:AA:2191:A:H2'	2.33	0.43
1:AA:2200:C:H4'	3:AC:47:LYS:HZ3	1.80	0.43
1:AA:1757:C:O2'	1:AA:2868:C:N3	2.44	0.43
1:AA:734:C:H2'	1:AA:735:U:O4'	2.18	0.43
1:AA:967:G:H4'	1:AA:2281:A:C5	2.53	0.43
1:AA:989:G:H5''	1:AA:990:A:H5'	2.00	0.43
3:AC:194:ILE:HD11	3:AC:227:PRO:HB2	1.99	0.43
4:AD:13:ARG:HD2	4:AD:13:ARG:HA	1.69	0.43
4:AD:72:LYS:HD3	4:AD:97:TYR:CE2	2.53	0.43
7:AG:103:LEU:HA	7:AG:103:LEU:HD23	1.78	0.43
8:AH:92:ILE:HA	8:AH:92:ILE:HD13	1.74	0.43
10:AL:23:VAL:HG13	10:AL:27:LEU:HD22	2.00	0.43
10:AL:40:ALA:HB3	10:AL:67:PHE:HZ	1.82	0.43
34:BA:1133:G:H2'	34:BA:1134:G:H8	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:1229:A:H8	34:BA:1229:A:O5'	2.01	0.43
34:BA:1429:C:H2'	34:BA:1430:C:H6	1.81	0.43
34:BA:408:A:H2'	34:BA:409:G:H8	1.82	0.43
34:BA:413:G:H1'	34:BA:428:G:N2	2.32	0.43
34:BA:604:G:N2	34:BA:635:G:C4	2.86	0.43
34:BA:689:C:OP2	44:BK:55:LYS:NZ	2.51	0.43
34:BA:696:A:O2'	34:BA:697:U:H5'	2.18	0.43
34:BA:735:C:H2'	34:BA:736:C:C6	2.53	0.43
37:BD:57:ARG:HE	37:BD:205:GLU:HB2	1.83	0.43
45:BL:26:ALA:HB3	45:BL:98:TYR:CZ	2.53	0.43
46:BM:17:VAL:O	46:BM:20:THR:OG1	2.28	0.43
49:BP:2:VAL:O	49:BP:64:ALA:HA	2.17	0.43
49:BP:36:ILE:HD12	49:BP:56:ALA:HB2	1.99	0.43
57:BZ:225:GLU:HG2	57:BZ:225:GLU:H	1.63	0.43
57:BZ:413:ILE:HA	57:BZ:476:VAL:HG12	1.99	0.43
57:BZ:673:PHE:HZ	57:BZ:676:TYR:CD2	2.35	0.43
25:C1:25:LYS:C	25:C1:27:GLU:N	2.70	0.43
1:CA:1290:C:H2'	1:CA:1291:C:C6	2.54	0.43
1:CA:2360:A:H8	1:CA:2360:A:O5'	2.01	0.43
1:CA:2784:C:H1'	5:CE:37:ARG:HH12	1.82	0.43
1:CA:590:A:H2'	1:CA:591:C:C6	2.53	0.43
1:CA:76:C:O3'	26:C2:59:ARG:HG3	2.17	0.43
3:CC:39:ASP:O	3:CC:178:LYS:HE3	2.17	0.43
14:CQ:29:PHE:HB2	14:CQ:105:GLU:OE2	2.18	0.43
17:CT:99:LEU:HA	17:CT:99:LEU:HD23	1.79	0.43
34:DA:1074:G:O2'	34:DA:1101:A:N1	2.44	0.43
34:DA:130:A:H1'	34:DA:263:A:O2'	2.18	0.43
34:DA:1348:U:H2'	34:DA:1349:A:H8	1.83	0.43
34:DA:195:A:OP2	63:DA:3299:HOH:O	2.21	0.43
34:DA:960:U:O2	34:DA:960:U:H2'	2.18	0.43
36:DC:54:ARG:NH1	36:DC:56:ASP:OD2	2.51	0.43
40:DG:71:PRO:HG3	40:DG:103:TRP:CH2	2.53	0.43
51:DR:22:VAL:HG23	51:DR:55:ARG:O	2.18	0.43
52:DS:15:LEU:HD21	52:DS:32:LYS:O	2.17	0.43
57:DZ:170:ARG:HA	57:DZ:170:ARG:HD3	1.68	0.43
57:DZ:238:THR:HG22	57:DZ:241:GLU:OE2	2.17	0.43
57:DZ:316:ILE:CD1	57:DZ:326:THR:HG23	2.48	0.43
57:DZ:629:GLY:HA2	57:DZ:648:PRO:HD3	2.00	0.43
1:AA:2101:U:O3'	25:A1:35:THR:OG1	2.34	0.43
25:A1:54:ALA:O	25:A1:56:GLN:N	2.50	0.43
1:AA:2274:U:P	24:A0:19:LYS:HZ3	2.41	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2327:G:C6	1:AA:2328:C:N4	2.86	0.43
1:AA:2444:A:H2'	1:AA:2445:A:C8	2.53	0.43
1:AA:424:G:H1'	1:AA:2243:C:O2'	2.19	0.43
1:AA:601:A:OP1	1:AA:1301:U:H4'	2.18	0.43
10:AL:21:PRO:HG3	10:AL:25:PRO:HD3	2.00	0.43
15:AR:8:ARG:NH1	15:AR:39:PRO:HB3	2.33	0.43
21:AX:31:HIS:HD2	21:AX:33:LYS:N	1.94	0.43
34:BA:1299:A:O2'	34:BA:1301:U:H5'	2.18	0.43
37:BD:13:ARG:HB2	37:BD:40:PRO:HD3	2.01	0.43
37:BD:206:PHE:C	37:BD:208:SER:H	2.21	0.43
41:BH:35:ILE:O	41:BH:38:ILE:HB	2.17	0.43
42:BI:53:VAL:O	42:BI:55:ALA:N	2.51	0.43
45:BL:44:THR:HA	45:BL:45:PRO:HD3	1.79	0.43
34:BA:189(F):U:C2	50:BQ:72:ARG:NH2	2.86	0.43
51:BR:70:ILE:HG23	51:BR:79:LEU:HD13	1.99	0.43
56:BW:28:G:H1	56:BW:42:C:N4	2.08	0.43
57:BZ:187:THR:HG22	57:BZ:199:ILE:HG12	2.00	0.43
57:BZ:-34:ARG:O	57:BZ:-32:LEU:N	2.51	0.43
1:CA:1056:G:H4'	1:CA:1086:A:C8	2.53	0.43
1:CA:921:G:H4'	1:CA:2269:A:C5	2.53	0.43
1:CA:2652:C:H2'	1:CA:2653:U:O4'	2.18	0.43
1:CA:2717:G:C6	1:CA:2718:G:C5	3.06	0.43
1:CA:2854:G:C2	1:CA:2864:G:C2	3.05	0.43
1:CA:322:A:H5'	1:CA:340:A:H1'	1.99	0.43
1:CA:748:G:C8	20:CW:89:ALA:HB1	2.53	0.43
1:CA:981:A:OP2	1:CA:982:C:N4	2.48	0.43
2:CB:24:G:N3	2:CB:27:C:N4	2.55	0.43
3:CC:54:ARG:HE	3:CC:57:GLN:HG2	1.83	0.43
6:CF:148:LEU:HD11	6:CF:193:VAL:HG21	2.00	0.43
10:CL:10:LEU:HB2	10:CL:55:VAL:HG13	2.00	0.43
16:CS:14:VAL:HG23	16:CS:15:ARG:N	2.33	0.43
16:CS:41:ASP:OD2	16:CS:44:LYS:HE2	2.18	0.43
34:DA:1075:C:H5''	35:DB:179:LYS:HZ1	1.82	0.43
34:DA:1095:U:H2'	34:DA:1096:C:O4'	2.18	0.43
34:DA:941:G:C2	34:DA:1343:G:C2	3.06	0.43
34:DA:1374:A:C4	34:DA:1375:A:C8	3.06	0.43
34:DA:630:G:H2'	34:DA:631:G:H8	1.84	0.43
45:DL:119:LYS:HB2	45:DL:120:TYR:HD2	1.83	0.43
50:DQ:45:HIS:HB3	50:DQ:72:ARG:HG3	1.99	0.43
51:DR:33:ASP:OD2	51:DR:36:ASN:HB2	2.18	0.43
57:DZ:272:LEU:HD12	57:DZ:275:ALA:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:38:ARG:NH1	57:DZ:270:GLN:HE21	2.16	0.43
57:DZ:638:GLY:C	57:DZ:640:ALA:HB3	2.38	0.43
25:A1:85:LEU:HA	25:A1:85:LEU:HD23	1.66	0.43
29:A5:48:GLU:C	29:A5:60:VAL:HG11	2.39	0.43
1:AA:1099:C:C2	1:AA:1153:G:N2	2.86	0.43
1:AA:1414:G:C2	1:AA:1415:G:C8	3.06	0.43
1:AA:1821:C:H2'	1:AA:1822:A:C5	2.54	0.43
1:AA:253:C:O2'	1:AA:254:A:H2'	2.18	0.43
1:AA:2579:G:H2'	1:AA:2580:C:C6	2.53	0.43
1:AA:966:G:C2'	1:AA:967:G:H5'	2.49	0.43
3:AC:60:ARG:HG3	3:AC:165:ARG:HB2	2.01	0.43
12:AO:35:VAL:HA	12:AO:62:VAL:HG12	2.00	0.43
13:AP:19:VAL:HB	13:AP:31:ALA:HB1	2.01	0.43
34:BA:1273:G:H3'	34:BA:1274:G:H8	1.83	0.43
34:BA:313:A:C6	34:BA:314:C:C4	3.07	0.43
34:BA:345:C:H4'	34:BA:346:G:C5	2.53	0.43
34:BA:866:C:C4	34:BA:867:G:H1'	2.54	0.43
34:BA:872:A:C2	34:BA:874:G:C6	3.06	0.43
35:BB:81:VAL:HG12	35:BB:215:LEU:HD11	1.98	0.43
36:BC:148:GLY:HA3	36:BC:172:ARG:O	2.18	0.43
49:BP:6:LEU:HD12	49:BP:19:ILE:HB	1.99	0.43
57:BZ:5:LEU:HD23	57:BZ:5:LEU:HA	1.67	0.43
1:CA:1540:U:C2'	1:CA:1541:G:H5'	2.48	0.43
1:CA:209:C:H2'	1:CA:210:C:O4'	2.18	0.43
1:CA:2351:G:O5'	1:CA:2351:G:H8	2.02	0.43
1:CA:2391:G:O2'	1:CA:2424:C:N4	2.41	0.43
1:CA:9:U:N3	1:CA:2629:A:H2	2.11	0.43
1:CA:2679:A:O2'	1:CA:2680:C:H5'	2.18	0.43
1:CA:2728:U:OP1	12:CO:70:LYS:NZ	2.51	0.43
1:CA:277:C:O2	1:CA:277:C:H2'	2.18	0.43
1:CA:71:A:H5''	1:CA:73:A:C8	2.54	0.43
1:CA:797:C:O5'	1:CA:797:C:H6	2.01	0.43
2:CB:29:A:C2	2:CB:30:C:C2	3.06	0.43
5:CE:14:ILE:HB	17:CT:14:TYR:CZ	2.54	0.43
6:CF:102:PRO:HB2	6:CF:105:VAL:HB	2.01	0.43
6:CF:13:SER:OG	6:CF:15:SER:HB2	2.18	0.43
17:CT:18:ASP:N	17:CT:18:ASP:OD1	2.41	0.43
34:DA:1015:A:N3	34:DA:1218:C:O2'	2.44	0.43
34:DA:1171:G:H8	34:DA:1171:G:OP2	2.01	0.43
34:DA:195:A:OP1	53:DT:68:LYS:NZ	2.50	0.43
34:DA:509:A:C6	34:DA:510:A:N1	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:517:G:N1	34:DA:533:A:OP2	2.50	0.43
34:DA:543:C:H2'	34:DA:544:G:O4'	2.19	0.43
34:DA:577:G:C4	34:DA:816:A:C2	3.06	0.43
34:DA:826:C:H2'	34:DA:827:U:C6	2.53	0.43
35:DB:218:ALA:O	35:DB:222:ILE:HG23	2.18	0.43
41:DH:13:ILE:O	41:DH:17:THR:HG23	2.17	0.43
48:DO:56:LEU:O	48:DO:60:VAL:HG23	2.18	0.43
48:DO:5:LYS:O	48:DO:9:GLN:HG2	2.18	0.43
56:DW:44:G:C5	56:DW:45:U:C4	3.06	0.43
56:DW:18:G:H4'	56:DW:60:U:C5	2.53	0.43
57:DZ:530:VAL:HG13	57:DZ:533:VAL:HG21	2.00	0.43
32:A8:26:LYS:HB3	63:A8:6310:HOH:O	2.19	0.43
33:A9:27:CYS:HB3	33:A9:32:HIS:HB2	1.99	0.43
1:AA:1478:C:H2'	1:AA:1479:U:O4'	2.18	0.43
1:AA:2564:U:H2'	1:AA:2566:U:H5''	2.01	0.43
1:AA:289:G:H2'	1:AA:290:G:C8	2.53	0.43
1:AA:310:C:H2'	1:AA:311:C:C6	2.54	0.43
1:AA:956:A:H2'	1:AA:957:A:C8	2.53	0.43
4:AD:35:LYS:HB2	4:AD:36:PRO:HD2	2.00	0.43
7:AG:26:GLN:HA	7:AG:26:GLN:HE21	1.82	0.43
34:BA:1079:G:H2'	34:BA:1080:A:C8	2.54	0.43
34:BA:625:G:H2'	34:BA:626:U:C6	2.53	0.43
34:BA:837:G:N2	34:BA:850:U:H1'	2.32	0.43
37:BD:101:LEU:C	37:BD:101:LEU:HD12	2.38	0.43
37:BD:173:TRP:CZ3	37:BD:174:LEU:HD11	2.53	0.43
38:BE:41:VAL:HG22	38:BE:69:VAL:HG11	2.00	0.43
41:BH:51:VAL:HG12	41:BH:52:ASP:N	2.28	0.43
34:BA:1225:A:OP1	46:BM:102:ARG:HA	2.19	0.43
53:BT:40:ALA:HB2	53:BT:55:ILE:HG22	2.01	0.43
57:BZ:590:ILE:HD13	57:BZ:593:ALA:HB3	2.00	0.43
1:CA:1268:A:C2'	1:CA:1269:A:O5'	2.66	0.43
1:CA:1530:C:N4	1:CA:1539:G:H1	2.17	0.43
1:CA:1860:G:C8	1:CA:1860:G:OP2	2.72	0.43
1:CA:2275:C:O2	14:CQ:85:LYS:HD2	2.18	0.43
1:CA:2747:G:O6	1:CA:2755:C:H5''	2.19	0.43
1:CA:2758:A:C2	1:CA:2759:G:H1'	2.52	0.43
1:CA:637:A:H2'	13:CP:117:GLU:OE2	2.19	0.43
1:CA:857:C:O2'	1:CA:858:U:H5'	2.18	0.43
1:CA:956:G:OP2	14:CQ:14:ARG:NH2	2.50	0.43
4:CD:94:LEU:HD23	4:CD:94:LEU:HA	1.83	0.43
5:CE:128:SER:OG	5:CE:129:HIS:N	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2053:G:OP1	5:CE:144:ARG:HG2	2.18	0.43
5:CE:169:ASN:HD22	5:CE:203:LYS:CB	2.31	0.43
7:CG:136:ARG:HD2	7:CG:137:GLU:H	1.84	0.43
7:CG:37:VAL:HB	7:CG:94:LEU:HB2	2.00	0.43
1:CA:875:G:H5''	23:CZ:149:SER:OG	2.18	0.43
34:DA:1154:G:C6	34:DA:1155:G:C6	3.06	0.43
34:DA:1256:A:OP2	36:DC:26:LYS:NZ	2.46	0.43
34:DA:533:A:N6	34:DA:536:C:O2	2.51	0.43
34:DA:547:A:H4'	34:DA:548:G:O5'	2.18	0.43
34:DA:841:U:H6	34:DA:841:U:P	2.42	0.43
34:DA:919:A:O2'	34:DA:920:U:H5'	2.19	0.43
36:DC:134:ILE:HD11	36:DC:153:VAL:HG23	2.00	0.43
37:DD:32:ALA:HB3	60:DD:501:SF4:S2	2.58	0.43
40:DG:78:ARG:HH21	40:DG:79:ARG:HH22	1.66	0.43
43:DJ:63:PHE:HZ	47:DN:49:HIS:CD2	2.36	0.43
57:DZ:655:TYR:C	57:DZ:657:THR:H	2.22	0.43
26:A2:51:ARG:O	26:A2:55:ARG:HG3	2.19	0.43
28:A4:57:GLU:HA	28:A4:58:ARG:HA	1.43	0.43
1:AA:1147:U:H2'	1:AA:1148:C:C6	2.53	0.43
1:AA:1464:G:O5'	1:AA:1464:G:H8	2.01	0.43
1:AA:2904:U:H2'	1:AA:2905:C:C6	2.54	0.43
1:AA:309:C:H2'	1:AA:310:C:C6	2.52	0.43
1:AA:887:C:H2'	1:AA:888:A:H8	1.83	0.43
1:AA:990:A:OP2	63:AA:4348:HOH:O	2.21	0.43
10:AL:51:ALA:HB2	10:AL:76:TYR:CE2	2.54	0.43
11:AN:67:LEU:HB3	11:AN:88:GLU:HG3	2.01	0.43
13:AP:105:LEU:H	13:AP:105:LEU:HD12	1.84	0.43
20:AW:54:ALA:CB	20:AW:107:LEU:HD22	2.49	0.43
23:AZ:157:LEU:HD11	23:AZ:163:LEU:HD13	1.99	0.43
34:BA:1075:C:OP1	35:BB:179:LYS:HE3	2.18	0.43
1:AA:1935:A:C6	34:BA:1494:G:H5'	2.53	0.43
34:BA:162:A:C8	34:BA:163:C:H1'	2.53	0.43
34:BA:457:C:N3	34:BA:475:G:C2	2.87	0.43
34:BA:583:A:H2'	34:BA:584:G:O4'	2.19	0.43
34:BA:941:G:C2	34:BA:942:G:H1'	2.53	0.43
40:BG:91:VAL:HG12	40:BG:95:ARG:HB3	1.99	0.43
42:BI:4:TYR:CE2	42:BI:88:TYR:HD1	2.36	0.43
1:CA:857:C:H4'	24:C0:23:VAL:HG21	1.99	0.43
27:C3:7:LYS:O	27:C3:54:VAL:HG22	2.18	0.43
32:C8:50:LEU:HA	32:C8:50:LEU:HD23	1.63	0.43
1:CA:1099:G:C2	1:CA:1100:C:H1'	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1489:U:HO2'	1:CA:1490:A:H8	1.64	0.43
1:CA:226:G:C2	1:CA:227:A:C6	3.06	0.43
1:CA:2835:A:N6	1:CA:2878:U:H3'	2.33	0.43
1:CA:2842:G:C2	1:CA:2876:G:C2	3.06	0.43
1:CA:471:A:O5'	1:CA:471:A:H8	2.01	0.43
1:CA:827:U:H5'	1:CA:828:U:O5'	2.19	0.43
4:CD:157:ARG:O	4:CD:196:VAL:HG11	2.19	0.43
5:CE:52:LEU:HA	5:CE:52:LEU:HD12	1.68	0.43
6:CF:152:GLU:HA	6:CF:190:GLU:OE2	2.19	0.43
13:CP:38:GLN:HG2	13:CP:45:LEU:N	2.29	0.43
1:CA:411:G:C5	13:CP:72:PRO:HB3	2.54	0.43
16:CS:24:LEU:HD23	16:CS:24:LEU:HA	1.73	0.43
23:CZ:59:LEU:HD11	23:CZ:88:PHE:CD1	2.54	0.43
34:DA:1008:C:H2'	34:DA:1009:G:O4'	2.19	0.43
34:DA:1064:G:N2	34:DA:1190:G:H2'	2.33	0.43
34:DA:1171:G:H2'	34:DA:1172:C:C6	2.54	0.43
34:DA:1226:C:H4'	34:DA:1227:A:OP1	2.17	0.43
34:DA:193:C:H2'	34:DA:194:C:C6	2.53	0.43
34:DA:410:G:C2	34:DA:429:U:C2	3.06	0.43
34:DA:448:A:P	34:DA:485:G:H22	2.40	0.43
34:DA:563:A:H2'	34:DA:567:G:C8	2.53	0.43
46:DM:5:ALA:HB1	46:DM:66:LEU:HD22	1.99	0.43
43:DJ:63:PHE:CE1	47:DN:58:LYS:HG2	2.53	0.43
49:DP:19:ILE:N	49:DP:37:GLY:O	2.50	0.43
55:DV:17:U:C2	56:DW:36:A:H2	2.36	0.43
57:DZ:550:MET:HB3	57:DZ:560:VAL:HB	2.00	0.43
57:DZ:655:TYR:CE2	57:DZ:659:LEU:HG	2.54	0.43
25:A1:63:ALA:O	25:A1:64:ALA:C	2.56	0.43
1:AA:1007:G:C5	1:AA:1008:U:C5	3.06	0.43
1:AA:1016:C:H2'	1:AA:1017:G:H5'	2.00	0.43
1:AA:1099:C:C4	1:AA:1100:A:C2	3.06	0.43
1:AA:1210:G:H2'	1:AA:1211:U:C6	2.54	0.43
1:AA:139:A:C8	1:AA:1454:C:O2'	2.69	0.43
1:AA:1480:A:H61	1:AA:1605:A:H61	1.61	0.43
1:AA:1657:C:OP1	63:AA:5027:HOH:O	2.21	0.43
1:AA:174:U:H2'	1:AA:175:G:H8	1.83	0.43
1:AA:2274:U:OP1	1:AA:2399:U:O2'	2.24	0.43
1:AA:2299:A:C4	1:AA:2301:G:C8	3.06	0.43
1:AA:668:A:O2'	1:AA:669:A:H5'	2.19	0.43
1:AA:702:A:H8	1:AA:703:G:C1'	2.32	0.43
2:AB:11:C:H3'	2:AB:12:C:H6	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:AN:73:THR:HG21	63:AN:3101:HOH:O	2.18	0.43
1:AA:1002:A:P	14:AQ:77:LYS:HG3	2.59	0.43
16:AS:36:TYR:N	16:AS:36:TYR:CD1	2.86	0.43
19:AV:4:ILE:HD12	19:AV:39:LEU:HB3	2.01	0.43
23:AZ:151:HIS:HD1	23:AZ:151:HIS:N	2.17	0.43
34:BA:1048:G:OP1	47:BN:4:LYS:HB2	2.18	0.43
34:BA:136:C:O3'	49:BP:65:GLN:NE2	2.52	0.43
34:BA:1508:G:H2'	34:BA:1509:C:O4'	2.19	0.43
34:BA:258:G:H2'	34:BA:259:G:H8	1.83	0.43
34:BA:259:G:O2'	34:BA:260:G:H5'	2.18	0.43
34:BA:949:A:C6	34:BA:950:U:C4	3.06	0.43
36:BC:108:ASN:HB3	36:BC:111:LEU:HG	1.99	0.43
37:BD:102:ASP:HA	37:BD:121:VAL:HG21	2.01	0.43
34:BA:1525:G:P	44:BK:120:ARG:NH2	2.91	0.43
49:BP:49:LEU:HD23	49:BP:76:GLN:HG2	2.00	0.43
51:BR:56:THR:HB	51:BR:58:LEU:HD22	2.00	0.43
24:C0:21:LEU:HD11	24:C0:41:ARG:HG2	1.99	0.43
1:CA:1091:G:H2'	1:CA:1092:C:O4'	2.19	0.43
1:CA:1167:U:H2'	1:CA:1168:G:C8	2.52	0.43
1:CA:1388:G:H2'	1:CA:1389:G:H8	1.84	0.43
1:CA:1448:G:N2	1:CA:1449:A:N6	2.67	0.43
1:CA:1932:A:N6	1:CA:1968:G:H1'	2.33	0.43
1:CA:2397:G:N2	1:CA:2420:C:H1'	2.34	0.43
1:CA:2718:G:C6	1:CA:2719:G:C5	3.06	0.43
1:CA:299:A:N1	1:CA:322:A:O2'	2.38	0.43
1:CA:874:G:O2'	23:CZ:120:ILE:HD11	2.19	0.43
2:CB:99:G:H8	2:CB:99:G:O5'	2.01	0.43
3:CC:11:LEU:HD11	3:CC:35:THR:HG23	2.01	0.43
4:CD:68:LYS:C	4:CD:70:TRP:H	2.22	0.43
6:CF:124:LEU:HB3	6:CF:193:VAL:HG22	1.99	0.43
15:CR:87:TYR:OH	15:CR:116:LEU:HB3	2.18	0.43
34:DA:1001:A:H2'	34:DA:1001(A):G:H8	1.83	0.43
34:DA:1040:U:H6	34:DA:1040:U:O5'	2.02	0.43
34:DA:1062:U:H2'	34:DA:1063:C:C6	2.53	0.43
34:DA:1183:A:H1'	34:DA:1184:G:OP1	2.19	0.43
34:DA:1401:G:C2	34:DA:1402:C:H1'	2.54	0.43
34:DA:409:G:OP1	37:DD:25:ARG:N	2.47	0.43
34:DA:872:A:C8	34:DA:874:G:C8	3.07	0.43
40:DG:75:VAL:HA	40:DG:87:VAL:O	2.18	0.43
42:DI:26:VAL:HG22	42:DI:61:ALA:HB3	2.00	0.43
46:DM:29:ARG:HG3	46:DM:64:TRP:CZ2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:750:G:O2'	48:DO:21:ASP:OD1	2.37	0.43
57:DZ:20:HIS:CG	57:DZ:21:ILE:H	2.37	0.43
57:DZ:417:THR:HA	57:DZ:418:LYS:HG2	2.00	0.43
57:DZ:423:LYS:NZ	57:DZ:471:LYS:HD3	2.33	0.43
57:DZ:487:ILE:HB	57:DZ:597:GLY:O	2.19	0.43
57:DZ:91:THR:O	57:DZ:92:ILE:HG22	2.18	0.43
27:A3:9:VAL:HG12	27:A3:32:GLN:HE22	1.84	0.43
1:AA:1885:A:C8	1:AA:1886:G:C8	3.06	0.43
1:AA:1885:A:C2	1:AA:2109:G:N3	2.87	0.43
1:AA:239:G:H2'	1:AA:240:A:C8	2.54	0.43
1:AA:2529:C:C6	1:AA:2554:A:N7	2.87	0.43
1:AA:649:C:O5'	1:AA:649:C:H6	2.02	0.43
4:AD:266:SER:O	4:AD:270:ILE:HD12	2.19	0.43
6:AF:34:TRP:CZ3	13:AP:8:PRO:HB3	2.53	0.43
20:AW:37:ARG:HD2	20:AW:38:TYR:CE2	2.53	0.43
23:AZ:146:ILE:HA	23:AZ:174:VAL:HG12	2.00	0.43
34:BA:1434:A:H2'	34:BA:1435:G:O4'	2.19	0.43
4:AD:202:LYS:NZ	34:BA:774:G:OP1	2.49	0.43
34:BA:872:A:C4	34:BA:874:G:N7	2.87	0.43
35:BB:97:TRP:CZ2	35:BB:102:LEU:HD13	2.54	0.43
35:BB:97:TRP:HZ3	35:BB:176:GLU:OE2	2.02	0.43
45:BL:122:THR:HG22	45:BL:123:LYS:O	2.19	0.43
52:BS:45:VAL:HG11	52:BS:64:GLU:HG2	1.99	0.43
57:BZ:153:MET:O	57:BZ:157:LEU:HD12	2.17	0.43
1:CA:2359:C:O2'	32:C8:54:GLU:HG3	2.19	0.43
1:CA:1266:G:C5	20:CW:15:ARG:NH1	2.86	0.43
1:CA:2774:C:H2'	1:CA:2775:A:O4'	2.18	0.43
1:CA:339:U:H6	1:CA:339:U:O5'	2.01	0.43
1:CA:391:G:C6	1:CA:411:G:C2	3.07	0.43
2:CB:28:C:OP1	16:CS:36:TYR:OH	2.33	0.43
3:CC:60:ARG:HG3	3:CC:165:ARG:HB2	2.01	0.43
5:CE:116:VAL:O	5:CE:122:PHE:HB2	2.18	0.43
6:CF:119:ARG:HG2	6:CF:119:ARG:O	2.18	0.43
6:CF:178:PRO:HB2	6:CF:201:VAL:HG21	2.01	0.43
13:CP:19:VAL:CG2	13:CP:31:ALA:HB1	2.49	0.43
23:CZ:61:LEU:HD12	23:CZ:62:PRO:HD2	2.00	0.43
34:DA:1001(A):G:C4	34:DA:1002:G:H1'	2.54	0.43
34:DA:1122:U:O4	34:DA:1123:A:N6	2.51	0.43
34:DA:1139:G:N2	34:DA:1142:G:O6	2.44	0.43
34:DA:396:G:H5''	57:DZ:349:LYS:NZ	2.34	0.43
34:DA:403:C:O2'	37:DD:122:ARG:NH2	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:435:C:H2'	34:DA:436:C:H6	1.80	0.43
34:DA:796:C:O5'	34:DA:796:C:H6	2.02	0.43
36:DC:113:ALA:O	36:DC:116:VAL:N	2.50	0.43
36:DC:12:LEU:HD23	36:DC:16:ARG:HB3	2.00	0.43
37:DD:104:VAL:HG21	37:DD:140:VAL:HG21	1.99	0.43
34:DA:1291:G:O2'	42:DI:38:GLN:HG3	2.18	0.43
46:DM:52:GLU:HG2	46:DM:55:ARG:NH2	2.33	0.43
56:DW:38:A:H2'	56:DW:39:PSU:O4'	2.18	0.43
30:A6:18:ARG:HD3	30:A6:42:TRP:CD1	2.53	0.43
1:AA:1101:G:H5'	1:AA:1102:G:OP2	2.18	0.43
1:AA:1405:A:H61	1:AA:1418:U:H3	1.67	0.43
1:AA:1462:G:HO2'	1:AA:1463:C:H5	1.66	0.43
1:AA:1541:A:O2'	1:AA:1542:A:H5'	2.19	0.43
1:AA:2219:U:C5	1:AA:2236:G:C6	3.07	0.43
1:AA:2359:C:H2'	1:AA:2360:U:C6	2.54	0.43
1:AA:402:C:H2'	1:AA:403:C:C6	2.54	0.43
6:AF:39:TRP:O	6:AF:40:GLN:C	2.57	0.43
7:AG:121:ASN:HA	7:AG:122:PRO:HD3	1.81	0.43
8:AH:6:ARG:HE	8:AH:6:ARG:HB3	1.39	0.43
1:AA:1110:C:H4'	10:AL:89:HIS:HA	1.99	0.43
1:AA:1233:U:C4'	19:AV:79:VAL:HG22	2.48	0.43
20:AW:75:TYR:CZ	20:AW:104:THR:HG21	2.54	0.43
34:BA:13:U:OP1	63:BA:5218:HOH:O	2.22	0.43
34:BA:284:G:H2'	34:BA:285:G:H8	1.82	0.43
34:BA:358:U:H2'	34:BA:359:U:H6	1.83	0.43
34:BA:939:G:H2'	34:BA:940:C:C6	2.54	0.43
34:BA:957:U:H4'	52:BS:79:THR:O	2.18	0.43
36:BC:56:ASP:O	36:BC:66:VAL:HA	2.19	0.43
37:BD:206:PHE:O	37:BD:208:SER:N	2.52	0.43
41:BH:83:ILE:HA	41:BH:136:GLU:O	2.19	0.43
53:BT:57:ARG:HH22	53:BT:100:ILE:HG13	1.84	0.43
57:BZ:182:ARG:O	57:BZ:184:LYS:HG3	2.19	0.43
57:BZ:-36:LEU:HD21	57:BZ:-29:LEU:HD22	2.00	0.43
57:BZ:420:ASP:CG	57:BZ:423:LYS:HE3	2.39	0.43
1:CA:1957:C:H2'	1:CA:1958:C:C6	2.54	0.43
1:CA:2243:U:H2'	1:CA:2244:U:C6	2.54	0.43
1:CA:2274:A:C2	1:CA:2276:G:H1'	2.54	0.43
1:CA:333:G:H5''	1:CA:334:C:OP2	2.18	0.43
1:CA:672:C:H2'	1:CA:673:C:C6	2.54	0.43
4:CD:73:VAL:HG13	4:CD:120:GLY:HA2	2.00	0.43
5:CE:47:VAL:HG23	5:CE:49:LEU:HD13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:CE:58:ARG:HA	5:CE:58:ARG:HD3	1.62	0.43
7:CG:171:ALA:O	7:CG:175:LEU:HB2	2.19	0.43
12:CO:10:VAL:HG21	12:CO:86:ILE:HD11	2.00	0.43
23:CZ:160:GLY:N	23:CZ:161:VAL:HB	2.34	0.43
34:DA:1251:A:H2'	34:DA:1252:A:O4'	2.19	0.43
34:DA:179:A:H2'	34:DA:180:U:H6	1.80	0.43
34:DA:355:C:C2	34:DA:356:A:C8	3.07	0.43
34:DA:393:A:C2	34:DA:394:G:C8	3.07	0.43
34:DA:44:G:C2	34:DA:45:U:H1'	2.53	0.43
34:DA:601:C:H2'	34:DA:602:A:C8	2.53	0.43
34:DA:680:C:O5'	34:DA:680:C:H6	2.02	0.43
35:DB:155:LEU:HA	35:DB:155:LEU:HD23	1.82	0.43
38:DE:102:ALA:HB1	38:DE:106:PRO:HG2	2.01	0.43
42:DI:63:ILE:HG21	42:DI:77:ILE:HG12	2.01	0.43
57:DZ:174:PHE:O	57:DZ:267:LYS:HE2	2.18	0.43
57:DZ:355:LEU:HB3	57:DZ:366:VAL:HG23	2.01	0.43
57:DZ:38:ARG:HH12	57:DZ:270:GLN:NE2	2.17	0.43
57:DZ:507:TYR:O	57:DZ:577:SER:OG	2.26	0.43
1:AA:1147:U:H2'	1:AA:1148:C:H6	1.84	0.43
1:AA:12:U:H2'	1:AA:12:U:O2	2.19	0.43
1:AA:1596:C:H2'	1:AA:1597:C:O4'	2.18	0.43
1:AA:1888:G:C6	1:AA:1889:G:C6	3.05	0.43
1:AA:2308:U:H4'	1:AA:2309:C:OP1	2.19	0.43
1:AA:602:G:H2'	1:AA:603:C:C6	2.54	0.43
1:AA:2154:U:C6	3:AC:6:LYS:HB2	2.52	0.43
4:AD:52:ARG:NH2	63:AD:412:HOH:O	2.50	0.43
5:AE:181:LEU:HA	5:AE:181:LEU:HD12	1.68	0.43
23:AZ:183:LEU:HD12	23:AZ:183:LEU:HA	1.68	0.43
34:BA:1203:C:H2'	34:BA:1204:A:C8	2.53	0.43
34:BA:1285:A:H4'	34:BA:1286:A:O5'	2.19	0.43
34:BA:1315:U:H2'	34:BA:1316:G:O4'	2.19	0.43
34:BA:152:A:N6	34:BA:170:U:C2	2.87	0.43
34:BA:319:G:H2'	34:BA:320:C:O4'	2.19	0.43
34:BA:502:G:N1	34:BA:503:C:N3	2.67	0.43
34:BA:567:G:H2'	34:BA:568:G:O4'	2.19	0.43
34:BA:636:U:H2'	34:BA:637:G:C8	2.54	0.43
34:BA:767:A:H2'	34:BA:768:A:O4'	2.19	0.43
35:BB:87:ARG:NH1	35:BB:220:ASP:OD1	2.49	0.43
37:BD:101:LEU:O	37:BD:102:ASP:C	2.57	0.43
37:BD:68:TYR:CD1	37:BD:97:LEU:HD22	2.53	0.43
41:BH:10:LEU:HD23	41:BH:10:LEU:N	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
42:BI:7:THR:O	42:BI:83:ARG:HD2	2.19	0.43
48:BO:26:GLU:HG2	48:BO:26:GLU:H	1.31	0.43
51:BR:61:LYS:HE3	51:BR:61:LYS:HB2	1.76	0.43
56:BW:44:G:H8	56:BW:44:G:OP2	2.02	0.43
57:BZ:210:ARG:O	57:BZ:213:HIS:HB3	2.18	0.43
57:BZ:494:GLU:CD	57:BZ:511:LYS:HE2	2.39	0.43
1:CA:210:C:OP2	31:C7:29:LYS:HE3	2.19	0.43
1:CA:493:G:H2'	1:CA:494:G:O4'	2.19	0.43
1:CA:602:G:O2'	1:CA:655:A:N6	2.51	0.43
3:CC:206:LYS:HB3	3:CC:206:LYS:HZ3	1.82	0.43
3:CC:6:LYS:N	3:CC:9:ARG:HH12	2.17	0.43
5:CE:116:VAL:HG13	5:CE:122:PHE:CD2	2.53	0.43
5:CE:18:ASP:HB3	17:CT:82:LEU:HD21	2.00	0.43
8:CH:12:PRO:O	8:CH:15:VAL:HG12	2.19	0.43
12:CO:31:LYS:HB3	12:CO:32:TYR:CD2	2.53	0.43
15:CR:13:HIS:O	15:CR:14:SER:C	2.57	0.43
34:DA:1133:G:N2	34:DA:1141:C:O2	2.51	0.43
34:DA:169:C:H2'	34:DA:170:U:C6	2.54	0.43
34:DA:473:G:C2	34:DA:474:G:N7	2.87	0.43
34:DA:938:A:C6	34:DA:939:G:C5	3.06	0.43
36:DC:123:GLN:O	36:DC:128:PHE:HB2	2.19	0.43
36:DC:155:GLY:O	36:DC:157:ILE:HG13	2.18	0.43
36:DC:6:HIS:NE2	36:DC:184:TYR:HE2	2.17	0.43
51:DR:59:SER:H	51:DR:62:GLU:HB2	1.83	0.43
52:DS:50:ALA:HB1	52:DS:57:HIS:HB3	2.01	0.43
55:DV:16:U:O5'	55:DV:16:U:H6	2.01	0.43
57:DZ:182:ARG:HG3	57:DZ:182:ARG:H	1.63	0.43
57:DZ:529:ILE:HA	57:DZ:529:ILE:HD13	1.75	0.43
1:AA:1221:G:H1'	1:AA:1222:A:C5'	2.49	0.43
1:AA:1216:G:N2	1:AA:1225:C:O2	2.52	0.43
1:AA:1314:A:H2'	1:AA:1315:A:O4'	2.18	0.43
1:AA:1374:G:O6	63:AA:4164:HOH:O	2.21	0.43
1:AA:1550:C:H2'	1:AA:1551:C:C6	2.52	0.43
1:AA:1688:A:H2'	1:AA:1689:G:O4'	2.18	0.43
1:AA:174:U:H2'	1:AA:175:G:C8	2.54	0.43
1:AA:1857:G:H2'	1:AA:1858:C:C6	2.54	0.43
1:AA:1539:C:H5	1:AA:2227:G:O2'	2.02	0.43
1:AA:2564:U:C2	1:AA:2566:U:H5'	2.54	0.43
1:AA:261:A:N7	1:AA:283:G:N2	2.62	0.43
1:AA:1790:A:C8	1:AA:2708:U:H1'	2.54	0.43
1:AA:609:A:N1	1:AA:856:G:O2'	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AD:26:LYS:HD3	4:AD:83:GLU:OE2	2.19	0.43
8:AH:105:LEU:HA	8:AH:105:LEU:HD12	1.79	0.43
8:AH:25:LYS:HE3	8:AH:27:LYS:HE2	2.00	0.43
15:AR:38:VAL:HG23	15:AR:110:PRO:O	2.19	0.43
18:AU:36:ARG:HD2	18:AU:40:PHE:CZ	2.54	0.43
23:AZ:111:VAL:HG12	23:AZ:112:ARG:N	2.34	0.43
34:BA:146:G:N2	34:BA:177:C:N3	2.66	0.43
34:BA:818:G:HO2'	34:BA:820:U:H6	1.62	0.43
34:BA:977:A:C8	34:BA:1223:C:C4	3.07	0.43
39:BF:11:ASN:O	39:BF:14:LEU:HG	2.19	0.43
40:BG:113:GLU:HG3	40:BG:119:ARG:HG2	2.01	0.43
41:BH:6:ILE:O	41:BH:7:ALA:C	2.57	0.43
51:BR:59:SER:N	51:BR:62:GLU:HG3	2.31	0.43
53:BT:14:LYS:O	53:BT:17:ARG:HB2	2.18	0.43
61:BZ:703:FUA:H322	61:BZ:703:FUA:H16	1.74	0.43
1:CA:1090:U:O2	1:CA:1102:C:H1'	2.19	0.43
1:CA:10:G:H2'	1:CA:11:G:C8	2.54	0.43
1:CA:1047:G:O2'	1:CA:1110:G:O6	2.31	0.43
1:CA:1338:G:C6	1:CA:1339:G:C5	3.06	0.43
1:CA:2052:G:N3	5:CE:149:ARG:HA	2.34	0.43
1:CA:2515:C:O2'	1:CA:2516:G:H5'	2.18	0.43
1:CA:2574:G:H2'	1:CA:2575:C:C6	2.53	0.43
1:CA:2776:A:C6	1:CA:2778:A:C6	3.07	0.43
1:CA:727:A:O2'	1:CA:728:G:H5'	2.19	0.43
8:CH:117:PRO:HA	8:CH:118:PRO:HD2	1.90	0.43
13:CP:136:GLU:O	13:CP:140:ALA:HB3	2.19	0.43
34:DA:1004:A:H5'	34:DA:1024:G:H22	1.83	0.43
34:DA:1357:A:N6	34:DA:1358:U:O4	2.52	0.43
34:DA:283:C:H2'	34:DA:284:G:O4'	2.18	0.43
34:DA:580:U:H2'	34:DA:581:G:O4'	2.19	0.43
36:DC:148:GLY:HA3	36:DC:203:PHE:HB3	2.01	0.43
43:DJ:42:THR:HG21	43:DJ:66:ARG:HB3	2.00	0.43
43:DJ:16:LEU:HD23	43:DJ:94:VAL:HG22	2.01	0.43
54:DU:15:ARG:HH11	54:DU:15:ARG:HB2	1.84	0.43
56:DW:40:C:O2'	56:DY:36:A:P	2.77	0.43
57:DZ:6:GLU:HA	57:DZ:9:LEU:CD2	2.49	0.43
30:A6:50:ARG:HB2	30:A6:50:ARG:HE	1.61	0.42
31:A7:34:ARG:NH1	31:A7:41:ARG:O	2.52	0.42
1:AA:1088:G:C6	1:AA:1089:C:C4	3.07	0.42
1:AA:1417:G:H2'	1:AA:1418:U:C5	2.53	0.42
1:AA:1568:G:H2'	1:AA:1569:U:O4'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:181:C:C2'	1:AA:182:U:H5'	2.46	0.42
1:AA:223:C:H2'	1:AA:224:U:C6	2.54	0.42
1:AA:526:A:C6	1:AA:527:A:C6	3.07	0.42
12:AO:64:ARG:HG2	12:AO:79:PHE:CB	2.49	0.42
17:AT:117:ASP:O	17:AT:121:ILE:HG13	2.19	0.42
23:AZ:136:PHE:O	23:AZ:137:ILE:HG13	2.19	0.42
2:AB:73:A:N1	23:AZ:34:ASN:ND2	2.67	0.42
34:BA:1072:G:H2'	34:BA:1073:U:C6	2.54	0.42
34:BA:1104:G:C2	34:BA:1105:A:C4	3.07	0.42
34:BA:16:A:N1	34:BA:919:A:H2	2.17	0.42
34:BA:176:C:H2'	34:BA:177:C:C6	2.54	0.42
34:BA:66:G:O4'	34:BA:173:U:C4	2.72	0.42
36:BC:50:ALA:HB1	36:BC:70:VAL:HG21	2.00	0.42
37:BD:30:LYS:HA	37:BD:35:ARG:HH11	1.83	0.42
34:BA:1291:G:H5'	40:BG:37:ASN:HD21	1.84	0.42
40:BG:79:ARG:HB3	40:BG:80:VAL:H	1.56	0.42
42:BI:48:GLU:OE1	42:BI:51:ARG:NH1	2.52	0.42
44:BK:48:ILE:HD12	44:BK:63:LEU:CB	2.49	0.42
47:BN:3:ARG:HB3	47:BN:4:LYS:H	1.69	0.42
57:BZ:89:ASP:CG	57:BZ:457:LEU:HB2	2.39	0.42
28:C4:40:HIS:O	28:C4:44:THR:N	2.30	0.42
1:CA:1195:G:O2'	1:CA:1196:C:H5'	2.19	0.42
1:CA:1340:U:H4'	1:CA:1341:U:OP2	2.18	0.42
1:CA:1364:G:OP2	25:C1:3:LYS:HG3	2.19	0.42
1:CA:1666:G:C2'	1:CA:1667:G:H5'	2.48	0.42
1:CA:1750:G:N3	1:CA:2860:A:H2	2.17	0.42
1:CA:528:A:C2	1:CA:2042:A:H2'	2.53	0.42
1:CA:2046:G:H2'	1:CA:2047:U:C6	2.53	0.42
1:CA:2318:G:H21	16:CS:3:ARG:HH12	1.67	0.42
1:CA:2469:A:C2	1:CA:2470:G:H1'	2.54	0.42
1:CA:489:G:H2'	1:CA:491:G:O4'	2.19	0.42
1:CA:2122:U:O2	3:CC:167:ASP:HB3	2.19	0.42
4:CD:238:GLY:O	4:CD:239:ARG:CB	2.63	0.42
8:CH:140:LYS:HE3	8:CH:140:LYS:HB2	1.74	0.42
13:CP:39:LYS:HA	13:CP:45:LEU:HG	2.00	0.42
13:CP:68:GLN:OE1	13:CP:68:GLN:HA	2.19	0.42
11:CN:42:TRP:CE3	18:CU:63:VAL:HG11	2.54	0.42
1:CA:1225:G:H4'	19:CV:84:LYS:HG2	2.01	0.42
34:DA:1085:U:H3'	34:DA:1086:U:H5	1.84	0.42
34:DA:157:G:H2'	34:DA:158:G:C8	2.50	0.42
34:DA:355:C:C4	34:DA:356:A:N7	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:987:G:O5'	34:DA:987:G:H8	2.01	0.42
38:DE:47:LYS:HB2	38:DE:47:LYS:HE2	1.64	0.42
40:DG:26:PHE:HB2	40:DG:62:PHE:HZ	1.84	0.42
42:DI:106:ALA:O	42:DI:108:VAL:N	2.52	0.42
42:DI:110:GLU:OE2	42:DI:113:LYS:NZ	2.52	0.42
42:DI:45:ALA:O	42:DI:48:GLU:HB2	2.19	0.42
34:DA:1254:C:OP1	43:DJ:45:ARG:HA	2.19	0.42
45:DL:31:PRO:HB2	45:DL:32:PHE:CD2	2.53	0.42
51:DR:31:LEU:HD13	51:DR:31:LEU:HA	1.80	0.42
34:DA:790:A:H5'	56:DW:39:PSU:OP1	2.18	0.42
57:DZ:431:LEU:HA	57:DZ:434:GLU:HB3	2.00	0.42
28:A4:63:TYR:N	28:A4:64:GLY:HA2	2.33	0.42
1:AA:1098:C:O2'	1:AA:1099:C:H5'	2.19	0.42
1:AA:1815:A:H4'	1:AA:1816:A:C5'	2.50	0.42
1:AA:2050:U:O4	63:AA:4233:HOH:O	2.19	0.42
1:AA:2216:G:H2'	1:AA:2217:C:O4'	2.19	0.42
1:AA:510:C:H2'	1:AA:511:C:C6	2.54	0.42
1:AA:556:C:H4'	1:AA:557:A:H5''	2.01	0.42
4:AD:159:ALA:HB1	4:AD:198:ASN:O	2.18	0.42
4:AD:261:LYS:NZ	4:AD:263:ARG:NH2	2.67	0.42
5:AE:150:VAL:CG1	5:AE:154:LYS:HG3	2.49	0.42
6:AF:103:LYS:O	6:AF:106:ARG:HG2	2.19	0.42
6:AF:13:SER:N	6:AF:16:GLY:O	2.45	0.42
10:AL:72:PRO:HA	10:AL:73:PRO:HD3	1.86	0.42
1:AA:1185:C:O3'	11:AN:25:ARG:NH1	2.52	0.42
17:AT:39:ARG:NH2	34:BA:345:C:C5	2.87	0.42
21:AX:29:TRP:CZ3	21:AX:78:LYS:HG2	2.53	0.42
23:AZ:110:GLY:C	23:AZ:111:VAL:O	2.56	0.42
23:AZ:99:TYR:CZ	23:AZ:125:LEU:HD13	2.54	0.42
34:BA:1041:A:H2'	34:BA:1042:G:O4'	2.19	0.42
34:BA:1112:C:O2	36:BC:179:ARG:HG2	2.19	0.42
34:BA:1274:G:H22	34:BA:1275:A:H62	1.66	0.42
34:BA:445:G:C4	34:BA:446:G:C8	3.07	0.42
34:BA:454:C:N4	34:BA:479:C:N3	2.67	0.42
34:BA:57:G:H2'	34:BA:58:C:O4'	2.19	0.42
36:BC:45:LYS:HD3	36:BC:46:GLU:HG2	2.01	0.42
36:BC:6:HIS:CD2	36:BC:7:PRO:HD2	2.54	0.42
57:BZ:544:LYS:O	57:BZ:548:GLU:HB2	2.19	0.42
1:CA:1668:A:H5''	12:CO:5:GLN:HG2	2.01	0.42
1:CA:18:C:H2'	1:CA:19:C:H6	1.84	0.42
1:CA:2393:A:H2'	1:CA:2394:C:H6	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:2853:C:H2'	1:CA:2854:G:C8	2.53	0.42
1:CA:311:A:C6	1:CA:328:U:C4	3.07	0.42
1:CA:993:G:OP2	18:CU:51:LYS:NZ	2.52	0.42
4:CD:96:HIS:CD2	4:CD:102:LYS:HG2	2.55	0.42
1:CA:1287:A:OP1	15:CR:105:ARG:HB3	2.18	0.42
15:CR:18:LEU:HA	15:CR:18:LEU:HD23	1.86	0.42
17:CT:15:VAL:HG12	17:CT:16:ARG:N	2.34	0.42
20:CW:35:ILE:HG22	20:CW:36:LEU:N	2.34	0.42
22:CY:52:SER:HB2	22:CY:53:PRO:HD2	2.01	0.42
34:DA:1074:G:C6	34:DA:1075:C:C4	3.07	0.42
34:DA:56:U:H2'	34:DA:57:G:C8	2.54	0.42
35:DB:133:LYS:O	35:DB:137:ARG:HG3	2.19	0.42
37:DD:59:ARG:O	37:DD:60:GLU:C	2.57	0.42
38:DE:139:LEU:C	38:DE:141:GLN:H	2.21	0.42
38:DE:81:GLU:HB3	38:DE:90:VAL:HG13	2.00	0.42
41:DH:94:TYR:CE1	41:DH:132:GLU:HB2	2.54	0.42
51:DR:65:ILE:O	51:DR:69:THR:HG23	2.19	0.42
57:DZ:605:ILE:HD13	57:DZ:675:HIS:CE1	2.55	0.42
1:AA:1968:U:H2'	1:AA:1969:C:C6	2.54	0.42
1:AA:2007:G:OP2	63:AA:4910:HOH:O	2.22	0.42
1:AA:553:A:C2	1:AA:2065:C:H4'	2.54	0.42
1:AA:2331:G:H1	16:AS:3:ARG:HD3	1.83	0.42
1:AA:424:G:O2'	1:AA:2243:C:H1'	2.19	0.42
1:AA:702:A:H2'	1:AA:703:G:O4'	2.19	0.42
8:AH:154:PRO:HD3	8:AH:162:ILE:O	2.19	0.42
19:AV:89:GLN:HA	19:AV:89:GLN:OE1	2.19	0.42
34:BA:1000:U:H2'	34:BA:1001:A:C8	2.53	0.42
34:BA:1026:G:H2'	34:BA:1026:G:N3	2.33	0.42
34:BA:1095:U:P	34:BA:1108:G:H1	2.41	0.42
34:BA:342:C:C2	34:BA:348:G:N2	2.87	0.42
34:BA:358:U:H2'	34:BA:359:U:C6	2.54	0.42
34:BA:444:C:C2	34:BA:445:G:C8	3.08	0.42
34:BA:522:C:C2'	34:BA:523:A:H5'	2.49	0.42
34:BA:558:G:C5	34:BA:559:A:C2	3.07	0.42
41:BH:56:LYS:O	41:BH:58:TYR:HD1	2.01	0.42
49:BP:57:ARG:HG2	49:BP:79:VAL:HG12	2.00	0.42
52:BS:41:VAL:N	52:BS:44:MET:SD	2.81	0.42
56:BY:9:A:H1'	56:BY:45:U:H2'	2.00	0.42
57:BZ:151:ARG:O	57:BZ:155:GLU:HB2	2.19	0.42
1:CA:1076:C:H2'	1:CA:1077:A:O4'	2.20	0.42
1:CA:530:G:O4'	1:CA:530:G:N3	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:553:G:C5	1:CA:554:U:C5	3.07	0.42
1:CA:565:C:H2'	1:CA:566:U:O4'	2.18	0.42
2:CB:117:G:H5'	16:CS:55:ALA:HB2	2.01	0.42
3:CC:195:ARG:HH11	3:CC:195:ARG:HG3	1.83	0.42
6:CF:169:ASN:HD22	6:CF:169:ASN:HA	1.51	0.42
1:CA:1257:C:H5'	6:CF:75:HIS:CE1	2.55	0.42
7:CG:117:PHE:HE1	7:CG:119:GLY:HA2	1.80	0.42
9:CK:23:SER:HA	9:CK:117:LEU:O	2.18	0.42
11:CN:33:LEU:HD13	11:CN:33:LEU:HA	1.80	0.42
1:CA:1141:U:H2'	11:CN:63:THR:HG21	2.00	0.42
16:CS:15:ARG:HB3	16:CS:19:LYS:NZ	2.34	0.42
18:CU:36:ARG:HD2	18:CU:40:PHE:CZ	2.55	0.42
20:CW:78:GLU:OE1	20:CW:99:ARG:NH1	2.49	0.42
23:CZ:130:PRO:O	23:CZ:133:ILE:HG12	2.20	0.42
34:DA:1072:G:H2'	34:DA:1073:U:O4'	2.19	0.42
34:DA:1222:G:C2	34:DA:1223:C:C2	3.07	0.42
34:DA:298:A:H5''	34:DA:299:G:OP2	2.19	0.42
34:DA:481:G:H21	34:DA:482:A:H62	1.67	0.42
34:DA:797:C:O2'	34:DA:798:G:H5'	2.19	0.42
34:DA:955:U:H2'	34:DA:956:U:O4'	2.19	0.42
37:DD:209:ARG:OXT	37:DD:209:ARG:HG2	2.18	0.42
38:DE:91:LEU:HA	38:DE:91:LEU:HD12	1.79	0.42
38:DE:99:GLY:O	38:DE:117:ASP:HA	2.19	0.42
56:DW:61:C:H2'	56:DW:62:C:C6	2.54	0.42
56:DY:35:A:N6	56:DY:36:A:N1	2.68	0.42
56:DY:51:U:O2	56:DY:63:G:N2	2.33	0.42
25:A1:91:LYS:HG2	25:A1:95:LEU:HD22	2.01	0.42
1:AA:1064:C:C2'	1:AA:1065:U:H5'	2.49	0.42
1:AA:1476:C:H2'	1:AA:1477:U:H6	1.83	0.42
1:AA:1549:U:H2'	1:AA:1550:C:C6	2.55	0.42
1:AA:1913:G:C6	1:AA:1914:C:C4	3.08	0.42
1:AA:2019:G:O2'	1:AA:2020:G:H5'	2.19	0.42
1:AA:2305:C:H2'	1:AA:2306:C:C6	2.54	0.42
1:AA:308:U:H2'	1:AA:309:C:C6	2.53	0.42
1:AA:974:G:O5'	1:AA:974:G:H8	2.02	0.42
3:AC:195:ARG:NH1	3:AC:195:ARG:HG3	2.34	0.42
3:AC:6:LYS:N	3:AC:9:ARG:HH12	2.17	0.42
5:AE:94:GLU:HG2	5:AE:94:GLU:H	1.59	0.42
6:AF:192:LEU:HD22	6:AF:194:MET:HG3	2.02	0.42
7:AG:112:PRO:C	7:AG:114:ILE:H	2.23	0.42
10:AL:30:HIS:HB2	10:AL:32:ALA:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:AL:41:PHE:HE1	10:AL:53:VAL:HB	1.83	0.42
11:AN:102:ALA:O	11:AN:106:MET:HG3	2.20	0.42
18:AU:113:ALA:O	18:AU:117:GLN:HG2	2.19	0.42
20:AW:29:LEU:O	20:AW:33:ARG:HG3	2.20	0.42
34:BA:1118:C:C1'	34:BA:1179:A:C4	2.99	0.42
34:BA:308:C:H2'	34:BA:309:G:C8	2.51	0.42
34:BA:542:G:H2'	34:BA:543:C:H6	1.85	0.42
34:BA:654:G:C2	34:BA:753:A:C4	3.07	0.42
34:BA:987:G:H2'	34:BA:988:G:C8	2.55	0.42
36:BC:152:ILE:HB	36:BC:199:LYS:HB2	2.02	0.42
42:BI:53:VAL:HG21	42:BI:92:TYR:CE1	2.54	0.42
52:BS:32:LYS:HE2	52:BS:57:HIS:CD2	2.54	0.42
56:BY:72:C:H2'	56:BY:73:A:O4'	2.18	0.42
57:BZ:-34:ARG:HB3	57:BZ:-34:ARG:HE	1.57	0.42
57:BZ:517:LEU:HA	57:BZ:517:LEU:HD12	1.72	0.42
57:BZ:621:ILE:HG13	57:BZ:643:ILE:HD13	2.01	0.42
1:CA:1000:A:C6	1:CA:1001:A:C6	3.06	0.42
1:CA:1115:G:H2'	1:CA:1116:C:H6	1.82	0.42
1:CA:1162:G:O3'	19:CV:24:LYS:NZ	2.51	0.42
1:CA:1287:A:H5"	1:CA:1288:U:OP2	2.19	0.42
1:CA:1557:C:P	1:CA:1558:A:HO2'	2.39	0.42
1:CA:2108:C:H2'	1:CA:2109:U:C6	2.53	0.42
1:CA:2861:G:C4	1:CA:2862:G:C8	3.07	0.42
3:CC:195:ARG:NH1	3:CC:195:ARG:HG3	2.35	0.42
3:CC:225:ILE:O	3:CC:227:PRO:HD3	2.19	0.42
5:CE:5:LEU:N	5:CE:5:LEU:HD23	2.34	0.42
5:CE:75:VAL:HG13	5:CE:77:ILE:H	1.84	0.42
6:CF:170:LEU:HG	6:CF:172:TRP:NE1	2.34	0.42
9:CK:69:PRO:O	9:CK:71:LEU:N	2.52	0.42
10:CL:41:PHE:CZ	10:CL:45:THR:HG21	2.54	0.42
12:CO:66:LYS:HE2	12:CO:80:ASP:O	2.20	0.42
13:CP:95:VAL:HA	13:CP:99:LEU:HD21	2.02	0.42
1:CA:958:U:H5"	14:CQ:14:ARG:HD3	2.02	0.42
20:CW:80:PRO:O	20:CW:100:THR:HB	2.20	0.42
34:DA:117:G:OP2	63:DA:3239:HOH:O	2.21	0.42
34:DA:1291:G:C6	34:DA:1292:U:C4	3.07	0.42
34:DA:203:U:H2'	34:DA:203:U:OP2	2.18	0.42
34:DA:705:U:C5	34:DA:706:A:C5	3.07	0.42
35:DB:178:ARG:HH22	41:DH:68:ARG:HH12	1.66	0.42
34:DA:1057:G:H1'	36:DC:195:VAL:HG12	2.02	0.42
43:DJ:7:LYS:HG2	43:DJ:71:LEU:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:888:C:H5'	46:DM:93:ARG:HD3	2.02	0.42
34:DA:264:U:H4'	50:DQ:63:ARG:HD3	2.01	0.42
56:DW:14:A:C4	56:DW:22:G:C2	3.07	0.42
57:DZ:164:MET:O	57:DZ:180:VAL:HG22	2.19	0.42
57:DZ:303:PRO:O	57:DZ:305:PRO:HD3	2.20	0.42
57:DZ:553:GLY:H	57:DZ:557:GLY:CA	2.24	0.42
1:AA:1888:G:H8	1:AA:1888:G:O5'	2.03	0.42
1:AA:2235:G:OP1	4:AD:172:TYR:OH	2.28	0.42
1:AA:2242:G:O3'	25:A1:43:TYR:HB2	2.19	0.42
1:AA:252:C:H2'	1:AA:253:C:O4'	2.19	0.42
1:AA:2638:C:H2'	1:AA:2639:G:C8	2.54	0.42
3:AC:195:ARG:HH11	3:AC:195:ARG:HG3	1.83	0.42
3:AC:225:ILE:O	3:AC:227:PRO:HD3	2.19	0.42
3:AC:54:ARG:HE	3:AC:57:GLN:HG2	1.83	0.42
4:AD:53:PHE:HB3	4:AD:218:ARG:O	2.20	0.42
7:AG:50:ALA:O	7:AG:52:ILE:N	2.53	0.42
10:AL:37:PHE:CE1	10:AL:41:PHE:HB2	2.55	0.42
11:AN:36:GLY:HA2	11:AN:38:HIS:CE1	2.54	0.42
14:AQ:2:LEU:HG	14:AQ:69:PHE:CE2	2.54	0.42
34:BA:1033:G:H2'	34:BA:1034:G:H8	1.84	0.42
34:BA:1289:A:N1	34:BA:1371:G:O2'	2.30	0.42
34:BA:325:A:H2'	34:BA:326:G:O4'	2.20	0.42
34:BA:430:A:OP1	37:BD:9:CYS:HB2	2.20	0.42
34:BA:895:G:H2'	34:BA:896:C:C6	2.55	0.42
35:BB:97:TRP:HZ2	35:BB:102:LEU:HD13	1.83	0.42
48:BO:18:PHE:HB2	48:BO:19:PRO:HD2	2.01	0.42
53:BT:49:ALA:O	53:BT:53:LEU:HG	2.20	0.42
57:BZ:100:VAL:H	57:BZ:100:VAL:HG23	1.63	0.42
57:BZ:168:ILE:N	57:BZ:176:GLY:O	2.49	0.42
57:BZ:277:VAL:HG12	57:BZ:278:ASP:N	2.34	0.42
27:C3:5:LYS:O	27:C3:56:VAL:HA	2.19	0.42
32:C8:34:TRP:CE3	32:C8:35:GLN:HG2	2.55	0.42
1:CA:1539:G:H2'	1:CA:1540:U:O4'	2.19	0.42
1:CA:1834:U:H4'	1:CA:1969:A:C6	2.54	0.42
1:CA:1893:C:C5	1:CA:1894:C:C5	3.07	0.42
1:CA:1916:A:O5'	1:CA:1916:A:H8	2.02	0.42
1:CA:2845:G:H2'	1:CA:2846:G:H8	1.83	0.42
1:CA:370:G:OP2	63:CA:3748:HOH:O	2.21	0.42
4:CD:123:ALA:HA	4:CD:124:PRO:HD3	1.85	0.42
5:CE:119:ARG:HD2	5:CE:120:TRP:NE1	2.35	0.42
7:CG:103:LEU:HD22	7:CG:178:PHE:HZ	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CH:71:LEU:HA	8:CH:71:LEU:HD12	1.82	0.42
12:CO:88:ASN:OD1	12:CO:92:GLU:HB2	2.19	0.42
18:CU:61:TRP:CZ3	18:CU:93:LYS:HB2	2.55	0.42
34:DA:1090:U:H2'	34:DA:1091:U:C6	2.55	0.42
34:DA:144:G:H2'	34:DA:145:G:C8	2.53	0.42
34:DA:1530:G:H2'	34:DA:1531:A:O4'	2.19	0.42
34:DA:779:C:H2'	34:DA:780:A:O4'	2.19	0.42
36:DC:175:LEU:H	36:DC:175:LEU:HD12	1.84	0.42
40:DG:147:ALA:C	40:DG:149:ARG:H	2.23	0.42
49:DP:49:LEU:HD12	49:DP:50:LYS:N	2.34	0.42
57:DZ:13:ARG:NH2	57:DZ:277:VAL:O	2.52	0.42
1:AA:493:G:OP1	31:A7:33:ARG:HD2	2.19	0.42
32:A8:6:THR:HB	32:A8:8:LYS:HE2	2.00	0.42
1:AA:1202:A:C8	18:AU:51:LYS:HD2	2.54	0.42
1:AA:170:A:H2'	1:AA:171:A:C8	2.54	0.42
1:AA:1849:U:H2'	4:AD:157:ARG:HG3	2.01	0.42
1:AA:2228:G:O2'	1:AA:2229:A:OP1	2.30	0.42
1:AA:2321:A:H8	1:AA:2321:A:O5'	2.02	0.42
1:AA:583:C:H2'	1:AA:584:G:O4'	2.20	0.42
1:AA:64:C:H2'	1:AA:65:C:C6	2.54	0.42
2:AB:54:G:H2'	2:AB:55:U:H6	1.84	0.42
4:AD:215:LEU:HA	4:AD:215:LEU:HD23	1.81	0.42
16:AS:61:ASN:O	16:AS:65:VAL:HG23	2.20	0.42
1:AA:572:A:H62	19:AV:19:LYS:NZ	2.16	0.42
19:AV:20:LEU:HD12	19:AV:21:ARG:H	1.84	0.42
20:AW:19:LEU:HB3	29:A5:25:LEU:CD1	2.50	0.42
34:BA:1270:C:H2'	34:BA:1271:G:H8	1.83	0.42
34:BA:1405:G:H1'	34:BA:1519:A:H4'	2.01	0.42
34:BA:189(K):U:H2'	34:BA:189(L):G:C8	2.54	0.42
34:BA:266:G:C8	34:BA:266:G:H5'	2.54	0.42
34:BA:410:G:C2	34:BA:429:U:C2	3.08	0.42
34:BA:578:C:O4'	34:BA:729:A:H1'	2.19	0.42
34:BA:750:G:H1'	48:BO:22:THR:O	2.20	0.42
35:BB:83:MET:H	35:BB:83:MET:HG2	1.63	0.42
41:BH:119:LEU:HB3	41:BH:123:GLU:HB3	2.01	0.42
44:BK:45:GLY:O	44:BK:50:TYR:HB2	2.19	0.42
34:BA:279:A:H5'	50:BQ:95:TYR:OH	2.19	0.42
52:BS:3:ARG:HG2	52:BS:4:SER:H	1.85	0.42
57:BZ:538:TYR:HB3	57:BZ:582:PHE:CD1	2.54	0.42
26:C2:49:LYS:HB3	26:C2:49:LYS:NZ	2.34	0.42
28:C4:59:PHE:HA	28:C4:60:GLN:C	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1005:C:OP2	1:CA:1011:G:O2'	2.21	0.42
1:CA:1189:A:OP2	63:CA:4135:HOH:O	2.21	0.42
1:CA:1545:A:H2'	1:CA:1546:C:O4'	2.19	0.42
1:CA:2258:C:H4'	1:CA:2259:G:OP2	2.20	0.42
1:CA:2488:A:H2'	1:CA:2489:G:C8	2.55	0.42
1:CA:2532:G:H2'	1:CA:2533:A:O4'	2.20	0.42
1:CA:2542:A:H4'	1:CA:2543:G:H8	1.84	0.42
1:CA:2572:A:C6	5:CE:144:ARG:NH1	2.87	0.42
1:CA:310:A:C4	1:CA:312:G:C8	3.07	0.42
1:CA:516:C:H6	1:CA:516:C:O5'	2.03	0.42
1:CA:569:U:C4	1:CA:570:G:C6	3.07	0.42
1:CA:833:U:O2	13:CP:55:ARG:NH2	2.53	0.42
4:CD:68:LYS:HB3	4:CD:70:TRP:CE2	2.54	0.42
8:CH:72:ILE:O	8:CH:75:ALA:HB3	2.19	0.42
18:CU:69:CYS:SG	18:CU:79:PHE:HB2	2.59	0.42
34:DA:1000:U:O2	34:DA:1042:G:N2	2.52	0.42
34:DA:1404:C:H2'	34:DA:1405:G:C8	2.55	0.42
34:DA:374:A:C6	34:DA:375:U:C4	3.06	0.42
34:DA:533:A:C6	34:DA:536:C:C2	3.07	0.42
45:DL:69:TYR:HD2	45:DL:99:HIS:CE1	2.37	0.42
46:DM:5:ALA:HB1	46:DM:66:LEU:HD13	2.01	0.42
49:DP:28:ARG:HG3	49:DP:29:ASP:OD1	2.19	0.42
39:DF:98:LEU:HD23	51:DR:30:ASP:HA	2.02	0.42
51:DR:76:LEU:HA	51:DR:76:LEU:HD12	1.80	0.42
53:DT:39:LYS:O	53:DT:43:LEU:HG	2.20	0.42
56:DY:46:7MG:H81	56:DY:46:7MG:H2'	1.69	0.42
30:A6:11:LEU:HB2	30:A6:21:TYR:HB2	2.02	0.42
30:A6:40:CYS:HA	30:A6:41:PRO:HD3	1.85	0.42
1:AA:118:U:OP2	63:AA:3919:HOH:O	2.21	0.42
1:AA:1619:A:H5'	63:AA:4021:HOH:O	2.18	0.42
1:AA:1834:A:H4'	4:AD:259:THR:CG2	2.50	0.42
1:AA:1954:A:H2'	1:AA:1955:G:O4'	2.19	0.42
1:AA:1952:G:O2'	1:AA:1990:G:O6	2.31	0.42
1:AA:2228:G:H2'	1:AA:2229:A:C2	2.54	0.42
1:AA:701:A:O2'	1:AA:702:A:H5'	2.20	0.42
2:AB:111:G:C6	2:AB:112:U:C4	3.08	0.42
2:AB:29:A:C2	2:AB:56:G:C2	3.08	0.42
4:AD:69:ARG:C	4:AD:71:ASP:N	2.72	0.42
5:AE:176:ILE:HA	5:AE:177:PRO:HD2	1.85	0.42
5:AE:56:PRO:C	5:AE:58:ARG:H	2.23	0.42
11:AN:62:VAL:CG1	11:AN:66:LYS:HB2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:AR:96:ARG:O	15:AR:96:ARG:HG3	2.18	0.42
22:AY:56:PRO:C	22:AY:58:GLY:H	2.23	0.42
34:BA:1103:C:H5'	35:BB:98:LEU:HD13	2.01	0.42
34:BA:189(B):C:N3	34:BA:189(J):G:C2	2.88	0.42
34:BA:923:A:H2'	34:BA:924:C:C6	2.55	0.42
37:BD:204:ILE:HG21	38:BE:98:THR:O	2.19	0.42
40:BG:78:ARG:HG3	40:BG:156:TRP:CE3	2.54	0.42
41:BH:12:ARG:HH21	41:BH:27:PRO:HD3	1.84	0.42
43:BJ:31:GLY:HA2	43:BJ:32:ALA:HA	1.58	0.42
34:BA:35:G:N2	45:BL:118:SER:OG	2.38	0.42
47:BN:6:LEU:HG	47:BN:23:ARG:NH2	2.34	0.42
57:BZ:182:ARG:HG3	57:BZ:182:ARG:H	1.62	0.42
57:BZ:300:GLU:C	57:BZ:301:ILE:HG12	2.39	0.42
26:C2:14:ARG:HA	26:C2:63:VAL:HG11	2.01	0.42
28:C4:16:CYS:SG	28:C4:17:GLY:N	2.93	0.42
30:C6:39:TYR:HA	30:C6:46:HIS:HA	2.01	0.42
30:C6:40:CYS:HA	30:C6:41:PRO:HD3	1.80	0.42
1:CA:1068:G:N2	1:CA:1096:A:H5'	2.33	0.42
1:CA:1314:C:C2	1:CA:1339:G:N2	2.87	0.42
1:CA:1408:C:C2	1:CA:1595:G:N2	2.88	0.42
1:CA:1805:U:H2'	1:CA:1806:C:C6	2.55	0.42
1:CA:2517:C:C2	1:CA:2542:A:N6	2.88	0.42
1:CA:262:A:H2'	1:CA:263:C:O4'	2.20	0.42
1:CA:627:A:C6	1:CA:637:A:C8	3.08	0.42
1:CA:985:C:H2'	1:CA:986:C:H6	1.84	0.42
6:CF:183:VAL:HG23	6:CF:183:VAL:O	2.19	0.42
8:CH:75:ALA:O	8:CH:79:VAL:HG22	2.20	0.42
10:CL:99:ILE:HD12	10:CL:103:GLN:HB3	2.01	0.42
14:CQ:135:ASP:CB	14:CQ:137:TYR:HB2	2.49	0.42
14:CQ:85:LYS:HG3	24:C0:8:GLY:O	2.20	0.42
17:CT:102:ILE:HB	17:CT:110:ILE:HG12	2.01	0.42
34:DA:1256:A:N6	34:DA:1278:U:O2	2.52	0.42
34:DA:7:G:H5'	34:DA:298:A:O4'	2.19	0.42
34:DA:339:C:H2'	34:DA:340:U:H6	1.84	0.42
34:DA:731:G:OP1	34:DA:766:A:H1'	2.20	0.42
34:DA:799:G:O6	34:DA:800:G:C2	2.73	0.42
36:DC:150:LYS:HD2	36:DC:201:TYR:HD1	1.84	0.42
36:DC:37:GLN:HA	36:DC:40:ARG:HG3	2.01	0.42
37:DD:150:GLU:HA	37:DD:153:ARG:NE	2.34	0.42
37:DD:169:LYS:HB3	37:DD:169:LYS:NZ	2.35	0.42
34:DA:8:A:N6	37:DD:209:ARG:HB2	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
39:DF:15:ASP:OD1	39:DF:18:GLN:N	2.38	0.42
45:DL:105:TYR:C	45:DL:107:ALA:H	2.23	0.42
45:DL:6:THR:HG23	45:DL:9:GLN:OE1	2.20	0.42
57:DZ:20:HIS:CE1	57:DZ:21:ILE:HG12	2.55	0.42
57:DZ:265:LYS:O	57:DZ:267:LYS:HG3	2.19	0.42
57:DZ:290:LYS:HG2	57:DZ:400:GLU:OE1	2.18	0.42
34:DA:396:G:P	57:DZ:349:LYS:HZ2	2.43	0.42
57:DZ:309:LEU:HB3	57:DZ:391:GLY:N	2.35	0.42
29:A5:15:ARG:HD3	29:A5:15:ARG:HH11	1.64	0.42
1:AA:1236:G:O5'	1:AA:1236:G:H8	2.03	0.42
1:AA:1743:G:C6	1:AA:1744:G:C4	3.07	0.42
1:AA:2284:U:H5''	1:AA:2285:A:OP1	2.20	0.42
1:AA:228:U:H2'	1:AA:229:G:O4'	2.18	0.42
1:AA:2710:U:H2'	1:AA:2711:C:C6	2.54	0.42
1:AA:478:G:H2'	1:AA:479:C:H6	1.85	0.42
7:AG:136:ARG:HG3	7:AG:137:GLU:HG3	2.01	0.42
15:AR:54:LEU:HD12	15:AR:54:LEU:HA	1.59	0.42
1:AA:1042:A:H4'	18:AU:91:ASP:OD2	2.18	0.42
19:AV:81:TYR:C	19:AV:82:ARG:HD2	2.40	0.42
21:AX:61:GLY:HA3	21:AX:73:ARG:O	2.19	0.42
17:AT:118:ARG:HG2	34:BA:1442(A):G:C8	2.55	0.42
34:BA:148:G:H2'	34:BA:149:A:H8	1.84	0.42
34:BA:681:C:H6	34:BA:681:C:O5'	2.02	0.42
34:BA:685:G:O2'	34:BA:686:U:H5'	2.20	0.42
37:BD:104:VAL:HG12	37:BD:105:VAL:N	2.35	0.42
43:BJ:48:THR:O	47:BN:34:TYR:OH	2.36	0.42
44:BK:97:ALA:O	44:BK:101:SER:HB3	2.20	0.42
53:BT:12:ALA:O	53:BT:15:ARG:HB2	2.20	0.42
57:BZ:201:ILE:HB	57:BZ:206:LEU:HD13	2.01	0.42
57:BZ:549:ALA:HB2	57:BZ:587:SER:HB2	2.00	0.42
1:CA:1831:G:H2'	1:CA:1832:C:C6	2.55	0.42
1:CA:2167:U:OP1	1:CA:2167:U:H4'	2.20	0.42
1:CA:2697:G:C2	1:CA:2711:A:C2	3.08	0.42
1:CA:2784:C:H6	1:CA:2784:C:O5'	2.03	0.42
1:CA:878:A:C6	1:CA:900:A:C8	3.07	0.42
1:CA:897:C:OP2	1:CA:897:C:H6	2.03	0.42
1:CA:921:G:C6	1:CA:922:U:C4	3.08	0.42
5:CE:176:ILE:O	5:CE:176:ILE:HG22	2.19	0.42
5:CE:171:GLU:O	5:CE:184:VAL:HG23	2.19	0.42
7:CG:164:GLU:CD	7:CG:164:GLU:H	2.20	0.42
7:CG:86:MET:HA	7:CG:87:PRO:HD2	1.95	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1063:G:H1'	10:CL:134:MET:HA	2.01	0.42
11:CN:39:ARG:HB3	11:CN:41:ASP:OD1	2.19	0.42
12:CO:107:ARG:NH2	17:CT:36:GLU:OE2	2.53	0.42
17:CT:41:ARG:NH2	34:DA:345:C:H3'	2.34	0.42
17:CT:56:GLY:N	17:CT:59:THR:HG22	2.35	0.42
34:DA:1327:C:H2'	34:DA:1328:C:H6	1.85	0.42
34:DA:1512:U:O2'	34:DA:1513:A:H5'	2.19	0.42
34:DA:304:U:H2'	34:DA:305:G:C8	2.54	0.42
34:DA:786:G:C2	34:DA:797:C:C2	3.08	0.42
35:DB:100:GLY:N	35:DB:176:GLU:OE2	2.51	0.42
37:DD:94:LEU:HA	37:DD:94:LEU:HD23	1.59	0.42
38:DE:8:GLU:OE2	38:DE:63:ARG:NH2	2.52	0.42
39:DF:78:GLU:C	39:DF:80:ARG:N	2.72	0.42
42:DI:127:LYS:O	42:DI:128:ARG:HB3	2.19	0.42
34:DA:972:C:O2'	43:DJ:55:LYS:O	2.35	0.42
44:DK:48:ILE:HG21	44:DK:63:LEU:HD12	2.01	0.42
51:DR:74:ARG:HG3	51:DR:74:ARG:H	1.68	0.42
57:DZ:328:ILE:H	57:DZ:328:ILE:HD13	1.85	0.42
26:A2:53:LEU:HA	26:A2:53:LEU:HD23	1.71	0.42
1:AA:1007:G:C6	1:AA:1008:U:C4	3.07	0.42
1:AA:1756:U:H1'	1:AA:2870:A:N3	2.35	0.42
1:AA:1841:A:H2'	1:AA:1842:G:O4'	2.20	0.42
1:AA:239:G:C6	1:AA:240:A:C6	3.08	0.42
1:AA:464:G:H2'	1:AA:465:G:O4'	2.20	0.42
1:AA:746:A:H2'	1:AA:747:G:O4'	2.19	0.42
2:AB:73:A:H5'	2:AB:74:U:OP2	2.19	0.42
3:AC:42:VAL:O	3:AC:216:THR:C	2.58	0.42
3:AC:48:LEU:CD2	3:AC:59:VAL:HG21	2.50	0.42
4:AD:34:VAL:HA	4:AD:62:TYR:O	2.20	0.42
6:AF:139:PHE:HB2	6:AF:166:ALA:HB1	2.01	0.42
12:AO:24:VAL:O	12:AO:24:VAL:HG22	2.20	0.42
1:AA:2575:U:H4'	12:AO:28:SER:HA	2.02	0.42
14:AQ:59:ARG:O	23:AZ:180:VAL:HB	2.20	0.42
16:AS:111:GLU:O	16:AS:112:PHE:HB3	2.20	0.42
17:AT:15:VAL:HG13	17:AT:79:HIS:HE1	1.85	0.42
1:AA:1043:G:OP1	18:AU:92:ARG:HB2	2.19	0.42
21:AX:44:GLU:HG2	21:AX:49:VAL:O	2.20	0.42
23:AZ:151:HIS:ND1	23:AZ:151:HIS:N	2.68	0.42
34:BA:1182:G:C4'	34:BA:1183:A:H5'	2.44	0.42
34:BA:1513:A:H2'	34:BA:1514:C:C6	2.55	0.42
34:BA:21:G:H2'	34:BA:22:G:C8	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:271:C:H2'	34:BA:272:C:H6	1.85	0.42
34:BA:658:G:H2'	34:BA:659:U:C6	2.55	0.42
34:BA:715:A:H2'	34:BA:716:A:C8	2.55	0.42
36:BC:19:GLU:O	36:BC:56:ASP:HA	2.20	0.42
37:BD:202:LEU:HD23	37:BD:202:LEU:HA	1.83	0.42
40:BG:73:MET:HG2	40:BG:145:ALA:HB1	2.02	0.42
41:BH:17:THR:HG22	41:BH:63:LEU:HG	2.02	0.42
42:BI:5:TYR:HA	42:BI:17:VAL:O	2.20	0.42
44:BK:48:ILE:O	44:BK:48:ILE:HG12	2.18	0.42
44:BK:92:GLU:O	44:BK:96:ARG:HB2	2.20	0.42
57:BZ:227:ILE:HD11	57:BZ:241:GLU:HG3	2.01	0.42
57:BZ:68:ALA:HB3	57:BZ:327:PHE:CD1	2.54	0.42
61:BZ:703:FUA:C15	61:BZ:703:FUA:H323	2.49	0.42
25:C1:91:LYS:HA	25:C1:94:LEU:HD12	2.02	0.42
1:CA:1197:G:O2'	1:CA:1198:U:H5'	2.20	0.42
1:CA:1221:C:N4	1:CA:1229:G:H1	2.18	0.42
1:CA:585:G:H2'	1:CA:1251:C:H42	1.85	0.42
1:CA:2815:C:H2'	1:CA:2816:C:O4'	2.20	0.42
1:CA:449:A:O2'	18:CU:3:ARG:NH1	2.52	0.42
1:CA:64:A:H2'	1:CA:65:C:O4'	2.20	0.42
1:CA:856:C:H2'	1:CA:857:C:C6	2.55	0.42
1:CA:867:C:O2	1:CA:913:U:H5'	2.20	0.42
1:CA:949:C:H2'	1:CA:950:G:C8	2.55	0.42
2:CB:58:A:N7	2:CB:59:A:N7	2.67	0.42
10:CL:134:MET:HG2	10:CL:134:MET:H	1.61	0.42
11:CN:18:ALA:HB2	11:CN:54:VAL:CG1	2.49	0.42
13:CP:45:LEU:HA	13:CP:45:LEU:HD23	1.60	0.42
5:CE:110:GLY:O	15:CR:3:HIS:HE1	2.03	0.42
34:DA:1239:A:H4'	34:DA:1240:U:H5''	2.00	0.42
34:DA:1244:C:C2	34:DA:1294:G:N2	2.87	0.42
34:DA:195:A:C6	34:DA:196:A:N1	2.87	0.42
34:DA:377:G:O2'	34:DA:378:G:H5'	2.19	0.42
34:DA:920:U:N3	34:DA:921:U:C4	2.88	0.42
38:DE:145:LYS:HE2	38:DE:145:LYS:HB3	1.77	0.42
46:DM:22:ILE:HB	46:DM:25:ILE:HB	2.01	0.42
56:DW:63:G:H2'	56:DW:64:A:C8	2.55	0.42
57:DZ:251:ILE:HD13	57:DZ:285:ASP:HB3	2.02	0.42
24:A0:32:ARG:O	24:A0:33:ALA:C	2.58	0.42
25:A1:86:SER:OG	25:A1:89:GLU:HG2	2.20	0.42
31:A7:38:GLY:O	31:A7:39:ARG:C	2.59	0.42
31:A7:5:TRP:CD1	31:A7:7:PRO:HD3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:1605:A:H4'	1:AA:1606:G:H2'	2.00	0.42
1:AA:2753:A:C6	1:AA:2777:A:C8	3.08	0.42
1:AA:304:C:H2'	1:AA:305:G:O4'	2.20	0.42
1:AA:864:C:O2'	1:AA:886:U:H5''	2.20	0.42
1:AA:915:U:C4	1:AA:916:G:N7	2.88	0.42
2:AB:46:A:C5	2:AB:47:C:C4	3.07	0.42
3:AC:11:LEU:HD11	3:AC:35:THR:HG23	2.01	0.42
5:AE:84:PHE:CZ	5:AE:86:PRO:HB3	2.55	0.42
8:AH:11:VAL:HA	8:AH:12:PRO:HD3	1.88	0.42
12:AO:98:VAL:HG13	12:AO:117:LEU:HB3	2.01	0.42
12:AO:64:ARG:HD2	12:AO:81:ASP:HB3	2.02	0.42
23:AZ:61:LEU:HD22	23:AZ:61:LEU:H	1.85	0.42
34:BA:988:G:C2	34:BA:1218:C:O2	2.73	0.42
34:BA:1452:C:O2'	34:BA:1457:G:C8	2.73	0.42
34:BA:741:G:C2	34:BA:742:G:C4	3.08	0.42
34:BA:921:U:H2'	34:BA:922:G:O4'	2.20	0.42
35:BB:19:HIS:HA	35:BB:39:ILE:HG21	2.02	0.42
40:BG:151:TYR:OH	44:BK:54:ARG:NH1	2.53	0.42
41:BH:65:TYR:HA	41:BH:79:VAL:HG23	2.02	0.42
44:BK:20:TYR:CD1	44:BK:83:ILE:HB	2.54	0.42
45:BL:27:LEU:HD13	45:BL:98:TYR:HE2	1.83	0.42
49:BP:23:ASP:OD1	49:BP:24:ALA:N	2.53	0.42
57:BZ:99:ARG:HB2	57:BZ:128:TYR:CD1	2.55	0.42
57:BZ:82:ILE:HD12	57:BZ:101:LEU:HD11	2.02	0.42
1:CA:577:G:O2'	1:CA:1254:A:OP1	2.37	0.42
1:CA:129:C:H2'	1:CA:130:C:H6	1.85	0.42
1:CA:2136:C:O2'	1:CA:2137:C:P	2.78	0.42
1:CA:2190:G:H2'	1:CA:2191:G:H8	1.85	0.42
1:CA:554:U:C4	1:CA:555:U:C4	3.08	0.42
1:CA:876:C:H2'	1:CA:877:U:O4'	2.20	0.42
1:CA:909:A:O2'	1:CA:910:A:H5'	2.19	0.42
1:CA:948:G:C6	1:CA:949:C:C4	3.07	0.42
3:CC:42:VAL:O	3:CC:216:THR:C	2.58	0.42
4:CD:124:PRO:HD2	4:CD:129:ASN:HD22	1.85	0.42
16:CS:101:LEU:HD23	16:CS:101:LEU:O	2.20	0.42
21:CX:44:GLU:HG2	21:CX:49:VAL:O	2.20	0.42
34:DA:1096:C:HO2'	34:DA:1170:A:HO2'	1.66	0.42
34:DA:114:U:H2'	34:DA:115:G:C8	2.54	0.42
34:DA:1402:C:O2	34:DA:1500:A:N1	2.52	0.42
34:DA:319:G:C2	34:DA:320:C:C2	3.08	0.42
34:DA:513:C:N3	34:DA:539:A:C2	2.88	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:DC:181:ASN:OD1	36:DC:204:LEU:HD12	2.20	0.42
34:DA:1298:C:N4	40:DG:114:ARG:HB3	2.35	0.42
44:DK:67:ASP:O	44:DK:71:LYS:HG3	2.19	0.42
50:DQ:27:PHE:CE2	50:DQ:36:ILE:HD11	2.55	0.42
34:DA:1400:C:C2	56:DW:34:G:C2	3.07	0.42
56:DW:41:C:O5'	56:DY:36:A:OP1	2.38	0.42
57:DZ:28:THR:O	57:DZ:32:ILE:HG13	2.19	0.42
57:DZ:330:VAL:HG21	57:DZ:369:LEU:HB3	2.02	0.42
57:DZ:80:ASN:HD22	57:DZ:374:LEU:HB2	1.84	0.42
27:A3:5:LYS:HZ1	27:A3:55:ARG:HH12	1.67	0.41
28:A4:68:ARG:H	28:A4:68:ARG:HG2	1.57	0.41
1:AA:2092:G:H2'	1:AA:2093:A:O4'	2.20	0.41
1:AA:2123:G:H1	1:AA:2210:C:H42	1.68	0.41
1:AA:2285:A:H2'	1:AA:2286:A:C8	2.54	0.41
1:AA:2371:C:O5'	1:AA:2371:C:H6	2.03	0.41
1:AA:2085:C:O2	1:AA:2462:A:N1	2.53	0.41
1:AA:274:U:H6	1:AA:274:U:H2'	1.67	0.41
2:AB:12:C:H2'	24:A0:73:GLY:HA3	2.02	0.41
8:AH:17:VAL:HG11	8:AH:50:VAL:HG21	2.02	0.41
1:AA:1126:C:O2'	10:AL:126:MET:HG3	2.20	0.41
10:AL:78:ILE:HD13	10:AL:134:MET:SD	2.60	0.41
10:AL:84:LEU:HD21	10:AL:96:VAL:HB	2.02	0.41
11:AN:40:PRO:O	18:AU:64:ARG:HD3	2.19	0.41
14:AQ:2:LEU:HA	14:AQ:2:LEU:HD12	1.74	0.41
15:AR:104:ARG:NH1	15:AR:107:ASP:OD2	2.53	0.41
17:AT:118:ARG:HG3	17:AT:118:ARG:HH11	1.85	0.41
34:BA:1112:C:O2	36:BC:179:ARG:N	2.45	0.41
34:BA:142:G:H2'	34:BA:143:A:C8	2.55	0.41
34:BA:354:G:C2'	34:BA:355:C:H5'	2.49	0.41
34:BA:452:A:O2'	34:BA:453:A:OP2	2.29	0.41
34:BA:548:G:H2'	34:BA:549:C:C6	2.54	0.41
34:BA:565:U:C4	34:BA:566:G:C5	3.07	0.41
34:BA:657:G:O2'	34:BA:658:G:H5'	2.19	0.41
34:BA:900:A:O5'	34:BA:900:A:H8	2.03	0.41
34:BA:987:G:H2'	34:BA:988:G:H8	1.84	0.41
39:BF:41:GLU:O	39:BF:43:LEU:HD12	2.19	0.41
57:BZ:181:LEU:HD23	57:BZ:182:ARG:CG	2.50	0.41
57:BZ:541:ALA:HB3	57:BZ:579:GLU:HG2	2.01	0.41
24:C0:23:VAL:HG22	24:C0:38:VAL:HG22	2.01	0.41
27:C3:4:LEU:HA	27:C3:4:LEU:HD23	1.77	0.41
1:CA:1233:C:C2'	1:CA:1234:U:H5'	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:1363:C:H2'	1:CA:1364:G:H8	1.84	0.41
1:CA:1469:A:H2'	1:CA:1470:G:O4'	2.19	0.41
1:CA:171:G:H2'	1:CA:172:C:C6	2.55	0.41
1:CA:1851:U:H2'	1:CA:1852:C:O4'	2.20	0.41
1:CA:1857:G:C6	1:CA:1858:G:N1	2.88	0.41
1:CA:1794:U:H1'	1:CA:1900:A:N3	2.35	0.41
1:CA:1914:C:O2	1:CA:1914:C:H2'	2.20	0.41
1:CA:2689:U:H2'	1:CA:2689:U:O2	2.20	0.41
1:CA:2690:C:N4	1:CA:2713:A:H1'	2.35	0.41
1:CA:384:U:H2'	1:CA:385:C:H6	1.85	0.41
1:CA:470:A:H8	1:CA:470:A:C5'	2.33	0.41
1:CA:844:C:C5	1:CA:845:G:C6	3.07	0.41
5:CE:25:VAL:HG21	17:CT:7:ILE:HD13	2.02	0.41
6:CF:178:PRO:HB2	6:CF:201:VAL:CG2	2.50	0.41
10:CL:119:ASP:HB3	10:CL:120:LEU:H	1.63	0.41
12:CO:98:VAL:HG13	12:CO:117:LEU:HB3	2.02	0.41
1:CA:252:G:P	13:CP:50:ARG:HH12	2.43	0.41
16:CS:15:ARG:HD3	16:CS:25:ARG:NH2	2.34	0.41
21:CX:12:VAL:HG22	21:CX:29:TRP:CD1	2.54	0.41
34:DA:300:A:H2'	34:DA:301:G:O4'	2.19	0.41
34:DA:451:A:N6	34:DA:480:U:H2'	2.35	0.41
35:DB:113:HIS:HA	35:DB:116:GLU:HG2	2.01	0.41
36:DC:87:LEU:O	36:DC:91:LEU:HB2	2.20	0.41
42:DI:6:GLY:O	42:DI:17:VAL:HG12	2.20	0.41
46:DM:121:LYS:HB2	57:DZ:507:TYR:OH	2.20	0.41
57:DZ:90:PHE:O	57:DZ:93:GLU:HB2	2.20	0.41
1:AA:1220:U:H1'	1:AA:1221:G:OP1	2.19	0.41
1:AA:1463:C:H4'	1:AA:1633:A:H2	1.85	0.41
1:AA:1475:G:O2'	1:AA:1476:C:H5'	2.20	0.41
1:AA:207:A:C2	1:AA:224:U:H4'	2.55	0.41
1:AA:2310:A:H2'	1:AA:2311:G:O4'	2.20	0.41
1:AA:507:G:C4	1:AA:532:A:C2	3.08	0.41
1:AA:520:G:C6	1:AA:521:G:C5	3.08	0.41
1:AA:648:G:H2'	1:AA:649:C:C6	2.55	0.41
8:AH:103:LEU:HG	8:AH:105:LEU:HD13	2.02	0.41
12:AO:14:THR:HG21	12:AO:86:ILE:HB	2.03	0.41
12:AO:71:ARG:HA	12:AO:72:PRO:HD3	1.95	0.41
18:AU:20:LEU:HA	18:AU:20:LEU:HD23	1.79	0.41
22:AY:92:ASN:HB3	22:AY:94:LYS:HG3	2.02	0.41
34:BA:1221:G:OP1	34:BA:1320:C:N4	2.40	0.41
34:BA:547:A:H4'	34:BA:548:G:O5'	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:748:C:H6	34:BA:748:C:H2'	1.65	0.41
34:BA:838:G:N2	34:BA:849:C:C2	2.88	0.41
35:BB:48:MET:HA	35:BB:51:LEU:HD12	2.02	0.41
36:BC:113:ALA:HB3	36:BC:114:PRO:HD3	2.02	0.41
37:BD:18:LYS:HG2	37:BD:33:MET:CG	2.48	0.41
37:BD:50:ARG:HA	37:BD:51:PRO:HD2	1.82	0.41
45:BL:55:VAL:HG12	45:BL:56:ALA:N	2.34	0.41
49:BP:21:VAL:CG1	49:BP:34:GLU:HB3	2.50	0.41
52:BS:58:VAL:HA	52:BS:59:PRO:HD2	1.78	0.41
57:BZ:-62:LEU:H	57:BZ:-62:LEU:HD12	1.85	0.41
57:BZ:89:ASP:O	57:BZ:454:MET:HB3	2.19	0.41
26:C2:66:GLU:HA	26:C2:69:ARG:NH1	2.35	0.41
30:C6:11:LEU:HA	30:C6:11:LEU:HD23	1.83	0.41
1:CA:1157:G:C2	1:CA:1158:C:C2	3.08	0.41
1:CA:2690:C:H6	1:CA:2690:C:OP2	2.03	0.41
1:CA:2830:G:N3	1:CA:2883:A:H2	2.18	0.41
1:CA:317:G:N2	1:CA:334:C:O2	2.48	0.41
1:CA:448:U:C4	1:CA:583:G:H1'	2.55	0.41
1:CA:830:G:H5''	63:CA:3984:HOH:O	2.20	0.41
1:CA:897:C:H3'	1:CA:898:C:H6	1.85	0.41
4:CD:70:TRP:CZ2	4:CD:150:LYS:HA	2.55	0.41
5:CE:8:LYS:HD2	5:CE:188:VAL:HG12	2.01	0.41
6:CF:158:THR:HG1	6:CF:160:ASN:H	1.60	0.41
8:CH:94:TYR:HA	8:CH:106:THR:O	2.20	0.41
10:CL:99:ILE:HG23	10:CL:103:GLN:CB	2.51	0.41
13:CP:127:ALA:O	13:CP:148:LEU:HD23	2.19	0.41
13:CP:56:SER:OG	13:CP:61:ARG:HD2	2.20	0.41
14:CQ:1:MET:HB2	14:CQ:2:LEU:H	1.71	0.41
15:CR:61:HIS:CD2	15:CR:65:LEU:HD21	2.55	0.41
1:CA:995:C:C2	18:CU:57:PHE:CE1	3.08	0.41
20:CW:23:LEU:HA	20:CW:23:LEU:HD12	1.81	0.41
34:DA:1486:G:H2'	34:DA:1487:G:C8	2.55	0.41
34:DA:317:G:C6	34:DA:318:G:N7	2.88	0.41
34:DA:438:G:N1	34:DA:495:A:OP2	2.42	0.41
34:DA:515:G:C4	34:DA:537:G:N2	2.89	0.41
34:DA:681:C:C2	34:DA:710:G:C2	3.07	0.41
34:DA:976:G:H22	34:DA:1362:C:H2'	1.86	0.41
37:DD:31:CYS:SG	37:DD:33:MET:HB2	2.60	0.41
48:DO:43:LEU:HA	48:DO:43:LEU:HD23	1.82	0.41
57:DZ:17:ILE:H	57:DZ:17:ILE:HD12	1.85	0.41
57:DZ:354:ARG:HB3	57:DZ:378:VAL:HB	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:403:GLU:O	57:DZ:404:VAL:HG23	2.20	0.41
57:DZ:546:ILE:HG13	57:DZ:590:ILE:HB	2.02	0.41
32:A8:52:LYS:HB3	32:A8:52:LYS:HE2	1.81	0.41
1:AA:1356:G:C2'	1:AA:1357:G:H5'	2.50	0.41
1:AA:1544:C:O4'	1:AA:1624:C:H4'	2.20	0.41
1:AA:1841:A:H2'	1:AA:1842:G:H5'	2.02	0.41
1:AA:1944:G:H2'	1:AA:1945:U:O4'	2.21	0.41
1:AA:2803:A:H2'	1:AA:2803:A:N3	2.35	0.41
1:AA:842:C:H2'	1:AA:843:C:H6	1.83	0.41
3:AC:60:ARG:NH2	3:AC:165:ARG:HH21	2.18	0.41
5:AE:101:ARG:HD2	5:AE:169:ASN:OD1	2.20	0.41
5:AE:176:ILE:HB	5:AE:181:LEU:HB2	2.03	0.41
7:AG:178:PHE:O	7:AG:180:PHE:CD2	2.73	0.41
12:AO:63:VAL:HG11	12:AO:85:VAL:HG23	2.01	0.41
13:AP:20:GLY:HA2	13:AP:28:GLY:HA2	2.01	0.41
16:AS:94:TYR:CE1	16:AS:99:LYS:HG3	2.55	0.41
22:AY:81:LYS:HB3	22:AY:81:LYS:HE2	1.97	0.41
23:AZ:111:VAL:C	23:AZ:113:ALA:N	2.73	0.41
34:BA:1104:G:OP1	35:BB:111:ARG:HD2	2.20	0.41
34:BA:1164:G:C2'	34:BA:1165:C:H5'	2.50	0.41
34:BA:1411:C:H2'	34:BA:1412:C:H6	1.84	0.41
34:BA:175:C:C2	34:BA:176:C:C5	3.08	0.41
34:BA:425:G:C2	34:BA:426:G:C8	3.09	0.41
34:BA:495:A:H4'	34:BA:496:A:OP1	2.21	0.41
34:BA:922:G:H3'	34:BA:923:A:C8	2.55	0.41
34:BA:955:U:H2'	34:BA:956:U:C6	2.52	0.41
37:BD:15:GLU:HG3	37:BD:63:LYS:HD3	2.02	0.41
37:BD:194:LEU:HD12	37:BD:195:ALA:H	1.84	0.41
46:BM:20:THR:C	46:BM:22:ILE:H	2.24	0.41
48:BO:62:GLN:HA	48:BO:65:ARG:NH1	2.34	0.41
1:CA:1154:G:OP2	18:CU:58:ARG:NH2	2.50	0.41
1:CA:1465:G:N1	1:CA:1466:G:C5	2.88	0.41
1:CA:1477:A:H2'	1:CA:1478:G:O4'	2.20	0.41
1:CA:2166:G:N2	1:CA:2167:U:O4	2.54	0.41
1:CA:224:G:H2'	1:CA:225:A:O4'	2.20	0.41
1:CA:2294:C:H2'	1:CA:2295:C:H6	1.85	0.41
1:CA:2854:G:C4	1:CA:2864:G:N2	2.88	0.41
1:CA:710:G:H2'	1:CA:711:G:C8	2.55	0.41
4:CD:68:LYS:O	4:CD:70:TRP:N	2.53	0.41
5:CE:170:LEU:HB3	5:CE:184:VAL:CG2	2.50	0.41
11:CN:94:HIS:O	11:CN:97:ARG:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:CO:2:ILE:CD1	12:CO:6:THR:HG21	2.47	0.41
15:CR:50:HIS:CE1	15:CR:54:LEU:HD11	2.55	0.41
20:CW:79:GLY:CA	20:CW:100:THR:HG22	2.50	0.41
23:CZ:45:ASP:OD1	23:CZ:49:ARG:HG3	2.20	0.41
23:CZ:53:ILE:HG22	23:CZ:71:VAL:HB	2.01	0.41
35:DB:163:PHE:HA	35:DB:185:ILE:O	2.20	0.41
36:DC:110:ASN:ND2	36:DC:144:SER:OG	2.52	0.41
37:DD:171:GLY:HA2	37:DD:172:PRO:HD2	1.83	0.41
38:DE:71:LEU:O	38:DE:72:GLN:HG2	2.20	0.41
39:DF:61:LEU:HD23	39:DF:63:TYR:OH	2.21	0.41
42:DI:17:VAL:HG23	42:DI:63:ILE:HG12	2.02	0.41
45:DL:27:LEU:HD13	45:DL:98:TYR:CE1	2.55	0.41
49:DP:65:GLN:HA	49:DP:66:PRO:HD2	1.94	0.41
50:DQ:80:GLY:O	50:DQ:82:MET:HG2	2.19	0.41
51:DR:51:LEU:HA	51:DR:52:PRO:HD3	1.85	0.41
52:DS:27:GLU:HB3	52:DS:28:LYS:HG2	2.01	0.41
56:DW:46:7MG:H81	56:DW:46:7MG:H2'	1.85	0.41
56:DW:65:G:H2'	56:DW:66:U:H6	1.86	0.41
57:DZ:110:SER:HB2	57:DZ:144:ALA:HB1	2.03	0.41
57:DZ:157:LEU:HD12	57:DZ:158:GLY:N	2.34	0.41
57:DZ:20:HIS:HB2	57:DZ:118:SER:HB2	2.03	0.41
57:DZ:647:VAL:HG11	57:DZ:652:MET:SD	2.60	0.41
1:AA:1065:U:O2'	1:AA:1067:A:C2	2.62	0.41
1:AA:2099:A:O2'	1:AA:2100:C:H5'	2.20	0.41
1:AA:291:G:N2	1:AA:395:C:C2	2.89	0.41
6:AF:149:ASP:OD2	6:AF:151:SER:N	2.46	0.41
1:AA:470:C:H4'	6:AF:49:ALA:HB2	2.01	0.41
6:AF:72:ARG:HE	6:AF:72:ARG:HB3	1.52	0.41
1:AA:2315:G:O2'	7:AG:132:ASN:HB2	2.20	0.41
7:AG:44:GLY:C	7:AG:46:ALA:H	2.22	0.41
8:AH:164:TYR:O	8:AH:167:GLU:HB3	2.21	0.41
8:AH:41:MET:HE3	8:AH:41:MET:HB3	1.90	0.41
10:AL:95:LYS:HA	10:AL:135:GLY:O	2.20	0.41
11:AN:107:LEU:HD23	11:AN:107:LEU:HA	1.80	0.41
18:AU:52:ARG:HH11	18:AU:52:ARG:HG3	1.85	0.41
23:AZ:31:ARG:HD2	23:AZ:31:ARG:HH11	1.72	0.41
34:BA:1423:G:C6	34:BA:1424:C:C4	3.09	0.41
34:BA:1516:G:H2'	34:BA:1518:A:OP2	2.20	0.41
34:BA:291:C:O2'	34:BA:292:G:H5'	2.19	0.41
34:BA:345:C:H4'	34:BA:346:G:N1	2.35	0.41
34:BA:405:U:H3'	34:BA:406:G:H5'	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:522:C:H2'	34:BA:523:A:H5'	2.01	0.41
34:BA:69:G:C2	34:BA:70:G:C5	3.08	0.41
35:BB:76:GLN:H	35:BB:76:GLN:HG2	1.51	0.41
34:BA:1112:C:H1'	36:BC:179:ARG:HE	1.84	0.41
37:BD:108:LEU:HD11	37:BD:174:LEU:HB3	2.02	0.41
37:BD:5:ILE:HD13	37:BD:5:ILE:O	2.21	0.41
38:BE:7:GLU:CD	38:BE:37:ARG:HH21	2.23	0.41
43:BJ:5:ARG:HB3	43:BJ:5:ARG:HE	1.51	0.41
47:BN:13:THR:HA	47:BN:14:PRO:HD3	1.80	0.41
53:BT:66:ALA:HB1	53:BT:71:THR:HG21	2.01	0.41
57:BZ:358:MET:HE1	57:BZ:363:ARG:HG2	2.01	0.41
27:C3:23:LEU:HB3	27:C3:28:LEU:O	2.21	0.41
1:CA:1467:C:C2	1:CA:1526:G:N2	2.89	0.41
1:CA:1932:A:H61	1:CA:1968:G:H1'	1.84	0.41
1:CA:2273:A:H2'	1:CA:2274:A:C8	2.55	0.41
1:CA:242:G:C8	32:C8:5:LYS:HG2	2.56	0.41
1:CA:2563:U:O2	1:CA:2565:A:C8	2.74	0.41
2:CB:16:G:C6	2:CB:69:G:C2	3.08	0.41
1:CA:614(C):A:C4	6:CF:180:GLY:HA3	2.56	0.41
7:CG:60:LEU:HD22	7:CG:63:ILE:HD11	2.01	0.41
8:CH:154:PRO:HB3	8:CH:163:TYR:CE2	2.55	0.41
1:CA:1061:U:O4	10:CL:73:PRO:HG3	2.20	0.41
11:CN:62:VAL:HG12	11:CN:67:LEU:HD22	2.03	0.41
13:CP:138:LEU:HD23	13:CP:145:PRO:HG3	2.03	0.41
21:CX:44:GLU:O	21:CX:46:ALA:N	2.53	0.41
23:CZ:14:LYS:HA	23:CZ:15:PRO:HD3	1.89	0.41
34:DA:1000:U:H2'	34:DA:1001:A:C8	2.55	0.41
34:DA:1077:G:N1	34:DA:1081:G:C6	2.88	0.41
34:DA:1154:G:C6	34:DA:1155:G:C5	3.08	0.41
34:DA:1366:C:O2'	43:DJ:60:ARG:NH2	2.54	0.41
34:DA:922:G:O2'	34:DA:1398:A:N1	2.38	0.41
34:DA:1394:A:C6	34:DA:1501:C:H4'	2.56	0.41
34:DA:334:C:H2'	34:DA:335:C:H6	1.85	0.41
34:DA:561:U:O2'	34:DA:562:C:P	2.79	0.41
34:DA:600:C:H42	34:DA:638:G:H1	1.68	0.41
34:DA:690:G:H2'	34:DA:691:G:C8	2.55	0.41
35:DB:41:ILE:HA	35:DB:41:ILE:HD13	1.86	0.41
46:DM:91:ARG:O	46:DM:110:ARG:NH1	2.52	0.41
50:DQ:81:ARG:HA	50:DQ:81:ARG:HD2	1.86	0.41
57:DZ:346:LYS:HE2	57:DZ:384:ILE:HG23	2.02	0.41
57:DZ:409:ILE:HG23	57:DZ:459:LEU:HD12	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
24:A0:65:GLY:HA3	24:A0:81:VAL:HG12	2.03	0.41
29:A5:58:LEU:HD22	29:A5:60:VAL:HG23	2.03	0.41
1:AA:1281:G:C6	1:AA:1282:G:N1	2.89	0.41
1:AA:2418:U:C2	13:AP:72:PRO:HG2	2.56	0.41
1:AA:2825:C:O5'	1:AA:2825:C:H6	2.03	0.41
1:AA:2879:G:H2'	1:AA:2880:C:C6	2.56	0.41
1:AA:449:A:C6	1:AA:450:A:C6	3.09	0.41
1:AA:604:C:H2'	1:AA:605:G:H8	1.84	0.41
3:AC:11:LEU:HD22	3:AC:11:LEU:H	1.86	0.41
4:AD:218:ARG:HB3	4:AD:219:PRO:HD2	2.02	0.41
2:AB:41:U:H5	7:AG:70:VAL:HB	1.86	0.41
1:AA:2679:C:O2	8:AH:109:PHE:HB3	2.19	0.41
19:AV:40:LEU:HA	19:AV:40:LEU:HD23	1.72	0.41
19:AV:19:LYS:HA	19:AV:94:LEU:O	2.19	0.41
34:BA:960:U:H2'	34:BA:1225:A:H62	1.84	0.41
34:BA:199:G:N2	34:BA:219:C:O2	2.53	0.41
34:BA:266:G:N2	63:BA:5136:HOH:O	2.53	0.41
34:BA:416:G:C6	34:BA:417:C:N3	2.88	0.41
34:BA:627:G:O2'	34:BA:628:G:H5'	2.21	0.41
34:BA:604:G:C2	34:BA:635:G:C4	3.09	0.41
35:BB:223:ILE:H	35:BB:223:ILE:HG12	1.35	0.41
39:BF:1:MET:HE3	39:BF:66:GLU:HB3	2.03	0.41
45:BL:84:LEU:HB2	45:BL:105:TYR:CD2	2.56	0.41
52:BS:70:LYS:HD3	52:BS:70:LYS:HA	1.95	0.41
57:BZ:123:ARG:O	57:BZ:126:GLU:HB2	2.20	0.41
57:BZ:443:HIS:HA	57:BZ:444:PRO:HD2	1.86	0.41
1:CA:54:G:O2'	31:C7:35:ARG:HD3	2.21	0.41
1:CA:1401:G:C5	1:CA:1402:C:C4	3.08	0.41
1:CA:1488:G:N2	1:CA:1502:C:C2	2.89	0.41
1:CA:14:A:N1	1:CA:2044:C:O2'	2.40	0.41
1:CA:1478:G:O2'	1:CA:1558:A:C2	2.70	0.41
1:CA:1575:C:H2'	1:CA:1576:U:O4'	2.21	0.41
1:CA:1271:G:N2	1:CA:1617:C:O4'	2.53	0.41
1:CA:189:G:H2'	1:CA:205:G:N2	2.36	0.41
1:CA:319:C:H2'	1:CA:320:A:O4'	2.20	0.41
5:CE:163:GLU:HG2	5:CE:164:ARG:H	1.86	0.41
10:CL:73:PRO:HB2	10:CL:76:TYR:H	1.85	0.41
16:CS:61:ASN:O	16:CS:65:VAL:HG23	2.21	0.41
19:CV:98:GLU:OE1	19:CV:100:ARG:NH1	2.48	0.41
20:CW:59:VAL:HG22	20:CW:64:MET:O	2.20	0.41
21:CX:12:VAL:CG2	21:CX:27:THR:HG22	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:1012:U:H2'	34:DA:1013:G:C8	2.55	0.41
34:DA:1055:A:C6	34:DA:1206:G:C5	3.09	0.41
34:DA:1278:U:H5'	34:DA:1279:A:C5'	2.51	0.41
34:DA:348:G:O2'	34:DA:349:A:H5'	2.20	0.41
34:DA:358:U:H2'	34:DA:359:U:C6	2.55	0.41
34:DA:539:A:N6	34:DA:540:G:O6	2.53	0.41
34:DA:778:G:C6	34:DA:779:C:N3	2.89	0.41
34:DA:863:U:H2'	34:DA:865:A:OP2	2.21	0.41
34:DA:961:U:H2'	34:DA:962:C:O4'	2.21	0.41
35:DB:95:GLN:HB2	35:DB:96:ARG:H	1.73	0.41
38:DE:11:ILE:HD13	38:DE:105:VAL:HG13	2.02	0.41
50:DQ:6:LEU:HA	50:DQ:6:LEU:HD12	1.86	0.41
57:DZ:19:ALA:HB3	57:DZ:25:LYS:HG3	2.02	0.41
57:DZ:631:ILE:HG13	57:DZ:632:LEU:N	2.36	0.41
61:DZ:703:FUA:C12	61:DZ:703:FUA:H231	2.45	0.41
1:AA:2102:G:OP1	25:A1:35:THR:HG21	2.20	0.41
30:A6:8:LYS:HE2	32:A8:34:TRP:CZ3	2.55	0.41
1:AA:1355:G:P	31:A7:9:ARG:HD3	2.60	0.41
1:AA:1668:G:C2	1:AA:1669:G:C8	3.09	0.41
1:AA:2169:G:H2'	1:AA:2170:G:O4'	2.20	0.41
1:AA:2211:U:H2'	1:AA:2212:G:H8	1.86	0.41
1:AA:278:G:C2	1:AA:279:G:C8	3.08	0.41
1:AA:718:C:N4	63:AA:4823:HOH:O	2.53	0.41
3:AC:167:ASP:OD1	3:AC:169:THR:OG1	2.38	0.41
3:AC:44:VAL:HG23	3:AC:176:VAL:CG2	2.51	0.41
4:AD:25:THR:HG21	4:AD:113:VAL:HG11	2.02	0.41
5:AE:98:PRO:HD3	5:AE:175:VAL:HG13	2.03	0.41
6:AF:164:ARG:O	6:AF:168:ARG:HB2	2.19	0.41
6:AF:185:ASP:OD1	6:AF:188:ARG:NH1	2.49	0.41
11:AN:112:LEU:HD12	11:AN:112:LEU:O	2.21	0.41
23:AZ:121:HIS:HE1	23:AZ:169:GLU:OE2	2.02	0.41
34:BA:116:A:H61	34:BA:313:A:H1'	1.84	0.41
34:BA:1207:G:H2'	34:BA:1208:C:H6	1.86	0.41
34:BA:1428:A:H2'	34:BA:1429:C:O4'	2.19	0.41
34:BA:389:A:C5	34:BA:390:C:H1'	2.56	0.41
34:BA:690:G:C6	34:BA:691:G:C6	3.09	0.41
34:BA:753:A:H4'	34:BA:754:C:C5'	2.49	0.41
34:BA:954:G:H2'	34:BA:955:U:C6	2.56	0.41
35:BB:28:PHE:CD1	35:BB:190:THR:HG22	2.55	0.41
35:BB:71:VAL:HG13	35:BB:93:VAL:HG23	2.03	0.41
37:BD:107:ARG:HH22	37:BD:194:LEU:CD1	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
38:BE:108:ALA:O	38:BE:110:LEU:N	2.53	0.41
45:BL:43:VAL:HG12	45:BL:44:THR:N	2.36	0.41
49:BP:4:ILE:HG21	49:BP:36:ILE:HD11	2.02	0.41
49:BP:73:LEU:HA	49:BP:76:GLN:HB3	2.02	0.41
56:BY:56:C:H2'	56:BY:57:G:O4'	2.20	0.41
57:BZ:339:SER:HB2	57:BZ:352:VAL:HG13	2.02	0.41
57:BZ:623:ASP:HB2	57:BZ:662:LYS:HD2	2.02	0.41
1:CA:1168:G:C2	1:CA:1182:A:C2	3.08	0.41
1:CA:1478:G:O2'	1:CA:1558:A:N1	2.50	0.41
1:CA:1773:A:C5	1:CA:1829:A:H1'	2.56	0.41
1:CA:2527:C:H2'	1:CA:2528:U:O4'	2.20	0.41
1:CA:2809:A:C2'	1:CA:2810:A:H5'	2.50	0.41
1:CA:357:A:H2'	1:CA:358:U:O4'	2.21	0.41
1:CA:464:U:H4'	31:C7:5:TRP:CZ3	2.56	0.41
2:CB:28:C:H2'	2:CB:29:A:H8	1.85	0.41
12:CO:20:MET:CE	12:CO:44:LYS:HE3	2.50	0.41
14:CQ:29:PHE:HB3	14:CQ:65:PHE:CE2	2.54	0.41
15:CR:24:GLN:OE1	15:CR:36:THR:HG21	2.19	0.41
16:CS:62:LYS:O	16:CS:65:VAL:HB	2.20	0.41
19:CV:7:THR:HB	19:CV:35:LEU:HD12	2.02	0.41
22:CY:15:VAL:HG21	22:CY:42:VAL:HG11	2.03	0.41
23:CZ:97:GLU:HA	23:CZ:126:VAL:O	2.20	0.41
34:DA:511:C:O3'	37:DD:43:HIS:CE1	2.74	0.41
34:DA:562:C:H4'	34:DA:563:A:O5'	2.20	0.41
35:DB:163:PHE:HA	35:DB:185:ILE:HG12	2.01	0.41
37:DD:23:GLY:HA3	37:DD:112:VAL:O	2.20	0.41
38:DE:105:VAL:HG21	38:DE:128:PRO:HB3	2.01	0.41
45:DL:125:PRO:HB2	45:DL:126:LYS:H	1.71	0.41
46:DM:10:PRO:HG2	46:DM:21:TYR:HD1	1.85	0.41
47:DN:41:ARG:HG3	47:DN:42:ILE:N	2.35	0.41
57:DZ:20:HIS:HB2	57:DZ:118:SER:OG	2.21	0.41
57:DZ:343:ASN:ND2	57:DZ:383:THR:HG23	2.36	0.41
61:DZ:703:FUA:C15	61:DZ:703:FUA:H323	2.49	0.41
25:A1:17:SER:HB2	25:A1:40:ARG:HG2	2.02	0.41
27:A3:26:LEU:HA	27:A3:26:LEU:HD23	1.80	0.41
33:A9:16:VAL:O	33:A9:17:ILE:HD12	2.21	0.41
1:AA:1373:C:H2'	1:AA:1374:G:O4'	2.21	0.41
1:AA:1537:G:C4	1:AA:1546:G:N2	2.89	0.41
1:AA:2022:G:C6	1:AA:2023:A:N7	2.89	0.41
1:AA:2047:C:H2'	1:AA:2048:C:H6	1.86	0.41
1:AA:2291:G:N7	24:A0:14:ARG:NH1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:2769:U:OP2	33:A9:19:ARG:NE	2.39	0.41
1:AA:27:G:C2	1:AA:537:G:N3	2.88	0.41
1:AA:616:G:N2	1:AA:712:C:C2	2.89	0.41
2:AB:73:A:N3	2:AB:73:A:H2'	2.36	0.41
1:AA:1537:G:O2'	4:AD:101:GLU:HB2	2.21	0.41
5:AE:116:VAL:HG13	5:AE:122:PHE:CB	2.50	0.41
6:AF:93:LYS:HA	6:AF:93:LYS:HD3	1.77	0.41
2:AB:41:U:C5	7:AG:70:VAL:HB	2.55	0.41
9:AK:48:GLY:C	9:AK:90:ALA:HB1	2.40	0.41
13:AP:97:PRO:HD3	13:AP:126:VAL:O	2.20	0.41
14:AQ:56:ARG:NH1	14:AQ:56:ARG:CG	2.82	0.41
17:AT:37:GLY:C	17:AT:39:ARG:H	2.23	0.41
14:AQ:62:GLY:O	23:AZ:178:GLU:HG2	2.20	0.41
34:BA:1024:G:H2'	34:BA:1025:U:H5''	2.03	0.41
34:BA:1127:G:H5'	34:BA:1280:A:O2'	2.21	0.41
34:BA:1299:A:H5''	34:BA:1299:A:N3	2.36	0.41
34:BA:1374:A:C4	34:BA:1375:A:C8	3.09	0.41
37:BD:79:PHE:CZ	37:BD:207:TYR:HD2	2.37	0.41
39:BF:92:LYS:HB2	39:BF:92:LYS:HE3	1.59	0.41
41:BH:78:GLN:HE21	41:BH:78:GLN:HB2	1.59	0.41
42:BI:28:VAL:HA	42:BI:63:ILE:O	2.21	0.41
49:BP:7:ALA:HB2	49:BP:20:VAL:HG11	2.03	0.41
50:BQ:41:LYS:HZ2	50:BQ:92:ARG:HH21	1.68	0.41
57:BZ:203:GLU:H	57:BZ:203:GLU:CD	2.24	0.41
57:BZ:271:LEU:HD12	57:BZ:271:LEU:H	1.85	0.41
57:BZ:555:LEU:HD11	57:BZ:599:PRO:HB2	2.02	0.41
57:BZ:590:ILE:O	57:BZ:594:VAL:HG23	2.21	0.41
24:C0:46:LYS:O	24:C0:78:TYR:HA	2.20	0.41
1:CA:1688:U:O2	1:CA:1700:A:H5'	2.19	0.41
1:CA:1823:G:OP1	4:CD:54:ARG:NH1	2.54	0.41
1:CA:2734:A:C8	1:CA:2735:G:C8	3.08	0.41
1:CA:777:A:C2	1:CA:778:G:C4	3.08	0.41
3:CC:48:LEU:CD2	3:CC:59:VAL:HG21	2.50	0.41
5:CE:182:LEU:HD12	5:CE:182:LEU:C	2.41	0.41
6:CF:162:LEU:HD12	6:CF:162:LEU:HA	1.79	0.41
6:CF:184:TYR:CE1	13:CP:3:LEU:HD21	2.56	0.41
17:CT:55:ASN:H	17:CT:59:THR:HG22	1.86	0.41
34:DA:1085:U:H3'	34:DA:1086:U:C5	2.56	0.41
34:DA:343:U:O3'	34:DA:344:A:H8	2.02	0.41
34:DA:411:A:C6	34:DA:429:U:C4	3.09	0.41
35:DB:72:GLY:O	35:DB:94:ASN:ND2	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
36:DC:105:GLU:HG3	36:DC:105:GLU:H	1.72	0.41
43:DJ:6:ILE:O	43:DJ:71:LEU:HD12	2.20	0.41
34:DA:981:U:H5'	47:DN:21:TYR:CZ	2.55	0.41
24:A0:5:LYS:HE3	24:A0:5:LYS:HB2	1.80	0.41
1:AA:1525:G:H2'	1:AA:1526:G:C8	2.55	0.41
1:AA:1700:G:H3'	15:AR:2:ARG:CD	2.51	0.41
1:AA:1739:U:H2'	1:AA:1741:C:C5	2.56	0.41
1:AA:2097:U:H2'	1:AA:2250:G:N2	2.36	0.41
1:AA:2845:A:C6	1:AA:2889:C:C6	3.09	0.41
1:AA:831:A:C8	1:AA:839:G:C5	3.09	0.41
3:AC:194:ILE:CD1	3:AC:227:PRO:CB	2.99	0.41
5:AE:111:ARG:H	5:AE:111:ARG:HG2	1.43	0.41
5:AE:2:LYS:HA	5:AE:84:PHE:CD1	2.56	0.41
8:AH:23:ARG:CZ	8:AH:23:ARG:HB2	2.50	0.41
12:AO:9:GLU:O	12:AO:83:ALA:HA	2.20	0.41
20:AW:83:LYS:C	20:AW:84:ARG:HD3	2.41	0.41
34:BA:1381:U:H1'	40:BG:79:ARG:HG3	2.02	0.41
34:BA:337:C:H2'	34:BA:338:A:C8	2.55	0.41
34:BA:370:C:O2'	34:BA:371:G:H5'	2.20	0.41
34:BA:40:C:C2	34:BA:41:G:C8	3.09	0.41
34:BA:557:G:N1	34:BA:558:G:C2	2.89	0.41
34:BA:685:G:N1	34:BA:686:U:O4	2.54	0.41
34:BA:665:A:N3	34:BA:732:C:H2'	2.35	0.41
34:BA:827:U:C4	34:BA:870:U:N3	2.89	0.41
34:BA:872:A:C2	34:BA:874:G:C5	3.08	0.41
34:BA:1060:C:C5	36:BC:2:GLY:HA3	2.55	0.41
40:BG:18:TYR:HD2	40:BG:59:LEU:HD22	1.86	0.41
49:BP:17:TYR:CE2	49:BP:41:PRO:HG3	2.55	0.41
46:BM:94:ARG:CZ	52:BS:80:TYR:HD2	2.34	0.41
57:BZ:443:HIS:CG	57:BZ:446:THR:HG22	2.55	0.41
1:CA:1149:G:H2'	1:CA:1150:C:C6	2.54	0.41
1:CA:1406:U:H2'	1:CA:1407:C:C6	2.56	0.41
1:CA:13:A:H2	1:CA:14:A:N6	2.19	0.41
1:CA:1507:A:O5'	1:CA:1507:A:H8	2.03	0.41
1:CA:1662:C:H2'	1:CA:1663:C:O4'	2.20	0.41
1:CA:2856:C:N3	1:CA:2862:G:C2	2.89	0.41
1:CA:285:C:O2'	1:CA:286:C:H5'	2.21	0.41
1:CA:470:A:OP1	6:CF:59:TYR:HE1	2.03	0.41
1:CA:7:G:H2'	1:CA:8:A:O4'	2.20	0.41
12:CO:4:PRO:O	12:CO:5:GLN:CB	2.68	0.41
15:CR:51:LEU:HD21	15:CR:69:ASP:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:CR:55:ALA:HA	15:CR:80:PHE:CE2	2.55	0.41
22:CY:42:VAL:HG12	22:CY:43:ASN:N	2.36	0.41
23:CZ:134:PRO:C	23:CZ:136:PHE:H	2.23	0.41
34:DA:1052:U:O2'	34:DA:1055:A:OP2	2.33	0.41
34:DA:1117:G:N2	34:DA:1180:A:O2'	2.54	0.41
34:DA:1418:A:H5''	34:DA:1419:G:OP2	2.20	0.41
34:DA:232:G:H2'	34:DA:233:C:O4'	2.21	0.41
34:DA:523:A:H61	45:DL:53:ARG:HH12	1.69	0.41
34:DA:678:U:H1'	34:DA:777:A:O3'	2.20	0.41
34:DA:785:G:N2	34:DA:798:G:C4	2.89	0.41
34:DA:857:C:H2'	34:DA:858:G:O4'	2.21	0.41
35:DB:71:VAL:HG12	35:DB:93:VAL:HG22	2.01	0.41
36:DC:16:ARG:HD2	36:DC:16:ARG:HA	1.81	0.41
36:DC:54:ARG:O	36:DC:69:HIS:ND1	2.39	0.41
45:DL:10:LEU:HA	45:DL:10:LEU:HD23	1.86	0.41
51:DR:44:LEU:HD21	51:DR:70:ILE:HD13	2.02	0.41
56:DY:75:C:HO2'	56:DY:76:A:H8	1.66	0.41
1:CA:2422:A:C6	56:DY:76:A:C2	3.09	0.41
57:DZ:159:ALA:O	57:DZ:161:PRO:HD3	2.21	0.41
57:DZ:237:PRO:HB3	57:DZ:241:GLU:OE1	2.21	0.41
57:DZ:31:ARG:O	57:DZ:34:TYR:HB3	2.20	0.41
57:DZ:669:PHE:H	57:DZ:669:PHE:HD1	1.69	0.41
57:DZ:88:VAL:O	57:DZ:91:THR:N	2.53	0.41
1:AA:1016:C:C2'	1:AA:1017:G:H5'	2.51	0.41
1:AA:1370:G:C5	1:AA:1374:G:O6	2.74	0.41
1:AA:1702:A:H4'	5:AE:115:GLY:N	2.36	0.41
1:AA:2294:G:OP1	1:AA:2295:C:H1'	2.21	0.41
1:AA:2342:G:H2'	1:AA:2343:G:O4'	2.21	0.41
1:AA:2418:U:H6	1:AA:2418:U:H2'	1.66	0.41
1:AA:2550:C:H2'	1:AA:2551:C:C6	2.55	0.41
1:AA:275:C:H2'	1:AA:276:C:C6	2.56	0.41
1:AA:2787:C:H2'	1:AA:2788:A:O4'	2.21	0.41
1:AA:585:U:C4	1:AA:2058:C:O4'	2.74	0.41
1:AA:721:G:O2'	6:AF:74:ARG:HD3	2.21	0.41
1:AA:1851:U:O2	4:AD:201:HIS:HB3	2.19	0.41
6:AF:53:THR:HB	6:AF:56:GLU:OE2	2.21	0.41
1:AA:1289:G:H4'	13:AP:7:ARG:NH2	2.36	0.41
16:AS:58:LEU:HA	16:AS:58:LEU:HD23	1.71	0.41
20:AW:51:LEU:HA	20:AW:51:LEU:HD23	1.86	0.41
21:AX:60:ARG:HH12	31:A7:47:ARG:NH2	2.19	0.41
22:AY:38:ILE:HD13	22:AY:66:PRO:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
23:AZ:105:VAL:O	23:AZ:141:VAL:HG22	2.21	0.41
34:BA:1277:C:H1'	34:BA:1282:C:O2	2.21	0.41
34:BA:186:C:C2	34:BA:187:C:C5	3.08	0.41
34:BA:872:A:C5	34:BA:874:G:C8	3.09	0.41
35:BB:150:SER:OG	35:BB:151:GLY:N	2.53	0.41
35:BB:197:VAL:HG12	35:BB:199:TYR:H	1.86	0.41
36:BC:124:ILE:HG22	36:BC:130:VAL:HG22	2.02	0.41
34:BA:620:C:C2	37:BD:135:LEU:HG	2.56	0.41
42:BI:92:TYR:HA	42:BI:92:TYR:HD1	1.77	0.41
50:BQ:62:SER:OG	50:BQ:72:ARG:HD2	2.20	0.41
52:BS:11:VAL:HG11	52:BS:16:LEU:HB2	2.02	0.41
56:BY:60:U:H5''	56:BY:61:C:C5	2.45	0.41
57:BZ:96:ARG:O	57:BZ:100:VAL:HG23	2.20	0.41
1:CA:1041:C:H5'	1:CA:1042:G:OP2	2.21	0.41
1:CA:1155:A:C4	1:CA:1157:G:C8	3.09	0.41
1:CA:1292:U:H2'	1:CA:1293:C:H6	1.83	0.41
1:CA:1368:G:O2'	1:CA:1369:G:H5'	2.21	0.41
1:CA:2092:U:H4'	1:CA:2093:G:O5'	2.21	0.41
1:CA:251:A:C5	1:CA:252:G:H1'	2.56	0.41
1:CA:264:C:O2'	1:CA:265:A:H2'	2.21	0.41
1:CA:2699:C:H2'	1:CA:2700:C:O4'	2.20	0.41
1:CA:501:A:C6	1:CA:502:A:C6	3.08	0.41
1:CA:705:A:H1'	4:CD:9:TYR:CE2	2.56	0.41
1:CA:818:G:H5'	1:CA:839:U:OP1	2.20	0.41
2:CB:28:C:H2'	2:CB:29:A:C8	2.56	0.41
4:CD:213:ARG:HD2	4:CD:213:ARG:HA	1.70	0.41
4:CD:37:LEU:HD12	4:CD:62:TYR:HB2	2.03	0.41
12:CO:103:ALA:O	12:CO:106:LEU:HB2	2.21	0.41
14:CQ:26:TYR:CE1	14:CQ:28:ALA:HB2	2.56	0.41
14:CQ:68:ILE:HD13	14:CQ:103:MET:HB3	2.03	0.41
20:CW:29:LEU:HD12	20:CW:29:LEU:O	2.21	0.41
23:CZ:150:LEU:HD12	23:CZ:171:ILE:HD11	2.02	0.41
23:CZ:69:THR:HG22	23:CZ:90:VAL:HA	2.01	0.41
34:DA:623:C:H2'	34:DA:624:C:H6	1.84	0.41
34:DA:687:A:O2'	34:DA:701:C:N4	2.53	0.41
34:DA:860:A:H2'	34:DA:861:G:O4'	2.20	0.41
34:DA:926:G:C6	34:DA:1505:G:C5	3.09	0.41
34:DA:953:G:H5'	34:DA:965:A:N6	2.35	0.41
36:DC:52:LEU:HD13	36:DC:118:GLN:HE22	1.86	0.41
43:DJ:46:ARG:HG3	43:DJ:64:GLU:HB3	2.02	0.41
44:DK:18:ARG:O	44:DK:32:ILE:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
52:DS:36:ARG:HG2	52:DS:51:VAL:HG13	2.02	0.41
52:DS:3:ARG:HG2	52:DS:4:SER:H	1.86	0.41
57:DZ:29:THR:O	57:DZ:33:LEU:HD23	2.20	0.41
57:DZ:419:ALA:O	57:DZ:422:GLU:HB3	2.21	0.41
57:DZ:468:ARG:HB3	57:DZ:468:ARG:HE	1.59	0.41
57:DZ:610:VAL:HG23	57:DZ:612:THR:HG22	2.03	0.41
57:DZ:647:VAL:HA	57:DZ:648:PRO:HD3	1.96	0.41
1:AA:2343:G:O3'	24:A0:43:THR:HG22	2.21	0.41
26:A2:9:GLN:OE1	26:A2:56:GLN:HG2	2.21	0.41
33:A9:10:ILE:HG21	33:A9:32:HIS:CD2	2.56	0.41
1:AA:1075:A:O5'	1:AA:1075:A:H8	2.04	0.41
1:AA:2303:U:OP1	1:AA:2393:C:H5'	2.21	0.41
1:AA:2385:G:H2'	1:AA:2386:C:C6	2.56	0.41
1:AA:280:C:H2'	1:AA:281:G:H8	1.84	0.41
1:AA:2845:A:C5	1:AA:2889:C:C5	3.09	0.41
1:AA:967:G:H2'	1:AA:968:U:C6	2.56	0.41
4:AD:146:GLU:HG2	4:AD:152:GLY:C	2.42	0.41
17:AT:39:ARG:NH2	34:BA:345:C:H5	2.19	0.41
34:BA:658:G:H2'	34:BA:659:U:H6	1.85	0.41
34:BA:828:A:H2'	34:BA:829:G:O5'	2.21	0.41
35:BB:221:LEU:HA	35:BB:221:LEU:HD22	1.84	0.41
36:BC:120:VAL:HB	36:BC:198:VAL:HG11	2.03	0.41
36:BC:129:ALA:HB3	36:BC:132:ARG:HB2	2.03	0.41
41:BH:116:LYS:HD2	41:BH:129:VAL:HG11	2.03	0.41
34:BA:599:C:H5''	41:BH:96:GLY:HA2	2.02	0.41
34:BA:966:G:H21	42:BI:127:LYS:NZ	2.19	0.41
42:BI:6:GLY:O	42:BI:17:VAL:HG12	2.21	0.41
43:BJ:55:LYS:HB3	43:BJ:55:LYS:HE2	1.89	0.41
47:BN:23:ARG:HD2	47:BN:30:ALA:HB2	2.03	0.41
48:BO:82:ILE:H	48:BO:82:ILE:HG13	1.75	0.41
57:BZ:183:MET:SD	57:BZ:213:HIS:CD2	3.14	0.41
26:C2:3:LEU:HD23	26:C2:3:LEU:HA	1.79	0.41
1:CA:108:U:H2'	1:CA:109:G:C8	2.54	0.41
1:CA:1321:A:H2'	1:CA:1322:A:O4'	2.20	0.41
1:CA:1379:A:H4'	1:CA:1380:G:OP2	2.21	0.41
1:CA:140:G:O4'	1:CA:141:A:H2	2.04	0.41
1:CA:1913:A:C6	57:DZ:580:MET:HE2	2.56	0.41
1:CA:2299:G:N2	1:CA:2318:G:N7	2.69	0.41
1:CA:2320:A:C2	1:CA:2333:A:C8	3.09	0.41
1:CA:239:U:H2'	1:CA:240:G:O4'	2.21	0.41
1:CA:2836:U:C4	1:CA:2883:A:N6	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:453:C:OP1	63:CA:4582:HOH:O	2.22	0.41
1:CA:589:C:H2'	1:CA:590:A:C8	2.55	0.41
2:CB:6:C:C2	2:CB:116:G:N2	2.89	0.41
5:CE:13:ARG:HD2	5:CE:20:ALA:HB1	2.03	0.41
5:CE:144:ARG:HB3	5:CE:145:LYS:H	1.56	0.41
8:CH:69:ARG:HG3	8:CH:70:THR:N	2.34	0.41
11:CN:62:VAL:HG13	11:CN:66:LYS:HB2	2.02	0.41
18:CU:79:PHE:CE1	18:CU:83:LEU:HD22	2.55	0.41
34:DA:401:C:OP2	37:DD:73:ARG:NE	2.53	0.41
34:DA:473:G:H8	34:DA:473:G:O5'	2.04	0.41
34:DA:518:C:O2'	34:DA:530:G:N2	2.54	0.41
34:DA:541:G:N2	34:DA:542:G:H1'	2.36	0.41
34:DA:995:C:H2'	34:DA:996:A:H8	1.86	0.41
35:DB:84:GLU:OE1	35:DB:87:ARG:NH1	2.54	0.41
40:DG:29:LYS:HD3	40:DG:29:LYS:HA	1.90	0.41
41:DH:113:SER:HB2	41:DH:134:ILE:HD11	2.03	0.41
45:DL:75:HIS:CD2	45:DL:77:LEU:HB2	2.56	0.41
46:DM:20:THR:C	46:DM:22:ILE:H	2.25	0.41
48:DO:21:ASP:OD2	48:DO:24:SER:HB3	2.20	0.41
49:DP:73:LEU:HD23	49:DP:73:LEU:HA	1.93	0.41
53:DT:31:SER:O	53:DT:34:LYS:HB2	2.21	0.41
56:DW:11:C:N4	56:DW:12:U:O4	2.54	0.41
56:DY:55:PSU:HN1	56:DY:57:G:C5'	2.33	0.41
57:DZ:349:LYS:HB2	57:DZ:349:LYS:HE3	1.90	0.41
57:DZ:74:TRP:HE1	57:DZ:273:LEU:C	2.24	0.41
24:A0:26:TYR:N	24:A0:29:GLN:OE1	2.40	0.41
24:A0:73:GLY:O	24:A0:74:ARG:C	2.60	0.41
1:AA:1046:A:C6	1:AA:1201:A:C8	3.09	0.41
1:AA:540:A:H2	1:AA:1306:G:N3	2.18	0.41
1:AA:1431:G:H4'	1:AA:1432:C:OP1	2.21	0.41
1:AA:153:C:OP2	25:A1:92:LYS:NZ	2.47	0.41
1:AA:2181:G:H2'	1:AA:2182:G:H8	1.86	0.41
1:AA:2556:G:H2'	1:AA:2557:G:O4'	2.21	0.41
1:AA:536:U:H5''	1:AA:537:G:OP2	2.21	0.41
1:AA:955:A:C6	1:AA:958:C:C2	3.09	0.41
2:AB:28:C:H2'	2:AB:29:A:O4'	2.21	0.41
4:AD:3:VAL:HG22	4:AD:18:VAL:O	2.21	0.41
6:AF:132:VAL:HG22	6:AF:132:VAL:O	2.21	0.41
8:AH:158:HIS:O	8:AH:160:LYS:N	2.54	0.41
10:AL:119:ASP:HB3	10:AL:122:ALA:HB3	2.02	0.41
10:AL:95:LYS:HG3	10:AL:135:GLY:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:AO:70:LYS:C	12:AO:71:ARG:HG3	2.41	0.41
15:AR:24:GLN:OE1	15:AR:36:THR:HG21	2.21	0.41
22:AY:92:ASN:N	22:AY:93:GLY:HA2	2.36	0.41
34:BA:1122:U:H2'	34:BA:1123:A:O4'	2.21	0.41
34:BA:314:C:O2'	34:BA:315:A:H5'	2.21	0.41
34:BA:731:G:H2'	34:BA:732:C:H6	1.86	0.41
34:BA:935:A:C2	34:BA:936:C:C2	3.09	0.41
34:BA:953:G:H2'	34:BA:954:G:O4'	2.21	0.41
34:BA:965:A:H5'	34:BA:969:A:O4'	2.20	0.41
39:BF:65:VAL:HG21	39:BF:67:MET:HE2	2.02	0.41
41:BH:17:THR:HB	41:BH:78:GLN:OE1	2.21	0.41
41:BH:41:ARG:HH22	41:BH:123:GLU:CD	2.24	0.41
46:BM:92:HIS:CE1	46:BM:98:VAL:HG21	2.56	0.41
47:BN:26:ARG:HB2	47:BN:39:LEU:HD22	2.03	0.41
49:BP:8:ARG:HG2	49:BP:9:PHE:N	2.36	0.41
56:BW:63:G:H2'	56:BW:64:A:O4'	2.21	0.41
57:BZ:172:ASP:OD2	57:BZ:173:THR:N	2.54	0.41
1:CA:2262:U:P	24:C0:19:LYS:HZ2	2.44	0.41
25:C1:53:VAL:HG22	25:C1:74:VAL:HG13	2.03	0.41
1:CA:1332:G:N3	1:CA:1332:G:H2'	2.36	0.41
1:CA:1394:U:H2'	1:CA:1395:A:O4'	2.20	0.41
1:CA:1526:G:C6	1:CA:1527:G:C2	3.09	0.41
1:CA:1802:A:N1	1:CA:1822:G:H1'	2.36	0.41
1:CA:1926:U:O2	1:CA:1928:A:C8	2.74	0.41
1:CA:2052:G:H21	5:CE:149:ARG:HA	1.85	0.41
1:CA:475:U:H1'	1:CA:509:C:C2	2.56	0.41
1:CA:539:G:C6	1:CA:540:C:C4	3.08	0.41
1:CA:606:U:O3'	1:CA:607:U:H4'	2.21	0.41
1:CA:780:G:C2	1:CA:782:A:C2	3.09	0.41
3:CC:11:LEU:CD1	3:CC:33:LEU:HA	2.50	0.41
3:CC:60:ARG:NH2	3:CC:165:ARG:HH21	2.18	0.41
6:CF:110:LEU:HD12	6:CF:110:LEU:HA	1.85	0.41
7:CG:121:ASN:HA	7:CG:122:PRO:HD3	1.75	0.41
7:CG:138:GLN:HB3	7:CG:153:ARG:O	2.21	0.41
8:CH:136:ILE:HG13	8:CH:136:ILE:H	1.43	0.41
13:CP:21:ARG:HD3	13:CP:21:ARG:HA	1.47	0.41
16:CS:30:ARG:HB2	16:CS:35:ILE:HD11	2.03	0.41
18:CU:63:VAL:HG13	63:CU:304:HOH:O	2.20	0.41
19:CV:25:LEU:HD23	19:CV:25:LEU:HA	1.67	0.41
20:CW:82:LEU:HD23	20:CW:82:LEU:HA	1.93	0.41
23:CZ:156:LYS:HB3	23:CZ:156:LYS:HE2	1.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:1401:G:H5''	34:DA:1402:C:OP2	2.21	0.41
17:CT:41:ARG:HH22	34:DA:346:G:P	2.44	0.41
34:DA:378:G:C6	34:DA:379:C:C4	3.08	0.41
34:DA:56:U:H2'	34:DA:57:G:H8	1.86	0.41
34:DA:942:G:H2'	34:DA:943:U:C6	2.55	0.41
46:DM:123:ALA:HB3	57:DZ:573:HIS:CG	2.56	0.41
57:DZ:-27:THR:O	57:DZ:-26:GLU:C	2.59	0.41
57:DZ:293:THR:HA	57:DZ:397:VAL:HG12	2.02	0.41
32:A8:23:VAL:HG13	32:A8:47:LYS:HB3	2.02	0.40
32:A8:31:HIS:O	32:A8:32:LEU:HB2	2.19	0.40
1:AA:2063:U:H2'	1:AA:2064:A:O5'	2.21	0.40
1:AA:451:G:N7	63:AA:4827:HOH:O	2.37	0.40
2:AB:37:C:C5	2:AB:38:C:C5	3.08	0.40
2:AB:49:C:C2'	2:AB:50:G:H5'	2.50	0.40
3:AC:20:VAL:O	3:AC:21:TYR:CB	2.58	0.40
7:AG:131:TYR:HB3	7:AG:159:VAL:CG1	2.51	0.40
14:AQ:21:THR:HG21	14:AQ:101:ARG:CB	2.50	0.40
19:AV:91:TYR:C	19:AV:91:TYR:CD1	2.94	0.40
34:BA:105:G:C6	34:BA:106:C:C4	3.09	0.40
34:BA:1311:G:N2	34:BA:1327:C:C2	2.89	0.40
34:BA:374:A:C6	34:BA:375:U:C4	3.09	0.40
34:BA:422:C:H2'	34:BA:422:C:OP2	2.21	0.40
34:BA:693:G:C6	34:BA:694:A:C6	3.10	0.40
34:BA:865:A:H2'	34:BA:866:C:O4'	2.21	0.40
34:BA:1112:C:N3	36:BC:178:LEU:HB2	2.35	0.40
39:BF:53:ALA:O	39:BF:54:LYS:HB2	2.21	0.40
40:BG:40:ALA:O	40:BG:43:PHE:HB3	2.21	0.40
42:BI:112:LYS:HD2	42:BI:118:LYS:O	2.21	0.40
48:BO:84:LYS:O	48:BO:84:LYS:HD3	2.20	0.40
49:BP:38:TYR:HB2	49:BP:39:TYR:H	1.71	0.40
50:BQ:22:LEU:HD12	50:BQ:23:VAL:N	2.36	0.40
57:BZ:-10:ARG:NH1	57:BZ:-10:ARG:HB2	2.35	0.40
57:BZ:236:GLU:HA	57:BZ:237:PRO:HD3	1.89	0.40
28:C4:15:ILE:HD13	28:C4:21:VAL:HG22	2.02	0.40
1:CA:1270:C:O2'	1:CA:1648:C:OP2	2.28	0.40
1:CA:2040:C:H2'	1:CA:2041:U:O4'	2.20	0.40
1:CA:2809:A:N1	1:CA:2892:A:C4	2.89	0.40
1:CA:443:A:N7	6:CF:45:ARG:HG2	2.36	0.40
1:CA:458:G:C8	31:C7:37:LYS:HG2	2.56	0.40
1:CA:783:A:C5	1:CA:785:G:H1'	2.56	0.40
1:CA:848:G:C2	1:CA:849:A:C5	3.09	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:CB:31:C:H2'	2:CB:32:C:H5'	2.03	0.40
5:CE:126:PRO:HB2	5:CE:131:ALA:HB2	2.03	0.40
17:CT:80:SER:HA	17:CT:81:PRO:HD2	1.83	0.40
21:CX:31:HIS:HA	21:CX:32:PRO:HD3	1.76	0.40
34:DA:1004:A:N6	34:DA:1037:C:C2	2.89	0.40
34:DA:1315:U:H2'	34:DA:1316:G:O4'	2.21	0.40
34:DA:1321:C:H5''	34:DA:1322:C:H2'	2.02	0.40
34:DA:168:G:O2'	34:DA:169:C:H5'	2.20	0.40
34:DA:187:C:O2'	53:DT:89:ARG:NH2	2.54	0.40
34:DA:416:G:C6	34:DA:417:C:N3	2.89	0.40
34:DA:373:A:C2	34:DA:482:A:C6	3.09	0.40
35:DB:30:ARG:HG3	35:DB:31:TYR:CD1	2.56	0.40
36:DC:12:LEU:HD23	36:DC:12:LEU:HA	1.97	0.40
37:DD:155:LEU:HD23	37:DD:156:GLU:H	1.86	0.40
41:DH:81:HIS:HB2	41:DH:138:TRP:OXT	2.20	0.40
42:DI:127:LYS:HE3	56:DW:34:G:OP2	2.21	0.40
43:DJ:5:ARG:HA	43:DJ:73:ASP:OD1	2.21	0.40
43:DJ:76:ASN:HA	43:DJ:77:PRO:HD2	1.90	0.40
50:DQ:34:LYS:O	50:DQ:36:ILE:HG23	2.21	0.40
56:DW:66:U:C4	56:DW:67:C:C4	3.09	0.40
57:DZ:-29:LEU:HB3	57:DZ:-27:THR:HG23	2.04	0.40
57:DZ:350:GLU:OE2	57:DZ:383:THR:OG1	2.24	0.40
57:DZ:356:LEU:HD23	57:DZ:358:MET:HE1	2.03	0.40
57:DZ:456:GLU:HB3	57:DZ:457:LEU:H	1.69	0.40
1:AA:1405:A:N3	1:AA:1405:A:O4'	2.53	0.40
1:AA:2152:U:H1'	1:AA:2180:A:N1	2.37	0.40
1:AA:2154:U:C6	3:AC:6:LYS:HB3	2.56	0.40
1:AA:2299:A:N3	1:AA:2301:G:C8	2.89	0.40
1:AA:2444:A:OP1	63:AA:4190:HOH:O	2.21	0.40
1:AA:2701:U:OP2	1:AA:2732:G:N2	2.46	0.40
1:AA:2145:G:O2'	3:AC:173:HIS:HB2	2.20	0.40
4:AD:218:ARG:HD3	4:AD:218:ARG:HH11	1.77	0.40
4:AD:44:ASN:OD1	4:AD:46:GLN:HB2	2.22	0.40
5:AE:85:ASN:HA	5:AE:86:PRO:HD2	1.88	0.40
6:AF:68:LYS:HE3	6:AF:68:LYS:HB3	1.87	0.40
1:AA:1104:G:N2	10:AL:126:MET:SD	2.94	0.40
12:AO:26:LYS:NZ	12:AO:37:ASP:OD2	2.49	0.40
14:AQ:114:ALA:O	14:AQ:118:LEU:HG	2.20	0.40
15:AR:46:GLY:HA2	15:AR:49:ASP:HB2	2.04	0.40
21:AX:35:THR:HG22	21:AX:38:GLU:HB3	2.03	0.40
21:AX:60:ARG:NH1	31:A7:47:ARG:NH2	2.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:BA:537:G:H2'	34:BA:538:G:H8	1.87	0.40
35:BB:25:ASN:HA	35:BB:26:PRO:HD3	1.87	0.40
36:BC:155:GLY:O	36:BC:157:ILE:N	2.54	0.40
38:BE:132:ALA:C	38:BE:134:ALA:N	2.75	0.40
38:BE:136:MET:O	38:BE:139:LEU:N	2.54	0.40
40:BG:78:ARG:HH21	40:BG:156:TRP:HB3	1.86	0.40
41:BH:114:THR:OG1	41:BH:117:GLY:O	2.32	0.40
43:BJ:30:SER:O	43:BJ:81:THR:HG21	2.21	0.40
45:BL:84:LEU:HD23	45:BL:105:TYR:CE2	2.57	0.40
47:BN:37:PHE:CE1	47:BN:53:LEU:HD13	2.56	0.40
50:BQ:76:LEU:HD11	50:BQ:78:GLU:O	2.21	0.40
57:BZ:-29:LEU:HB2	57:BZ:-27:THR:HG23	2.01	0.40
57:BZ:404:VAL:HG22	57:BZ:405:PRO:N	2.36	0.40
57:BZ:-7:GLU:C	57:BZ:-6:ARG:HH11	2.25	0.40
1:CA:1234:U:H2'	1:CA:1235:G:O4'	2.21	0.40
1:CA:2646:C:H2'	1:CA:2647:U:O4'	2.21	0.40
1:CA:56:A:H2'	1:CA:57:C:O4'	2.21	0.40
1:CA:607:U:OP1	6:CF:102:PRO:HA	2.22	0.40
1:CA:953:A:OP2	14:CQ:16:ARG:NE	2.54	0.40
1:CA:997:G:H2'	1:CA:998:C:H6	1.85	0.40
16:CS:34:HIS:HB3	16:CS:35:ILE:H	1.65	0.40
17:CT:105:LEU:HA	17:CT:105:LEU:HD23	1.66	0.40
22:CY:67:LEU:CD2	22:CY:71:LYS:HD3	2.51	0.40
34:DA:1323:G:H2'	34:DA:1324:A:C8	2.57	0.40
34:DA:641:U:H1'	34:DA:642:A:N7	2.36	0.40
34:DA:90:U:H2'	34:DA:91:C:H6	1.87	0.40
42:DI:65:VAL:HG21	42:DI:73:GLN:HB3	2.03	0.40
34:DA:1198:G:O2'	43:DJ:55:LYS:HE2	2.21	0.40
56:DY:28:G:H2'	56:DY:29:G:H8	1.85	0.40
57:DZ:637:ARG:HB3	57:DZ:638:GLY:H	1.71	0.40
24:A0:14:ARG:O	24:A0:15:ASP:HB2	2.21	0.40
24:A0:2:ALA:N	63:A0:201:HOH:O	2.55	0.40
28:A4:10:VAL:CG2	28:A4:29:PRO:HG3	2.48	0.40
1:AA:859:C:H1'	1:AA:1296:G:C2	2.56	0.40
1:AA:1485:A:C2	1:AA:1600:A:C5	3.09	0.40
1:AA:2418:U:C6	1:AA:2418:U:H5'	2.56	0.40
1:AA:2653:G:C8	1:AA:2653:G:H5''	2.56	0.40
1:AA:326:C:H2'	1:AA:327:U:C6	2.55	0.40
1:AA:509:A:C8	1:AA:510:C:C5	3.10	0.40
1:AA:592:U:C4	1:AA:593:G:C6	3.10	0.40
1:AA:605:G:H2'	1:AA:606:G:C8	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:874:U:H6	63:AA:4779:HOH:O	2.02	0.40
5:AE:21:VAL:O	5:AE:23:VAL:HG13	2.21	0.40
10:AL:53:VAL:HA	10:AL:54:PRO:HD3	1.75	0.40
13:AP:112:LEU:HD23	13:AP:112:LEU:HA	1.87	0.40
19:AV:94:LEU:HA	19:AV:94:LEU:HD23	1.82	0.40
21:AX:35:THR:HG22	21:AX:38:GLU:HB2	2.02	0.40
23:AZ:98:MET:CE	23:AZ:133:ILE:HD12	2.51	0.40
34:BA:613:C:H42	34:BA:627:G:H1	1.69	0.40
34:BA:938:A:C5	34:BA:939:G:C8	3.10	0.40
35:BB:111:ARG:O	35:BB:114:ARG:HB3	2.21	0.40
35:BB:121:LEU:HA	35:BB:121:LEU:HD13	1.90	0.40
37:BD:200:GLU:HG2	37:BD:200:GLU:H	1.73	0.40
43:BJ:75:ILE:O	43:BJ:77:PRO:HD3	2.21	0.40
45:BL:102:ARG:H	45:BL:102:ARG:HG2	1.65	0.40
46:BM:10:PRO:HG3	46:BM:21:TYR:CD1	2.57	0.40
48:BO:31:LEU:HD23	48:BO:31:LEU:HA	1.53	0.40
49:BP:40:ASP:OD2	49:BP:44:THR:OG1	2.25	0.40
53:BT:88:VAL:O	53:BT:92:LEU:HG	2.22	0.40
56:BY:49:C:H2'	56:BY:50:U:H6	1.86	0.40
57:BZ:239:GLU:O	57:BZ:242:LEU:N	2.48	0.40
57:BZ:-27:THR:O	57:BZ:-26:GLU:C	2.60	0.40
57:BZ:399:LEU:HD12	57:BZ:399:LEU:HA	1.80	0.40
57:BZ:573:HIS:HD2	57:BZ:576:ASP:H	1.63	0.40
1:CA:2884:U:C2	29:C5:52:TYR:CE1	3.08	0.40
1:CA:467:G:OP1	31:C7:33:ARG:NH1	2.55	0.40
1:CA:11:G:C2'	1:CA:12:U:H5''	2.45	0.40
1:CA:1721:G:H5''	1:CA:1721:G:N3	2.36	0.40
1:CA:2029:G:H2'	1:CA:2031:A:OP1	2.21	0.40
1:CA:2228:G:C5	1:CA:2229:C:C4	3.09	0.40
1:CA:2274:A:C5	1:CA:2276:G:C8	3.09	0.40
1:CA:2719:G:O2'	1:CA:2720:U:H5'	2.22	0.40
1:CA:986:C:H2'	1:CA:987:G:H5'	2.02	0.40
2:CB:54:G:C4	2:CB:55:U:C5	3.10	0.40
1:CA:2305:A:N3	7:CG:136:ARG:HA	2.37	0.40
11:CN:57:ALA:HB3	11:CN:124:ALA:HA	2.03	0.40
12:CO:25:LEU:O	12:CO:26:LYS:HB2	2.21	0.40
14:CQ:69:PHE:CD1	14:CQ:70:PRO:HD2	2.56	0.40
18:CU:65:ILE:CD1	18:CU:95:LEU:HB3	2.51	0.40
18:CU:80:ILE:HA	18:CU:80:ILE:HD13	1.94	0.40
23:CZ:152:ALA:HB1	23:CZ:163:LEU:HD21	2.03	0.40
34:DA:101:A:C5	34:DA:102:G:N7	2.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
34:DA:1248:A:H2'	34:DA:1249:C:C6	2.57	0.40
34:DA:460:G:C6	34:DA:470:C:H5''	2.57	0.40
35:DB:77:ALA:HB2	35:DB:211:ILE:HD13	2.03	0.40
37:DD:171:GLY:O	37:DD:173:TRP:N	2.54	0.40
37:DD:58:LEU:O	37:DD:61:LYS:HB3	2.21	0.40
38:DE:139:LEU:HA	38:DE:142:LEU:HD12	2.02	0.40
38:DE:33:VAL:HG22	38:DE:34:VAL:H	1.86	0.40
38:DE:80:ILE:HD12	38:DE:80:ILE:HA	1.93	0.40
46:DM:70:LEU:HA	46:DM:70:LEU:HD23	1.89	0.40
50:DQ:37:LYS:O	50:DQ:38:ARG:HD3	2.21	0.40
51:DR:73:ALA:HB3	51:DR:79:LEU:HD12	2.03	0.40
57:DZ:435:ASP:C	57:DZ:437:THR:H	2.25	0.40
25:A1:19:GLN:O	25:A1:35:THR:HG22	2.21	0.40
26:A2:48:HIS:O	26:A2:52:ASP:HB2	2.21	0.40
29:A5:16:ARG:NH1	29:A5:16:ARG:HG2	2.29	0.40
1:AA:1053:C:OP1	11:AN:35:ARG:NH1	2.53	0.40
1:AA:125:A:H5''	1:AA:126:C:C6	2.57	0.40
1:AA:2227:G:O2'	1:AA:2228:G:OP1	2.35	0.40
1:AA:348:A:H2'	1:AA:349:G:O5'	2.21	0.40
1:AA:504:A:C6	1:AA:506:A:C6	3.09	0.40
1:AA:593:G:H2'	1:AA:2052:A:N7	2.36	0.40
1:AA:804:U:H2'	1:AA:805:C:O4'	2.21	0.40
6:AF:11:VAL:HB	6:AF:18:ARG:HB3	2.03	0.40
1:AA:709:G:H5''	13:AP:16:ARG:HG3	2.03	0.40
18:AU:3:ARG:HD3	18:AU:3:ARG:HH11	1.61	0.40
23:AZ:110:GLY:CA	23:AZ:145:GLU:HA	2.51	0.40
34:BA:185:A:C2	34:BA:193:C:C2	3.09	0.40
34:BA:736:C:C2	34:BA:737:A:C8	3.10	0.40
34:BA:865:A:H2	34:BA:918:A:H4'	1.86	0.40
35:BB:155:LEU:HD11	35:BB:159:PRO:HG3	2.04	0.40
35:BB:15:VAL:O	35:BB:16:HIS:ND1	2.54	0.40
37:BD:121:VAL:HA	37:BD:126:ILE:HG13	2.02	0.40
37:BD:173:TRP:CD1	37:BD:173:TRP:N	2.89	0.40
38:BE:122:GLU:HB2	38:BE:126:ARG:HD3	2.03	0.40
42:BI:95:LYS:O	42:BI:96:LEU:HD23	2.21	0.40
47:BN:23:ARG:NH1	47:BN:30:ALA:HB2	2.36	0.40
48:BO:41:GLU:HA	48:BO:44:LYS:HD2	2.03	0.40
48:BO:5:LYS:H	48:BO:5:LYS:HD2	1.85	0.40
50:BQ:60:ILE:HG12	50:BQ:61:GLU:N	2.36	0.40
56:BW:37:MIA:H3'	56:BW:38:A:H8	1.86	0.40
56:BY:44:G:H8	56:BY:44:G:OP2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
56:BY:5:G:H1'	56:BY:69:G:N2	2.36	0.40
57:BZ:577:SER:OG	57:BZ:578:SER:N	2.55	0.40
27:C3:6:VAL:HG13	27:C3:56:VAL:HG13	2.03	0.40
32:C8:60:LEU:HA	32:C8:60:LEU:HD23	1.92	0.40
1:CA:1171:G:N2	1:CA:1179:C:C2	2.84	0.40
1:CA:1653:G:OP1	1:CA:2822:G:N1	2.52	0.40
1:CA:2079:U:H2'	1:CA:2080:G:O4'	2.22	0.40
1:CA:2233:U:H2'	1:CA:2234:G:C8	2.57	0.40
1:CA:2311:A:H1'	7:CG:88:ILE:HD12	2.03	0.40
1:CA:2521:C:H2'	1:CA:2522:U:O4'	2.21	0.40
1:CA:2815:C:H2'	1:CA:2816:C:H6	1.87	0.40
1:CA:303:U:H2'	1:CA:304:G:O4'	2.20	0.40
1:CA:701:G:N2	1:CA:732:C:C2	2.89	0.40
1:CA:886:C:H1'	1:CA:890:A:H61	1.85	0.40
2:CB:15:A:H1'	2:CB:110:G:N7	2.36	0.40
5:CE:116:VAL:HG13	5:CE:122:PHE:CB	2.51	0.40
5:CE:112:GLY:O	5:CE:159:HIS:HA	2.21	0.40
6:CF:170:LEU:HA	6:CF:170:LEU:HD12	1.87	0.40
1:CA:2744:G:N2	8:CH:143:GLN:OE1	2.53	0.40
11:CN:128:HIS:HA	11:CN:129:PRO:HD3	1.69	0.40
18:CU:20:LEU:HD23	18:CU:20:LEU:HA	1.81	0.40
34:DA:1030:C:H42	34:DA:1031:G:H1	1.68	0.40
34:DA:1060:C:C5'	43:DJ:51:ARG:HB3	2.50	0.40
34:DA:192:U:H4'	53:DT:57:ARG:HD3	2.04	0.40
34:DA:474:G:H2'	34:DA:475:G:C8	2.55	0.40
34:DA:491:G:C4	34:DA:492:G:C8	3.09	0.40
34:DA:502:G:N2	34:DA:544:G:N3	2.70	0.40
34:DA:543:C:C2	34:DA:544:G:C8	3.09	0.40
34:DA:543:C:O2'	34:DA:544:G:H5'	2.21	0.40
34:DA:692:U:O2'	34:DA:694:A:N7	2.47	0.40
34:DA:1240:U:OP2	40:DG:115:ARG:HA	2.21	0.40
40:DG:20:ASP:OD2	40:DG:23:VAL:HG23	2.22	0.40
45:DL:75:HIS:HD2	45:DL:77:LEU:N	2.18	0.40
50:DQ:13:ASP:OD1	50:DQ:13:ASP:N	2.54	0.40
50:DQ:11:VAL:HB	50:DQ:88:TYR:CE2	2.57	0.40
51:DR:29:PHE:CD2	51:DR:39:VAL:HG11	2.57	0.40
56:DW:19:G:H5''	56:DW:60:U:O4	2.21	0.40
56:DW:63:G:H2'	56:DW:64:A:H8	1.86	0.40
57:DZ:114:VAL:HB	57:DZ:152:THR:HB	2.04	0.40
57:DZ:174:PHE:HD2	57:DZ:267:LYS:HD3	1.86	0.40
57:DZ:25:LYS:HZ3	57:DZ:25:LYS:HG3	1.75	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
57:DZ:511:LYS:HB2	57:DZ:511:LYS:HE3	1.85	0.40
25:A1:7:ILE:HD12	25:A1:98:LEU:HD11	2.03	0.40
26:A2:63:VAL:O	26:A2:66:GLU:HB3	2.22	0.40
33:A9:3:VAL:HA	33:A9:35:ARG:O	2.21	0.40
1:AA:1137:G:N2	1:AA:1147:U:H1'	2.37	0.40
1:AA:1266:C:H2'	1:AA:1267:C:H6	1.87	0.40
1:AA:1882:U:H2'	1:AA:1883:C:O4'	2.21	0.40
1:AA:2690:C:H2'	1:AA:2691:A:O4'	2.21	0.40
1:AA:2865:C:O2'	1:AA:2866:C:H5'	2.21	0.40
1:AA:505:A:H4'	1:AA:506:A:OP1	2.22	0.40
1:AA:761:U:O2'	1:AA:763:A:N7	2.38	0.40
2:AB:78:A:C2	2:AB:100:A:C4	3.09	0.40
2:AB:33:G:N2	2:AB:50:G:C4	2.89	0.40
3:AC:54:ARG:CZ	3:AC:55:SER:O	2.69	0.40
3:AC:6:LYS:CA	3:AC:9:ARG:NH1	2.85	0.40
7:AG:133:LEU:HD12	7:AG:135:LEU:CD1	2.52	0.40
10:AL:117:THR:OG1	10:AL:118:THR:N	2.55	0.40
10:AL:38:VAL:O	10:AL:42:ASN:HB2	2.21	0.40
10:AL:78:ILE:HD11	10:AL:136:VAL:HG11	2.04	0.40
11:AN:46:VAL:CG2	11:AN:48:MET:HG2	2.51	0.40
19:AV:49:THR:O	19:AV:49:THR:HG22	2.20	0.40
1:AA:795:G:C8	20:AW:89:ALA:HB1	2.56	0.40
34:BA:1005:A:H5''	34:BA:1006:C:OP2	2.21	0.40
34:BA:1468:A:H8	34:BA:1468:A:O5'	2.04	0.40
34:BA:143:A:H2	34:BA:220:G:H1	1.69	0.40
34:BA:416:G:H2'	34:BA:417:C:O4'	2.22	0.40
34:BA:697:U:C5	34:BA:698:G:C8	3.10	0.40
34:BA:814:A:H2'	34:BA:816:A:C5'	2.49	0.40
35:BB:126:GLU:HB3	35:BB:127:ILE:H	1.65	0.40
35:BB:193:ASP:HB3	35:BB:196:LEU:HB2	2.02	0.40
36:BC:20:SER:O	47:BN:54:PRO:HB3	2.20	0.40
37:BD:70:ILE:HG13	37:BD:100:ARG:NH2	2.36	0.40
38:BE:127:ASN:HA	38:BE:128:PRO:HD3	1.87	0.40
38:BE:82:VAL:HB	38:BE:138:ALA:HB2	2.04	0.40
46:BM:79:LYS:NZ	46:BM:83:ASP:OD1	2.50	0.40
49:BP:39:TYR:CD2	49:BP:41:PRO:HD3	2.57	0.40
49:BP:8:ARG:HD3	49:BP:8:ARG:HH11	1.74	0.40
57:BZ:681:LYS:HE3	57:BZ:681:LYS:HB3	1.57	0.40
24:C0:48:GLY:N	24:C0:79:VAL:O	2.54	0.40
26:C2:9:GLN:OE1	26:C2:56:GLN:HG2	2.21	0.40
33:C9:9:ARG:NH1	33:C9:16:VAL:HG23	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CA:143(A):C:H4'	21:CX:38:GLU:OE2	2.22	0.40
1:CA:1568:G:H5'	4:CD:59:LYS:O	2.22	0.40
1:CA:2046:G:H2'	1:CA:2047:U:H6	1.86	0.40
1:CA:2120:G:H22	3:CC:169:THR:CG2	2.35	0.40
1:CA:2107:C:N4	1:CA:2182:G:H1	2.11	0.40
1:CA:2556:C:H2'	1:CA:2557:G:O4'	2.21	0.40
1:CA:2725:A:C4	1:CA:2727:G:C8	3.10	0.40
1:CA:2862:G:O2'	1:CA:2863:C:H5'	2.22	0.40
1:CA:89:G:H3'	1:CA:90:U:H5''	2.04	0.40
1:CA:927:G:H2'	1:CA:928:G:O4'	2.21	0.40
2:CB:16:G:H2'	2:CB:17:C:C6	2.57	0.40
4:CD:68:LYS:O	4:CD:69:ARG:HB2	2.21	0.40
5:CE:24:THR:HG21	5:CE:187:ALA:HA	2.03	0.40
8:CH:158:HIS:O	8:CH:160:LYS:N	2.55	0.40
13:CP:52:GLU:OE1	13:CP:55:ARG:NH1	2.47	0.40
15:CR:9:LYS:HA	15:CR:17:ARG:NE	2.37	0.40
16:CS:3:ARG:HB2	16:CS:3:ARG:CZ	2.51	0.40
23:CZ:99:TYR:HB3	23:CZ:123:ASP:OD1	2.22	0.40
23:CZ:78:LYS:HB3	23:CZ:78:LYS:NZ	2.36	0.40
34:DA:1077:G:N2	34:DA:1081:G:C4	2.90	0.40
34:DA:552:U:O3'	45:DL:87:GLY:HA3	2.21	0.40
34:DA:575:G:H4'	34:DA:576:G:O5'	2.21	0.40
34:DA:694:A:H2'	34:DA:695:A:O4'	2.21	0.40
38:DE:129:ILE:O	38:DE:132:ALA:HB3	2.21	0.40
42:DI:16:ARG:HB2	42:DI:64:THR:CG2	2.50	0.40
44:DK:38:ASN:HA	44:DK:39:PRO:HD2	1.88	0.40
46:DM:20:THR:HA	46:DM:25:ILE:O	2.22	0.40
46:DM:3:ARG:HG2	46:DM:8:GLU:OE1	2.21	0.40
34:DA:1401:G:OP2	55:DV:18:C:C5	2.75	0.40
56:DY:36:A:H2'	56:DY:37:MIA:O4'	2.22	0.40
57:DZ:126:GLU:OE2	57:DZ:132:ARG:NH2	2.54	0.40
57:DZ:455:GLY:O	57:DZ:458:HIS:HB3	2.21	0.40

There are no symmetry-related clashes.

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	
3	AC	133/228 (58%)	90 (68%)	25 (19%)	18 (14%)	0	1
3	CC	133/228 (58%)	90 (68%)	25 (19%)	18 (14%)	0	1
4	AD	273/276 (99%)	249 (91%)	20 (7%)	4 (2%)	12	37
4	CD	273/276 (99%)	242 (89%)	26 (10%)	5 (2%)	10	32
5	AE	202/206 (98%)	191 (95%)	9 (4%)	2 (1%)	18	50
5	CE	202/206 (98%)	174 (86%)	19 (9%)	9 (4%)	3	9
6	AF	201/210 (96%)	187 (93%)	11 (6%)	3 (2%)	12	37
6	CF	201/210 (96%)	186 (92%)	12 (6%)	3 (2%)	12	37
7	AG	179/182 (98%)	143 (80%)	25 (14%)	11 (6%)	2	4
7	CG	179/182 (98%)	148 (83%)	20 (11%)	11 (6%)	2	4
8	AH	172/180 (96%)	150 (87%)	20 (12%)	2 (1%)	15	44
8	CH	172/180 (96%)	148 (86%)	19 (11%)	5 (3%)	5	18
9	AK	128/173 (74%)	68 (53%)	33 (26%)	27 (21%)	0	0
9	CK	128/173 (74%)	69 (54%)	24 (19%)	35 (27%)	0	0
10	AL	137/147 (93%)	105 (77%)	23 (17%)	9 (7%)	1	4
10	CL	137/147 (93%)	95 (69%)	33 (24%)	9 (7%)	1	4
11	AN	138/140 (99%)	133 (96%)	4 (3%)	1 (1%)	25	59
11	CN	138/140 (99%)	125 (91%)	11 (8%)	2 (1%)	13	39
12	AO	120/122 (98%)	114 (95%)	4 (3%)	2 (2%)	11	34
12	CO	120/122 (98%)	105 (88%)	12 (10%)	3 (2%)	6	22
13	AP	147/150 (98%)	132 (90%)	13 (9%)	2 (1%)	13	39
13	CP	147/150 (98%)	128 (87%)	16 (11%)	3 (2%)	9	28
14	AQ	139/141 (99%)	124 (89%)	13 (9%)	2 (1%)	13	39
14	CQ	139/141 (99%)	121 (87%)	15 (11%)	3 (2%)	8	26
15	AR	116/118 (98%)	100 (86%)	13 (11%)	3 (3%)	6	21
15	CR	116/118 (98%)	92 (79%)	16 (14%)	8 (7%)	1	3
16	AS	108/112 (96%)	92 (85%)	12 (11%)	4 (4%)	4	13
16	CS	108/112 (96%)	86 (80%)	17 (16%)	5 (5%)	3	9
17	AT	129/146 (88%)	118 (92%)	10 (8%)	1 (1%)	22	55

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
17	CT	129/146 (88%)	109 (84%)	15 (12%)	5 (4%)	3	12
18	AU	114/118 (97%)	111 (97%)	3 (3%)	0	100	100
18	CU	114/118 (97%)	103 (90%)	10 (9%)	1 (1%)	20	52
19	AV	99/101 (98%)	92 (93%)	5 (5%)	2 (2%)	9	28
19	CV	99/101 (98%)	87 (88%)	9 (9%)	3 (3%)	5	17
20	AW	110/113 (97%)	106 (96%)	4 (4%)	0	100	100
20	CW	110/113 (97%)	100 (91%)	10 (9%)	0	100	100
21	AX	93/96 (97%)	88 (95%)	5 (5%)	0	100	100
21	CX	93/96 (97%)	80 (86%)	11 (12%)	2 (2%)	8	26
22	AY	105/110 (96%)	91 (87%)	11 (10%)	3 (3%)	5	18
22	CY	105/110 (96%)	90 (86%)	13 (12%)	2 (2%)	9	30
23	AZ	183/206 (89%)	146 (80%)	24 (13%)	13 (7%)	1	3
23	CZ	183/206 (89%)	140 (76%)	31 (17%)	12 (7%)	1	4
24	A0	81/85 (95%)	72 (89%)	8 (10%)	1 (1%)	15	44
24	C0	81/85 (95%)	70 (86%)	11 (14%)	0	100	100
25	A1	95/98 (97%)	86 (90%)	7 (7%)	2 (2%)	8	27
25	C1	95/98 (97%)	87 (92%)	4 (4%)	4 (4%)	3	10
26	A2	68/72 (94%)	62 (91%)	5 (7%)	1 (2%)	12	37
26	C2	68/72 (94%)	63 (93%)	4 (6%)	1 (2%)	12	37
27	A3	57/60 (95%)	55 (96%)	2 (4%)	0	100	100
27	C3	57/60 (95%)	52 (91%)	5 (9%)	0	100	100
28	A4	67/71 (94%)	41 (61%)	19 (28%)	7 (10%)	0	1
28	C4	67/71 (94%)	53 (79%)	11 (16%)	3 (4%)	3	9
29	A5	57/60 (95%)	53 (93%)	4 (7%)	0	100	100
29	C5	57/60 (95%)	51 (90%)	5 (9%)	1 (2%)	10	32
30	A6	51/54 (94%)	48 (94%)	2 (4%)	1 (2%)	9	28
30	C6	51/54 (94%)	46 (90%)	4 (8%)	1 (2%)	9	28
31	A7	46/49 (94%)	44 (96%)	2 (4%)	0	100	100
31	C7	46/49 (94%)	43 (94%)	1 (2%)	2 (4%)	3	10
32	A8	62/65 (95%)	60 (97%)	2 (3%)	0	100	100
32	C8	62/65 (95%)	59 (95%)	2 (3%)	1 (2%)	11	36

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
33	A9	35/37 (95%)	35 (100%)	0	0	100	100
33	C9	35/37 (95%)	33 (94%)	2 (6%)	0	100	100
35	BB	229/256 (90%)	176 (77%)	42 (18%)	11 (5%)	2	8
35	DB	229/256 (90%)	177 (77%)	37 (16%)	15 (7%)	1	4
36	BC	204/239 (85%)	169 (83%)	28 (14%)	7 (3%)	4	15
36	DC	204/239 (85%)	175 (86%)	26 (13%)	3 (2%)	12	37
37	BD	206/209 (99%)	154 (75%)	34 (16%)	18 (9%)	1	2
37	DD	206/209 (99%)	161 (78%)	31 (15%)	14 (7%)	1	3
38	BE	146/162 (90%)	111 (76%)	26 (18%)	9 (6%)	2	4
38	DE	146/162 (90%)	122 (84%)	21 (14%)	3 (2%)	8	27
39	BF	98/101 (97%)	80 (82%)	15 (15%)	3 (3%)	5	16
39	DF	98/101 (97%)	82 (84%)	15 (15%)	1 (1%)	18	50
40	BG	153/156 (98%)	131 (86%)	17 (11%)	5 (3%)	4	15
40	DG	153/156 (98%)	133 (87%)	14 (9%)	6 (4%)	3	12
41	BH	135/138 (98%)	111 (82%)	18 (13%)	6 (4%)	3	9
41	DH	135/138 (98%)	122 (90%)	13 (10%)	0	100	100
42	BI	125/128 (98%)	104 (83%)	16 (13%)	5 (4%)	3	11
42	DI	125/128 (98%)	104 (83%)	18 (14%)	3 (2%)	7	23
43	BJ	95/105 (90%)	80 (84%)	12 (13%)	3 (3%)	5	16
43	DJ	94/105 (90%)	76 (81%)	9 (10%)	9 (10%)	1	1
44	BK	112/129 (87%)	96 (86%)	12 (11%)	4 (4%)	4	13
44	DK	112/129 (87%)	98 (88%)	11 (10%)	3 (3%)	6	20
45	BL	120/132 (91%)	111 (92%)	7 (6%)	2 (2%)	11	34
45	DL	120/132 (91%)	103 (86%)	13 (11%)	4 (3%)	4	15
46	BM	115/126 (91%)	89 (77%)	23 (20%)	3 (3%)	6	21
46	DM	120/126 (95%)	100 (83%)	12 (10%)	8 (7%)	1	4
47	BN	58/61 (95%)	47 (81%)	9 (16%)	2 (3%)	4	15
47	DN	58/61 (95%)	53 (91%)	5 (9%)	0	100	100
48	BO	86/89 (97%)	72 (84%)	8 (9%)	6 (7%)	1	3
48	DO	86/89 (97%)	75 (87%)	8 (9%)	3 (4%)	4	14
49	BP	80/88 (91%)	52 (65%)	19 (24%)	9 (11%)	0	1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
49	DP	80/88 (91%)	65 (81%)	12 (15%)	3 (4%)	4	12
50	BQ	97/105 (92%)	82 (84%)	9 (9%)	6 (6%)	2	4
50	DQ	97/105 (92%)	87 (90%)	7 (7%)	3 (3%)	5	16
51	BR	66/88 (75%)	60 (91%)	6 (9%)	0	100	100
51	DR	66/88 (75%)	57 (86%)	7 (11%)	2 (3%)	5	17
52	BS	82/93 (88%)	66 (80%)	14 (17%)	2 (2%)	7	23
52	DS	81/93 (87%)	68 (84%)	8 (10%)	5 (6%)	2	4
53	BT	94/106 (89%)	77 (82%)	10 (11%)	7 (7%)	1	3
53	DT	94/106 (89%)	81 (86%)	10 (11%)	3 (3%)	5	16
54	BU	21/27 (78%)	20 (95%)	1 (5%)	0	100	100
54	DU	21/27 (78%)	19 (90%)	1 (5%)	1 (5%)	2	8
57	BZ	726/758 (96%)	569 (78%)	106 (15%)	51 (7%)	1	3
57	DZ	726/758 (96%)	554 (76%)	121 (17%)	51 (7%)	1	3
All	All	13389/14444 (93%)	11230 (84%)	1582 (12%)	577 (4%)	3	10

All (577) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	AC	42	VAL
3	AC	47	LYS
3	AC	68	GLY
3	AC	180	SER
3	AC	181	PHE
4	AD	275	LYS
7	AG	43	LEU
7	AG	47	LYS
7	AG	51	ARG
7	AG	181	ARG
9	AK	29	TYR
9	AK	47	ASN
9	AK	71	LEU
9	AK	74	LEU
9	AK	77	PRO
9	AK	80	VAL
9	AK	91	LYS
9	AK	104	ILE
9	AK	105	PRO

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Mol	Chain	Res	Type
9	AK	107	VAL
9	AK	125	LEU
9	AK	128	LEU
10	AL	16	LYS
16	AS	59	LYS
23	AZ	136	PHE
23	AZ	154	ASP
23	AZ	158	PRO
23	AZ	177	PRO
23	AZ	178	GLU
23	AZ	183	LEU
23	AZ	184	ALA
28	A4	4	GLY
28	A4	34	GLU
28	A4	59	PHE
28	A4	62	ARG
30	A6	29	ASN
35	BB	10	LEU
35	BB	13	ALA
35	BB	17	PHE
35	BB	125	PRO
36	BC	65	ALA
37	BD	5	ILE
37	BD	31	CYS
37	BD	42	GLN
37	BD	101	LEU
37	BD	102	ASP
37	BD	178	VAL
37	BD	179	GLU
37	BD	207	TYR
38	BE	98	THR
38	BE	140	ARG
39	BF	42	GLU
40	BG	79	ARG
40	BG	80	VAL
42	BI	41	VAL
42	BI	54	ASP
43	BJ	31	GLY
44	BK	106	LYS
47	BN	4	LYS
48	BO	21	ASP
49	BP	66	PRO

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Mol	Chain	Res	Type
53	BT	10	LEU
53	BT	100	ILE
57	BZ	-57	GLU
57	BZ	-25	SER
57	BZ	-23	LEU
57	BZ	39	ILE
57	BZ	87	HIS
57	BZ	88	VAL
57	BZ	100	VAL
57	BZ	183	MET
57	BZ	239	GLU
57	BZ	240	GLU
57	BZ	324	ARG
57	BZ	402	ILE
57	BZ	404	VAL
57	BZ	446	THR
57	BZ	469	GLU
57	BZ	504	ARG
3	CC	42	VAL
3	CC	47	LYS
3	CC	68	GLY
3	CC	180	SER
3	CC	181	PHE
4	CD	239	ARG
5	CE	17	ASP
5	CE	74	PRO
6	CF	130	ALA
7	CG	14	GLU
7	CG	47	LYS
7	CG	126	ASP
7	CG	181	ARG
8	CH	126	PRO
9	CK	68	LEU
9	CK	69	PRO
9	CK	71	LEU
9	CK	74	LEU
9	CK	75	GLN
9	CK	77	PRO
9	CK	80	VAL
9	CK	85	ASP
9	CK	90	ALA
9	CK	93	LEU

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Mol	Chain	Res	Type
9	CK	107	VAL
9	CK	128	LEU
10	CL	115	LEU
12	CO	26	LYS
14	CQ	28	ALA
15	CR	14	SER
15	CR	58	GLY
15	CR	86	ARG
16	CS	57	LYS
17	CT	100	TYR
19	CV	79	VAL
23	CZ	154	ASP
23	CZ	158	PRO
23	CZ	161	VAL
23	CZ	183	LEU
23	CZ	184	ALA
25	C1	3	LYS
25	C1	26	ARG
25	C1	85	LEU
30	C6	44	ARG
35	DB	10	LEU
35	DB	17	PHE
35	DB	21	ARG
35	DB	24	TRP
35	DB	74	LYS
37	DD	136	PRO
37	DD	154	ASN
38	DE	140	ARG
39	DF	79	LEU
40	DG	80	VAL
42	DI	54	ASP
42	DI	107	ARG
43	DJ	56	HIS
43	DJ	75	ILE
46	DM	106	ASN
48	DO	19	PRO
50	DQ	74	LEU
51	DR	52	PRO
53	DT	100	ILE
57	DZ	-57	GLU
57	DZ	-25	SER
57	DZ	-12	ALA

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Mol	Chain	Res	Type
57	DZ	89	ASP
57	DZ	92	ILE
57	DZ	160	ARG
57	DZ	183	MET
57	DZ	290	LYS
57	DZ	303	PRO
57	DZ	402	ILE
57	DZ	416	LYS
57	DZ	472	VAL
57	DZ	528	ALA
3	AC	53	ARG
3	AC	161	ARG
3	AC	179	ALA
6	AF	130	ALA
7	AG	21	ARG
7	AG	78	SER
7	AG	140	ILE
9	AK	49	ALA
9	AK	75	GLN
9	AK	119	ALA
9	AK	123	GLU
9	AK	132	ASP
10	AL	89	HIS
15	AR	42	LYS
17	AT	129	ARG
19	AV	79	VAL
22	AY	78	ALA
24	A0	74	ARG
35	BB	16	HIS
35	BB	37	ASN
35	BB	106	LYS
36	BC	66	VAL
36	BC	70	VAL
37	BD	172	PRO
38	BE	68	GLU
38	BE	85	GLY
38	BE	103	GLY
38	BE	133	TYR
38	BE	146	ALA
41	BH	41	ARG
41	BH	51	VAL
41	BH	104	ARG

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Mol	Chain	Res	Type
41	BH	133	LEU
42	BI	43	ALA
46	BM	101	GLN
47	BN	52	GLN
49	BP	78	GLY
50	BQ	34	LYS
50	BQ	49	GLU
50	BQ	87	LYS
53	BT	47	GLY
53	BT	96	GLY
57	BZ	-33	GLY
57	BZ	-8	ALA
57	BZ	-4	ALA
57	BZ	86	GLY
57	BZ	171	GLU
57	BZ	172	ASP
57	BZ	199	ILE
57	BZ	235	GLU
57	BZ	400	GLU
57	BZ	416	LYS
57	BZ	418	LYS
57	BZ	468	ARG
57	BZ	472	VAL
57	BZ	656	ALA
57	BZ	671	MET
3	CC	53	ARG
3	CC	161	ARG
3	CC	179	ALA
5	CE	69	LYS
6	CF	21	ALA
7	CG	51	ARG
7	CG	81	LYS
9	CK	33	PRO
9	CK	50	ARG
9	CK	53	VAL
9	CK	70	GLU
9	CK	73	GLY
9	CK	100	ASN
9	CK	104	ILE
9	CK	110	GLY
10	CL	13	PRO
10	CL	89	HIS

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Mol	Chain	Res	Type
11	CN	2	LYS
13	CP	140	ALA
14	CQ	3	MET
14	CQ	55	VAL
15	CR	4	LEU
15	CR	71	GLN
16	CS	82	ILE
17	CT	10	VAL
21	CX	94	GLY
23	CZ	60	GLU
23	CZ	64	GLY
23	CZ	182	LYS
28	C4	45	GLY
28	C4	46	GLN
29	C5	37	LYS
31	C7	46	VAL
35	DB	26	PRO
36	DC	26	LYS
37	DD	3	ARG
37	DD	56	VAL
37	DD	76	ARG
37	DD	110	PHE
37	DD	171	GLY
37	DD	182	LYS
38	DE	150	ARG
40	DG	79	ARG
40	DG	148	ASN
43	DJ	29	ARG
43	DJ	55	LYS
43	DJ	77	PRO
43	DJ	79	ARG
44	DK	49	GLY
46	DM	100	GLY
48	DO	88	ARG
49	DP	78	GLY
50	DQ	33	GLY
52	DS	25	LYS
57	DZ	-1	GLU
57	DZ	39	ILE
57	DZ	85	PRO
57	DZ	182	ARG
57	DZ	403	GLU

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Mol	Chain	Res	Type
57	DZ	446	THR
57	DZ	456	GLU
57	DZ	468	ARG
57	DZ	533	VAL
3	AC	30	VAL
3	AC	43	GLU
3	AC	52	PRO
3	AC	69	LEU
3	AC	184	GLU
3	AC	202	PRO
3	AC	209	PHE
4	AD	156	ALA
5	AE	52	LEU
6	AF	25	PRO
7	AG	74	LYS
7	AG	79	ASN
9	AK	30	GLN
9	AK	33	PRO
9	AK	93	LEU
9	AK	112	LEU
12	AO	5	GLN
19	AV	31	ALA
22	AY	54	LYS
23	AZ	159	PRO
23	AZ	161	VAL
25	A1	3	LYS
35	BB	19	HIS
37	BD	3	ARG
37	BD	109	GLY
37	BD	142	PRO
38	BE	132	ALA
41	BH	6	ILE
44	BK	49	GLY
44	BK	96	ARG
44	BK	100	ALA
45	BL	50	SER
49	BP	56	ALA
50	BQ	82	MET
50	BQ	94	ASN
57	BZ	85	PRO
57	BZ	170	ARG
57	BZ	274	ASP

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Mol	Chain	Res	Type
57	BZ	486	THR
57	BZ	521	SER
57	BZ	638	GLY
3	CC	30	VAL
3	CC	43	GLU
3	CC	52	PRO
3	CC	69	LEU
3	CC	184	GLU
3	CC	202	PRO
3	CC	209	PHE
4	CD	3	VAL
5	CE	103	ASP
8	CH	159	GLU
8	CH	169	VAL
9	CK	20	ALA
9	CK	30	GLN
9	CK	56	ASN
9	CK	62	ALA
9	CK	63	LEU
9	CK	114	GLY
9	CK	119	ALA
9	CK	132	ASP
18	CU	72	HIS
19	CV	29	PRO
22	CY	43	ASN
23	CZ	178	GLU
26	C2	22	GLU
28	C4	68	ARG
31	C7	45	ALA
35	DB	20	GLU
35	DB	77	ALA
35	DB	134	GLU
35	DB	232	PRO
37	DD	47	ARG
40	DG	33	ASP
40	DG	55	GLY
44	DK	106	LYS
45	DL	106	ASP
45	DL	125	PRO
46	DM	67	GLU
46	DM	118	ALA
52	DS	30	LEU

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Mol	Chain	Res	Type
52	DS	42	PRO
53	DT	102	GLY
54	DU	7	ARG
57	DZ	-65	LYS
57	DZ	-8	ALA
57	DZ	9	LEU
57	DZ	115	GLU
57	DZ	153	MET
57	DZ	235	GLU
57	DZ	404	VAL
57	DZ	444	PRO
57	DZ	457	LEU
57	DZ	636	PRO
57	DZ	671	MET
3	AC	16	ASP
4	AD	169	GLU
8	AH	159	GLU
9	AK	5	ARG
9	AK	110	GLY
10	AL	33	ASN
10	AL	42	ASN
13	AP	29	LYS
13	AP	122	PRO
15	AR	45	ARG
23	AZ	137	ILE
23	AZ	157	LEU
23	AZ	164	ALA
28	A4	41	PRO
36	BC	61	ALA
36	BC	81	GLY
36	BC	88	ARG
37	BD	105	VAL
37	BD	182	LYS
38	BE	77	PRO
39	BF	70	ASP
41	BH	7	ALA
42	BI	97	LYS
46	BM	12	ASN
48	BO	44	LYS
48	BO	78	TYR
49	BP	18	ARG
49	BP	49	LEU

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Mol	Chain	Res	Type
49	BP	64	ALA
52	BS	81	ARG
53	BT	102	GLY
57	BZ	-26	GLU
57	BZ	22	ASP
57	BZ	181	LEU
57	BZ	269	VAL
57	BZ	473	ASP
57	BZ	500	GLN
57	BZ	640	ALA
3	CC	16	ASP
5	CE	162	ALA
7	CG	32	PRO
7	CG	43	LEU
7	CG	50	ALA
8	CH	65	HIS
9	CK	47	ASN
9	CK	101	PRO
10	CL	72	PRO
12	CO	5	GLN
13	CP	29	LYS
13	CP	122	PRO
16	CS	84	GLN
17	CT	82	LEU
35	DB	123	ALA
35	DB	125	PRO
35	DB	131	PRO
36	DC	79	ARG
36	DC	156	ARG
37	DD	101	LEU
37	DD	172	PRO
38	DE	69	VAL
42	DI	118	LYS
43	DJ	78	ASN
45	DL	120	TYR
46	DM	5	ALA
46	DM	6	GLY
50	DQ	81	ARG
51	DR	41	LYS
52	DS	12	ASP
52	DS	29	ARG
57	DZ	-24	ASN

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Mol	Chain	Res	Type
57	DZ	88	VAL
57	DZ	170	ARG
57	DZ	199	ILE
57	DZ	247	ARG
57	DZ	640	ALA
57	DZ	656	ALA
3	AC	21	TYR
4	AD	146	GLU
5	AE	57	LYS
6	AF	207	GLY
7	AG	113	ARG
8	AH	127	GLU
9	AK	85	ASP
10	AL	13	PRO
10	AL	51	ALA
11	AN	14	VAL
12	AO	25	LEU
14	AQ	13	GLN
14	AQ	40	ALA
16	AS	82	ILE
26	A2	68	ARG
28	A4	49	PHE
35	BB	231	GLU
37	BD	199	ASN
40	BG	6	ARG
40	BG	93	PRO
42	BI	90	PRO
45	BL	88	GLY
49	BP	39	TYR
52	BS	67	VAL
53	BT	52	ALA
57	BZ	243	VAL
57	BZ	353	ALA
57	BZ	444	PRO
3	CC	21	TYR
4	CD	125	ILE
4	CD	275	LYS
5	CE	182	LEU
9	CK	86	PRO
9	CK	120	LYS
10	CL	3	LYS
10	CL	50	ASP

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Mol	Chain	Res	Type
10	CL	119	ASP
11	CN	87	LEU
12	CO	35	VAL
15	CR	45	ARG
15	CR	49	ASP
15	CR	107	ASP
16	CS	96	GLY
19	CV	43	GLU
21	CX	45	THR
22	CY	66	PRO
23	CZ	155	LEU
23	CZ	157	LEU
23	CZ	160	GLY
25	C1	45	ASN
35	DB	124	SER
37	DD	5	ILE
37	DD	59	ARG
43	DJ	36	GLY
46	DM	68	GLY
48	DO	47	LYS
57	DZ	380	LEU
57	DZ	532	GLY
3	AC	221	PRO
16	AS	72	ALA
23	AZ	156	LYS
25	A1	55	GLY
35	BB	78	GLN
35	BB	126	GLU
36	BC	101	LEU
37	BD	104	VAL
43	BJ	79	ARG
46	BM	67	GLU
48	BO	79	ARG
48	BO	86	GLY
53	BT	63	ILE
57	BZ	209	ALA
57	BZ	281	PRO
3	CC	221	PRO
5	CE	52	LEU
6	CF	132	VAL
7	CG	52	ILE
8	CH	168	PRO

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Mol	Chain	Res	Type
9	CK	84	GLU
16	CS	22	GLY
17	CT	20	PRO
17	CT	106	SER
32	C8	17	THR
35	DB	227	GLY
57	DZ	154	GLN
57	DZ	575	VAL
57	DZ	637	ARG
9	AK	22	GLY
9	AK	114	GLY
16	AS	60	GLY
28	A4	45	GLY
40	BG	55	GLY
5	CE	102	VAL
53	DT	63	ILE
57	DZ	502	GLY
7	AG	32	PRO
9	AK	86	PRO
10	AL	24	GLY
10	AL	113	PRO
57	BZ	307	GLY
9	CK	105	PRO
10	CL	21	PRO
37	DD	37	PRO
43	DJ	24	VAL
49	DP	51	VAL
10	AL	21	PRO
37	BD	37	PRO
39	BF	6	VAL
48	BO	36	ILE
49	BP	19	ILE
57	BZ	-16	ILE
4	CD	236	GLY
7	CG	63	ILE
9	CK	129	PRO
10	CL	24	GLY
45	DL	88	GLY
57	DZ	237	PRO
22	AY	103	GLY
37	BD	136	PRO
43	BJ	77	PRO

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Mol	Chain	Res	Type
49	BP	41	PRO
50	BQ	33	GLY
5	CE	71	GLY
40	DG	19	GLY
44	DK	39	PRO
46	DM	4	ILE
49	DP	53	VAL
57	DZ	86	GLY
57	DZ	163	VAL
57	DZ	638	GLY
15	AR	92	GLY
57	DZ	384	ILE

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	
3	AC	111/180 (62%)	104 (94%)	7 (6%)	21	51
3	CC	111/180 (62%)	103 (93%)	8 (7%)	17	43
4	AD	215/218 (99%)	181 (84%)	34 (16%)	3	9
4	CD	216/218 (99%)	179 (83%)	37 (17%)	2	7
5	AE	164/166 (99%)	130 (79%)	34 (21%)	1	4
5	CE	164/166 (99%)	130 (79%)	34 (21%)	1	4
6	AF	160/166 (96%)	128 (80%)	32 (20%)	1	4
6	CF	159/166 (96%)	124 (78%)	35 (22%)	1	3
7	AG	143/156 (92%)	114 (80%)	29 (20%)	1	4
7	CG	142/156 (91%)	111 (78%)	31 (22%)	1	3
8	AH	144/148 (97%)	121 (84%)	23 (16%)	3	8
8	CH	144/148 (97%)	124 (86%)	20 (14%)	4	12
10	AL	104/111 (94%)	83 (80%)	21 (20%)	1	4
10	CL	104/111 (94%)	84 (81%)	20 (19%)	1	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
11	AN	118/119 (99%)	97 (82%)	21 (18%)	2	6
11	CN	118/119 (99%)	92 (78%)	26 (22%)	1	3
12	AO	100/100 (100%)	88 (88%)	12 (12%)	6	18
12	CO	100/100 (100%)	82 (82%)	18 (18%)	2	6
13	AP	116/116 (100%)	90 (78%)	26 (22%)	1	3
13	CP	115/116 (99%)	95 (83%)	20 (17%)	2	7
14	AQ	111/111 (100%)	99 (89%)	12 (11%)	7	22
14	CQ	111/111 (100%)	91 (82%)	20 (18%)	2	6
15	AR	101/101 (100%)	81 (80%)	20 (20%)	1	4
15	CR	101/101 (100%)	78 (77%)	23 (23%)	1	3
16	AS	87/88 (99%)	70 (80%)	17 (20%)	1	4
16	CS	85/88 (97%)	66 (78%)	19 (22%)	1	3
17	AT	115/127 (91%)	98 (85%)	17 (15%)	3	10
17	CT	113/127 (89%)	90 (80%)	23 (20%)	1	4
18	AU	93/94 (99%)	79 (85%)	14 (15%)	3	10
18	CU	93/94 (99%)	80 (86%)	13 (14%)	4	12
19	AV	80/82 (98%)	63 (79%)	17 (21%)	1	3
19	CV	80/82 (98%)	68 (85%)	12 (15%)	3	10
20	AW	90/92 (98%)	77 (86%)	13 (14%)	4	11
20	CW	90/92 (98%)	82 (91%)	8 (9%)	11	32
21	AX	77/78 (99%)	69 (90%)	8 (10%)	8	24
21	CX	77/78 (99%)	69 (90%)	8 (10%)	8	24
22	AY	85/91 (93%)	71 (84%)	14 (16%)	2	8
22	CY	85/91 (93%)	69 (81%)	16 (19%)	2	5
23	AZ	156/179 (87%)	121 (78%)	35 (22%)	1	3
23	CZ	156/179 (87%)	129 (83%)	27 (17%)	2	7
24	A0	65/67 (97%)	61 (94%)	4 (6%)	21	52
24	C0	65/67 (97%)	60 (92%)	5 (8%)	15	39
25	A1	80/83 (96%)	70 (88%)	10 (12%)	5	16
25	C1	80/83 (96%)	66 (82%)	14 (18%)	2	6
26	A2	65/67 (97%)	55 (85%)	10 (15%)	3	9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
26	C2	65/67 (97%)	51 (78%)	14 (22%)	1	3
27	A3	51/52 (98%)	41 (80%)	10 (20%)	1	4
27	C3	50/52 (96%)	42 (84%)	8 (16%)	3	8
28	A4	60/63 (95%)	42 (70%)	18 (30%)	0	1
28	C4	53/63 (84%)	41 (77%)	12 (23%)	1	3
29	A5	50/52 (96%)	43 (86%)	7 (14%)	4	12
29	C5	50/52 (96%)	42 (84%)	8 (16%)	3	8
30	A6	51/52 (98%)	39 (76%)	12 (24%)	1	2
30	C6	50/52 (96%)	39 (78%)	11 (22%)	1	3
31	A7	41/42 (98%)	34 (83%)	7 (17%)	2	7
31	C7	41/42 (98%)	32 (78%)	9 (22%)	1	3
32	A8	54/55 (98%)	46 (85%)	8 (15%)	3	10
32	C8	54/55 (98%)	44 (82%)	10 (18%)	2	5
33	A9	34/34 (100%)	32 (94%)	2 (6%)	23	54
33	C9	34/34 (100%)	32 (94%)	2 (6%)	23	54
35	BB	192/220 (87%)	153 (80%)	39 (20%)	1	4
35	DB	187/220 (85%)	151 (81%)	36 (19%)	1	5
36	BC	143/188 (76%)	129 (90%)	14 (10%)	9	27
36	DC	141/188 (75%)	116 (82%)	25 (18%)	2	6
37	BD	170/181 (94%)	143 (84%)	27 (16%)	3	9
37	DD	174/181 (96%)	137 (79%)	37 (21%)	1	3
38	BE	113/123 (92%)	89 (79%)	24 (21%)	1	3
38	DE	114/123 (93%)	92 (81%)	22 (19%)	1	5
39	BF	84/90 (93%)	68 (81%)	16 (19%)	2	5
39	DF	86/90 (96%)	75 (87%)	11 (13%)	5	15
40	BG	119/127 (94%)	102 (86%)	17 (14%)	4	11
40	DG	120/127 (94%)	99 (82%)	21 (18%)	2	6
41	BH	114/119 (96%)	96 (84%)	18 (16%)	3	9
41	DH	114/119 (96%)	92 (81%)	22 (19%)	1	5
42	BI	91/99 (92%)	75 (82%)	16 (18%)	2	6
42	DI	89/99 (90%)	76 (85%)	13 (15%)	3	11

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
43	BJ	66/92 (72%)	61 (92%)	5 (8%)	15	40
43	DJ	69/92 (75%)	65 (94%)	4 (6%)	23	55
44	BK	83/99 (84%)	68 (82%)	15 (18%)	2	6
44	DK	83/99 (84%)	77 (93%)	6 (7%)	17	43
45	BL	97/109 (89%)	83 (86%)	14 (14%)	4	11
45	DL	97/109 (89%)	82 (84%)	15 (16%)	3	9
46	BM	91/101 (90%)	70 (77%)	21 (23%)	1	2
46	DM	92/101 (91%)	79 (86%)	13 (14%)	4	12
47	BN	49/50 (98%)	40 (82%)	9 (18%)	2	5
47	DN	49/50 (98%)	40 (82%)	9 (18%)	2	5
48	BO	78/80 (98%)	69 (88%)	9 (12%)	6	20
48	DO	78/80 (98%)	68 (87%)	10 (13%)	5	15
49	BP	69/74 (93%)	53 (77%)	16 (23%)	1	2
49	DP	68/74 (92%)	56 (82%)	12 (18%)	2	6
50	BQ	94/97 (97%)	74 (79%)	20 (21%)	1	3
50	DQ	94/97 (97%)	85 (90%)	9 (10%)	10	28
51	BR	59/77 (77%)	48 (81%)	11 (19%)	2	5
51	DR	59/77 (77%)	47 (80%)	12 (20%)	1	4
52	BS	70/80 (88%)	60 (86%)	10 (14%)	4	11
52	DS	67/80 (84%)	59 (88%)	8 (12%)	6	18
53	BT	70/82 (85%)	56 (80%)	14 (20%)	1	4
53	DT	71/82 (87%)	63 (89%)	8 (11%)	7	20
54	BU	18/22 (82%)	16 (89%)	2 (11%)	7	21
54	DU	18/22 (82%)	18 (100%)	0	100	100
57	BZ	609/636 (96%)	485 (80%)	124 (20%)	1	4
57	DZ	609/636 (96%)	474 (78%)	135 (22%)	1	3
All	All	10785/11672 (92%)	8911 (83%)	1874 (17%)	2	7

All (1874) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	AC	28	ARG
3	AC	32	GLU

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Mol	Chain	Res	Type
3	AC	48	LEU
3	AC	50	ILE
3	AC	53	ARG
3	AC	54	ARG
3	AC	203	GLU
4	AD	3	VAL
4	AD	12	SER
4	AD	13	ARG
4	AD	14	ARG
4	AD	25	THR
4	AD	35	LYS
4	AD	38	LYS
4	AD	61	LEU
4	AD	78	LYS
4	AD	89	SER
4	AD	94	LEU
4	AD	99	ASP
4	AD	103	ARG
4	AD	106	ILE
4	AD	111	LEU
4	AD	113	VAL
4	AD	126	GLN
4	AD	141	VAL
4	AD	142	VAL
4	AD	157	ARG
4	AD	173	VAL
4	AD	200	ASP
4	AD	202	LYS
4	AD	211	ARG
4	AD	218	ARG
4	AD	221	VAL
4	AD	229	VAL
4	AD	242	ARG
4	AD	246	PRO
4	AD	253	GLN
4	AD	257	LEU
4	AD	259	THR
4	AD	265	PRO
4	AD	274	ARG
5	AE	1	MET
5	AE	7	VAL
5	AE	12	THR

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Mol	Chain	Res	Type
5	AE	21	VAL
5	AE	33	VAL
5	AE	34	VAL
5	AE	40	GLU
5	AE	41	LYS
5	AE	47	VAL
5	AE	49	LEU
5	AE	76	ARG
5	AE	77	ILE
5	AE	78	LEU
5	AE	81	ILE
5	AE	82	ARG
5	AE	93	VAL
5	AE	111	ARG
5	AE	113	PHE
5	AE	116	VAL
5	AE	119	ARG
5	AE	121	ASN
5	AE	128	SER
5	AE	144	ARG
5	AE	145	LYS
5	AE	154	LYS
5	AE	163	GLU
5	AE	170	LEU
5	AE	175	VAL
5	AE	178	GLU
5	AE	181	LEU
5	AE	184	VAL
5	AE	185	LYS
5	AE	188	VAL
5	AE	200	GLU
6	AF	8	GLN
6	AF	12	LEU
6	AF	13	SER
6	AF	15	SER
6	AF	20	LEU
6	AF	24	LEU
6	AF	33	LEU
6	AF	38	ARG
6	AF	43	LYS
6	AF	50	SER
6	AF	53	THR

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Mol	Chain	Res	Type
6	AF	57	VAL
6	AF	72	ARG
6	AF	82	ILE
6	AF	95	ARG
6	AF	106	ARG
6	AF	110	LEU
6	AF	112	MET
6	AF	116	ASP
6	AF	125	LEU
6	AF	127	GLU
6	AF	132	VAL
6	AF	133	ASN
6	AF	140	LEU
6	AF	158	THR
6	AF	162	LEU
6	AF	170	LEU
6	AF	183	VAL
6	AF	191	ARG
6	AF	192	LEU
6	AF	197	ASP
6	AF	205	ARG
7	AG	3	LEU
7	AG	5	VAL
7	AG	7	LEU
7	AG	26	GLN
7	AG	31	VAL
7	AG	32	PRO
7	AG	37	VAL
7	AG	41	GLN
7	AG	43	LEU
7	AG	45	GLU
7	AG	79	ASN
7	AG	81	LYS
7	AG	82	LEU
7	AG	91	ARG
7	AG	103	LEU
7	AG	104	GLU
7	AG	116	ASP
7	AG	126	ASP
7	AG	128	ARG
7	AG	130	ASN
7	AG	133	LEU

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Mol	Chain	Res	Type
7	AG	136	ARG
7	AG	137	GLU
7	AG	153	ARG
7	AG	161	THR
7	AG	162	THR
7	AG	170	ARG
7	AG	175	LEU
7	AG	181	ARG
8	AH	6	ARG
8	AH	13	LYS
8	AH	15	VAL
8	AH	27	LYS
8	AH	56	SER
8	AH	59	ARG
8	AH	60	ARG
8	AH	69	ARG
8	AH	71	LEU
8	AH	81	GLU
8	AH	84	SER
8	AH	88	LEU
8	AH	89	ILE
8	AH	92	ILE
8	AH	95	ARG
8	AH	98	LEU
8	AH	105	LEU
8	AH	122	THR
8	AH	124	GLU
8	AH	136	ILE
8	AH	149	ARG
8	AH	153	LYS
8	AH	159	GLU
10	AL	2	LYS
10	AL	3	LYS
10	AL	4	VAL
10	AL	30	HIS
10	AL	34	ILE
10	AL	35	MET
10	AL	45	THR
10	AL	50	ASP
10	AL	52	ILE
10	AL	58	THR
10	AL	59	ILE

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Mol	Chain	Res	Type
10	AL	70	LYS
10	AL	85	GLU
10	AL	86	LYS
10	AL	95	LYS
10	AL	104	VAL
10	AL	106	GLU
10	AL	116	ASN
10	AL	117	THR
10	AL	118	THR
10	AL	136	VAL
11	AN	14	VAL
11	AN	28	THR
11	AN	33	LEU
11	AN	34	LEU
11	AN	38	HIS
11	AN	48	MET
11	AN	58	ASP
11	AN	61	ARG
11	AN	62	VAL
11	AN	67	LEU
11	AN	73	THR
11	AN	83	LYS
11	AN	87	LEU
11	AN	89	LYS
11	AN	97	ARG
11	AN	99	LEU
11	AN	120	LEU
11	AN	121	LYS
11	AN	133	GLN
11	AN	137	LYS
11	AN	140	VAL
12	AO	7	TYR
12	AO	8	LEU
12	AO	10	VAL
12	AO	17	ARG
12	AO	23	ARG
12	AO	24	VAL
12	AO	28	SER
12	AO	52	VAL
12	AO	61	VAL
12	AO	69	ILE
12	AO	86	ILE

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Mol	Chain	Res	Type
12	AO	94	ARG
13	AP	2	LYS
13	AP	3	LEU
13	AP	21	ARG
13	AP	42	SER
13	AP	55	ARG
13	AP	58	THR
13	AP	59	LEU
13	AP	64	LYS
13	AP	65	ARG
13	AP	70	GLN
13	AP	76	LYS
13	AP	77	ARG
13	AP	83	VAL
13	AP	98	GLU
13	AP	101	VAL
13	AP	105	LEU
13	AP	106	LEU
13	AP	107	LYS
13	AP	112	LEU
13	AP	119	GLU
13	AP	125	VAL
13	AP	126	VAL
13	AP	139	LYS
13	AP	147	LEU
13	AP	148	LEU
13	AP	149	GLU
14	AQ	2	LEU
14	AQ	5	ARG
14	AQ	7	MET
14	AQ	8	LYS
14	AQ	10	ARG
14	AQ	35	VAL
14	AQ	42	ILE
14	AQ	45	GLN
14	AQ	55	VAL
14	AQ	66	ILE
14	AQ	75	THR
14	AQ	109	VAL
15	AR	15	SER
15	AR	17	ARG
15	AR	18	LEU

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Mol	Chain	Res	Type
15	AR	20	LEU
15	AR	27	SER
15	AR	28	LEU
15	AR	29	LEU
15	AR	33	ARG
15	AR	36	THR
15	AR	44	LEU
15	AR	54	LEU
15	AR	60	LEU
15	AR	63	ARG
15	AR	65	LEU
15	AR	67	LEU
15	AR	79	LEU
15	AR	91	GLN
15	AR	100	LEU
15	AR	111	LEU
15	AR	114	VAL
16	AS	3	ARG
16	AS	4	LEU
16	AS	11	LYS
16	AS	14	VAL
16	AS	17	ARG
16	AS	20	ARG
16	AS	26	LEU
16	AS	36	TYR
16	AS	44	LYS
16	AS	48	LEU
16	AS	50	SER
16	AS	52	SER
16	AS	54	LEU
16	AS	69	VAL
16	AS	76	LYS
16	AS	83	LYS
16	AS	110	LEU
17	AT	6	LEU
17	AT	9	LEU
17	AT	11	GLU
17	AT	12	SER
17	AT	17	THR
17	AT	23	ARG
17	AT	28	VAL
17	AT	35	LYS

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Mol	Chain	Res	Type
17	AT	49	VAL
17	AT	59	THR
17	AT	78	LEU
17	AT	85	LYS
17	AT	89	VAL
17	AT	96	ARG
17	AT	115	ARG
17	AT	118	ARG
17	AT	128	GLU
18	AU	5	LYS
18	AU	8	VAL
18	AU	19	LYS
18	AU	36	ARG
18	AU	56	ASP
18	AU	59	ARG
18	AU	74	LEU
18	AU	84	LYS
18	AU	92	ARG
18	AU	95	LEU
18	AU	101	ARG
18	AU	104	GLN
18	AU	108	GLU
18	AU	111	GLU
19	AV	6	LYS
19	AV	15	GLU
19	AV	18	LEU
19	AV	21	ARG
19	AV	28	GLU
19	AV	32	THR
19	AV	43	GLU
19	AV	46	VAL
19	AV	52	VAL
19	AV	56	SER
19	AV	61	VAL
19	AV	72	VAL
19	AV	73	SER
19	AV	79	VAL
19	AV	95	LEU
19	AV	98	GLU
19	AV	100	ARG
20	AW	4	LYS
20	AW	11	ARG

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Mol	Chain	Res	Type
20	AW	12	ILE
20	AW	17	VAL
20	AW	19	LEU
20	AW	23	LEU
20	AW	35	ILE
20	AW	41	LYS
20	AW	51	LEU
20	AW	65	LEU
20	AW	67	ASP
20	AW	90	ARG
20	AW	107	LEU
21	AX	33	LYS
21	AX	35	THR
21	AX	45	THR
21	AX	57	LEU
21	AX	65	ARG
21	AX	66	LEU
21	AX	70	LEU
21	AX	72	LYS
22	AY	1	MET
22	AY	7	VAL
22	AY	8	LYS
22	AY	23	ARG
22	AY	26	LYS
22	AY	47	LYS
22	AY	50	ARG
22	AY	55	TYR
22	AY	72	VAL
22	AY	73	ARG
22	AY	81	LYS
22	AY	90	LEU
22	AY	91	GLU
22	AY	96	ILE
23	AZ	5	LEU
23	AZ	16	SER
23	AZ	18	LEU
23	AZ	19	ARG
23	AZ	28	MET
23	AZ	31	ARG
23	AZ	37	VAL
23	AZ	40	ASP
23	AZ	42	VAL

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Mol	Chain	Res	Type
23	AZ	46	LYS
23	AZ	52	SER
23	AZ	56	VAL
23	AZ	65	GLN
23	AZ	72	ARG
23	AZ	73	GLN
23	AZ	74	VAL
23	AZ	77	ASP
23	AZ	78	LYS
23	AZ	86	VAL
23	AZ	87	ASP
23	AZ	91	LEU
23	AZ	97	GLU
23	AZ	98	MET
23	AZ	124	ILE
23	AZ	129	SER
23	AZ	136	PHE
23	AZ	149	SER
23	AZ	150	LEU
23	AZ	151	HIS
23	AZ	156	LYS
23	AZ	157	LEU
23	AZ	163	LEU
23	AZ	170	THR
23	AZ	180	VAL
23	AZ	183	LEU
24	A0	7	LEU
24	A0	20	ARG
24	A0	43	THR
24	A0	55	ARG
25	A1	21	ARG
25	A1	30	VAL
25	A1	32	LYS
25	A1	40	ARG
25	A1	51	VAL
25	A1	59	THR
25	A1	75	GLU
25	A1	80	LEU
25	A1	92	LYS
25	A1	95	LEU
26	A2	8	LYS
26	A2	32	LEU

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Mol	Chain	Res	Type
26	A2	40	SER
26	A2	41	ILE
26	A2	52	ASP
26	A2	53	LEU
26	A2	62	THR
26	A2	64	LEU
26	A2	67	LYS
26	A2	70	GLN
27	A3	5	LYS
27	A3	8	LEU
27	A3	18	ASP
27	A3	23	LEU
27	A3	29	ARG
27	A3	32	GLN
27	A3	34	GLU
27	A3	35	ARG
27	A3	54	VAL
27	A3	58	VAL
28	A4	1	MET
28	A4	8	LYS
28	A4	22	ILE
28	A4	27	THR
28	A4	31	ILE
28	A4	34	GLU
28	A4	35	VAL
28	A4	36	CYS
28	A4	39	CYS
28	A4	46	GLN
28	A4	48	ARG
28	A4	49	PHE
28	A4	56	VAL
28	A4	57	GLU
28	A4	58	ARG
28	A4	61	ARG
28	A4	67	TYR
28	A4	68	ARG
29	A5	6	VAL
29	A5	16	ARG
29	A5	29	THR
29	A5	33	CYS
29	A5	40	LYS
29	A5	55	ARG

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Mol	Chain	Res	Type
29	A5	60	VAL
30	A6	4	GLU
30	A6	6	ARG
30	A6	7	ILE
30	A6	14	THR
30	A6	24	GLU
30	A6	25	LYS
30	A6	33	LYS
30	A6	44	ARG
30	A6	47	THR
30	A6	48	VAL
30	A6	50	ARG
30	A6	54	ILE
31	A7	1	MET
31	A7	9	ARG
31	A7	10	ARG
31	A7	14	LYS
31	A7	24	THR
31	A7	43	THR
31	A7	46	VAL
32	A8	13	ARG
32	A8	14	VAL
32	A8	30	ARG
32	A8	31	HIS
32	A8	32	LEU
32	A8	43	GLN
32	A8	46	ARG
32	A8	52	LYS
33	A9	4	ARG
33	A9	17	ILE
35	BB	8	LYS
35	BB	11	LEU
35	BB	15	VAL
35	BB	16	HIS
35	BB	17	PHE
35	BB	19	HIS
35	BB	20	GLU
35	BB	21	ARG
35	BB	24	TRP
35	BB	37	ASN
35	BB	41	ILE
35	BB	48	MET

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Mol	Chain	Res	Type
35	BB	56	ARG
35	BB	64	ARG
35	BB	76	GLN
35	BB	78	GLN
35	BB	80	ILE
35	BB	83	MET
35	BB	98	LEU
35	BB	107	THR
35	BB	110	GLN
35	BB	111	ARG
35	BB	112	VAL
35	BB	113	HIS
35	BB	127	ILE
35	BB	142	LEU
35	BB	145	LEU
35	BB	160	ASP
35	BB	168	THR
35	BB	169	LYS
35	BB	187	LEU
35	BB	195	ASP
35	BB	200	ILE
35	BB	208	ILE
35	BB	217	ARG
35	BB	221	LEU
35	BB	223	ILE
35	BB	224	GLN
35	BB	229	VAL
36	BC	3	ASN
36	BC	15	THR
36	BC	28	GLN
36	BC	29	TYR
36	BC	45	LYS
36	BC	67	THR
36	BC	108	ASN
36	BC	131	ARG
36	BC	143	GLU
36	BC	178	LEU
36	BC	181	ASN
36	BC	192	THR
36	BC	196	LEU
36	BC	207	VAL
37	BD	5	ILE

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Mol	Chain	Res	Type
37	BD	13	ARG
37	BD	15	GLU
37	BD	31	CYS
37	BD	58	LEU
37	BD	61	LYS
37	BD	70	ILE
37	BD	73	ARG
37	BD	76	ARG
37	BD	101	LEU
37	BD	103	ASN
37	BD	107	ARG
37	BD	119	GLN
37	BD	126	ILE
37	BD	127	THR
37	BD	134	ASP
37	BD	135	LEU
37	BD	139	ARG
37	BD	140	VAL
37	BD	142	PRO
37	BD	158	ILE
37	BD	168	ARG
37	BD	173	TRP
37	BD	182	LYS
37	BD	193	ASP
37	BD	196	LEU
37	BD	200	GLU
38	BE	10	MET
38	BE	11	ILE
38	BE	18	ARG
38	BE	20	GLN
38	BE	25	ARG
38	BE	27	ARG
38	BE	28	PHE
38	BE	31	LEU
38	BE	34	VAL
38	BE	38	GLN
38	BE	40	ARG
38	BE	41	VAL
38	BE	47	LYS
38	BE	76	ILE
38	BE	78	HIS
38	BE	79	GLU

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Mol	Chain	Res	Type
38	BE	91	LEU
38	BE	121	LYS
38	BE	137	GLU
38	BE	140	ARG
38	BE	147	ASP
38	BE	148	VAL
38	BE	149	GLU
38	BE	150	ARG
39	BF	30	LEU
39	BF	37	VAL
39	BF	40	VAL
39	BF	61	LEU
39	BF	64	GLN
39	BF	69	GLU
39	BF	71	ARG
39	BF	72	VAL
39	BF	73	ASN
39	BF	75	LEU
39	BF	82	ARG
39	BF	83	ASP
39	BF	86	ARG
39	BF	92	LYS
39	BF	94	GLN
39	BF	98	LEU
40	BG	8	GLU
40	BG	12	LEU
40	BG	15	ASP
40	BG	16	LEU
40	BG	29	LYS
40	BG	38	LEU
40	BG	41	ARG
40	BG	51	GLN
40	BG	61	VAL
40	BG	73	MET
40	BG	75	VAL
40	BG	78	ARG
40	BG	98	SER
40	BG	104	LEU
40	BG	113	GLU
40	BG	114	ARG
40	BG	115	ARG
41	BH	2	LEU

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Mol	Chain	Res	Type
41	BH	10	LEU
41	BH	18	ARG
41	BH	21	LYS
41	BH	24	THR
41	BH	25	ASP
41	BH	26	VAL
41	BH	51	VAL
41	BH	52	ASP
41	BH	54	ASP
41	BH	75	ARG
41	BH	77	GLU
41	BH	78	GLN
41	BH	91	ARG
41	BH	98	LYS
41	BH	112	LEU
41	BH	122	ARG
41	BH	133	LEU
42	BI	3	GLN
42	BI	23	ASN
42	BI	25	LYS
42	BI	27	THR
42	BI	53	VAL
42	BI	64	THR
42	BI	75	ASP
42	BI	81	ILE
42	BI	92	TYR
42	BI	93	ARG
42	BI	102	LEU
42	BI	104	ARG
42	BI	110	GLU
42	BI	113	LYS
42	BI	127	LYS
42	BI	128	ARG
43	BJ	16	LEU
43	BJ	21	GLN
43	BJ	65	LEU
43	BJ	94	VAL
43	BJ	100	THR
44	BK	14	VAL
44	BK	16	SER
44	BK	18	ARG
44	BK	31	THR

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Mol	Chain	Res	Type
44	BK	40	ILE
44	BK	48	ILE
44	BK	51	LYS
44	BK	83	ILE
44	BK	84	VAL
44	BK	87	THR
44	BK	96	ARG
44	BK	98	LEU
44	BK	104	GLN
44	BK	109	VAL
44	BK	112	THR
45	BL	7	ILE
45	BL	18	VAL
45	BL	23	LYS
45	BL	27	LEU
45	BL	33	ARG
45	BL	36	VAL
45	BL	46	LYS
45	BL	53	ARG
45	BL	57	LYS
45	BL	60	LEU
45	BL	67	THR
45	BL	83	VAL
45	BL	97	ARG
45	BL	100	ILE
46	BM	3	ARG
46	BM	4	ILE
46	BM	11	ARG
46	BM	15	VAL
46	BM	16	ASP
46	BM	17	VAL
46	BM	19	LEU
46	BM	32	GLU
46	BM	37	THR
46	BM	39	ILE
46	BM	44	ARG
46	BM	50	GLU
46	BM	57	ARG
46	BM	63	THR
46	BM	66	LEU
46	BM	70	LEU
46	BM	84	ILE

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Mol	Chain	Res	Type
46	BM	102	ARG
46	BM	105	THR
46	BM	110	ARG
46	BM	116	THR
47	BN	3	ARG
47	BN	7	ILE
47	BN	18	VAL
47	BN	23	ARG
47	BN	33	VAL
47	BN	41	ARG
47	BN	44	LEU
47	BN	49	HIS
47	BN	57	ARG
48	BO	3	ILE
48	BO	5	LYS
48	BO	26	GLU
48	BO	35	ARG
48	BO	39	LEU
48	BO	41	GLU
48	BO	42	HIS
48	BO	66	LEU
48	BO	88	ARG
49	BP	1	MET
49	BP	2	VAL
49	BP	4	ILE
49	BP	5	ARG
49	BP	8	ARG
49	BP	11	SER
49	BP	16	HIS
49	BP	18	ARG
49	BP	19	ILE
49	BP	29	ASP
49	BP	50	LYS
49	BP	54	GLU
49	BP	69	THR
49	BP	71	ARG
49	BP	74	LEU
49	BP	76	GLN
50	BQ	5	VAL
50	BQ	6	LEU
50	BQ	9	VAL
50	BQ	36	ILE

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Mol	Chain	Res	Type
50	BQ	43	LEU
50	BQ	45	HIS
50	BQ	49	GLU
50	BQ	53	LEU
50	BQ	60	ILE
50	BQ	62	SER
50	BQ	68	ARG
50	BQ	70	ARG
50	BQ	72	ARG
50	BQ	73	VAL
50	BQ	74	LEU
50	BQ	77	VAL
50	BQ	78	GLU
50	BQ	90	ILE
50	BQ	91	ARG
50	BQ	92	ARG
51	BR	26	LEU
51	BR	29	PHE
51	BR	31	LEU
51	BR	35	ARG
51	BR	37	VAL
51	BR	51	LEU
51	BR	53	ARG
51	BR	55	ARG
51	BR	58	LEU
51	BR	76	LEU
51	BR	85	LEU
52	BS	4	SER
52	BS	12	ASP
52	BS	28	LYS
52	BS	34	TRP
52	BS	65	ASN
52	BS	66	MET
52	BS	78	ARG
52	BS	81	ARG
52	BS	83	HIS
52	BS	85	LYS
53	BT	8	ARG
53	BT	13	LEU
53	BT	15	ARG
53	BT	31	SER
53	BT	45	GLN

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Mol	Chain	Res	Type
53	BT	58	LYS
53	BT	60	GLU
53	BT	61	SER
53	BT	62	LEU
53	BT	70	SER
53	BT	75	ASN
53	BT	84	LEU
53	BT	93	GLU
53	BT	100	ILE
54	BU	9	ARG
54	BU	10	ARG
57	BZ	-64	VAL
57	BZ	-58	LEU
57	BZ	-56	ASN
57	BZ	-52	VAL
57	BZ	-49	VAL
57	BZ	-45	LYS
57	BZ	-42	TYR
57	BZ	-32	LEU
57	BZ	-30	VAL
57	BZ	-29	LEU
57	BZ	-25	SER
57	BZ	-22	LYS
57	BZ	-20	LEU
57	BZ	-10	ARG
57	BZ	-6	ARG
57	BZ	-3	GLU
57	BZ	1	LEU
57	BZ	2	LYS
57	BZ	8	ASP
57	BZ	12	LEU
57	BZ	13	ARG
57	BZ	21	ILE
57	BZ	30	GLU
57	BZ	33	LEU
57	BZ	39	ILE
57	BZ	70	THR
57	BZ	79	ILE
57	BZ	81	ILE
57	BZ	83	ASP
57	BZ	88	VAL
57	BZ	91	THR

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Mol	Chain	Res	Type
57	BZ	92	ILE
57	BZ	97	SER
57	BZ	107	VAL
57	BZ	115	GLU
57	BZ	117	GLN
57	BZ	130	VAL
57	BZ	132	ARG
57	BZ	146	LEU
57	BZ	152	THR
57	BZ	157	LEU
57	BZ	160	ARG
57	BZ	171	GLU
57	BZ	172	ASP
57	BZ	182	ARG
57	BZ	186	TYR
57	BZ	196	ILE
57	BZ	198	GLU
57	BZ	210	ARG
57	BZ	238	THR
57	BZ	240	GLU
57	BZ	247	ARG
57	BZ	252	ASP
57	BZ	255	ILE
57	BZ	256	THR
57	BZ	264	LEU
57	BZ	269	VAL
57	BZ	271	LEU
57	BZ	279	TYR
57	BZ	284	LEU
57	BZ	285	ASP
57	BZ	286	ILE
57	BZ	289	ILE
57	BZ	297	GLU
57	BZ	298	VAL
57	BZ	301	ILE
57	BZ	312	LEU
57	BZ	324	ARG
57	BZ	328	ILE
57	BZ	329	ARG
57	BZ	352	VAL
57	BZ	354	ARG
57	BZ	356	LEU

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Mol	Chain	Res	Type
57	BZ	361	ASN
57	BZ	363	ARG
57	BZ	368	GLU
57	BZ	377	VAL
57	BZ	385	THR
57	BZ	396	ARG
57	BZ	399	LEU
57	BZ	409	ILE
57	BZ	420	ASP
57	BZ	421	GLN
57	BZ	422	GLU
57	BZ	424	LEU
57	BZ	425	SER
57	BZ	426	GLN
57	BZ	431	LEU
57	BZ	437	THR
57	BZ	438	PHE
57	BZ	442	THR
57	BZ	445	GLU
57	BZ	457	LEU
57	BZ	461	ILE
57	BZ	464	ASP
57	BZ	468	ARG
57	BZ	473	ASP
57	BZ	478	LYS
57	BZ	481	VAL
57	BZ	484	ARG
57	BZ	488	THR
57	BZ	491	VAL
57	BZ	498	ILE
57	BZ	506	GLN
57	BZ	510	VAL
57	BZ	512	ILE
57	BZ	517	LEU
57	BZ	526	VAL
57	BZ	529	ILE
57	BZ	566	THR
57	BZ	587	SER
57	BZ	590	ILE
57	BZ	592	GLU
57	BZ	610	VAL
57	BZ	614	GLU

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Mol	Chain	Res	Type
57	BZ	624	LEU
57	BZ	631	ILE
57	BZ	634	MET
57	BZ	641	GLN
57	BZ	644	ARG
57	BZ	651	GLU
57	BZ	659	LEU
57	BZ	677	GLN
57	BZ	679	VAL
3	CC	28	ARG
3	CC	32	GLU
3	CC	48	LEU
3	CC	50	ILE
3	CC	53	ARG
3	CC	54	ARG
3	CC	203	GLU
3	CC	208	THR
4	CD	3	VAL
4	CD	10	THR
4	CD	12	SER
4	CD	14	ARG
4	CD	27	THR
4	CD	54	ARG
4	CD	61	LEU
4	CD	69	ARG
4	CD	71	ASP
4	CD	89	SER
4	CD	94	LEU
4	CD	98	VAL
4	CD	101	GLU
4	CD	103	ARG
4	CD	105	ILE
4	CD	106	ILE
4	CD	111	LEU
4	CD	134	ARG
4	CD	140	THR
4	CD	142	VAL
4	CD	147	LEU
4	CD	173	VAL
4	CD	193	VAL
4	CD	211	ARG
4	CD	217	ARG

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Mol	Chain	Res	Type
4	CD	221	VAL
4	CD	229	VAL
4	CD	242	ARG
4	CD	253	GLN
4	CD	257	LEU
4	CD	259	THR
4	CD	260	ARG
4	CD	261	LYS
4	CD	262	ARG
4	CD	265	PRO
4	CD	274	ARG
4	CD	276	LYS
5	CE	2	LYS
5	CE	9	VAL
5	CE	12	THR
5	CE	19	ARG
5	CE	21	VAL
5	CE	24	THR
5	CE	33	VAL
5	CE	34	VAL
5	CE	38	THR
5	CE	40	GLU
5	CE	52	LEU
5	CE	58	ARG
5	CE	63	LEU
5	CE	73	GLU
5	CE	74	PRO
5	CE	75	VAL
5	CE	82	ARG
5	CE	84	PHE
5	CE	85	ASN
5	CE	92	THR
5	CE	93	VAL
5	CE	101	ARG
5	CE	116	VAL
5	CE	119	ARG
5	CE	144	ARG
5	CE	145	LYS
5	CE	154	LYS
5	CE	165	VAL
5	CE	179	GLU
5	CE	181	LEU

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Mol	Chain	Res	Type
5	CE	182	LEU
5	CE	195	LEU
5	CE	196	VAL
5	CE	202	LYS
6	CF	19	GLU
6	CF	20	LEU
6	CF	23	ASP
6	CF	24	LEU
6	CF	27	GLU
6	CF	28	ILE
6	CF	43	LYS
6	CF	48	THR
6	CF	50	SER
6	CF	53	THR
6	CF	57	VAL
6	CF	66	PRO
6	CF	70	THR
6	CF	74	ARG
6	CF	78	ILE
6	CF	82	ILE
6	CF	88	VAL
6	CF	102	PRO
6	CF	104	LYS
6	CF	106	ARG
6	CF	119	ARG
6	CF	135	LYS
6	CF	137	LYS
6	CF	140	LEU
6	CF	158	THR
6	CF	162	LEU
6	CF	169	ASN
6	CF	170	LEU
6	CF	175	THR
6	CF	176	LEU
6	CF	181	LEU
6	CF	192	LEU
6	CF	200	GLU
6	CF	203	GLN
6	CF	205	ARG
7	CG	3	LEU
7	CG	5	VAL
7	CG	7	LEU

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Mol	Chain	Res	Type
7	CG	9	ARG
7	CG	16	ARG
7	CG	18	GLU
7	CG	20	ILE
7	CG	32	PRO
7	CG	43	LEU
7	CG	60	LEU
7	CG	63	ILE
7	CG	70	VAL
7	CG	71	THR
7	CG	75	LYS
7	CG	108	ASN
7	CG	115	ARG
7	CG	126	ASP
7	CG	128	ARG
7	CG	136	ARG
7	CG	138	GLN
7	CG	143	GLU
7	CG	148	MET
7	CG	150	ASP
7	CG	153	ARG
7	CG	161	THR
7	CG	164	GLU
7	CG	165	THR
7	CG	170	ARG
7	CG	173	LEU
7	CG	174	GLU
7	CG	175	LEU
8	CH	3	ARG
8	CH	4	ILE
8	CH	6	ARG
8	CH	7	LEU
8	CH	33	LEU
8	CH	41	MET
8	CH	49	VAL
8	CH	59	ARG
8	CH	69	ARG
8	CH	71	LEU
8	CH	80	SER
8	CH	105	LEU
8	CH	106	THR
8	CH	122	THR

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Mol	Chain	Res	Type
8	CH	125	VAL
8	CH	130	ARG
8	CH	136	ILE
8	CH	139	GLN
8	CH	149	ARG
8	CH	171	LEU
10	CL	2	LYS
10	CL	4	VAL
10	CL	29	GLN
10	CL	30	HIS
10	CL	38	VAL
10	CL	59	ILE
10	CL	65	PHE
10	CL	75	SER
10	CL	77	LEU
10	CL	86	LYS
10	CL	93	ARG
10	CL	95	LYS
10	CL	96	VAL
10	CL	102	GLU
10	CL	105	LEU
10	CL	112	MET
10	CL	117	THR
10	CL	118	THR
10	CL	134	MET
10	CL	136	VAL
11	CN	10	GLU
11	CN	16	ILE
11	CN	22	THR
11	CN	25	ARG
11	CN	26	LEU
11	CN	29	LYS
11	CN	32	THR
11	CN	33	LEU
11	CN	34	LEU
11	CN	38	HIS
11	CN	46	VAL
11	CN	48	MET
11	CN	58	ASP
11	CN	59	LYS
11	CN	63	THR
11	CN	70	LYS

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Mol	Chain	Res	Type
11	CN	76	SER
11	CN	87	LEU
11	CN	97	ARG
11	CN	99	LEU
11	CN	109	LYS
11	CN	120	LEU
11	CN	127	ASP
11	CN	133	GLN
11	CN	137	LYS
11	CN	138	LEU
12	CO	1	MET
12	CO	17	ARG
12	CO	18	LYS
12	CO	19	ILE
12	CO	22	ILE
12	CO	23	ARG
12	CO	24	VAL
12	CO	35	VAL
12	CO	52	VAL
12	CO	69	ILE
12	CO	77	ILE
12	CO	78	ARG
12	CO	86	ILE
12	CO	91	LEU
12	CO	96	THR
12	CO	98	VAL
12	CO	113	LYS
12	CO	116	SER
13	CP	2	LYS
13	CP	3	LEU
13	CP	15	ARG
13	CP	21	ARG
13	CP	42	SER
13	CP	55	ARG
13	CP	57	THR
13	CP	59	LEU
13	CP	65	ARG
13	CP	75	ILE
13	CP	86	LYS
13	CP	92	GLU
13	CP	95	VAL
13	CP	106	LEU

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Mol	Chain	Res	Type
13	CP	112	LEU
13	CP	125	VAL
13	CP	132	LYS
13	CP	135	LEU
13	CP	147	LEU
13	CP	148	LEU
14	CQ	1	MET
14	CQ	3	MET
14	CQ	6	ARG
14	CQ	12	GLN
14	CQ	16	ARG
14	CQ	21	THR
14	CQ	35	VAL
14	CQ	45	GLN
14	CQ	57	HIS
14	CQ	60	ARG
14	CQ	63	LYS
14	CQ	75	THR
14	CQ	94	VAL
14	CQ	109	VAL
14	CQ	110	THR
14	CQ	111	GLU
14	CQ	112	GLU
14	CQ	127	ILE
14	CQ	128	LYS
14	CQ	133	ARG
15	CR	1	MET
15	CR	4	LEU
15	CR	9	LYS
15	CR	14	SER
15	CR	18	LEU
15	CR	22	ARG
15	CR	28	LEU
15	CR	29	LEU
15	CR	37	THR
15	CR	40	LYS
15	CR	43	GLU
15	CR	44	LEU
15	CR	54	LEU
15	CR	60	LEU
15	CR	63	ARG
15	CR	65	LEU

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Mol	Chain	Res	Type
15	CR	75	LEU
15	CR	79	LEU
15	CR	86	ARG
15	CR	88	ARG
15	CR	90	ARG
15	CR	102	GLU
15	CR	114	VAL
16	CS	3	ARG
16	CS	4	LEU
16	CS	12	PHE
16	CS	13	ARG
16	CS	19	LYS
16	CS	20	ARG
16	CS	26	LEU
16	CS	29	PHE
16	CS	31	SER
16	CS	35	ILE
16	CS	36	TYR
16	CS	49	VAL
16	CS	50	SER
16	CS	52	SER
16	CS	58	LEU
16	CS	62	LYS
16	CS	67	ARG
16	CS	78	LEU
16	CS	110	LEU
17	CT	6	LEU
17	CT	8	LYS
17	CT	13	ARG
17	CT	18	ASP
17	CT	19	LEU
17	CT	23	ARG
17	CT	39	ARG
17	CT	42	ILE
17	CT	51	ARG
17	CT	53	ARG
17	CT	59	THR
17	CT	64	ARG
17	CT	65	LYS
17	CT	67	SER
17	CT	74	ARG
17	CT	78	LEU

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Mol	Chain	Res	Type
17	CT	85	LYS
17	CT	87	ASP
17	CT	96	ARG
17	CT	98	LYS
17	CT	115	ARG
17	CT	118	ARG
17	CT	123	GLN
18	CU	5	LYS
18	CU	31	SER
18	CU	36	ARG
18	CU	52	ARG
18	CU	59	ARG
18	CU	69	CYS
18	CU	74	LEU
18	CU	84	LYS
18	CU	93	LYS
18	CU	95	LEU
18	CU	100	VAL
18	CU	104	GLN
18	CU	108	GLU
19	CV	6	LYS
19	CV	13	ARG
19	CV	15	GLU
19	CV	18	LEU
19	CV	52	VAL
19	CV	61	VAL
19	CV	62	LEU
19	CV	69	LYS
19	CV	72	VAL
19	CV	73	SER
19	CV	95	LEU
19	CV	99	ILE
20	CW	6	ILE
20	CW	11	ARG
20	CW	19	LEU
20	CW	23	LEU
20	CW	29	LEU
20	CW	60	ASN
20	CW	100	THR
20	CW	107	LEU
21	CX	9	LEU
21	CX	35	THR

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Mol	Chain	Res	Type
21	CX	45	THR
21	CX	52	VAL
21	CX	57	LEU
21	CX	65	ARG
21	CX	82	GLN
21	CX	87	GLN
22	CY	6	HIS
22	CY	7	VAL
22	CY	23	ARG
22	CY	34	LYS
22	CY	43	ASN
22	CY	45	VAL
22	CY	49	VAL
22	CY	64	GLU
22	CY	67	LEU
22	CY	72	VAL
22	CY	85	VAL
22	CY	88	LYS
22	CY	91	GLU
22	CY	95	LYS
22	CY	102	CYS
22	CY	107	ASP
23	CZ	6	LYS
23	CZ	16	SER
23	CZ	19	ARG
23	CZ	33	LEU
23	CZ	37	VAL
23	CZ	42	VAL
23	CZ	61	LEU
23	CZ	72	ARG
23	CZ	76	LEU
23	CZ	78	LYS
23	CZ	80	ARG
23	CZ	81	ARG
23	CZ	87	ASP
23	CZ	89	PHE
23	CZ	91	LEU
23	CZ	98	MET
23	CZ	107	THR
23	CZ	111	VAL
23	CZ	124	ILE
23	CZ	136	PHE

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Mol	Chain	Res	Type
23	CZ	149	SER
23	CZ	150	LEU
23	CZ	153	SER
23	CZ	154	ASP
23	CZ	156	LYS
23	CZ	161	VAL
23	CZ	170	THR
24	C0	3	HIS
24	C0	11	ARG
24	C0	20	ARG
24	C0	74	ARG
24	C0	82	ARG
25	C1	4	VAL
25	C1	30	VAL
25	C1	37	ILE
25	C1	38	SER
25	C1	40	ARG
25	C1	41	ARG
25	C1	52	ARG
25	C1	59	THR
25	C1	62	VAL
25	C1	69	LYS
25	C1	78	LYS
25	C1	80	LEU
25	C1	94	LEU
25	C1	95	LEU
26	C2	2	LYS
26	C2	9	GLN
26	C2	14	ARG
26	C2	28	LYS
26	C2	30	ARG
26	C2	32	LEU
26	C2	41	ILE
26	C2	51	ARG
26	C2	52	ASP
26	C2	53	LEU
26	C2	59	ARG
26	C2	62	THR
26	C2	63	VAL
26	C2	70	GLN
27	C3	6	VAL
27	C3	8	LEU

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Mol	Chain	Res	Type
27	C3	23	LEU
27	C3	24	LYS
27	C3	30	ARG
27	C3	36	VAL
27	C3	50	VAL
27	C3	56	VAL
28	C4	5	ILE
28	C4	34	GLU
28	C4	35	VAL
28	C4	50	VAL
28	C4	56	VAL
28	C4	58	ARG
28	C4	59	PHE
28	C4	61	ARG
28	C4	63	TYR
28	C4	67	TYR
28	C4	68	ARG
28	C4	69	LYS
29	C5	6	VAL
29	C5	9	LYS
29	C5	12	SER
29	C5	16	ARG
29	C5	20	ARG
29	C5	36	CYS
29	C5	40	LYS
29	C5	57	VAL
30	C6	5	VAL
30	C6	6	ARG
30	C6	7	ILE
30	C6	23	THR
30	C6	25	LYS
30	C6	32	ASN
30	C6	38	LYS
30	C6	40	CYS
30	C6	44	ARG
30	C6	48	VAL
30	C6	50	ARG
31	C7	1	MET
31	C7	4	THR
31	C7	9	ARG
31	C7	14	LYS
31	C7	29	LYS

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Mol	Chain	Res	Type
31	C7	32	LYS
31	C7	43	THR
31	C7	47	ARG
31	C7	48	LYS
32	C8	3	LYS
32	C8	6	THR
32	C8	14	VAL
32	C8	26	LYS
32	C8	30	ARG
32	C8	31	HIS
32	C8	32	LEU
32	C8	34	TRP
32	C8	36	LYS
32	C8	46	ARG
33	C9	19	ARG
33	C9	27	CYS
35	DB	7	VAL
35	DB	12	GLU
35	DB	24	TRP
35	DB	27	LYS
35	DB	44	LEU
35	DB	47	THR
35	DB	51	LEU
35	DB	53	ARG
35	DB	56	ARG
35	DB	58	ILE
35	DB	67	THR
35	DB	68	ILE
35	DB	80	ILE
35	DB	87	ARG
35	DB	94	ASN
35	DB	107	THR
35	DB	108	ILE
35	DB	113	HIS
35	DB	115	LEU
35	DB	117	GLU
35	DB	119	GLU
35	DB	122	PHE
35	DB	128	GLU
35	DB	154	LEU
35	DB	155	LEU
35	DB	157	ARG

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Mol	Chain	Res	Type
35	DB	170	GLU
35	DB	179	LYS
35	DB	185	ILE
35	DB	187	LEU
35	DB	189	ASP
35	DB	210	SER
35	DB	212	GLN
35	DB	217	ARG
35	DB	224	GLN
35	DB	229	VAL
36	DC	3	ASN
36	DC	6	HIS
36	DC	16	ARG
36	DC	26	LYS
36	DC	40	ARG
36	DC	43	LEU
36	DC	47	LEU
36	DC	52	LEU
36	DC	56	ASP
36	DC	63	ASN
36	DC	102	ASN
36	DC	104	GLN
36	DC	105	GLU
36	DC	108	ASN
36	DC	115	LEU
36	DC	120	VAL
36	DC	131	ARG
36	DC	152	ILE
36	DC	178	LEU
36	DC	188	LEU
36	DC	190	ARG
36	DC	191	THR
36	DC	195	VAL
36	DC	202	ILE
36	DC	207	VAL
37	DD	5	ILE
37	DD	8	VAL
37	DD	12	CYS
37	DD	13	ARG
37	DD	19	LEU
37	DD	28	SER
37	DD	31	CYS

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Mol	Chain	Res	Type
37	DD	34	GLU
37	DD	47	ARG
37	DD	57	ARG
37	DD	58	LEU
37	DD	61	LYS
37	DD	65	ARG
37	DD	73	ARG
37	DD	76	ARG
37	DD	83	SER
37	DD	85	LYS
37	DD	91	SER
37	DD	103	ASN
37	DD	115	ARG
37	DD	120	LEU
37	DD	126	ILE
37	DD	127	THR
37	DD	129	ASN
37	DD	132	ARG
37	DD	135	LEU
37	DD	139	ARG
37	DD	140	VAL
37	DD	150	GLU
37	DD	153	ARG
37	DD	155	LEU
37	DD	162	LEU
37	DD	170	VAL
37	DD	187	ARG
37	DD	188	LEU
37	DD	194	LEU
37	DD	208	SER
38	DE	11	ILE
38	DE	24	ARG
38	DE	25	ARG
38	DE	31	LEU
38	DE	41	VAL
38	DE	47	LYS
38	DE	51	VAL
38	DE	55	VAL
38	DE	68	GLU
38	DE	69	VAL
38	DE	71	LEU
38	DE	75	THR

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Mol	Chain	Res	Type
38	DE	78	HIS
38	DE	87	SER
38	DE	89	ILE
38	DE	90	VAL
38	DE	91	LEU
38	DE	107	ARG
38	DE	116	THR
38	DE	135	THR
38	DE	137	GLU
38	DE	150	ARG
39	DF	25	ILE
39	DF	27	GLN
39	DF	28	ARG
39	DF	48	LEU
39	DF	69	GLU
39	DF	70	ASP
39	DF	72	VAL
39	DF	79	LEU
39	DF	82	ARG
39	DF	86	ARG
39	DF	87	ARG
40	DG	8	GLU
40	DG	10	ARG
40	DG	11	GLN
40	DG	24	THR
40	DG	32	ARG
40	DG	41	ARG
40	DG	47	CYS
40	DG	57	GLU
40	DG	67	GLU
40	DG	72	ARG
40	DG	73	MET
40	DG	76	ARG
40	DG	85	TYR
40	DG	87	VAL
40	DG	90	GLU
40	DG	104	LEU
40	DG	114	ARG
40	DG	115	ARG
40	DG	140	ASP
40	DG	153	HIS
40	DG	154	TYR

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Mol	Chain	Res	Type
41	DH	2	LEU
41	DH	21	LYS
41	DH	24	THR
41	DH	25	ASP
41	DH	26	VAL
41	DH	34	GLU
41	DH	37	ARG
41	DH	51	VAL
41	DH	68	ARG
41	DH	78	GLN
41	DH	84	ARG
41	DH	85	ARG
41	DH	91	ARG
41	DH	95	VAL
41	DH	98	LYS
41	DH	99	GLU
41	DH	104	ARG
41	DH	111	ILE
41	DH	112	LEU
41	DH	120	THR
41	DH	133	LEU
41	DH	137	VAL
42	DI	7	THR
42	DI	20	ARG
42	DI	23	ASN
42	DI	27	THR
42	DI	41	VAL
42	DI	53	VAL
42	DI	66	ARG
42	DI	81	ILE
42	DI	92	TYR
42	DI	102	LEU
42	DI	103	THR
42	DI	104	ARG
42	DI	109	VAL
43	DJ	17	ASP
43	DJ	38	ILE
43	DJ	58	ASP
43	DJ	72	VAL
44	DK	28	THR
44	DK	30	VAL
44	DK	33	THR

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Mol	Chain	Res	Type
44	DK	54	ARG
44	DK	96	ARG
44	DK	114	VAL
45	DL	8	ASN
45	DL	24	VAL
45	DL	27	LEU
45	DL	33	ARG
45	DL	38	THR
45	DL	39	VAL
45	DL	47	LYS
45	DL	55	VAL
45	DL	59	ARG
45	DL	83	VAL
45	DL	104	VAL
45	DL	106	ASP
45	DL	113	ARG
45	DL	114	LYS
45	DL	118	SER
46	DM	3	ARG
46	DM	15	VAL
46	DM	29	ARG
46	DM	56	LEU
46	DM	66	LEU
46	DM	70	LEU
46	DM	73	GLU
46	DM	91	ARG
46	DM	103	THR
46	DM	106	ASN
46	DM	108	ARG
46	DM	110	ARG
46	DM	114	ARG
47	DN	12	ARG
47	DN	15	LYS
47	DN	22	THR
47	DN	33	VAL
47	DN	35	ARG
47	DN	41	ARG
47	DN	42	ILE
47	DN	44	LEU
47	DN	58	LYS
48	DO	3	ILE
48	DO	4	THR

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Mol	Chain	Res	Type
48	DO	22	THR
48	DO	26	GLU
48	DO	38	ARG
48	DO	39	LEU
48	DO	48	LYS
48	DO	58	MET
48	DO	66	LEU
48	DO	68	ARG
49	DP	2	VAL
49	DP	4	ILE
49	DP	5	ARG
49	DP	8	ARG
49	DP	11	SER
49	DP	20	VAL
49	DP	21	VAL
49	DP	25	ARG
49	DP	27	LYS
49	DP	60	LEU
49	DP	62	VAL
49	DP	69	THR
50	DQ	6	LEU
50	DQ	13	ASP
50	DQ	37	LYS
50	DQ	50	LYS
50	DQ	52	LYS
50	DQ	69	LYS
50	DQ	70	ARG
50	DQ	72	ARG
50	DQ	76	LEU
51	DR	21	LYS
51	DR	25	THR
51	DR	31	LEU
51	DR	32	ARG
51	DR	37	VAL
51	DR	41	LYS
51	DR	47	THR
51	DR	53	ARG
51	DR	69	THR
51	DR	76	LEU
51	DR	84	LYS
51	DR	85	LEU
52	DS	12	ASP

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Mol	Chain	Res	Type
52	DS	15	LEU
52	DS	16	LEU
52	DS	33	THR
52	DS	36	ARG
52	DS	38	SER
52	DS	43	GLU
52	DS	78	ARG
53	DT	8	ARG
53	DT	23	ARG
53	DT	56	MET
53	DT	61	SER
53	DT	72	LEU
53	DT	79	ARG
53	DT	80	ARG
53	DT	84	LEU
57	DZ	-58	LEU
57	DZ	-52	VAL
57	DZ	-44	PRO
57	DZ	-32	LEU
57	DZ	-30	VAL
57	DZ	-29	LEU
57	DZ	-22	LYS
57	DZ	-6	ARG
57	DZ	-3	GLU
57	DZ	6	GLU
57	DZ	9	LEU
57	DZ	10	LYS
57	DZ	14	ASN
57	DZ	15	ILE
57	DZ	21	ILE
57	DZ	22	ASP
57	DZ	28	THR
57	DZ	36	THR
57	DZ	75	LYS
57	DZ	76	ASP
57	DZ	79	ILE
57	DZ	81	ILE
57	DZ	87	HIS
57	DZ	92	ILE
57	DZ	100	VAL
57	DZ	101	LEU
57	DZ	105	ILE

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Mol	Chain	Res	Type
57	DZ	111	SER
57	DZ	114	VAL
57	DZ	117	GLN
57	DZ	118	SER
57	DZ	123	ARG
57	DZ	130	VAL
57	DZ	132	ARG
57	DZ	133	ILE
57	DZ	142	THR
57	DZ	146	LEU
57	DZ	157	LEU
57	DZ	165	GLN
57	DZ	170	ARG
57	DZ	181	LEU
57	DZ	182	ARG
57	DZ	184	LYS
57	DZ	186	TYR
57	DZ	187	THR
57	DZ	192	LEU
57	DZ	196	ILE
57	DZ	201	ILE
57	DZ	203	GLU
57	DZ	207	ASP
57	DZ	212	TYR
57	DZ	213	HIS
57	DZ	215	LYS
57	DZ	217	VAL
57	DZ	219	VAL
57	DZ	222	ASP
57	DZ	225	GLU
57	DZ	227	ILE
57	DZ	228	MET
57	DZ	229	LEU
57	DZ	232	LEU
57	DZ	236	GLU
57	DZ	240	GLU
57	DZ	242	LEU
57	DZ	255	ILE
57	DZ	258	VAL
57	DZ	262	SER
57	DZ	279	TYR
57	DZ	284	LEU

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Mol	Chain	Res	Type
57	DZ	285	ASP
57	DZ	292	THR
57	DZ	297	GLU
57	DZ	298	VAL
57	DZ	299	VAL
57	DZ	302	HIS
57	DZ	312	LEU
57	DZ	322	VAL
57	DZ	328	ILE
57	DZ	345	THR
57	DZ	352	VAL
57	DZ	355	LEU
57	DZ	357	ARG
57	DZ	361	ASN
57	DZ	363	ARG
57	DZ	364	GLU
57	DZ	368	GLU
57	DZ	377	VAL
57	DZ	385	THR
57	DZ	389	LEU
57	DZ	399	LEU
57	DZ	402	ILE
57	DZ	403	GLU
57	DZ	408	VAL
57	DZ	417	THR
57	DZ	421	GLN
57	DZ	424	LEU
57	DZ	428	LEU
57	DZ	434	GLU
57	DZ	440	VAL
57	DZ	445	GLU
57	DZ	451	ILE
57	DZ	452	SER
57	DZ	457	LEU
57	DZ	461	ILE
57	DZ	466	LEU
57	DZ	468	ARG
57	DZ	480	GLN
57	DZ	487	ILE
57	DZ	494	GLU
57	DZ	506	GLN
57	DZ	512	ILE

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Mol	Chain	Res	Type
57	DZ	523	PHE
57	DZ	542	VAL
57	DZ	561	VAL
57	DZ	571	SER
57	DZ	576	ASP
57	DZ	590	ILE
57	DZ	610	VAL
57	DZ	612	THR
57	DZ	614	GLU
57	DZ	615	GLU
57	DZ	621	ILE
57	DZ	623	ASP
57	DZ	625	ASN
57	DZ	634	MET
57	DZ	639	ASN
57	DZ	643	ILE
57	DZ	649	LEU
57	DZ	651	GLU
57	DZ	657	THR
57	DZ	659	LEU
57	DZ	661	SER
57	DZ	669	PHE
57	DZ	670	VAL
57	DZ	674	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (175) such sidechains are listed below:

Mol	Chain	Res	Type
3	AC	45	HIS
3	AC	67	HIS
3	AC	173	HIS
3	AC	189	ASN
3	AC	200	HIS
4	AD	253	GLN
5	AE	54	GLN
5	AE	85	ASN
6	AF	69	HIS
6	AF	203	GLN
7	AG	26	GLN
10	AL	29	GLN
10	AL	47	ASN
13	AP	38	GLN

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Mol	Chain	Res	Type
17	AT	43	GLN
17	AT	123	GLN
18	AU	81	HIS
18	AU	104	GLN
21	AX	31	HIS
21	AX	82	GLN
22	AY	92	ASN
23	AZ	34	ASN
23	AZ	121	HIS
24	A0	35	ASN
26	A2	38	GLN
27	A3	32	GLN
28	A4	46	GLN
32	A8	35	GLN
33	A9	36	GLN
35	BB	40	HIS
35	BB	76	GLN
36	BC	28	GLN
36	BC	37	GLN
36	BC	69	HIS
36	BC	98	ASN
36	BC	136	GLN
36	BC	181	ASN
37	BD	77	ASN
37	BD	119	GLN
37	BD	123	HIS
37	BD	125	HIS
37	BD	161	ASN
38	BE	38	GLN
38	BE	56	GLN
38	BE	73	ASN
38	BE	141	GLN
39	BF	73	ASN
40	BG	28	ASN
40	BG	86	GLN
42	BI	23	ASN
42	BI	31	GLN
42	BI	73	GLN
42	BI	89	ASN
42	BI	124	GLN
43	BJ	21	GLN
43	BJ	56	HIS

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Mol	Chain	Res	Type
44	BK	93	GLN
44	BK	104	GLN
45	BL	99	HIS
46	BM	12	ASN
47	BN	49	HIS
48	BO	9	GLN
48	BO	13	GLN
48	BO	28	GLN
48	BO	62	GLN
50	BQ	16	GLN
52	BS	65	ASN
52	BS	69	HIS
52	BS	83	HIS
53	BT	26	ASN
53	BT	90	GLN
57	BZ	-50	GLN
57	BZ	-24	ASN
57	BZ	165	GLN
57	BZ	213	HIS
57	BZ	270	GLN
57	BZ	361	ASN
57	BZ	426	GLN
57	BZ	448	GLN
57	BZ	475	ASN
57	BZ	509	HIS
57	BZ	573	HIS
57	BZ	641	GLN
57	BZ	675	HIS
3	CC	67	HIS
3	CC	189	ASN
3	CC	200	HIS
4	CD	87	ASN
4	CD	126	GLN
4	CD	253	GLN
5	CE	85	ASN
5	CE	143	ASN
5	CE	169	ASN
6	CF	29	ASN
6	CF	69	HIS
6	CF	75	HIS
6	CF	133	ASN
6	CF	169	ASN

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Mol	Chain	Res	Type
6	CF	203	GLN
7	CG	40	ASN
7	CG	108	ASN
8	CH	158	HIS
10	CL	30	HIS
10	CL	42	ASN
10	CL	103	GLN
13	CP	38	GLN
14	CQ	57	HIS
14	CQ	89	ASN
15	CR	50	HIS
16	CS	68	GLN
17	CT	43	GLN
17	CT	58	ASN
17	CT	123	GLN
18	CU	72	HIS
19	CV	64	HIS
21	CX	31	HIS
21	CX	82	GLN
22	CY	43	ASN
23	CZ	34	ASN
23	CZ	50	GLN
23	CZ	55	HIS
23	CZ	65	GLN
24	C0	29	GLN
24	C0	50	ASN
26	C2	46	GLN
33	C9	36	GLN
35	DB	45	GLN
35	DB	76	GLN
35	DB	224	GLN
36	DC	104	GLN
36	DC	110	ASN
36	DC	118	GLN
36	DC	123	GLN
37	DD	45	GLN
37	DD	77	ASN
37	DD	123	HIS
37	DD	125	HIS
37	DD	129	ASN
37	DD	160	GLN
38	DE	20	GLN

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Mol	Chain	Res	Type
38	DE	73	ASN
38	DE	130	ASN
39	DF	100	ASN
40	DG	28	ASN
40	DG	51	GLN
40	DG	97	GLN
42	DI	31	GLN
43	DJ	21	GLN
43	DJ	68	HIS
44	DK	93	GLN
44	DK	117	ASN
45	DL	49	ASN
45	DL	75	HIS
45	DL	78	GLN
45	DL	99	HIS
46	DM	77	ASN
46	DM	106	ASN
48	DO	13	GLN
48	DO	28	GLN
50	DQ	16	GLN
52	DS	83	HIS
53	DT	16	HIS
57	DZ	-50	GLN
57	DZ	-24	ASN
57	DZ	77	HIS
57	DZ	117	GLN
57	DZ	266	ASN
57	DZ	270	GLN
57	DZ	302	HIS
57	DZ	361	ASN
57	DZ	362	HIS
57	DZ	421	GLN
57	DZ	480	GLN
57	DZ	506	GLN
57	DZ	641	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	AA	2845/2915 (97%)	527 (18%)	56 (1%)
1	CA	2839/2915 (97%)	579 (20%)	39 (1%)

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	AB	119/121 (98%)	23 (19%)	0
2	CB	119/121 (98%)	21 (17%)	0
34	BA	1491/1521 (98%)	310 (20%)	22 (1%)
34	DA	1498/1521 (98%)	303 (20%)	24 (1%)
55	BV	6/18 (33%)	1 (16%)	0
55	DV	5/18 (27%)	1 (20%)	0
56	BW	74/76 (97%)	15 (20%)	0
56	BY	71/76 (93%)	23 (32%)	2 (2%)
56	DW	74/76 (97%)	19 (25%)	2 (2%)
56	DY	69/76 (90%)	21 (30%)	1 (1%)
All	All	9210/9454 (97%)	1843 (20%)	146 (1%)

All (1843) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	AA	9	U
1	AA	12	U
1	AA	13	A
1	AA	34	C
1	AA	45	C
1	AA	56	C
1	AA	62	U
1	AA	68	C
1	AA	70	A
1	AA	73	A
1	AA	74	G
1	AA	77	A
1	AA	83	A
1	AA	95	G
1	AA	116	A
1	AA	117	A
1	AA	118	U
1	AA	119	G
1	AA	120	G
1	AA	123	G
1	AA	129	G
1	AA	131	C
1	AA	146	G
1	AA	149	A
1	AA	170	A
1	AA	171	A
1	AA	185	A

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Mol	Chain	Res	Type
1	AA	188	A
1	AA	189	U
1	AA	190	C
1	AA	194	G
1	AA	202	A
1	AA	203	G
1	AA	204	G
1	AA	205	A
1	AA	208	G
1	AA	211	A
1	AA	214	A
1	AA	217	A
1	AA	218	A
1	AA	222	A
1	AA	237	G
1	AA	239	G
1	AA	254	A
1	AA	265	U
1	AA	269	G
1	AA	271	U
1	AA	272	U
1	AA	273	G
1	AA	274	U
1	AA	275	C
1	AA	276	C
1	AA	279	G
1	AA	289	G
1	AA	295	C
1	AA	296	U
1	AA	299	G
1	AA	303	C
1	AA	311	C
1	AA	334	A
1	AA	335	A
1	AA	348	A
1	AA	349	G
1	AA	353	G
1	AA	354	A
1	AA	358	C
1	AA	376	G
1	AA	387	G
1	AA	397	G

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Mol	Chain	Res	Type
1	AA	407	U
1	AA	413	G
1	AA	423	G
1	AA	431	C
1	AA	434	G
1	AA	438	G
1	AA	439	A
1	AA	455	A
1	AA	464	G
1	AA	469	A
1	AA	470	C
1	AA	474	U
1	AA	477	C
1	AA	482	C
1	AA	483	A
1	AA	496	A
1	AA	501	U
1	AA	502	G
1	AA	507	G
1	AA	511	C
1	AA	512	C
1	AA	519	G
1	AA	529	U
1	AA	530	A
1	AA	534	C
1	AA	535	C
1	AA	543	G
1	AA	553	A
1	AA	554	A
1	AA	555	G
1	AA	556	C
1	AA	557	A
1	AA	558	G
1	AA	573	G
1	AA	574	G
1	AA	579	G
1	AA	586	G
1	AA	596	G
1	AA	597	C
1	AA	598	A
1	AA	609	A
1	AA	610	C

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Mol	Chain	Res	Type
1	AA	626	A
1	AA	627	G
1	AA	630	U
1	AA	638	U
1	AA	639	G
1	AA	641	G
1	AA	659	C
1	AA	662	A
1	AA	671	A
1	AA	697	C
1	AA	698	G
1	AA	702	A
1	AA	703	G
1	AA	716	G
1	AA	724	A
1	AA	725	C
1	AA	733	G
1	AA	777	C
1	AA	787	U
1	AA	811	A
1	AA	812	G
1	AA	821	A
1	AA	822	G
1	AA	823	G
1	AA	829	A
1	AA	831	A
1	AA	832	G
1	AA	837	C
1	AA	839	G
1	AA	852	G
1	AA	859	C
1	AA	874	U
1	AA	875	U
1	AA	906	G
1	AA	913	A
1	AA	914	C
1	AA	926	G
1	AA	927	G
1	AA	932	C
1	AA	933	C
1	AA	934	A
1	AA	935	C

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Mol	Chain	Res	Type
1	AA	936	C
1	AA	937	A
1	AA	938	G
1	AA	939	C
1	AA	940	C
1	AA	942	A
1	AA	943	C
1	AA	945	A
1	AA	953	U
1	AA	954	C
1	AA	956	A
1	AA	967	G
1	AA	968	U
1	AA	977	G
1	AA	986	A
1	AA	990	A
1	AA	991	G
1	AA	992	G
1	AA	1004	A
1	AA	1006	C
1	AA	1019	G
1	AA	1020	C
1	AA	1029	A
1	AA	1036	A
1	AA	1040	C
1	AA	1042	A
1	AA	1048	G
1	AA	1051	C
1	AA	1058	U
1	AA	1059	C
1	AA	1068	G
1	AA	1072	U
1	AA	1073	A
1	AA	1079	U
1	AA	1084	C
1	AA	1087	C
1	AA	1093	G
1	AA	1097	G
1	AA	1099	C
1	AA	1100	A
1	AA	1105	G
1	AA	1107	U

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Mol	Chain	Res	Type
1	AA	1108	G
1	AA	1112	U
1	AA	1116	A
1	AA	1119	A
1	AA	1120	G
1	AA	1121	C
1	AA	1122	C
1	AA	1128	U
1	AA	1129	U
1	AA	1132	A
1	AA	1134	A
1	AA	1135	G
1	AA	1141	A
1	AA	1147	U
1	AA	1155	C
1	AA	1156	G
1	AA	1161	G
1	AA	1195	G
1	AA	1217	G
1	AA	1218	G
1	AA	1219	A
1	AA	1220	U
1	AA	1221	G
1	AA	1222	A
1	AA	1223	C
1	AA	1255	A
1	AA	1256	U
1	AA	1265	A
1	AA	1299	A
1	AA	1302	G
1	AA	1317	G
1	AA	1318	A
1	AA	1319	U
1	AA	1321	A
1	AA	1346	U
1	AA	1347	A
1	AA	1349	G
1	AA	1359	U
1	AA	1367	A
1	AA	1398	U
1	AA	1405	A
1	AA	1406	A

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Mol	Chain	Res	Type
1	AA	1411	A
1	AA	1416	C
1	AA	1422	C
1	AA	1423	G
1	AA	1424	A
1	AA	1426	G
1	AA	1430	A
1	AA	1431	G
1	AA	1432	C
1	AA	1462	G
1	AA	1463	C
1	AA	1466	U
1	AA	1467	G
1	AA	1474	C
1	AA	1477	U
1	AA	1491	A
1	AA	1493	C
1	AA	1496	A
1	AA	1497	G
1	AA	1503	G
1	AA	1506	G
1	AA	1514	C
1	AA	1518	A
1	AA	1529	G
1	AA	1539	C
1	AA	1554	A
1	AA	1555	C
1	AA	1556	A
1	AA	1569	U
1	AA	1578	C
1	AA	1589	A
1	AA	1590	C
1	AA	1605	A
1	AA	1607	G
1	AA	1613	A
1	AA	1616	A
1	AA	1625	U
1	AA	1628	G
1	AA	1629	C
1	AA	1631	C
1	AA	1632	A
1	AA	1652	G

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Mol	Chain	Res	Type
1	AA	1653	C
1	AA	1654	A
1	AA	1655	A
1	AA	1656	A
1	AA	1663	C
1	AA	1671	C
1	AA	1688	A
1	AA	1694	G
1	AA	1695	C
1	AA	1700	G
1	AA	1701	A
1	AA	1721	G
1	AA	1743	G
1	AA	1746	G
1	AA	1747	A
1	AA	1748	A
1	AA	1765	U
1	AA	1766	G
1	AA	1767	A
1	AA	1768	U
1	AA	1769	G
1	AA	1776	G
1	AA	1777	G
1	AA	1787	G
1	AA	1793	A
1	AA	1794	G
1	AA	1795	G
1	AA	1804	A
1	AA	1811	A
1	AA	1812	C
1	AA	1813	C
1	AA	1822	A
1	AA	1831	C
1	AA	1832	G
1	AA	1833	A
1	AA	1845	G
1	AA	1847	G
1	AA	1848	G
1	AA	1851	U
1	AA	1859	G
1	AA	1860	A
1	AA	1870	G

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Mol	Chain	Res	Type
1	AA	1878	A
1	AA	1886	G
1	AA	1889	G
1	AA	1892	G
1	AA	1898	A
1	AA	1911	A
1	AA	1922	A
1	AA	1928	G
1	AA	1937	U
1	AA	1951	G
1	AA	1952	G
1	AA	1953	U
1	AA	1954	A
1	AA	1960	A
1	AA	1966	U
1	AA	1977	U
1	AA	1985	U
1	AA	1986	G
1	AA	1987	C
1	AA	1989	C
1	AA	1992	A
1	AA	1993	A
1	AA	1994	A
1	AA	2010	C
1	AA	2014	G
1	AA	2015	U
1	AA	2018	C
1	AA	2019	G
1	AA	2045	G
1	AA	2052	A
1	AA	2053	A
1	AA	2055	A
1	AA	2061	C
1	AA	2064	A
1	AA	2065	C
1	AA	2071	G
1	AA	2077	C
1	AA	2078	G
1	AA	2082	A
1	AA	2083	G
1	AA	2084	A
1	AA	2091	G

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Mol	Chain	Res	Type
1	AA	2102	G
1	AA	2104	A
1	AA	2122	G
1	AA	2133	C
1	AA	2134	G
1	AA	2135	U
1	AA	2139	A
1	AA	2141	A
1	AA	2149	G
1	AA	2151	C
1	AA	2154	U
1	AA	2155	G
1	AA	2156	A
1	AA	2157	A
1	AA	2158	C
1	AA	2160	C
1	AA	2161	C
1	AA	2162	C
1	AA	2163	G
1	AA	2164	C
1	AA	2167	C
1	AA	2168	C
1	AA	2169	G
1	AA	2172	U
1	AA	2173	G
1	AA	2175	G
1	AA	2178	G
1	AA	2179	G
1	AA	2180	A
1	AA	2181	G
1	AA	2182	G
1	AA	2188	G
1	AA	2189	U
1	AA	2190	G
1	AA	2191	A
1	AA	2192	A
1	AA	2194	U
1	AA	2195	A
1	AA	2197	C
1	AA	2204	G
1	AA	2206	G
1	AA	2210	C

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Mol	Chain	Res	Type
1	AA	2211	U
1	AA	2213	G
1	AA	2214	G
1	AA	2217	C
1	AA	2218	C
1	AA	2220	A
1	AA	2221	A
1	AA	2227	G
1	AA	2228	G
1	AA	2229	A
1	AA	2237	A
1	AA	2280	A
1	AA	2281	A
1	AA	2285	A
1	AA	2287	C
1	AA	2291	G
1	AA	2295	C
1	AA	2298	A
1	AA	2299	A
1	AA	2301	G
1	AA	2307	C
1	AA	2317	A
1	AA	2320	G
1	AA	2324	U
1	AA	2327	G
1	AA	2332	A
1	AA	2337	G
1	AA	2346	G
1	AA	2347	A
1	AA	2348	A
1	AA	2352	G
1	AA	2355	C
1	AA	2356	U
1	AA	2359	C
1	AA	2361	G
1	AA	2362	C
1	AA	2369	U
1	AA	2376	C
1	AA	2377	G
1	AA	2395	G
1	AA	2397	C
1	AA	2418	U

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Mol	Chain	Res	Type
1	AA	2419	G
1	AA	2426	G
1	AA	2434	A
1	AA	2435	U
1	AA	2437	A
1	AA	2440	G
1	AA	2441	G
1	AA	2442	A
1	AA	2445	A
1	AA	2446	A
1	AA	2451	A
1	AA	2453	C
1	AA	2460	A
1	AA	2461	U
1	AA	2482	G
1	AA	2488	A
1	AA	2490	A
1	AA	2503	U
1	AA	2514	G
1	AA	2517	G
1	AA	2518	U
1	AA	2519	C
1	AA	2530	A
1	AA	2541	G
1	AA	2547	G
1	AA	2566	U
1	AA	2567	U
1	AA	2578	A
1	AA	2579	G
1	AA	2584	A
1	AA	2585	C
1	AA	2596	U
1	AA	2597	U
1	AA	2600	G
1	AA	2613	C
1	AA	2614	A
1	AA	2621	U
1	AA	2622	C
1	AA	2623	U
1	AA	2624	C
1	AA	2642	G
1	AA	2653	G

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Mol	Chain	Res	Type
1	AA	2670	C
1	AA	2674	A
1	AA	2681	G
1	AA	2691	A
1	AA	2701	U
1	AA	2702	C
1	AA	2715	C
1	AA	2725	A
1	AA	2726	A
1	AA	2727	G
1	AA	2730	G
1	AA	2738	A
1	AA	2739	U
1	AA	2746	A
1	AA	2757	G
1	AA	2770	A
1	AA	2771	A
1	AA	2778	A
1	AA	2779	G
1	AA	2782	C
1	AA	2791	A
1	AA	2803	A
1	AA	2804	C
1	AA	2807	C
1	AA	2813	G
1	AA	2814	C
1	AA	2818	U
1	AA	2830	A
1	AA	2831	A
1	AA	2843	G
1	AA	2845	A
1	AA	2876	U
1	AA	2882	G
1	AA	2883	A
1	AA	2892	A
1	AA	2902	G
1	AA	2903	G
1	AA	2906	U
2	AB	2	C
2	AB	5	C
2	AB	10	C
2	AB	12	C

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Mol	Chain	Res	Type
2	AB	13	A
2	AB	15	A
2	AB	47	C
2	AB	50	G
2	AB	53	A
2	AB	54	G
2	AB	56	G
2	AB	57	A
2	AB	58	A
2	AB	59	A
2	AB	67	G
2	AB	73	A
2	AB	76	G
2	AB	89	G
2	AB	90	A
2	AB	95	C
2	AB	110	G
2	AB	111	G
2	AB	113	G
34	BA	5	U
34	BA	7	G
34	BA	9	G
34	BA	10	A
34	BA	11	G
34	BA	32	A
34	BA	39	G
34	BA	43	C
34	BA	44	G
34	BA	47	C
34	BA	48	C
34	BA	50	A
34	BA	51	A
34	BA	61	G
34	BA	71	C
34	BA	77	G
34	BA	78	G
34	BA	79	G
34	BA	97	G
34	BA	101	A
34	BA	105	G
34	BA	116	A
34	BA	121	C

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Mol	Chain	Res	Type
34	BA	127	G
34	BA	131	C
34	BA	137	C
34	BA	163	C
34	BA	174	C
34	BA	182	U
34	BA	189(F)	U
34	BA	189(G)	G
34	BA	189(J)	G
34	BA	194	C
34	BA	195	A
34	BA	197	A
34	BA	199	G
34	BA	202	U
34	BA	204	U
34	BA	216	G
34	BA	220	G
34	BA	231	G
34	BA	236	G
34	BA	240	C
34	BA	243	A
34	BA	244	U
34	BA	247	G
34	BA	251	G
34	BA	259	G
34	BA	261	U
34	BA	266	G
34	BA	267	C
34	BA	270	A
34	BA	289	G
34	BA	298	A
34	BA	318	G
34	BA	321	A
34	BA	328	C
34	BA	329	A
34	BA	331	G
34	BA	332	G
34	BA	347	G
34	BA	348	G
34	BA	350	G
34	BA	351	G
34	BA	352	C

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Mol	Chain	Res	Type
34	BA	353	A
34	BA	354	G
34	BA	355	C
34	BA	359	U
34	BA	367	U
34	BA	372	C
34	BA	384	G
34	BA	397	A
34	BA	398	C
34	BA	411	A
34	BA	412	A
34	BA	413	G
34	BA	422	C
34	BA	424	G
34	BA	429	U
34	BA	432	A
34	BA	439	A
34	BA	442	C
34	BA	443	C
34	BA	446	G
34	BA	452	A
34	BA	470	C
34	BA	484	G
34	BA	485	G
34	BA	496	A
34	BA	498	U
34	BA	499	A
34	BA	504	C
34	BA	505	G
34	BA	509	A
34	BA	510	A
34	BA	511	C
34	BA	514	C
34	BA	516	U
34	BA	518	C
34	BA	521	G
34	BA	527	G
34	BA	531	U
34	BA	532	A
34	BA	533	A
34	BA	536	C
34	BA	547	A

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Mol	Chain	Res	Type
34	BA	558	G
34	BA	559	A
34	BA	560	U
34	BA	561	U
34	BA	564	C
34	BA	570	G
34	BA	572	A
34	BA	573	A
34	BA	576	G
34	BA	581	G
34	BA	618	C
34	BA	630	G
34	BA	631	G
34	BA	633	G
34	BA	649	G
34	BA	653	A
34	BA	665	A
34	BA	687	A
34	BA	688	G
34	BA	693	G
34	BA	694	A
34	BA	697	U
34	BA	705	U
34	BA	714	G
34	BA	717	C
34	BA	723	U
34	BA	724	G
34	BA	728	A
34	BA	731	G
34	BA	734	G
34	BA	738	C
34	BA	747	C
34	BA	749	C
34	BA	752	G
34	BA	755	G
34	BA	774	G
34	BA	777	A
34	BA	792	A
34	BA	793	U
34	BA	794	A
34	BA	815	A
34	BA	816	A

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Mol	Chain	Res	Type
34	BA	817	C
34	BA	828	A
34	BA	829	G
34	BA	840	C
34	BA	841	U
34	BA	848	C
34	BA	851	G
34	BA	859	A
34	BA	870	U
34	BA	872	A
34	BA	874	G
34	BA	902	G
34	BA	905	U
34	BA	910	C
34	BA	913	A
34	BA	914	A
34	BA	922	G
34	BA	923	A
34	BA	926	G
34	BA	927	G
34	BA	934	C
34	BA	936	C
34	BA	941	G
34	BA	942	G
34	BA	960	U
34	BA	961	U
34	BA	964	A
34	BA	968	A
34	BA	969	A
34	BA	971	G
34	BA	973	G
34	BA	974	A
34	BA	975	A
34	BA	976	G
34	BA	977	A
34	BA	992	U
34	BA	993	G
34	BA	1003	G
34	BA	1004	A
34	BA	1005	A
34	BA	1009	G
34	BA	1016	A

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Mol	Chain	Res	Type
34	BA	1019	C
34	BA	1021	G
34	BA	1022	G
34	BA	1023	G
34	BA	1025	U
34	BA	1026	G
34	BA	1028	C
34	BA	1029	C
34	BA	1030	C
34	BA	1030(A)	G
34	BA	1030(C)	G
34	BA	1031	G
34	BA	1033	G
34	BA	1042	G
34	BA	1045	C
34	BA	1053	G
34	BA	1054	C
34	BA	1055	A
34	BA	1056	U
34	BA	1063	C
34	BA	1065	U
34	BA	1066	C
34	BA	1067	A
34	BA	1068	G
34	BA	1070	U
34	BA	1081	G
34	BA	1086	U
34	BA	1094	G
34	BA	1095	U
34	BA	1101	A
34	BA	1108	G
34	BA	1123	A
34	BA	1124	G
34	BA	1125	U
34	BA	1126	U
34	BA	1130	A
34	BA	1136	U
34	BA	1137	C
34	BA	1139	G
34	BA	1140	C
34	BA	1141	C
34	BA	1146	A

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Mol	Chain	Res	Type
34	BA	1150	U
34	BA	1152	A
34	BA	1157	A
34	BA	1159	U
34	BA	1165	C
34	BA	1166	G
34	BA	1182	G
34	BA	1183	A
34	BA	1184	G
34	BA	1189	C
34	BA	1190	G
34	BA	1196	U
34	BA	1197	G
34	BA	1200	C
34	BA	1201	A
34	BA	1202	G
34	BA	1213	A
34	BA	1214	C
34	BA	1227	A
34	BA	1228	C
34	BA	1236	A
34	BA	1238	A
34	BA	1240	U
34	BA	1241	G
34	BA	1253	G
34	BA	1256	A
34	BA	1257	U
34	BA	1258	G
34	BA	1260	C
34	BA	1278	U
34	BA	1279	A
34	BA	1280	A
34	BA	1286	A
34	BA	1287	A
34	BA	1299	A
34	BA	1300	G
34	BA	1305	G
34	BA	1320	C
34	BA	1322	C
34	BA	1338	G
34	BA	1340	A
34	BA	1346	A

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Mol	Chain	Res	Type
34	BA	1347	G
34	BA	1353	G
34	BA	1359	C
34	BA	1360	A
34	BA	1363	C
34	BA	1363(A)	A
34	BA	1379	G
34	BA	1389	C
34	BA	1397	C
34	BA	1398	A
34	BA	1400	C
34	BA	1402	C
34	BA	1419	G
34	BA	1442	G
34	BA	1442(A)	G
34	BA	1446	U
34	BA	1447	A
34	BA	1452	C
34	BA	1460	A
34	BA	1468	A
34	BA	1471	G
34	BA	1479	C
34	BA	1487	G
34	BA	1491	G
34	BA	1496	C
34	BA	1497	G
34	BA	1502	A
34	BA	1503	A
34	BA	1504	G
34	BA	1506	U
34	BA	1507	A
34	BA	1517	G
34	BA	1529	G
34	BA	1530	G
34	BA	1531	A
55	BV	13	A
56	BW	8	4SU
56	BW	13	C
56	BW	14	A
56	BW	17	C
56	BW	18	G
56	BW	20	U

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Mol	Chain	Res	Type
56	BW	22	G
56	BW	31	A
56	BW	42	C
56	BW	43	C
56	BW	45	U
56	BW	47	U
56	BW	48	C
56	BW	72	C
56	BW	76	A
56	BY	5	G
56	BY	6	G
56	BY	9	A
56	BY	13	C
56	BY	14	A
56	BY	20	U
56	BY	21	A
56	BY	23	A
56	BY	26	A
56	BY	34	G
56	BY	36	A
56	BY	41	C
56	BY	42	C
56	BY	44	G
56	BY	45	U
56	BY	46	7MG
56	BY	47	U
56	BY	48	C
56	BY	49	C
56	BY	57	G
56	BY	59	U
56	BY	60	U
56	BY	68	C
1	CA	10	G
1	CA	12	U
1	CA	14	A
1	CA	15	G
1	CA	34	C
1	CA	35	G
1	CA	36	G
1	CA	45	C
1	CA	51	G
1	CA	54	G

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Mol	Chain	Res	Type
1	CA	55	G
1	CA	71	A
1	CA	74	A
1	CA	75	G
1	CA	78	A
1	CA	84	A
1	CA	90	U
1	CA	95	G
1	CA	100	G
1	CA	102	G
1	CA	118	A
1	CA	119	A
1	CA	120	U
1	CA	133	C
1	CA	139(A)	G
1	CA	141	A
1	CA	149	A
1	CA	154(A)	C
1	CA	157	U
1	CA	181	A
1	CA	196	A
1	CA	197	A
1	CA	199	A
1	CA	205	G
1	CA	214	G
1	CA	215	G
1	CA	216	A
1	CA	221	A
1	CA	222	A
1	CA	225	A
1	CA	226	G
1	CA	228	A
1	CA	229	A
1	CA	232	G
1	CA	248	G
1	CA	271(A)	A
1	CA	271(I)	G
1	CA	271(K)	U
1	CA	271(L)	U
1	CA	271(M)	G
1	CA	271(P)	C
1	CA	272(A)	U

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Mol	Chain	Res	Type
1	CA	272(B)	G
1	CA	272(J)	C
1	CA	277	C
1	CA	278	A
1	CA	283	A
1	CA	286	C
1	CA	289	A
1	CA	298	G
1	CA	299	A
1	CA	310	A
1	CA	311	A
1	CA	317	G
1	CA	324	A
1	CA	327	G
1	CA	329	G
1	CA	330	A
1	CA	333	G
1	CA	338	G
1	CA	339	U
1	CA	352	G
1	CA	362	U
1	CA	363	G
1	CA	363(C)	G
1	CA	363(E)	U
1	CA	386	G
1	CA	407	G
1	CA	411	G
1	CA	412	A
1	CA	416	C
1	CA	422	A
1	CA	428	A
1	CA	443	A
1	CA	444	C
1	CA	451	C
1	CA	455	C
1	CA	456	C
1	CA	457	A
1	CA	470	A
1	CA	471	A
1	CA	475	U
1	CA	477	A
1	CA	481	G

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Mol	Chain	Res	Type
1	CA	504	U
1	CA	505	A
1	CA	507	A
1	CA	508	G
1	CA	509	C
1	CA	521	G
1	CA	528	A
1	CA	529	A
1	CA	530	G
1	CA	531	C
1	CA	532	A
1	CA	533	G
1	CA	545	G
1	CA	557	U
1	CA	563	G
1	CA	573	G
1	CA	575	A
1	CA	586	A
1	CA	588	U
1	CA	592	G
1	CA	603	A
1	CA	604	G
1	CA	606	U
1	CA	607	U
1	CA	614(A)	U
1	CA	614(B)	G
1	CA	614(C)	A
1	CA	615	G
1	CA	620	G
1	CA	623	G
1	CA	627	A
1	CA	631	A
1	CA	637	A
1	CA	645	C
1	CA	646	A
1	CA	652(B)	A
1	CA	652(C)	G
1	CA	652(U)	G
1	CA	653	A
1	CA	669	G
1	CA	686	G
1	CA	730	C

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Mol	Chain	Res	Type
1	CA	732	C
1	CA	743	G
1	CA	762	U
1	CA	764	A
1	CA	765	G
1	CA	774	A
1	CA	775	G
1	CA	776	G
1	CA	782	A
1	CA	784	A
1	CA	785	G
1	CA	792	G
1	CA	794	G
1	CA	802	A
1	CA	805	G
1	CA	812	C
1	CA	819	A
1	CA	822	U
1	CA	827	U
1	CA	828	U
1	CA	831	G
1	CA	847	U
1	CA	857	C
1	CA	859	G
1	CA	866	A
1	CA	879	G
1	CA	880	G
1	CA	884	C
1	CA	886	C
1	CA	887	A
1	CA	888	C
1	CA	889	C
1	CA	890	A
1	CA	893	C
1	CA	896	A
1	CA	897	C
1	CA	898	C
1	CA	900	A
1	CA	901	A
1	CA	907	U
1	CA	910	A
1	CA	914	C

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Mol	Chain	Res	Type
1	CA	915	C
1	CA	917	A
1	CA	926	A
1	CA	932	G
1	CA	938	G
1	CA	941	A
1	CA	944	G
1	CA	945	A
1	CA	946	G
1	CA	958	U
1	CA	959	A
1	CA	961	C
1	CA	974	G
1	CA	975	C
1	CA	983	A
1	CA	996	A
1	CA	1005	C
1	CA	1009	A
1	CA	1012	U
1	CA	1013	C
1	CA	1018	C
1	CA	1020	A
1	CA	1022	G
1	CA	1025	G
1	CA	1027	A
1	CA	1033	U
1	CA	1038	C
1	CA	1039	G
1	CA	1041	C
1	CA	1042	G
1	CA	1046	A
1	CA	1047	G
1	CA	1048	A
1	CA	1049	C
1	CA	1050	A
1	CA	1055	G
1	CA	1057	A
1	CA	1058	G
1	CA	1060	U
1	CA	1061	U
1	CA	1062	G
1	CA	1070	A

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Mol	Chain	Res	Type
1	CA	1073	A
1	CA	1075	C
1	CA	1076	C
1	CA	1079	C
1	CA	1083	U
1	CA	1088	A
1	CA	1090	U
1	CA	1101	U
1	CA	1106	G
1	CA	1110	G
1	CA	1111	A
1	CA	1112	G
1	CA	1114	G
1	CA	1116	C
1	CA	1144	G
1	CA	1149	G
1	CA	1171	G
1	CA	1180	C
1	CA	1188	U
1	CA	1205	U
1	CA	1220	A
1	CA	1221	C
1	CA	1223	G
1	CA	1229	G
1	CA	1236	G
1	CA	1241	A
1	CA	1248	G
1	CA	1250	G
1	CA	1253	A
1	CA	1256	G
1	CA	1268	A
1	CA	1269	A
1	CA	1271	G
1	CA	1272	A
1	CA	1287	A
1	CA	1300	U
1	CA	1301	A
1	CA	1305	C
1	CA	1314	C
1	CA	1321	A
1	CA	1342	A
1	CA	1348	G

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Mol	Chain	Res	Type
1	CA	1351	C
1	CA	1359	A
1	CA	1360	A
1	CA	1365	A
1	CA	1368	G
1	CA	1370	C
1	CA	1384	A
1	CA	1385	G
1	CA	1388	G
1	CA	1390	U
1	CA	1416	G
1	CA	1417	C
1	CA	1419	A
1	CA	1420	U
1	CA	1421	G
1	CA	1422	G
1	CA	1426	G
1	CA	1427	A
1	CA	1428	C
1	CA	1437	C
1	CA	1445	A
1	CA	1449	A
1	CA	1450	G
1	CA	1452	A
1	CA	1455	G
1	CA	1459	G
1	CA	1460	A
1	CA	1461	G
1	CA	1467	C
1	CA	1471	A
1	CA	1473	G
1	CA	1482	G
1	CA	1489	U
1	CA	1490	A
1	CA	1493	C
1	CA	1496	A
1	CA	1497	U
1	CA	1509	C
1	CA	1509(A)	A
1	CA	1525	G
1	CA	1531	C
1	CA	1542	A

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Mol	Chain	Res	Type
1	CA	1543	C
1	CA	1544	A
1	CA	1547	C
1	CA	1558	A
1	CA	1559	G
1	CA	1560	G
1	CA	1566	A
1	CA	1569	A
1	CA	1578	U
1	CA	1580	A
1	CA	1583	A
1	CA	1584	C
1	CA	1586	A
1	CA	1608	A
1	CA	1609	A
1	CA	1612	C
1	CA	1640	C
1	CA	1647	G
1	CA	1648	C
1	CA	1653	G
1	CA	1654	A
1	CA	1670	C
1	CA	1674	G
1	CA	1675	C
1	CA	1682	G
1	CA	1696	G
1	CA	1700	A
1	CA	1701	A
1	CA	1703	G
1	CA	1722	A
1	CA	1742	G
1	CA	1746	G
1	CA	1756	G
1	CA	1758	G
1	CA	1762	A
1	CA	1763	G
1	CA	1764	G
1	CA	1769	G
1	CA	1773	A
1	CA	1774	C
1	CA	1780	A
1	CA	1782	C

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Mol	Chain	Res	Type
1	CA	1791	A
1	CA	1800	C
1	CA	1801	G
1	CA	1816	G
1	CA	1829	A
1	CA	1835	G
1	CA	1847	A
1	CA	1848	A
1	CA	1859	A
1	CA	1877	A
1	CA	1878	G
1	CA	1886	C
1	CA	1900	A
1	CA	1906	G
1	CA	1913	A
1	CA	1914	C
1	CA	1929	G
1	CA	1930	G
1	CA	1934	C
1	CA	1937	A
1	CA	1938	A
1	CA	1955	U
1	CA	1962	C
1	CA	1963	U
1	CA	1967	C
1	CA	1970	A
1	CA	1971	A
1	CA	1972	A
1	CA	1984	G
1	CA	1992	G
1	CA	1993	U
1	CA	1994	C
1	CA	1997	G
1	CA	2020	A
1	CA	2023	G
1	CA	2027	G
1	CA	2030	A
1	CA	2031	A
1	CA	2032	G
1	CA	2033	A
1	CA	2043	C
1	CA	2046	G

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Mol	Chain	Res	Type
1	CA	2052	G
1	CA	2055	C
1	CA	2056	G
1	CA	2060	A
1	CA	2061	G
1	CA	2062	A
1	CA	2069	G
1	CA	2082	A
1	CA	2099	U
1	CA	2102	U
1	CA	2105	C
1	CA	2106	G
1	CA	2110	G
1	CA	2111	C
1	CA	2113	U
1	CA	2115	G
1	CA	2116	G
1	CA	2117	A
1	CA	2119	A
1	CA	2122	U
1	CA	2125	G
1	CA	2126	A
1	CA	2127	G
1	CA	2129	C
1	CA	2131	G
1	CA	2132	U
1	CA	2133	G
1	CA	2134	A
1	CA	2135	A
1	CA	2136	C
1	CA	2137	C
1	CA	2142	C
1	CA	2144	U
1	CA	2146	C
1	CA	2150	U
1	CA	2153	G
1	CA	2154	G
1	CA	2157	G
1	CA	2158	A
1	CA	2162	G
1	CA	2164	C
1	CA	2165	G

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Mol	Chain	Res	Type
1	CA	2167	U
1	CA	2168	G
1	CA	2169	A
1	CA	2172	U
1	CA	2175	C
1	CA	2178	C
1	CA	2181	G
1	CA	2184	G
1	CA	2185	C
1	CA	2186	G
1	CA	2188	C
1	CA	2189	U
1	CA	2190	G
1	CA	2192	G
1	CA	2195	C
1	CA	2198	A
1	CA	2206	G
1	CA	2207	G
1	CA	2208	A
1	CA	2218	U
1	CA	2225	A
1	CA	2238	G
1	CA	2240	C
1	CA	2256	G
1	CA	2268	A
1	CA	2273	A
1	CA	2275	C
1	CA	2278	A
1	CA	2283	C
1	CA	2287	A
1	CA	2288	A
1	CA	2289	G
1	CA	2305	A
1	CA	2306	C
1	CA	2308	G
1	CA	2312	U
1	CA	2318	G
1	CA	2319	G
1	CA	2320	A
1	CA	2325	G
1	CA	2327	A
1	CA	2334	G

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Mol	Chain	Res	Type
1	CA	2336	A
1	CA	2343	C
1	CA	2347	C
1	CA	2350	C
1	CA	2353	G
1	CA	2354	G
1	CA	2383	G
1	CA	2385	C
1	CA	2406	U
1	CA	2410	G
1	CA	2422	A
1	CA	2423	U
1	CA	2425	A
1	CA	2428	G
1	CA	2429	G
1	CA	2430	A
1	CA	2432	A
1	CA	2434	A
1	CA	2435	A
1	CA	2439	A
1	CA	2440	C
1	CA	2441	C
1	CA	2448	A
1	CA	2459	A
1	CA	2460	U
1	CA	2465	C
1	CA	2468	G
1	CA	2474	C
1	CA	2476	A
1	CA	2502	G
1	CA	2505	G
1	CA	2506	U
1	CA	2507	C
1	CA	2518	A
1	CA	2520	C
1	CA	2525	G
1	CA	2529	G
1	CA	2531	A
1	CA	2533	A
1	CA	2554	U
1	CA	2556	C
1	CA	2562	U

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Mol	Chain	Res	Type
1	CA	2564	A
1	CA	2566	A
1	CA	2567	G
1	CA	2573	C
1	CA	2579	C
1	CA	2582	G
1	CA	2596	U
1	CA	2602	A
1	CA	2609	U
1	CA	2610	C
1	CA	2611	U
1	CA	2612	C
1	CA	2621	A
1	CA	2630	G
1	CA	2632	A
1	CA	2634	G
1	CA	2647	U
1	CA	2653	U
1	CA	2654	A
1	CA	2663	G
1	CA	2673	G
1	CA	2674	G
1	CA	2689	U
1	CA	2690	C
1	CA	2699	C
1	CA	2702	U
1	CA	2712(A)	A
1	CA	2713	A
1	CA	2720	U
1	CA	2721	A
1	CA	2726	U
1	CA	2727	G
1	CA	2733	A
1	CA	2745	C
1	CA	2748	A
1	CA	2751	G
1	CA	2757	A
1	CA	2758	A
1	CA	2760	C
1	CA	2765	A
1	CA	2766	G
1	CA	2767	C

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Mol	Chain	Res	Type
1	CA	2775	A
1	CA	2778	A
1	CA	2802	G
1	CA	2803	C
1	CA	2807	G
1	CA	2809	A
1	CA	2810	A
1	CA	2818	G
1	CA	2820	A
1	CA	2821	A
1	CA	2833	G
1	CA	2835	A
1	CA	2839	G
1	CA	2872	G
1	CA	2874	C
1	CA	2875	C
1	CA	2877	G
1	CA	2879	C
1	CA	2892	A
1	CA	2893	G
1	CA	2894	G
1	CA	2896	C
1	CA	2897	U
2	CB	2	C
2	CB	5	C
2	CB	7	G
2	CB	12	C
2	CB	13	A
2	CB	15	A
2	CB	20	C
2	CB	21	G
2	CB	32	C
2	CB	34	U
2	CB	42	C
2	CB	52	A
2	CB	53	A
2	CB	56	G
2	CB	67	G
2	CB	73	A
2	CB	85	G
2	CB	94	C
2	CB	110	G

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Mol	Chain	Res	Type
2	CB	116	G
2	CB	118	G
34	DA	5	U
34	DA	9	G
34	DA	30	U
34	DA	32	A
34	DA	39	G
34	DA	47	C
34	DA	48	C
34	DA	49	U
34	DA	51	A
34	DA	54	C
34	DA	62	U
34	DA	73	G
34	DA	80	G
34	DA	89	C
34	DA	97	G
34	DA	101	A
34	DA	116	A
34	DA	121	C
34	DA	131	C
34	DA	143	A
34	DA	163	C
34	DA	174	C
34	DA	182	U
34	DA	189(E)	U
34	DA	195	A
34	DA	197	A
34	DA	201	C
34	DA	203	U
34	DA	204	U
34	DA	216	G
34	DA	231	G
34	DA	243	A
34	DA	245	C
34	DA	247	G
34	DA	251	G
34	DA	252	U
34	DA	258	G
34	DA	266	G
34	DA	267	C
34	DA	269	C

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Mol	Chain	Res	Type
34	DA	281	G
34	DA	289	G
34	DA	290	C
34	DA	298	A
34	DA	301	G
34	DA	306	G
34	DA	308	C
34	DA	318	G
34	DA	328	C
34	DA	329	A
34	DA	332	G
34	DA	341	C
34	DA	342	C
34	DA	343	U
34	DA	344	A
34	DA	346	G
34	DA	347	G
34	DA	351	G
34	DA	352	C
34	DA	353	A
34	DA	354	G
34	DA	367	U
34	DA	372	C
34	DA	373	A
34	DA	382	A
34	DA	384	G
34	DA	398	C
34	DA	406	G
34	DA	412	A
34	DA	413	G
34	DA	424	G
34	DA	429	U
34	DA	430	A
34	DA	439	A
34	DA	442	C
34	DA	443	C
34	DA	452	A
34	DA	485	G
34	DA	496	A
34	DA	498	U
34	DA	499	A
34	DA	500	G

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Mol	Chain	Res	Type
34	DA	505	G
34	DA	509	A
34	DA	510	A
34	DA	511	C
34	DA	518	C
34	DA	527	G
34	DA	531	U
34	DA	532	A
34	DA	533	A
34	DA	536	C
34	DA	545	C
34	DA	547	A
34	DA	559	A
34	DA	560	U
34	DA	561	U
34	DA	562	C
34	DA	564	C
34	DA	567	G
34	DA	572	A
34	DA	573	A
34	DA	576	G
34	DA	577	G
34	DA	581	G
34	DA	592	G
34	DA	596	C
34	DA	600	C
34	DA	601	C
34	DA	619	U
34	DA	630	G
34	DA	632	A
34	DA	651	C
34	DA	653	A
34	DA	665	A
34	DA	675	A
34	DA	687	A
34	DA	688	G
34	DA	693	G
34	DA	695	A
34	DA	698	G
34	DA	702	A
34	DA	717	C
34	DA	721	G

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Mol	Chain	Res	Type
34	DA	723	U
34	DA	728	A
34	DA	731	G
34	DA	749	C
34	DA	755	G
34	DA	766	A
34	DA	768	A
34	DA	774	G
34	DA	777	A
34	DA	788	U
34	DA	792	A
34	DA	793	U
34	DA	794	A
34	DA	816	A
34	DA	817	C
34	DA	821	G
34	DA	825	G
34	DA	827	U
34	DA	828	A
34	DA	829	G
34	DA	834	C
34	DA	840	C
34	DA	841	U
34	DA	848	C
34	DA	851	G
34	DA	859	A
34	DA	873	A
34	DA	874	G
34	DA	876	G
34	DA	880	C
34	DA	884	U
34	DA	887	G
34	DA	902	G
34	DA	913	A
34	DA	914	A
34	DA	916	G
34	DA	922	G
34	DA	926	G
34	DA	927	G
34	DA	934	C
34	DA	935	A
34	DA	936	C

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Mol	Chain	Res	Type
34	DA	959	A
34	DA	961	U
34	DA	966	G
34	DA	968	A
34	DA	969	A
34	DA	971	G
34	DA	974	A
34	DA	975	A
34	DA	976	G
34	DA	977	A
34	DA	981	U
34	DA	989	C
34	DA	992	U
34	DA	993	G
34	DA	999	C
34	DA	1003	G
34	DA	1005	A
34	DA	1006	C
34	DA	1007	C
34	DA	1016	A
34	DA	1017	G
34	DA	1022	G
34	DA	1025	U
34	DA	1026	G
34	DA	1027	C
34	DA	1028	C
34	DA	1030	C
34	DA	1030(A)	G
34	DA	1030(B)	C
34	DA	1033	G
34	DA	1064	G
34	DA	1065	U
34	DA	1066	C
34	DA	1067	A
34	DA	1081	G
34	DA	1094	G
34	DA	1095	U
34	DA	1097	C
34	DA	1101	A
34	DA	1105	A
34	DA	1117	G
34	DA	1121	U

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Mol	Chain	Res	Type
34	DA	1122	U
34	DA	1125	U
34	DA	1127	G
34	DA	1130	A
34	DA	1136	U
34	DA	1137	C
34	DA	1138	G
34	DA	1139	G
34	DA	1140	C
34	DA	1146	A
34	DA	1152	A
34	DA	1154	G
34	DA	1155	G
34	DA	1159	U
34	DA	1161	C
34	DA	1164	G
34	DA	1166	G
34	DA	1181	G
34	DA	1183	A
34	DA	1184	G
34	DA	1190	G
34	DA	1193	G
34	DA	1195	C
34	DA	1196	U
34	DA	1197	G
34	DA	1201	A
34	DA	1202	G
34	DA	1211	U
34	DA	1212	U
34	DA	1213	A
34	DA	1214	C
34	DA	1220	G
34	DA	1225	A
34	DA	1227	A
34	DA	1228	C
34	DA	1236	A
34	DA	1238	A
34	DA	1240	U
34	DA	1241	G
34	DA	1249	C
34	DA	1252	A
34	DA	1253	G

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Mol	Chain	Res	Type
34	DA	1254	C
34	DA	1256	A
34	DA	1257	U
34	DA	1258	G
34	DA	1261	A
34	DA	1279	A
34	DA	1280	A
34	DA	1281	U
34	DA	1282	C
34	DA	1285	A
34	DA	1286	A
34	DA	1287	A
34	DA	1297	C
34	DA	1298	C
34	DA	1301	U
34	DA	1305	G
34	DA	1320	C
34	DA	1322	C
34	DA	1326	C
34	DA	1340	A
34	DA	1346	A
34	DA	1347	G
34	DA	1355	G
34	DA	1360	A
34	DA	1363	C
34	DA	1363(A)	A
34	DA	1364	U
34	DA	1368	G
34	DA	1381	U
34	DA	1397	C
34	DA	1398	A
34	DA	1401	G
34	DA	1419	G
34	DA	1442	G
34	DA	1442(A)	G
34	DA	1442(B)	A
34	DA	1446	U
34	DA	1447	A
34	DA	1452	C
34	DA	1456	G
34	DA	1469	G
34	DA	1487	G

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Mol	Chain	Res	Type
34	DA	1497	G
34	DA	1503	A
34	DA	1504	G
34	DA	1506	U
34	DA	1517	G
34	DA	1520	G
34	DA	1525	G
34	DA	1529	G
34	DA	1530	G
34	DA	1531	A
34	DA	1532	U
55	DV	14	A
56	DW	8	4SU
56	DW	14	A
56	DW	16	U
56	DW	17	C
56	DW	18	G
56	DW	19	G
56	DW	20	U
56	DW	21	A
56	DW	40	C
56	DW	41	C
56	DW	43	C
56	DW	45	U
56	DW	46	7MG
56	DW	47	U
56	DW	48	C
56	DW	49	C
56	DW	62	C
56	DW	73	A
56	DW	76	A
56	DY	9	A
56	DY	13	C
56	DY	14	A
56	DY	19	G
56	DY	26	A
56	DY	33	U
56	DY	34	G
56	DY	39	PSU
56	DY	45	U
56	DY	46	7MG
56	DY	47	U

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Mol	Chain	Res	Type
56	DY	48	C
56	DY	49	C
56	DY	52	G
56	DY	54	5MU
56	DY	55	PSU
56	DY	57	G
56	DY	58	A
56	DY	59	U
56	DY	65	G
56	DY	70	G

All (146) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	AA	184	A
1	AA	185	A
1	AA	188	A
1	AA	204	G
1	AA	210	A
1	AA	238	C
1	AA	302	A
1	AA	334	A
1	AA	528	A
1	AA	572	A
1	AA	596	G
1	AA	645	G
1	AA	716	G
1	AA	732	A
1	AA	793	A
1	AA	811	A
1	AA	821	A
1	AA	823	G
1	AA	837	C
1	AA	906	G
1	AA	990	A
1	AA	1019	G
1	AA	1057	G
1	AA	1072	U
1	AA	1098	C
1	AA	1154	U
1	AA	1188	A
1	AA	1219	A

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Mol	Chain	Res	Type
1	AA	1220	U
1	AA	1221	G
1	AA	1255	A
1	AA	1425	A
1	AA	1442	U
1	AA	1466	U
1	AA	1654	A
1	AA	1655	A
1	AA	1700	G
1	AA	1793	A
1	AA	1811	A
1	AA	2014	G
1	AA	2019	G
1	AA	2203	G
1	AA	2209	G
1	AA	2227	G
1	AA	2287	C
1	AA	2347	A
1	AA	2358	A
1	AA	2418	U
1	AA	2434	A
1	AA	2451	A
1	AA	2459	G
1	AA	2518	U
1	AA	2623	U
1	AA	2701	U
1	AA	2769	U
1	AA	2902	G
34	BA	115	G
34	BA	243	A
34	BA	250	A
34	BA	266	G
34	BA	347	G
34	BA	509	A
34	BA	510	A
34	BA	530	G
34	BA	560	U
34	BA	687	A
34	BA	748	C
34	BA	793	U
34	BA	839	U
34	BA	884	U

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Mol	Chain	Res	Type
34	BA	913	A
34	BA	991	U
34	BA	1065	U
34	BA	1067	A
34	BA	1165	C
34	BA	1201	A
34	BA	1285	A
34	BA	1442	G
56	BY	19	G
56	BY	58	A
1	CA	27	G
1	CA	195	A
1	CA	196	A
1	CA	271(K)	U
1	CA	277	C
1	CA	310	A
1	CA	503	A
1	CA	645	C
1	CA	685	A
1	CA	746	A
1	CA	764	A
1	CA	774	A
1	CA	776	G
1	CA	830	G
1	CA	856	C
1	CA	900	A
1	CA	974	G
1	CA	1026	U
1	CA	1057	A
1	CA	1240	U
1	CA	1300	U
1	CA	1379	A
1	CA	1420	U
1	CA	1427	A
1	CA	1558	A
1	CA	1608	A
1	CA	1653	G
1	CA	1913	A
1	CA	1992	G
1	CA	1997	G
1	CA	2110	G
1	CA	2288	A

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Mol	Chain	Res	Type
1	CA	2318	G
1	CA	2406	U
1	CA	2439	A
1	CA	2566	A
1	CA	2689	U
1	CA	2726	U
1	CA	2756	U
34	DA	115	G
34	DA	119	A
34	DA	251	G
34	DA	266	G
34	DA	428	G
34	DA	429	U
34	DA	499	A
34	DA	509	A
34	DA	535	A
34	DA	560	U
34	DA	687	A
34	DA	748	C
34	DA	793	U
34	DA	840	C
34	DA	873	A
34	DA	884	U
34	DA	913	A
34	DA	991	U
34	DA	1064	G
34	DA	1065	U
34	DA	1183	A
34	DA	1201	A
34	DA	1279	A
34	DA	1442	G
56	DW	13	C
56	DW	45	U
56	DY	46	7MG

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

28 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$
56	PSU	BW	32	56	16,21,22	1.07	2 (12%)	20,30,33	3.67	7 (35%)
56	MIA	BW	37	56	23,31,32	1.89	4 (17%)	25,44,47	2.79	6 (24%)
56	PSU	BW	39	56	16,21,22	1.68	1 (6%)	20,30,33	3.66	7 (35%)
56	7MG	BW	46	56	20,26,27	1.56	2 (10%)	22,39,42	2.86	5 (22%)
56	5MU	BW	54	56	14,22,23	0.81	1 (7%)	16,32,35	2.02	3 (18%)
56	PSU	BW	55	56	16,21,22	1.35	1 (6%)	20,30,33	3.68	6 (30%)
56	4SU	BW	8	56	14,21,22	1.14	1 (7%)	15,30,33	1.47	2 (13%)
56	PSU	BY	32	56	16,21,22	1.12	1 (6%)	20,30,33	3.54	6 (30%)
56	MIA	BY	37	56	18,24,32	1.28	2 (11%)	17,35,47	1.91	3 (17%)
56	PSU	BY	39	56	16,21,22	1.21	1 (6%)	20,30,33	3.67	6 (30%)
56	7MG	BY	46	56	20,26,27	1.68	2 (10%)	22,39,42	2.79	6 (27%)
56	5MU	BY	54	56	14,22,23	0.77	0	16,32,35	2.34	2 (12%)
56	PSU	BY	55	56	16,21,22	1.25	1 (6%)	20,30,33	3.56	6 (30%)
56	4SU	BY	8	56	14,21,22	1.21	1 (7%)	15,30,33	1.69	3 (20%)
56	PSU	DW	32	56	16,21,22	0.86	0	20,30,33	4.11	8 (40%)
56	MIA	DW	37	56	23,31,32	2.01	3 (13%)	25,44,47	1.23	4 (16%)
56	PSU	DW	39	56	16,21,22	1.37	2 (12%)	20,30,33	3.81	8 (40%)
56	7MG	DW	46	56	20,26,27	1.69	2 (10%)	22,39,42	2.76	5 (22%)
56	5MU	DW	54	56	14,22,23	0.90	1 (7%)	16,32,35	2.29	3 (18%)
56	PSU	DW	55	56	16,21,22	1.12	1 (6%)	20,30,33	3.78	7 (35%)
56	4SU	DW	8	56	14,21,22	1.23	1 (7%)	15,30,33	1.59	2 (13%)
56	PSU	DY	32	56	16,21,22	1.14	1 (6%)	20,30,33	3.56	6 (30%)
56	MIA	DY	37	56	18,24,32	1.17	2 (11%)	17,35,47	1.83	2 (11%)
56	PSU	DY	39	56	16,21,22	1.34	2 (12%)	20,30,33	3.77	6 (30%)
56	7MG	DY	46	56	20,26,27	1.65	3 (15%)	22,39,42	2.98	7 (31%)
56	5MU	DY	54	56	14,22,23	0.76	0	16,32,35	2.30	3 (18%)
56	PSU	DY	55	56	16,21,22	1.35	2 (12%)	20,30,33	3.53	6 (30%)
56	4SU	DY	8	56	14,21,22	1.29	1 (7%)	15,30,33	1.52	2 (13%)

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
56	PSU	BW	32	56	-	0/7/25/26	0/2/2/2
56	MIA	BW	37	56	-	0/11/33/34	0/3/3/3
56	PSU	BW	39	56	-	0/7/25/26	0/2/2/2
56	7MG	BW	46	56	-	0/7/37/38	0/3/3/3
56	5MU	BW	54	56	-	0/3/25/26	0/2/2/2
56	PSU	BW	55	56	-	0/7/25/26	0/2/2/2
56	4SU	BW	8	56	-	0/3/25/26	0/2/2/2
56	PSU	BY	32	56	-	0/7/25/26	0/2/2/2
56	MIA	BY	37	56	-	0/3/25/34	0/3/3/3
56	PSU	BY	39	56	-	0/7/25/26	0/2/2/2
56	7MG	BY	46	56	-	0/7/37/38	0/3/3/3
56	5MU	BY	54	56	-	0/3/25/26	0/2/2/2
56	PSU	BY	55	56	-	0/7/25/26	0/2/2/2
56	4SU	BY	8	56	-	0/3/25/26	0/2/2/2
56	PSU	DW	32	56	-	0/7/25/26	0/2/2/2
56	MIA	DW	37	56	-	0/11/33/34	0/3/3/3
56	PSU	DW	39	56	-	0/7/25/26	0/2/2/2
56	7MG	DW	46	56	-	0/7/37/38	0/3/3/3
56	5MU	DW	54	56	-	0/3/25/26	0/2/2/2
56	PSU	DW	55	56	-	0/7/25/26	0/2/2/2
56	4SU	DW	8	56	-	0/3/25/26	0/2/2/2
56	PSU	DY	32	56	-	0/7/25/26	0/2/2/2
56	MIA	DY	37	56	-	0/3/25/34	0/3/3/3
56	PSU	DY	39	56	-	0/7/25/26	0/2/2/2
56	7MG	DY	46	56	-	0/7/37/38	0/3/3/3
56	5MU	DY	54	56	-	0/3/25/26	0/2/2/2
56	PSU	DY	55	56	-	0/7/25/26	0/2/2/2
56	4SU	DY	8	56	-	0/3/25/26	0/2/2/2

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
56	DW	37	MIA	C2-S10	-7.89	1.69	1.75
56	BW	37	MIA	C2-S10	-7.17	1.69	1.75
56	BW	39	PSU	C5-C1'	-5.72	1.47	1.52
56	BW	55	PSU	C5-C1'	-4.01	1.48	1.52
56	DY	8	4SU	C4-S4	-3.88	1.60	1.67
56	DY	55	PSU	C5-C1'	-3.73	1.49	1.52
56	DW	39	PSU	C5-C1'	-3.70	1.49	1.52
56	DW	8	4SU	C4-S4	-3.59	1.60	1.67
56	BY	8	4SU	C4-S4	-3.57	1.60	1.67
56	BY	55	PSU	C5-C1'	-3.50	1.49	1.52
56	BW	8	4SU	C4-S4	-3.48	1.60	1.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
56	DY	39	PSU	C5-C1'	-3.44	1.49	1.52
56	BY	39	PSU	C5-C1'	-3.31	1.49	1.52
56	BW	37	MIA	O5'-C5'	-3.27	1.40	1.44
56	DY	32	PSU	C5-C1'	-2.93	1.49	1.52
56	BW	32	PSU	C2-N3	-2.77	1.32	1.38
56	DW	55	PSU	C5-C1'	-2.75	1.49	1.52
56	BY	32	PSU	C5-C1'	-2.58	1.50	1.52
56	DW	37	MIA	O5'-C5'	-2.16	1.41	1.44
56	DW	54	5MU	O5'-C5'	-2.15	1.41	1.44
56	DW	39	PSU	O5'-C5'	-2.10	1.41	1.44
56	BW	32	PSU	C5-C1'	-2.09	1.50	1.52
56	DY	39	PSU	O4'-C1'	-2.09	1.41	1.44
56	DY	55	PSU	C2-N1	-2.08	1.34	1.38
56	BW	54	5MU	O5'-C5'	-2.01	1.41	1.44
56	DY	46	7MG	CM7-N7	2.16	1.49	1.46
56	BW	37	MIA	C6-N1	2.30	1.36	1.33
56	DY	37	MIA	C2-N3	2.38	1.36	1.32
56	BY	37	MIA	C2-N3	2.48	1.36	1.32
56	BW	37	MIA	C5-C4	2.67	1.46	1.40
56	BW	46	7MG	C5-C4	2.70	1.46	1.39
56	BY	46	7MG	C5-C4	3.13	1.47	1.39
56	DY	46	7MG	C5-C4	3.18	1.47	1.39
56	DW	46	7MG	C5-C4	3.24	1.47	1.39
56	DW	37	MIA	C5-C4	3.27	1.47	1.40
56	DY	37	MIA	C5-C4	3.34	1.48	1.40
56	BY	37	MIA	C5-C4	3.60	1.48	1.40
56	DY	46	7MG	C6-C5	5.65	1.48	1.41
56	BW	46	7MG	C6-C5	5.87	1.48	1.41
56	BY	46	7MG	C6-C5	6.06	1.48	1.41
56	DW	46	7MG	C6-C5	6.23	1.48	1.41

All (137) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
56	DW	32	PSU	N1-C2-N3	-11.82	119.90	128.40
56	BW	37	MIA	C11-S10-C2	-11.82	93.56	102.29
56	BW	32	PSU	N1-C2-N3	-11.71	119.97	128.40
56	BW	55	PSU	N1-C2-N3	-10.33	120.97	128.40
56	DW	55	PSU	N1-C2-N3	-10.14	121.10	128.40
56	DY	32	PSU	N1-C2-N3	-9.96	121.23	128.40
56	BY	39	PSU	N1-C2-N3	-9.71	121.41	128.40
56	DY	39	PSU	C5-C4-N3	-9.71	117.47	125.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
56	BY	32	PSU	N1-C2-N3	-9.68	121.44	128.40
56	DY	39	PSU	N1-C2-N3	-9.48	121.58	128.40
56	DY	55	PSU	N1-C2-N3	-9.41	121.63	128.40
56	DW	39	PSU	N1-C2-N3	-9.39	121.64	128.40
56	BY	55	PSU	N1-C2-N3	-9.36	121.67	128.40
56	BW	39	PSU	N1-C2-N3	-9.23	121.76	128.40
56	BW	39	PSU	C5-C4-N3	-8.92	118.11	125.43
56	DW	55	PSU	C5-C4-N3	-8.76	118.24	125.43
56	BY	39	PSU	C5-C4-N3	-8.71	118.28	125.43
56	BY	55	PSU	C5-C4-N3	-8.69	118.30	125.43
56	DY	55	PSU	C5-C4-N3	-8.30	118.62	125.43
56	BY	32	PSU	C5-C4-N3	-8.12	118.77	125.43
56	DW	39	PSU	C5-C4-N3	-8.11	118.77	125.43
56	DY	32	PSU	C5-C4-N3	-8.05	118.82	125.43
56	BW	55	PSU	C5-C4-N3	-7.67	119.14	125.43
56	DW	32	PSU	C5-C4-N3	-6.97	119.71	125.43
56	BY	37	MIA	N3-C2-N1	-6.43	123.26	128.86
56	DY	37	MIA	N3-C2-N1	-6.22	123.44	128.86
56	BY	54	5MU	C5-C4-N3	-5.90	118.73	125.24
56	BW	46	7MG	C5-C4-N3	-5.87	116.67	126.47
56	DY	54	5MU	C5-C4-N3	-5.58	119.09	125.24
56	DY	46	7MG	C5-C4-N3	-5.45	117.37	126.47
56	DW	54	5MU	C5-C4-N3	-5.38	119.31	125.24
56	BY	46	7MG	C5-C4-N3	-5.23	117.75	126.47
56	DW	46	7MG	C5-C4-N3	-5.13	117.91	126.47
56	DW	46	7MG	C5-C6-N1	-4.92	115.65	123.37
56	DY	46	7MG	C5-C6-N1	-4.80	115.83	123.37
56	BY	46	7MG	C5-C6-N1	-4.64	116.10	123.37
56	BW	32	PSU	C5-C6-N1	-4.54	118.51	124.39
56	BW	54	5MU	C5-C4-N3	-4.41	120.37	125.24
56	DW	39	PSU	C5-C6-N1	-4.37	118.72	124.39
56	BW	39	PSU	C5-C6-N1	-4.31	118.80	124.39
56	BW	55	PSU	C5-C6-N1	-4.07	119.12	124.39
56	BW	32	PSU	C5-C4-N3	-3.99	122.15	125.43
56	BW	37	MIA	C4-C5-N7	-3.96	105.58	109.41
56	BW	32	PSU	C5-C1'-C2'	-3.90	108.82	115.55
56	BY	8	4SU	C5-C4-N3	-3.88	118.82	123.73
56	DW	8	4SU	C5-C4-N3	-3.87	118.84	123.73
56	DW	55	PSU	C5-C6-N1	-3.87	119.37	124.39
56	DY	55	PSU	C5-C6-N1	-3.84	119.41	124.39
56	DW	39	PSU	C5-C1'-C2'	-3.78	109.03	115.55
56	BW	46	7MG	C5-C6-N1	-3.68	117.59	123.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
56	BY	32	PSU	C5-C6-N1	-3.66	119.65	124.39
56	DW	32	PSU	C5-C1'-C2'	-3.53	109.46	115.55
56	BY	55	PSU	C5-C6-N1	-3.47	119.90	124.39
56	BY	37	MIA	C4-C5-N7	-3.28	106.24	109.41
56	BY	39	PSU	C5-C6-N1	-3.28	120.13	124.39
56	BW	55	PSU	C5-C1'-C2'	-3.27	109.90	115.55
56	DW	54	5MU	C5-C6-N1	-3.24	118.64	122.15
56	DY	39	PSU	O4'-C1'-C5	-3.22	104.94	109.93
56	DY	32	PSU	C5-C6-N1	-3.20	120.24	124.39
56	DY	8	4SU	C5-C4-N3	-3.16	119.73	123.73
56	DY	37	MIA	C4-C5-N7	-3.09	106.43	109.41
56	BW	37	MIA	C5-C6-N1	-3.08	117.57	120.64
56	BY	39	PSU	C5-C1'-C2'	-3.07	110.25	115.55
56	DW	39	PSU	O2'-C2'-C1'	-2.92	105.60	112.21
56	BW	39	PSU	O4'-C1'-C5	-2.88	105.47	109.93
56	DW	32	PSU	C5-C6-N1	-2.79	120.77	124.39
56	DY	39	PSU	C5-C6-N1	-2.77	120.80	124.39
56	BW	8	4SU	C5-C4-N3	-2.72	120.29	123.73
56	DW	37	MIA	C4-C5-N7	-2.64	106.86	109.41
56	DY	55	PSU	C5-C1'-C2'	-2.55	111.14	115.55
56	BY	55	PSU	C5-C1'-C2'	-2.54	111.16	115.55
56	BW	37	MIA	N3-C2-N1	-2.52	122.32	126.85
56	DY	54	5MU	C5-C6-N1	-2.51	119.43	122.15
56	DW	55	PSU	C5-C1'-C2'	-2.42	111.38	115.55
56	BW	54	5MU	C5-C6-N1	-2.23	119.74	122.15
56	DY	46	7MG	C5-C4-N9	-2.23	103.07	106.31
56	DY	32	PSU	C5-C1'-C2'	-2.22	111.72	115.55
56	DW	37	MIA	C11-S10-C2	-2.20	100.67	102.29
56	BW	39	PSU	C5-C1'-C2'	-2.12	111.88	115.55
56	DW	37	MIA	C5-C6-N1	-2.09	118.55	120.64
56	BY	46	7MG	C5-C4-N9	-2.06	103.32	106.31
56	BY	8	4SU	C6-N1-C2	-2.04	117.98	121.28
56	DW	55	PSU	O2'-C2'-C1'	-2.01	107.67	112.21
56	DW	46	7MG	C2-N3-C4	2.02	119.63	113.95
56	BY	37	MIA	C2-N1-C6	2.03	122.32	118.77
56	BY	32	PSU	O4'-C1'-C2'	2.10	107.83	104.45
56	DY	46	7MG	C4-N9-C1'	2.14	131.76	126.58
56	DY	46	7MG	C2-N3-C4	2.20	120.12	113.95
56	BY	46	7MG	C2-N3-C4	2.24	120.25	113.95
56	BW	37	MIA	N6-C6-N1	2.38	121.51	118.54
56	DW	37	MIA	C2-N1-C6	2.59	121.09	113.47
56	BW	37	MIA	C2-N1-C6	2.86	121.89	113.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
56	BW	46	7MG	C2-N3-C4	2.88	122.04	113.95
56	BW	32	PSU	O4'-C1'-C5	2.90	114.42	109.93
56	DW	32	PSU	C4-C5-C1'	2.90	126.76	121.15
56	BW	46	7MG	C6-N1-C2	3.64	121.29	116.06
56	DY	39	PSU	C6-N1-C2	3.76	121.38	115.36
56	BW	8	4SU	C2-N3-C4	3.85	120.79	115.11
56	BY	55	PSU	C6-N1-C2	4.03	121.81	115.36
56	DY	8	4SU	C2-N3-C4	4.08	121.13	115.11
56	BW	39	PSU	C6-N1-C2	4.13	121.97	115.36
56	BY	39	PSU	C6-N1-C2	4.13	121.97	115.36
56	DY	32	PSU	C6-N1-C2	4.20	122.09	115.36
56	DW	8	4SU	C2-N3-C4	4.25	121.39	115.11
56	BY	32	PSU	C6-N1-C2	4.35	122.33	115.36
56	DY	55	PSU	C6-N1-C2	4.46	122.49	115.36
56	BY	8	4SU	C2-N3-C4	4.51	121.76	115.11
56	DY	46	7MG	C6-N1-C2	4.51	122.55	116.06
56	BY	46	7MG	C6-N1-C2	4.55	122.61	116.06
56	DW	39	PSU	C6-N1-C2	4.57	122.67	115.36
56	BW	55	PSU	C6-N1-C2	4.61	122.74	115.36
56	DW	46	7MG	C6-N1-C2	4.61	122.70	116.06
56	DW	55	PSU	C6-N1-C2	4.67	122.84	115.36
56	DW	32	PSU	C6-N1-C2	4.89	123.18	115.36
56	BW	32	PSU	C4-N3-C2	5.39	119.88	115.16
56	DW	39	PSU	O4'-C1'-C5	5.40	118.29	109.93
56	BW	32	PSU	C6-N1-C2	5.70	124.48	115.36
56	DW	32	PSU	O4'-C1'-C5	5.77	118.87	109.93
56	BW	54	5MU	C4-N3-C2	5.81	120.24	115.16
56	DW	39	PSU	C4-N3-C2	5.99	120.40	115.16
56	DW	54	5MU	C4-N3-C2	6.08	120.48	115.16
56	DY	55	PSU	C4-N3-C2	6.35	120.72	115.16
56	BW	55	PSU	C4-N3-C2	6.44	120.80	115.16
56	DY	54	5MU	C4-N3-C2	6.45	120.80	115.16
56	BY	54	5MU	C4-N3-C2	6.52	120.86	115.16
56	BW	39	PSU	C4-N3-C2	6.70	121.02	115.16
56	BY	32	PSU	C4-N3-C2	6.72	121.04	115.16
56	BY	55	PSU	C4-N3-C2	6.78	121.09	115.16
56	DY	32	PSU	C4-N3-C2	6.85	121.15	115.16
56	DW	55	PSU	C4-N3-C2	6.94	121.23	115.16
56	BY	39	PSU	C4-N3-C2	7.02	121.30	115.16
56	DW	32	PSU	C4-N3-C2	7.39	121.62	115.16
56	DY	39	PSU	C4-N3-C2	7.48	121.70	115.16
56	DW	46	7MG	N3-C4-N9	9.01	138.48	126.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
56	BY	46	7MG	N3-C4-N9	9.35	138.93	126.98
56	BW	46	7MG	N3-C4-N9	9.75	139.44	126.98
56	DY	46	7MG	N3-C4-N9	9.84	139.55	126.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
56	BW	37	MIA	1	0
56	BY	8	4SU	1	0
56	DW	37	MIA	2	0
56	DW	39	PSU	4	0
56	DW	46	7MG	2	0
56	DW	54	5MU	1	0
56	DW	55	PSU	2	0
56	DY	37	MIA	1	0
56	DY	46	7MG	2	0
56	DY	55	PSU	3	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2058 ligands modelled in this entry, 2052 are monoatomic - leaving 6 for Mogul analysis.

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	$\# Z > 2$
60	SF4	BD	501	37	0,12,12	0.00	-	0,24,24	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
61	FUA	BZ	703	-	37,40,40	1.70	6 (16%)	45,64,64	1.72	7 (15%)
62	GDP	BZ	704	58	25,30,30	1.09	2 (8%)	26,47,47	1.92	6 (23%)
60	SF4	DD	501	37	0,12,12	0.00	-	0,24,24	0.00	-
61	FUA	DZ	703	-	37,40,40	1.70	6 (16%)	45,64,64	1.72	7 (15%)
62	GDP	DZ	704	58	25,30,30	1.20	2 (8%)	26,47,47	2.15	6 (23%)

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
60	SF4	BD	501	37	-	0/0/48/48	0/6/5/5
61	FUA	BZ	703	-	-	0/10/92/92	0/4/4/4
62	GDP	BZ	704	58	-	0/12/32/32	0/3/3/3
60	SF4	DD	501	37	-	0/0/48/48	0/6/5/5
61	FUA	DZ	703	-	-	0/10/92/92	0/4/4/4
62	GDP	DZ	704	58	-	0/12/32/32	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
61	BZ	703	FUA	C23-C22	-6.40	1.39	1.51
61	DZ	703	FUA	C23-C22	-6.35	1.40	1.51
61	BZ	703	FUA	C23-C24	-4.14	1.39	1.53
61	DZ	703	FUA	C23-C24	-4.13	1.39	1.53
61	DZ	703	FUA	C24-C25	-3.13	1.39	1.50
61	BZ	703	FUA	C24-C25	-3.13	1.39	1.50
61	BZ	703	FUA	C14-C8	-2.89	1.53	1.58
61	DZ	703	FUA	C14-C8	-2.87	1.53	1.58
61	BZ	703	FUA	C10-C9	-2.10	1.53	1.57
61	DZ	703	FUA	C10-C9	-2.09	1.53	1.57
61	BZ	703	FUA	C25-C26	2.33	1.39	1.32
61	DZ	703	FUA	C25-C26	2.35	1.39	1.32
62	BZ	704	GDP	C5-C4	2.61	1.46	1.40
62	BZ	704	GDP	C6-C5	2.74	1.46	1.41
62	DZ	704	GDP	C5-C4	3.37	1.48	1.40
62	DZ	704	GDP	C6-C5	4.22	1.49	1.41

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
62	DZ	704	GDP	C5-C6-N1	-4.65	116.87	123.48
61	BZ	703	FUA	C13-C12-C11	-4.43	105.80	111.97
61	DZ	703	FUA	C13-C12-C11	-4.40	105.85	111.97
62	DZ	704	GDP	C6-C5-C4	-3.92	116.95	120.84
61	DZ	703	FUA	C16-O2-C31	-3.87	111.10	117.10
61	BZ	703	FUA	C16-O2-C31	-3.86	111.12	117.10
61	BZ	703	FUA	C8-C9-C10	-3.54	112.76	116.44
62	BZ	704	GDP	C6-C5-C4	-3.50	117.36	120.84
61	DZ	703	FUA	C8-C9-C10	-3.50	112.81	116.44
62	BZ	704	GDP	C5-C6-N1	-3.35	118.71	123.48
62	DZ	704	GDP	N3-C2-N1	-3.05	123.00	127.46
62	BZ	704	GDP	N3-C2-N1	-2.97	123.12	127.46
62	DZ	704	GDP	C4-C5-N7	-2.78	106.72	109.41
61	BZ	703	FUA	C12-C13-C14	-2.01	110.45	114.72
61	DZ	703	FUA	C12-C13-C14	-2.01	110.45	114.72
62	BZ	704	GDP	C2'-C3'-C4'	2.10	106.70	102.62
61	BZ	703	FUA	C28-C26-C27	2.21	119.75	114.60
61	DZ	703	FUA	C28-C26-C27	2.21	119.77	114.60
61	BZ	703	FUA	O2-C31-C32	2.88	116.52	111.10
61	DZ	703	FUA	O2-C31-C32	2.90	116.54	111.10
62	BZ	704	GDP	C6-N1-C2	3.37	120.91	116.06
62	DZ	704	GDP	C2-N3-C4	4.65	120.58	115.16
62	BZ	704	GDP	C2-N3-C4	4.91	120.90	115.16
61	DZ	703	FUA	C24-C23-C22	5.09	124.09	112.50
61	BZ	703	FUA	C24-C23-C22	5.09	124.11	112.50
62	DZ	704	GDP	C6-N1-C2	5.20	123.54	116.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 45 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
60	BD	501	SF4	1	0
61	BZ	703	FUA	11	0
62	BZ	704	GDP	5	0
60	DD	501	SF4	2	0
61	DZ	703	FUA	17	0
62	DZ	704	GDP	9	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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	AA	2852/2915 (97%)	0.32	112 (3%) 40 29	14, 34, 139, 364	4 (0%)
1	CA	2848/2915 (97%)	0.37	128 (4%) 34 24	27, 57, 180, 356	0
2	AB	120/121 (99%)	-0.17	0 100 100	23, 50, 73, 110	0
2	CB	120/121 (99%)	0.09	1 (0%) 86 81	64, 92, 120, 168	0
3	AC	137/228 (60%)	10.34	133 (97%) 0 0	258, 289, 307, 313	0
3	CC	137/228 (60%)	11.69	133 (97%) 0 0	281, 312, 331, 336	0
4	AD	275/276 (99%)	-0.33	2 (0%) 87 83	13, 35, 59, 137	1 (0%)
4	CD	275/276 (99%)	-0.27	2 (0%) 87 83	19, 48, 74, 130	2 (0%)
5	AE	204/206 (99%)	-0.39	0 100 100	5, 33, 57, 80	3 (1%)
5	CE	204/206 (99%)	0.00	4 (1%) 65 56	21, 63, 107, 134	0
6	AF	203/210 (96%)	-0.28	0 100 100	10, 35, 78, 174	0
6	CF	203/210 (96%)	-0.27	0 100 100	21, 64, 107, 155	0
7	AG	181/182 (99%)	0.03	4 (2%) 62 52	34, 78, 134, 212	1 (0%)
7	CG	181/182 (99%)	0.70	21 (11%) 5 3	73, 112, 177, 207	0
8	AH	174/180 (96%)	-0.36	1 (0%) 89 86	26, 46, 70, 112	0
8	CH	174/180 (96%)	1.30	44 (25%) 1 0	65, 112, 161, 200	0
9	AK	130/173 (75%)	0.74	17 (13%) 4 2	48, 105, 170, 232	0
9	CK	130/173 (75%)	2.60	68 (52%) 0 0	75, 162, 204, 231	0
10	AL	139/147 (94%)	3.50	92 (66%) 0 0	96, 173, 233, 253	0
10	CL	139/147 (94%)	5.83	120 (86%) 0 0	127, 196, 252, 287	1 (0%)
11	AN	140/140 (100%)	-0.38	0 100 100	14, 28, 61, 97	1 (0%)
11	CN	140/140 (100%)	0.15	3 (2%) 64 54	32, 72, 108, 150	0
12	AO	122/122 (100%)	-0.33	0 100 100	17, 37, 63, 79	1 (0%)
12	CO	122/122 (100%)	-0.22	0 100 100	36, 59, 85, 106	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	AP	149/150 (99%)	-0.07	0 100 100	11, 42, 81, 107	1 (0%)
13	CP	149/150 (99%)	0.24	4 (2%) 55 44	30, 68, 117, 137	0
14	AQ	141/141 (100%)	-0.32	0 100 100	11, 34, 54, 81	0
14	CQ	141/141 (100%)	-0.17	2 (1%) 75 69	38, 71, 101, 120	0
15	AR	118/118 (100%)	-0.34	0 100 100	16, 29, 45, 56	0
15	CR	118/118 (100%)	-0.12	0 100 100	33, 56, 90, 106	0
16	AS	110/112 (98%)	-0.22	0 100 100	29, 51, 81, 94	0
16	CS	110/112 (98%)	0.35	5 (4%) 34 24	46, 85, 120, 152	0
17	AT	131/146 (89%)	-0.29	1 (0%) 86 81	24, 41, 92, 165	0
17	CT	131/146 (89%)	-0.09	1 (0%) 86 81	43, 65, 105, 143	0
18	AU	116/118 (98%)	-0.33	0 100 100	9, 22, 38, 90	1 (0%)
18	CU	116/118 (98%)	-0.05	0 100 100	27, 65, 93, 107	0
19	AV	101/101 (100%)	-0.45	0 100 100	9, 28, 50, 75	0
19	CV	101/101 (100%)	0.17	1 (0%) 82 77	36, 80, 113, 171	0
20	AW	112/113 (99%)	-0.36	0 100 100	13, 26, 43, 112	1 (0%)
20	CW	112/113 (99%)	-0.12	0 100 100	27, 50, 81, 119	0
21	AX	95/96 (98%)	-0.30	1 (1%) 80 74	16, 35, 67, 99	1 (1%)
21	CX	95/96 (98%)	-0.02	2 (2%) 64 54	38, 62, 86, 107	0
22	AY	107/110 (97%)	-0.24	1 (0%) 84 79	24, 44, 87, 161	0
22	CY	107/110 (97%)	0.48	10 (9%) 9 5	46, 76, 115, 167	0
23	AZ	185/206 (89%)	-0.38	0 100 100	29, 57, 92, 148	0
23	CZ	185/206 (89%)	0.65	20 (10%) 6 3	61, 106, 149, 213	0
24	A0	83/85 (97%)	-0.08	4 (4%) 31 21	12, 35, 83, 225	1 (1%)
24	C0	83/85 (97%)	0.74	9 (10%) 6 3	42, 66, 122, 228	0
25	A1	97/98 (98%)	-0.14	2 (2%) 64 54	19, 43, 80, 101	1 (1%)
25	C1	97/98 (98%)	-0.17	1 (1%) 82 77	31, 52, 91, 125	0
26	A2	70/72 (97%)	-0.28	2 (2%) 52 41	25, 44, 69, 123	1 (1%)
26	C2	70/72 (97%)	-0.01	2 (2%) 52 41	49, 71, 101, 117	0
27	A3	59/60 (98%)	-0.32	0 100 100	14, 30, 56, 101	1 (1%)
27	C3	59/60 (98%)	0.65	8 (13%) 3 2	46, 73, 112, 150	0
28	A4	69/71 (97%)	1.01	17 (24%) 1 0	59, 118, 206, 239	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
28	C4	69/71 (97%)	1.49	24 (34%) 0 0	81, 159, 207, 228	0
29	A5	59/60 (98%)	-0.30	0 100 100	8, 26, 40, 52	0
29	C5	59/60 (98%)	-0.09	0 100 100	27, 52, 90, 104	0
30	A6	53/54 (98%)	-0.42	0 100 100	23, 41, 55, 74	0
30	C6	53/54 (98%)	-0.33	0 100 100	41, 61, 80, 106	0
31	A7	48/49 (97%)	-0.24	1 (2%) 64 54	14, 24, 69, 134	1 (2%)
31	C7	48/49 (97%)	0.01	2 (4%) 37 26	26, 40, 96, 119	0
32	A8	64/65 (98%)	-0.30	0 100 100	16, 29, 45, 65	1 (1%)
32	C8	64/65 (98%)	-0.10	0 100 100	37, 52, 73, 85	0
33	A9	37/37 (100%)	-0.06	0 100 100	23, 35, 57, 68	1 (2%)
33	C9	37/37 (100%)	0.87	5 (13%) 3 2	45, 79, 96, 127	0
34	BA	1495/1521 (98%)	0.47	95 (6%) 20 12	31, 85, 186, 337	0
34	DA	1501/1521 (98%)	0.57	131 (8%) 11 6	39, 90, 196, 346	0
35	BB	231/256 (90%)	0.67	34 (14%) 3 1	43, 106, 173, 223	0
35	DB	231/256 (90%)	0.81	34 (14%) 3 1	71, 125, 176, 215	0
36	BC	206/239 (86%)	1.14	43 (20%) 1 1	56, 119, 174, 197	0
36	DC	206/239 (86%)	1.42	54 (26%) 1 0	69, 136, 182, 212	0
37	BD	208/209 (99%)	0.32	9 (4%) 36 26	44, 87, 138, 196	0
37	DD	208/209 (99%)	0.22	3 (1%) 75 69	59, 87, 136, 201	0
38	BE	148/162 (91%)	-0.07	1 (0%) 87 83	35, 73, 105, 128	0
38	DE	148/162 (91%)	0.04	2 (1%) 75 69	50, 81, 117, 182	0
39	BF	100/101 (99%)	-0.13	0 100 100	56, 86, 117, 138	0
39	DF	100/101 (99%)	-0.01	2 (2%) 65 56	48, 87, 115, 134	0
40	BG	155/156 (99%)	1.20	34 (21%) 1 1	68, 113, 183, 226	0
40	DG	155/156 (99%)	1.77	44 (28%) 1 0	72, 131, 194, 222	0
41	BH	137/138 (99%)	0.09	3 (2%) 62 52	47, 73, 99, 119	0
41	DH	137/138 (99%)	0.09	1 (0%) 87 83	57, 81, 111, 140	0
42	BI	127/128 (99%)	1.90	51 (40%) 0 0	65, 125, 167, 199	0
42	DI	127/128 (99%)	2.47	70 (55%) 0 0	91, 146, 193, 216	0
43	BJ	97/105 (92%)	1.64	35 (36%) 0 0	83, 131, 186, 215	0
43	DJ	96/105 (91%)	2.20	47 (48%) 0 0	92, 151, 200, 234	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
44	BK	114/129 (88%)	0.15	0 100 100	35, 79, 125, 151	0
44	DK	114/129 (88%)	0.23	2 (1%) 69 60	52, 91, 116, 177	0
45	BL	122/132 (92%)	-0.24	0 100 100	37, 60, 78, 113	0
45	DL	122/132 (92%)	-0.09	1 (0%) 86 81	46, 72, 96, 117	0
46	BM	117/126 (92%)	1.67	36 (30%) 0 0	77, 134, 182, 211	0
46	DM	122/126 (96%)	2.08	44 (36%) 0 0	94, 151, 201, 275	0
47	BN	60/61 (98%)	1.09	12 (20%) 1 1	67, 113, 146, 168	0
47	DN	60/61 (98%)	1.88	23 (38%) 0 0	98, 137, 179, 200	0
48	BO	88/89 (98%)	-0.12	1 (1%) 80 74	36, 71, 106, 121	0
48	DO	88/89 (98%)	0.02	1 (1%) 80 74	47, 71, 107, 153	0
49	BP	82/88 (93%)	0.42	1 (1%) 79 72	49, 80, 119, 171	0
49	DP	82/88 (93%)	0.48	5 (6%) 22 14	54, 78, 111, 153	0
50	BQ	99/105 (94%)	0.06	1 (1%) 82 77	44, 73, 99, 124	0
50	DQ	99/105 (94%)	0.03	0 100 100	44, 78, 104, 119	0
51	BR	68/88 (77%)	0.34	5 (7%) 15 8	42, 81, 123, 136	0
51	DR	68/88 (77%)	0.30	1 (1%) 74 67	53, 83, 128, 145	0
52	BS	84/93 (90%)	3.22	53 (63%) 0 0	97, 145, 198, 212	0
52	DS	83/93 (89%)	3.31	52 (62%) 0 0	90, 165, 216, 226	0
53	BT	96/106 (90%)	0.21	2 (2%) 64 54	62, 85, 122, 162	0
53	DT	96/106 (90%)	0.35	3 (3%) 49 38	58, 86, 135, 157	0
54	BU	23/27 (85%)	2.27	10 (43%) 0 0	62, 119, 158, 177	0
54	DU	23/27 (85%)	2.25	12 (52%) 0 0	92, 134, 173, 189	0
55	BV	7/18 (38%)	2.79	4 (57%) 0 0	53, 88, 211, 226	0
55	DV	6/18 (33%)	3.00	4 (66%) 0 0	84, 106, 214, 225	0
56	BW	69/76 (90%)	0.89	5 (7%) 16 9	38, 72, 106, 212	0
56	BY	67/76 (88%)	8.76	66 (98%) 0 0	82, 289, 329, 354	0
56	DW	69/76 (90%)	1.18	10 (14%) 3 2	54, 98, 141, 254	0
56	DY	66/76 (86%)	10.03	66 (100%) 0 0	213, 296, 333, 355	0
57	BZ	730/758 (96%)	0.17	36 (4%) 30 20	36, 79, 135, 190	0
57	DZ	730/758 (96%)	0.64	97 (13%) 4 2	36, 102, 169, 225	0
All	All	22825/23898 (95%)	0.61	2256 (9%) 8 4	5, 71, 189, 364	25 (0%)

All (2256) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	AC	159	ALA	36.4
3	CC	27	ALA	33.8
3	AC	57	GLN	33.4
3	CC	68	GLY	33.3
3	CC	172	ILE	32.3
3	CC	171	ALA	31.0
56	BY	35	A	29.7
56	DY	38	A	29.5
3	CC	67	HIS	29.4
3	AC	58	ASN	27.4
3	CC	180	SER	27.3
3	AC	200	HIS	27.1
3	AC	174	ALA	26.9
3	AC	59	VAL	26.6
3	CC	227	PRO	25.9
3	AC	173	HIS	24.1
3	CC	173	HIS	24.0
3	CC	181	PHE	23.6
3	CC	159	ALA	23.4
3	AC	166	ASN	23.4
56	BY	36	A	22.8
3	AC	67	HIS	22.7
3	CC	175	PRO	22.2
3	CC	164	PHE	22.0
3	CC	178	LYS	22.0
3	AC	39	ASP	22.0
46	DM	123	ALA	21.4
3	CC	41	THR	21.2
3	CC	39	ASP	21.1
3	CC	166	ASN	21.1
3	AC	164	PHE	21.0
3	CC	170	GLY	20.9
3	CC	57	GLN	20.8
10	CL	19	PRO	20.3
3	AC	177	GLY	20.2
3	CC	70	GLY	20.0
3	CC	179	ALA	20.0
3	AC	66	PRO	20.0
3	CC	4	HIS	19.5
40	DG	83	ALA	19.3
56	DY	29	G	19.2
3	CC	204	GLY	19.2

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Mol	Chain	Res	Type	RSRZ
1	CA	2123	G	18.9
10	CL	14	ALA	18.8
3	AC	172	ILE	18.7
3	AC	163	GLU	18.6
3	CC	226	ASN	18.6
3	AC	196	ALA	18.5
1	CA	2124	G	18.3
56	BY	33	U	18.1
1	CA	2179	C	18.0
56	DY	36	A	17.9
3	AC	63	VAL	17.8
1	CA	2122	U	17.7
3	CC	213	VAL	17.6
3	CC	200	HIS	17.5
3	CC	183	PRO	17.5
3	AC	176	VAL	17.5
3	CC	189	ASN	17.3
10	CL	13	PRO	17.3
1	AA	2152	U	17.2
3	CC	199	ALA	17.2
3	CC	60	ARG	17.1
1	AA	2181	G	17.1
3	CC	177	GLY	17.0
3	AC	25	GLU	16.9
1	CA	2115	G	16.9
3	CC	52	PRO	16.9
3	CC	71	LYS	16.9
3	CC	212	SER	16.9
56	DY	34	G	16.8
3	AC	23	ILE	16.8
3	AC	44	VAL	16.8
56	BY	34	G	16.7
3	CC	69	LEU	16.5
10	CL	24	GLY	16.5
3	CC	44	VAL	16.5
1	AA	2138	G	16.5
1	CA	2146	C	16.4
3	CC	165	ARG	16.4
3	AC	56	ASP	16.2
3	CC	160	GLY	16.2
3	AC	52	PRO	16.2
1	AA	2182	G	16.1

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Mol	Chain	Res	Type	RSRZ
10	CL	20	ALA	16.1
34	BA	1030(B)	C	16.0
3	CC	202	PRO	16.0
1	CA	2121	G	15.9
3	AC	167	ASP	15.8
1	AA	2131	U	15.8
56	BY	42	C	15.6
10	CL	11	GLN	15.6
3	CC	66	PRO	15.6
1	CA	2110	G	15.5
3	CC	185	LYS	15.5
3	AC	55	SER	15.5
3	CC	182	PRO	15.5
10	CL	10	LEU	15.5
3	AC	198	GLU	15.5
9	CK	50	ARG	15.4
3	AC	186	LEU	15.4
1	CA	2112	G	15.3
3	AC	171	ALA	15.2
56	DY	42	C	15.1
3	CC	59	VAL	15.1
3	CC	163	GLU	15.1
56	DY	62	C	15.0
3	AC	170	GLY	15.0
10	CL	12	LEU	15.0
3	AC	27	ALA	14.9
3	AC	26	ALA	14.9
3	AC	187	ALA	14.8
3	CC	56	ASP	14.8
3	AC	162	ILE	14.8
1	AA	2188	G	14.8
1	CA	2127	G	14.7
1	CA	2168	G	14.7
3	AC	189	ASN	14.6
3	CC	21	TYR	14.6
1	AA	2137	G	14.6
3	CC	186	LEU	14.6
46	DM	124	PRO	14.6
3	CC	192	ALA	14.5
1	AA	2132	G	14.4
3	AC	190	ILE	14.2
3	AC	183	PRO	14.2

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Mol	Chain	Res	Type	RSRZ
56	DY	1	G	14.2
10	CL	37	PHE	14.1
1	AA	2136	A	14.1
3	CC	23	ILE	14.1
1	AA	2151	C	14.1
3	CC	176	VAL	14.0
1	CA	2159	G	13.9
56	DY	72	C	13.9
56	BY	1	G	13.9
3	CC	20	VAL	13.9
56	DY	40	C	13.9
1	AA	2135	U	13.8
56	DY	64	A	13.8
10	CL	115	LEU	13.7
1	AA	2190	G	13.7
1	CA	2117	A	13.7
1	AA	2145	G	13.7
34	DA	1030(B)	C	13.6
3	CC	167	ASP	13.6
56	DY	30	G	13.5
1	CA	2139	C	13.4
56	DY	2	C	13.4
56	DY	56	C	13.4
1	AA	2201	C	13.3
1	CA	2111	C	13.2
1	AA	2167	C	13.1
56	BY	24	G	13.1
56	BY	23	A	13.0
3	CC	25	GLU	13.0
3	CC	203	GLU	12.9
56	BY	38	A	12.9
3	CC	174	ALA	12.9
3	CC	65	LEU	12.9
56	DY	33	U	12.8
10	CL	5	VAL	12.8
1	AA	2202	U	12.8
3	AC	191	ARG	12.8
3	CC	42	VAL	12.8
3	AC	192	ALA	12.8
56	DY	41	C	12.7
10	AL	52	ILE	12.7
3	AC	60	ARG	12.7

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Mol	Chain	Res	Type	RSRZ
3	CC	225	ILE	12.7
3	AC	195	ARG	12.7
52	DS	30	LEU	12.6
3	AC	207	GLY	12.6
1	AA	2163	G	12.6
56	DY	65	G	12.5
10	AL	7	VAL	12.5
3	CC	195	ARG	12.4
56	BY	70	G	12.4
3	AC	70	GLY	12.4
1	CA	2147	G	12.4
56	DY	28	G	12.3
3	AC	65	LEU	12.2
1	AA	2156	A	12.2
3	AC	204	GLY	12.2
3	AC	41	THR	12.2
56	DY	63	G	12.2
3	AC	69	LEU	12.2
10	CL	28	GLY	12.1
1	CA	2135	A	12.1
1	CA	2104	G	12.1
1	AA	2164	C	12.1
56	DY	3	C	12.1
3	AC	203	GLU	12.1
1	AA	2126	G	12.0
3	AC	199	ALA	12.0
10	CL	8	VAL	12.0
56	DY	61	C	11.9
3	CC	196	ALA	11.8
56	DY	74	C	11.8
3	AC	64	SER	11.8
56	DY	35	A	11.7
1	CA	2113	U	11.7
56	DY	53	G	11.7
3	AC	21	TYR	11.7
3	AC	68	GLY	11.7
1	AA	2168	C	11.6
10	CL	15	GLY	11.6
3	AC	188	ASP	11.6
3	AC	185	LYS	11.5
56	BY	29	G	11.5
3	AC	178	LYS	11.5

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Mol	Chain	Res	Type	RSRZ
56	DY	52	G	11.5
40	DG	79	ARG	11.4
56	BY	2	C	11.4
3	AC	4	HIS	11.4
1	CA	2109	U	11.4
10	CL	3	LYS	11.4
1	AA	2165	C	11.3
1	CA	2170	A	11.3
3	AC	226	ASN	11.3
10	CL	9	LYS	11.3
40	DG	78	ARG	11.3
1	AA	2134	G	11.3
10	CL	22	PRO	11.3
56	BY	63	G	11.3
3	CC	9	ARG	11.3
57	DZ	576	ASP	11.3
10	CL	114	ASP	11.2
3	AC	24	ASP	11.2
34	BA	1003	G	11.2
34	DA	1001(A)	G	11.2
1	AA	2169	G	11.2
56	DY	31	A	11.2
42	DI	30	GLY	11.1
10	CL	25	PRO	11.1
3	AC	20	VAL	11.0
3	CC	58	ASN	11.0
3	AC	165	ARG	11.0
46	DM	121	LYS	11.0
3	CC	61	GLY	10.9
3	AC	201	LYS	10.9
1	CA	2145	C	10.8
1	AA	2139	A	10.8
10	CL	21	PRO	10.8
46	DM	120	LYS	10.8
9	CK	89	ALA	10.8
56	BY	62	C	10.8
56	DY	6	G	10.7
3	AC	182	PRO	10.7
24	C0	3	HIS	10.7
3	CC	51	ASP	10.6
3	CC	31	LYS	10.6
56	BY	73	A	10.6

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Mol	Chain	Res	Type	RSRZ
9	AK	51	LEU	10.6
1	CA	2134	A	10.5
56	DY	71	G	10.5
1	CA	2105	C	10.5
56	BY	71	G	10.5
10	CL	136	VAL	10.5
1	CA	2166	G	10.5
1	CA	2120	G	10.4
3	AC	205	ALA	10.3
10	CL	18	THR	10.3
56	DY	73	A	10.3
56	BY	26	A	10.3
56	DY	75	C	10.2
9	CK	53	VAL	10.2
1	AA	2203	G	10.2
40	DG	156	TRP	10.2
10	CL	7	VAL	10.1
1	AA	2162	C	10.1
3	CC	207	GLY	10.1
56	BY	22	G	10.1
3	CC	210	LEU	10.1
1	CA	2177	C	10.1
3	CC	40	GLU	10.1
3	AC	51	ASP	10.1
3	CC	46	ALA	10.0
1	CA	2155	G	10.0
3	CC	191	ARG	9.9
10	AL	2	LYS	9.9
56	DY	70	G	9.9
56	DY	4	C	9.9
3	AC	197	LEU	9.9
56	BY	61	C	9.9
10	AL	64	SER	9.9
10	AL	13	PRO	9.9
56	DY	51	U	9.9
9	CK	115	GLN	9.8
56	DY	57	G	9.8
1	CA	2165	G	9.8
3	AC	28	ARG	9.8
3	CC	62	THR	9.8
56	DY	5	G	9.8
56	DY	58	A	9.8

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Mol	Chain	Res	Type	RSRZ
10	AL	60	TYR	9.7
56	BY	56	C	9.7
56	DY	24	G	9.7
1	CA	2160	G	9.7
3	CC	10	ALA	9.7
56	BY	28	G	9.7
10	CL	50	ASP	9.6
10	CL	31	GLY	9.6
3	AC	160	GLY	9.6
3	CC	54	ARG	9.6
10	CL	29	GLN	9.6
1	AA	2191	A	9.6
56	BY	72	C	9.5
1	AA	935	C	9.5
1	CA	2136	C	9.5
3	CC	162	ILE	9.5
3	CC	28	ARG	9.5
56	DY	18	G	9.5
3	CC	188	ASP	9.5
57	DZ	575	VAL	9.4
34	DA	1030	C	9.4
56	DY	19	G	9.3
3	CC	64	SER	9.3
34	DA	1030(D)	A	9.3
56	BY	27	G	9.3
1	AA	2154	U	9.2
3	AC	48	LEU	9.2
1	CA	2180	U	9.2
3	CC	193	PHE	9.2
56	BY	19	G	9.2
1	CA	2178	C	9.2
10	CL	2	LYS	9.2
1	AA	2187	G	9.1
56	BY	57	G	9.1
52	BS	56	GLN	9.1
3	AC	169	THR	9.1
10	CL	4	VAL	9.1
3	AC	224	ARG	9.1
3	CC	194	ILE	9.0
3	CC	205	ALA	9.0
1	AA	2166	U	9.0
3	CC	5	GLY	9.0

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Mol	Chain	Res	Type	RSRZ
52	DS	63	THR	9.0
3	AC	211	ARG	8.9
3	CC	198	GLU	8.9
3	CC	211	ARG	8.9
10	CL	122	ALA	8.9
56	DY	14	A	8.8
22	AY	1	MET	8.8
10	CL	127	ILE	8.8
56	BY	30	G	8.8
56	DY	60	U	8.8
10	CL	99	ILE	8.8
3	CC	184	GLU	8.8
10	AL	14	ALA	8.8
1	AA	2130	C	8.8
10	AL	49	GLY	8.8
3	AC	218	THR	8.7
56	DY	22	G	8.7
56	BY	40	C	8.7
1	CA	2169	A	8.6
10	AL	21	PRO	8.6
1	CA	2181	G	8.6
1	AA	2157	A	8.6
1	CA	2133	G	8.6
3	AC	175	PRO	8.5
3	AC	161	ARG	8.5
10	AL	15	GLY	8.5
10	CL	17	ALA	8.5
10	AL	16	LYS	8.5
3	CC	197	LEU	8.5
57	DZ	507	TYR	8.5
52	BS	57	HIS	8.5
1	CA	2128	C	8.5
10	CL	69	THR	8.5
1	CA	2167	U	8.4
1	AA	2155	G	8.4
40	DG	84	ASN	8.4
3	AC	202	PRO	8.4
34	DA	1036	G	8.4
3	AC	219	MET	8.4
10	CL	41	PHE	8.3
10	CL	138	VAL	8.4
34	DA	1030(A)	G	8.3

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Mol	Chain	Res	Type	RSRZ
3	CC	55	SER	8.3
3	CC	24	ASP	8.3
56	DY	26	A	8.3
1	CA	2154	G	8.3
3	CC	63	VAL	8.3
42	BI	80	GLY	8.3
56	DY	76	A	8.3
10	CL	96	VAL	8.2
56	BY	64	A	8.2
56	BY	53	G	8.2
57	DZ	538	TYR	8.2
40	DG	80	VAL	8.2
3	AC	61	GLY	8.2
3	CC	38	PHE	8.1
9	CK	96	PHE	8.1
3	CC	190	ILE	8.1
56	BY	74	C	8.1
10	CL	49	GLY	8.1
1	CA	2129	C	8.1
52	DS	29	ARG	8.1
34	DA	1033	G	8.1
3	AC	210	LEU	8.0
34	BA	1002	G	8.0
40	DG	82	GLY	8.0
56	DY	15	G	8.0
3	AC	71	LYS	8.0
3	CC	53	ARG	8.0
24	C0	2	ALA	7.9
9	CK	51	LEU	7.9
10	AL	23	VAL	7.9
10	AL	27	LEU	7.9
52	DS	69	HIS	7.9
56	DY	50	U	7.8
46	DM	6	GLY	7.8
3	CC	26	ALA	7.8
3	CC	8	TYR	7.8
56	BY	14	A	7.8
56	BY	58	A	7.8
3	CC	187	ALA	7.8
10	CL	54	PRO	7.8
56	BY	5	G	7.7
1	CA	2143	C	7.7

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Mol	Chain	Res	Type	RSRZ
9	CK	49	ALA	7.7
34	BA	1030	C	7.7
3	AC	181	PHE	7.7
1	AA	2142	G	7.7
1	CA	2144	U	7.7
34	BA	1029	C	7.7
1	AA	2133	C	7.7
1	AA	2199	C	7.7
47	DN	25	VAL	7.7
46	BM	2	ALA	7.7
34	BA	1036	G	7.7
3	CC	219	MET	7.6
10	CL	48	MET	7.6
10	CL	57	ILE	7.6
1	CA	2126	A	7.6
3	CC	6	LYS	7.6
3	AC	193	PHE	7.6
9	CK	114	GLY	7.6
40	BG	156	TRP	7.6
1	CA	2137	C	7.6
52	BS	71	LEU	7.6
10	CL	63	ARG	7.6
56	BY	43	C	7.5
3	CC	22	THR	7.5
10	AL	12	LEU	7.5
9	AK	53	VAL	7.5
9	AK	49	ALA	7.5
1	AA	2161	C	7.5
9	AK	90	ALA	7.5
10	CL	110	GLN	7.5
40	BG	79	ARG	7.4
34	DA	1030(C)	G	7.4
34	DA	1257	U	7.4
46	BM	24	GLY	7.4
56	BY	75	C	7.4
10	CL	123	ALA	7.4
56	DY	43	C	7.4
3	AC	208	THR	7.4
1	AA	2192	A	7.3
9	CK	94	VAL	7.3
3	AC	54	ARG	7.3
52	BS	4	SER	7.3

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Mol	Chain	Res	Type	RSRZ
1	AA	2153	G	7.3
56	BY	65	G	7.3
1	AA	2177	G	7.3
56	BY	41	C	7.3
56	DY	27	G	7.3
56	BY	45	U	7.2
34	DA	1031	G	7.2
10	AL	4	VAL	7.2
9	AK	50	ARG	7.2
52	BS	62	ILE	7.2
1	AA	2180	A	7.2
52	DS	31	ILE	7.2
56	BW	44	G	7.2
43	DJ	74	ILE	7.2
10	CL	30	HIS	7.2
34	DA	1001	A	7.2
42	BI	30	GLY	7.1
56	BY	3	C	7.1
52	BS	51	VAL	7.1
1	AA	2186	C	7.1
56	BY	69	G	7.1
56	DY	21	A	7.1
1	AA	2183	C	7.1
1	CA	2138	C	7.1
3	AC	206	LYS	7.1
56	DY	13	C	7.1
1	CA	2156	G	7.1
3	AC	53	ARG	7.1
10	CL	61	ALA	7.1
43	DJ	27	ALA	7.1
3	CC	208	THR	7.1
3	CC	50	ILE	7.0
1	AA	2189	U	7.0
10	CL	26	ALA	7.0
46	DM	119	GLY	7.0
56	BY	15	G	7.0
3	CC	43	GLU	7.0
40	BG	81	GLY	7.0
1	AA	2170	G	7.0
1	CA	2174	C	7.0
9	AK	104	ILE	6.9
42	DI	62	TYR	6.9

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Mol	Chain	Res	Type	RSRZ
1	CA	2114	A	6.9
10	CL	62	ASP	6.9
34	DA	1032	G	6.9
42	DI	49	PRO	6.9
57	DZ	574	GLU	6.9
56	BY	4	C	6.9
56	DY	44	G	6.9
10	CL	116	ASN	6.9
3	AC	179	ALA	6.8
10	CL	65	PHE	6.8
34	BA	1030(A)	G	6.8
34	DA	1002	G	6.8
1	AA	2141	A	6.8
1	AA	2200	C	6.8
3	AC	209	PHE	6.8
56	BY	52	G	6.8
56	BY	20	U	6.8
28	A4	65	ASP	6.8
1	CA	2141	G	6.8
46	DM	122	LYS	6.8
35	DB	228	GLY	6.7
1	CA	2140	C	6.7
56	DY	25	C	6.7
57	BZ	538	TYR	6.7
10	AL	65	PHE	6.7
10	AL	66	THR	6.7
3	AC	194	ILE	6.7
34	BA	1028	C	6.7
3	CC	7	ARG	6.6
9	AK	88	ALA	6.6
3	AC	31	LYS	6.6
3	AC	220	GLY	6.6
36	DC	155	GLY	6.6
34	BA	1030(C)	G	6.6
42	DI	105	ASP	6.6
56	BY	21	A	6.6
3	CC	49	GLY	6.6
56	DY	66	U	6.6
52	DS	4	SER	6.5
10	AL	62	ASP	6.5
52	BS	64	GLU	6.5
9	CK	125	LEU	6.5

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Mol	Chain	Res	Type	RSRZ
10	AL	58	THR	6.5
52	DS	56	GLN	6.5
8	CH	29	PRO	6.5
52	DS	82	GLY	6.5
28	C4	62	ARG	6.5
1	CA	2142	C	6.5
1	AA	2143	G	6.5
56	BY	11	C	6.4
56	BY	6	G	6.4
57	DZ	530	VAL	6.4
42	DI	5	TYR	6.4
1	CA	2108	C	6.4
10	AL	67	PHE	6.4
42	DI	18	PHE	6.4
1	AA	2173	G	6.4
57	DZ	573	HIS	6.3
23	CZ	156	LYS	6.3
56	BY	44	G	6.3
1	CA	888	C	6.3
40	DG	37	ASN	6.3
10	CL	137	GLU	6.3
1	AA	2160	C	6.3
43	BJ	99	LYS	6.3
56	BY	12	U	6.3
3	AC	22	THR	6.3
54	BU	17	THR	6.3
24	C0	5	LYS	6.3
10	CL	56	GLU	6.2
34	BA	1035	A	6.2
35	BB	228	GLY	6.2
10	AL	48	MET	6.2
10	AL	56	GLU	6.2
10	AL	61	ALA	6.2
40	BG	78	ARG	6.2
52	DS	67	VAL	6.2
3	CC	47	LYS	6.2
3	CC	201	LYS	6.2
10	AL	3	LYS	6.2
52	DS	48	THR	6.2
3	AC	10	ALA	6.2
9	CK	117	LEU	6.2
52	DS	49	ILE	6.2

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Mol	Chain	Res	Type	RSRZ
10	AL	25	PRO	6.2
35	BB	136	VAL	6.1
1	AA	2176	G	6.1
56	BY	13	C	6.1
1	CA	2116	G	6.1
56	DY	23	A	6.1
3	AC	29	LEU	6.1
52	BS	30	LEU	6.1
24	A0	3	HIS	6.1
10	CL	128	ALA	6.1
47	DN	2	ALA	6.1
10	AL	124	ALA	6.1
56	DY	47	U	6.1
10	CL	52	ILE	6.1
1	AA	2196	C	6.1
52	BS	29	ARG	6.1
56	BY	9	A	6.1
1	AA	2158	C	6.0
3	AC	180	SER	6.0
40	BG	34	GLY	6.0
3	AC	227	PRO	6.0
3	AC	35	THR	6.0
56	DY	67	C	6.0
1	CA	229	A	6.0
34	BA	1286	A	6.0
36	DC	87	LEU	6.0
3	CC	18	ASN	6.0
52	DS	62	ILE	6.0
40	DG	154	TYR	6.0
8	CH	27	LYS	6.0
24	C0	4	LYS	6.0
3	CC	35	THR	5.9
10	CL	68	VAL	5.9
43	BJ	98	ILE	5.9
34	BA	1001(A)	G	5.9
52	BS	65	ASN	5.9
52	BS	84	GLY	5.9
3	AC	38	PHE	5.9
10	AL	10	LEU	5.9
46	DM	82	MET	5.9
9	CK	100	ASN	5.8
3	AC	45	HIS	5.8

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Mol	Chain	Res	Type	RSRZ
28	C4	68	ARG	5.8
55	BV	12	A	5.8
56	BY	7	A	5.8
56	DY	69	G	5.8
52	DS	61	TYR	5.8
54	BU	18	TYR	5.8
10	CL	104	VAL	5.8
47	DN	38	GLY	5.8
3	AC	62	THR	5.8
54	DU	17	THR	5.8
40	BG	80	VAL	5.7
10	AL	37	PHE	5.7
52	DS	66	MET	5.7
28	C4	64	GLY	5.7
3	AC	225	ILE	5.7
9	CK	90	ALA	5.7
1	AA	1555	C	5.7
43	DJ	100	THR	5.7
56	DY	12	U	5.7
10	CL	55	VAL	5.7
52	BS	74	PHE	5.7
34	BA	1257	U	5.7
56	BY	25	C	5.6
9	CK	77	PRO	5.6
47	DN	17	LYS	5.6
1	CA	2152	G	5.6
22	CY	1	MET	5.6
1	CA	2164	C	5.6
34	BA	1004	A	5.6
55	DV	14	A	5.6
52	BS	59	PRO	5.6
36	DC	39	ILE	5.6
56	BY	10	G	5.6
1	CA	2106	G	5.6
3	AC	212	SER	5.6
43	DJ	39	PRO	5.6
34	DA	1286	A	5.6
35	BB	135	GLN	5.6
8	CH	13	LYS	5.6
56	BY	49	C	5.5
1	CA	2182	G	5.5
34	BA	1033	G	5.5

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Mol	Chain	Res	Type	RSRZ
56	BY	67	C	5.5
34	BA	1030(D)	A	5.5
52	BS	72	GLY	5.5
9	CK	97	ALA	5.5
43	DJ	20	ALA	5.5
52	DS	43	GLU	5.5
34	DA	1137	C	5.5
42	DI	9	ARG	5.5
10	CL	126	MET	5.5
56	BY	31	A	5.5
10	CL	98	ARG	5.4
46	DM	93	ARG	5.4
46	BM	95	GLY	5.4
10	AL	46	ALA	5.4
56	BY	18	G	5.4
3	AC	9	ARG	5.4
9	CK	25	PHE	5.4
28	A4	68	ARG	5.4
10	CL	59	ILE	5.4
1	AA	2210	C	5.4
10	AL	22	PRO	5.4
52	DS	79	THR	5.4
56	BY	47	U	5.4
46	DM	98	VAL	5.3
10	AL	8	VAL	5.3
1	AA	1221	G	5.3
46	BM	83	ASP	5.3
10	CL	135	GLY	5.3
34	DA	1256	A	5.3
35	BB	66	GLY	5.3
35	DB	222	ILE	5.3
9	CK	44	LEU	5.3
28	C4	51	ASP	5.3
10	AL	28	GLY	5.3
52	DS	42	PRO	5.3
34	DA	1034	G	5.3
1	CA	2125	G	5.3
52	DS	35	SER	5.3
1	AA	2123	G	5.3
1	AA	2175	G	5.3
1	AA	2184	G	5.3
1	CA	889	C	5.2

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Mol	Chain	Res	Type	RSRZ
56	DY	45	U	5.2
1	CA	2153	G	5.2
46	DM	87	TYR	5.2
3	AC	40	GLU	5.2
1	CA	2157	G	5.2
10	CL	134	MET	5.2
10	CL	93	ARG	5.2
36	BC	2	GLY	5.2
1	AA	696	C	5.2
3	AC	47	LYS	5.2
52	DS	28	LYS	5.2
42	DI	8	GLY	5.2
3	AC	34	ALA	5.2
10	AL	63	ARG	5.2
28	A4	55	ARG	5.2
1	AA	2144	U	5.2
36	DC	36	ASP	5.2
1	AA	934	A	5.1
1	AA	2179	G	5.1
57	DZ	231	TYR	5.1
56	BY	60	U	5.1
57	BZ	570	GLY	5.1
43	DJ	6	ILE	5.1
1	CA	2175	C	5.1
42	DI	19	LEU	5.1
9	CK	52	PHE	5.1
43	DJ	10	GLY	5.1
42	DI	104	ARG	5.1
1	AA	2125	C	5.1
36	BC	206	GLU	5.1
56	BY	50	U	5.1
34	DA	1035	A	5.1
36	BC	39	ILE	5.1
43	DJ	85	LEU	5.1
34	BA	1034	G	5.1
43	DJ	73	ASP	5.1
1	CA	2158	A	5.1
42	DI	17	VAL	5.0
1	CA	2151	G	5.0
9	CK	26	LEU	5.0
57	DZ	567	LEU	5.0
10	AL	136	VAL	5.0

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Mol	Chain	Res	Type	RSRZ
36	BC	154	SER	5.0
10	AL	20	ALA	5.0
42	DI	43	ALA	5.0
52	DS	3	ARG	5.0
1	AA	933	C	5.0
56	BY	51	U	5.0
40	DG	77	SER	5.0
36	BC	155	GLY	5.0
52	DS	12	ASP	5.0
10	CL	70	LYS	5.0
34	DA	1021	G	5.0
57	BZ	495	GLY	5.0
56	DY	11	C	5.0
34	BA	1026	G	4.9
10	AL	26	ALA	4.9
10	AL	50	ASP	4.9
34	BA	999	C	4.9
57	BZ	499	ARG	4.9
3	AC	50	ILE	4.9
28	A4	59	PHE	4.9
3	CC	218	THR	4.9
1	AA	2159	C	4.9
1	AA	2178	G	4.9
34	BA	1031	G	4.9
52	BS	41	VAL	4.9
52	BS	22	LEU	4.9
57	DZ	536	LYS	4.9
46	DM	45	VAL	4.9
56	DY	7	A	4.9
10	AL	19	PRO	4.9
36	DC	23	TYR	4.9
36	DC	64	VAL	4.9
10	CL	60	TYR	4.9
7	CG	48	GLU	4.9
3	AC	223	VAL	4.8
46	BM	93	ARG	4.8
46	BM	6	GLY	4.8
1	AA	932	C	4.8
10	CL	42	ASN	4.8
3	AC	42	VAL	4.8
8	CH	105	LEU	4.8
3	CC	169	THR	4.8

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Mol	Chain	Res	Type	RSRZ
56	DY	59	U	4.8
57	DZ	600	VAL	4.8
8	CH	97	ARG	4.8
35	BB	122	PHE	4.8
52	BS	31	ILE	4.8
52	DS	40	ILE	4.8
3	CC	168	LYS	4.8
34	DA	1042	G	4.8
46	DM	23	TYR	4.8
3	AC	5	GLY	4.7
1	AA	2127	C	4.7
1	AA	2207	C	4.7
46	DM	92	HIS	4.7
43	BJ	24	VAL	4.7
51	BR	85	LEU	4.7
10	AL	59	ILE	4.7
28	C4	69	LYS	4.7
10	CL	35	MET	4.7
40	BG	54	THR	4.7
52	BS	33	THR	4.7
34	BA	1138	G	4.7
56	DW	44	G	4.7
10	AL	9	LYS	4.7
10	CL	140	GLY	4.7
17	AT	38	ASN	4.7
43	DJ	96	ILE	4.7
1	AA	2172	U	4.7
3	CC	17	PRO	4.7
42	DI	42	ARG	4.7
52	BS	61	TYR	4.7
28	A4	54	GLY	4.7
36	BC	199	LYS	4.7
42	BI	61	ALA	4.6
42	DI	27	THR	4.6
1	CA	2176	A	4.6
52	DS	14	HIS	4.6
34	BA	1447	A	4.6
56	DY	9	A	4.6
46	BM	36	LYS	4.6
43	DJ	70	ARG	4.6
52	DS	7	LYS	4.6
1	AA	2208	G	4.6

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Mol	Chain	Res	Type	RSRZ
1	CA	2130	U	4.6
34	DA	1026	G	4.6
28	A4	66	SER	4.6
10	CL	125	ARG	4.6
57	BZ	530	VAL	4.6
40	BG	84	ASN	4.6
42	DI	81	ILE	4.6
9	CK	86	PRO	4.5
43	BJ	69	ASN	4.5
43	DJ	81	THR	4.5
1	AA	2149	G	4.5
10	AL	30	HIS	4.5
10	CL	85	GLU	4.5
34	DA	1043	C	4.5
42	DI	86	VAL	4.5
43	BJ	28	ARG	4.5
3	AC	46	ALA	4.5
35	BB	123	ALA	4.5
1	CA	1509	C	4.5
34	DA	1027	C	4.5
52	BS	32	LYS	4.5
52	BS	40	ILE	4.5
35	DB	232	PRO	4.5
42	DI	93	ARG	4.5
36	BC	89	GLU	4.5
52	BS	83	HIS	4.5
10	CL	139	VAL	4.5
57	BZ	229	LEU	4.5
36	DC	190	ARG	4.5
40	BG	55	GLY	4.5
36	BC	32	LEU	4.5
10	CL	23	VAL	4.5
1	CA	887	A	4.5
55	BV	13	A	4.5
1	CA	652(D)	C	4.5
10	CL	51	ALA	4.5
35	BB	118	LEU	4.5
10	CL	81	ALA	4.4
34	BA	1043	C	4.4
42	DI	31	GLN	4.4
46	DM	5	ALA	4.4
28	A4	69	LYS	4.4

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Mol	Chain	Res	Type	RSRZ
55	DV	13	A	4.4
9	CK	57	THR	4.4
37	DD	23	GLY	4.4
42	BI	98	PRO	4.4
1	CA	2101	G	4.4
40	DG	16	LEU	4.4
57	DZ	539	ILE	4.4
3	AC	217	THR	4.4
56	DW	17	C	4.4
56	DY	49	C	4.4
34	DA	1265	G	4.4
10	AL	132	ARG	4.4
10	CL	107	ILE	4.4
28	C4	49	PHE	4.4
56	DY	48	C	4.4
9	CK	131	MET	4.4
42	BI	17	VAL	4.4
28	C4	45	GLY	4.3
1	CA	652(B)	A	4.3
34	BA	1274	G	4.3
42	BI	63	ILE	4.3
42	DI	6	GLY	4.3
46	DM	80	ARG	4.3
1	AA	2204	G	4.3
3	AC	32	GLU	4.3
56	BY	66	U	4.3
47	BN	17	LYS	4.3
3	CC	214	TYR	4.3
43	DJ	8	LEU	4.3
54	DU	18	TYR	4.3
3	AC	222	SER	4.3
44	DK	17	GLY	4.3
10	AL	137	GLU	4.3
35	DB	135	GLN	4.3
47	DN	15	LYS	4.3
56	BY	48	C	4.3
10	CL	6	ALA	4.3
3	CC	11	LEU	4.3
35	DB	233	SER	4.3
46	BM	98	VAL	4.3
52	BS	60	VAL	4.3
10	AL	127	ILE	4.3

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Mol	Chain	Res	Type	RSRZ
37	BD	23	GLY	4.3
42	BI	81	ILE	4.3
34	BA	841	U	4.3
34	BA	1032	G	4.3
10	AL	6	ALA	4.2
9	CK	85	ASP	4.2
46	BM	23	TYR	4.2
3	CC	29	LEU	4.2
36	DC	63	ASN	4.2
36	DC	85	ARG	4.2
52	DS	41	VAL	4.2
9	CK	11	ALA	4.2
35	BB	233	SER	4.2
34	BA	1042	G	4.2
52	BS	58	VAL	4.2
42	DI	99	LEU	4.2
1	AA	2128	G	4.2
46	DM	84	ILE	4.2
49	DP	19	ILE	4.2
42	DI	98	PRO	4.2
34	DA	1370	G	4.2
57	BZ	541	ALA	4.2
42	BI	102	LEU	4.2
10	CL	100	THR	4.2
52	BS	45	VAL	4.2
24	C0	6	GLY	4.2
25	C1	2	SER	4.2
42	DI	95	LYS	4.2
10	AL	35	MET	4.2
10	CL	66	THR	4.2
10	CL	117	THR	4.2
52	DS	32	LYS	4.2
10	AL	33	ASN	4.2
34	BA	1136	U	4.2
34	DA	1040	U	4.2
34	BA	1021	G	4.1
36	BC	152	ILE	4.1
34	DA	1287	A	4.1
46	DM	51	ALA	4.1
40	BG	82	GLY	4.1
52	BS	50	ALA	4.1
1	AA	2174	G	4.1

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Mol	Chain	Res	Type	RSRZ
1	CA	2100	G	4.1
40	DG	18	TYR	4.1
57	DZ	411	VAL	4.1
54	BU	13	ILE	4.1
1	CA	2132	U	4.1
34	BA	1037	C	4.1
42	DI	66	ARG	4.1
3	AC	214	TYR	4.1
10	AL	76	TYR	4.1
54	DU	5	ASP	4.1
40	DG	27	ILE	4.1
34	DA	79	G	4.1
46	BM	28	ALA	4.1
40	DG	85	TYR	4.1
10	AL	24	GLY	4.0
1	CA	652(U)	G	4.0
46	DM	85	GLY	4.0
46	DM	86	CYS	4.0
52	BS	12	ASP	4.0
1	CA	2190	G	4.0
10	CL	132	ARG	4.0
34	DA	1018	C	4.0
42	BI	38	GLN	4.0
40	DG	17	VAL	4.0
52	BS	49	ILE	4.0
34	BA	1000	U	4.0
53	BT	9	ASN	4.0
52	BS	66	MET	4.0
42	BI	15	ALA	4.0
57	DZ	531	GLY	4.0
57	BZ	540	PRO	4.0
36	DC	101	LEU	4.0
52	BS	63	THR	4.0
46	BM	94	ARG	4.0
52	DS	50	ALA	4.0
1	AA	694	G	4.0
57	DZ	255	ILE	4.0
3	CC	220	GLY	4.0
9	CK	119	ALA	4.0
33	C9	37	GLY	4.0
46	BM	80	ARG	4.0
10	AL	31	GLY	4.0

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Mol	Chain	Res	Type	RSRZ
57	DZ	508	GLY	4.0
56	DY	68	C	3.9
43	DJ	78	ASN	3.9
36	BC	164	ARG	3.9
23	CZ	9	TYR	3.9
8	CH	26	VAL	3.9
34	BA	1137	C	3.9
9	CK	116	ILE	3.9
52	BS	8	GLY	3.9
3	CC	13	GLU	3.9
56	BY	59	U	3.9
1	CA	1847	A	3.9
34	DA	1029	C	3.9
55	DV	15	A	3.9
10	CL	89	HIS	3.9
1	AA	2150	C	3.9
34	BA	1027	C	3.9
40	DG	81	GLY	3.9
28	A4	57	GLU	3.9
43	BJ	97	GLU	3.9
9	CK	58	LEU	3.9
57	BZ	527	ASN	3.9
8	CH	113	VAL	3.9
10	CL	94	GLU	3.9
52	BS	2	PRO	3.9
40	DG	155	ARG	3.9
10	CL	58	THR	3.9
40	BG	89	MET	3.9
46	DM	117	VAL	3.9
43	BJ	41	PRO	3.9
3	AC	184	GLU	3.9
10	CL	32	ALA	3.9
34	DA	1045	C	3.9
1	AA	2171	G	3.9
56	DY	10	G	3.9
28	C4	50	VAL	3.8
43	DJ	98	ILE	3.8
36	BC	87	LEU	3.8
42	DI	7	THR	3.8
1	AA	2197	C	3.8
56	BY	68	C	3.8
57	DZ	430	ARG	3.8

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Mol	Chain	Res	Type	RSRZ
36	DC	77	ILE	3.8
52	BS	52	TYR	3.8
36	BC	91	LEU	3.8
1	CA	2148	G	3.8
34	DA	1024	G	3.8
42	BI	6	GLY	3.8
43	BJ	29	ARG	3.8
10	CL	45	THR	3.8
36	BC	57	ILE	3.8
43	DJ	72	VAL	3.8
46	BM	45	VAL	3.8
28	C4	63	TYR	3.8
35	DB	33	TYR	3.8
57	BZ	232	LEU	3.8
9	CK	130	THR	3.8
57	DZ	529	ILE	3.8
57	DZ	565	VAL	3.8
34	BA	1025	U	3.8
10	CL	105	LEU	3.8
52	BS	26	GLY	3.8
34	DA	1044	A	3.8
54	DU	16	GLY	3.8
3	CC	161	ARG	3.8
52	DS	57	HIS	3.7
3	CC	222	SER	3.7
46	BM	85	GLY	3.7
57	BZ	502	GLY	3.7
9	CK	54	ALA	3.7
43	BJ	34	VAL	3.7
57	DZ	537	GLU	3.7
1	AA	697	C	3.7
1	CA	1079	C	3.7
1	CA	2119	A	3.7
35	DB	139	LYS	3.7
42	DI	87	GLN	3.7
40	BG	47	CYS	3.7
1	CA	2185	C	3.7
34	BA	1044	A	3.7
3	CC	30	VAL	3.7
52	DS	60	VAL	3.7
1	CA	2189	U	3.7
3	CC	32	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
46	BM	62	ASN	3.7
7	CG	42	GLY	3.7
52	DS	26	GLY	3.7
57	DZ	633	GLY	3.7
36	DC	7	PRO	3.7
52	BS	48	THR	3.7
3	CC	15	VAL	3.7
35	DB	229	VAL	3.7
36	DC	57	ILE	3.7
10	CL	111	LYS	3.7
57	DZ	403	GLU	3.7
34	DA	1266	G	3.7
46	BM	26	GLY	3.7
22	CY	65	ALA	3.7
46	DM	43	THR	3.7
3	AC	43	GLU	3.7
10	CL	34	ILE	3.7
40	BG	57	GLU	3.7
3	AC	12	LEU	3.7
1	CA	886	C	3.7
34	DA	999	C	3.7
52	BS	85	LYS	3.7
42	DI	29	ASN	3.7
55	BV	14	A	3.7
10	AL	120	LEU	3.7
42	DI	103	THR	3.7
43	DJ	99	LYS	3.6
57	DZ	506	GLN	3.6
10	CL	43	ALA	3.6
10	CL	27	LEU	3.6
3	CC	37	LYS	3.6
42	BI	75	ASP	3.6
52	DS	46	GLY	3.6
1	AA	2122	G	3.6
10	CL	73	PRO	3.6
23	CZ	155	LEU	3.6
34	DA	1531	A	3.6
57	DZ	492	ASP	3.6
40	BG	56	GLN	3.6
43	BJ	20	ALA	3.6
9	CK	132	ASP	3.6
3	CC	34	ALA	3.6

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Mol	Chain	Res	Type	RSRZ
57	DZ	543	GLN	3.6
34	DA	1039	C	3.6
35	DB	230	VAL	3.6
1	CA	2103	C	3.6
57	DZ	491	VAL	3.6
57	DZ	540	PRO	3.6
1	AA	698	G	3.6
40	DG	76	ARG	3.6
10	AL	116	ASN	3.6
42	DI	73	GLN	3.6
46	DM	41	PRO	3.6
3	AC	30	VAL	3.6
10	AL	96	VAL	3.6
57	DZ	601	ILE	3.6
42	DI	75	ASP	3.6
46	DM	83	ASP	3.6
4	AD	276	LYS	3.5
9	CK	109	SER	3.5
57	DZ	234	GLY	3.5
57	DZ	566	THR	3.5
34	DA	1041	A	3.5
35	BB	131	PRO	3.5
42	BI	28	VAL	3.5
34	BA	1212	U	3.5
43	BJ	40	LEU	3.5
57	BZ	231	TYR	3.5
24	C0	65	GLY	3.5
36	BC	145	GLY	3.5
38	DE	22	GLY	3.5
9	CK	56	ASN	3.5
10	CL	87	GLY	3.5
46	BM	100	GLY	3.5
7	CG	152	LEU	3.5
34	DA	1264	C	3.5
9	CK	24	PHE	3.5
10	AL	70	LYS	3.5
40	DG	153	HIS	3.5
9	AK	131	MET	3.5
43	BJ	18	ALA	3.5
36	DC	131	ARG	3.5
43	BJ	35	SER	3.5
43	DJ	36	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
28	C4	67	TYR	3.5
23	CZ	60	GLU	3.5
36	BC	163	ALA	3.5
35	BB	37	ASN	3.5
3	CC	12	LEU	3.5
46	BM	82	MET	3.5
34	BA	1039	C	3.5
8	AH	2	SER	3.4
23	CZ	153	SER	3.4
3	CC	19	LYS	3.4
8	CH	102	ALA	3.4
9	CK	104	ILE	3.4
35	DB	123	ALA	3.4
42	DI	82	ALA	3.4
57	BZ	539	ILE	3.4
57	DZ	526	VAL	3.4
25	A1	2	SER	3.4
42	BI	19	LEU	3.4
42	BI	84	ALA	3.4
22	CY	55	TYR	3.4
8	CH	48	GLY	3.4
8	CH	95	ARG	3.4
36	BC	127	ARG	3.4
52	DS	64	GLU	3.4
34	DA	1037	C	3.4
1	CA	900	A	3.4
10	CL	16	LYS	3.4
36	DC	37	GLN	3.4
43	BJ	62	HIS	3.4
28	C4	29	PRO	3.4
10	AL	11	GLN	3.4
10	CL	64	SER	3.4
57	DZ	578	SER	3.4
10	AL	51	ALA	3.4
36	DC	121	ALA	3.4
36	BC	181	ASN	3.4
1	CA	2833	G	3.4
36	DC	129	ALA	3.4
46	BM	25	ILE	3.4
42	BI	97	LYS	3.4
52	BS	55	LYS	3.4
52	DS	34	TRP	3.4

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Mol	Chain	Res	Type	RSRZ
34	DA	1532	U	3.4
42	DI	102	LEU	3.4
47	DN	39	LEU	3.4
36	DC	189	ALA	3.4
43	BJ	10	GLY	3.4
57	DZ	681	LYS	3.4
9	CK	12	THR	3.4
8	CH	88	LEU	3.4
34	DA	204	U	3.4
34	DA	1164	G	3.4
8	CH	16	SER	3.4
10	CL	92	GLY	3.4
35	DB	130	ARG	3.3
47	DN	26	ARG	3.3
35	DB	231	GLU	3.3
42	BI	31	GLN	3.3
36	BC	200	ALA	3.3
34	BA	998	G	3.3
54	DU	6	ARG	3.3
1	AA	2211	U	3.3
34	BA	1353	G	3.3
57	DZ	224	ASP	3.3
57	DZ	577	SER	3.3
57	DZ	579	GLU	3.3
34	DA	1211	U	3.3
1	CA	2102	U	3.3
3	CC	217	THR	3.3
36	DC	158	GLY	3.3
42	BI	64	THR	3.3
42	BI	74	ILE	3.3
34	DA	1020	U	3.3
46	DM	95	GLY	3.3
1	AA	2193	A	3.3
16	CS	56	LEU	3.3
42	BI	20	ARG	3.3
17	CT	38	ASN	3.3
42	BI	45	ALA	3.3
1	CA	1099	G	3.3
34	DA	1003	G	3.3
34	DA	1022	G	3.3
9	CK	121	ASP	3.3
48	BO	89	GLY	3.3

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Mol	Chain	Res	Type	RSRZ
57	DZ	527	ASN	3.3
57	DZ	417	THR	3.3
42	DI	56	LEU	3.3
42	DI	67	GLY	3.3
35	BB	207	ALA	3.2
52	BS	75	ALA	3.2
28	A4	62	ARG	3.2
40	DG	89	MET	3.2
52	BS	39	THR	3.2
52	BS	44	MET	3.2
36	DC	91	LEU	3.2
42	DI	50	LEU	3.2
3	AC	19	LYS	3.2
36	DC	157	ILE	3.2
8	CH	17	VAL	3.2
8	CH	76	VAL	3.2
10	CL	38	VAL	3.2
40	BG	83	ALA	3.2
1	CA	878	A	3.2
52	BS	46	GLY	3.2
52	DS	71	LEU	3.2
57	BZ	494	GLU	3.2
35	DB	140	HIS	3.2
36	BC	202	ILE	3.2
43	BJ	38	ILE	3.2
52	DS	45	VAL	3.2
57	DZ	594	VAL	3.2
34	DA	1353	G	3.2
10	AL	134	MET	3.2
34	BA	1275	A	3.2
52	BS	42	PRO	3.2
9	CK	99	SER	3.2
10	CL	101	TRP	3.2
27	C3	6	VAL	3.2
36	DC	24	ALA	3.2
42	BI	29	ASN	3.2
46	DM	13	LYS	3.2
1	CA	2191	G	3.2
47	DN	28	GLY	3.2
28	C4	61	ARG	3.2
57	DZ	40	HIS	3.2
42	BI	41	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
10	CL	47	ASN	3.2
40	DG	28	ASN	3.2
10	AL	36	GLU	3.2
43	DJ	84	GLN	3.2
7	AG	49	ASP	3.2
57	DZ	197	ARG	3.2
9	AK	52	PHE	3.2
10	AL	45	THR	3.2
43	DJ	26	ALA	3.2
26	C2	7	ARG	3.2
47	BN	15	LYS	3.2
1	CA	2118	U	3.2
3	CC	36	ALA	3.2
10	AL	17	ALA	3.2
36	DC	80	GLY	3.2
35	DB	132	LYS	3.2
43	BJ	8	LEU	3.2
46	BM	84	ILE	3.2
10	AL	131	ALA	3.2
1	CA	883	G	3.2
47	DN	19	ARG	3.2
47	DN	29	ARG	3.2
22	CY	63	LYS	3.2
36	DC	32	LEU	3.2
34	DA	1019	C	3.2
8	CH	49	VAL	3.1
34	BA	1041	A	3.1
43	DJ	71	LEU	3.1
47	BN	16	PHE	3.1
52	DS	2	PRO	3.1
9	CK	60	ARG	3.1
43	BJ	36	GLY	3.1
3	AC	18	ASN	3.1
10	AL	47	ASN	3.1
57	BZ	507	TYR	3.1
34	DA	80	G	3.1
34	DA	1138	G	3.1
49	DP	59	TRP	3.1
1	CA	2163	C	3.1
8	CH	107	VAL	3.1
56	DW	4	C	3.1
57	DZ	404	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
23	CZ	112	ARG	3.1
1	AA	2124	U	3.1
3	CC	215	VAL	3.1
10	CL	97	GLY	3.1
40	DG	34	GLY	3.1
22	CY	91	GLU	3.1
28	A4	53	GLU	3.1
21	CX	69	TYR	3.1
36	DC	193	TYR	3.1
35	DB	124	SER	3.1
34	BA	1023	G	3.1
34	BA	1258	G	3.1
34	DA	723	U	3.1
43	BJ	72	VAL	3.1
52	BS	67	VAL	3.1
42	BI	78	LYS	3.1
1	CA	645	C	3.1
9	CK	74	LEU	3.1
42	DI	88	TYR	3.1
3	AC	168	LYS	3.1
31	C7	48	LYS	3.1
10	AL	98	ARG	3.1
42	DI	46	ALA	3.1
24	A0	7	LEU	3.1
34	BA	78	G	3.1
34	DA	1023	G	3.1
28	C4	59	PHE	3.1
57	BZ	393	ASP	3.1
3	AC	6	LYS	3.1
42	DI	63	ILE	3.1
3	AC	213	VAL	3.1
36	DC	89	GLU	3.1
1	CA	885	C	3.1
57	BZ	572	TYR	3.1
10	AL	34	ILE	3.1
43	BJ	5	ARG	3.1
43	DJ	59	SER	3.1
53	BT	45	GLN	3.1
9	AK	24	PHE	3.1
9	AK	74	LEU	3.1
10	AL	39	LYS	3.1
23	CZ	157	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
36	DC	72	LYS	3.1
52	DS	10	PHE	3.1
10	AL	110	GLN	3.1
34	DA	1291	G	3.1
34	DA	90	U	3.0
34	DA	1025	U	3.0
42	DI	79	LEU	3.0
34	BA	1006	C	3.0
34	DA	1270	C	3.0
36	DC	184	TYR	3.0
57	BZ	531	GLY	3.0
10	AL	86	LYS	3.0
35	BB	137	ARG	3.0
35	BB	138	LEU	3.0
34	BA	1001	A	3.0
34	BA	1246	C	3.0
36	BC	98	ASN	3.0
43	DJ	68	HIS	3.0
52	DS	47	HIS	3.0
7	CG	159	VAL	3.0
10	CL	53	VAL	3.0
34	BA	1038	C	3.0
35	DB	131	PRO	3.0
46	BM	37	THR	3.0
43	BJ	25	GLU	3.0
43	DJ	25	GLU	3.0
13	CP	137	LYS	3.0
40	DG	36	LYS	3.0
34	BA	1350	A	3.0
35	BB	132	LYS	3.0
35	DB	122	PHE	3.0
43	DJ	65	LEU	3.0
34	DA	1013	G	3.0
46	BM	110	ARG	3.0
10	AL	90	LYS	3.0
5	CE	78	LEU	3.0
34	DA	1209	C	3.0
36	DC	192	THR	3.0
54	DU	4	GLY	3.0
37	BD	144	ASP	3.0
3	CC	223	VAL	3.0
8	CH	19	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
9	CK	55	LYS	3.0
35	DB	112	VAL	3.0
51	BR	21	LYS	3.0
34	DA	998	G	3.0
42	DI	12	GLU	3.0
43	DJ	64	GLU	3.0
57	DZ	626	ALA	3.0
36	BC	38	ARG	3.0
10	CL	46	ALA	3.0
34	DA	1446	U	3.0
1	AA	2185	C	3.0
28	A4	64	GLY	3.0
34	BA	1174	G	3.0
8	CH	106	THR	3.0
10	CL	90	LYS	3.0
54	BU	22	ARG	2.9
57	DZ	602	LEU	2.9
34	DA	1314	C	2.9
56	BW	17	C	2.9
10	AL	57	ILE	2.9
28	C4	31	ILE	2.9
52	DS	11	VAL	2.9
23	CZ	150	LEU	2.9
34	BA	202	U	2.9
43	DJ	21	GLN	2.9
34	BA	1277	C	2.9
36	DC	191	THR	2.9
3	AC	37	LYS	2.9
8	CH	169	VAL	2.9
34	DA	1275	A	2.9
9	AK	105	PRO	2.9
46	DM	25	ILE	2.9
46	DM	88	ARG	2.9
26	C2	1	MET	2.9
35	DB	136	VAL	2.9
34	BA	1024	G	2.9
34	DA	1202	G	2.9
46	DM	118	ALA	2.9
1	CA	890	A	2.9
34	DA	1136	U	2.9
46	DM	40	ASN	2.9
7	CG	49	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	AA	931	C	2.9
10	AL	43	ALA	2.9
35	DB	227	GLY	2.9
40	DG	39	ALA	2.9
43	BJ	27	ALA	2.9
34	DA	1190	G	2.9
36	BC	23	TYR	2.9
54	DU	7	ARG	2.9
10	AL	112	MET	2.9
34	DA	1028	C	2.9
42	DI	14	VAL	2.9
36	BC	101	LEU	2.9
1	AA	271	U	2.9
34	DA	1212	U	2.9
57	DZ	683	VAL	2.9
36	BC	197	GLY	2.9
23	CZ	69	THR	2.9
42	DI	64	THR	2.9
34	BA	204	U	2.9
43	BJ	73	ASP	2.9
57	DZ	180	VAL	2.9
16	CS	58	LEU	2.9
49	DP	48	TRP	2.9
3	CC	16	ASP	2.9
7	CG	39	ILE	2.9
37	BD	20	TYR	2.9
42	DI	92	TYR	2.9
57	BZ	576	ASP	2.9
34	BA	1287	A	2.9
35	BB	227	GLY	2.9
1	AA	2209	G	2.9
36	DC	49	SER	2.8
34	DA	91	C	2.8
52	DS	33	THR	2.8
42	BI	88	TYR	2.8
3	CC	209	PHE	2.8
7	AG	80	PHE	2.8
8	CH	15	VAL	2.8
10	CL	67	PHE	2.8
36	BC	128	PHE	2.8
52	DS	51	VAL	2.8
1	CA	1083	U	2.8

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Mol	Chain	Res	Type	RSRZ
9	AK	115	GLN	2.8
35	BB	130	ARG	2.8
35	BB	222	ILE	2.8
46	BM	4	ILE	2.8
52	BS	80	TYR	2.8
4	CD	167	GLY	2.8
40	BG	130	GLY	2.8
36	DC	86	VAL	2.8
42	BI	59	PHE	2.8
7	CG	81	LYS	2.8
1	CA	2150	U	2.8
28	A4	52	THR	2.8
34	DA	1008	C	2.8
46	BM	55	ARG	2.8
52	DS	39	THR	2.8
57	DZ	563	ILE	2.8
38	DE	10	MET	2.8
1	AA	2803	A	2.8
34	DA	1248	A	2.8
1	CA	1044	G	2.8
40	BG	12	LEU	2.8
23	CZ	21	ALA	2.8
36	DC	160	ALA	2.8
5	CE	58	ARG	2.8
27	C3	3	ARG	2.8
8	CH	129	THR	2.8
46	BM	92	HIS	2.8
52	BS	10	PHE	2.8
46	BM	87	TYR	2.8
54	DU	14	TRP	2.8
57	DZ	390	VAL	2.8
10	AL	42	ASN	2.8
42	BI	46	ALA	2.8
52	BS	27	GLU	2.8
8	CH	25	LYS	2.8
34	BA	1040	U	2.8
28	C4	19	GLY	2.8
40	DG	130	GLY	2.8
42	BI	21	PRO	2.8
34	DA	1236	A	2.8
34	DA	1321	C	2.8
46	BM	74	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
47	DN	13	THR	2.8
1	CA	652(T)	C	2.8
10	AL	93	ARG	2.8
57	DZ	511	LYS	2.8
11	CN	71	ILE	2.8
35	BB	226	ARG	2.8
46	DM	16	ASP	2.8
40	DG	38	LEU	2.8
34	BA	1045	C	2.8
34	DA	1288	A	2.8
40	DG	56	GLN	2.8
34	DA	1274	G	2.8
34	DA	1316	G	2.8
54	BU	19	GLY	2.8
9	CK	69	PRO	2.8
36	DC	8	ILE	2.8
10	AL	125	ARG	2.8
40	DG	32	ARG	2.8
54	DU	15	ARG	2.8
1	CA	2804	C	2.8
42	DI	4	TYR	2.8
9	CK	48	GLY	2.7
42	DI	72	GLY	2.7
35	DB	144	ARG	2.7
42	BI	95	LYS	2.7
57	DZ	518	PRO	2.7
23	CZ	140	ASP	2.7
23	CZ	128	VAL	2.7
42	DI	58	HIS	2.7
47	DN	18	VAL	2.7
10	CL	118	THR	2.7
24	A0	2	ALA	2.7
52	BS	53	ASN	2.7
52	DS	44	MET	2.7
9	CK	4	LYS	2.7
28	A4	58	ARG	2.7
43	DJ	46	ARG	2.7
52	DS	37	ARG	2.7
57	DZ	226	ASN	2.7
9	CK	122	VAL	2.7
31	A7	48	LYS	2.7
57	DZ	188	TYR	2.7

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Mol	Chain	Res	Type	RSRZ
10	CL	129	GLY	2.7
42	DI	53	VAL	2.7
51	DR	58	LEU	2.7
8	CH	51	ARG	2.7
42	DI	120	ARG	2.7
10	AL	97	GLY	2.7
19	CV	63	GLY	2.7
36	BC	191	THR	2.7
42	DI	22	GLY	2.7
57	DZ	233	GLU	2.7
9	AK	25	PHE	2.7
1	AA	218	A	2.7
10	AL	68	VAL	2.7
10	CL	133	SER	2.7
40	DG	125	MET	2.7
34	DA	947	G	2.7
47	DN	16	PHE	2.7
52	DS	65	ASN	2.7
56	DW	40	C	2.7
43	BJ	85	LEU	2.7
51	BR	22	VAL	2.7
54	BU	5	ASP	2.7
57	BZ	575	VAL	2.7
10	CL	102	GLU	2.7
28	C4	30	GLU	2.7
52	DS	8	GLY	2.7
57	DZ	541	ALA	2.7
1	CA	1078	U	2.7
7	CG	182	LYS	2.7
34	BA	1020	U	2.7
1	AA	936	C	2.7
34	DA	1326	C	2.7
3	CC	221	PRO	2.7
8	CH	96	ALA	2.7
36	DC	22	TRP	2.7
36	DC	65	ALA	2.7
9	CK	78	SER	2.7
57	DZ	229	LEU	2.7
34	DA	1173	G	2.7
40	DG	55	GLY	2.7
42	DI	106	ALA	2.7
57	DZ	174	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
43	DJ	75	ILE	2.7
9	CK	124	ALA	2.7
42	BI	76	ALA	2.7
35	DB	113	HIS	2.7
7	CG	23	PHE	2.6
42	DI	33	PHE	2.6
34	DA	1261	A	2.6
34	DA	1349	A	2.6
40	BG	27	ILE	2.6
35	BB	231	GLU	2.6
57	BZ	536	LYS	2.6
43	DJ	87	THR	2.6
35	DB	137	ARG	2.6
36	DC	83	ARG	2.6
47	DN	35	ARG	2.6
1	AA	938	G	2.6
34	DA	1011	G	2.6
34	DA	1057	G	2.6
1	AA	1878	A	2.6
7	CG	146	TYR	2.6
43	BJ	23	ILE	2.6
57	DZ	568	TYR	2.6
36	DC	162	GLN	2.6
42	BI	47	LEU	2.6
41	BH	128	GLY	2.6
1	CA	1080	C	2.6
34	BA	1140	C	2.6
40	DG	40	ALA	2.6
9	CK	123	GLU	2.6
36	DC	182	ILE	2.6
1	AA	2147	G	2.6
24	C0	7	LEU	2.6
34	BA	79	G	2.6
34	DA	1142	G	2.6
1	CA	901	A	2.6
46	BM	7	VAL	2.6
46	DM	110	ARG	2.6
28	C4	52	THR	2.6
10	CL	106	GLU	2.6
34	BA	76	C	2.6
56	BW	43	C	2.6
57	DZ	225	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
43	DJ	38	ILE	2.6
28	C4	48	ARG	2.6
36	BC	131	ARG	2.6
40	BG	140	ASP	2.6
57	DZ	542	VAL	2.6
1	AA	693	G	2.6
34	DA	1293	G	2.6
55	BV	15	A	2.6
10	AL	123	ALA	2.6
48	DO	15	PHE	2.6
10	AL	99	ILE	2.6
33	C9	18	ARG	2.6
3	AC	8	TYR	2.6
36	DC	194	GLY	2.6
40	BG	44	TYR	2.6
47	DN	60	SER	2.6
57	DZ	175	SER	2.6
57	DZ	401	SER	2.6
3	AC	221	PRO	2.6
40	DG	62	PHE	2.6
42	DI	20	ARG	2.6
4	AD	275	LYS	2.6
34	BA	1247	U	2.6
42	BI	67	GLY	2.6
35	BB	33	TYR	2.6
36	BC	82	GLU	2.6
23	CZ	4	ARG	2.6
27	C3	30	ARG	2.6
56	DW	71	G	2.6
56	BW	45	U	2.6
28	C4	32	TYR	2.6
40	DG	33	ASP	2.6
10	CL	113	PRO	2.6
43	DJ	69	ASN	2.6
34	DA	1182	G	2.6
10	AL	138	VAL	2.6
42	BI	26	VAL	2.6
16	CS	44	LYS	2.6
57	DZ	432	ALA	2.6
9	CK	73	GLY	2.5
42	BI	8	GLY	2.5
1	AA	2195	A	2.5

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Mol	Chain	Res	Type	RSRZ
8	CH	44	VAL	2.5
40	DG	35	LYS	2.5
9	CK	34	ALA	2.5
43	DJ	18	ALA	2.5
52	DS	59	PRO	2.5
9	CK	68	LEU	2.5
36	DC	35	GLU	2.5
4	CD	262	ARG	2.5
7	AG	51	ARG	2.5
35	BB	19	HIS	2.5
36	DC	132	ARG	2.5
57	DZ	299	VAL	2.5
7	CG	25	TYR	2.5
13	CP	140	ALA	2.5
47	BN	2	ALA	2.5
42	DI	24	GLY	2.5
46	DM	78	ILE	2.5
47	BN	12	ARG	2.5
8	CH	24	VAL	2.5
10	CL	71	THR	2.5
47	DN	22	THR	2.5
57	DZ	476	VAL	2.5
34	BA	1317	C	2.5
34	DA	1320	C	2.5
42	BI	43	ALA	2.5
46	DM	42	ALA	2.5
35	BB	78	GLN	2.5
36	DC	74	GLY	2.5
40	DG	112	PRO	2.5
46	DM	97	PRO	2.5
28	A4	61	ARG	2.5
43	DJ	9	ARG	2.5
36	DC	204	LEU	2.5
42	DI	96	LEU	2.5
9	CK	27	VAL	2.5
3	CC	45	HIS	2.5
22	CY	46	LYS	2.5
9	CK	95	GLN	2.5
34	DA	1249	C	2.5
47	DN	32	SER	2.5
34	DA	1313	U	2.5
27	C3	5	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
43	DJ	11	PHE	2.5
34	BA	1352	C	2.5
43	DJ	30	SER	2.5
3	AC	14	LYS	2.5
1	AA	2146	G	2.5
34	DA	1273	G	2.5
43	DJ	47	PHE	2.5
57	DZ	422	GLU	2.5
42	BI	106	ALA	2.5
42	BI	62	TYR	2.5
42	DI	114	TYR	2.5
34	DA	1038	C	2.5
34	DA	1165	C	2.5
28	C4	8	LYS	2.5
57	BZ	496	LYS	2.5
8	CH	40	GLU	2.5
57	DZ	514	VAL	2.5
36	DC	62	ASP	2.5
46	DM	63	THR	2.5
57	DZ	684	GLN	2.5
38	BE	9	LYS	2.5
7	CG	2	PRO	2.5
8	CH	18	GLU	2.5
46	DM	17	VAL	2.5
1	AA	2148	A	2.5
10	CL	82	ALA	2.5
37	DD	2	GLY	2.5
47	BN	20	ALA	2.5
8	CH	94	TYR	2.5
34	BA	630	G	2.5
34	DA	1175	G	2.5
35	DB	121	LEU	2.5
36	BC	34	LEU	2.5
9	CK	8	GLU	2.4
46	BM	97	PRO	2.4
43	DJ	76	ASN	2.4
23	CZ	143	GLY	2.4
34	DA	1245	A	2.4
36	DC	28	GLN	2.4
36	DC	145	GLY	2.4
47	BN	13	THR	2.4
10	CL	79	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
36	BC	156	ARG	2.4
54	BU	9	ARG	2.4
28	C4	11	PRO	2.4
1	CA	2131	G	2.4
1	CA	2162	G	2.4
36	DC	154	SER	2.4
3	AC	15	VAL	2.4
34	BA	1367	C	2.4
35	BB	62	ALA	2.4
1	AA	2614	A	2.4
31	C7	47	ARG	2.4
57	BZ	504	ARG	2.4
37	BD	37	PRO	2.4
8	CH	63	SER	2.4
10	AL	41	PHE	2.4
33	C9	6	SER	2.4
35	BB	229	VAL	2.4
34	DA	1134	G	2.4
42	DI	57	GLY	2.4
53	DT	47	GLY	2.4
57	BZ	543	GLN	2.4
10	AL	32	ALA	2.4
42	DI	15	ALA	2.4
7	CG	90	LEU	2.4
33	C9	17	ILE	2.4
34	BA	1288	A	2.4
43	BJ	71	LEU	2.4
42	BI	5	TYR	2.4
52	DS	58	VAL	2.4
34	BA	1211	U	2.4
34	DA	1000	U	2.4
42	DI	71	SER	2.4
14	CQ	6	ARG	2.4
34	BA	1282	C	2.4
55	DV	18	C	2.4
10	AL	128	ALA	2.4
21	CX	92	LEU	2.4
40	BG	16	LEU	2.4
43	BJ	96	ILE	2.4
1	AA	700	A	2.4
42	DI	90	PRO	2.4
1	CA	2172	U	2.4

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Mol	Chain	Res	Type	RSRZ
7	CG	116	ASP	2.4
9	AK	89	ALA	2.4
9	CK	126	ALA	2.4
1	CA	652(C)	G	2.4
1	CA	879	G	2.4
47	DN	49	HIS	2.4
56	DW	53	G	2.4
8	CH	54	ARG	2.4
57	DZ	533	VAL	2.4
3	AC	16	ASP	2.4
44	DK	77	MET	2.4
35	DB	37	ASN	2.4
57	BZ	7	ASN	2.4
47	BN	39	LEU	2.4
34	DA	1272	G	2.4
3	AC	17	PRO	2.4
8	CH	43	VAL	2.4
9	CK	76	GLY	2.4
10	CL	36	GLU	2.4
36	DC	171	GLY	2.4
40	DG	54	THR	2.4
9	CK	28	ASN	2.4
10	AL	133	SER	2.4
52	BS	38	SER	2.4
1	AA	2129	C	2.4
9	CK	19	ARG	2.4
42	DI	78	LYS	2.4
36	BC	86	VAL	2.4
37	BD	163	GLU	2.4
28	A4	63	TYR	2.4
43	BJ	37	PRO	2.4
34	BA	73	G	2.4
34	DA	1310	G	2.4
54	DU	8	THR	2.4
34	BA	1318	A	2.4
34	DA	1004	A	2.4
3	CC	14	LYS	2.4
35	BB	133	LYS	2.4
57	DZ	99	ARG	2.4
9	CK	84	GLU	2.3
57	DZ	189	GLY	2.3
40	DG	11	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
1	AA	2206	G	2.3
36	BC	63	ASN	2.3
42	DI	89	ASN	2.3
57	BZ	563	ILE	2.3
8	CH	47	GLU	2.3
1	CA	2107	C	2.3
42	BI	33	PHE	2.3
57	BZ	453	GLY	2.3
35	DB	110	GLN	2.3
47	DN	34	TYR	2.3
54	BU	20	LYS	2.3
8	CH	23	ARG	2.3
57	BZ	419	ALA	2.3
34	BA	1385	G	2.3
34	DA	157	G	2.3
34	DA	1331	G	2.3
46	BM	32	GLU	2.3
40	BG	26	PHE	2.3
51	BR	29	PHE	2.3
7	CG	160	VAL	2.3
34	DA	1149	C	2.3
35	BB	165	VAL	2.3
41	DH	129	VAL	2.3
52	BS	25	LYS	2.3
23	CZ	93	ASP	2.3
54	BU	23	PRO	2.3
40	BG	154	TYR	2.3
46	DM	28	ALA	2.3
7	CG	77	ILE	2.3
9	CK	23	SER	2.3
36	DC	181	ASN	2.3
1	CA	2184	G	2.3
1	CA	2793	G	2.3
34	BA	1009	G	2.3
34	BA	1276	G	2.3
57	DZ	596	LYS	2.3
36	BC	64	VAL	2.3
1	CA	652(V)	C	2.3
1	CA	2161	C	2.3
34	BA	1314	C	2.3
40	DG	41	ARG	2.3
40	DG	44	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
39	DF	92	LYS	2.3
13	CP	93	GLY	2.3
23	CZ	66	SER	2.3
9	CK	21	GLN	2.3
42	BI	65	VAL	2.3
1	CA	882	G	2.3
1	CA	1170	G	2.3
28	C4	18	CYS	2.3
34	BA	1271	G	2.3
34	DA	1300	G	2.3
42	DI	54	ASP	2.3
40	BG	151	TYR	2.3
42	BI	92	TYR	2.3
43	DJ	40	LEU	2.3
42	BI	104	ARG	2.3
10	AL	29	GLN	2.3
46	BM	64	TRP	2.3
35	BB	117	GLU	2.3
34	BA	91	C	2.3
34	BA	1261	A	2.3
34	BA	1362	C	2.3
34	DA	1141	C	2.3
34	DA	1367	C	2.3
35	DB	218	ALA	2.3
40	BG	124	LEU	2.3
40	DG	73	MET	2.3
10	AL	79	ARG	2.3
35	DB	32	ILE	2.3
43	BJ	43	ARG	2.3
46	DM	39	ILE	2.3
57	DZ	582	PHE	2.3
8	CH	99	VAL	2.3
36	DC	75	VAL	2.3
43	BJ	19	SER	2.3
57	BZ	204	GLU	2.3
40	BG	153	HIS	2.3
43	DJ	62	HIS	2.3
23	CZ	62	PRO	2.3
8	CH	98	LEU	2.3
28	A4	67	TYR	2.3
34	BA	1363	C	2.3
34	DA	1168	A	2.3

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Mol	Chain	Res	Type	RSRZ
36	BC	196	LEU	2.3
39	DF	35	ALA	2.3
42	BI	77	ILE	2.3
46	BM	89	GLY	2.3
26	A2	12	GLU	2.3
28	C4	23	GLU	2.3
37	DD	161	ASN	2.3
42	DI	28	VAL	2.3
57	DZ	195	ASP	2.3
57	DZ	504	ARG	2.3
8	CH	33	LEU	2.3
40	BG	59	LEU	2.3
1	CA	1042	G	2.3
34	DA	1224	G	2.3
34	DA	1304	G	2.3
10	AL	55	VAL	2.2
57	BZ	594	VAL	2.2
33	C9	22	ARG	2.2
1	CA	1033	U	2.2
56	DW	45	U	2.2
3	AC	36	ALA	2.2
10	CL	108	ALA	2.2
51	BR	24	ALA	2.2
53	DT	40	ALA	2.2
57	DZ	516	PRO	2.2
57	DZ	513	LYS	2.2
1	AA	699	C	2.2
41	BH	58	TYR	2.2
42	BI	101	PHE	2.2
8	CH	58	GLU	2.2
46	BM	35	GLU	2.2
22	CY	45	VAL	2.2
36	BC	85	ARG	2.2
34	DA	1174	G	2.2
34	BA	723	U	2.2
34	DA	1056	U	2.2
27	C3	28	LEU	2.2
35	BB	61	LEU	2.2
36	BC	74	GLY	2.2
42	DI	85	LEU	2.2
43	DJ	91	PRO	2.2
42	DI	59	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
57	DZ	204	GLU	2.2
1	CA	1043	C	2.2
34	DA	1363	C	2.2
56	DW	3	C	2.2
57	BZ	491	VAL	2.2
34	BA	162	A	2.2
35	DB	133	LYS	2.2
40	BG	24	THR	2.2
10	CL	40	ALA	2.2
35	DB	19	HIS	2.2
27	C3	60	GLU	2.2
42	DI	52	ALA	2.2
28	A4	18	CYS	2.2
57	BZ	523	PHE	2.2
57	DZ	525	PHE	2.2
42	BI	3	GLN	2.2
42	DI	125	TYR	2.2
57	DZ	519	ARG	2.2
57	DZ	630	GLN	2.2
5	CE	33	VAL	2.2
11	CN	103	VAL	2.2
47	BN	11	LYS	2.2
52	DS	70	LYS	2.2
9	CK	16	ASN	2.2
10	AL	117	THR	2.2
14	CQ	30	GLY	2.2
34	BA	1360	A	2.2
10	CL	44	ALA	2.2
40	BG	99	LEU	2.2
43	BJ	39	PRO	2.2
2	CB	89	G	2.2
5	CE	61	ARG	2.2
10	CL	112	MET	2.2
34	DA	1178	G	2.2
46	DM	99	ARG	2.2
47	DN	57	ARG	2.2
57	DZ	550	MET	2.2
40	DG	151	TYR	2.2
1	CA	2188	C	2.2
34	DA	1116	C	2.2
43	BJ	94	VAL	2.2
46	DM	7	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
35	DB	11	LEU	2.2
42	BI	60	ASP	2.2
10	AL	122	ALA	2.2
36	DC	177	THR	2.2
42	DI	55	ALA	2.2
45	DL	94	PRO	2.2
47	BN	19	ARG	2.2
54	DU	9	ARG	2.2
9	CK	45	LYS	2.2
10	CL	95	LYS	2.2
9	AK	87	VAL	2.2
1	AA	692	C	2.2
35	DB	128	GLU	2.2
57	DZ	685	GLU	2.2
1	AA	2194	U	2.2
7	CG	43	LEU	2.2
23	CZ	131	ARG	2.2
57	DZ	619	ASP	2.2
52	BS	28	LYS	2.2
10	AL	91	PRO	2.2
36	BC	162	GLN	2.2
43	DJ	48	THR	2.2
49	DP	2	VAL	2.2
9	CK	41	ARG	2.2
34	BA	1181	G	2.2
42	BI	39	GLY	2.2
47	BN	29	ARG	2.2
52	DS	25	LYS	2.2
57	BZ	471	LYS	2.2
57	DZ	221	ALA	2.2
47	BN	14	PRO	2.2
10	AL	5	VAL	2.2
36	DC	103	VAL	2.2
57	DZ	-26	GLU	2.2
37	BD	3	ARG	2.2
42	DI	83	ARG	2.2
52	BS	3	ARG	2.2
3	AC	11	LEU	2.2
24	C0	71	ASP	2.2
34	BA	1311	G	2.2
34	BA	1320	C	2.2
34	DA	1161	C	2.2

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Mol	Chain	Res	Type	RSRZ
40	DG	12	LEU	2.2
46	DM	56	LEU	2.2
53	DT	9	ASN	2.2
56	DW	5	G	2.2
40	BG	50	ILE	2.2
22	CY	52	SER	2.2
57	BZ	578	SER	2.2
42	BI	27	THR	2.2
27	C3	59	VAL	2.2
36	DC	38	ARG	2.2
37	BD	132	ARG	2.2
47	DN	45	ARG	2.2
28	C4	4	GLY	2.1
34	DA	975	A	2.2
57	DZ	608	VAL	2.2
57	DZ	545	GLY	2.1
34	DA	1210	C	2.1
57	DZ	487	ILE	2.1
1	CA	2187	G	2.1
8	CH	60	ARG	2.1
8	CH	80	SER	2.1
35	BB	140	HIS	2.1
36	BC	21	ARG	2.1
42	DI	48	GLU	2.1
43	BJ	100	THR	2.1
8	CH	35	VAL	2.1
7	CG	131	TYR	2.1
36	BC	123	GLN	2.1
40	DG	131	LYS	2.1
43	DJ	45	ARG	2.1
9	CK	37	THR	2.1
40	BG	69	VAL	2.1
24	C0	78	TYR	2.1
9	CK	128	LEU	2.1
36	BC	204	LEU	2.1
43	DJ	63	PHE	2.1
43	DJ	89	ASP	2.1
46	BM	48	LEU	2.1
52	BS	5	LEU	2.1
34	BA	1005	A	2.1
34	BA	1126	U	2.1
52	DS	24	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	CA	898	C	2.1
56	BW	56	C	2.1
56	DW	72	C	2.1
46	BM	34	LEU	2.1
46	BM	96	LEU	2.1
36	BC	132	ARG	2.1
49	DP	9	PHE	2.1
26	A2	11	GLU	2.1
34	DA	1324	A	2.1
37	BD	26	CYS	2.1
47	DN	27	CYS	2.1
7	CG	155	MET	2.1
9	CK	103	GLY	2.1
34	BA	217	C	2.1
54	BU	14	TRP	2.1
57	DZ	490	PRO	2.1
57	DZ	638	GLY	2.1
10	AL	18	THR	2.1
3	CC	48	LEU	2.1
42	BI	79	LEU	2.1
42	BI	93	ARG	2.1
1	AA	2212	G	2.1
34	BA	1173	G	2.1
34	DA	1120	G	2.1
8	CH	89	ILE	2.1
34	DA	1148	U	2.1
22	CY	34	LYS	2.1
9	CK	22	GLY	2.1
34	DA	1357	A	2.1
35	BB	234	PRO	2.1
57	DZ	535	PRO	2.1
3	AC	215	VAL	2.1
34	BA	1007	C	2.1
34	DA	1246	C	2.1
37	BD	140	VAL	2.1
40	BG	141	VAL	2.1
42	DI	51	ARG	2.1
52	BS	69	HIS	2.1
57	DZ	339	SER	2.1
10	CL	103	GLN	2.1
36	BC	62	ASP	2.1
42	BI	73	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
34	DA	1235	U	2.1
1	CA	1058	G	2.1
1	CA	1533	G	2.1
34	BA	93	G	2.1
35	BB	147	LYS	2.1
34	BA	1046	A	2.1
34	DA	1170	A	2.1
40	BG	4	ARG	2.1
47	DN	23	ARG	2.1
34	DA	980	C	2.1
10	CL	119	ASP	2.1
34	DA	1205	U	2.1
35	BB	237	ALA	2.1
57	DZ	196	ILE	2.1
34	BA	77	G	2.1
34	DA	1365	G	2.1
57	DZ	494	GLU	2.1
1	CA	884	C	2.1
3	AC	33	LEU	2.1
11	CN	87	LEU	2.1
57	DZ	509	HIS	2.1
7	CG	157	ILE	2.1
49	BP	22	THR	2.1
52	DS	36	ARG	2.1
57	DZ	599	PRO	2.1
1	CA	1046	A	2.0
34	BA	1349	A	2.0
34	DA	1179	A	2.0
41	BH	52	ASP	2.0
43	DJ	28	ARG	2.0
7	CG	65	GLY	2.0
46	DM	50	GLU	2.0
27	C3	54	VAL	2.0
57	BZ	-48	VAL	2.0
46	DM	77	ASN	2.0
8	CH	103	LEU	2.0
57	DZ	497	PHE	2.0
7	AG	78	SER	2.0
23	CZ	80	ARG	2.0
34	DA	1317	C	2.0
43	BJ	17	ASP	2.0
9	CK	70	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
16	CS	22	GLY	2.0
34	DA	1447	A	2.0
35	DB	97	TRP	2.0
22	CY	62	GLU	2.0
35	BB	129	GLU	2.0
43	BJ	7	LYS	2.0
50	BQ	20	THR	2.0
10	CL	91	PRO	2.0
23	CZ	1	MET	2.0
13	CP	147	LEU	2.0
8	CH	69	ARG	2.0
21	AX	68	ARG	2.0
40	BG	51	GLN	2.0
40	DG	42	ILE	2.0
34	DA	1260	C	2.0
35	DB	116	GLU	2.0
36	BC	60	ALA	2.0
1	AA	691	G	2.0
16	CS	3	ARG	2.0
25	A1	98	LEU	2.0
9	AK	103	GLY	2.0
10	AL	85	GLU	2.0
35	BB	16	HIS	2.0
24	A0	6	GLY	2.0
54	DU	19	GLY	2.0
57	BZ	508	GLY	2.0
1	AA	1137	G	2.0
1	CA	652(E)	G	2.0
34	DA	1157	A	2.0
43	DJ	80	LYS	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
56	4SU	BY	8	20/21	0.26	0.50	-	300,300,300,300	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
56	5MU	DY	54	21/22	0.34	0.74	-	305,305,305,305	1
56	PSU	DW	39	20/21	0.87	0.35	-	93,93,93,93	3
56	7MG	DW	46	24/25	0.88	0.29	-	114,114,114,114	2
56	MIA	DW	37	29/30	0.91	0.27	-	94,94,94,94	0
56	5MU	BW	54	21/22	0.95	0.25	-	74,74,74,74	1
56	4SU	DY	8	20/21	0.17	0.51	-	275,275,275,275	0
56	PSU	BY	39	20/21	0.06	1.19	-	316,316,316,316	0
56	PSU	BW	55	20/21	0.88	0.24	-	74,74,74,74	5
56	PSU	BY	55	20/21	0.04	0.67	-	302,302,302,302	1
56	PSU	DY	32	20/21	0.18	1.61	-	268,268,268,268	0
56	PSU	DW	55	20/21	0.78	0.23	-	106,106,106,106	2
56	PSU	BY	32	20/21	0.34	0.89	-	254,254,254,254	1
56	5MU	BY	54	21/22	0.23	0.82	-	315,315,315,315	0
56	7MG	BW	46	24/25	0.94	0.20	-	63,63,63,63	5
56	PSU	DY	55	20/21	0.42	0.71	-	246,246,246,246	0
56	4SU	DW	8	20/21	0.93	0.20	-	88,88,88,88	3
56	PSU	DY	39	20/21	0.28	1.29	-	284,284,284,284	0
56	PSU	BW	39	20/21	0.95	0.25	-	65,65,65,65	3
56	PSU	BW	32	20/21	0.92	0.17	-	81,81,81,81	1
56	PSU	DW	32	20/21	0.88	0.21	-	106,106,106,106	1
56	5MU	DW	54	21/22	0.88	0.33	-	114,114,114,114	1
56	7MG	DY	46	24/25	0.52	0.53	-	302,302,302,302	0
56	7MG	BY	46	24/25	0.25	0.45	-	302,302,302,302	0
56	MIA	DY	37	22/30	0.12	1.85	-	319,319,319,319	1
56	MIA	BW	37	29/30	0.92	0.26	-	79,79,79,79	2
56	MIA	BY	37	22/30	0.05	1.39	-	284,284,284,284	0
56	4SU	BW	8	20/21	0.94	0.15	-	51,51,51,51	6

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
58	MG	AA	3018	1/1	0.83	1.37	126.81	78,78,78,78	0
58	MG	CA	3106	1/1	0.86	0.74	100.56	55,55,55,55	0
58	MG	CA	3043	1/1	0.59	0.83	64.26	102,102,102,102	0
58	MG	AA	3135	1/1	0.94	0.66	58.03	62,62,62,62	1
58	MG	AA	3171	1/1	0.96	0.49	53.19	25,25,25,25	1
58	MG	BA	1711	1/1	0.64	0.60	47.54	71,71,71,71	0
58	MG	CA	3185	1/1	0.91	0.48	46.01	59,59,59,59	0
58	MG	CA	3467	1/1	0.87	0.64	45.66	80,80,80,80	0
58	MG	CA	3314	1/1	0.74	0.53	41.67	77,77,77,77	0
58	MG	AA	3604	1/1	0.84	0.39	37.61	38,38,38,38	1
58	MG	BA	1801	1/1	0.89	0.46	34.36	69,69,69,69	0
58	MG	CA	3088	1/1	0.76	0.53	33.45	75,75,75,75	0
58	MG	CA	3619	1/1	0.89	0.44	33.17	47,47,47,47	1
58	MG	CA	3290	1/1	0.75	0.47	32.36	75,75,75,75	0
58	MG	CA	3114	1/1	0.93	0.45	32.24	39,39,39,39	0
58	MG	CA	3073	1/1	0.84	0.60	31.06	91,91,91,91	0
58	MG	BA	1657	1/1	0.86	0.41	30.60	73,73,73,73	0
58	MG	A7	103	1/1	0.82	0.50	30.26	38,38,38,38	1
58	MG	BA	1738	1/1	0.96	0.43	30.16	56,56,56,56	0
58	MG	CA	3168	1/1	0.86	0.52	29.98	56,56,56,56	0
58	MG	CA	3218	1/1	0.95	0.50	27.35	54,54,54,54	0
58	MG	CA	3223	1/1	0.83	0.58	27.21	65,65,65,65	0
58	MG	CA	3618	1/1	0.69	0.36	27.08	65,65,65,65	0
58	MG	CA	3498	1/1	0.94	0.49	26.75	68,68,68,68	0
58	MG	AH	3002	1/1	0.82	0.72	26.53	74,74,74,74	0
58	MG	CA	3542	1/1	0.94	0.39	26.28	68,68,68,68	0
58	MG	DA	1651	1/1	0.95	0.41	26.07	59,59,59,59	0
58	MG	CA	3230	1/1	0.93	0.42	26.04	51,51,51,51	0
58	MG	AA	3162	1/1	0.93	0.37	26.00	47,47,47,47	0
58	MG	DA	1743	1/1	0.92	0.35	25.94	72,72,72,72	0
58	MG	AB	3023	1/1	0.82	0.48	25.90	76,76,76,76	0
58	MG	AA	3211	1/1	0.92	0.57	25.62	42,42,42,42	1
58	MG	AA	3822	1/1	0.92	0.55	25.50	65,65,65,65	0
58	MG	C7	101	1/1	0.89	0.67	25.03	42,42,42,42	1
58	MG	DA	1668	1/1	0.97	0.42	24.86	62,62,62,62	0
58	MG	AA	3453	1/1	0.94	0.33	24.32	56,56,56,56	0
58	MG	BA	1664	1/1	0.88	0.39	24.13	59,59,59,59	0
58	MG	AA	3663	1/1	0.88	0.45	23.95	62,62,62,62	0
58	MG	AA	3132	1/1	0.88	0.28	23.76	27,27,27,27	1
58	MG	AA	3116	1/1	0.92	0.39	23.49	51,51,51,51	0
58	MG	AA	3768	1/1	0.72	0.40	23.47	99,99,99,99	0
58	MG	AA	3835	1/1	0.82	0.60	23.27	111,111,111,111	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
58	MG	AA	3212	1/1	0.94	0.41	23.24	34,34,34,34	1
58	MG	AA	3112	1/1	0.76	0.60	22.98	98,98,98,98	0
58	MG	CA	3441	1/1	0.91	0.35	22.41	77,77,77,77	0
58	MG	CA	3035	1/1	0.84	0.49	21.69	60,60,60,60	0
58	MG	AA	3462	1/1	0.94	0.45	21.41	71,71,71,71	0
58	MG	AA	3770	1/1	0.94	0.39	21.01	37,37,37,37	0
58	MG	BA	1629	1/1	0.91	0.49	20.83	64,64,64,64	0
58	MG	AA	3739	1/1	0.56	0.51	20.54	94,94,94,94	0
58	MG	AA	3117	1/1	0.85	0.32	20.44	30,30,30,30	1
58	MG	AA	3040	1/1	0.96	0.30	20.34	40,40,40,40	1
58	MG	CA	3588	1/1	0.95	0.35	20.31	63,63,63,63	0
58	MG	AA	3035	1/1	0.86	0.44	19.89	57,57,57,57	0
58	MG	AA	3196	1/1	0.94	0.39	19.69	52,52,52,52	0
58	MG	AA	3061	1/1	0.91	0.30	19.64	27,27,27,27	0
58	MG	DA	1684	1/1	0.85	0.54	19.53	72,72,72,72	0
58	MG	CA	3660	1/1	0.44	0.66	19.19	101,101,101,101	0
58	MG	AA	3101	1/1	0.90	0.36	19.02	52,52,52,52	0
58	MG	AA	3297	1/1	0.96	0.31	18.70	20,20,20,20	1
58	MG	AN	3001	1/1	0.31	0.89	18.66	85,85,85,85	0
58	MG	AA	3051	1/1	0.88	0.35	18.60	36,36,36,36	0
58	MG	CA	3166	1/1	0.96	0.38	18.35	44,44,44,44	0
58	MG	AA	3206	1/1	0.83	0.28	18.19	39,39,39,39	0
58	MG	AU	203	1/1	0.95	0.45	18.12	62,62,62,62	0
58	MG	AA	3185	1/1	0.89	0.29	18.11	41,41,41,41	0
58	MG	AA	3221	1/1	0.88	0.34	18.07	56,56,56,56	0
58	MG	AA	3133	1/1	0.79	0.37	17.72	69,69,69,69	0
58	MG	CA	3348	1/1	0.88	0.27	17.69	54,54,54,54	0
58	MG	AA	3134	1/1	0.97	0.39	17.48	59,59,59,59	1
58	MG	CA	3182	1/1	0.97	0.36	17.45	27,27,27,27	0
58	MG	BA	1756	1/1	0.78	0.37	17.44	68,68,68,68	0
58	MG	AA	3301	1/1	0.93	0.33	17.11	23,23,23,23	0
58	MG	DA	1672	1/1	0.61	0.47	16.84	73,73,73,73	0
58	MG	DA	1606	1/1	0.67	1.00	16.78	84,84,84,84	0
58	MG	AA	3184	1/1	0.81	0.34	16.74	68,68,68,68	0
58	MG	DA	1647	1/1	0.95	0.33	16.73	58,58,58,58	0
58	MG	CA	3322	1/1	0.99	0.28	16.68	45,45,45,45	0
58	MG	AW	3003	1/1	0.88	0.45	16.57	52,52,52,52	0
58	MG	CA	3500	1/1	0.64	0.43	16.55	64,64,64,64	0
58	MG	CF	301	1/1	0.67	0.40	16.46	63,63,63,63	0
58	MG	AA	3436	1/1	0.96	0.34	16.37	38,38,38,38	0
58	MG	AA	3606	1/1	0.89	0.33	16.09	61,61,61,61	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3028	1/1	0.97	0.55	16.00	35,35,35,35	1
58	MG	AA	3702	1/1	0.96	0.37	15.80	35,35,35,35	1
58	MG	AA	3708	1/1	0.90	0.52	15.56	53,53,53,53	1
58	MG	AA	3829	1/1	0.93	0.60	15.40	88,88,88,88	0
58	MG	AX	101	1/1	0.74	0.44	15.36	75,75,75,75	0
58	MG	AA	3168	1/1	0.80	0.38	15.05	63,63,63,63	0
58	MG	CA	3276	1/1	0.95	0.31	14.95	50,50,50,50	0
58	MG	CA	3038	1/1	0.90	0.37	14.89	48,48,48,48	0
58	MG	CA	3201	1/1	0.98	0.38	14.71	59,59,59,59	0
58	MG	AA	3525	1/1	0.95	0.33	14.67	40,40,40,40	0
58	MG	AA	3110	1/1	0.90	0.32	14.45	79,79,79,79	0
58	MG	CA	3226	1/1	0.84	0.35	14.36	69,69,69,69	0
58	MG	BA	1686	1/1	0.96	0.32	14.33	52,52,52,52	0
58	MG	AA	3267	1/1	0.91	0.42	14.05	63,63,63,63	0
58	MG	CA	3229	1/1	0.96	0.33	14.00	51,51,51,51	0
58	MG	CA	3597	1/1	0.94	0.28	13.92	39,39,39,39	0
58	MG	CU	201	1/1	0.89	0.56	13.82	64,64,64,64	0
58	MG	AA	3223	1/1	0.84	0.42	13.79	35,35,35,35	0
58	MG	AA	3823	1/1	0.97	0.24	13.46	37,37,37,37	1
58	MG	AA	3771	1/1	0.88	0.24	13.23	31,31,31,31	1
58	MG	CA	3212	1/1	0.83	0.33	13.18	69,69,69,69	0
58	MG	DA	1636	1/1	0.91	0.41	13.09	70,70,70,70	0
58	MG	AA	3819	1/1	0.92	0.26	13.05	25,25,25,25	1
58	MG	CA	3313	1/1	0.95	0.33	12.88	50,50,50,50	0
58	MG	AA	3173	1/1	0.96	0.24	12.82	60,60,60,60	0
58	MG	AA	3372	1/1	0.86	0.31	12.69	63,63,63,63	0
58	MG	CE	301	1/1	0.95	0.37	12.67	66,66,66,66	0
58	MG	CA	3428	1/1	0.93	0.29	12.49	54,54,54,54	1
58	MG	CA	3361	1/1	0.97	0.31	12.25	58,58,58,58	0
58	MG	AA	3138	1/1	0.94	0.37	12.22	50,50,50,50	0
58	MG	CA	3654	1/1	0.85	0.39	12.10	51,51,51,51	0
58	MG	CA	3661	1/1	0.91	0.26	12.01	74,74,74,74	0
58	MG	CA	3169	1/1	0.95	0.28	11.95	34,34,34,34	0
58	MG	AD	301	1/1	0.82	0.49	11.78	70,70,70,70	0
58	MG	AA	3331	1/1	0.95	0.28	11.75	34,34,34,34	0
58	MG	DA	1697	1/1	0.93	0.31	11.70	62,62,62,62	0
58	MG	AA	3400	1/1	0.98	0.36	11.68	39,39,39,39	0
58	MG	CA	3030	1/1	0.89	0.47	11.67	57,57,57,57	1
58	MG	AA	3316	1/1	0.97	0.29	11.65	60,60,60,60	0
58	MG	CA	3110	1/1	0.95	0.35	11.12	63,63,63,63	0
58	MG	AA	3179	1/1	0.97	0.31	11.04	71,71,71,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3084	1/1	0.92	0.34	11.01	59,59,59,59	1
58	MG	CA	3324	1/1	0.94	0.32	10.99	40,40,40,40	0
58	MG	CA	3014	1/1	0.89	0.44	10.96	62,62,62,62	0
58	MG	CD	303	1/1	0.97	0.52	10.87	37,37,37,37	0
58	MG	AA	3420	1/1	0.93	0.23	10.86	26,26,26,26	0
58	MG	DA	1618	1/1	0.91	0.45	10.86	65,65,65,65	0
58	MG	BA	1623	1/1	0.92	0.27	10.84	65,65,65,65	0
58	MG	CF	303	1/1	0.92	0.39	10.84	62,62,62,62	0
58	MG	AA	3282	1/1	0.90	0.49	10.72	40,40,40,40	0
58	MG	CA	3137	1/1	0.89	0.36	10.65	73,73,73,73	0
58	MG	AA	3150	1/1	0.96	0.31	10.51	15,15,15,15	0
58	MG	AA	3773	1/1	0.94	0.34	10.48	30,30,30,30	1
58	MG	AA	3620	1/1	0.95	0.32	10.45	40,40,40,40	0
58	MG	CA	3455	1/1	0.95	0.27	10.29	47,47,47,47	0
58	MG	BA	1616	1/1	0.67	0.61	10.19	134,134,134,134	0
58	MG	CA	3027	1/1	0.85	0.36	10.16	44,44,44,44	0
58	MG	DA	1638	1/1	0.91	0.33	10.13	83,83,83,83	0
58	MG	AA	3210	1/1	0.95	0.33	9.99	24,24,24,24	1
58	MG	AA	3698	1/1	0.94	0.28	9.93	32,32,32,32	1
58	MG	BA	1783	1/1	0.90	0.33	9.62	69,69,69,69	0
58	MG	AH	3001	1/1	0.93	0.30	9.59	52,52,52,52	0
58	MG	AA	3130	1/1	0.94	0.26	9.55	37,37,37,37	0
58	MG	AA	3602	1/1	0.94	0.27	9.54	37,37,37,37	0
58	MG	CA	3375	1/1	0.92	0.32	9.52	71,71,71,71	0
58	MG	AA	3039	1/1	0.95	0.32	9.51	39,39,39,39	1
58	MG	CA	3326	1/1	0.92	0.24	9.50	34,34,34,34	0
58	MG	AA	3418	1/1	0.92	0.24	9.47	43,43,43,43	0
58	MG	CA	3607	1/1	0.84	0.28	9.46	97,97,97,97	0
58	MG	AD	302	1/1	0.87	0.38	9.43	19,19,19,19	0
58	MG	AA	3559	1/1	0.96	0.22	9.30	39,39,39,39	0
58	MG	BA	1683	1/1	0.94	0.30	9.22	69,69,69,69	0
58	MG	BA	1721	1/1	0.91	0.23	9.15	60,60,60,60	0
58	MG	CA	3163	1/1	0.90	0.34	9.12	40,40,40,40	0
58	MG	AA	3174	1/1	0.89	0.31	9.02	63,63,63,63	0
58	MG	BA	1630	1/1	0.70	0.31	9.00	63,63,63,63	0
58	MG	CA	3489	1/1	0.89	0.25	8.92	80,80,80,80	0
58	MG	DA	1649	1/1	0.71	0.85	8.91	93,93,93,93	0
58	MG	AB	3008	1/1	0.94	0.47	8.80	51,51,51,51	0
58	MG	CA	3463	1/1	0.85	0.28	8.66	49,49,49,49	0
58	MG	AA	3805	1/1	0.94	0.33	8.64	35,35,35,35	1
58	MG	AA	3249	1/1	0.87	0.34	8.63	24,24,24,24	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3432	1/1	0.91	0.28	8.57	61,61,61,61	0
58	MG	CA	3119	1/1	0.98	0.31	8.51	128,128,128,128	0
58	MG	BA	1755	1/1	0.98	0.31	8.51	38,38,38,38	0
58	MG	CA	3213	1/1	0.90	0.27	8.46	44,44,44,44	0
58	MG	AA	3816	1/1	0.94	0.29	8.37	43,43,43,43	0
58	MG	AA	3081	1/1	0.85	0.29	8.35	40,40,40,40	0
58	MG	CA	3013	1/1	0.79	0.31	8.28	63,63,63,63	0
58	MG	AA	3617	1/1	0.89	0.23	8.25	49,49,49,49	0
58	MG	CA	3420	1/1	0.86	0.29	8.22	71,71,71,71	0
58	MG	AA	3311	1/1	0.89	0.21	8.03	33,33,33,33	0
58	MG	CA	3626	1/1	0.91	0.28	8.02	75,75,75,75	0
58	MG	AA	3354	1/1	0.83	0.31	8.01	60,60,60,60	0
58	MG	AD	309	1/1	0.74	0.30	7.94	57,57,57,57	0
58	MG	AA	3250	1/1	0.85	0.30	7.94	46,46,46,46	0
58	MG	AA	3190	1/1	0.92	0.25	7.93	45,45,45,45	0
58	MG	AA	3565	1/1	0.88	0.28	7.92	29,29,29,29	0
58	MG	DT	3001	1/1	0.87	0.53	7.92	67,67,67,67	0
58	MG	AA	3706	1/1	0.95	0.24	7.90	27,27,27,27	1
58	MG	AF	303	1/1	0.96	0.32	7.82	50,50,50,50	0
58	MG	AU	202	1/1	0.97	0.30	7.80	29,29,29,29	1
58	MG	AA	3817	1/1	0.90	0.32	7.79	61,61,61,61	0
58	MG	AA	3833	1/1	0.92	0.40	7.70	49,49,49,49	0
58	MG	AA	3037	1/1	0.96	0.31	7.69	45,45,45,45	0
58	MG	AD	304	1/1	0.91	0.31	7.64	38,38,38,38	1
58	MG	AA	3381	1/1	0.94	0.27	7.64	27,27,27,27	0
58	MG	DA	1694	1/1	0.91	0.29	7.62	60,60,60,60	0
58	MG	AA	3120	1/1	0.91	0.25	7.61	46,46,46,46	0
58	MG	AU	201	1/1	0.93	0.30	7.60	44,44,44,44	0
58	MG	AA	3824	1/1	0.89	0.27	7.56	45,45,45,45	0
58	MG	CA	3221	1/1	0.86	0.27	7.45	54,54,54,54	0
58	MG	DA	1768	1/1	0.94	0.43	7.31	73,73,73,73	0
58	MG	AA	3048	1/1	0.97	0.26	7.26	34,34,34,34	0
58	MG	CA	3530	1/1	0.95	0.26	7.07	59,59,59,59	0
58	MG	CA	3124	1/1	0.93	0.26	6.82	48,48,48,48	0
58	MG	CA	3396	1/1	0.75	0.24	6.74	58,58,58,58	0
58	MG	AA	3508	1/1	0.94	0.27	6.70	49,49,49,49	0
58	MG	AA	3564	1/1	0.98	0.22	6.66	19,19,19,19	0
58	MG	AD	308	1/1	0.93	0.39	6.62	42,42,42,42	0
58	MG	AA	3507	1/1	0.97	0.25	6.57	14,14,14,14	0
58	MG	AA	3253	1/1	0.94	0.27	6.56	29,29,29,29	1
58	MG	DA	1680	1/1	0.92	0.31	6.53	62,62,62,62	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3383	1/1	0.96	0.25	6.49	44,44,44,44	0
58	MG	AA	3113	1/1	0.91	0.32	6.48	64,64,64,64	0
58	MG	AA	3798	1/1	0.98	0.27	6.45	35,35,35,35	0
58	MG	CA	3217	1/1	0.80	0.30	6.38	52,52,52,52	0
58	MG	CA	3490	1/1	0.97	0.26	6.38	67,67,67,67	0
58	MG	AA	3240	1/1	0.95	0.34	6.37	30,30,30,30	0
58	MG	AA	3059	1/1	0.77	0.26	6.27	51,51,51,51	0
58	MG	CE	303	1/1	0.90	0.34	6.26	54,54,54,54	0
58	MG	CA	3650	1/1	0.95	0.27	6.23	27,27,27,27	0
58	MG	AA	3506	1/1	0.96	0.27	6.23	32,32,32,32	0
58	MG	CA	3190	1/1	0.91	0.37	6.20	83,83,83,83	0
58	MG	AA	3589	1/1	0.93	0.26	6.14	21,21,21,21	1
58	MG	CA	3091	1/1	0.92	0.35	6.04	111,111,111,111	0
58	MG	AA	3488	1/1	0.98	0.23	6.04	20,20,20,20	0
58	MG	CA	3330	1/1	0.94	0.24	5.99	43,43,43,43	0
58	MG	CA	3109	1/1	0.86	0.25	5.96	54,54,54,54	0
58	MG	AA	3443	1/1	0.93	0.21	5.94	65,65,65,65	0
58	MG	CA	3011	1/1	0.95	0.23	5.94	47,47,47,47	0
58	MG	CA	3127	1/1	0.97	0.25	5.90	63,63,63,63	0
58	MG	CA	3476	1/1	0.95	0.25	5.83	55,55,55,55	0
58	MG	AA	3791	1/1	0.99	0.26	5.73	16,16,16,16	0
58	MG	CA	3452	1/1	0.96	0.24	5.71	61,61,61,61	0
58	MG	AA	3596	1/1	0.87	0.23	5.69	40,40,40,40	0
58	MG	AA	3247	1/1	0.87	0.31	5.64	63,63,63,63	0
58	MG	BA	1648	1/1	0.85	0.22	5.62	37,37,37,37	0
58	MG	AV	202	1/1	0.89	0.30	5.60	55,55,55,55	1
58	MG	CA	3458	1/1	0.90	0.24	5.59	49,49,49,49	0
58	MG	A8	5001	1/1	0.88	0.35	5.58	59,59,59,59	0
58	MG	CA	3603	1/1	0.80	0.25	5.56	51,51,51,51	0
58	MG	DA	1669	1/1	0.89	0.40	5.49	84,84,84,84	0
58	MG	AD	310	1/1	0.93	0.31	5.47	58,58,58,58	0
58	MG	AA	3329	1/1	0.98	0.23	5.43	17,17,17,17	0
58	MG	BA	1811	1/1	0.81	0.32	5.35	75,75,75,75	0
58	MG	AA	3034	1/1	0.90	0.29	5.34	57,57,57,57	0
58	MG	AA	3023	1/1	0.98	0.31	5.33	33,33,33,33	1
58	MG	AA	3142	1/1	0.94	0.18	5.29	26,26,26,26	1
58	MG	AA	3717	1/1	0.96	0.24	5.27	47,47,47,47	0
58	MG	AA	3014	1/1	0.91	0.20	5.19	45,45,45,45	0
58	MG	AA	3187	1/1	0.93	0.26	5.12	32,32,32,32	0
58	MG	CQ	201	1/1	0.66	0.28	5.08	62,62,62,62	0
58	MG	AA	3551	1/1	0.93	0.23	5.00	52,52,52,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3036	1/1	0.92	0.19	4.95	51,51,51,51	0
58	MG	CA	3037	1/1	0.85	0.23	4.95	58,58,58,58	0
58	MG	DA	1722	1/1	0.86	0.26	4.92	65,65,65,65	0
58	MG	CA	3023	1/1	0.96	0.25	4.91	46,46,46,46	0
58	MG	AD	305	1/1	0.77	0.40	4.88	53,53,53,53	1
58	MG	AA	3439	1/1	0.97	0.23	4.80	17,17,17,17	0
58	MG	AA	3357	1/1	0.95	0.22	4.80	27,27,27,27	0
58	MG	AA	3820	1/1	0.94	0.29	4.79	40,40,40,40	0
58	MG	AA	3830	1/1	0.95	0.26	4.75	45,45,45,45	0
58	MG	AA	3395	1/1	0.96	0.21	4.74	18,18,18,18	0
58	MG	AA	3183	1/1	0.88	0.25	4.74	35,35,35,35	1
58	MG	CA	3353	1/1	0.97	0.23	4.71	48,48,48,48	0
58	MG	AA	3440	1/1	0.96	0.22	4.61	31,31,31,31	0
58	MG	AA	3806	1/1	0.81	0.25	4.61	61,61,61,61	0
58	MG	CA	3041	1/1	0.94	0.27	4.60	31,31,31,31	0
58	MG	CA	3263	1/1	0.93	0.25	4.58	57,57,57,57	0
58	MG	AA	3309	1/1	0.92	0.20	4.50	44,44,44,44	0
58	MG	CA	3086	1/1	0.94	0.24	4.50	36,36,36,36	0
58	MG	CA	3358	1/1	0.93	0.29	4.49	45,45,45,45	0
58	MG	AB	3003	1/1	0.88	0.24	4.46	51,51,51,51	0
58	MG	AA	3704	1/1	0.88	0.27	4.44	59,59,59,59	0
58	MG	BA	1671	1/1	0.81	0.23	4.41	75,75,75,75	0
58	MG	BA	1723	1/1	0.89	0.31	4.37	71,71,71,71	0
58	MG	BA	1678	1/1	0.87	0.26	4.35	69,69,69,69	0
58	MG	AA	3827	1/1	0.96	0.22	4.35	40,40,40,40	0
58	MG	DA	1610	1/1	0.86	0.22	4.35	75,75,75,75	0
58	MG	CA	3409	1/1	0.91	0.24	4.25	40,40,40,40	0
58	MG	CV	202	1/1	0.91	0.28	4.23	85,85,85,85	0
58	MG	CA	3364	1/1	0.97	0.24	4.21	29,29,29,29	0
58	MG	DA	1689	1/1	0.87	0.22	4.15	58,58,58,58	0
58	MG	BA	1626	1/1	0.88	0.30	4.13	75,75,75,75	0
58	MG	AA	3711	1/1	0.93	0.29	4.11	34,34,34,34	1
58	MG	CA	3555	1/1	0.74	0.26	4.11	83,83,83,83	0
58	MG	CA	3133	1/1	0.91	0.24	4.07	69,69,69,69	0
58	MG	AA	3082	1/1	0.97	0.20	3.98	23,23,23,23	1
58	MG	AA	3102	1/1	0.97	0.19	3.92	49,49,49,49	0
58	MG	AA	3811	1/1	0.93	0.29	3.85	58,58,58,58	0
58	MG	CA	3309	1/1	0.97	0.21	3.78	29,29,29,29	0
58	MG	AA	3231	1/1	0.85	0.22	3.74	53,53,53,53	0
58	MG	AA	3442	1/1	0.81	0.30	3.61	49,49,49,49	0
58	MG	AA	3812	1/1	0.91	0.22	3.59	42,42,42,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AF	302	1/1	0.91	0.29	3.56	41,41,41,41	0
58	MG	AA	3109	1/1	0.88	0.22	3.54	50,50,50,50	0
58	MG	AA	3621	1/1	0.79	0.18	3.46	46,46,46,46	0
58	MG	AA	3020	1/1	0.96	0.20	3.40	25,25,25,25	0
58	MG	AA	3388	1/1	0.99	0.23	3.37	25,25,25,25	0
58	MG	CA	3002	1/1	0.43	0.28	3.37	114,114,114,114	0
58	MG	AE	304	1/1	0.89	0.27	3.33	30,30,30,30	0
58	MG	CA	3596	1/1	0.72	0.23	3.29	72,72,72,72	0
58	MG	AA	3012	1/1	0.80	0.23	3.26	34,34,34,34	0
58	MG	AA	3456	1/1	0.91	0.18	3.25	30,30,30,30	0
58	MG	AA	3410	1/1	0.91	0.22	3.18	30,30,30,30	0
58	MG	BA	1695	1/1	0.70	0.21	3.15	98,98,98,98	0
58	MG	CA	3491	1/1	0.91	0.20	3.15	51,51,51,51	0
58	MG	BA	1693	1/1	0.70	0.28	3.14	67,67,67,67	0
58	MG	AA	3831	1/1	0.89	0.22	3.11	65,65,65,65	0
58	MG	AA	3828	1/1	0.92	0.28	3.08	37,37,37,37	1
58	MG	AA	3735	1/1	0.67	0.23	3.01	35,35,35,35	0
58	MG	DF	3001	1/1	0.79	0.22	2.97	49,49,49,49	0
58	MG	AA	3581	1/1	0.92	0.22	2.95	40,40,40,40	0
58	MG	BA	1690	1/1	0.81	0.30	2.90	89,89,89,89	0
58	MG	BA	1763	1/1	0.91	0.27	2.90	62,62,62,62	0
58	MG	BT	3001	1/1	0.95	0.37	2.86	62,62,62,62	0
58	MG	AA	3726	1/1	0.92	0.19	2.82	67,67,67,67	0
58	MG	CA	3178	1/1	0.93	0.22	2.82	57,57,57,57	0
58	MG	AA	3128	1/1	0.86	0.28	2.80	59,59,59,59	0
58	MG	AA	3044	1/1	0.94	0.20	2.76	34,34,34,34	0
58	MG	CA	3328	1/1	0.86	0.26	2.76	55,55,55,55	0
58	MG	DA	1658	1/1	0.92	0.21	2.69	72,72,72,72	0
58	MG	AQ	202	1/1	0.97	0.21	2.68	31,31,31,31	0
58	MG	BA	1615	1/1	0.78	0.31	2.62	74,74,74,74	0
58	MG	AA	3272	1/1	0.74	0.45	2.62	52,52,52,52	0
58	MG	CA	3251	1/1	0.94	0.19	2.57	56,56,56,56	0
58	MG	CA	3567	1/1	0.94	0.20	2.50	56,56,56,56	0
58	MG	CA	3159	1/1	0.83	0.41	2.50	69,69,69,69	0
58	MG	CA	3392	1/1	0.85	0.21	2.49	35,35,35,35	0
58	MG	AA	3257	1/1	0.88	0.19	2.47	14,14,14,14	0
58	MG	CA	3302	1/1	0.86	0.21	2.45	68,68,68,68	0
58	MG	AA	3485	1/1	0.92	0.21	2.45	14,14,14,14	0
58	MG	AA	3832	1/1	0.97	0.22	2.43	38,38,38,38	0
58	MG	CA	3281	1/1	0.96	0.26	2.42	51,51,51,51	0
58	MG	AA	3043	1/1	0.96	0.22	2.40	31,31,31,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3736	1/1	0.96	0.21	2.40	78,78,78,78	0
58	MG	AA	3815	1/1	0.98	0.17	2.31	29,29,29,29	1
58	MG	AA	3623	1/1	0.67	0.20	2.28	74,74,74,74	0
58	MG	AA	3314	1/1	0.97	0.20	2.25	28,28,28,28	0
58	MG	CA	3177	1/1	0.96	0.22	2.25	36,36,36,36	0
58	MG	C3	3001	1/1	0.95	0.33	2.13	69,69,69,69	0
58	MG	BA	1740	1/1	0.93	0.19	2.12	50,50,50,50	0
58	MG	CA	3266	1/1	0.96	0.21	2.10	69,69,69,69	0
58	MG	CA	3146	1/1	0.93	0.24	2.07	60,60,60,60	0
58	MG	CA	3340	1/1	0.88	0.22	2.07	48,48,48,48	0
58	MG	AA	3793	1/1	0.87	0.22	2.07	28,28,28,28	0
58	MG	AA	3499	1/1	0.95	0.18	2.07	51,51,51,51	1
58	MG	AA	3741	1/1	0.93	0.20	2.01	34,34,34,34	1
58	MG	AA	3519	1/1	0.87	0.19	2.00	21,21,21,21	0
58	MG	BA	1773	1/1	0.93	0.24	1.99	78,78,78,78	0
58	MG	CA	3214	1/1	0.95	0.19	1.95	40,40,40,40	0
58	MG	DA	1765	1/1	0.84	0.19	1.95	95,95,95,95	0
58	MG	CA	3362	1/1	0.93	0.19	1.92	44,44,44,44	0
58	MG	BA	1603	1/1	0.68	0.23	1.90	67,67,67,67	0
58	MG	AB	3014	1/1	0.94	0.18	1.86	67,67,67,67	0
58	MG	AA	3334	1/1	0.91	0.21	1.85	57,57,57,57	0
58	MG	AB	3016	1/1	0.96	0.17	1.81	34,34,34,34	0
58	MG	AA	3721	1/1	0.91	0.22	1.80	10,10,10,10	0
58	MG	AA	3518	1/1	0.87	0.22	1.79	33,33,33,33	0
58	MG	AA	3356	1/1	0.89	0.20	1.78	35,35,35,35	0
58	MG	CA	3318	1/1	0.99	0.22	1.74	33,33,33,33	0
58	MG	AA	3394	1/1	0.96	0.18	1.74	27,27,27,27	0
58	MG	BA	1752	1/1	0.97	0.21	1.73	48,48,48,48	0
58	MG	BF	3001	1/1	0.91	0.26	1.72	74,74,74,74	0
58	MG	AB	3017	1/1	0.69	0.18	1.72	77,77,77,77	0
58	MG	CV	201	1/1	0.79	0.22	1.70	100,100,100,100	0
58	MG	AA	3278	1/1	0.92	0.18	1.70	36,36,36,36	0
58	MG	AA	3543	1/1	0.83	0.20	1.66	52,52,52,52	1
58	MG	AP	201	1/1	0.96	0.21	1.62	28,28,28,28	1
58	MG	CW	201	1/1	0.94	0.26	1.58	39,39,39,39	0
58	MG	CA	3657	1/1	0.84	0.21	1.58	67,67,67,67	0
58	MG	AA	3714	1/1	0.94	0.20	1.52	56,56,56,56	0
58	MG	CA	3400	1/1	0.92	0.23	1.52	62,62,62,62	0
58	MG	CA	3521	1/1	0.93	0.20	1.48	61,61,61,61	0
58	MG	CA	3243	1/1	0.70	0.19	1.47	78,78,78,78	0
58	MG	CA	3433	1/1	0.95	0.16	1.47	71,71,71,71	0
58	MG	AA	3505	1/1	0.94	0.17	1.43	31,31,31,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	BA	1700	1/1	0.85	0.24	1.43	61,61,61,61	0
58	MG	CB	3007	1/1	0.76	0.19	1.42	65,65,65,65	0
58	MG	AA	3401	1/1	0.95	0.19	1.42	33,33,33,33	0
58	MG	DA	1739	1/1	0.87	0.20	1.40	79,79,79,79	0
61	FUA	BZ	703	37/37	0.85	0.27	1.36	69,69,69,69	0
58	MG	AA	3821	1/1	0.97	0.20	1.29	41,41,41,41	1
58	MG	AA	3290	1/1	0.92	0.18	1.28	63,63,63,63	0
58	MG	AD	307	1/1	0.87	0.16	1.25	37,37,37,37	0
58	MG	DA	1682	1/1	0.95	0.20	1.24	47,47,47,47	0
58	MG	AD	303	1/1	0.88	0.17	1.24	63,63,63,63	0
58	MG	CA	3662	1/1	0.87	0.23	1.22	55,55,55,55	0
58	MG	CA	3487	1/1	0.92	0.21	1.21	70,70,70,70	0
58	MG	AA	3433	1/1	0.97	0.21	1.20	28,28,28,28	0
58	MG	AA	3181	1/1	0.92	0.18	1.19	56,56,56,56	0
58	MG	AA	3459	1/1	0.98	0.19	1.14	18,18,18,18	0
58	MG	CG	3001	1/1	0.69	0.31	1.10	83,83,83,83	0
58	MG	AA	3050	1/1	0.95	0.19	1.08	53,53,53,53	0
58	MG	CA	3655	1/1	0.98	0.23	1.03	70,70,70,70	0
58	MG	AA	3045	1/1	0.98	0.19	1.03	43,43,43,43	0
58	MG	CA	3550	1/1	0.97	0.18	1.00	33,33,33,33	0
58	MG	CA	3332	1/1	0.92	0.22	0.99	42,42,42,42	0
61	FUA	DZ	703	37/37	0.87	0.23	0.98	85,85,85,85	0
58	MG	CA	3413	1/1	0.94	0.21	0.97	39,39,39,39	0
58	MG	AA	3723	1/1	0.98	0.20	0.95	21,21,21,21	0
58	MG	AA	3511	1/1	0.92	0.21	0.91	12,12,12,12	0
58	MG	AA	3213	1/1	0.84	0.17	0.87	58,58,58,58	0
58	MG	BA	1624	1/1	0.63	0.19	0.85	87,87,87,87	0
58	MG	BK	201	1/1	0.87	0.17	0.85	57,57,57,57	0
58	MG	DE	201	1/1	0.86	0.18	0.82	84,84,84,84	0
58	MG	AA	3473	1/1	0.98	0.21	0.80	15,15,15,15	0
58	MG	DA	1720	1/1	0.47	0.19	0.76	72,72,72,72	0
58	MG	BA	1704	1/1	0.87	0.23	0.74	61,61,61,61	0
58	MG	CA	3453	1/1	0.97	0.20	0.68	39,39,39,39	0
58	MG	AA	3169	1/1	0.94	0.20	0.67	39,39,39,39	0
58	MG	BA	1662	1/1	0.85	0.17	0.65	53,53,53,53	0
58	MG	CA	3457	1/1	0.96	0.21	0.62	47,47,47,47	0
58	MG	AA	3300	1/1	0.95	0.19	0.60	50,50,50,50	0
58	MG	AA	3145	1/1	0.95	0.17	0.59	33,33,33,33	0
58	MG	AD	306	1/1	0.91	0.16	0.59	73,73,73,73	0
58	MG	AA	3622	1/1	0.96	0.17	0.59	45,45,45,45	0
58	MG	DA	1637	1/1	0.89	0.25	0.57	76,76,76,76	0
58	MG	AA	3469	1/1	0.95	0.16	0.57	43,43,43,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	BA	1748	1/1	0.86	0.30	0.55	84,84,84,84	0
58	MG	CA	3267	1/1	0.88	0.17	0.53	56,56,56,56	0
58	MG	AA	3546	1/1	0.92	0.20	0.50	32,32,32,32	0
58	MG	AA	3404	1/1	0.95	0.19	0.50	19,19,19,19	0
58	MG	CA	3315	1/1	0.97	0.16	0.46	60,60,60,60	0
58	MG	AA	3389	1/1	0.99	0.19	0.46	17,17,17,17	0
58	MG	AA	3041	1/1	0.77	0.17	0.44	75,75,75,75	0
58	MG	AA	3235	1/1	0.57	0.16	0.39	64,64,64,64	0
58	MG	CF	302	1/1	0.95	0.19	0.31	56,56,56,56	0
58	MG	CA	3417	1/1	0.98	0.21	0.25	37,37,37,37	0
58	MG	AA	3007	1/1	0.91	0.16	0.25	21,21,21,21	0
58	MG	AA	3315	1/1	0.96	0.18	0.24	34,34,34,34	0
58	MG	AA	3535	1/1	0.95	0.19	0.20	15,15,15,15	0
58	MG	CF	305	1/1	0.77	0.16	0.18	51,51,51,51	0
58	MG	CA	3448	1/1	0.96	0.19	0.17	43,43,43,43	0
58	MG	CA	3564	1/1	0.99	0.19	0.14	40,40,40,40	1
58	MG	AA	3275	1/1	0.93	0.22	0.14	47,47,47,47	1
58	MG	BA	1800	1/1	0.65	0.21	0.13	84,84,84,84	0
62	GDP	DZ	704	28/28	0.93	0.17	0.13	80,80,80,80	3
58	MG	BA	1810	1/1	0.85	0.26	0.12	66,66,66,66	0
58	MG	BA	1685	1/1	0.90	0.18	0.12	50,50,50,50	0
58	MG	BA	1640	1/1	0.90	0.19	0.12	55,55,55,55	0
58	MG	CA	3579	1/1	0.88	0.18	0.12	51,51,51,51	0
58	MG	AA	3540	1/1	0.97	0.19	0.05	29,29,29,29	0
58	MG	AA	3377	1/1	0.99	0.19	0.05	20,20,20,20	0
58	MG	CA	3424	1/1	0.96	0.17	0.03	55,55,55,55	0
58	MG	CA	3003	1/1	0.92	0.20	0.02	45,45,45,45	0
58	MG	BA	1702	1/1	0.78	0.18	0.00	89,89,89,89	0
58	MG	AA	3686	1/1	0.93	0.18	-0.01	73,73,73,73	0
58	MG	CA	3273	1/1	0.87	0.17	-0.02	69,69,69,69	0
58	MG	AA	3341	1/1	0.92	0.21	-0.04	15,15,15,15	0
58	MG	AA	3217	1/1	0.96	0.17	-0.09	6,6,6,6	0
58	MG	AA	3687	1/1	0.93	0.18	-0.12	45,45,45,45	0
58	MG	AA	3258	1/1	0.82	0.18	-0.14	22,22,22,22	0
58	MG	DA	1715	1/1	0.90	0.24	-0.17	76,76,76,76	0
58	MG	CA	3656	1/1	0.92	0.23	-0.19	61,61,61,61	0
58	MG	AA	3378	1/1	0.95	0.17	-0.20	19,19,19,19	0
58	MG	AA	3517	1/1	0.97	0.19	-0.21	19,19,19,19	0
58	MG	CA	3311	1/1	0.92	0.15	-0.22	48,48,48,48	0
58	MG	AA	3649	1/1	0.91	0.14	-0.27	92,92,92,92	0
58	MG	AA	3383	1/1	0.90	0.19	-0.34	54,54,54,54	0
58	MG	CA	3047	1/1	0.88	0.16	-0.37	61,61,61,61	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CQ	203	1/1	0.86	0.15	-0.41	67,67,67,67	0
58	MG	CA	3615	1/1	0.96	0.18	-0.44	38,38,38,38	0
58	MG	AA	3583	1/1	0.99	0.16	-0.45	13,13,13,13	0
58	MG	AA	3202	1/1	0.74	0.15	-0.46	47,47,47,47	0
58	MG	CA	3020	1/1	0.76	0.17	-0.48	63,63,63,63	0
58	MG	AA	3524	1/1	0.94	0.17	-0.48	28,28,28,28	0
58	MG	DA	1709	1/1	0.89	0.16	-0.48	72,72,72,72	0
58	MG	CA	3439	1/1	0.91	0.20	-0.49	38,38,38,38	0
58	MG	AA	3515	1/1	0.95	0.17	-0.51	20,20,20,20	0
58	MG	CA	3069	1/1	0.94	0.18	-0.52	81,81,81,81	0
58	MG	CA	3192	1/1	0.92	0.14	-0.52	58,58,58,58	0
58	MG	AA	3734	1/1	0.98	0.17	-0.55	22,22,22,22	0
58	MG	AV	201	1/1	0.98	0.19	-0.59	38,38,38,38	0
58	MG	CB	3004	1/1	0.93	0.15	-0.60	68,68,68,68	0
62	GDP	BZ	704	28/28	0.97	0.15	-0.60	53,53,53,53	1
58	MG	CA	3337	1/1	0.93	0.17	-0.60	41,41,41,41	0
58	MG	AA	3792	1/1	0.92	0.17	-0.61	45,45,45,45	0
58	MG	CA	3664	1/1	0.92	0.15	-0.63	54,54,54,54	0
58	MG	AA	3542	1/1	0.95	0.15	-0.63	63,63,63,63	0
58	MG	CA	3160	1/1	0.91	0.13	-0.68	42,42,42,42	0
58	MG	CA	3057	1/1	0.93	0.17	-0.70	49,49,49,49	0
58	MG	AA	3794	1/1	0.91	0.17	-0.76	58,58,58,58	1
58	MG	DA	1630	1/1	0.81	0.16	-0.80	61,61,61,61	0
58	MG	AA	3393	1/1	0.98	0.18	-0.83	21,21,21,21	0
59	ZN	A5	101	1/1	1.00	0.12	-0.95	36,36,36,36	0
58	MG	BA	1679	1/1	0.96	0.15	-0.99	36,36,36,36	0
58	MG	AG	202	1/1	0.93	0.14	-1.02	73,73,73,73	0
58	MG	DA	1695	1/1	0.93	0.14	-1.04	66,66,66,66	0
58	MG	DA	1666	1/1	0.94	0.17	-1.05	66,66,66,66	0
58	MG	BA	1675	1/1	0.91	0.16	-1.05	57,57,57,57	0
58	MG	BA	1789	1/1	0.85	0.13	-1.06	68,68,68,68	0
58	MG	CA	3410	1/1	0.96	0.19	-1.12	31,31,31,31	0
58	MG	DA	1601	1/1	0.89	0.14	-1.12	61,61,61,61	0
58	MG	AA	3725	1/1	0.93	0.17	-1.13	39,39,39,39	0
59	ZN	C5	102	1/1	0.99	0.09	-1.17	68,68,68,68	0
58	MG	AA	3818	1/1	0.94	0.16	-1.26	19,19,19,19	0
58	MG	BM	201	1/1	0.87	0.12	-1.34	57,57,57,57	0
58	MG	CA	3105	1/1	0.61	0.12	-1.38	80,80,80,80	0
59	ZN	A4	501	1/1	0.96	0.06	-1.39	137,137,137,137	0
58	MG	AA	3009	1/1	0.94	0.14	-1.41	24,24,24,24	0
58	MG	DA	1721	1/1	0.92	0.13	-1.42	67,67,67,67	0
58	MG	AA	3571	1/1	0.89	0.15	-1.43	48,48,48,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
58	MG	AA	3826	1/1	0.96	0.17	-1.44	20,20,20,20	0
59	ZN	A9	501	1/1	1.00	0.10	-1.44	42,42,42,42	0
58	MG	CA	3346	1/1	0.95	0.17	-1.45	31,31,31,31	0
58	MG	CA	3427	1/1	0.86	0.14	-1.53	55,55,55,55	0
58	MG	AA	3200	1/1	0.73	0.11	-1.54	91,91,91,91	0
58	MG	CA	3227	1/1	0.96	0.15	-1.55	53,53,53,53	0
59	ZN	AY	501	1/1	0.99	0.08	-1.55	65,65,65,65	0
59	ZN	C6	501	1/1	0.99	0.10	-1.57	61,61,61,61	0
58	MG	AA	3072	1/1	0.94	0.14	-1.58	26,26,26,26	0
58	MG	AA	3099	1/1	0.87	0.13	-1.58	62,62,62,62	0
58	MG	AA	3529	1/1	0.97	0.16	-1.59	16,16,16,16	0
58	MG	AA	3619	1/1	0.96	0.15	-1.61	42,42,42,42	0
58	MG	CA	3261	1/1	0.96	0.17	-1.62	29,29,29,29	0
58	MG	DE	202	1/1	0.80	0.13	-1.66	94,94,94,94	0
58	MG	DA	1670	1/1	0.97	0.14	-1.68	75,75,75,75	0
60	SF4	DD	501	8/8	0.98	0.10	-1.70	90,90,90,90	1
58	MG	CA	3104	1/1	0.95	0.15	-1.72	48,48,48,48	0
58	MG	CA	3101	1/1	0.66	0.14	-1.73	78,78,78,78	0
58	MG	AA	3413	1/1	0.98	0.16	-1.78	20,20,20,20	0
58	MG	CA	3339	1/1	0.98	0.14	-1.78	34,34,34,34	0
58	MG	AA	3670	1/1	0.96	0.13	-1.79	33,33,33,33	0
60	SF4	BD	501	8/8	0.97	0.09	-1.79	80,80,80,80	0
59	ZN	C4	501	1/1	0.59	0.06	-1.85	189,189,189,189	0
58	MG	DA	1652	1/1	0.75	0.13	-1.86	71,71,71,71	0
58	MG	CA	3462	1/1	0.95	0.13	-1.89	49,49,49,49	0
58	MG	CA	3658	1/1	0.94	0.12	-1.90	50,50,50,50	0
58	MG	CA	3019	1/1	0.95	0.15	-1.92	28,28,28,28	0
58	MG	CA	3425	1/1	0.96	0.14	-1.93	53,53,53,53	0
58	MG	CA	3600	1/1	0.92	0.12	-1.93	50,50,50,50	0
58	MG	CA	3501	1/1	0.96	0.16	-1.94	45,45,45,45	1
59	ZN	A6	102	1/1	1.00	0.11	-1.96	46,46,46,46	0
58	MG	AA	3532	1/1	0.96	0.16	-2.00	25,25,25,25	0
58	MG	BA	1617	1/1	0.94	0.13	-2.01	73,73,73,73	0
58	MG	CA	3189	1/1	0.90	0.10	-2.07	68,68,68,68	0
58	MG	DA	1730	1/1	0.83	0.15	-2.09	88,88,88,88	0
58	MG	AA	3251	1/1	0.98	0.14	-2.11	42,42,42,42	0
58	MG	CA	3050	1/1	0.98	0.10	-2.12	44,44,44,44	0
59	ZN	C9	501	1/1	0.99	0.07	-2.14	94,94,94,94	0
58	MG	AA	3084	1/1	0.96	0.11	-2.15	32,32,32,32	0
58	MG	CA	3617	1/1	0.99	0.13	-2.18	31,31,31,31	0
58	MG	DA	1686	1/1	0.77	0.15	-2.19	53,53,53,53	0
58	MG	CA	3310	1/1	0.94	0.12	-2.19	48,48,48,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
59	ZN	DN	501	1/1	0.96	0.06	-2.29	129,129,129,129	0
58	MG	BA	1733	1/1	0.93	0.15	-2.35	62,62,62,62	0
58	MG	AA	3069	1/1	0.95	0.13	-2.38	34,34,34,34	0
58	MG	CA	3157	1/1	0.92	0.15	-2.38	55,55,55,55	0
58	MG	AA	3384	1/1	0.98	0.15	-2.39	34,34,34,34	0
58	MG	AA	3313	1/1	0.97	0.13	-2.40	35,35,35,35	0
59	ZN	BN	501	1/1	0.95	0.05	-2.50	123,123,123,123	0
58	MG	AA	3614	1/1	0.94	0.12	-2.50	56,56,56,56	0
58	MG	CA	3274	1/1	0.98	0.12	-2.53	50,50,50,50	1
58	MG	AA	3528	1/1	0.97	0.15	-2.55	28,28,28,28	0
58	MG	DA	1617	1/1	0.89	0.10	-2.56	48,48,48,48	0
58	MG	AA	3574	1/1	0.95	0.16	-2.58	30,30,30,30	1
58	MG	DA	1624	1/1	0.87	0.13	-2.61	44,44,44,44	0
59	ZN	CY	501	1/1	0.98	0.04	-2.66	93,93,93,93	0
58	MG	AA	3307	1/1	0.98	0.15	-2.67	6,6,6,6	0
58	MG	CA	3138	1/1	0.85	0.11	-2.70	63,63,63,63	0
58	MG	CA	3306	1/1	0.93	0.12	-2.72	41,41,41,41	0
58	MG	BA	1607	1/1	0.91	0.12	-2.75	62,62,62,62	0
58	MG	AA	3613	1/1	0.62	0.12	-2.76	104,104,104,104	0
58	MG	AA	3408	1/1	0.99	0.15	-2.78	20,20,20,20	0
58	MG	AA	3514	1/1	0.98	0.14	-2.82	18,18,18,18	0
58	MG	AA	3322	1/1	0.98	0.13	-2.84	32,32,32,32	1
58	MG	CE	302	1/1	0.97	0.13	-2.85	47,47,47,47	0
58	MG	CA	3372	1/1	0.99	0.13	-2.92	42,42,42,42	0
58	MG	CA	3320	1/1	0.98	0.15	-3.03	30,30,30,30	0
58	MG	AB	3020	1/1	0.93	0.11	-3.07	55,55,55,55	0
58	MG	CA	3010	1/1	0.84	0.08	-3.07	41,41,41,41	0
58	MG	AA	3526	1/1	0.98	0.16	-3.08	20,20,20,20	0
58	MG	CA	3360	1/1	0.94	0.12	-3.09	49,49,49,49	0
58	MG	AA	3344	1/1	0.96	0.10	-3.10	84,84,84,84	0
58	MG	AA	3396	1/1	0.94	0.14	-3.13	22,22,22,22	0
58	MG	CA	3419	1/1	0.92	0.11	-3.24	60,60,60,60	0
58	MG	AA	3398	1/1	0.99	0.14	-3.26	15,15,15,15	0
58	MG	AA	3022	1/1	0.92	0.16	-3.28	9,9,9,9	0
58	MG	BA	1680	1/1	0.98	0.11	-3.29	53,53,53,53	0
58	MG	AA	3547	1/1	0.88	0.10	-3.29	31,31,31,31	0
58	MG	CA	3211	1/1	0.94	0.11	-3.32	40,40,40,40	0
58	MG	BA	1620	1/1	0.96	0.09	-3.42	37,37,37,37	0
58	MG	DA	1687	1/1	0.67	0.13	-3.42	100,100,100,100	0
58	MG	CA	3370	1/1	0.94	0.13	-3.43	47,47,47,47	0
58	MG	CA	3464	1/1	0.94	0.14	-3.43	46,46,46,46	0
58	MG	BA	1654	1/1	0.95	0.09	-3.46	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	BB	3001	1/1	0.91	0.12	-3.49	75,75,75,75	0
58	MG	AG	201	1/1	0.95	0.10	-3.49	52,52,52,52	0
58	MG	AA	3750	1/1	0.91	0.13	-3.49	24,24,24,24	0
58	MG	AA	3486	1/1	0.95	0.16	-3.54	27,27,27,27	0
58	MG	AA	3572	1/1	0.93	0.14	-3.56	32,32,32,32	0
58	MG	CA	3278	1/1	0.87	0.15	-3.57	37,37,37,37	0
58	MG	CA	3488	1/1	0.91	0.10	-3.58	51,51,51,51	0
58	MG	CA	3265	1/1	0.96	0.11	-3.59	40,40,40,40	0
58	MG	CA	3560	1/1	0.98	0.13	-3.60	36,36,36,36	0
58	MG	AA	3435	1/1	0.96	0.16	-3.61	20,20,20,20	0
58	MG	AA	3324	1/1	0.96	0.12	-3.76	33,33,33,33	0
58	MG	AA	3642	1/1	0.95	0.14	-3.76	49,49,49,49	0
58	MG	AA	3387	1/1	0.99	0.14	-3.80	17,17,17,17	0
58	MG	AA	3038	1/1	0.99	0.13	-3.83	11,11,11,11	0
58	MG	CA	3584	1/1	0.95	0.13	-3.87	46,46,46,46	0
58	MG	AA	3775	1/1	0.89	0.11	-3.94	45,45,45,45	0
58	MG	BA	1674	1/1	0.94	0.05	-3.96	68,68,68,68	0
58	MG	AA	3556	1/1	0.94	0.15	-3.99	39,39,39,39	0
58	MG	AA	3054	1/1	0.94	0.14	-4.07	38,38,38,38	0
58	MG	DA	1645	1/1	0.98	0.12	-4.12	64,64,64,64	0
58	MG	CA	3282	1/1	0.98	0.12	-4.16	50,50,50,50	0
58	MG	CA	3343	1/1	0.91	0.13	-4.18	36,36,36,36	0
58	MG	AA	3825	1/1	0.94	0.13	-4.22	17,17,17,17	1
58	MG	BA	1741	1/1	0.96	0.09	-4.33	46,46,46,46	0
58	MG	AB	3007	1/1	0.93	0.09	-4.38	45,45,45,45	0
58	MG	AA	3777	1/1	0.98	0.10	-4.50	19,19,19,19	0
58	MG	AA	3299	1/1	0.98	0.13	-4.53	20,20,20,20	0
58	MG	BA	1611	1/1	0.97	0.13	-4.71	31,31,31,31	0
58	MG	DA	1654	1/1	0.97	0.10	-4.76	30,30,30,30	0
58	MG	AA	3385	1/1	0.98	0.13	-5.03	28,28,28,28	0
58	MG	AA	3403	1/1	0.96	0.09	-5.09	28,28,28,28	0
58	MG	AA	3539	1/1	0.96	0.14	-5.15	28,28,28,28	0
58	MG	BA	1612	1/1	0.91	0.10	-5.23	75,75,75,75	0
58	MG	CA	3526	1/1	0.94	0.10	-5.34	40,40,40,40	0
58	MG	CA	3321	1/1	0.97	0.12	-5.48	31,31,31,31	0
58	MG	CA	3317	1/1	0.92	0.12	-5.54	44,44,44,44	0
58	MG	CA	3438	1/1	0.92	0.15	-5.71	46,46,46,46	0
58	MG	AA	3434	1/1	0.93	0.12	-5.77	17,17,17,17	0
58	MG	AA	3616	1/1	0.94	0.10	-5.81	28,28,28,28	0
58	MG	AA	3374	1/1	0.98	0.08	-5.89	18,18,18,18	0
58	MG	AA	3003	1/1	0.98	0.09	-6.00	20,20,20,20	0
58	MG	AA	3502	1/1	0.97	0.12	-6.14	51,51,51,51	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
58	MG	AA	3021	1/1	0.92	0.11	-6.28	40,40,40,40	0
58	MG	CA	3647	1/1	0.91	0.07	-6.69	66,66,66,66	0
58	MG	AA	3449	1/1	0.97	0.11	-6.79	15,15,15,15	0
58	MG	AA	3754	1/1	0.98	0.08	-6.82	29,29,29,29	0
58	MG	CA	3029	1/1	0.94	0.11	-6.84	33,33,33,33	0
58	MG	AA	3342	1/1	0.99	0.14	-7.04	5,5,5,5	0
58	MG	AA	3684	1/1	0.97	0.10	-7.16	29,29,29,29	0
58	MG	AA	3669	1/1	0.97	0.06	-7.28	33,33,33,33	0
58	MG	CA	3242	1/1	0.91	0.12	-7.99	41,41,41,41	0
58	MG	BA	1643	1/1	0.94	0.07	-8.03	58,58,58,58	0
58	MG	BA	1728	1/1	0.97	0.11	-8.14	52,52,52,52	0
58	MG	AA	3562	1/1	0.97	0.06	-8.28	48,48,48,48	1
58	MG	BA	1613	1/1	0.90	0.08	-9.17	101,101,101,101	0
58	MG	CA	3628	1/1	0.96	0.13	-9.52	66,66,66,66	0
58	MG	CA	3291	1/1	0.97	0.12	-9.95	27,27,27,27	0
58	MG	AA	3493	1/1	0.95	0.13	-10.19	30,30,30,30	1
58	MG	AA	3498	1/1	0.99	0.12	-10.33	37,37,37,37	0
58	MG	AA	3011	1/1	0.97	0.09	-11.13	40,40,40,40	0
58	MG	BA	1758	1/1	0.98	0.06	-12.63	57,57,57,57	0
58	MG	CA	3021	1/1	0.96	0.12	-13.57	29,29,29,29	0
58	MG	DA	1621	1/1	0.94	0.12	-15.63	44,44,44,44	0
58	MG	BW	503	1/1	0.89	0.15	-	45,45,45,45	0
58	MG	CA	3423	1/1	0.88	0.19	-	50,50,50,50	0
58	MG	CA	3048	1/1	0.94	0.18	-	86,86,86,86	0
58	MG	BA	1794	1/1	0.96	0.17	-	77,77,77,77	0
58	MG	CA	3454	1/1	0.92	0.18	-	81,81,81,81	0
58	MG	CA	3074	1/1	0.92	0.50	-	53,53,53,53	0
58	MG	BA	1757	1/1	0.87	0.28	-	65,65,65,65	0
58	MG	AA	3452	1/1	0.95	0.14	-	69,69,69,69	0
58	MG	AQ	203	1/1	0.94	0.43	-	41,41,41,41	0
58	MG	AA	3656	1/1	0.81	0.22	-	56,56,56,56	0
58	MG	AY	502	1/1	0.91	0.31	-	58,58,58,58	0
58	MG	AA	3146	1/1	0.95	0.23	-	33,33,33,33	1
58	MG	CA	3545	1/1	0.91	0.13	-	68,68,68,68	0
58	MG	CA	3644	1/1	0.65	0.25	-	84,84,84,84	0
58	MG	BA	1644	1/1	0.81	0.33	-	75,75,75,75	0
58	MG	AA	3164	1/1	0.89	0.23	-	58,58,58,58	0
58	MG	DA	1700	1/1	0.40	0.28	-	124,124,124,124	0
58	MG	CA	3056	1/1	0.76	0.25	-	85,85,85,85	0
58	MG	AA	3699	1/1	0.92	0.37	-	46,46,46,46	1
58	MG	DA	1702	1/1	0.95	0.10	-	65,65,65,65	0
58	MG	AA	3013	1/1	0.95	0.18	-	35,35,35,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3121	1/1	0.89	0.15	-	45,45,45,45	0
58	MG	BA	1786	1/1	0.97	0.16	-	66,66,66,66	0
58	MG	CA	3236	1/1	0.78	0.45	-	87,87,87,87	0
58	MG	CA	3087	1/1	0.86	0.21	-	67,67,67,67	0
58	MG	CA	3188	1/1	0.91	0.56	-	58,58,58,58	0
58	MG	CA	3444	1/1	0.80	0.15	-	91,91,91,91	0
58	MG	CA	3365	1/1	0.97	0.20	-	48,48,48,48	0
58	MG	DA	1626	1/1	0.94	0.50	-	49,49,49,49	0
58	MG	AA	3071	1/1	0.95	0.75	-	41,41,41,41	0
58	MG	CA	3547	1/1	0.96	0.14	-	61,61,61,61	0
58	MG	AA	3359	1/1	0.79	0.19	-	51,51,51,51	0
58	MG	DA	1605	1/1	0.92	0.23	-	73,73,73,73	0
58	MG	CA	3286	1/1	0.87	0.21	-	64,64,64,64	0
58	MG	CA	3352	1/1	0.98	0.18	-	68,68,68,68	0
58	MG	AA	3005	1/1	0.92	0.21	-	62,62,62,62	0
58	MG	CA	3573	1/1	0.79	0.20	-	80,80,80,80	0
58	MG	CB	3009	1/1	0.91	0.16	-	67,67,67,67	0
58	MG	CA	3234	1/1	0.96	0.30	-	54,54,54,54	0
58	MG	DA	1690	1/1	0.89	0.53	-	82,82,82,82	0
58	MG	CA	3629	1/1	0.96	0.17	-	55,55,55,55	0
58	MG	AB	3021	1/1	0.72	0.21	-	61,61,61,61	0
58	MG	BA	1622	1/1	0.70	1.05	-	75,75,75,75	0
58	MG	AA	3479	1/1	0.88	0.26	-	54,54,54,54	0
58	MG	AA	3705	1/1	0.98	0.24	-	24,24,24,24	1
58	MG	AA	3087	1/1	0.93	0.25	-	72,72,72,72	0
58	MG	BA	1781	1/1	0.96	0.39	-	62,62,62,62	0
58	MG	AA	3073	1/1	0.94	0.27	-	58,58,58,58	0
58	MG	CA	3553	1/1	0.79	0.21	-	81,81,81,81	0
58	MG	AB	3009	1/1	0.94	0.08	-	56,56,56,56	0
58	MG	CA	3341	1/1	0.97	0.34	-	71,71,71,71	0
58	MG	CA	3275	1/1	0.95	0.23	-	42,42,42,42	0
58	MG	BV	101	1/1	0.82	0.34	-	110,110,110,110	0
58	MG	AA	3789	1/1	0.95	0.18	-	54,54,54,54	0
58	MG	AA	3137	1/1	0.82	0.29	-	56,56,56,56	0
58	MG	DA	1633	1/1	0.87	0.49	-	73,73,73,73	0
58	MG	CE	305	1/1	0.96	0.26	-	43,43,43,43	0
58	MG	AA	3321	1/1	0.95	0.32	-	70,70,70,70	0
58	MG	AA	3157	1/1	0.98	0.11	-	34,34,34,34	0
58	MG	CA	3514	1/1	0.72	0.80	-	105,105,105,105	0
58	MG	AA	3283	1/1	0.82	0.56	-	62,62,62,62	0
58	MG	CA	3152	1/1	0.68	0.28	-	56,56,56,56	0
58	MG	AZ	301	1/1	0.73	0.34	-	98,98,98,98	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	DA	1685	1/1	0.94	0.37	-	65,65,65,65	0
58	MG	CA	3203	1/1	0.85	0.12	-	60,60,60,60	0
58	MG	AA	3780	1/1	0.92	0.32	-	42,42,42,42	0
58	MG	AA	3254	1/1	0.96	0.19	-	42,42,42,42	0
58	MG	CA	3634	1/1	0.86	0.10	-	81,81,81,81	0
58	MG	AA	3659	1/1	0.89	0.24	-	14,14,14,14	0
58	MG	AA	3609	1/1	0.98	0.19	-	53,53,53,53	0
58	MG	CA	3225	1/1	0.85	0.47	-	65,65,65,65	0
58	MG	AA	3057	1/1	0.87	0.46	-	57,57,57,57	0
58	MG	CA	3331	1/1	0.94	0.36	-	52,52,52,52	0
58	MG	AA	3182	1/1	0.90	0.22	-	76,76,76,76	0
58	MG	CA	3518	1/1	0.88	0.32	-	86,86,86,86	0
58	MG	AA	3733	1/1	0.90	0.20	-	68,68,68,68	0
58	MG	CA	3430	1/1	0.89	0.33	-	53,53,53,53	0
58	MG	DA	1619	1/1	0.78	0.31	-	71,71,71,71	0
58	MG	AA	3167	1/1	0.96	0.14	-	65,65,65,65	0
58	MG	BA	1731	1/1	0.96	0.27	-	63,63,63,63	0
58	MG	BA	1676	1/1	0.96	0.25	-	44,44,44,44	0
58	MG	DA	1740	1/1	0.92	0.09	-	81,81,81,81	0
58	MG	CA	3296	1/1	0.97	0.32	-	42,42,42,42	0
58	MG	CP	203	1/1	0.82	0.25	-	67,67,67,67	0
58	MG	BA	1659	1/1	0.82	0.42	-	84,84,84,84	0
58	MG	AA	3339	1/1	0.87	0.29	-	43,43,43,43	0
58	MG	AA	3246	1/1	0.69	0.41	-	72,72,72,72	0
58	MG	CA	3297	1/1	0.98	0.34	-	56,56,56,56	0
58	MG	AA	3047	1/1	0.89	0.35	-	34,34,34,34	0
58	MG	CA	3044	1/1	0.91	0.35	-	52,52,52,52	0
58	MG	AA	3495	1/1	0.93	0.17	-	59,59,59,59	0
58	MG	BA	1809	1/1	0.91	0.22	-	68,68,68,68	0
58	MG	AA	3679	1/1	0.88	0.28	-	65,65,65,65	0
58	MG	AA	3423	1/1	0.98	0.15	-	22,22,22,22	0
58	MG	BA	1775	1/1	0.72	0.21	-	75,75,75,75	0
58	MG	CA	3621	1/1	0.88	0.31	-	73,73,73,73	0
58	MG	CA	3445	1/1	0.93	0.32	-	40,40,40,40	0
58	MG	AA	3326	1/1	0.97	0.12	-	58,58,58,58	0
58	MG	DA	1703	1/1	0.95	0.29	-	74,74,74,74	0
58	MG	BA	1701	1/1	0.96	0.46	-	58,58,58,58	0
58	MG	CA	3036	1/1	0.85	0.19	-	44,44,44,44	0
58	MG	DA	1677	1/1	0.87	0.13	-	74,74,74,74	0
58	MG	AA	3067	1/1	0.97	0.41	-	55,55,55,55	0
58	MG	BA	1697	1/1	0.90	0.38	-	78,78,78,78	0
58	MG	AA	3579	1/1	0.93	0.14	-	38,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
58	MG	AA	3799	1/1	0.90	0.19	-	47,47,47,47	0
58	MG	CA	3529	1/1	0.96	0.08	-	57,57,57,57	0
58	MG	AA	3497	1/1	0.93	0.04	-	46,46,46,46	0
58	MG	CA	3052	1/1	0.91	0.43	-	69,69,69,69	0
58	MG	CA	3536	1/1	0.87	0.12	-	71,71,71,71	0
58	MG	AA	3744	1/1	0.67	0.29	-	86,86,86,86	0
58	MG	CA	3195	1/1	0.91	0.15	-	60,60,60,60	0
58	MG	AA	3143	1/1	0.87	0.33	-	47,47,47,47	0
58	MG	AA	3237	1/1	0.86	0.14	-	71,71,71,71	0
58	MG	AA	3380	1/1	0.93	0.14	-	15,15,15,15	0
58	MG	AA	3266	1/1	0.93	0.50	-	50,50,50,50	0
58	MG	AA	3161	1/1	0.86	0.48	-	60,60,60,60	0
58	MG	CA	3630	1/1	0.91	0.12	-	63,63,63,63	0
58	MG	CA	3480	1/1	0.88	0.29	-	55,55,55,55	0
58	MG	CA	3541	1/1	0.90	0.17	-	71,71,71,71	0
58	MG	AA	3718	1/1	0.88	0.11	-	43,43,43,43	0
58	MG	AA	3063	1/1	0.80	0.45	-	67,67,67,67	0
58	MG	AA	3634	1/1	0.91	0.32	-	62,62,62,62	0
58	MG	AA	3630	1/1	0.87	0.23	-	71,71,71,71	0
58	MG	AA	3802	1/1	0.86	0.38	-	55,55,55,55	0
58	MG	A0	102	1/1	0.87	0.23	-	40,40,40,40	0
58	MG	AA	3118	1/1	0.90	0.62	-	64,64,64,64	0
58	MG	AF	301	1/1	0.93	0.17	-	43,43,43,43	0
58	MG	AA	3302	1/1	0.91	0.21	-	56,56,56,56	0
58	MG	BA	1706	1/1	0.94	0.43	-	62,62,62,62	0
58	MG	BA	1651	1/1	0.80	0.16	-	102,102,102,102	0
58	MG	AA	3064	1/1	0.90	0.17	-	29,29,29,29	0
58	MG	BA	1771	1/1	0.19	0.94	-	115,115,115,115	0
58	MG	CA	3175	1/1	0.99	0.28	-	42,42,42,42	0
58	MG	AA	3373	1/1	0.82	0.29	-	59,59,59,59	0
58	MG	CA	3299	1/1	0.98	0.16	-	61,61,61,61	0
58	MG	CA	3652	1/1	0.95	0.17	-	53,53,53,53	0
58	MG	AA	3160	1/1	0.94	0.17	-	57,57,57,57	0
58	MG	CA	3148	1/1	0.98	0.29	-	65,65,65,65	0
58	MG	CB	3002	1/1	0.72	0.27	-	66,66,66,66	0
58	MG	CA	3082	1/1	0.73	0.31	-	70,70,70,70	0
58	MG	AA	3763	1/1	0.96	0.36	-	65,65,65,65	0
58	MG	BA	1798	1/1	0.91	0.15	-	69,69,69,69	0
58	MG	BA	1673	1/1	0.78	0.96	-	77,77,77,77	0
58	MG	CA	3571	1/1	0.66	0.14	-	65,65,65,65	0
58	MG	AA	3238	1/1	0.93	0.35	-	61,61,61,61	0
58	MG	BA	1734	1/1	0.68	0.40	-	81,81,81,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3033	1/1	0.96	0.23	-	29,29,29,29	1
58	MG	CA	3215	1/1	0.86	0.09	-	73,73,73,73	0
58	MG	DA	1748	1/1	0.92	0.16	-	78,78,78,78	0
58	MG	AA	3585	1/1	0.69	0.17	-	65,65,65,65	0
58	MG	CA	3477	1/1	0.89	0.11	-	74,74,74,74	0
58	MG	CB	3010	1/1	0.93	0.23	-	55,55,55,55	0
58	MG	DA	1712	1/1	0.93	0.14	-	81,81,81,81	0
58	MG	AA	3446	1/1	0.91	0.51	-	61,61,61,61	0
58	MG	AA	3580	1/1	0.79	0.14	-	23,23,23,23	0
58	MG	AA	3428	1/1	0.94	0.17	-	35,35,35,35	0
58	MG	CA	3659	1/1	0.95	0.11	-	77,77,77,77	0
58	MG	AA	3318	1/1	0.96	0.25	-	51,51,51,51	1
58	MG	BA	1791	1/1	0.89	0.14	-	63,63,63,63	0
58	MG	AA	3364	1/1	0.88	0.34	-	81,81,81,81	0
58	MG	AA	3232	1/1	0.91	0.30	-	79,79,79,79	0
58	MG	CA	3640	1/1	0.86	0.29	-	59,59,59,59	0
58	MG	CA	3602	1/1	0.89	0.17	-	66,66,66,66	0
58	MG	AA	3567	1/1	0.97	0.21	-	51,51,51,51	0
58	MG	DA	1691	1/1	0.76	0.20	-	85,85,85,85	0
58	MG	AA	3633	1/1	0.88	0.25	-	48,48,48,48	1
58	MG	DA	1644	1/1	0.96	0.16	-	57,57,57,57	0
58	MG	CA	3220	1/1	0.88	0.24	-	78,78,78,78	0
58	MG	AA	3086	1/1	0.95	0.15	-	47,47,47,47	0
58	MG	CA	3531	1/1	0.91	0.21	-	51,51,51,51	0
58	MG	CA	3107	1/1	0.36	0.47	-	108,108,108,108	0
58	MG	AA	3431	1/1	0.98	0.10	-	25,25,25,25	0
58	MG	AA	3159	1/1	0.94	0.23	-	46,46,46,46	1
58	MG	AA	3628	1/1	0.89	0.28	-	80,80,80,80	0
58	MG	DA	1736	1/1	0.95	0.11	-	79,79,79,79	0
58	MG	CA	3509	1/1	0.83	0.18	-	76,76,76,76	0
58	MG	CA	3538	1/1	0.89	0.07	-	71,71,71,71	0
58	MG	AA	3749	1/1	0.98	0.14	-	14,14,14,14	0
58	MG	CA	3204	1/1	0.74	0.37	-	74,74,74,74	0
58	MG	CA	3565	1/1	0.98	0.39	-	51,51,51,51	0
58	MG	A0	103	1/1	0.93	0.12	-	37,37,37,37	0
58	MG	AA	3426	1/1	0.91	0.15	-	50,50,50,50	0
58	MG	CA	3401	1/1	0.93	0.23	-	69,69,69,69	0
58	MG	AA	3358	1/1	0.88	0.17	-	63,63,63,63	0
58	MG	AA	3214	1/1	0.93	0.16	-	34,34,34,34	0
58	MG	AA	3491	1/1	0.88	0.29	-	35,35,35,35	0
58	MG	AA	3809	1/1	0.89	0.22	-	62,62,62,62	0
58	MG	CA	3376	1/1	0.94	0.08	-	66,66,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	BA	1694	1/1	0.86	0.07	-	80,80,80,80	0
58	MG	AA	3095	1/1	0.77	0.47	-	82,82,82,82	0
58	MG	DA	1757	1/1	0.89	0.33	-	75,75,75,75	0
58	MG	AA	3340	1/1	0.98	0.10	-	59,59,59,59	0
58	MG	AA	3647	1/1	0.91	0.21	-	72,72,72,72	0
58	MG	DA	1657	1/1	0.40	0.20	-	93,93,93,93	0
58	MG	CA	3164	1/1	0.96	0.41	-	41,41,41,41	0
58	MG	AA	3558	1/1	0.83	0.14	-	51,51,51,51	0
58	MG	AA	3441	1/1	0.93	0.18	-	51,51,51,51	1
58	MG	CA	3139	1/1	0.84	0.33	-	123,123,123,123	0
58	MG	AA	3355	1/1	0.86	0.17	-	58,58,58,58	0
58	MG	CA	3510	1/1	0.92	0.14	-	65,65,65,65	0
58	MG	BA	1646	1/1	0.83	0.80	-	75,75,75,75	0
58	MG	AA	3075	1/1	0.97	0.15	-	9,9,9,9	0
58	MG	DA	1764	1/1	0.92	0.08	-	55,55,55,55	0
58	MG	BA	1747	1/1	0.94	0.40	-	68,68,68,68	0
58	MG	AA	3607	1/1	0.97	0.09	-	30,30,30,30	0
58	MG	AA	3371	1/1	0.96	0.34	-	62,62,62,62	0
58	MG	AA	3730	1/1	0.96	0.27	-	30,30,30,30	0
58	MG	CA	3125	1/1	0.88	0.50	-	73,73,73,73	0
58	MG	CA	3651	1/1	0.86	0.24	-	31,31,31,31	0
58	MG	CA	3134	1/1	0.55	1.10	-	86,86,86,86	0
58	MG	AA	3294	1/1	0.88	0.25	-	66,66,66,66	0
58	MG	AA	3225	1/1	0.92	0.26	-	27,27,27,27	1
58	MG	CA	3009	1/1	0.85	0.51	-	67,67,67,67	0
58	MG	AA	3320	1/1	0.95	0.19	-	24,24,24,24	0
58	MG	CA	3351	1/1	0.91	0.14	-	46,46,46,46	0
58	MG	CA	3594	1/1	0.37	0.21	-	80,80,80,80	0
58	MG	CA	3055	1/1	0.97	0.51	-	39,39,39,39	0
58	MG	DA	1653	1/1	0.90	0.19	-	60,60,60,60	0
58	MG	AA	3636	1/1	0.84	0.25	-	65,65,65,65	0
58	MG	CA	3627	1/1	0.93	0.18	-	63,63,63,63	0
58	MG	CA	3248	1/1	0.81	0.65	-	77,77,77,77	0
58	MG	CA	3636	1/1	0.82	0.19	-	80,80,80,80	0
58	MG	DA	1761	1/1	0.97	0.29	-	66,66,66,66	0
58	MG	AR	201	1/1	0.96	0.28	-	32,32,32,32	0
58	MG	DA	1719	1/1	0.92	0.12	-	74,74,74,74	0
58	MG	CA	3645	1/1	0.85	0.14	-	82,82,82,82	0
58	MG	AA	3476	1/1	0.82	0.24	-	69,69,69,69	0
58	MG	CA	3136	1/1	0.90	0.35	-	53,53,53,53	0
58	MG	AA	3070	1/1	0.89	0.40	-	60,60,60,60	0
58	MG	DA	1648	1/1	0.96	0.22	-	50,50,50,50	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3067	1/1	0.84	0.34	-	72,72,72,72	0
58	MG	AA	3289	1/1	0.94	0.11	-	27,27,27,27	0
58	MG	AA	3584	1/1	0.96	0.08	-	65,65,65,65	0
58	MG	AB	3012	1/1	0.99	0.22	-	29,29,29,29	1
58	MG	BA	1797	1/1	0.80	0.21	-	63,63,63,63	0
58	MG	CA	3354	1/1	0.97	0.25	-	61,61,61,61	0
58	MG	CA	3200	1/1	0.89	0.47	-	54,54,54,54	0
58	MG	AA	3480	1/1	0.91	0.08	-	54,54,54,54	0
58	MG	AA	3586	1/1	0.90	0.14	-	62,62,62,62	0
58	MG	CA	3623	1/1	0.90	0.26	-	64,64,64,64	0
58	MG	CA	3240	1/1	0.91	0.15	-	64,64,64,64	0
58	MG	AA	3566	1/1	0.96	0.18	-	27,27,27,27	0
58	MG	AA	3561	1/1	0.88	0.26	-	58,58,58,58	0
58	MG	AA	3415	1/1	0.96	0.07	-	56,56,56,56	0
58	MG	CA	3012	1/1	0.97	0.26	-	59,59,59,59	0
58	MG	CA	3172	1/1	0.79	0.33	-	83,83,83,83	0
58	MG	AA	3406	1/1	0.91	0.16	-	57,57,57,57	0
58	MG	AA	3695	1/1	0.93	0.51	-	78,78,78,78	0
58	MG	AA	3001	1/1	0.96	0.10	-	37,37,37,37	0
58	MG	DA	1750	1/1	0.86	0.17	-	68,68,68,68	0
58	MG	BA	1614	1/1	0.83	0.32	-	88,88,88,88	0
58	MG	AA	3697	1/1	0.80	0.15	-	63,63,63,63	0
58	MG	CA	3199	1/1	0.90	0.35	-	74,74,74,74	0
58	MG	AA	3330	1/1	0.97	0.13	-	69,69,69,69	0
58	MG	AA	3748	1/1	0.96	0.28	-	56,56,56,56	0
58	MG	AA	3738	1/1	0.80	0.26	-	75,75,75,75	0
58	MG	AA	3555	1/1	0.97	0.17	-	38,38,38,38	0
58	MG	AA	3197	1/1	0.96	0.38	-	37,37,37,37	0
58	MG	DA	1604	1/1	0.81	0.37	-	76,76,76,76	0
58	MG	AA	3188	1/1	0.98	0.18	-	13,13,13,13	0
58	MG	AA	3530	1/1	0.98	0.12	-	53,53,53,53	0
58	MG	BA	1805	1/1	0.64	0.18	-	83,83,83,83	0
58	MG	CA	3610	1/1	0.81	0.16	-	98,98,98,98	0
58	MG	AA	3761	1/1	0.88	0.46	-	50,50,50,50	0
58	MG	AA	3764	1/1	0.97	0.20	-	55,55,55,55	0
58	MG	DA	1763	1/1	0.85	0.20	-	77,77,77,77	0
58	MG	AA	3227	1/1	0.92	0.26	-	55,55,55,55	0
58	MG	AA	3277	1/1	0.80	0.23	-	79,79,79,79	0
58	MG	CA	3080	1/1	0.61	0.22	-	87,87,87,87	0
58	MG	CA	3103	1/1	0.85	0.49	-	62,62,62,62	0
58	MG	CA	3582	1/1	0.93	0.19	-	44,44,44,44	0
58	MG	AA	3678	1/1	0.97	0.24	-	31,31,31,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	DW	501	1/1	0.91	0.26	-	74,74,74,74	0
58	MG	AA	3693	1/1	0.96	0.21	-	48,48,48,48	0
58	MG	AA	3189	1/1	0.95	0.40	-	40,40,40,40	0
58	MG	BA	1739	1/1	0.86	0.14	-	93,93,93,93	0
58	MG	CA	3018	1/1	0.78	0.32	-	64,64,64,64	0
58	MG	DA	1718	1/1	0.74	0.15	-	77,77,77,77	0
58	MG	BW	502	1/1	0.96	0.11	-	53,53,53,53	0
58	MG	CA	3556	1/1	0.86	0.08	-	48,48,48,48	0
58	MG	AA	3611	1/1	0.97	0.15	-	51,51,51,51	0
58	MG	DA	1681	1/1	0.95	0.17	-	55,55,55,55	0
58	MG	CA	3349	1/1	0.97	0.26	-	41,41,41,41	0
58	MG	CA	3347	1/1	0.93	0.13	-	59,59,59,59	0
58	MG	AA	3463	1/1	0.94	0.30	-	46,46,46,46	0
58	MG	BA	1672	1/1	0.88	0.26	-	61,61,61,61	0
58	MG	CA	3099	1/1	0.88	0.15	-	92,92,92,92	0
58	MG	BA	1790	1/1	0.86	0.10	-	72,72,72,72	0
58	MG	AA	3361	1/1	0.98	0.15	-	29,29,29,29	0
58	MG	CA	3186	1/1	0.81	0.35	-	69,69,69,69	0
58	MG	DA	1701	1/1	0.85	0.28	-	63,63,63,63	0
58	MG	AA	3668	1/1	0.93	0.19	-	40,40,40,40	0
58	MG	CA	3319	1/1	0.93	0.19	-	67,67,67,67	0
58	MG	CA	3193	1/1	0.94	0.23	-	64,64,64,64	0
58	MG	CA	3015	1/1	0.70	0.54	-	85,85,85,85	0
58	MG	AA	3467	1/1	0.98	0.12	-	44,44,44,44	0
58	MG	CA	3591	1/1	0.95	0.14	-	83,83,83,83	0
58	MG	CA	3061	1/1	0.92	0.46	-	67,67,67,67	0
58	MG	BA	1602	1/1	0.92	0.11	-	53,53,53,53	0
58	MG	AA	3492	1/1	0.88	0.36	-	45,45,45,45	0
58	MG	AA	3046	1/1	0.96	0.23	-	34,34,34,34	0
58	MG	AA	3681	1/1	0.91	0.37	-	65,65,65,65	0
58	MG	CA	3595	1/1	0.95	0.10	-	53,53,53,53	0
58	MG	CA	3173	1/1	0.69	0.46	-	65,65,65,65	0
58	MG	AA	3287	1/1	0.97	0.40	-	47,47,47,47	0
58	MG	AA	3545	1/1	0.89	0.10	-	51,51,51,51	1
58	MG	BA	1726	1/1	0.98	0.30	-	52,52,52,52	0
58	MG	CA	3006	1/1	0.73	0.26	-	67,67,67,67	0
58	MG	CB	3008	1/1	0.89	0.12	-	59,59,59,59	0
58	MG	AA	3800	1/1	0.94	0.13	-	30,30,30,30	0
58	MG	BA	1714	1/1	0.85	0.26	-	88,88,88,88	0
58	MG	CA	3483	1/1	0.93	0.48	-	69,69,69,69	0
58	MG	BA	1806	1/1	0.70	0.28	-	81,81,81,81	0
58	MG	CA	3141	1/1	0.76	0.30	-	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3760	1/1	0.68	0.20	-	27,27,27,27	0
58	MG	AA	3106	1/1	0.93	0.28	-	52,52,52,52	0
58	MG	AA	3527	1/1	0.95	0.18	-	26,26,26,26	0
58	MG	CA	3170	1/1	0.94	0.34	-	47,47,47,47	0
58	MG	AA	3573	1/1	0.88	0.14	-	50,50,50,50	0
58	MG	CA	3062	1/1	0.64	0.25	-	68,68,68,68	0
58	MG	AA	3363	1/1	0.93	0.34	-	28,28,28,28	0
58	MG	AA	3430	1/1	0.93	0.14	-	44,44,44,44	0
58	MG	DW	502	1/1	0.78	0.11	-	84,84,84,84	0
58	MG	CA	3469	1/1	0.86	0.13	-	69,69,69,69	0
58	MG	AA	3349	1/1	0.85	0.29	-	40,40,40,40	0
58	MG	AA	3487	1/1	0.93	0.17	-	39,39,39,39	0
58	MG	BA	1605	1/1	0.52	0.28	-	67,67,67,67	0
58	MG	CA	3031	1/1	0.82	0.08	-	76,76,76,76	0
58	MG	DD	502	1/1	0.96	0.63	-	62,62,62,62	0
58	MG	AA	3031	1/1	0.93	0.25	-	10,10,10,10	1
58	MG	CA	3562	1/1	0.96	0.21	-	76,76,76,76	0
58	MG	CA	3180	1/1	0.75	0.49	-	108,108,108,108	0
58	MG	BA	1633	1/1	0.83	0.35	-	62,62,62,62	0
58	MG	AA	3027	1/1	0.93	0.48	-	77,77,77,77	0
58	MG	AA	3370	1/1	0.94	0.28	-	58,58,58,58	0
58	MG	AA	3066	1/1	0.97	0.29	-	50,50,50,50	0
58	MG	AA	3170	1/1	0.94	0.44	-	53,53,53,53	0
58	MG	CA	3357	1/1	0.92	0.06	-	80,80,80,80	0
58	MG	AA	3201	1/1	0.86	0.37	-	65,65,65,65	0
58	MG	AV	203	1/1	0.97	0.35	-	38,38,38,38	0
58	MG	CA	3241	1/1	0.61	0.55	-	107,107,107,107	0
58	MG	AA	3366	1/1	0.95	0.22	-	53,53,53,53	0
58	MG	CA	3268	1/1	0.95	0.30	-	69,69,69,69	0
58	MG	AZ	302	1/1	0.88	0.21	-	68,68,68,68	0
58	MG	BA	1634	1/1	0.68	0.51	-	100,100,100,100	0
58	MG	AA	3233	1/1	0.82	0.36	-	55,55,55,55	0
58	MG	AA	3513	1/1	0.97	0.11	-	35,35,35,35	0
58	MG	AA	3451	1/1	0.94	0.24	-	48,48,48,48	0
58	MG	AA	3627	1/1	0.98	0.17	-	54,54,54,54	0
58	MG	AA	3305	1/1	0.96	0.24	-	55,55,55,55	0
58	MG	CA	3416	1/1	0.98	0.17	-	48,48,48,48	0
58	MG	DA	1725	1/1	0.78	0.20	-	70,70,70,70	0
58	MG	CA	3112	1/1	0.88	0.26	-	61,61,61,61	0
58	MG	AA	3336	1/1	0.94	0.17	-	54,54,54,54	0
58	MG	BA	1699	1/1	0.94	0.17	-	75,75,75,75	0
58	MG	AA	3643	1/1	0.99	0.17	-	49,49,49,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	BA	1799	1/1	0.95	0.21	-	65,65,65,65	0
58	MG	CA	3232	1/1	0.96	0.09	-	60,60,60,60	0
58	MG	CA	3004	1/1	0.87	0.42	-	64,64,64,64	0
58	MG	AA	3291	1/1	0.88	0.15	-	44,44,44,44	0
58	MG	AA	3807	1/1	0.77	0.30	-	62,62,62,62	1
58	MG	BA	1649	1/1	0.89	0.24	-	68,68,68,68	0
58	MG	AA	3756	1/1	0.96	0.16	-	40,40,40,40	1
58	MG	BA	1727	1/1	0.92	0.10	-	45,45,45,45	0
58	MG	BA	1735	1/1	0.89	0.28	-	72,72,72,72	0
58	MG	CA	3187	1/1	0.91	0.28	-	37,37,37,37	0
58	MG	AA	3516	1/1	0.97	0.17	-	18,18,18,18	0
58	MG	BA	1684	1/1	0.89	0.13	-	81,81,81,81	0
58	MG	CA	3585	1/1	0.85	0.21	-	78,78,78,78	0
58	MG	AA	3834	1/1	0.80	0.22	-	58,58,58,58	0
58	MG	AA	3575	1/1	0.96	0.12	-	69,69,69,69	0
58	MG	CA	3356	1/1	0.92	0.16	-	57,57,57,57	0
58	MG	CA	3049	1/1	0.97	0.13	-	81,81,81,81	0
58	MG	AA	3477	1/1	0.88	0.17	-	58,58,58,58	0
58	MG	AA	3369	1/1	0.96	0.24	-	27,27,27,27	0
58	MG	CA	3499	1/1	0.93	0.24	-	83,83,83,83	0
58	MG	AA	3788	1/1	0.91	0.27	-	58,58,58,58	1
58	MG	AA	3207	1/1	0.94	0.18	-	22,22,22,22	1
58	MG	AA	3522	1/1	0.95	0.22	-	28,28,28,28	0
58	MG	AA	3457	1/1	0.91	0.18	-	67,67,67,67	0
58	MG	AA	3481	1/1	0.88	0.15	-	51,51,51,51	0
58	MG	AA	3582	1/1	0.96	0.41	-	66,66,66,66	0
58	MG	CA	3345	1/1	0.99	0.22	-	38,38,38,38	0
58	MG	AA	3598	1/1	0.94	0.23	-	54,54,54,54	0
58	MG	BA	1670	1/1	0.94	0.23	-	92,92,92,92	0
58	MG	BA	1708	1/1	0.76	0.20	-	79,79,79,79	0
58	MG	CA	3040	1/1	0.92	0.15	-	66,66,66,66	0
58	MG	CA	3171	1/1	0.94	0.33	-	55,55,55,55	0
58	MG	BA	1618	1/1	0.74	0.30	-	57,57,57,57	0
58	MG	CA	3460	1/1	0.70	0.47	-	109,109,109,109	0
58	MG	AA	3786	1/1	0.89	0.37	-	59,59,59,59	0
58	MG	CA	3609	1/1	0.88	0.23	-	76,76,76,76	0
58	MG	CA	3633	1/1	0.96	0.23	-	68,68,68,68	0
58	MG	AA	3242	1/1	0.90	0.27	-	72,72,72,72	0
58	MG	AA	3779	1/1	0.91	0.20	-	62,62,62,62	0
58	MG	AA	3646	1/1	0.89	0.13	-	58,58,58,58	0
58	MG	AA	3576	1/1	0.97	0.28	-	38,38,38,38	0
58	MG	BA	1639	1/1	0.92	0.37	-	49,49,49,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3312	1/1	0.97	0.14	-	51,51,51,51	0
58	MG	BA	1770	1/1	0.98	0.38	-	56,56,56,56	0
58	MG	CA	3151	1/1	0.91	0.17	-	50,50,50,50	0
58	MG	AA	3444	1/1	0.89	0.27	-	63,63,63,63	0
58	MG	AA	3790	1/1	0.96	0.06	-	49,49,49,49	0
58	MG	CA	3308	1/1	0.96	0.11	-	41,41,41,41	0
58	MG	AA	3151	1/1	0.81	0.21	-	50,50,50,50	0
58	MG	AB	3018	1/1	0.93	0.14	-	84,84,84,84	0
58	MG	AA	3376	1/1	0.97	0.17	-	19,19,19,19	0
58	MG	DA	1679	1/1	0.95	0.42	-	58,58,58,58	0
58	MG	AA	3368	1/1	0.93	0.21	-	37,37,37,37	0
58	MG	AA	3650	1/1	0.92	0.07	-	60,60,60,60	0
58	MG	CA	3598	1/1	0.80	0.09	-	73,73,73,73	0
58	MG	AA	3348	1/1	0.93	0.22	-	32,32,32,32	0
58	MG	AA	3629	1/1	0.94	0.17	-	61,61,61,61	0
58	MG	AA	3122	1/1	0.82	0.28	-	54,54,54,54	0
58	MG	AA	3554	1/1	0.89	0.15	-	49,49,49,49	0
58	MG	CA	3250	1/1	0.94	0.16	-	52,52,52,52	0
58	MG	CA	3147	1/1	0.84	0.35	-	76,76,76,76	0
58	MG	DA	1665	1/1	0.91	0.49	-	61,61,61,61	0
58	MG	DA	1671	1/1	0.90	0.59	-	83,83,83,83	0
58	MG	CA	3574	1/1	0.74	0.15	-	72,72,72,72	0
58	MG	AA	3333	1/1	0.94	0.15	-	66,66,66,66	0
58	MG	AA	3379	1/1	0.95	0.14	-	30,30,30,30	0
58	MG	AA	3653	1/1	0.88	0.15	-	68,68,68,68	0
58	MG	AA	3654	1/1	0.97	0.29	-	66,66,66,66	0
58	MG	AA	3720	1/1	0.85	0.72	-	77,77,77,77	0
58	MG	CA	3631	1/1	0.91	0.13	-	77,77,77,77	0
58	MG	AA	3203	1/1	0.91	0.33	-	46,46,46,46	1
58	MG	BA	1698	1/1	0.62	0.28	-	74,74,74,74	0
58	MG	AA	3407	1/1	0.98	0.12	-	49,49,49,49	0
58	MG	BA	1760	1/1	0.85	0.13	-	89,89,89,89	0
58	MG	AA	3347	1/1	0.97	0.10	-	38,38,38,38	0
58	MG	AA	3538	1/1	0.85	0.26	-	61,61,61,61	1
58	MG	AA	3494	1/1	0.97	0.23	-	50,50,50,50	0
58	MG	DZ	702	1/1	0.99	0.28	-	57,57,57,57	0
58	MG	CA	3338	1/1	0.97	0.17	-	63,63,63,63	0
58	MG	A2	102	1/1	0.79	0.37	-	54,54,54,54	0
58	MG	AA	3375	1/1	0.89	0.25	-	57,57,57,57	0
58	MG	CA	3022	1/1	0.96	0.20	-	38,38,38,38	0
58	MG	AA	3127	1/1	0.98	0.37	-	71,71,71,71	0
58	MG	DA	1759	1/1	0.82	0.35	-	76,76,76,76	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3198	1/1	0.91	0.13	-	37,37,37,37	0
58	MG	DA	1655	1/1	0.96	0.42	-	58,58,58,58	0
58	MG	CA	3113	1/1	0.81	0.30	-	92,92,92,92	0
58	MG	AA	3125	1/1	0.93	0.22	-	23,23,23,23	1
58	MG	CA	3207	1/1	0.96	0.15	-	71,71,71,71	0
58	MG	AA	3470	1/1	0.96	0.16	-	29,29,29,29	0
58	MG	DA	1692	1/1	0.87	0.15	-	76,76,76,76	0
58	MG	AA	3450	1/1	0.91	0.27	-	53,53,53,53	0
58	MG	AA	3068	1/1	0.92	0.48	-	65,65,65,65	0
58	MG	AA	3610	1/1	0.50	0.18	-	51,51,51,51	1
58	MG	CA	3005	1/1	0.92	0.23	-	59,59,59,59	0
58	MG	AB	3013	1/1	0.98	0.18	-	54,54,54,54	0
58	MG	A7	101	1/1	0.85	0.16	-	49,49,49,49	1
58	MG	AA	3552	1/1	0.97	0.15	-	63,63,63,63	0
58	MG	CA	3007	1/1	0.97	0.10	-	28,28,28,28	0
58	MG	DA	1635	1/1	0.86	0.35	-	65,65,65,65	0
58	MG	AA	3114	1/1	0.93	0.21	-	17,17,17,17	0
58	MG	CA	3492	1/1	0.68	0.54	-	105,105,105,105	0
58	MG	AA	3425	1/1	0.95	0.24	-	18,18,18,18	0
58	MG	CA	3572	1/1	0.93	0.16	-	54,54,54,54	0
58	MG	CA	3046	1/1	-0.24	0.35	-	114,114,114,114	0
58	MG	CA	3118	1/1	0.82	0.23	-	57,57,57,57	0
58	MG	AA	3521	1/1	0.98	0.16	-	29,29,29,29	0
58	MG	AA	3490	1/1	0.82	0.09	-	50,50,50,50	0
58	MG	CA	3486	1/1	0.95	0.27	-	69,69,69,69	0
58	MG	CA	3373	1/1	0.85	0.50	-	58,58,58,58	0
58	MG	DA	1662	1/1	0.80	0.23	-	64,64,64,64	0
58	MG	AA	3591	1/1	0.96	0.25	-	65,65,65,65	0
58	MG	AA	3703	1/1	0.84	0.10	-	76,76,76,76	0
58	MG	AA	3796	1/1	0.95	0.26	-	50,50,50,50	0
58	MG	AA	3256	1/1	0.97	0.16	-	21,21,21,21	1
58	MG	CA	3442	1/1	0.93	0.46	-	67,67,67,67	0
58	MG	BA	1749	1/1	0.95	0.12	-	48,48,48,48	0
58	MG	CA	3108	1/1	0.92	0.18	-	60,60,60,60	0
58	MG	AA	3808	1/1	0.93	0.20	-	28,28,28,28	1
58	MG	CA	3336	1/1	0.93	0.16	-	69,69,69,69	0
58	MG	AA	3165	1/1	0.83	0.42	-	57,57,57,57	0
58	MG	BA	1707	1/1	0.88	0.26	-	50,50,50,50	0
58	MG	AA	3746	1/1	0.94	0.15	-	64,64,64,64	0
58	MG	CA	3608	1/1	0.91	0.22	-	56,56,56,56	0
58	MG	CA	3149	1/1	0.92	0.08	-	66,66,66,66	0
58	MG	AA	3317	1/1	0.97	0.14	-	58,58,58,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	BA	1636	1/1	0.96	0.40	-	64,64,64,64	0
58	MG	DA	1642	1/1	0.93	0.20	-	66,66,66,66	0
58	MG	AA	3466	1/1	0.95	0.12	-	63,63,63,63	0
58	MG	CA	3161	1/1	0.94	0.19	-	66,66,66,66	0
58	MG	CA	3570	1/1	0.86	0.21	-	77,77,77,77	0
58	MG	BA	1713	1/1	0.96	0.27	-	55,55,55,55	0
58	MG	AA	3665	1/1	0.85	0.34	-	85,85,85,85	0
58	MG	BZ	701	1/1	0.22	0.30	-	137,137,137,137	0
58	MG	CA	3515	1/1	0.85	0.20	-	79,79,79,79	0
58	MG	CA	3551	1/1	0.82	0.17	-	88,88,88,88	0
58	MG	DA	1676	1/1	0.83	0.09	-	78,78,78,78	0
58	MG	CA	3075	1/1	0.84	0.40	-	71,71,71,71	0
58	MG	DA	1646	1/1	0.95	0.34	-	62,62,62,62	0
58	MG	DA	1678	1/1	0.71	0.37	-	82,82,82,82	0
58	MG	AA	3327	1/1	0.97	0.14	-	13,13,13,13	0
58	MG	CQ	202	1/1	0.93	0.66	-	64,64,64,64	0
58	MG	AA	3219	1/1	0.92	0.29	-	58,58,58,58	0
58	MG	CA	3092	1/1	0.98	0.16	-	70,70,70,70	0
58	MG	CA	3495	1/1	0.85	0.15	-	65,65,65,65	0
58	MG	AA	3016	1/1	0.76	0.40	-	59,59,59,59	0
58	MG	CA	3174	1/1	0.78	0.54	-	61,61,61,61	0
58	MG	DA	1693	1/1	0.90	0.22	-	54,54,54,54	0
58	MG	AA	3382	1/1	0.82	0.17	-	36,36,36,36	1
58	MG	AA	3104	1/1	0.96	0.17	-	28,28,28,28	0
58	MG	CA	3316	1/1	0.94	0.17	-	50,50,50,50	0
58	MG	CA	3643	1/1	0.93	0.09	-	76,76,76,76	0
58	MG	CA	3581	1/1	0.92	0.12	-	81,81,81,81	0
58	MG	BA	1807	1/1	0.94	0.15	-	83,83,83,83	0
58	MG	AA	3261	1/1	0.85	0.36	-	51,51,51,51	0
58	MG	BA	1729	1/1	0.97	0.18	-	49,49,49,49	0
58	MG	DA	1622	1/1	0.89	0.54	-	77,77,77,77	0
58	MG	BA	1796	1/1	0.88	0.11	-	70,70,70,70	0
58	MG	CE	304	1/1	0.85	0.76	-	68,68,68,68	0
58	MG	BA	1762	1/1	0.90	0.07	-	52,52,52,52	1
58	MG	AA	3195	1/1	0.98	0.19	-	50,50,50,50	0
58	MG	AA	3155	1/1	0.84	0.31	-	64,64,64,64	0
58	MG	DA	1760	1/1	0.95	0.37	-	66,66,66,66	0
58	MG	AA	3058	1/1	0.90	0.20	-	35,35,35,35	0
58	MG	CA	3535	1/1	0.93	0.30	-	69,69,69,69	0
58	MG	DA	1710	1/1	0.97	0.20	-	70,70,70,70	0
58	MG	CA	3440	1/1	0.98	0.31	-	57,57,57,57	0
58	MG	AA	3810	1/1	0.91	0.24	-	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3115	1/1	0.72	0.45	-	67,67,67,67	1
58	MG	AA	3141	1/1	0.92	0.47	-	40,40,40,40	0
58	MG	BA	1720	1/1	0.98	0.45	-	62,62,62,62	0
58	MG	AA	3192	1/1	0.18	0.65	-	76,76,76,76	0
58	MG	AA	3148	1/1	0.89	0.34	-	68,68,68,68	0
58	MG	CB	3001	1/1	0.89	0.20	-	99,99,99,99	0
58	MG	DA	1628	1/1	0.86	0.49	-	74,74,74,74	0
58	MG	AA	3060	1/1	0.88	0.69	-	65,65,65,65	0
58	MG	AA	3010	1/1	0.76	0.62	-	68,68,68,68	0
58	MG	BA	1746	1/1	0.89	0.17	-	83,83,83,83	0
58	MG	CA	3394	1/1	0.96	0.16	-	55,55,55,55	0
58	MG	DA	1659	1/1	0.91	0.08	-	64,64,64,64	0
58	MG	AA	3096	1/1	0.85	0.14	-	63,63,63,63	0
58	MG	CA	3167	1/1	0.93	0.10	-	60,60,60,60	0
58	MG	CA	3465	1/1	0.95	0.34	-	66,66,66,66	0
58	MG	BA	1724	1/1	0.81	0.20	-	64,64,64,64	0
58	MG	CA	3543	1/1	0.90	0.20	-	63,63,63,63	0
58	MG	BA	1779	1/1	0.91	0.15	-	46,46,46,46	1
58	MG	CA	3065	1/1	0.89	0.12	-	52,52,52,52	0
58	MG	AB	3019	1/1	0.93	0.18	-	65,65,65,65	0
58	MG	CA	3095	1/1	0.90	0.31	-	64,64,64,64	0
58	MG	CA	3447	1/1	0.82	0.44	-	80,80,80,80	0
58	MG	BA	1737	1/1	0.95	0.27	-	63,63,63,63	0
58	MG	AA	3696	1/1	0.88	0.16	-	69,69,69,69	0
58	MG	AA	3269	1/1	0.77	0.46	-	84,84,84,84	0
58	MG	AA	3747	1/1	0.92	0.39	-	58,58,58,58	0
58	MG	CA	3247	1/1	0.92	0.66	-	66,66,66,66	0
58	MG	CA	3583	1/1	0.76	0.30	-	78,78,78,78	1
58	MG	CA	3344	1/1	0.95	0.08	-	87,87,87,87	0
58	MG	CA	3429	1/1	0.92	0.36	-	74,74,74,74	0
58	MG	DA	1698	1/1	0.93	0.40	-	68,68,68,68	0
58	MG	AA	3076	1/1	0.96	0.25	-	92,92,92,92	0
58	MG	AA	3391	1/1	0.96	0.15	-	45,45,45,45	0
58	MG	AA	3312	1/1	0.96	0.19	-	55,55,55,55	0
58	MG	CA	3398	1/1	0.85	0.10	-	65,65,65,65	0
58	MG	CA	3287	1/1	0.94	0.15	-	52,52,52,52	0
58	MG	DA	1738	1/1	0.94	0.59	-	80,80,80,80	0
58	MG	AA	3732	1/1	0.69	0.23	-	68,68,68,68	0
58	MG	CA	3256	1/1	0.97	0.18	-	40,40,40,40	0
58	MG	AA	3688	1/1	0.91	0.19	-	25,25,25,25	1
58	MG	AA	3700	1/1	0.95	0.22	-	35,35,35,35	0
58	MG	AA	3199	1/1	0.89	0.18	-	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3639	1/1	0.93	0.32	-	55,55,55,55	0
58	MG	AA	3615	1/1	0.85	0.20	-	35,35,35,35	1
58	MG	DA	1767	1/1	0.90	0.14	-	74,74,74,74	0
58	MG	AA	3597	1/1	0.90	0.09	-	63,63,63,63	0
58	MG	DA	1675	1/1	0.92	0.36	-	74,74,74,74	0
58	MG	AA	3496	1/1	0.87	0.17	-	52,52,52,52	0
58	MG	BA	1709	1/1	0.51	0.21	-	104,104,104,104	0
58	MG	AA	3236	1/1	0.69	0.21	-	57,57,57,57	0
58	MG	AA	3119	1/1	0.93	0.34	-	40,40,40,40	1
58	MG	CA	3484	1/1	0.96	0.27	-	76,76,76,76	0
58	MG	AA	3280	1/1	0.88	0.34	-	47,47,47,47	0
58	MG	AA	3178	1/1	0.44	0.53	-	78,78,78,78	0
58	MG	CA	3262	1/1	0.92	0.17	-	61,61,61,61	0
58	MG	DA	1733	1/1	0.70	0.17	-	83,83,83,83	0
58	MG	DA	1747	1/1	0.87	0.17	-	70,70,70,70	0
58	MG	BA	1767	1/1	0.72	0.09	-	58,58,58,58	0
58	MG	AA	3216	1/1	0.90	0.21	-	51,51,51,51	0
58	MG	AA	3194	1/1	0.93	0.48	-	44,44,44,44	0
58	MG	BA	1667	1/1	0.39	0.21	-	89,89,89,89	0
58	MG	CA	3369	1/1	0.98	0.13	-	59,59,59,59	0
58	MG	DA	1661	1/1	0.86	0.41	-	70,70,70,70	0
58	MG	DZ	701	1/1	0.57	0.61	-	111,111,111,111	0
58	MG	AA	3782	1/1	0.94	0.27	-	74,74,74,74	0
58	MG	CA	3246	1/1	0.71	0.50	-	59,59,59,59	0
58	MG	DA	1731	1/1	0.93	0.51	-	82,82,82,82	0
58	MG	AA	3405	1/1	0.93	0.40	-	46,46,46,46	0
58	MG	CA	3305	1/1	0.98	0.18	-	62,62,62,62	0
58	MG	DA	1745	1/1	0.94	0.18	-	61,61,61,61	0
58	MG	AB	3005	1/1	0.97	0.28	-	69,69,69,69	0
58	MG	CA	3434	1/1	0.92	0.18	-	28,28,28,28	0
58	MG	AA	3601	1/1	0.97	0.39	-	47,47,47,47	0
58	MG	CR	201	1/1	0.85	0.45	-	51,51,51,51	0
58	MG	CA	3085	1/1	0.79	0.54	-	66,66,66,66	0
58	MG	CA	3143	1/1	0.92	0.71	-	57,57,57,57	0
58	MG	AA	3801	1/1	0.82	0.07	-	88,88,88,88	0
58	MG	BA	1669	1/1	0.67	0.30	-	73,73,73,73	0
58	MG	CA	3228	1/1	0.86	0.26	-	63,63,63,63	0
58	MG	AA	3549	1/1	0.97	0.08	-	54,54,54,54	0
58	MG	AA	3474	1/1	0.97	0.28	-	53,53,53,53	0
58	MG	CD	302	1/1	0.80	0.45	-	95,95,95,95	0
58	MG	BA	1635	1/1	0.96	0.23	-	72,72,72,72	0
58	MG	AA	3568	1/1	0.94	0.19	-	16,16,16,16	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3208	1/1	0.81	0.44	-	84,84,84,84	0
58	MG	AA	3399	1/1	0.97	0.10	-	18,18,18,18	0
58	MG	CA	3391	1/1	0.97	0.05	-	63,63,63,63	0
58	MG	CA	3601	1/1	0.81	0.08	-	75,75,75,75	0
58	MG	CA	3293	1/1	0.97	0.06	-	71,71,71,71	0
58	MG	AA	3360	1/1	0.88	0.12	-	114,114,114,114	0
58	MG	DA	1758	1/1	0.94	0.16	-	71,71,71,71	0
58	MG	CA	3576	1/1	0.95	0.10	-	36,36,36,36	0
58	MG	AA	3264	1/1	0.91	0.08	-	62,62,62,62	0
58	MG	CA	3132	1/1	0.86	0.20	-	30,30,30,30	0
58	MG	CA	3549	1/1	0.93	0.06	-	57,57,57,57	0
58	MG	BA	1606	1/1	0.89	0.17	-	74,74,74,74	0
58	MG	CA	3076	1/1	0.74	0.41	-	84,84,84,84	0
58	MG	CA	3466	1/1	0.96	0.48	-	57,57,57,57	0
58	MG	CA	3561	1/1	0.58	0.19	-	95,95,95,95	0
58	MG	CA	3179	1/1	0.90	0.53	-	75,75,75,75	0
58	MG	AA	3218	1/1	0.96	0.16	-	38,38,38,38	0
58	MG	AA	3644	1/1	0.94	0.18	-	56,56,56,56	0
58	MG	AA	3310	1/1	0.90	0.26	-	58,58,58,58	0
58	MG	CA	3334	1/1	0.96	0.27	-	44,44,44,44	0
58	MG	AA	3753	1/1	0.89	0.18	-	41,41,41,41	0
58	MG	AA	3803	1/1	0.86	0.32	-	62,62,62,62	0
58	MG	DA	1683	1/1	0.95	0.37	-	54,54,54,54	0
58	MG	AA	3065	1/1	0.87	0.54	-	48,48,48,48	0
58	MG	BA	1780	1/1	0.87	0.20	-	81,81,81,81	0
58	MG	DA	1616	1/1	0.95	0.24	-	64,64,64,64	0
58	MG	AA	3758	1/1	0.78	0.36	-	81,81,81,81	0
58	MG	CA	3115	1/1	0.91	0.23	-	76,76,76,76	0
58	MG	CA	3016	1/1	0.89	0.56	-	80,80,80,80	0
58	MG	AA	3715	1/1	0.92	0.08	-	66,66,66,66	0
58	MG	AA	3489	1/1	0.93	0.09	-	64,64,64,64	0
58	MG	CO	201	1/1	0.94	0.16	-	64,64,64,64	0
58	MG	AA	3599	1/1	0.83	0.41	-	58,58,58,58	0
58	MG	CA	3523	1/1	0.86	0.32	-	40,40,40,40	0
58	MG	CA	3001	1/1	0.75	0.40	-	73,73,73,73	0
58	MG	AA	3751	1/1	0.78	0.64	-	56,56,56,56	1
58	MG	AA	3285	1/1	0.91	0.38	-	45,45,45,45	0
58	MG	BA	1722	1/1	0.97	0.44	-	55,55,55,55	0
58	MG	DA	1728	1/1	0.90	0.16	-	63,63,63,63	0
58	MG	BL	201	1/1	0.88	0.27	-	80,80,80,80	0
58	MG	AA	3691	1/1	0.75	0.21	-	89,89,89,89	0
58	MG	CA	3155	1/1	0.73	0.25	-	86,86,86,86	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3675	1/1	0.96	0.10	-	38,38,38,38	0
58	MG	AA	3421	1/1	0.97	0.18	-	12,12,12,12	0
58	MG	AA	3397	1/1	0.94	0.13	-	15,15,15,15	0
58	MG	AN	3002	1/1	0.97	0.16	-	27,27,27,27	0
58	MG	AA	3094	1/1	0.89	0.75	-	111,111,111,111	0
58	MG	CA	3620	1/1	0.51	0.27	-	69,69,69,69	0
58	MG	CA	3034	1/1	0.88	0.58	-	101,101,101,101	0
58	MG	AA	3509	1/1	0.96	0.16	-	49,49,49,49	0
58	MG	AA	3024	1/1	0.78	0.14	-	57,57,57,57	0
58	MG	BA	1754	1/1	0.91	0.09	-	98,98,98,98	0
58	MG	AA	3772	1/1	0.95	0.28	-	17,17,17,17	1
58	MG	CA	3238	1/1	0.90	0.26	-	59,59,59,59	0
58	MG	BA	1601	1/1	0.88	0.27	-	93,93,93,93	0
58	MG	BA	1788	1/1	0.87	0.12	-	75,75,75,75	0
58	MG	AA	3745	1/1	0.92	0.18	-	29,29,29,29	0
58	MG	CA	3497	1/1	0.94	0.32	-	73,73,73,73	0
58	MG	CA	3456	1/1	0.98	0.13	-	60,60,60,60	0
58	MG	CA	3231	1/1	0.96	0.38	-	57,57,57,57	0
58	MG	CA	3382	1/1	0.95	0.19	-	40,40,40,40	0
58	MG	CA	3525	1/1	0.93	0.30	-	83,83,83,83	0
58	MG	AA	3042	1/1	0.85	0.43	-	38,38,38,38	0
58	MG	AA	3198	1/1	0.95	0.07	-	58,58,58,58	0
58	MG	AA	3129	1/1	0.91	0.50	-	66,66,66,66	1
58	MG	AA	3740	1/1	0.92	0.17	-	45,45,45,45	0
58	MG	C5	101	1/1	0.95	0.64	-	65,65,65,65	0
58	MG	DA	1663	1/1	0.71	0.22	-	72,72,72,72	0
58	MG	CA	3120	1/1	0.96	0.52	-	62,62,62,62	0
58	MG	DA	1754	1/1	0.57	0.37	-	81,81,81,81	0
58	MG	AA	3759	1/1	0.91	0.29	-	65,65,65,65	0
58	MG	AA	3783	1/1	0.96	0.21	-	54,54,54,54	0
58	MG	DA	1711	1/1	0.93	0.32	-	45,45,45,45	0
58	MG	AA	3288	1/1	0.92	0.17	-	24,24,24,24	0
58	MG	AA	3153	1/1	0.89	0.30	-	59,59,59,59	0
58	MG	BA	1803	1/1	0.93	0.22	-	64,64,64,64	0
58	MG	DA	1753	1/1	0.96	0.40	-	70,70,70,70	0
58	MG	BA	1628	1/1	0.89	0.56	-	55,55,55,55	0
58	MG	CA	3077	1/1	0.90	0.39	-	66,66,66,66	0
58	MG	CA	3519	1/1	0.89	0.28	-	79,79,79,79	0
58	MG	DA	1727	1/1	0.89	0.17	-	66,66,66,66	0
58	MG	AB	3001	1/1	0.76	0.57	-	87,87,87,87	0
58	MG	CA	3479	1/1	0.95	0.22	-	46,46,46,46	0
58	MG	AA	3784	1/1	0.48	0.38	-	74,74,74,74	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	DA	1752	1/1	0.96	0.22	-	74,74,74,74	0
58	MG	CA	3418	1/1	0.88	0.27	-	41,41,41,41	0
58	MG	DA	1735	1/1	0.89	0.57	-	83,83,83,83	0
58	MG	CA	3548	1/1	0.87	0.14	-	48,48,48,48	1
58	MG	AA	3259	1/1	0.98	0.33	-	20,20,20,20	1
58	MG	AA	3512	1/1	0.93	0.34	-	60,60,60,60	0
58	MG	BA	1689	1/1	0.53	0.77	-	91,91,91,91	0
58	MG	CA	3563	1/1	0.69	0.20	-	91,91,91,91	0
58	MG	AA	3648	1/1	0.97	0.17	-	40,40,40,40	0
58	MG	CA	3323	1/1	0.88	0.23	-	67,67,67,67	0
58	MG	AA	3353	1/1	0.91	0.08	-	76,76,76,76	0
58	MG	AA	3645	1/1	0.93	0.51	-	79,79,79,79	0
58	MG	AA	3652	1/1	0.91	0.28	-	77,77,77,77	0
58	MG	DA	1744	1/1	0.89	0.10	-	79,79,79,79	0
58	MG	AA	3062	1/1	0.89	0.35	-	67,67,67,67	0
58	MG	AA	3338	1/1	0.95	0.15	-	30,30,30,30	0
58	MG	AA	3533	1/1	0.95	0.14	-	22,22,22,22	0
58	MG	BA	1619	1/1	0.91	0.19	-	52,52,52,52	0
58	MG	AA	3208	1/1	0.85	0.32	-	61,61,61,61	0
58	MG	BA	1665	1/1	0.89	0.38	-	55,55,55,55	0
58	MG	AA	3689	1/1	0.92	0.16	-	55,55,55,55	1
58	MG	CA	3191	1/1	0.93	0.29	-	65,65,65,65	0
58	MG	CA	3355	1/1	0.97	0.11	-	41,41,41,41	0
58	MG	AA	3464	1/1	0.81	0.15	-	63,63,63,63	0
58	MG	DA	1751	1/1	0.72	0.26	-	81,81,81,81	0
58	MG	CA	3380	1/1	0.95	0.20	-	71,71,71,71	0
58	MG	BA	1681	1/1	0.53	1.14	-	84,84,84,84	0
58	MG	BA	1785	1/1	0.69	0.33	-	87,87,87,87	0
58	MG	AA	3701	1/1	0.86	0.46	-	43,43,43,43	1
58	MG	CA	3255	1/1	0.88	0.48	-	67,67,67,67	0
58	MG	CA	3388	1/1	0.85	0.12	-	83,83,83,83	0
58	MG	AA	3742	1/1	0.94	0.25	-	82,82,82,82	0
58	MG	DA	1674	1/1	0.94	0.40	-	62,62,62,62	0
58	MG	AA	3124	1/1	0.95	0.63	-	63,63,63,63	0
58	MG	AA	3088	1/1	0.91	0.33	-	34,34,34,34	0
58	MG	CA	3605	1/1	0.93	0.27	-	73,73,73,73	0
58	MG	AA	3553	1/1	0.93	0.17	-	40,40,40,40	0
58	MG	AA	3690	1/1	0.90	0.25	-	71,71,71,71	0
58	MG	BA	1730	1/1	0.88	0.26	-	78,78,78,78	0
58	MG	AA	3149	1/1	0.90	0.27	-	62,62,62,62	0
58	MG	DA	1708	1/1	0.90	0.10	-	87,87,87,87	0
58	MG	AA	3813	1/1	0.93	0.59	-	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
58	MG	CA	3194	1/1	0.89	0.57	-	72,72,72,72	0
58	MG	AA	3220	1/1	0.82	0.28	-	62,62,62,62	0
58	MG	DA	1762	1/1	0.80	0.08	-	73,73,73,73	0
58	MG	AA	3241	1/1	0.94	0.21	-	69,69,69,69	0
58	MG	BA	1745	1/1	0.88	0.43	-	66,66,66,66	0
58	MG	CA	3284	1/1	0.95	0.20	-	92,92,92,92	0
58	MG	AA	3685	1/1	0.93	0.20	-	47,47,47,47	0
58	MG	AA	3682	1/1	0.96	0.30	-	53,53,53,53	0
58	MG	AA	3468	1/1	0.94	0.37	-	52,52,52,52	0
58	MG	AA	3390	1/1	0.89	0.22	-	35,35,35,35	0
58	MG	CA	3244	1/1	0.72	0.67	-	78,78,78,78	0
58	MG	AA	3683	1/1	0.96	0.37	-	67,67,67,67	0
58	MG	CA	3415	1/1	0.98	0.21	-	31,31,31,31	1
58	MG	CA	3070	1/1	0.63	0.81	-	87,87,87,87	0
58	MG	AA	3510	1/1	0.95	0.36	-	58,58,58,58	0
58	MG	CA	3625	1/1	0.85	0.54	-	79,79,79,79	0
58	MG	AA	3618	1/1	0.94	0.16	-	38,38,38,38	0
58	MG	AP	202	1/1	0.88	0.18	-	40,40,40,40	0
58	MG	CA	3421	1/1	0.57	0.25	-	76,76,76,76	0
58	MG	AA	3475	1/1	0.96	0.14	-	62,62,62,62	0
58	MG	AA	3716	1/1	0.96	0.17	-	57,57,57,57	0
58	MG	CA	3612	1/1	0.94	0.23	-	68,68,68,68	0
58	MG	AA	3710	1/1	0.79	0.27	-	75,75,75,75	0
58	MG	AA	3676	1/1	0.95	0.19	-	66,66,66,66	0
58	MG	DA	1607	1/1	0.84	0.32	-	61,61,61,61	0
58	MG	CA	3450	1/1	0.97	0.19	-	48,48,48,48	0
58	MG	AA	3177	1/1	0.94	0.34	-	61,61,61,61	0
58	MG	AA	3667	1/1	0.95	0.20	-	28,28,28,28	0
58	MG	C8	5001	1/1	0.90	0.36	-	51,51,51,51	0
58	MG	CA	3532	1/1	0.75	0.12	-	79,79,79,79	0
58	MG	BA	1725	1/1	0.90	0.11	-	59,59,59,59	0
58	MG	AA	3350	1/1	0.92	0.29	-	31,31,31,31	0
58	MG	BA	1772	1/1	0.81	0.18	-	70,70,70,70	0
58	MG	AA	3578	1/1	0.94	0.42	-	55,55,55,55	0
58	MG	AA	3478	1/1	0.98	0.16	-	40,40,40,40	0
58	MG	CA	3505	1/1	0.93	0.27	-	73,73,73,73	0
58	MG	BA	1776	1/1	0.96	0.26	-	64,64,64,64	0
58	MG	AF	305	1/1	0.94	0.22	-	55,55,55,55	0
58	MG	CA	3116	1/1	0.81	0.43	-	75,75,75,75	0
58	MG	DA	1756	1/1	0.93	0.24	-	68,68,68,68	0
58	MG	CA	3537	1/1	0.69	0.29	-	78,78,78,78	0
58	MG	AA	3461	1/1	0.87	0.50	-	66,66,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	BA	1660	1/1	0.73	0.21	-	70,70,70,70	0
58	MG	CA	3473	1/1	0.93	0.17	-	54,54,54,54	0
58	MG	CA	3637	1/1	0.98	0.36	-	59,59,59,59	0
58	MG	CA	3335	1/1	0.73	0.18	-	62,62,62,62	0
58	MG	CA	3363	1/1	0.86	0.24	-	66,66,66,66	0
58	MG	AA	3230	1/1	0.82	0.58	-	80,80,80,80	0
58	MG	AA	3657	1/1	0.90	0.24	-	43,43,43,43	1
58	MG	AA	3328	1/1	0.94	0.17	-	18,18,18,18	0
58	MG	DA	1734	1/1	0.86	0.25	-	83,83,83,83	0
58	MG	AA	3587	1/1	0.91	0.31	-	53,53,53,53	0
58	MG	CB	3003	1/1	0.95	0.13	-	77,77,77,77	0
58	MG	DA	1742	1/1	0.96	0.33	-	79,79,79,79	0
58	MG	CA	3379	1/1	0.90	0.29	-	65,65,65,65	0
58	MG	AA	3303	1/1	0.91	0.33	-	56,56,56,56	0
58	MG	AA	3471	1/1	0.95	0.26	-	34,34,34,34	0
58	MG	CA	3481	1/1	0.58	0.34	-	91,91,91,91	0
58	MG	AB	3002	1/1	0.98	0.18	-	59,59,59,59	0
58	MG	BA	1658	1/1	0.93	0.51	-	76,76,76,76	0
58	MG	AA	3417	1/1	0.97	0.17	-	25,25,25,25	0
58	MG	AA	3265	1/1	0.87	0.50	-	77,77,77,77	0
58	MG	DA	1634	1/1	0.59	0.34	-	90,90,90,90	0
58	MG	AA	3234	1/1	0.95	0.17	-	30,30,30,30	1
58	MG	CA	3181	1/1	0.94	0.24	-	47,47,47,47	0
58	MG	AA	3098	1/1	0.93	0.31	-	58,58,58,58	0
58	MG	CA	3653	1/1	0.78	0.39	-	95,95,95,95	0
58	MG	DA	1696	1/1	0.86	0.20	-	91,91,91,91	0
58	MG	CA	3592	1/1	0.95	0.61	-	76,76,76,76	0
58	MG	CA	3039	1/1	0.82	0.91	-	71,71,71,71	0
58	MG	AA	3204	1/1	0.72	0.39	-	57,57,57,57	0
58	MG	AA	3186	1/1	0.83	0.16	-	37,37,37,37	0
58	MG	AO	5001	1/1	0.87	0.18	-	55,55,55,55	0
58	MG	CA	3408	1/1	0.98	0.13	-	64,64,64,64	0
58	MG	AA	3308	1/1	0.81	0.13	-	30,30,30,30	0
58	MG	CA	3528	1/1	0.83	0.42	-	79,79,79,79	0
58	MG	CA	3233	1/1	0.91	0.46	-	71,71,71,71	0
58	MG	CA	3502	1/1	0.99	0.14	-	70,70,70,70	0
58	MG	AA	3176	1/1	0.90	0.31	-	50,50,50,50	0
58	MG	CA	3496	1/1	0.98	0.17	-	63,63,63,63	0
58	MG	AA	3172	1/1	0.85	0.75	-	71,71,71,71	0
58	MG	CA	3196	1/1	0.92	0.56	-	64,64,64,64	0
58	MG	AA	3484	1/1	0.92	0.23	-	35,35,35,35	0
58	MG	AA	3594	1/1	0.92	0.23	-	43,43,43,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3412	1/1	0.88	0.26	-	81,81,81,81	0
58	MG	DA	1716	1/1	0.90	0.37	-	78,78,78,78	0
58	MG	CA	3539	1/1	0.95	0.43	-	73,73,73,73	0
58	MG	CA	3058	1/1	0.72	0.42	-	77,77,77,77	0
58	MG	AA	3454	1/1	0.98	0.12	-	51,51,51,51	0
58	MG	CA	3622	1/1	0.84	0.24	-	55,55,55,55	0
58	MG	AA	3337	1/1	0.97	0.24	-	10,10,10,10	0
58	MG	AA	3319	1/1	0.90	0.18	-	69,69,69,69	0
58	MG	DW	503	1/1	0.62	0.18	-	84,84,84,84	0
58	MG	AA	3108	1/1	0.76	0.48	-	125,125,125,125	0
58	MG	CA	3260	1/1	0.97	0.23	-	65,65,65,65	0
58	MG	AA	3346	1/1	0.94	0.21	-	46,46,46,46	0
58	MG	AB	3015	1/1	0.96	0.18	-	40,40,40,40	0
58	MG	CA	3131	1/1	0.80	0.21	-	62,62,62,62	0
58	MG	CA	3117	1/1	0.80	0.31	-	68,68,68,68	0
58	MG	AA	3386	1/1	0.98	0.22	-	29,29,29,29	0
58	MG	CA	3474	1/1	0.83	0.33	-	76,76,76,76	0
58	MG	AW	3004	1/1	0.89	0.32	-	65,65,65,65	0
58	MG	AA	3612	1/1	0.86	0.21	-	56,56,56,56	0
58	MG	CA	3270	1/1	0.98	0.19	-	35,35,35,35	0
58	MG	CA	3130	1/1	0.91	0.68	-	73,73,73,73	0
58	MG	AA	3639	1/1	0.87	0.18	-	77,77,77,77	0
58	MG	BA	1743	1/1	0.95	0.18	-	52,52,52,52	0
58	MG	CA	3235	1/1	0.85	0.61	-	75,75,75,75	0
58	MG	CA	3368	1/1	0.92	0.21	-	59,59,59,59	0
58	MG	CA	3176	1/1	0.96	0.48	-	50,50,50,50	0
58	MG	CA	3402	1/1	0.93	0.12	-	70,70,70,70	0
58	MG	CA	3471	1/1	0.94	0.17	-	45,45,45,45	0
58	MG	BA	1736	1/1	0.85	0.12	-	73,73,73,73	0
58	MG	CA	3032	1/1	0.85	0.58	-	67,67,67,67	0
58	MG	CA	3300	1/1	0.80	0.40	-	86,86,86,86	0
58	MG	CA	3142	1/1	0.90	0.27	-	69,69,69,69	0
58	MG	CA	3294	1/1	0.75	0.21	-	83,83,83,83	0
58	MG	CA	3289	1/1	0.97	0.32	-	42,42,42,42	0
58	MG	AA	3776	1/1	0.94	0.12	-	40,40,40,40	0
58	MG	CA	3558	1/1	0.91	0.20	-	51,51,51,51	1
58	MG	AA	3166	1/1	0.94	0.17	-	31,31,31,31	0
58	MG	CA	3096	1/1	0.82	0.35	-	68,68,68,68	0
58	MG	AA	3465	1/1	0.94	0.21	-	39,39,39,39	0
58	MG	DA	1656	1/1	0.79	0.23	-	75,75,75,75	0
58	MG	CA	3406	1/1	0.86	0.13	-	70,70,70,70	0
58	MG	AA	3531	1/1	0.92	0.30	-	62,62,62,62	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CB	3006	1/1	0.85	0.13	-	83,83,83,83	0
58	MG	AA	3409	1/1	0.98	0.20	-	45,45,45,45	0
58	MG	CA	3072	1/1	0.86	0.28	-	56,56,56,56	0
58	MG	BA	1625	1/1	0.87	0.32	-	57,57,57,57	0
58	MG	AA	3640	1/1	0.72	0.44	-	77,77,77,77	0
58	MG	CA	3468	1/1	0.82	0.16	-	61,61,61,61	0
58	MG	AA	3030	1/1	0.93	0.32	-	26,26,26,26	1
58	MG	CA	3663	1/1	0.85	0.39	-	91,91,91,91	0
58	MG	CA	3504	1/1	0.88	0.08	-	62,62,62,62	0
58	MG	DA	1714	1/1	0.97	0.12	-	51,51,51,51	0
58	MG	CA	3210	1/1	0.93	0.31	-	75,75,75,75	0
58	MG	AA	3603	1/1	0.91	0.19	-	63,63,63,63	0
58	MG	CA	3552	1/1	0.97	0.14	-	69,69,69,69	0
58	MG	AW	3002	1/1	0.91	0.28	-	55,55,55,55	0
58	MG	AA	3180	1/1	0.89	0.31	-	94,94,94,94	0
58	MG	CA	3575	1/1	0.93	0.10	-	43,43,43,43	1
58	MG	AA	3008	1/1	0.95	0.17	-	19,19,19,19	0
58	MG	CA	3051	1/1	0.97	0.44	-	63,63,63,63	0
58	MG	CA	3414	1/1	0.93	0.20	-	50,50,50,50	0
58	MG	BA	1774	1/1	0.83	0.13	-	61,61,61,61	0
58	MG	BA	1642	1/1	0.85	0.42	-	69,69,69,69	0
58	MG	AA	3729	1/1	0.84	0.11	-	38,38,38,38	0
58	MG	AA	3279	1/1	0.88	0.35	-	53,53,53,53	0
58	MG	AA	3074	1/1	0.97	0.35	-	15,15,15,15	0
58	MG	CA	3437	1/1	0.97	0.11	-	48,48,48,48	0
58	MG	AA	3306	1/1	0.81	0.19	-	47,47,47,47	0
58	MG	BA	1804	1/1	0.92	0.18	-	67,67,67,67	0
58	MG	AA	3680	1/1	0.92	0.25	-	59,59,59,59	0
58	MG	CA	3292	1/1	0.89	0.08	-	73,73,73,73	0
58	MG	CA	3054	1/1	0.88	0.14	-	71,71,71,71	0
58	MG	CA	3071	1/1	0.32	0.51	-	97,97,97,97	0
58	MG	CA	3123	1/1	0.83	0.87	-	88,88,88,88	0
58	MG	AA	3123	1/1	0.92	0.37	-	37,37,37,37	1
58	MG	AA	3766	1/1	0.60	0.20	-	72,72,72,72	0
58	MG	AA	3273	1/1	0.83	0.33	-	90,90,90,90	0
58	MG	AA	3028	1/1	0.87	0.37	-	51,51,51,51	1
58	MG	DA	1615	1/1	0.98	0.54	-	53,53,53,53	0
58	MG	CA	3422	1/1	0.95	0.26	-	55,55,55,55	0
58	MG	BA	1682	1/1	0.89	0.11	-	69,69,69,69	0
58	MG	DA	1640	1/1	0.95	0.18	-	79,79,79,79	0
58	MG	AA	3362	1/1	0.87	0.48	-	69,69,69,69	0
58	MG	AA	3774	1/1	0.86	0.23	-	80,80,80,80	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	BA	1784	1/1	0.88	0.22	-	60,60,60,60	0
58	MG	CA	3206	1/1	0.93	0.59	-	56,56,56,56	0
58	MG	CA	3271	1/1	0.90	0.32	-	57,57,57,57	0
58	MG	CA	3079	1/1	0.94	0.34	-	57,57,57,57	0
58	MG	CA	3145	1/1	0.92	0.08	-	79,79,79,79	0
58	MG	AA	3091	1/1	0.89	0.39	-	38,38,38,38	1
58	MG	BA	1609	1/1	0.89	0.14	-	69,69,69,69	0
58	MG	CA	3649	1/1	0.78	0.35	-	85,85,85,85	0
58	MG	CD	301	1/1	0.74	0.47	-	81,81,81,81	0
58	MG	AA	3017	1/1	0.81	0.16	-	78,78,78,78	0
58	MG	AA	3737	1/1	0.94	0.15	-	29,29,29,29	0
58	MG	AA	3245	1/1	0.82	0.78	-	69,69,69,69	0
58	MG	CA	3303	1/1	0.93	0.47	-	54,54,54,54	0
58	MG	AA	3767	1/1	0.60	0.31	-	63,63,63,63	1
58	MG	AA	3077	1/1	0.92	0.34	-	50,50,50,50	0
58	MG	CA	3288	1/1	0.97	0.25	-	54,54,54,54	0
58	MG	AB	3004	1/1	0.79	0.30	-	89,89,89,89	0
58	MG	CA	3374	1/1	0.80	0.43	-	76,76,76,76	0
58	MG	CA	3042	1/1	0.60	0.73	-	95,95,95,95	0
58	MG	BA	1652	1/1	0.77	0.14	-	69,69,69,69	0
58	MG	AA	3447	1/1	0.80	0.18	-	61,61,61,61	0
58	MG	BM	202	1/1	0.86	0.39	-	65,65,65,65	0
58	MG	AA	3778	1/1	0.93	0.21	-	54,54,54,54	0
58	MG	BA	1691	1/1	0.86	0.56	-	74,74,74,74	0
58	MG	DA	1643	1/1	0.90	0.16	-	55,55,55,55	0
58	MG	AA	3228	1/1	0.96	0.19	-	51,51,51,51	0
58	MG	BA	1716	1/1	0.78	0.17	-	67,67,67,67	0
58	MG	AA	3154	1/1	0.91	0.35	-	46,46,46,46	0
58	MG	AA	3432	1/1	0.90	0.33	-	57,57,57,57	0
58	MG	CE	306	1/1	0.92	0.07	-	67,67,67,67	0
58	MG	AA	3605	1/1	0.86	0.28	-	68,68,68,68	0
58	MG	CB	3005	1/1	0.94	0.31	-	62,62,62,62	0
58	MG	CA	3397	1/1	0.98	0.13	-	59,59,59,59	0
58	MG	CN	5001	1/1	0.91	0.09	-	77,77,77,77	0
58	MG	CA	3154	1/1	0.80	0.31	-	72,72,72,72	0
58	MG	AA	3131	1/1	0.96	0.45	-	55,55,55,55	0
58	MG	AA	3097	1/1	0.98	0.20	-	26,26,26,26	0
58	MG	CA	3078	1/1	0.87	0.25	-	47,47,47,47	0
58	MG	BA	1715	1/1	0.74	0.21	-	83,83,83,83	0
58	MG	CA	3102	1/1	0.92	0.35	-	56,56,56,56	0
58	MG	CA	3557	1/1	0.89	0.11	-	76,76,76,76	0
58	MG	BA	1795	1/1	0.97	0.27	-	69,69,69,69	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3205	1/1	0.85	0.46	-	64,64,64,64	0
58	MG	CA	3377	1/1	0.97	0.19	-	80,80,80,80	0
58	MG	AA	3416	1/1	0.97	0.17	-	30,30,30,30	0
58	MG	BA	1637	1/1	0.88	0.23	-	66,66,66,66	0
58	MG	AA	3392	1/1	0.96	0.17	-	42,42,42,42	0
58	MG	CA	3472	1/1	0.93	0.60	-	72,72,72,72	0
58	MG	CA	3184	1/1	0.89	0.30	-	66,66,66,66	0
58	MG	CA	3081	1/1	0.82	0.16	-	68,68,68,68	0
58	MG	AA	3692	1/1	0.97	0.18	-	53,53,53,53	0
58	MG	DA	1704	1/1	0.96	0.07	-	69,69,69,69	0
58	MG	CB	3013	1/1	0.80	0.19	-	98,98,98,98	0
58	MG	CA	3053	1/1	0.96	0.75	-	58,58,58,58	0
58	MG	CA	3350	1/1	0.79	0.09	-	85,85,85,85	0
58	MG	AA	3448	1/1	0.89	0.05	-	78,78,78,78	0
58	MG	AA	3437	1/1	0.89	0.25	-	54,54,54,54	0
58	MG	AA	3424	1/1	0.86	0.15	-	65,65,65,65	0
58	MG	AA	3262	1/1	0.97	0.38	-	70,70,70,70	0
58	MG	BA	1650	1/1	0.71	0.36	-	72,72,72,72	0
58	MG	BA	1759	1/1	0.95	0.14	-	63,63,63,63	0
58	MG	AA	3638	1/1	0.71	0.41	-	72,72,72,72	0
58	MG	AA	3624	1/1	0.93	0.15	-	65,65,65,65	0
58	MG	AA	3608	1/1	0.86	0.18	-	73,73,73,73	0
58	MG	AA	3103	1/1	0.98	0.07	-	15,15,15,15	0
58	MG	BA	1761	1/1	0.93	0.18	-	55,55,55,55	0
58	MG	CA	3426	1/1	0.91	0.19	-	55,55,55,55	0
58	MG	BA	1808	1/1	0.93	0.14	-	50,50,50,50	0
58	MG	AA	3105	1/1	0.84	0.14	-	81,81,81,81	0
58	MG	DA	1602	1/1	0.92	0.12	-	80,80,80,80	0
58	MG	CA	3508	1/1	0.94	0.15	-	96,96,96,96	0
58	MG	CP	202	1/1	0.88	0.42	-	71,71,71,71	0
58	MG	AA	3728	1/1	0.92	0.21	-	61,61,61,61	0
58	MG	CA	3064	1/1	0.95	0.06	-	43,43,43,43	0
58	MG	CA	3578	1/1	0.85	0.26	-	80,80,80,80	0
58	MG	AA	3015	1/1	0.81	0.48	-	64,64,64,64	0
58	MG	DA	1611	1/1	0.96	0.09	-	38,38,38,38	0
58	MG	CA	3566	1/1	0.96	0.13	-	41,41,41,41	0
58	MG	AA	3156	1/1	0.95	0.34	-	33,33,33,33	1
58	MG	DA	1660	1/1	0.52	0.63	-	90,90,90,90	0
58	MG	AA	3085	1/1	0.84	0.32	-	53,53,53,53	0
58	MG	AA	3367	1/1	0.96	0.17	-	52,52,52,52	0
58	MG	AA	3274	1/1	0.92	0.29	-	55,55,55,55	0
58	MG	CA	3066	1/1	0.62	0.59	-	84,84,84,84	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3563	1/1	0.97	0.16	-	49,49,49,49	1
58	MG	DA	1705	1/1	0.92	0.32	-	62,62,62,62	0
58	MG	AA	3422	1/1	0.88	0.10	-	71,71,71,71	0
58	MG	CA	3387	1/1	0.95	0.34	-	70,70,70,70	0
58	MG	DJ	5001	1/1	0.53	0.48	-	105,105,105,105	0
58	MG	AA	3577	1/1	0.93	0.11	-	32,32,32,32	0
58	MG	AA	3332	1/1	0.95	0.14	-	46,46,46,46	0
58	MG	AA	3298	1/1	0.93	0.08	-	59,59,59,59	0
58	MG	AA	3724	1/1	0.97	0.28	-	40,40,40,40	0
58	MG	CA	3586	1/1	0.92	0.15	-	69,69,69,69	0
58	MG	AA	3343	1/1	0.95	0.14	-	65,65,65,65	0
58	MG	AA	3402	1/1	0.96	0.30	-	33,33,33,33	0
58	MG	CA	3520	1/1	0.92	0.18	-	59,59,59,59	0
58	MG	CA	3604	1/1	0.80	0.19	-	69,69,69,69	0
58	MG	AA	3365	1/1	0.91	0.31	-	57,57,57,57	0
58	MG	AA	3472	1/1	0.96	0.21	-	24,24,24,24	0
58	MG	CA	3366	1/1	0.96	0.24	-	61,61,61,61	0
58	MG	AA	3293	1/1	0.96	0.20	-	32,32,32,32	0
58	MG	AA	3004	1/1	0.92	0.16	-	25,25,25,25	0
58	MG	AA	3672	1/1	0.79	0.35	-	25,25,25,25	1
58	MG	CA	3307	1/1	0.97	0.30	-	52,52,52,52	0
58	MG	CA	3237	1/1	0.96	0.31	-	75,75,75,75	0
58	MG	AA	3052	1/1	0.86	0.64	-	65,65,65,65	0
58	MG	AA	3752	1/1	0.84	0.60	-	72,72,72,72	0
58	MG	AA	3352	1/1	0.89	0.25	-	51,51,51,51	0
58	MG	BA	1627	1/1	0.89	0.28	-	87,87,87,87	0
58	MG	CA	3059	1/1	0.77	0.37	-	60,60,60,60	0
58	MG	AA	3460	1/1	0.83	0.47	-	72,72,72,72	0
58	MG	CA	3257	1/1	0.94	0.48	-	57,57,57,57	0
58	MG	AA	3438	1/1	0.89	0.22	-	57,57,57,57	0
58	MG	AA	3229	1/1	0.89	0.31	-	43,43,43,43	0
58	MG	BA	1751	1/1	0.89	0.13	-	58,58,58,58	0
58	MG	DA	1620	1/1	0.81	0.14	-	58,58,58,58	0
58	MG	CA	3224	1/1	0.69	0.97	-	81,81,81,81	0
58	MG	CA	3219	1/1	0.98	0.21	-	31,31,31,31	0
58	MG	AA	3570	1/1	0.85	0.17	-	15,15,15,15	0
58	MG	CA	3589	1/1	0.85	0.07	-	79,79,79,79	0
58	MG	AA	3139	1/1	0.92	0.10	-	58,58,58,58	0
58	MG	DA	1673	1/1	0.65	0.37	-	100,100,100,100	0
58	MG	CA	3599	1/1	0.54	0.23	-	80,80,80,80	0
58	MG	AA	3804	1/1	0.93	0.40	-	68,68,68,68	0
58	MG	AA	3548	1/1	0.95	0.16	-	57,57,57,57	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3540	1/1	0.83	0.33	-	85,85,85,85	0
58	MG	CA	3371	1/1	0.95	0.21	-	55,55,55,55	0
58	MG	CA	3144	1/1	0.90	0.32	-	67,67,67,67	0
58	MG	AA	3080	1/1	0.80	0.40	-	57,57,57,57	0
58	MG	AA	3660	1/1	0.91	0.23	-	70,70,70,70	0
58	MG	AA	3323	1/1	0.86	0.12	-	22,22,22,22	0
58	MG	CA	3068	1/1	0.78	0.26	-	57,57,57,57	0
58	MG	CA	3060	1/1	0.91	0.41	-	77,77,77,77	0
58	MG	CA	3407	1/1	0.90	0.23	-	55,55,55,55	0
58	MG	AA	3419	1/1	0.96	0.14	-	31,31,31,31	0
58	MG	AA	3757	1/1	0.90	0.12	-	55,55,55,55	0
58	MG	CA	3205	1/1	0.09	0.69	-	105,105,105,105	0
58	MG	BA	1655	1/1	0.91	0.32	-	69,69,69,69	0
58	MG	AA	3557	1/1	0.97	0.17	-	19,19,19,19	0
58	MG	DK	5001	1/1	0.93	0.27	-	100,100,100,100	0
58	MG	CA	3386	1/1	0.90	0.28	-	65,65,65,65	0
58	MG	AA	3429	1/1	0.96	0.21	-	41,41,41,41	0
58	MG	CQ	205	1/1	0.55	0.51	-	81,81,81,81	0
58	MG	CA	3249	1/1	0.95	0.19	-	46,46,46,46	0
58	MG	CB	3012	1/1	0.95	0.35	-	76,76,76,76	0
58	MG	AA	3121	1/1	0.86	0.33	-	70,70,70,70	0
58	MG	BE	3001	1/1	0.94	0.12	-	60,60,60,60	0
58	MG	CQ	204	1/1	0.85	0.31	-	61,61,61,61	0
58	MG	CA	3435	1/1	0.97	0.11	-	52,52,52,52	0
58	MG	AB	3006	1/1	0.77	0.30	-	72,72,72,72	0
58	MG	CA	3197	1/1	0.90	0.48	-	64,64,64,64	0
58	MG	BA	1766	1/1	0.75	0.36	-	86,86,86,86	0
58	MG	AA	3658	1/1	0.99	0.15	-	62,62,62,62	0
58	MG	BA	1610	1/1	0.66	0.12	-	79,79,79,79	0
58	MG	CA	3239	1/1	0.72	0.39	-	75,75,75,75	0
58	MG	CA	3590	1/1	0.71	0.17	-	95,95,95,95	0
58	MG	CA	3642	1/1	0.83	0.99	-	80,80,80,80	0
58	MG	CA	3451	1/1	0.95	0.19	-	63,63,63,63	0
58	MG	CA	3269	1/1	0.94	0.13	-	86,86,86,86	0
58	MG	BA	1764	1/1	0.85	0.33	-	63,63,63,63	0
58	MG	BA	1638	1/1	0.79	0.63	-	78,78,78,78	0
58	MG	AA	3765	1/1	0.85	0.37	-	63,63,63,63	0
58	MG	AA	3713	1/1	0.79	0.41	-	52,52,52,52	1
58	MG	CA	3327	1/1	0.94	0.25	-	53,53,53,53	0
58	MG	AA	3727	1/1	0.88	0.15	-	49,49,49,49	0
58	MG	CA	3507	1/1	0.63	0.26	-	100,100,100,100	0
58	MG	BA	1692	1/1	0.82	0.30	-	86,86,86,86	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3632	1/1	0.85	0.17	-	79,79,79,79	0
58	MG	AA	3078	1/1	0.80	0.29	-	66,66,66,66	0
58	MG	CA	3129	1/1	0.85	0.54	-	69,69,69,69	0
58	MG	DA	1608	1/1	0.87	0.08	-	47,47,47,47	0
58	MG	BA	1631	1/1	0.89	0.11	-	48,48,48,48	0
58	MG	DA	1713	1/1	0.96	0.54	-	72,72,72,72	0
58	MG	DA	1749	1/1	0.85	0.43	-	80,80,80,80	0
58	MG	CA	3277	1/1	0.95	0.11	-	90,90,90,90	0
58	MG	BA	1792	1/1	0.95	0.19	-	75,75,75,75	0
58	MG	CA	3153	1/1	0.89	0.19	-	78,78,78,78	0
58	MG	AA	3193	1/1	0.85	0.23	-	62,62,62,62	0
58	MG	AA	3534	1/1	0.98	0.17	-	27,27,27,27	0
58	MG	AA	3286	1/1	0.94	0.21	-	52,52,52,52	0
58	MG	DA	1755	1/1	0.74	0.64	-	86,86,86,86	0
58	MG	CA	3098	1/1	0.60	0.39	-	83,83,83,83	0
58	MG	BA	1632	1/1	0.90	0.30	-	54,54,54,54	0
58	MG	DA	1625	1/1	0.92	0.55	-	73,73,73,73	0
58	MG	AA	3458	1/1	0.92	0.19	-	72,72,72,72	0
58	MG	BA	1663	1/1	0.94	0.09	-	79,79,79,79	0
58	MG	CA	3090	1/1	0.87	0.53	-	65,65,65,65	0
58	MG	AA	3536	1/1	0.94	0.10	-	35,35,35,35	0
58	MG	BA	1661	1/1	0.73	0.99	-	82,82,82,82	0
58	MG	CA	3280	1/1	0.90	0.18	-	30,30,30,30	0
58	MG	CA	3111	1/1	0.82	0.13	-	71,71,71,71	0
58	MG	DA	1766	1/1	0.85	0.12	-	58,58,58,58	0
58	MG	CA	3359	1/1	0.93	0.19	-	42,42,42,42	0
58	MG	AA	3093	1/1	0.91	0.28	-	27,27,27,27	1
58	MG	CA	3245	1/1	0.89	0.48	-	57,57,57,57	0
58	MG	DA	1639	1/1	0.87	0.16	-	75,75,75,75	0
58	MG	CA	3298	1/1	0.94	0.15	-	68,68,68,68	0
58	MG	AA	3655	1/1	0.98	0.16	-	61,61,61,61	0
58	MG	AA	3163	1/1	0.88	0.38	-	40,40,40,40	0
58	MG	AA	3503	1/1	0.94	0.15	-	64,64,64,64	0
58	MG	CA	3534	1/1	0.84	0.18	-	79,79,79,79	0
58	MG	CA	3593	1/1	0.68	0.26	-	73,73,73,73	0
58	MG	AA	3661	1/1	0.94	0.32	-	43,43,43,43	0
58	MG	AA	3588	1/1	0.90	0.19	-	38,38,38,38	0
58	MG	AA	3292	1/1	0.90	0.13	-	74,74,74,74	0
58	MG	AA	3029	1/1	0.87	0.29	-	53,53,53,53	0
58	MG	CA	3511	1/1	0.96	0.12	-	68,68,68,68	0
58	MG	AA	3666	1/1	0.95	0.07	-	64,64,64,64	0
58	MG	CA	3089	1/1	0.72	0.41	-	80,80,80,80	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3092	1/1	0.94	0.13	-	53,53,53,53	0
58	MG	AA	3635	1/1	0.99	0.11	-	23,23,23,23	0
58	MG	CA	3517	1/1	0.99	0.33	-	64,64,64,64	0
58	MG	DA	1729	1/1	0.88	0.13	-	57,57,57,57	0
58	MG	CA	3008	1/1	0.72	0.41	-	100,100,100,100	0
58	MG	CA	3638	1/1	0.91	0.32	-	55,55,55,55	0
58	MG	BA	1641	1/1	0.92	0.23	-	71,71,71,71	0
58	MG	CA	3527	1/1	0.47	0.13	-	78,78,78,78	0
58	MG	AA	3268	1/1	0.80	0.12	-	88,88,88,88	0
58	MG	CA	3641	1/1	0.79	0.40	-	67,67,67,67	0
58	MG	AF	304	1/1	0.92	0.32	-	62,62,62,62	0
58	MG	AA	3244	1/1	0.45	0.25	-	100,100,100,100	0
58	MG	CA	3524	1/1	0.79	0.09	-	77,77,77,77	0
58	MG	AA	3224	1/1	0.83	0.41	-	75,75,75,75	0
58	MG	BA	1765	1/1	0.97	0.11	-	61,61,61,61	0
58	MG	CA	3554	1/1	0.95	0.18	-	67,67,67,67	0
58	MG	AA	3595	1/1	0.96	0.25	-	55,55,55,55	0
58	MG	AA	3560	1/1	0.87	0.27	-	58,58,58,58	0
58	MG	BA	1787	1/1	0.88	0.24	-	90,90,90,90	0
58	MG	AA	3483	1/1	0.85	0.07	-	43,43,43,43	1
58	MG	DA	1737	1/1	0.91	0.16	-	72,72,72,72	0
58	MG	CA	3094	1/1	0.92	0.32	-	87,87,87,87	0
58	MG	DA	1629	1/1	0.93	0.41	-	58,58,58,58	0
58	MG	CP	201	1/1	0.85	0.82	-	65,65,65,65	0
58	MG	CA	3025	1/1	0.83	0.28	-	77,77,77,77	0
58	MG	DA	1667	1/1	0.96	0.24	-	49,49,49,49	0
58	MG	AA	3140	1/1	0.93	0.31	-	50,50,50,50	0
58	MG	CA	3478	1/1	0.94	0.14	-	58,58,58,58	0
58	MG	AA	3500	1/1	0.95	0.12	-	47,47,47,47	0
58	MG	AA	3055	1/1	0.95	0.28	-	35,35,35,35	0
58	MG	CA	3045	1/1	0.88	0.41	-	67,67,67,67	0
58	MG	AN	3003	1/1	0.91	0.13	-	47,47,47,47	0
58	MG	CA	3093	1/1	0.58	0.64	-	84,84,84,84	0
58	MG	AA	3781	1/1	0.76	0.33	-	44,44,44,44	1
58	MG	CA	3156	1/1	0.93	0.32	-	68,68,68,68	0
58	MG	BA	1777	1/1	0.90	0.21	-	79,79,79,79	0
58	MG	DA	1706	1/1	0.72	0.28	-	128,128,128,128	0
58	MG	AA	3281	1/1	0.84	0.40	-	61,61,61,61	0
58	MG	AA	3625	1/1	0.88	0.21	-	60,60,60,60	0
58	MG	DA	1627	1/1	0.80	0.23	-	70,70,70,70	0
58	MG	A4	502	1/1	0.63	0.65	-	123,123,123,123	0
58	MG	AA	3263	1/1	0.88	0.80	-	80,80,80,80	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	BL	202	1/1	0.94	0.35	-	67,67,67,67	0
58	MG	DA	1741	1/1	0.93	0.16	-	78,78,78,78	0
58	MG	BA	1769	1/1	0.97	0.28	-	63,63,63,63	0
58	MG	CA	3494	1/1	0.94	0.21	-	63,63,63,63	0
58	MG	AB	3010	1/1	0.92	0.18	-	47,47,47,47	1
58	MG	CA	3122	1/1	0.91	0.22	-	67,67,67,67	0
58	MG	CA	3516	1/1	0.96	0.12	-	62,62,62,62	0
58	MG	AA	3226	1/1	0.92	0.27	-	56,56,56,56	0
58	MG	CA	3587	1/1	0.99	0.12	-	34,34,34,34	0
58	MG	AA	3722	1/1	0.94	0.14	-	18,18,18,18	0
58	MG	CA	3026	1/1	0.97	0.25	-	81,81,81,81	0
58	MG	AA	3674	1/1	0.94	0.25	-	75,75,75,75	0
58	MG	BA	1712	1/1	0.93	0.14	-	61,61,61,61	0
58	MG	AA	3032	1/1	0.95	0.27	-	36,36,36,36	0
58	MG	AA	3662	1/1	0.89	0.22	-	59,59,59,59	0
58	MG	CA	3616	1/1	0.66	0.66	-	79,79,79,79	0
58	MG	CA	3254	1/1	0.96	0.21	-	42,42,42,42	0
58	MG	AA	3083	1/1	0.94	0.24	-	38,38,38,38	1
58	MG	AA	3243	1/1	0.98	0.24	-	24,24,24,24	1
58	MG	CA	3611	1/1	0.74	0.70	-	91,91,91,91	0
58	MG	AA	3276	1/1	0.99	0.28	-	47,47,47,47	1
58	MG	CA	3446	1/1	0.95	0.18	-	63,63,63,63	0
58	MG	DA	1732	1/1	0.89	0.36	-	76,76,76,76	0
58	MG	A2	101	1/1	0.93	0.20	-	35,35,35,35	0
58	MG	AA	3270	1/1	0.77	0.24	-	54,54,54,54	0
58	MG	AA	3053	1/1	0.97	0.16	-	14,14,14,14	0
58	MG	DA	1726	1/1	0.93	0.17	-	77,77,77,77	0
58	MG	AA	3296	1/1	0.95	0.12	-	17,17,17,17	0
58	MG	AA	3252	1/1	0.73	0.55	-	66,66,66,66	0
58	MG	AA	3255	1/1	0.89	0.35	-	53,53,53,53	0
58	MG	AA	3412	1/1	0.96	0.21	-	43,43,43,43	0
58	MG	AE	302	1/1	0.99	0.23	-	18,18,18,18	0
58	MG	AB	3022	1/1	0.98	0.05	-	58,58,58,58	0
58	MG	BA	1604	1/1	0.92	0.15	-	65,65,65,65	0
58	MG	AA	3144	1/1	0.97	0.12	-	40,40,40,40	0
58	MG	CA	3033	1/1	0.65	0.87	-	89,89,89,89	0
58	MG	AA	3147	1/1	0.88	0.27	-	69,69,69,69	0
58	MG	CA	3150	1/1	0.91	0.18	-	54,54,54,54	0
58	MG	BA	1645	1/1	0.77	0.64	-	61,61,61,61	0
58	MG	CA	3459	1/1	0.97	0.18	-	48,48,48,48	0
58	MG	CA	3461	1/1	0.94	0.18	-	43,43,43,43	0
58	MG	A7	102	1/1	0.95	0.10	-	43,43,43,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3632	1/1	0.93	0.14	-	76,76,76,76	0
58	MG	AB	3011	1/1	0.96	0.19	-	30,30,30,30	0
58	MG	CA	3389	1/1	0.69	0.49	-	75,75,75,75	0
58	MG	CA	3381	1/1	0.94	0.10	-	38,38,38,38	0
58	MG	AA	3191	1/1	0.94	0.26	-	42,42,42,42	0
58	MG	BA	1778	1/1	0.98	0.07	-	42,42,42,42	0
58	MG	CA	3395	1/1	0.96	0.33	-	65,65,65,65	0
58	MG	DA	1632	1/1	0.81	0.32	-	61,61,61,61	0
58	MG	AA	3482	1/1	0.98	0.15	-	65,65,65,65	0
58	MG	AA	3158	1/1	0.90	0.31	-	97,97,97,97	0
58	MG	CA	3162	1/1	0.97	0.47	-	46,46,46,46	0
58	MG	AA	3049	1/1	0.94	0.27	-	52,52,52,52	0
58	MG	CA	3503	1/1	0.83	0.17	-	62,62,62,62	0
58	MG	AA	3785	1/1	0.92	0.19	-	72,72,72,72	0
58	MG	CA	3613	1/1	0.92	0.34	-	74,74,74,74	0
58	MG	AA	3107	1/1	0.91	0.51	-	76,76,76,76	0
58	MG	CA	3279	1/1	0.94	0.16	-	34,34,34,34	0
58	MG	AA	3631	1/1	0.94	0.30	-	46,46,46,46	0
58	MG	CA	3580	1/1	0.88	0.07	-	100,100,100,100	0
58	MG	CA	3544	1/1	0.60	0.19	-	81,81,81,81	0
58	MG	CO	202	1/1	0.93	0.21	-	53,53,53,53	0
58	MG	BA	1742	1/1	0.67	0.21	-	79,79,79,79	0
58	MG	AA	3152	1/1	0.93	0.29	-	71,71,71,71	0
58	MG	CA	3470	1/1	0.93	0.35	-	72,72,72,72	0
58	MG	AA	3427	1/1	0.96	0.11	-	33,33,33,33	0
58	MG	CA	3304	1/1	0.76	0.13	-	93,93,93,93	0
58	MG	AA	3345	1/1	0.89	0.10	-	68,68,68,68	0
58	MG	CA	3097	1/1	0.82	0.26	-	80,80,80,80	0
58	MG	AA	3411	1/1	0.80	0.23	-	47,47,47,47	0
58	MG	AA	3504	1/1	0.94	0.20	-	58,58,58,58	0
58	MG	CA	3063	1/1	0.84	0.29	-	53,53,53,53	0
58	MG	CA	3209	1/1	0.20	0.65	-	93,93,93,93	0
58	MG	AA	3455	1/1	0.91	0.33	-	58,58,58,58	0
58	MG	AA	3673	1/1	0.89	0.17	-	67,67,67,67	0
58	MG	AA	3520	1/1	0.96	0.13	-	38,38,38,38	0
58	MG	CA	3384	1/1	0.87	0.23	-	71,71,71,71	0
58	MG	AA	3731	1/1	0.91	0.19	-	42,42,42,42	0
58	MG	AA	3814	1/1	0.76	0.37	-	93,93,93,93	0
58	MG	BN	502	1/1	0.95	0.24	-	66,66,66,66	0
58	MG	CA	3606	1/1	0.91	0.42	-	65,65,65,65	0
58	MG	AA	3351	1/1	0.97	0.14	-	30,30,30,30	0
58	MG	AA	3335	1/1	0.91	0.23	-	41,41,41,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3677	1/1	0.92	0.22	-	41,41,41,41	0
58	MG	AA	3709	1/1	0.98	0.19	-	29,29,29,29	1
58	MG	AA	3769	1/1	0.93	0.23	-	57,57,57,57	0
58	MG	CA	3648	1/1	0.95	0.33	-	53,53,53,53	0
58	MG	AA	3026	1/1	0.82	0.39	-	86,86,86,86	0
58	MG	CA	3569	1/1	0.94	0.26	-	79,79,79,79	0
58	MG	AA	3248	1/1	0.73	0.52	-	64,64,64,64	0
58	MG	CA	3522	1/1	0.90	0.34	-	56,56,56,56	0
58	MG	CA	3399	1/1	0.93	0.10	-	75,75,75,75	0
58	MG	AA	3592	1/1	0.90	0.25	-	52,52,52,52	0
58	MG	CA	3252	1/1	0.94	0.23	-	64,64,64,64	0
58	MG	BA	1793	1/1	0.80	0.66	-	86,86,86,86	0
58	MG	BA	1666	1/1	0.77	0.53	-	75,75,75,75	0
58	MG	DA	1723	1/1	0.97	0.30	-	66,66,66,66	0
58	MG	CA	3475	1/1	0.91	0.36	-	55,55,55,55	0
58	MG	AE	301	1/1	0.79	0.28	-	69,69,69,69	0
58	MG	AA	3025	1/1	0.93	0.39	-	35,35,35,35	1
58	MG	DA	1699	1/1	0.99	0.08	-	75,75,75,75	0
58	MG	DA	1664	1/1	0.94	0.15	-	64,64,64,64	0
58	MG	AE	303	1/1	0.97	0.25	-	41,41,41,41	0
58	MG	AA	3215	1/1	0.89	0.62	-	42,42,42,42	1
58	MG	A9	502	1/1	0.88	0.28	-	60,60,60,60	0
58	MG	CA	3624	1/1	0.53	0.17	-	104,104,104,104	0
58	MG	AA	3637	1/1	0.89	0.29	-	45,45,45,45	0
58	MG	A5	102	1/1	0.89	0.34	-	60,60,60,60	0
58	MG	AA	3671	1/1	0.93	0.22	-	58,58,58,58	0
58	MG	CA	3253	1/1	0.70	0.18	-	95,95,95,95	0
58	MG	DA	1641	1/1	0.92	0.09	-	77,77,77,77	0
58	MG	CA	3126	1/1	0.79	0.28	-	93,93,93,93	0
58	MG	AA	3089	1/1	0.90	0.31	-	47,47,47,47	1
58	MG	AA	3006	1/1	0.93	0.46	-	52,52,52,52	0
58	MG	AA	3569	1/1	0.96	0.17	-	19,19,19,19	0
58	MG	CA	3449	1/1	0.95	0.10	-	66,66,66,66	0
58	MG	BA	1703	1/1	0.93	0.25	-	78,78,78,78	0
58	MG	CA	3512	1/1	0.81	0.48	-	65,65,65,65	0
58	MG	DA	1688	1/1	0.97	0.20	-	66,66,66,66	0
58	MG	CA	3333	1/1	0.94	0.32	-	75,75,75,75	0
58	MG	AA	3550	1/1	0.94	0.21	-	47,47,47,47	0
58	MG	CA	3404	1/1	0.94	0.19	-	54,54,54,54	0
58	MG	AW	3001	1/1	0.90	0.31	-	52,52,52,52	0
58	MG	DA	1612	1/1	0.82	0.39	-	72,72,72,72	0
58	MG	CA	3533	1/1	0.79	0.20	-	81,81,81,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3301	1/1	0.98	0.35	-	47,47,47,47	0
58	MG	AA	3541	1/1	0.98	0.18	-	43,43,43,43	0
58	MG	BA	1608	1/1	0.89	0.50	-	60,60,60,60	0
58	MG	CA	3393	1/1	0.68	0.08	-	82,82,82,82	0
58	MG	CA	3017	1/1	0.93	0.58	-	46,46,46,46	0
58	MG	DA	1650	1/1	0.91	0.34	-	61,61,61,61	0
58	MG	CA	3158	1/1	0.70	0.32	-	70,70,70,70	0
58	MG	CA	3614	1/1	0.84	0.20	-	52,52,52,52	0
58	MG	CA	3259	1/1	0.96	0.22	-	47,47,47,47	0
58	MG	BA	1653	1/1	0.85	0.43	-	78,78,78,78	0
58	MG	AA	3797	1/1	0.96	0.26	-	15,15,15,15	1
58	MG	CA	3635	1/1	0.85	0.16	-	79,79,79,79	0
58	MG	CA	3222	1/1	0.96	0.26	-	75,75,75,75	0
58	MG	AA	3712	1/1	0.91	0.23	-	46,46,46,46	0
58	MG	A8	5002	1/1	0.97	0.24	-	31,31,31,31	0
58	MG	AA	3795	1/1	0.90	0.33	-	68,68,68,68	1
58	MG	AA	3002	1/1	0.86	0.21	-	55,55,55,55	0
58	MG	AA	3762	1/1	0.90	0.19	-	53,53,53,53	1
58	MG	AA	3271	1/1	0.93	0.36	-	69,69,69,69	0
58	MG	BA	1718	1/1	0.86	0.52	-	83,83,83,83	0
58	MG	AA	3743	1/1	0.88	0.25	-	80,80,80,80	0
58	MG	CA	3443	1/1	0.97	0.11	-	36,36,36,36	0
58	MG	CA	3272	1/1	0.84	0.48	-	75,75,75,75	0
58	MG	BA	1744	1/1	0.88	0.10	-	37,37,37,37	0
58	MG	CA	3367	1/1	0.95	0.23	-	65,65,65,65	0
58	MG	AA	3523	1/1	0.97	0.20	-	30,30,30,30	0
58	MG	BA	1668	1/1	0.86	0.25	-	83,83,83,83	0
58	MG	CA	3264	1/1	0.98	0.18	-	59,59,59,59	0
58	MG	DA	1609	1/1	0.94	0.30	-	46,46,46,46	0
58	MG	AA	3719	1/1	0.94	0.12	-	58,58,58,58	0
58	MG	AA	3651	1/1	0.84	0.24	-	52,52,52,52	0
58	MG	CA	3482	1/1	0.90	0.24	-	70,70,70,70	0
58	MG	AA	3414	1/1	0.82	0.16	-	36,36,36,36	0
58	MG	AA	3175	1/1	0.90	0.33	-	51,51,51,51	0
58	MG	BA	1647	1/1	0.83	0.15	-	75,75,75,75	0
58	MG	AA	3593	1/1	0.94	0.21	-	25,25,25,25	1
58	MG	CY	502	1/1	0.92	0.17	-	56,56,56,56	0
58	MG	AA	3126	1/1	0.88	0.42	-	50,50,50,50	0
58	MG	BA	1687	1/1	0.88	0.33	-	72,72,72,72	0
58	MG	BA	1677	1/1	0.86	0.16	-	87,87,87,87	0
58	MG	AA	3707	1/1	0.98	0.24	-	29,29,29,29	1
58	MG	CA	3405	1/1	0.84	0.23	-	91,91,91,91	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3787	1/1	0.97	0.25	-	53,53,53,53	0
58	MG	CA	3403	1/1	0.93	0.08	-	91,91,91,91	0
58	MG	CA	3295	1/1	0.73	0.20	-	84,84,84,84	0
58	MG	AA	3501	1/1	0.95	0.11	-	24,24,24,24	0
58	MG	AQ	201	1/1	0.77	0.51	-	62,62,62,62	0
58	MG	BA	1768	1/1	0.94	0.07	-	75,75,75,75	0
58	MG	AA	3590	1/1	0.92	0.21	-	69,69,69,69	0
58	MG	BA	1719	1/1	0.83	0.27	-	80,80,80,80	0
58	MG	BA	1656	1/1	0.86	0.13	-	90,90,90,90	0
58	MG	BD	502	1/1	0.83	0.63	-	82,82,82,82	0
58	MG	AA	3325	1/1	0.97	0.10	-	66,66,66,66	0
58	MG	DA	1724	1/1	0.91	0.31	-	61,61,61,61	0
58	MG	CA	3411	1/1	0.94	0.36	-	61,61,61,61	0
58	MG	CA	3506	1/1	0.89	0.17	-	58,58,58,58	0
58	MG	CA	3128	1/1	0.90	0.42	-	71,71,71,71	0
58	MG	CA	3342	1/1	0.98	0.17	-	45,45,45,45	0
58	MG	AA	3295	1/1	0.89	0.39	-	47,47,47,47	0
58	MG	BZ	702	1/1	0.95	0.34	-	46,46,46,46	0
58	MG	CA	3325	1/1	0.88	0.10	-	38,38,38,38	0
58	MG	CA	3165	1/1	0.92	0.17	-	62,62,62,62	0
58	MG	CA	3100	1/1	0.84	0.46	-	79,79,79,79	0
58	MG	CA	3285	1/1	0.95	0.37	-	57,57,57,57	0
58	MG	BA	1802	1/1	0.86	0.14	-	68,68,68,68	1
58	MG	DA	1613	1/1	0.91	0.25	-	72,72,72,72	0
58	MG	AA	3600	1/1	0.81	0.25	-	60,60,60,60	0
58	MG	AA	3111	1/1	0.89	0.43	-	48,48,48,48	0
58	MG	DA	1707	1/1	0.71	0.30	-	87,87,87,87	0
58	MG	CA	3258	1/1	0.86	0.36	-	70,70,70,70	0
58	MG	AA	3100	1/1	0.87	0.27	-	53,53,53,53	0
58	MG	CA	3546	1/1	0.63	0.11	-	119,119,119,119	0
58	MG	AA	3222	1/1	0.90	0.32	-	28,28,28,28	0
58	MG	AA	3079	1/1	0.92	0.10	-	34,34,34,34	0
58	MG	CA	3436	1/1	0.95	0.12	-	75,75,75,75	0
58	MG	CA	3485	1/1	0.76	0.23	-	83,83,83,83	0
58	MG	AA	3304	1/1	0.88	0.25	-	30,30,30,30	0
58	MG	AA	3019	1/1	0.94	0.30	-	58,58,58,58	0
58	MG	A0	101	1/1	0.98	0.14	-	43,43,43,43	0
58	MG	AA	3239	1/1	0.86	0.36	-	64,64,64,64	0
58	MG	CA	3568	1/1	0.89	0.20	-	41,41,41,41	0
58	MG	CA	3083	1/1	0.66	0.75	-	90,90,90,90	0
58	MG	DA	1623	1/1	0.55	0.13	-	117,117,117,117	0
58	MG	AA	3284	1/1	0.89	0.42	-	44,44,44,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	CA	3385	1/1	0.96	0.47	-	64,64,64,64	0
58	MG	AA	3641	1/1	0.98	0.22	-	41,41,41,41	0
58	MG	BA	1732	1/1	0.86	0.07	-	71,71,71,71	0
58	MG	DA	1631	1/1	0.95	0.20	-	70,70,70,70	0
58	MG	CA	3024	1/1	0.83	0.68	-	88,88,88,88	0
58	MG	AA	3260	1/1	0.75	0.39	-	71,71,71,71	0
58	MG	CA	3513	1/1	0.90	0.26	-	75,75,75,75	0
58	MG	AA	3056	1/1	0.92	0.26	-	63,63,63,63	0
58	MG	AA	3544	1/1	0.97	0.11	-	16,16,16,16	0
58	MG	BA	1621	1/1	0.93	0.41	-	49,49,49,49	0
58	MG	CB	3011	1/1	0.94	0.29	-	53,53,53,53	0
58	MG	CA	3431	1/1	0.95	0.29	-	100,100,100,100	0
58	MG	DA	1603	1/1	0.86	0.30	-	74,74,74,74	0
58	MG	BA	1705	1/1	0.48	0.21	-	92,92,92,92	0
58	MG	DA	1614	1/1	0.79	0.80	-	87,87,87,87	0
58	MG	CA	3390	1/1	0.95	0.14	-	64,64,64,64	0
58	MG	BA	1753	1/1	0.95	0.13	-	94,94,94,94	0
58	MG	BA	1717	1/1	0.96	0.18	-	44,44,44,44	0
58	MG	BA	1750	1/1	0.97	0.27	-	65,65,65,65	0
58	MG	CA	3378	1/1	0.81	0.12	-	84,84,84,84	0
58	MG	CA	3135	1/1	0.96	0.21	-	66,66,66,66	0
58	MG	AA	3445	1/1	0.65	0.23	-	75,75,75,75	0
58	MG	CA	3493	1/1	0.81	0.39	-	88,88,88,88	0
58	MG	AA	3694	1/1	0.89	0.15	-	45,45,45,45	0
58	MG	CA	3140	1/1	0.71	0.60	-	98,98,98,98	0
58	MG	CA	3202	1/1	0.13	0.83	-	77,77,77,77	0
58	MG	BW	501	1/1	0.93	0.27	-	48,48,48,48	0
58	MG	AA	3537	1/1	0.78	0.15	-	95,95,95,95	0
58	MG	DA	1746	1/1	0.81	0.19	-	91,91,91,91	0
58	MG	AA	3626	1/1	0.81	0.27	-	74,74,74,74	0
58	MG	A6	101	1/1	0.93	0.37	-	65,65,65,65	0
58	MG	CA	3646	1/1	0.62	0.20	-	95,95,95,95	0
58	MG	AA	3664	1/1	0.97	0.23	-	57,57,57,57	0
58	MG	CA	3559	1/1	0.98	0.26	-	52,52,52,52	1
58	MG	AA	3755	1/1	0.83	0.42	-	63,63,63,63	0
58	MG	AA	3136	1/1	0.72	0.20	-	52,52,52,52	0
58	MG	BA	1710	1/1	0.88	0.69	-	73,73,73,73	0
58	MG	CA	3283	1/1	0.89	0.19	-	60,60,60,60	0
58	MG	CA	3183	1/1	0.76	1.07	-	86,86,86,86	0
58	MG	DA	1717	1/1	0.94	0.30	-	52,52,52,52	0
58	MG	CA	3216	1/1	0.81	0.62	-	66,66,66,66	0
58	MG	BA	1782	1/1	0.91	0.19	-	50,50,50,50	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
58	MG	AA	3090	1/1	0.93	0.56	-	30,30,30,30	1
58	MG	BA	1696	1/1	0.72	0.40	-	68,68,68,68	0
58	MG	CF	304	1/1	0.98	0.36	-	65,65,65,65	0
58	MG	CA	3577	1/1	0.92	0.17	-	83,83,83,83	0
58	MG	BA	1688	1/1	0.89	0.71	-	70,70,70,70	0
58	MG	CA	3329	1/1	0.97	0.21	-	29,29,29,29	0
58	MG	AA	3209	1/1	0.93	0.33	-	63,63,63,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.