



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 20, 2017 – 06:27 AM EDT

PDB ID : 3CMU
Title : Mechanism of homologous recombination from the RecA-ssDNA/dsDNA structures
Authors : Pavletich, N.P.
Deposited on : unknown
Resolution : 4.20 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-20029824
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-20029824

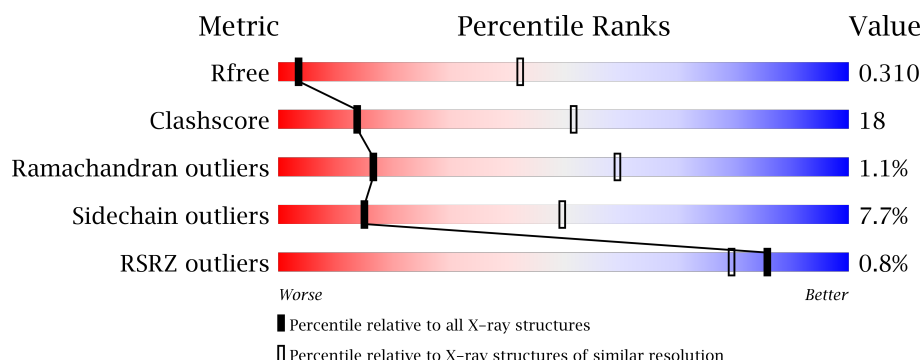
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.20 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	1177 (4.80-3.60)
Clashscore	112137	1025 (4.72-3.66)
Ramachandran outliers	110173	1024 (4.76-3.62)
Sidechain outliers	110143	1008 (4.76-3.62)
RSRZ outliers	101464	1188 (4.80-3.60)

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	B	18	
2	A	2050	

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
4	ALF	A	3501	-	-	-	X
4	ALF	A	501	-	-	X	-
4	ALF	A	5501	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15093 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 DNA chain called DNA (5'-D(*DTP*DTP*DTP*DTP*DTP*DTP*DTP*DT-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	15	Total	C	N	O	P	0	0	0
			297	150	30	103	14			

- Molecule 2 is a protein called Protein recA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	1937	Total	C	N	O	S	0	0	0
			14598	9178	2528	2824	68			

There are 75 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	GLY	-	LINKER	UNP P0A7G6
A	27	ALA	-	LINKER	UNP P0A7G6
A	28	HIS	-	LINKER	UNP P0A7G6
A	29	MET	-	LINKER	UNP P0A7G6
A	986	THR	-	LINKER	UNP P0A7G6
A	987	GLY	-	LINKER	UNP P0A7G6
A	988	SER	-	LINKER	UNP P0A7G6
A	989	THR	-	LINKER	UNP P0A7G6
A	990	GLY	-	LINKER	UNP P0A7G6
A	991	SER	-	LINKER	UNP P0A7G6
A	992	GLY	-	LINKER	UNP P0A7G6
A	993	THR	-	LINKER	UNP P0A7G6
A	994	THR	-	LINKER	UNP P0A7G6
A	995	GLY	-	LINKER	UNP P0A7G6
A	996	SER	-	LINKER	UNP P0A7G6
A	997	THR	-	LINKER	UNP P0A7G6
A	998	GLY	-	LINKER	UNP P0A7G6
A	999	SER	-	LINKER	UNP P0A7G6
A	1000	MET	-	LINKER	UNP P0A7G6

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	1986	THR	-	LINKER	UNP P0A7G6
A	1987	GLY	-	LINKER	UNP P0A7G6
A	1988	SER	-	LINKER	UNP P0A7G6
A	1989	THR	-	LINKER	UNP P0A7G6
A	1990	GLY	-	LINKER	UNP P0A7G6
A	1991	SER	-	LINKER	UNP P0A7G6
A	1992	MET	-	LINKER	UNP P0A7G6
A	1993	GLY	-	LINKER	UNP P0A7G6
A	1994	HIS	-	LINKER	UNP P0A7G6
A	1995	THR	-	LINKER	UNP P0A7G6
A	1996	THR	-	LINKER	UNP P0A7G6
A	1997	GLY	-	LINKER	UNP P0A7G6
A	1998	ALA	-	LINKER	UNP P0A7G6
A	1999	MET	-	LINKER	UNP P0A7G6
A	2000	SER	-	LINKER	UNP P0A7G6
A	2986	THR	-	LINKER	UNP P0A7G6
A	2987	GLY	-	LINKER	UNP P0A7G6
A	2988	SER	-	LINKER	UNP P0A7G6
A	2989	THR	-	LINKER	UNP P0A7G6
A	2990	GLY	-	LINKER	UNP P0A7G6
A	2991	SER	-	LINKER	UNP P0A7G6
A	2992	MET	-	LINKER	UNP P0A7G6
A	2993	ALA	-	LINKER	UNP P0A7G6
A	2994	SER	-	LINKER	UNP P0A7G6
A	2995	THR	-	LINKER	UNP P0A7G6
A	2996	GLY	-	LINKER	UNP P0A7G6
A	2997	SER	-	LINKER	UNP P0A7G6
A	2998	THR	-	LINKER	UNP P0A7G6
A	2999	GLY	-	LINKER	UNP P0A7G6
A	3000	SER	-	LINKER	UNP P0A7G6
A	3989	THR	-	LINKER	UNP P0A7G6
A	3990	GLY	-	LINKER	UNP P0A7G6
A	3991	ALA	-	LINKER	UNP P0A7G6
A	3992	THR	-	LINKER	UNP P0A7G6
A	3993	GLY	-	LINKER	UNP P0A7G6
A	3994	ALA	-	LINKER	UNP P0A7G6
A	3995	MET	-	LINKER	UNP P0A7G6
A	3996	SER	-	LINKER	UNP P0A7G6
A	3997	GLY	-	LINKER	UNP P0A7G6
A	3998	ARG	-	LINKER	UNP P0A7G6
A	3999	MET	-	LINKER	UNP P0A7G6
A	4000	SER	-	LINKER	UNP P0A7G6

Continued on next page...

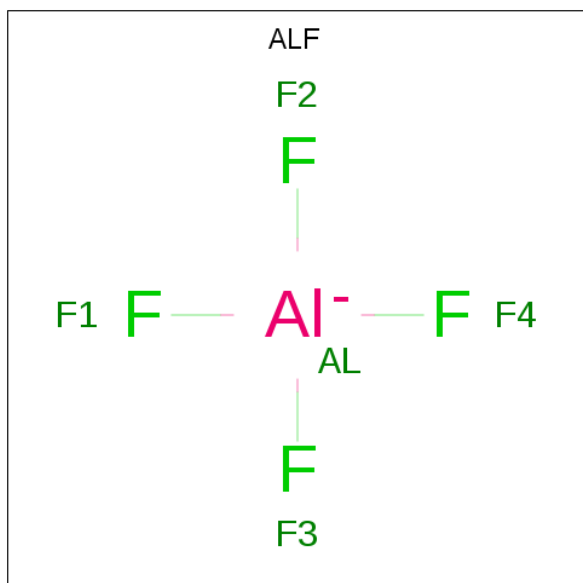
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	4987	THR	-	LINKER	UNP P0A7G6
A	4988	GLY	-	LINKER	UNP P0A7G6
A	4989	SER	-	LINKER	UNP P0A7G6
A	4990	THR	-	LINKER	UNP P0A7G6
A	4991	GLY	-	LINKER	UNP P0A7G6
A	4992	SER	-	LINKER	UNP P0A7G6
A	4993	GLY	-	LINKER	UNP P0A7G6
A	4994	SER	-	LINKER	UNP P0A7G6
A	4995	SER	-	LINKER	UNP P0A7G6
A	4996	THR	-	LINKER	UNP P0A7G6
A	4997	GLY	-	LINKER	UNP P0A7G6
A	4998	SER	-	LINKER	UNP P0A7G6
A	4999	MET	-	LINKER	UNP P0A7G6
A	5000	SER	-	LINKER	UNP P0A7G6

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	6	Total Mg 6 6	0	0

- Molecule 4 is TETRAFLUOROALUMINATE ION (three-letter code: ALF) (formula: AlF_4).



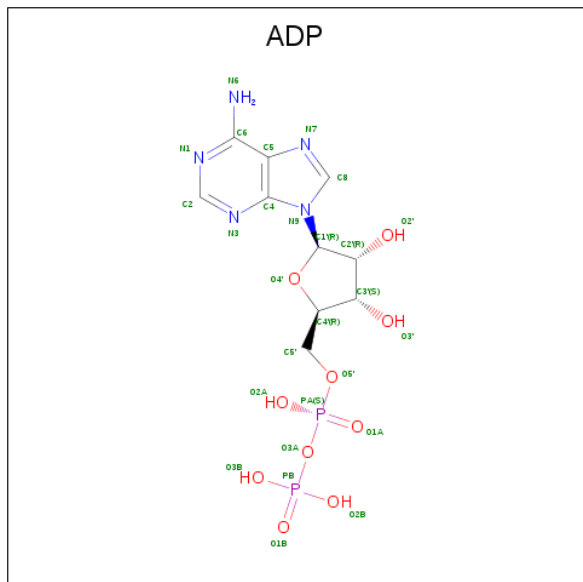
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Al F 5 1 4	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	Al	F	0	0
			5	1	4		
4	A	1	Total	Al	F	0	0
			5	1	4		
4	A	1	Total	Al	F	0	0
			5	1	4		
4	A	1	Total	Al	F	0	0
			5	1	4		
4	A	1	Total	Al	F	0	0
			5	1	4		

- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



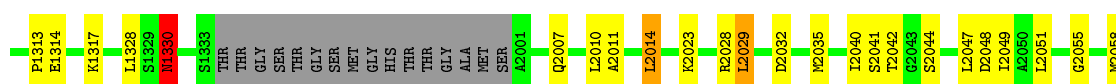
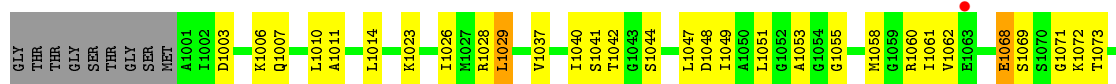
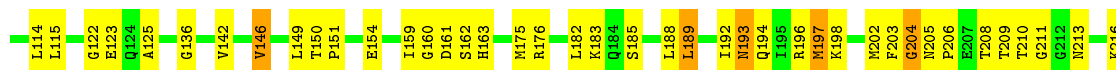
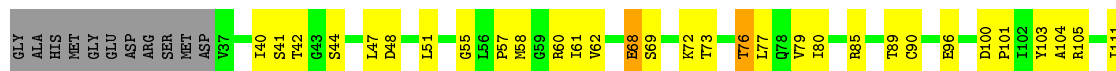
3 Residue-property plots

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 ($\text{RSRZ} > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- [illegible]



- Molecule 2: Protein recA



S5220	V5146	D5032	E4273	M4175	S4070	ALA	G3211	H3113	I3026	F2260	G2160
V5221	L5149	M5035	L4274	R4176	G4071	THR	G3212	L3114	I3027	Y2264	A2168
L5223	P5150	T5042	V4275	K4177	T4072	GLY	N3213	L3115	R3028	G2265	R2169
I5225	S5041	I5040	G4278	L4182	T4076	MET	L3215	G3123	I3029	I2268	Q2173
R5226	T5042	T5042	L4284	Q4184	L4077	GLY	R3216	E3122	G3030	G2271	A2174
E5235	E5045	G5045	E4285	S4185	Q4078	ARG	R3217	Q3124	E3031	G2272	M2175
E5235	S5044	S5044	E4285	MET	V4079	MET	R3218	A3125	D3032	L2274	R2176
K5248	L5047	L5047	Y4291	L4189	I4080	SER	A3219	G3136	M3035	V2275	K2177
N5249	D5048	D5048	S4292	I4192	R4085	A4001	V3221	V3142	I3036	L2275	L2182
K5250	L5051	L5051	I4298	J4193	R4085	D4003	R3222	V3143	V3037	G2278	K2183
I5251	L5051	L5051	I4298	Q4194	T4089	D4003	L3223	D3144	I3040	E2284	S2185
F5260	G5055	G5055	P4313	I4195	C4090	K4006	D3224	S3145	S3041	E2285	S2185
G5265	G5055	G5055	E4314	R4196	C4090	Q4007	R3226	V3146	T3042	L2284	L2189
G5265	M5058	M5058	K4317	M4197	H4097	L4010	K3248	L3149	G3043	Y2291	L2192
I5268	G5069	G5069	Y4323	K4198	A4098	A4011	I3249	T3150	S3044	S2292	I2193
Y5271	R5060	R5060	V4323	V4201	P4101	L4014	I3250	P3151	L3047	I2298	I2194
G5272	I5061	I5061	L4328	M4202	I4102	L4014	I3251	E3154	D3048	P2313	I2195
L5274	V5062	V5062	S4329	F4203	Y4103	F4021	A3252	I3159	I3051	E2314	R2196
L5274	E5068	E5068	N4330	G4204	A4104	A3253	A3253	G3160	G3052	T2315	M2197
V5275	K5072	K5072	S4333	N4205	R4105	K4023	P3254	I3169	A3053	G2316	V2201
N5181	L5077	L5077	THR	E4207	I4111	K4023	P3260	H3163	G3055	K2317	M2202
K5183	G5078	G5078	THR	T4208	D4112	I4026	Y3264	Q3173	M3058	L2328	G2204
Q5184	T5076	T5076	GLY	T4209	M4113	M4027	G3265	I3175	G3059	S2329	M2205
S5185	L5077	L5077	THR	G4211	L4114	R4028	I3268	R3176	R3060	M2330	P2206
L5189	G5078	G5078	GLY	G4212	L4115	G4030	I3268	R3177	I3061	E2207	E2207
I5192	V5079	V5079	SER	N4213	Q4118	S4034	Y3271	K3177	V3062	S2333	T2208
Q5194	I5080	I5080	GLY	K4216	G4122	M4035	G3272	G3180	E3068	THR	T2209
I5195	R5085	R5085	SER	F4217	E4123	T4040	L3274	H3181	S3069	GLY	G2211
R5196	T5088	T5088	GLY	Y4218	Q4124	S4041	I3275	L3182	S3070	THR	G2212
MET	C5090	C5090	GLY	S4220	A4125	T4042	G3278	K3183	G3071	SER	M2213
L5199	D5100	D5100	SER	V4221	G4136	G4043	G3278	Q3184	K3072	THR	A2214
I5102	P5101	P5101	GLY	R4222	V4142	S4044	L3283	S3185	T3073	GLY	L2215
Y5103	I5102	I5102	GLY	L4223	V4146	L4047	E3285	L3189	T3076	SER	K2216
R5105	Y5103	Y5103	GLY	D4224	V4146	D4048	E3285	I3192	T3077	ALA	Y2218
I5111	A5104	A5104	GLY	I4225	A4147	I4049	Y3291	N3193	G3078	SER	A2219
I5111	R5105	R5105	GLY	R4226	A4148	A4050	S3292	Q3194	V3079	THR	S2220
I5114	I5111	I5111	GLY	K4248	L4149	L4051	S3292	I3195	I3080	GLY	V2221
L5115	L5114	L5114	GLY	N4249	T4150	G4052	I3298	R3196	R3085	THR	L2223
G5122	G5122	G5122	GLY	K4250	P4151	A4053	P3313	M3197	T3089	GLY	I2224
Q5124	A5125	A5125	GLY	I4251	E4154	G4054	E3314	I3197	C3090	SER	R2225
A5125	L5115	L5115	GLY	P4254	I4159	M4088	K3317	V3201	A3001	ALA	E2227
G5136	G5136	G5136	GLY	F4260	G4160	G4059	I3317	M3202	Q3007	THR	K2248
V5142	V5142	V5142	GLY	G4265	H4163	R4060	N3330	F3203	I3102	GLY	K2249
K5216	K5216	K5216	GLY	I4268	H4163	I4061	S3333	N3205	Y3103	THR	M2249
F5217	F5217	F5217	GLY	Q4173	R4169	V4062	THR	P3206	A3104	GLY	K2250
				G4272	A4174	E4063	GLY	E3207	R3105	THR	I2251
						E4068		T3208	I3111	THR	P2254
						S4069		T3210	D3112	THR	K3023

GLOBAL-STATISTICS INFOmissingINFO

4 Model quality ⓘ

4.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ALF, MG, ADP

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	B	2.06	5/326 (1.5%)	2.67	39/502 (7.8%)
2	A	0.59	0/14765	0.67	0/19866
All	All	0.66	5/15091 (0.0%)	0.79	39/20368 (0.2%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1002	DT	N3-C4	6.64	1.44	1.38
1	B	1002	DT	C1'-N1	6.34	1.57	1.49
1	B	1012	DT	O3'-P	-6.04	1.53	1.61
1	B	1002	DT	N1-C6	5.66	1.42	1.38
1	B	1002	DT	N1-C2	5.58	1.42	1.38

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1014	DT	O4'-C1'-N1	11.00	115.70	108.00
1	B	1011	DT	O4'-C1'-N1	10.59	115.41	108.00
1	B	1005	DT	O4'-C1'-N1	9.55	114.69	108.00
1	B	1008	DT	O4'-C1'-N1	9.30	114.51	108.00
1	B	1002	DT	N3-C4-O4	9.05	125.33	119.90
1	B	1012	DT	C5-C4-O4	-8.23	119.14	124.90
1	B	1015	DT	C5-C4-O4	-7.85	119.40	124.90
1	B	1012	DT	P-O3'-C3'	-7.82	110.32	119.70
1	B	1009	DT	C5-C4-O4	-7.78	119.45	124.90
1	B	1003	DT	P-O3'-C3'	7.58	128.80	119.70
1	B	1006	DT	C5-C4-O4	-7.58	119.59	124.90
1	B	1003	DT	N3-C4-O4	7.38	124.33	119.90
1	B	1014	DT	P-O3'-C3'	6.99	128.09	119.70
1	B	1002	DT	C5-C4-O4	-6.94	120.04	124.90

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1012	DT	N3-C4-O4	6.91	124.04	119.90
1	B	1003	DT	C5-C4-O4	-6.87	120.09	124.90
1	B	1008	DT	P-O3'-C3'	6.65	127.68	119.70
1	B	1015	DT	N3-C4-O4	6.61	123.87	119.90
1	B	1009	DT	N3-C4-O4	6.59	123.85	119.90
1	B	1005	DT	P-O3'-C3'	6.42	127.40	119.70
1	B	1006	DT	N3-C4-O4	6.35	123.71	119.90
1	B	1002	DT	C5-C6-N1	6.23	127.44	123.70
1	B	1008	DT	C4'-C3'-C2'	-6.23	97.49	103.10
1	B	1011	DT	C4'-C3'-C2'	-6.20	97.52	103.10
1	B	1011	DT	P-O3'-C3'	6.14	127.07	119.70
1	B	1005	DT	C4'-C3'-C2'	-5.92	97.78	103.10
1	B	1002	DT	N1-C1'-C2'	5.89	123.80	112.60
1	B	1014	DT	C4'-C3'-C2'	-5.79	97.89	103.10
1	B	1003	DT	O4'-C1'-N1	5.67	111.97	108.00
1	B	1012	DT	O4'-C1'-N1	5.63	111.94	108.00
1	B	1002	DT	C2-N1-C1'	5.46	126.94	118.20
1	B	1006	DT	O3'-P-O5'	-5.39	93.76	104.00
1	B	1007	DT	N3-C4-O4	5.24	123.04	119.90
1	B	1009	DT	O4'-C1'-N1	5.20	111.64	108.00
1	B	1012	DT	OP2-P-O3'	5.14	116.50	105.20
1	B	1010	DT	N3-C4-O4	5.11	122.97	119.90
1	B	1015	DT	O4'-C1'-N1	5.11	111.57	108.00
1	B	1008	DT	C1'-O4'-C4'	-5.05	105.05	110.10
1	B	1006	DT	O4'-C1'-N1	5.01	111.51	108.00

There are no chirality outliers.

There are no planarity outliers.

4.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	297	0	182	25	0
2	A	14598	0	14995	523	1
3	A	6	0	0	0	0
4	A	30	0	0	8	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	162	0	72	24	0
All	All	15093	0	15249	538	1

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 18.

All (538) 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)
2:A:4194:GLN:HE21	2:A:4196:ARG:HH12	1.10	1.00
2:A:2194:GLN:HE21	2:A:2196:ARG:HH12	1.09	0.98
2:A:194:GLN:HE21	2:A:196:ARG:HH12	1.08	0.98
2:A:5194:GLN:HE21	2:A:5196:ARG:HH12	1.11	0.95
2:A:3194:GLN:HE21	2:A:3196:ARG:HH12	1.10	0.95
2:A:2068:GLU:HG2	2:A:3216:LYS:HB3	1.49	0.94
2:A:1194:GLN:HE21	2:A:1196:ARG:HH12	1.10	0.92
2:A:4160:GLY:H	2:A:5173:GLN:HE22	1.10	0.91
2:A:1160:GLY:H	2:A:2173:GLN:HE22	1.19	0.88
2:A:3194:GLN:NE2	2:A:3196:ARG:HH12	1.72	0.87
2:A:194:GLN:NE2	2:A:196:ARG:HH12	1.71	0.87
2:A:4194:GLN:NE2	2:A:4196:ARG:HH12	1.73	0.87
2:A:2194:GLN:NE2	2:A:2196:ARG:HH12	1.73	0.87
2:A:2159:ILE:HA	2:A:3173:GLN:NE2	1.89	0.86
2:A:3068:GLU:HG2	2:A:4216:LYS:HB3	1.58	0.85
2:A:5194:GLN:NE2	2:A:5196:ARG:HH12	1.74	0.85
2:A:5146:VAL:HA	2:A:5149:LEU:HD13	1.59	0.85
2:A:1194:GLN:NE2	2:A:1196:ARG:HH12	1.73	0.85
2:A:1146:VAL:HA	2:A:1149:LEU:HD13	1.59	0.84
2:A:3159:ILE:HA	2:A:4173:GLN:NE2	1.94	0.83
2:A:4146:VAL:HA	2:A:4149:LEU:HD13	1.61	0.82
2:A:146:VAL:HA	2:A:149:LEU:HD13	1.60	0.82
2:A:2069:SER:OG	2:A:3248:LYS:HG3	1.80	0.81
2:A:2146:VAL:HA	2:A:2149:LEU:HD13	1.62	0.81
2:A:154:GLU:O	2:A:1177:LYS:HE2	1.81	0.81
2:A:4115:LEU:HD22	2:A:5026:ILE:HG12	1.62	0.80
2:A:5068:GLU:HA	4:A:5501:ALF:F1	1.71	0.80
2:A:2111:ILE:HG22	2:A:3030:GLY:HA2	1.64	0.80
2:A:3146:VAL:HA	2:A:3149:LEU:HD13	1.61	0.80
2:A:3160:GLY:H	2:A:4173:GLN:HE22	1.30	0.80
2:A:2160:GLY:H	2:A:3173:GLN:HE22	1.28	0.79
2:A:2111:ILE:HG22	2:A:3030:GLY:CA	2.13	0.79

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:3069:SER:OG	2:A:4248:LYS:HG3	1.85	0.77
2:A:1069:SER:OG	2:A:2248:LYS:HG3	1.86	0.76
2:A:1068:GLU:HG2	2:A:2216:LYS:HB3	1.68	0.75
2:A:3068:GLU:HG2	2:A:4216:LYS:CB	2.17	0.75
2:A:2068:GLU:HG2	2:A:3216:LYS:CB	2.16	0.74
2:A:4160:GLY:N	2:A:5173:GLN:HE22	1.84	0.74
2:A:4111:ILE:HG22	2:A:5030:GLY:CA	2.18	0.74
2:A:2071:GLY:HA2	5:A:2502:ADP:H5'1	1.72	0.72
2:A:4154:GLU:O	2:A:5177:LYS:HE2	1.89	0.72
2:A:3254:PRO:HG3	5:A:2502:ADP:O2'	1.90	0.71
2:A:42:THR:HG21	2:A:47:LEU:HD22	1.73	0.71
2:A:1042:THR:HG21	2:A:1047:LEU:HD22	1.74	0.70
2:A:4112:ASP:O	2:A:5028:ARG:HG2	1.91	0.70
2:A:4159:ILE:HA	2:A:5173:GLN:NE2	2.08	0.69
2:A:96:GLU:OE2	4:A:501:ALF:F4	1.99	0.69
2:A:4042:THR:HG21	2:A:4047:LEU:HD22	1.74	0.68
2:A:1159:ILE:HG12	2:A:2177:LYS:HD3	1.74	0.68
2:A:1068:GLU:HG2	2:A:2216:LYS:CB	2.23	0.68
2:A:2071:GLY:HA2	5:A:2502:ADP:C5'	2.24	0.68
2:A:4111:ILE:HG22	2:A:5030:GLY:N	2.09	0.68
2:A:3042:THR:HG21	2:A:3047:LEU:HD22	1.75	0.68
2:A:5042:THR:HG21	2:A:5047:LEU:HD22	1.76	0.68
2:A:160:GLY:H	2:A:1173:GLN:HE22	1.42	0.67
2:A:5072:LYS:N	5:A:5502:ADP:O2B	2.27	0.67
2:A:2049:ILE:HA	2:A:2328:LEU:CD2	2.24	0.67
2:A:4072:LYS:N	5:A:4502:ADP:O2B	2.25	0.67
2:A:2159:ILE:HA	2:A:3173:GLN:HE22	1.57	0.67
2:A:3159:ILE:HG12	2:A:4177:LYS:HD3	1.77	0.67
2:A:1159:ILE:HA	2:A:2173:GLN:NE2	2.10	0.66
2:A:4159:ILE:HG12	2:A:5177:LYS:HD3	1.78	0.66
2:A:4291:TYR:O	2:A:4298:ILE:HG12	1.95	0.66
2:A:3291:TYR:O	2:A:3298:ILE:HG12	1.96	0.65
2:A:2042:THR:HG21	2:A:2047:LEU:HD22	1.76	0.65
2:A:3159:ILE:HA	2:A:4173:GLN:HE22	1.60	0.65
2:A:4193:ASN:HD22	2:A:4194:GLN:H	1.45	0.65
2:A:111:ILE:HG23	2:A:1029:LEU:HD13	1.79	0.65
2:A:1291:TYR:O	2:A:1298:ILE:HG12	1.97	0.64
2:A:1072:LYS:N	5:A:1502:ADP:O2B	2.31	0.64
2:A:3101:PRO:HD3	2:A:4035:MET:HG3	1.79	0.63
2:A:2291:TYR:O	2:A:2298:ILE:HG12	1.99	0.63
2:A:3159:ILE:HG23	2:A:4173:GLN:HE21	1.64	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1160:GLY:N	2:A:2173:GLN:HE22	1.92	0.63
2:A:291:TYR:O	2:A:298:ILE:HG12	1.98	0.63
2:A:2160:GLY:N	2:A:3173:GLN:HE22	1.96	0.63
2:A:5193:ASN:HD22	2:A:5194:GLN:H	1.47	0.62
2:A:5291:TYR:O	2:A:5298:ILE:HG12	1.98	0.62
2:A:2072:LYS:HB3	2:A:2192:ILE:HG23	1.80	0.62
2:A:4278:GLY:HA3	2:A:4284:ILE:HD12	1.82	0.62
2:A:193:ASN:HD22	2:A:194:GLN:H	1.47	0.62
2:A:5278:GLY:HA3	2:A:5284:ILE:HD12	1.82	0.61
2:A:3072:LYS:HB3	2:A:3192:ILE:HG23	1.83	0.61
2:A:3072:LYS:N	5:A:3502:ADP:O2B	2.31	0.61
2:A:1278:GLY:HA3	2:A:1284:ILE:HD12	1.82	0.61
2:A:4197:MET:HA	2:A:4206:PRO:O	2.01	0.61
2:A:5208:THR:HG22	2:A:5209:THR:H	1.65	0.61
2:A:2150:THR:OG1	2:A:3176:ARG:HG2	2.01	0.61
1:B:1003:DT:N3	2:A:198:LYS:HB3	2.15	0.61
2:A:1193:ASN:HD22	2:A:1194:GLN:H	1.49	0.61
2:A:5072:LYS:HB3	2:A:5192:ILE:HG23	1.82	0.60
2:A:72:LYS:HB3	2:A:192:ILE:HG23	1.82	0.60
2:A:4208:THR:HG22	2:A:4209:THR:H	1.65	0.60
2:A:2159:ILE:HG12	2:A:3177:LYS:HD3	1.82	0.60
2:A:3193:ASN:HD22	2:A:3194:GLN:H	1.48	0.60
2:A:4254:PRO:HG3	5:A:3502:ADP:O2'	2.01	0.60
2:A:5164:MET:O	2:A:5166:LEU:N	2.35	0.60
1:B:1003:DT:H3	2:A:198:LYS:HB3	1.66	0.60
2:A:1154:GLU:O	2:A:2177:LYS:HE2	2.02	0.60
2:A:1072:LYS:HB3	2:A:1192:ILE:HG23	1.82	0.60
2:A:4068:GLU:HB3	2:A:5217:PHE:HA	1.83	0.60
1:B:1004:DT:C6	1:B:1005:DT:H72	2.37	0.59
2:A:2193:ASN:HD22	2:A:2194:GLN:H	1.49	0.59
2:A:3160:GLY:N	2:A:4173:GLN:HE22	1.97	0.59
2:A:3071:GLY:HA2	5:A:3502:ADP:O1A	2.02	0.59
2:A:4072:LYS:HB3	2:A:4192:ILE:HG23	1.83	0.59
2:A:3208:THR:HG22	2:A:3209:THR:H	1.68	0.59
2:A:208:THR:HG22	2:A:209:THR:H	1.68	0.59
2:A:4068:GLU:HG2	2:A:5216:LYS:HB3	1.84	0.59
2:A:68:GLU:HG2	2:A:1216:LYS:HB3	1.85	0.58
2:A:73:THR:O	2:A:76:THR:HB	2.04	0.58
2:A:2111:ILE:HG22	2:A:3030:GLY:N	2.18	0.58
2:A:2154:GLU:O	2:A:3177:LYS:HE2	2.04	0.58
2:A:2254:PRO:HG3	5:A:1502:ADP:O2'	2.04	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:3250:LYS:HA	5:A:2502:ADP:N7	2.18	0.58
2:A:3115:LEU:HD22	2:A:4026:ILE:HG12	1.85	0.58
2:A:2208:THR:HG22	2:A:2209:THR:H	1.69	0.58
2:A:1197:MET:HA	2:A:1206:PRO:O	2.05	0.57
2:A:2049:ILE:HA	2:A:2328:LEU:HD21	1.84	0.57
1:B:1010:DT:C6	1:B:1011:DT:H72	2.40	0.57
2:A:4069:SER:OG	2:A:5248:LYS:HG3	2.04	0.57
2:A:1250:LYS:HE3	4:A:501:ALF:F3	1.95	0.57
2:A:3111:ILE:HG23	2:A:4029:LEU:HD13	1.86	0.57
2:A:3053:ALA:HA	2:A:3330:ASN:O	2.05	0.56
2:A:4154:GLU:HG3	2:A:4163:HIS:HD2	1.69	0.56
2:A:5250:LYS:HE3	4:A:4501:ALF:F3	1.94	0.56
2:A:4111:ILE:HG22	2:A:5030:GLY:HA2	1.85	0.56
2:A:5073:THR:O	2:A:5076:THR:HB	2.05	0.56
1:B:1007:DT:C6	1:B:1008:DT:H72	2.40	0.56
2:A:2048:ASP:HA	2:A:2051:LEU:HD12	1.87	0.56
2:A:5076:THR:HG22	2:A:5077:LEU:N	2.21	0.56
2:A:4048:ASP:HA	2:A:4051:LEU:HD12	1.87	0.56
2:A:1208:THR:HG22	2:A:1209:THR:H	1.71	0.56
2:A:1210:THR:HG22	2:A:1211:GLY:H	1.71	0.56
2:A:210:THR:HG22	2:A:211:GLY:H	1.71	0.56
2:A:2210:THR:HG22	2:A:2211:GLY:H	1.70	0.56
2:A:3197:MET:HA	2:A:3206:PRO:O	2.05	0.56
2:A:3278:GLY:HA3	2:A:3284:ILE:HD12	1.87	0.56
2:A:5048:ASP:HA	2:A:5051:LEU:HD12	1.88	0.56
1:B:1013:DT:C6	1:B:1014:DT:H72	2.40	0.56
2:A:3048:ASP:HA	2:A:3051:LEU:HD12	1.88	0.55
2:A:4193:ASN:ND2	2:A:4194:GLN:H	2.04	0.55
2:A:2278:GLY:HA3	2:A:2284:ILE:HD12	1.88	0.55
2:A:2073:THR:O	2:A:2076:THR:HB	2.06	0.55
2:A:3248:LYS:NZ	4:A:2501:ALF:F1	2.21	0.55
2:A:2285:GLU:HB3	2:A:2292:SER:HB2	1.89	0.55
2:A:5271:TYR:O	2:A:5275:VAL:HG23	2.06	0.55
2:A:40:ILE:HG22	2:A:41:SER:O	2.06	0.55
2:A:1271:TYR:O	2:A:1275:VAL:HG23	2.06	0.55
2:A:2136:GLY:HA2	2:A:2185:SER:HB3	1.89	0.55
2:A:4136:GLY:HA2	2:A:4185:SER:HB3	1.89	0.55
2:A:3111:ILE:HG22	2:A:4030:GLY:N	2.22	0.55
2:A:4073:THR:O	2:A:4076:THR:HB	2.07	0.55
2:A:5136:GLY:HA2	2:A:5185:SER:HB3	1.89	0.55
2:A:3073:THR:O	2:A:3076:THR:HB	2.07	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:5040:ILE:HG22	2:A:5041:SER:O	2.07	0.54
2:A:1273:GLU:O	2:A:1274:LEU:C	2.45	0.54
1:B:1003:DT:H3'	2:A:196:ARG:HB3	1.89	0.54
2:A:278:GLY:HA3	2:A:284:ILE:HD12	1.88	0.54
2:A:4210:THR:HG22	2:A:4211:GLY:H	1.72	0.54
2:A:3136:GLY:HA2	2:A:3185:SER:HB3	1.90	0.54
2:A:3210:THR:HG22	2:A:3211:GLY:H	1.72	0.54
2:A:1048:ASP:HA	2:A:1051:LEU:HD12	1.89	0.54
2:A:5210:THR:HG22	2:A:5211:GLY:H	1.72	0.54
2:A:1073:THR:O	2:A:1076:THR:HB	2.08	0.54
2:A:3150:THR:OG1	2:A:4176:ARG:HG2	2.08	0.54
2:A:3111:ILE:HG22	2:A:4030:GLY:CA	2.37	0.54
2:A:4040:ILE:HG22	2:A:4041:SER:O	2.08	0.54
2:A:4076:THR:HG22	2:A:4077:LEU:N	2.22	0.54
1:B:1015:DT:H6	1:B:1015:DT:O5'	1.91	0.54
2:A:1111:ILE:HG23	2:A:2029:LEU:HD13	1.90	0.54
2:A:2123:GLU:OE1	2:A:2151:PRO:HA	2.08	0.53
2:A:2154:GLU:OE2	2:A:3176:ARG:HD2	2.07	0.53
2:A:3076:THR:HG22	2:A:3077:LEU:N	2.23	0.53
2:A:4007:GLN:HA	2:A:4010:LEU:HB3	1.91	0.53
2:A:5193:ASN:ND2	2:A:5194:GLN:H	2.06	0.53
2:A:3040:ILE:HG12	2:A:3058:MET:SD	2.48	0.53
2:A:40:ILE:HG12	2:A:58:MET:SD	2.48	0.53
2:A:3273:GLU:O	2:A:3274:LEU:C	2.46	0.53
2:A:4271:TYR:O	2:A:4275:VAL:HG23	2.09	0.53
2:A:5040:ILE:HG12	2:A:5058:MET:SD	2.49	0.53
2:A:48:ASP:HA	2:A:51:LEU:HD12	1.90	0.53
1:B:1006:DT:O5'	1:B:1006:DT:H6	1.92	0.53
2:A:210:THR:HG22	2:A:211:GLY:N	2.24	0.53
2:A:76:THR:HG22	2:A:77:LEU:N	2.23	0.53
2:A:2264:TYR:CD2	2:A:3254:PRO:HD2	2.43	0.53
2:A:4053:ALA:HA	2:A:4330:ASN:O	2.09	0.53
2:A:1248:LYS:NZ	4:A:501:ALF:F1	2.17	0.53
2:A:1136:GLY:HA2	2:A:1185:SER:HB3	1.89	0.53
2:A:136:GLY:HA2	2:A:185:SER:HB3	1.90	0.53
2:A:2076:THR:HG22	2:A:2077:LEU:N	2.24	0.53
2:A:2040:ILE:HG22	2:A:2041:SER:O	2.10	0.52
2:A:2111:ILE:CG2	2:A:3030:GLY:HA2	2.38	0.52
2:A:2007:GLN:HA	2:A:2010:LEU:HB3	1.91	0.52
2:A:3040:ILE:HG22	2:A:3041:SER:O	2.08	0.52
2:A:1210:THR:HG22	2:A:1211:GLY:N	2.24	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:285:GLU:HB3	2:A:292:SER:HB2	1.90	0.52
2:A:4115:LEU:HD13	2:A:5026:ILE:HD11	1.91	0.52
2:A:2271:TYR:O	2:A:2275:VAL:HG23	2.10	0.52
2:A:4273:GLU:O	2:A:4274:LEU:C	2.48	0.52
2:A:3123:GLU:OE1	2:A:3151:PRO:HA	2.09	0.52
2:A:2210:THR:HG22	2:A:2211:GLY:N	2.25	0.52
2:A:1249:ASN:OD1	2:A:1251:ILE:HG22	2.10	0.52
2:A:4040:ILE:HG12	2:A:4058:MET:SD	2.50	0.52
2:A:5076:THR:O	2:A:5079:VAL:HG22	2.09	0.52
2:A:4210:THR:HG22	2:A:4211:GLY:N	2.25	0.52
2:A:5249:ASN:OD1	2:A:5251:ILE:HG22	2.10	0.52
2:A:1076:THR:HG22	2:A:1077:LEU:N	2.23	0.51
2:A:1076:THR:O	2:A:1079:VAL:HG22	2.10	0.51
2:A:193:ASN:ND2	2:A:194:GLN:H	2.08	0.51
2:A:5123:GLU:OE1	2:A:5151:PRO:HA	2.09	0.51
2:A:76:THR:O	2:A:79:VAL:HG22	2.10	0.51
2:A:273:GLU:O	2:A:274:LEU:C	2.48	0.51
2:A:2040:ILE:HG12	2:A:2058:MET:SD	2.50	0.51
2:A:3193:ASN:ND2	2:A:3194:GLN:H	2.07	0.51
2:A:4060:ARG:HB3	2:A:4220:SER:OG	2.10	0.51
2:A:4071:GLY:HA2	5:A:4502:ADP:H5'1	1.91	0.51
1:B:1009:DT:O5'	1:B:1009:DT:H6	1.93	0.51
2:A:2076:THR:O	2:A:2079:VAL:HG22	2.11	0.51
2:A:2197:MET:HA	2:A:2206:PRO:O	2.11	0.51
2:A:1040:ILE:HG22	2:A:1041:SER:O	2.09	0.51
2:A:3007:GLN:HA	2:A:3010:LEU:HB3	1.93	0.51
2:A:1007:GLN:HA	2:A:1010:LEU:HB3	1.92	0.51
2:A:159:ILE:HG12	2:A:1177:LYS:HD3	1.92	0.51
2:A:4154:GLU:HG3	2:A:4163:HIS:CD2	2.46	0.51
2:A:5210:THR:HG22	2:A:5211:GLY:N	2.25	0.51
1:B:1012:DT:O5'	1:B:1012:DT:H6	1.93	0.51
2:A:1193:ASN:ND2	2:A:1194:GLN:H	2.09	0.51
2:A:123:GLU:OE1	2:A:151:PRO:HA	2.10	0.51
2:A:4076:THR:O	2:A:4079:VAL:HG22	2.11	0.51
2:A:4123:GLU:OE1	2:A:4151:PRO:HA	2.11	0.51
2:A:2273:GLU:O	2:A:2274:LEU:C	2.47	0.50
2:A:5285:GLU:HB3	2:A:5292:SER:HB2	1.93	0.50
2:A:2193:ASN:ND2	2:A:2194:GLN:H	2.08	0.50
2:A:2069:SER:OG	2:A:3248:LYS:CG	2.56	0.50
2:A:1053:ALA:HA	2:A:1330:ASN:O	2.12	0.50
2:A:1193:ASN:HD21	2:A:1210:THR:HB	1.76	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:5103:TYR:O	2:A:5104:ALA:C	2.50	0.50
1:B:1011:DT:H2''	2:A:4169:ARG:HH11	1.76	0.50
2:A:2249:ASN:OD1	2:A:2251:ILE:HG22	2.10	0.50
2:A:3249:ASN:OD1	2:A:3251:ILE:HG22	2.11	0.50
2:A:5273:GLU:O	2:A:5274:LEU:C	2.48	0.50
2:A:2101:PRO:HD3	2:A:3035:MET:HG3	1.94	0.50
2:A:4249:ASN:OD1	2:A:4251:ILE:HG22	2.12	0.50
2:A:249:ASN:OD1	2:A:251:ILE:HG22	2.12	0.49
2:A:3111:ILE:HD12	2:A:4029:LEU:HD13	1.94	0.49
2:A:4097:HIS:NE2	2:A:5180:GLY:HA2	2.27	0.49
2:A:3210:THR:HG22	2:A:3211:GLY:N	2.26	0.49
2:A:1123:GLU:OE1	2:A:1151:PRO:HA	2.11	0.49
2:A:1077:LEU:O	2:A:1080:ILE:HB	2.13	0.49
2:A:1285:GLU:HB3	2:A:1292:SER:HB2	1.93	0.49
2:A:1060:ARG:HB3	2:A:1220:SER:OG	2.13	0.49
2:A:1062:VAL:HG22	2:A:1221:VAL:HB	1.95	0.49
2:A:3068:GLU:HB3	2:A:4217:PHE:HA	1.93	0.49
2:A:3193:ASN:HD21	2:A:3210:THR:HB	1.78	0.49
2:A:5074:THR:OG1	5:A:5502:ADP:H5'1	2.11	0.49
2:A:1103:TYR:O	2:A:1104:ALA:C	2.51	0.49
1:B:1003:DT:O5'	1:B:1003:DT:H6	1.95	0.49
2:A:4049:ILE:HA	2:A:4328:LEU:CD2	2.42	0.49
2:A:2011:ALA:HA	2:A:2014:LEU:HB2	1.94	0.49
2:A:2146:VAL:CA	2:A:2149:LEU:HD13	2.39	0.49
2:A:69:SER:OG	2:A:1248:LYS:HG3	2.13	0.49
2:A:1040:ILE:HG12	2:A:1058:MET:SD	2.53	0.49
2:A:5007:GLN:HA	2:A:5010:LEU:HB3	1.94	0.49
2:A:4285:GLU:HB3	2:A:4292:SER:HB2	1.95	0.48
2:A:5060:ARG:HB3	2:A:5220:SER:OG	2.12	0.48
2:A:1011:ALA:HA	2:A:1014:LEU:HB2	1.95	0.48
2:A:3285:GLU:HB3	2:A:3292:SER:HB2	1.94	0.48
2:A:2103:TYR:O	2:A:2104:ALA:C	2.51	0.48
2:A:3271:TYR:O	2:A:3275:VAL:HG23	2.13	0.48
2:A:4205:ASN:HA	2:A:4206:PRO:HD3	1.64	0.48
2:A:3076:THR:O	2:A:3079:VAL:HG22	2.13	0.48
2:A:1068:GLU:HB3	2:A:2217:PHE:HA	1.96	0.48
2:A:4103:TYR:O	2:A:4104:ALA:C	2.50	0.48
2:A:4193:ASN:HD21	2:A:4210:THR:HB	1.78	0.48
2:A:5077:LEU:O	2:A:5080:ILE:HB	2.13	0.48
2:A:193:ASN:HD21	2:A:210:THR:HB	1.78	0.48
2:A:3011:ALA:HA	2:A:3014:LEU:HB2	1.95	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:4044:SER:HB2	2:A:4268:ILE:HD13	1.96	0.48
2:A:4313:PRO:O	2:A:4317:LYS:HG2	2.14	0.48
2:A:2060:ARG:HB3	2:A:2220:SER:OG	2.13	0.48
2:A:4068:GLU:HG2	2:A:5216:LYS:CB	2.43	0.48
2:A:103:TYR:O	2:A:104:ALA:C	2.51	0.48
2:A:1044:SER:HB2	2:A:1268:ILE:HD13	1.96	0.48
2:A:271:TYR:O	2:A:275:VAL:HG23	2.13	0.48
2:A:3154:GLU:O	2:A:4177:LYS:HE2	2.14	0.48
2:A:2097:HIS:CE1	2:A:3180:GLY:HA2	2.49	0.48
2:A:90:CYS:HB2	2:A:114:LEU:HD23	1.96	0.48
2:A:60:ARG:HB3	2:A:220:SER:OG	2.14	0.47
2:A:3103:TYR:O	2:A:3104:ALA:C	2.50	0.47
2:A:3124:GLN:HG3	2:A:4021:PHE:CD1	2.48	0.47
2:A:4071:GLY:HA2	5:A:4502:ADP:C5'	2.43	0.47
2:A:5193:ASN:HD21	2:A:5210:THR:HB	1.79	0.47
2:A:3216:LYS:HA	2:A:3222:ARG:HD3	1.96	0.47
2:A:4077:LEU:O	2:A:4080:ILE:HB	2.14	0.47
2:A:216:LYS:HA	2:A:222:ARG:HD3	1.96	0.47
2:A:2194:GLN:HG2	2:A:3217:PHE:CD2	2.50	0.47
2:A:4208:THR:HG22	2:A:4209:THR:N	2.29	0.47
2:A:5216:LYS:HA	2:A:5222:ARG:HD3	1.97	0.47
2:A:2193:ASN:HD21	2:A:2210:THR:HB	1.79	0.47
2:A:3032:ASP:HB3	2:A:3035:MET:HB2	1.96	0.47
2:A:5208:THR:HG22	2:A:5209:THR:N	2.29	0.47
2:A:68:GLU:HG2	2:A:1216:LYS:CB	2.44	0.47
2:A:2118:GLN:NE2	2:A:3183:LYS:HG2	2.29	0.47
2:A:3111:ILE:HG22	2:A:4030:GLY:HA2	1.95	0.47
2:A:3040:ILE:O	2:A:3055:GLY:HA3	2.15	0.47
2:A:3090:CYS:HB2	2:A:3114:LEU:HD23	1.97	0.47
2:A:205:ASN:HA	2:A:206:PRO:HD3	1.63	0.47
2:A:2216:LYS:HA	2:A:2222:ARG:HD3	1.95	0.47
2:A:313:PRO:O	2:A:317:LYS:HG2	2.15	0.47
2:A:4146:VAL:CA	2:A:4149:LEU:HD13	2.39	0.47
2:A:1105:ARG:N	2:A:1111:ILE:HD11	2.30	0.47
2:A:3077:LEU:O	2:A:3080:ILE:HB	2.15	0.47
2:A:5122:GLY:O	2:A:5125:ALA:HB3	2.14	0.47
2:A:146:VAL:CA	2:A:149:LEU:HD13	2.38	0.47
2:A:3069:SER:OG	2:A:4248:LYS:CG	2.61	0.47
2:A:4115:LEU:HA	2:A:4115:LEU:HD23	1.73	0.46
2:A:105:ARG:N	2:A:111:ILE:HD11	2.31	0.46
2:A:1313:PRO:O	2:A:1317:LYS:HG2	2.15	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:2105:ARG:N	2:A:2111:ILE:HD11	2.31	0.46
2:A:3060:ARG:HB3	2:A:3220:SER:OG	2.15	0.46
1:B:1016:DT:H3'	2:A:5212:GLY:HA2	1.97	0.46
2:A:1159:ILE:HG23	2:A:2173:GLN:HE21	1.81	0.46
2:A:2159:ILE:CA	2:A:3173:GLN:HE22	2.28	0.46
2:A:4105:ARG:N	2:A:4111:ILE:HD11	2.31	0.46
2:A:2098:ALA:HA	2:A:3060:ARG:CZ	2.45	0.46
2:A:4090:CYS:HB2	2:A:4114:LEU:HD23	1.97	0.46
2:A:1090:CYS:HB2	2:A:1114:LEU:HD23	1.97	0.46
2:A:2077:LEU:O	2:A:2080:ILE:HB	2.15	0.46
2:A:44:SER:HB2	2:A:268:ILE:HD13	1.96	0.46
2:A:62:VAL:HG22	2:A:221:VAL:HB	1.98	0.46
2:A:3044:SER:HB2	2:A:3268:ILE:HD13	1.97	0.46
2:A:3159:ILE:CA	2:A:4173:GLN:HE22	2.28	0.46
2:A:40:ILE:O	2:A:55:GLY:HA3	2.16	0.46
2:A:4147:ALA:O	2:A:5176:ARG:HG3	2.16	0.46
2:A:77:LEU:O	2:A:80:ILE:HB	2.16	0.46
2:A:68:GLU:HB3	2:A:1217:PHE:HA	1.96	0.45
2:A:76:THR:CG2	2:A:142:VAL:HG21	2.46	0.45
2:A:3252:ALA:C	5:A:2502:ADP:C2	2.89	0.45
2:A:4076:THR:CG2	2:A:4142:VAL:HG21	2.46	0.45
2:A:5062:VAL:HG22	2:A:5221:VAL:HB	1.98	0.45
2:A:1071:GLY:HA2	5:A:1502:ADP:C5'	2.46	0.45
2:A:2044:SER:HB2	2:A:2268:ILE:HD13	1.98	0.45
2:A:2115:LEU:HD22	2:A:3026:ILE:HG12	1.98	0.45
2:A:1061:ILE:HD12	2:A:1061:ILE:N	2.32	0.45
2:A:154:GLU:OE2	2:A:1176:ARG:HD2	2.16	0.45
2:A:3105:ARG:N	2:A:3111:ILE:HD11	2.31	0.45
2:A:3122:GLY:O	2:A:3125:ALA:HB3	2.16	0.45
2:A:1040:ILE:O	2:A:1055:GLY:HA3	2.16	0.45
2:A:2072:LYS:HB3	2:A:2192:ILE:CG2	2.47	0.45
2:A:4160:GLY:H	2:A:5173:GLN:NE2	1.94	0.45
2:A:4216:LYS:HA	2:A:4222:ARG:HD3	1.97	0.45
2:A:4071:GLY:HA2	5:A:4502:ADP:O1A	2.16	0.45
2:A:2208:THR:HG22	2:A:2209:THR:N	2.32	0.45
2:A:3076:THR:CG2	2:A:3142:VAL:HG21	2.46	0.45
2:A:5032:ASP:HB3	2:A:5035:MET:HB2	1.99	0.45
2:A:1146:VAL:CA	2:A:1149:LEU:HD13	2.36	0.45
2:A:115:LEU:HD23	2:A:115:LEU:HA	1.73	0.45
2:A:1216:LYS:HA	2:A:1222:ARG:HD3	1.97	0.45
2:A:2076:THR:CG2	2:A:2142:VAL:HG21	2.47	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:3124:GLN:HG3	2:A:4021:PHE:CG	2.52	0.45
2:A:5044:SER:HB2	2:A:5268:ILE:HD13	1.98	0.45
1:B:1007:DT:H4'	2:A:2168:ALA:CB	2.47	0.45
1:B:1006:DT:H4'	2:A:2169:ARG:HA	1.99	0.45
2:A:4040:ILE:O	2:A:4055:GLY:HA3	2.17	0.45
2:A:4115:LEU:HD13	2:A:5026:ILE:CD1	2.47	0.45
2:A:5090:CYS:HB2	2:A:5114:LEU:HD23	1.98	0.45
2:A:5011:ALA:HA	2:A:5014:LEU:HB2	1.99	0.44
2:A:4062:VAL:HG22	2:A:4221:VAL:HB	1.98	0.44
2:A:2040:ILE:O	2:A:2055:GLY:HA3	2.16	0.44
2:A:208:THR:HG22	2:A:209:THR:N	2.31	0.44
2:A:2062:VAL:HG22	2:A:2221:VAL:HB	2.00	0.44
2:A:4193:ASN:HD22	2:A:4194:GLN:N	2.14	0.44
1:B:1014:DT:C7	1:B:1015:DT:O4	2.65	0.44
2:A:1071:GLY:HA2	5:A:1502:ADP:O1A	2.17	0.44
2:A:122:GLY:O	2:A:125:ALA:HB3	2.18	0.44
2:A:3071:GLY:HA2	5:A:3502:ADP:H5'1	1.99	0.44
2:A:3182:LEU:HD11	2:A:3189:LEU:HD12	1.99	0.44
2:A:4097:HIS:CD2	2:A:5180:GLY:HA2	2.53	0.44
2:A:1072:LYS:HB3	2:A:1192:ILE:CG2	2.48	0.44
2:A:1071:GLY:HA2	5:A:1502:ADP:H5'1	2.00	0.44
2:A:1154:GLU:HG3	2:A:1163:HIS:HD2	1.82	0.44
2:A:2042:THR:O	2:A:2079:VAL:HG12	2.18	0.44
2:A:3154:GLU:HG3	2:A:3163:HIS:HD2	1.82	0.44
2:A:4118:GLN:NE2	2:A:5183:LYS:HG2	2.32	0.44
1:B:1002:DT:H2'	1:B:1003:DT:C5	2.53	0.44
2:A:1003:ASP:HA	2:A:1006:LYS:HG3	1.99	0.44
2:A:2115:LEU:HA	2:A:2115:LEU:HD23	1.72	0.44
2:A:2122:GLY:O	2:A:2125:ALA:HB3	2.17	0.44
2:A:5040:ILE:O	2:A:5055:GLY:HA3	2.17	0.44
2:A:2182:LEU:HD11	2:A:2189:LEU:HD12	2.00	0.44
2:A:4103:TYR:HE1	2:A:4265:GLY:H	1.66	0.44
2:A:5115:LEU:HD23	2:A:5115:LEU:HA	1.73	0.44
2:A:5146:VAL:CA	2:A:5149:LEU:HD13	2.37	0.44
2:A:5225:ILE:O	2:A:5226:ARG:HG3	2.18	0.44
2:A:2090:CYS:HB2	2:A:2114:LEU:HD23	1.99	0.43
2:A:2100:ASP:OD2	2:A:3037:VAL:HG21	2.18	0.43
2:A:3146:VAL:CA	2:A:3149:LEU:HD13	2.38	0.43
2:A:5205:ASN:HA	2:A:5206:PRO:HD3	1.64	0.43
2:A:3208:THR:HG22	2:A:3209:THR:N	2.31	0.43
2:A:5061:ILE:N	2:A:5061:ILE:HD12	2.33	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1042:THR:O	2:A:1079:VAL:HG12	2.18	0.43
4:A:5501:ALF:F2	5:A:5502:ADP:O1B	2.26	0.43
2:A:1103:TYR:HE1	2:A:1265:GLY:H	1.67	0.43
2:A:1205:ASN:HA	2:A:1206:PRO:HD3	1.63	0.43
2:A:2097:HIS:NE2	2:A:3180:GLY:HA2	2.33	0.43
2:A:5182:LEU:HD11	2:A:5189:LEU:HD12	2.01	0.43
2:A:1122:GLY:O	2:A:1125:ALA:HB3	2.19	0.43
2:A:3071:GLY:HA2	5:A:3502:ADP:C5'	2.48	0.43
2:A:4011:ALA:HA	2:A:4014:LEU:HB2	1.99	0.43
2:A:4260:PHE:CD1	2:A:4268:ILE:HG23	2.53	0.43
2:A:5216:LYS:HG2	2:A:5222:ARG:CZ	2.49	0.43
2:A:2250:LYS:HE3	4:A:1501:ALF:F3	2.08	0.43
2:A:2118:GLN:HE22	2:A:3183:LYS:HG2	1.82	0.43
2:A:2194:GLN:NE2	2:A:3217:PHE:CG	2.87	0.43
2:A:3264:TYR:CD2	2:A:4254:PRO:HD2	2.54	0.43
2:A:5105:ARG:N	2:A:5111:ILE:HD11	2.33	0.43
2:A:1208:THR:HG22	2:A:1209:THR:N	2.33	0.43
2:A:197:MET:HA	2:A:206:PRO:O	2.19	0.43
2:A:2216:LYS:HG2	2:A:2222:ARG:CZ	2.49	0.43
2:A:103:TYR:HE1	2:A:265:GLY:H	1.66	0.43
2:A:4111:ILE:CG2	2:A:5030:GLY:HA2	2.49	0.43
2:A:4098:ALA:HA	2:A:5060:ARG:CZ	2.49	0.43
2:A:3103:TYR:HE1	2:A:3265:GLY:H	1.67	0.43
2:A:4122:GLY:O	2:A:4125:ALA:HB3	2.18	0.43
1:B:1004:DT:OP1	2:A:1213:ASN:N	2.51	0.43
2:A:3072:LYS:HB3	2:A:3192:ILE:CG2	2.49	0.43
2:A:3313:PRO:O	2:A:3317:LYS:HG2	2.19	0.43
2:A:4175:MET:O	2:A:4176:ARG:C	2.56	0.43
2:A:4182:LEU:HD11	2:A:4189:LEU:HD12	2.00	0.43
2:A:4209:THR:HG22	2:A:4216:LYS:HE2	2.01	0.43
1:B:1011:DT:C7	1:B:1012:DT:O4	2.67	0.43
2:A:3111:ILE:HG23	2:A:4029:LEU:HB3	2.01	0.42
2:A:3062:VAL:HG22	2:A:3221:VAL:HB	2.00	0.42
2:A:4076:THR:HG23	2:A:4142:VAL:HG21	2.01	0.42
2:A:1182:LEU:HD11	2:A:1189:LEU:HD12	2.00	0.42
2:A:3175:MET:HB3	2:A:3218:TYR:CG	2.54	0.42
2:A:4097:HIS:CE1	2:A:5180:GLY:HA2	2.54	0.42
2:A:1068:GLU:HG2	2:A:2216:LYS:HB2	1.99	0.42
2:A:1159:ILE:HA	2:A:2173:GLN:HE22	1.82	0.42
2:A:182:LEU:HD11	2:A:189:LEU:HD12	2.01	0.42
2:A:2072:LYS:NZ	5:A:2502:ADP:PB	2.92	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:4217:PHE:O	2:A:4248:LYS:NZ	2.52	0.42
2:A:1076:THR:CG2	2:A:1142:VAL:HG21	2.49	0.42
2:A:76:THR:HG23	2:A:142:VAL:HG21	2.01	0.42
2:A:3042:THR:O	2:A:3079:VAL:HG12	2.19	0.42
2:A:3216:LYS:HG2	2:A:3222:ARG:CZ	2.49	0.42
2:A:2032:ASP:HB3	2:A:2035:MET:HB2	2.01	0.42
2:A:2150:THR:HG23	2:A:3176:ARG:HD2	2.01	0.42
2:A:2313:PRO:O	2:A:2317:LYS:HG2	2.19	0.42
2:A:2068:GLU:CG	2:A:3216:LYS:HB3	2.35	0.42
2:A:100:ASP:HA	2:A:101:PRO:HD2	1.89	0.42
2:A:2098:ALA:HB1	2:A:3250:LYS:HD3	2.02	0.42
2:A:216:LYS:HG2	2:A:222:ARG:CZ	2.49	0.42
2:A:4061:ILE:HD12	2:A:4061:ILE:N	2.35	0.42
2:A:4042:THR:O	2:A:4079:VAL:HG12	2.19	0.42
2:A:4111:ILE:HG23	2:A:5029:LEU:HB3	2.01	0.42
2:A:3205:ASN:HA	2:A:3206:PRO:HD3	1.62	0.42
2:A:3260:PHE:CD1	2:A:3268:ILE:HG23	2.55	0.42
2:A:4225:ILE:O	2:A:4226:ARG:HG3	2.20	0.42
2:A:5076:THR:CG2	2:A:5142:VAL:HG21	2.50	0.42
2:A:1175:MET:HB3	2:A:1218:TYR:CG	2.55	0.42
2:A:2225:ILE:O	2:A:2226:ARG:HG3	2.19	0.42
2:A:3202:MET:H	2:A:3202:MET:HE2	1.84	0.42
2:A:2194:GLN:NE2	2:A:3217:PHE:HB2	2.34	0.42
2:A:3225:ILE:O	2:A:3226:ARG:HG3	2.19	0.42
2:A:3101:PRO:HB3	2:A:4029:LEU:HD22	2.00	0.42
2:A:3217:PHE:O	2:A:3248:LYS:NZ	2.52	0.42
2:A:5217:PHE:O	2:A:5248:LYS:NZ	2.53	0.42
2:A:203:PHE:CG	2:A:204:GLY:N	2.88	0.42
2:A:2159:ILE:HG23	2:A:3173:GLN:HE21	1.85	0.42
2:A:5042:THR:O	2:A:5079:VAL:HG12	2.20	0.42
2:A:5103:TYR:HE1	2:A:5265:GLY:H	1.67	0.42
1:B:1003:DT:H5'	2:A:1172:SER:CB	2.50	0.42
2:A:209:THR:HG22	2:A:216:LYS:HE2	2.00	0.41
2:A:217:PHE:O	2:A:248:LYS:NZ	2.52	0.41
2:A:1175:MET:O	2:A:1176:ARG:C	2.59	0.41
2:A:2154:GLU:OE2	2:A:3176:ARG:NH1	2.52	0.41
2:A:42:THR:O	2:A:79:VAL:HG12	2.19	0.41
2:A:3159:ILE:HG12	2:A:4177:LYS:CD	2.49	0.41
2:A:4003:ASP:HA	2:A:4006:LYS:HG3	2.01	0.41
2:A:4101:PRO:HG2	2:A:4102:ILE:HD12	2.03	0.41
2:A:115:LEU:HD22	2:A:1026:ILE:HG12	2.02	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:3224:ASP:OD1	2:A:3226:ARG:NE	2.52	0.41
2:A:4034:SER:O	2:A:4183:LYS:HD2	2.20	0.41
2:A:4216:LYS:HG2	2:A:4222:ARG:CZ	2.49	0.41
2:A:1100:ASP:HA	2:A:1101:PRO:HD2	1.90	0.41
1:B:1007:DT:H71	2:A:1197:MET:HG2	2.03	0.41
2:A:1225:ILE:O	2:A:1226:ARG:HG3	2.21	0.41
2:A:2103:TYR:HE1	2:A:2265:GLY:H	1.67	0.41
2:A:2209:THR:HG22	2:A:2216:LYS:HE2	2.03	0.41
2:A:3209:THR:HG22	2:A:3216:LYS:HE2	2.02	0.41
2:A:4072:LYS:HB3	2:A:4192:ILE:CG2	2.50	0.41
2:A:5101:PRO:HG2	2:A:5102:ILE:HD12	2.03	0.41
2:A:1260:PHE:CD1	2:A:1268:ILE:HG23	2.56	0.41
2:A:2217:PHE:O	2:A:2248:LYS:NZ	2.53	0.41
2:A:4175:MET:HB3	2:A:4218:TYR:CG	2.56	0.41
2:A:5100:ASP:HA	2:A:5101:PRO:HD2	1.88	0.41
2:A:1217:PHE:O	2:A:1248:LYS:NZ	2.54	0.41
2:A:175:MET:O	2:A:176:ARG:C	2.59	0.41
2:A:2175:MET:O	2:A:2176:ARG:C	2.59	0.41
2:A:2213:ASN:O	2:A:2214:ALA:C	2.59	0.41
2:A:4154:GLU:HA	2:A:4163:HIS:CD2	2.55	0.41
2:A:5072:LYS:HB3	2:A:5192:ILE:CG2	2.49	0.41
2:A:2175:MET:HB3	2:A:2218:TYR:CG	2.55	0.41
2:A:3193:ASN:HD22	2:A:3194:GLN:N	2.16	0.41
2:A:5193:ASN:HD22	2:A:5194:GLN:N	2.15	0.41
2:A:61:ILE:N	2:A:61:ILE:HD12	2.35	0.41
2:A:72:LYS:HB2	2:A:72:LYS:HE2	1.88	0.41
2:A:72:LYS:HB3	2:A:192:ILE:CG2	2.48	0.41
2:A:2205:ASN:HA	2:A:2206:PRO:HD3	1.62	0.41
2:A:3076:THR:HG23	2:A:3142:VAL:HG21	2.01	0.41
2:A:3175:MET:HB3	2:A:3218:TYR:CD1	2.56	0.41
1:B:1015:DT:N3	2:A:4198:LYS:HA	2.35	0.41
1:B:1008:DT:C7	1:B:1009:DT:O4	2.69	0.41
2:A:1154:GLU:HG3	2:A:1163:HIS:CD2	2.55	0.41
2:A:2076:THR:HG23	2:A:2142:VAL:HG21	2.02	0.41
2:A:4175:MET:HB3	2:A:4218:TYR:CD1	2.56	0.41
2:A:3042:THR:HA	2:A:3079:VAL:HG12	2.03	0.41
2:A:3112:ASP:O	2:A:4028:ARG:HG2	2.21	0.41
2:A:3213:ASN:O	2:A:3214:ALA:C	2.59	0.40
2:A:4049:ILE:HD11	2:A:4323:VAL:HG12	2.03	0.40
1:B:1004:DT:H2"	1:B:1005:DT:H6	1.86	0.40
2:A:4202:MET:H	2:A:4202:MET:HE2	1.86	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1049:ILE:HA	2:A:1328:LEU:CD2	2.51	0.40
2:A:1068:GLU:H	2:A:1068:GLU:HG3	1.57	0.40
2:A:57:PRO:O	2:A:188:LEU:HD13	2.21	0.40
2:A:175:MET:HB3	2:A:218:TYR:CG	2.57	0.40
2:A:3254:PRO:CG	5:A:2502:ADP:O2'	2.65	0.40
2:A:4224:ASP:OD1	2:A:4226:ARG:NE	2.49	0.40
2:A:5044:SER:OG	2:A:5273:GLU:OE2	2.27	0.40
2:A:5260:PHE:CD1	2:A:5268:ILE:HG23	2.55	0.40
2:A:5313:PRO:O	2:A:5317:LYS:HG2	2.21	0.40
2:A:1115:LEU:HA	2:A:1115:LEU:HD23	1.75	0.40
2:A:1272:GLY:O	2:A:1275:VAL:HB	2.21	0.40
2:A:225:ILE:O	2:A:226:ARG:HG3	2.21	0.40
2:A:2260:PHE:CD1	2:A:2268:ILE:HG23	2.56	0.40
2:A:3251:ILE:HD12	2:A:3251:ILE:HA	1.92	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1276:ASP:OD2	2:A:5235:GLU:OE2[4_556]	2.08	0.12

4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	A	1921/2050 (94%)	1719 (90%)	180 (9%)	22 (1%)	17 59

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	5165	GLY
2	A	1204	GLY

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	A	2204	GLY
2	A	3204	GLY
2	A	4204	GLY
2	A	161	ASP
2	A	1023	LYS
2	A	2023	LYS
2	A	3023	LYS
2	A	4023	LYS
2	A	5023	LYS
2	A	1037	VAL
2	A	1275	VAL
2	A	4275	VAL
2	A	275	VAL
2	A	2275	VAL
2	A	3275	VAL
2	A	5275	VAL
2	A	204	GLY
2	A	1330	ASN
2	A	3037	VAL
2	A	4330	ASN

4.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	A	1524/1606 (95%)	1406 (92%)	118 (8%)	15	50

All (118) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	68	GLU
2	A	76	THR
2	A	85	ARG
2	A	89	THR
2	A	146	VAL
2	A	150	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	A	162	SER
2	A	163	HIS
2	A	183	LYS
2	A	189	LEU
2	A	193	ASN
2	A	197	MET
2	A	202	MET
2	A	213	ASN
2	A	223	LEU
2	A	285	GLU
2	A	314	GLU
2	A	315	THR
2	A	1028	ARG
2	A	1029	LEU
2	A	1068	GLU
2	A	1076	THR
2	A	1085	ARG
2	A	1089	THR
2	A	1146	VAL
2	A	1150	THR
2	A	1183	LYS
2	A	1189	LEU
2	A	1193	ASN
2	A	1201	VAL
2	A	1202	MET
2	A	1203	PHE
2	A	1213	ASN
2	A	1223	LEU
2	A	1285	GLU
2	A	1314	GLU
2	A	1330	ASN
2	A	2014	LEU
2	A	2028	ARG
2	A	2029	LEU
2	A	2068	GLU
2	A	2076	THR
2	A	2079	VAL
2	A	2085	ARG
2	A	2089	THR
2	A	2146	VAL
2	A	2150	THR
2	A	2183	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	A	2189	LEU
2	A	2193	ASN
2	A	2201	VAL
2	A	2202	MET
2	A	2203	PHE
2	A	2213	ASN
2	A	2223	LEU
2	A	2251	ILE
2	A	2285	GLU
2	A	2314	GLU
2	A	2315	THR
2	A	2330	ASN
2	A	3027	MET
2	A	3028	ARG
2	A	3029	LEU
2	A	3068	GLU
2	A	3076	THR
2	A	3085	ARG
2	A	3089	THR
2	A	3146	VAL
2	A	3150	THR
2	A	3183	LYS
2	A	3189	LEU
2	A	3193	ASN
2	A	3201	VAL
2	A	3202	MET
2	A	3203	PHE
2	A	3213	ASN
2	A	3223	LEU
2	A	3285	GLU
2	A	3314	GLU
2	A	3330	ASN
2	A	4028	ARG
2	A	4029	LEU
2	A	4068	GLU
2	A	4076	THR
2	A	4079	VAL
2	A	4085	ARG
2	A	4089	THR
2	A	4146	VAL
2	A	4183	LYS
2	A	4189	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	A	4193	ASN
2	A	4201	VAL
2	A	4202	MET
2	A	4203	PHE
2	A	4213	ASN
2	A	4223	LEU
2	A	4251	ILE
2	A	4285	GLU
2	A	4314	GLU
2	A	4330	ASN
2	A	5014	LEU
2	A	5027	MET
2	A	5028	ARG
2	A	5029	LEU
2	A	5068	GLU
2	A	5076	THR
2	A	5085	ARG
2	A	5089	THR
2	A	5146	VAL
2	A	5155	ILE
2	A	5164	MET
2	A	5183	LYS
2	A	5189	LEU
2	A	5193	ASN
2	A	5213	ASN
2	A	5223	LEU
2	A	5285	GLU
2	A	5314	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	181	ASN
2	A	193	ASN
2	A	194	GLN
2	A	236	ASN
2	A	304	ASN
2	A	1118	GLN
2	A	1163	HIS
2	A	1173	GLN
2	A	1181	ASN
2	A	1193	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	A	1194	GLN
2	A	1236	ASN
2	A	1304	ASN
2	A	2118	GLN
2	A	2163	HIS
2	A	2173	GLN
2	A	2181	ASN
2	A	2193	ASN
2	A	2194	GLN
2	A	2236	ASN
2	A	2304	ASN
2	A	3097	HIS
2	A	3118	GLN
2	A	3163	HIS
2	A	3173	GLN
2	A	3181	ASN
2	A	3193	ASN
2	A	3194	GLN
2	A	3236	ASN
2	A	3304	ASN
2	A	4118	GLN
2	A	4163	HIS
2	A	4173	GLN
2	A	4181	ASN
2	A	4193	ASN
2	A	4194	GLN
2	A	4236	ASN
2	A	4304	ASN
2	A	5097	HIS
2	A	5173	GLN
2	A	5181	ASN
2	A	5193	ASN
2	A	5194	GLN
2	A	5236	ASN
2	A	5304	ASN

4.3.3 RNA ⓘ

There are no RNA molecules in this entry.

4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

4.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

4.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 6 are monoatomic - leaving 12 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$
4	ALF	A	1501	-	0,4,4	0.00	-	0,6,6	0.00	-
5	ADP	A	1502	3	25,29,29	1.15	2 (8%)	24,45,45	1.79	2 (8%)
4	ALF	A	2501	-	0,4,4	0.00	-	0,6,6	0.00	-
5	ADP	A	2502	3	25,29,29	1.05	1 (4%)	24,45,45	1.67	2 (8%)
4	ALF	A	3501	-	0,4,4	0.00	-	0,6,6	0.00	-
5	ADP	A	3502	3	25,29,29	1.06	1 (4%)	24,45,45	1.83	3 (12%)
4	ALF	A	4501	-	0,4,4	0.00	-	0,6,6	0.00	-
5	ADP	A	4502	3	25,29,29	1.03	2 (8%)	24,45,45	1.84	3 (12%)
4	ALF	A	501	-	0,4,4	0.00	-	0,6,6	0.00	-
5	ADP	A	502	3	25,29,29	1.03	1 (4%)	24,45,45	1.85	4 (16%)
4	ALF	A	5501	-	0,4,4	0.00	-	0,6,6	0.00	-
5	ADP	A	5502	3	25,29,29	1.10	2 (8%)	24,45,45	1.68	2 (8%)

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
4	ALF	A	1501	-	-	0/0/0/0	0/0/0/0
5	ADP	A	1502	3	-	0/12/32/32	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ALF	A	2501	-	-	0/0/0/0	0/0/0/0
5	ADP	A	2502	3	-	0/12/32/32	0/3/3/3
4	ALF	A	3501	-	-	0/0/0/0	0/0/0/0
5	ADP	A	3502	3	-	0/12/32/32	0/3/3/3
4	ALF	A	4501	-	-	0/0/0/0	0/0/0/0
5	ADP	A	4502	3	-	0/12/32/32	0/3/3/3
4	ALF	A	501	-	-	0/0/0/0	0/0/0/0
5	ADP	A	502	3	-	0/12/32/32	0/3/3/3
4	ALF	A	5501	-	-	0/0/0/0	0/0/0/0
5	ADP	A	5502	3	-	0/12/32/32	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	4502	ADP	C2-N3	2.03	1.35	1.32
5	A	5502	ADP	C2-N3	2.24	1.35	1.32
5	A	1502	ADP	C2-N3	2.27	1.36	1.32
5	A	502	ADP	C5-C4	3.10	1.47	1.40
5	A	2502	ADP	C5-C4	3.12	1.47	1.40
5	A	4502	ADP	C5-C4	3.13	1.47	1.40
5	A	5502	ADP	C5-C4	3.29	1.47	1.40
5	A	3502	ADP	C5-C4	3.30	1.47	1.40
5	A	1502	ADP	C5-C4	3.44	1.48	1.40

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	502	ADP	N3-C2-N1	-7.24	122.56	128.86
5	A	1502	ADP	N3-C2-N1	-7.14	122.64	128.86
5	A	4502	ADP	N3-C2-N1	-6.99	122.77	128.86
5	A	3502	ADP	N3-C2-N1	-6.92	122.83	128.86
5	A	2502	ADP	N3-C2-N1	-6.32	123.36	128.86
5	A	5502	ADP	N3-C2-N1	-5.78	123.83	128.86
5	A	5502	ADP	C4-C5-N7	-3.26	106.26	109.41
5	A	4502	ADP	C4-C5-N7	-3.23	106.29	109.41
5	A	3502	ADP	C4-C5-N7	-2.64	106.86	109.41
5	A	1502	ADP	C4-C5-N7	-2.63	106.87	109.41
5	A	2502	ADP	C4-C5-N7	-2.56	106.93	109.41
5	A	4502	ADP	O2B-PB-O1B	2.02	118.42	110.50
5	A	502	ADP	C2-N1-C6	2.09	122.42	118.77
5	A	3502	ADP	C2-N1-C6	2.10	122.45	118.77
5	A	502	ADP	O3B-PB-O2B	2.38	117.22	107.61

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
5	A	502	ADP	N6-C6-N1	2.48	123.69	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 31 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1501	ALF	1	0
5	A	1502	ADP	5	0
4	A	2501	ALF	1	0
5	A	2502	ADP	7	0
5	A	3502	ADP	5	0
4	A	4501	ALF	1	0
5	A	4502	ADP	4	0
4	A	501	ALF	3	0
4	A	5501	ALF	2	0
5	A	5502	ADP	3	0

4.7 Other polymers [i](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

5 Fit of model and data [i](#)

5.1 Protein, DNA and RNA chains [i](#)

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	B	15/18 (83%)	0.27	0 100 100	118, 147, 162, 174	0
2	A	1937/2050 (94%)	-0.27	16 (0%) 86 79	147, 196, 220, 220	0
All	All	1952/2068 (94%)	-0.26	16 (0%) 86 79	118, 195, 220, 220	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	1200	GLY	2.9
2	A	5205	ASN	2.8
2	A	5001	ALA	2.8
2	A	295	GLY	2.5
2	A	3283	LEU	2.5
2	A	3144	ASP	2.4
2	A	3193	ASN	2.4
2	A	4001	ALA	2.4
2	A	229	GLY	2.3
2	A	1063	GLU	2.3
2	A	4195	ILE	2.3
2	A	2227	ARG	2.1
2	A	2144	ASP	2.1
2	A	228	ILE	2.0
2	A	4063	GLU	2.0
2	A	5005	ASN	2.0

5.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.4 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
4	ALF	A	3501	5/5	0.98	0.52	0.89	181,181,182,183	0
5	ADP	A	502	27/27	0.93	0.37	0.62	169,190,193,193	0
4	ALF	A	1501	5/5	0.95	0.33	0.37	190,194,195,196	0
4	ALF	A	501	5/5	0.97	0.39	0.18	165,168,169,169	0
5	ADP	A	1502	27/27	0.92	0.25	-0.12	191,198,201,202	0
5	ADP	A	3502	27/27	0.91	0.29	-0.21	182,190,197,197	0
4	ALF	A	4501	5/5	0.99	0.26	-0.33	159,160,161,162	0
5	ADP	A	2502	27/27	0.93	0.20	-0.57	176,188,192,193	0
5	ADP	A	4502	27/27	0.97	0.18	-0.76	157,178,183,183	0
4	ALF	A	5501	5/5	0.95	0.17	-0.95	167,172,172,172	0
4	ALF	A	2501	5/5	0.94	0.23	-0.96	177,177,179,180	0
5	ADP	A	5502	27/27	0.96	0.13	-1.26	173,185,193,193	0
3	MG	A	4500	1/1	0.99	0.35	-	162,162,162,162	0
3	MG	A	1500	1/1	0.99	0.47	-	186,186,186,186	0
3	MG	A	3500	1/1	0.99	0.60	-	180,180,180,180	0
3	MG	A	500	1/1	0.98	0.55	-	162,162,162,162	0
3	MG	A	2500	1/1	0.97	0.32	-	177,177,177,177	0
3	MG	A	5500	1/1	0.98	0.10	-	53,53,53,53	0

5.5 Other polymers [i](#)

There are no such residues in this entry.