



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 11:43 PM GMT

PDB ID : 3I3Y
Title : Crystal structure of Ribokinase in Complex with D-Ribose from *Klebsiella pneumoniae*
Authors : Satyanarayana, L.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-07-01
Resolution : 2.15 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

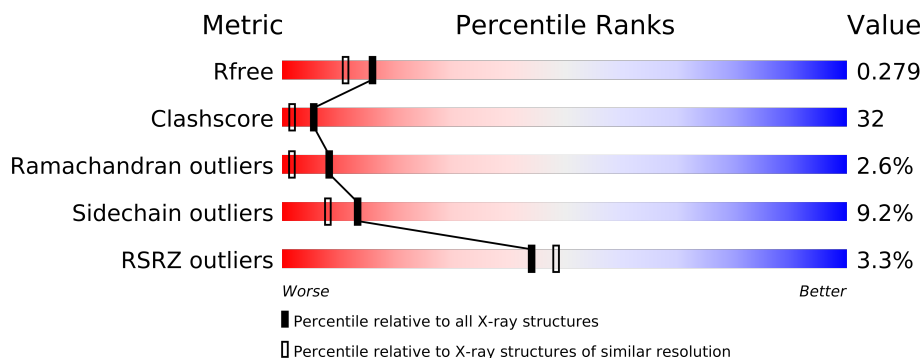
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.15 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1094 (2.18-2.14)
Clashscore	79885	1299 (2.18-2.14)
Ramachandran outliers	78287	1272 (2.18-2.14)
Sidechain outliers	78261	1272 (2.18-2.14)
RSRZ outliers	66119	1094 (2.18-2.14)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	299	
1	B	299	
1	C	299	
1	D	299	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	SO4	C	1	-	X
3	GOL	C	501	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8707 atoms, of which 0 are hydrogen and 0 are deuterium.

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 protein called Carbohydrate kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	Se	0	0	0
			2121	1333	372	408	3	5			
1	B	282	Total	C	N	O	S	Se	0	0	0
			2099	1322	368	401	3	5			
1	C	285	Total	C	N	O	S	Se	0	0	0
			2121	1333	372	408	3	5			
1	D	285	Total	C	N	O	S	Se	0	0	0
			2121	1333	372	408	3	5			

There are 44 discrepancies between the modelled and reference sequences:

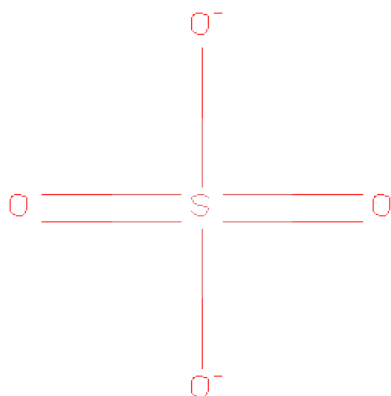
Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	MSE	-	EXPRESSION TAG	UNP A6T989
A	-2	SER	-	EXPRESSION TAG	UNP A6T989
A	-1	LEU	-	EXPRESSION TAG	UNP A6T989
A	290	GLU	-	EXPRESSION TAG	UNP A6T989
A	291	GLY	-	EXPRESSION TAG	UNP A6T989
A	292	HIS	-	EXPRESSION TAG	UNP A6T989
A	293	HIS	-	EXPRESSION TAG	UNP A6T989
A	294	HIS	-	EXPRESSION TAG	UNP A6T989
A	295	HIS	-	EXPRESSION TAG	UNP A6T989
A	296	HIS	-	EXPRESSION TAG	UNP A6T989
A	297	HIS	-	EXPRESSION TAG	UNP A6T989
B	-3	MSE	-	EXPRESSION TAG	UNP A6T989
B	-2	SER	-	EXPRESSION TAG	UNP A6T989
B	-1	LEU	-	EXPRESSION TAG	UNP A6T989
B	290	GLU	-	EXPRESSION TAG	UNP A6T989
B	291	GLY	-	EXPRESSION TAG	UNP A6T989
B	292	HIS	-	EXPRESSION TAG	UNP A6T989
B	293	HIS	-	EXPRESSION TAG	UNP A6T989
B	294	HIS	-	EXPRESSION TAG	UNP A6T989
B	295	HIS	-	EXPRESSION TAG	UNP A6T989
B	296	HIS	-	EXPRESSION TAG	UNP A6T989

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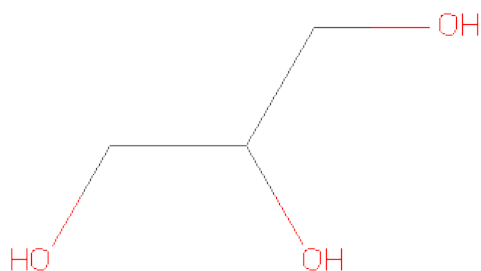
Chain	Residue	Modelled	Actual	Comment	Reference
B	297	HIS	-	EXPRESSION TAG	UNP A6T989
C	-3	MSE	-	EXPRESSION TAG	UNP A6T989
C	-2	SER	-	EXPRESSION TAG	UNP A6T989
C	-1	LEU	-	EXPRESSION TAG	UNP A6T989
C	290	GLU	-	EXPRESSION TAG	UNP A6T989
C	291	GLY	-	EXPRESSION TAG	UNP A6T989
C	292	HIS	-	EXPRESSION TAG	UNP A6T989
C	293	HIS	-	EXPRESSION TAG	UNP A6T989
C	294	HIS	-	EXPRESSION TAG	UNP A6T989
C	295	HIS	-	EXPRESSION TAG	UNP A6T989
C	296	HIS	-	EXPRESSION TAG	UNP A6T989
C	297	HIS	-	EXPRESSION TAG	UNP A6T989
D	-3	MSE	-	EXPRESSION TAG	UNP A6T989
D	-2	SER	-	EXPRESSION TAG	UNP A6T989
D	-1	LEU	-	EXPRESSION TAG	UNP A6T989
D	290	GLU	-	EXPRESSION TAG	UNP A6T989
D	291	GLY	-	EXPRESSION TAG	UNP A6T989
D	292	HIS	-	EXPRESSION TAG	UNP A6T989
D	293	HIS	-	EXPRESSION TAG	UNP A6T989
D	294	HIS	-	EXPRESSION TAG	UNP A6T989
D	295	HIS	-	EXPRESSION TAG	UNP A6T989
D	296	HIS	-	EXPRESSION TAG	UNP A6T989
D	297	HIS	-	EXPRESSION TAG	UNP A6T989

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



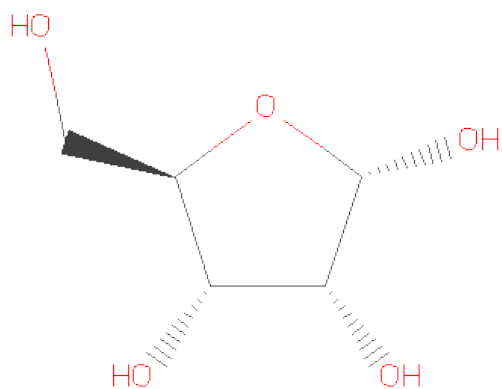
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is SUGAR (RIBOSE) (three-letter code: RIB) (formula: $C_5H_{10}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			10	5	5		

- Molecule 5 is water.

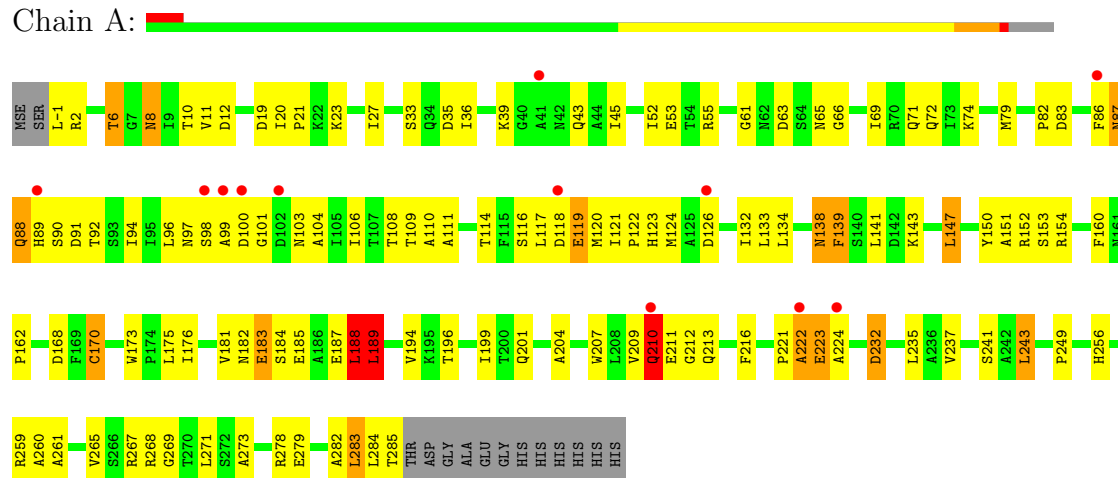
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	60	Total	O	0	0
			60	60		
5	B	41	Total	O	0	0
			41	41		
5	C	60	Total	O	0	0
			60	60		
5	D	38	Total	O	0	0
			38	38		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

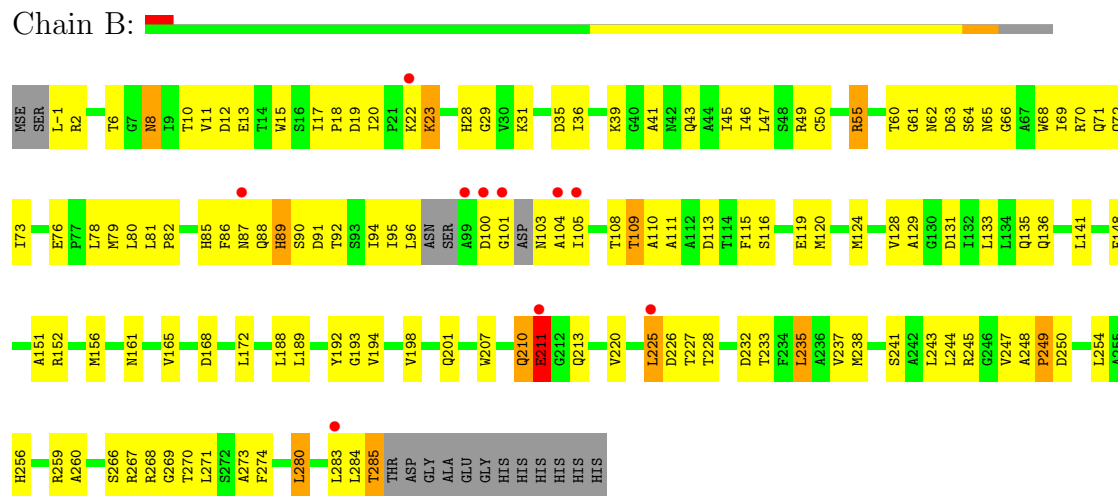
• Molecule 1: Carbohydrate kinase

Chain A:



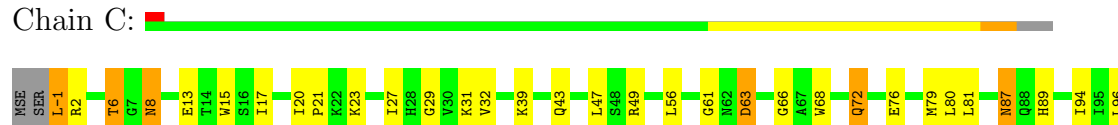
• Molecule 1: Carbohydrate kinase

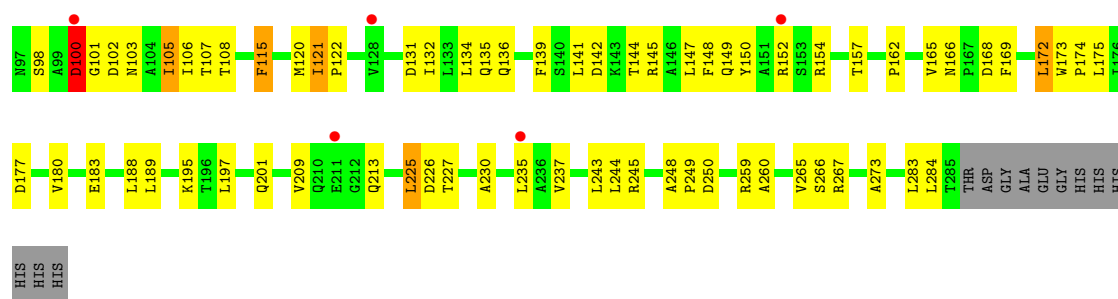
Chain B:



• Molecule 1: Carbohydrate kinase

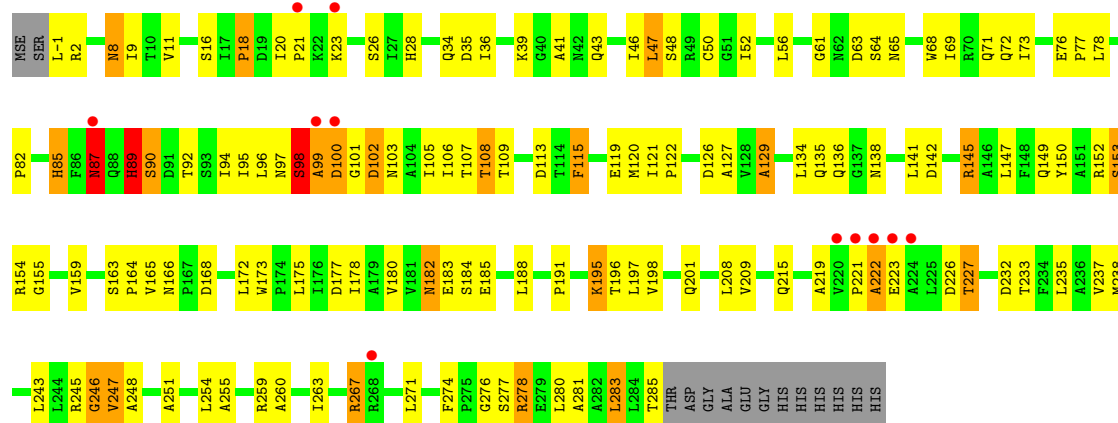
Chain C:





● Molecule 1: Carbohydrate kinase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.49Å 105.90Å 121.15Å 90.00° 99.45° 90.00°	Depositor
Resolution (Å)	34.97 – 2.15 34.97 – 2.05	Depositor EDS
% Data completeness (in resolution range)	91.3 (34.97-2.15) 91.3 (34.97-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.50 (at 2.05Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.244 , 0.276 0.251 , 0.279	Depositor DCC
R_{free} test set	2279 reflections (3.23%)	DCC
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.350	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 165981 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	8707	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, RIB, SO4

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	A	0.39	0/2155	0.72	0/2931
1	B	0.36	0/2131	0.66	0/2895
1	C	0.36	0/2155	0.68	0/2931
1	D	0.36	0/2155	0.69	0/2931
All	All	0.37	0/8596	0.69	0/11688

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2121	0	2122	165	0
1	B	2099	0	2105	152	0
1	C	2121	0	2122	100	0
1	D	2121	0	2122	163	0
2	B	15	0	0	0	0
2	C	10	0	0	0	0
2	D	5	0	0	0	0
3	C	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	10	0	10	0	0
5	A	60	0	0	12	0
5	B	41	0	0	7	0
5	C	60	0	0	1	0
5	D	38	0	0	6	0
All	All	8707	0	8489	544	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 32.

All (544) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:61:GLY:H	1:A:89:HIS:CD2	1.66	1.13
1:A:10:THR:HG23	1:A:91:ASP:HA	1.34	1.10
1:D:195:LYS:H	1:D:195:LYS:HD2	0.95	1.09
1:A:181:VAL:HG13	1:A:185:GLU:HB2	1.42	1.01
1:D:259:ARG:HB3	1:D:283:LEU:HD11	1.45	0.98
1:A:61:GLY:N	1:A:89:HIS:HD2	1.63	0.95
1:C:6:THR:HG21	1:C:147:LEU:HD13	1.43	0.95
1:A:6:THR:HG23	1:A:120:MSE:HE1	1.49	0.93
1:A:61:GLY:H	1:A:89:HIS:HD2	0.93	0.93
1:D:195:LYS:HD2	1:D:195:LYS:N	1.81	0.93
1:B:70:ARG:HD3	1:B:85:HIS:NE2	1.86	0.89
1:B:63:ASP:H	1:B:89:HIS:HE1	1.22	0.87
1:A:188:LEU:O	1:A:189:LEU:HB2	1.71	0.86
1:B:228:THR:HG21	1:B:270:THR:HG22	1.57	0.86
1:A:45:ILE:HD13	1:A:271:LEU:HD11	1.56	0.86
1:A:19:ASP:HA	1:A:97:ASN:HD22	1.40	0.86
1:C:20:ILE:O	1:C:20:ILE:HD12	1.78	0.83
1:A:100:ASP:HB3	1:C:23:LYS:NZ	1.93	0.83
1:D:43:GLN:NE2	1:D:136:GLN:HE21	1.77	0.83
1:D:195:LYS:CD	1:D:195:LYS:H	1.80	0.83
1:B:63:ASP:H	1:B:89:HIS:CE1	1.97	0.83
1:B:63:ASP:HB3	1:B:89:HIS:CE1	2.13	0.82
1:C:79:MSE:HE2	1:C:81:LEU:HD21	1.59	0.82
1:C:166:ASN:HD21	1:C:168:ASP:HB2	1.43	0.82
1:A:20:ILE:HG13	1:A:97:ASN:HD21	1.46	0.80
1:D:177:ASP:HA	1:D:195:LYS:HD3	1.62	0.80
1:A:209:VAL:O	1:A:210:GLN:HB2	1.79	0.80
1:D:141:LEU:HD11	1:D:172:LEU:HD11	1.65	0.79
1:A:66:GLY:CA	1:A:89:HIS:NE2	2.47	0.78
1:D:278:ARG:HG2	5:D:325:HOH:O	1.83	0.78

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:8:ASN:H	1:B:8:ASN:HD22	1.31	0.77
1:C:-1:LEU:HD12	1:C:2:ARG:N	2.00	0.77
1:C:152:ARG:CZ	1:C:175:LEU:HA	2.14	0.77
1:B:63:ASP:N	1:B:89:HIS:HE1	1.82	0.76
1:D:11:VAL:HG21	1:D:65:ASN:HB3	1.65	0.76
1:A:194:VAL:O	1:A:210:GLN:HG2	1.86	0.76
1:C:6:THR:HG23	1:C:139:PHE:HE2	1.49	0.76
1:D:145:ARG:HH11	1:D:145:ARG:CB	2.00	0.75
1:D:16:SER:HA	1:D:97:ASN:HB3	1.69	0.74
1:B:79:MSE:HE2	1:B:81:LEU:HD21	1.69	0.74
1:B:220:VAL:HG21	1:B:259:ARG:HG3	1.69	0.74
1:B:65:ASN:HB2	1:B:89:HIS:CD2	2.23	0.73
1:C:152:ARG:NH1	1:C:175:LEU:HA	2.03	0.73
1:A:-1:LEU:N	5:A:349:HOH:O	2.20	0.73
1:D:43:GLN:HE22	1:D:136:GLN:HE21	1.37	0.73
1:A:117:LEU:H	1:A:143:LYS:HZ2	1.37	0.73
1:B:39:LYS:HE2	1:B:232:ASP:OD1	1.86	0.73
1:B:23:LYS:HD3	1:D:101:GLY:HA3	1.71	0.73
1:B:141:LEU:HD11	1:B:172:LEU:HD11	1.71	0.72
1:B:68:TRP:O	1:B:72:GLN:HG2	1.89	0.72
1:C:68:TRP:HE1	1:C:72:GLN:NE2	1.88	0.72
1:B:124:MSE:HE3	1:B:156:MSE:HE1	1.72	0.71
1:B:63:ASP:HB3	1:B:89:HIS:NE2	2.06	0.71
1:B:28:HIS:CE1	1:D:138:ASN:HD22	2.09	0.70
1:B:23:LYS:NZ	1:D:101:GLY:HA3	2.06	0.70
1:D:238:MSE:HG3	1:D:254:LEU:HD12	1.72	0.70
1:A:65:ASN:CB	1:A:89:HIS:HE1	2.04	0.70
1:A:282:ALA:O	1:A:285:THR:HG22	1.92	0.70
1:B:210:GLN:HB2	5:B:307:HOH:O	1.91	0.70
1:A:117:LEU:C	1:A:119:GLU:H	1.94	0.70
1:D:71:GLN:NE2	1:D:72:GLN:HE21	1.90	0.70
1:D:141:LEU:HD22	1:D:168:ASP:HB3	1.74	0.69
1:D:82:PRO:HD3	1:D:115:PHE:CZ	2.26	0.69
1:A:224:ALA:HA	1:A:265:VAL:HG12	1.74	0.69
1:C:68:TRP:HE1	1:C:72:GLN:HE22	1.40	0.69
1:B:285:THR:HG22	5:B:328:HOH:O	1.92	0.69
1:B:95:ILE:HG23	1:B:104:ALA:HB3	1.75	0.68
1:B:70:ARG:HD3	1:B:85:HIS:CE1	2.28	0.68
1:A:196:THR:HG21	1:A:249:PRO:HG2	1.75	0.68
1:A:121:ILE:HD12	1:A:124:MSE:HE3	1.74	0.68
1:D:134:LEU:HA	1:D:159:VAL:HG13	1.74	0.68
1:C:49:ARG:HD2	1:C:76:GLU:OE1	1.93	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:105:ILE:C	1:C:105:ILE:HD12	2.14	0.67
1:C:225:LEU:HD21	1:C:266:SER:C	2.14	0.67
1:A:241:SER:OG	1:A:256:HIS:HD2	1.76	0.67
1:A:100:ASP:HB3	1:C:23:LYS:HZ2	1.58	0.67
1:A:36:ILE:HD11	1:A:72:GLN:HB2	1.74	0.67
1:A:183:GLU:CD	1:A:201:GLN:HG2	2.15	0.66
1:D:109:THR:HG23	1:D:113:ASP:OD2	1.95	0.66
1:B:63:ASP:CG	1:B:64:SER:H	1.99	0.66
1:D:121:ILE:HG13	1:D:122:PRO:CD	2.26	0.66
1:B:94:ILE:HD12	1:B:94:ILE:N	2.11	0.66
1:B:60:THR:HG23	1:B:85:HIS:HD1	1.60	0.65
1:A:66:GLY:N	1:A:89:HIS:NE2	2.43	0.65
1:A:117:LEU:HD22	1:A:143:LYS:NZ	2.12	0.65
1:A:210:GLN:O	1:A:211:GLU:HB2	1.97	0.65
1:D:99:ALA:O	1:D:100:ASP:HB3	1.96	0.65
1:D:61:GLY:HA3	1:D:89:HIS:O	1.97	0.65
1:C:17:ILE:O	1:C:98:SER:HA	1.97	0.65
1:B:100:ASP:HB3	1:D:23:LYS:NZ	2.12	0.65
1:B:247:VAL:O	1:B:248:ALA:HB3	1.96	0.64
1:B:19:ASP:HB3	1:B:100:ASP:OD1	1.97	0.64
1:C:154:ARG:HH11	1:C:154:ARG:HG2	1.62	0.64
1:B:70:ARG:CD	1:B:85:HIS:NE2	2.60	0.64
1:B:20:ILE:HD13	1:D:21:PRO:O	1.96	0.64
1:A:283:LEU:HD12	1:A:283:LEU:C	2.18	0.64
1:A:108:THR:HG23	1:C:15:TRP:CZ3	2.33	0.64
1:B:109:THR:HG22	1:D:28:HIS:HE1	1.63	0.64
1:A:23:LYS:NZ	1:C:100:ASP:HB2	2.13	0.63
1:A:43:GLN:HG3	1:A:232:ASP:HB2	1.80	0.63
1:C:225:LEU:CD2	1:C:225:LEU:H	2.12	0.63
1:A:100:ASP:CB	1:C:23:LYS:NZ	2.62	0.63
1:A:187:GLU:O	1:A:188:LEU:HG	1.98	0.63
1:B:11:VAL:HG11	1:B:65:ASN:CG	2.19	0.63
1:C:6:THR:O	1:C:6:THR:HG22	1.97	0.63
1:B:63:ASP:CG	1:B:64:SER:N	2.52	0.63
1:C:43:GLN:HE21	1:C:134:LEU:HD21	1.63	0.63
1:D:39:LYS:HB3	1:D:136:GLN:HG2	1.81	0.63
1:B:28:HIS:HE1	1:D:138:ASN:HD22	1.46	0.62
1:B:63:ASP:CA	1:B:89:HIS:HE1	2.11	0.62
1:A:139:PHE:HB2	5:A:314:HOH:O	1.98	0.62
1:C:115:PHE:HB3	1:C:120:MSE:HE3	1.82	0.62
1:A:61:GLY:N	1:A:89:HIS:CD2	2.47	0.62
1:A:6:THR:HG23	1:A:120:MSE:CE	2.27	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:36:ILE:HD11	1:B:72:GLN:HB2	1.81	0.62
1:B:109:THR:O	1:B:111:ALA:N	2.31	0.61
1:D:152:ARG:O	1:D:153:SER:HB3	1.98	0.61
1:B:165:VAL:HG21	1:B:188:LEU:HD23	1.82	0.61
1:B:60:THR:HG23	1:B:85:HIS:ND1	2.14	0.61
1:D:115:PHE:HB3	1:D:120:MSE:HE3	1.80	0.61
1:D:152:ARG:O	1:D:153:SER:CB	2.46	0.61
1:D:237:VAL:HG21	1:D:260:ALA:CB	2.29	0.61
1:B:124:MSE:CE	1:B:156:MSE:SE	2.99	0.61
1:D:247:VAL:HG22	1:D:248:ALA:H	1.64	0.61
1:D:121:ILE:HG13	1:D:122:PRO:N	2.15	0.61
1:A:104:ALA:HB1	1:C:21:PRO:HB2	1.83	0.61
1:A:121:ILE:HD11	1:A:150:TYR:CE1	2.36	0.61
1:B:66:GLY:H	1:B:89:HIS:CE1	2.18	0.61
1:B:17:ILE:HB	1:B:18:PRO:HD2	1.81	0.61
1:B:225:LEU:HD21	1:B:266:SER:O	2.01	0.60
1:D:8:ASN:HD22	1:D:8:ASN:H	1.47	0.60
1:A:138:ASN:O	1:A:139:PHE:CG	2.54	0.60
1:A:66:GLY:HA2	1:A:89:HIS:NE2	2.15	0.60
1:B:8:ASN:H	1:B:8:ASN:ND2	1.98	0.60
1:D:36:ILE:CD1	1:D:72:GLN:HG3	2.31	0.60
1:A:117:LEU:O	1:A:119:GLU:N	2.32	0.60
1:C:68:TRP:NE1	1:C:72:GLN:NE2	2.50	0.60
1:A:65:ASN:HB2	1:A:89:HIS:HE1	1.65	0.59
1:D:43:GLN:NE2	1:D:136:GLN:NE2	2.50	0.59
1:B:105:ILE:HD12	1:B:105:ILE:N	2.18	0.59
1:B:241:SER:OG	1:B:256:HIS:HD2	1.85	0.59
1:B:244:LEU:HD12	1:B:284:LEU:HD13	1.83	0.59
1:A:43:GLN:HG3	1:A:232:ASP:CB	2.32	0.59
1:C:141:LEU:HD22	1:C:168:ASP:HB3	1.84	0.59
1:B:17:ILE:HG22	1:B:29:GLY:HA3	1.84	0.59
1:C:-1:LEU:CD1	1:C:131:ASP:HA	2.32	0.58
1:B:28:HIS:NE2	1:D:164:PRO:HG3	2.18	0.58
1:A:12:ASP:OD1	1:A:92:THR:HG22	2.03	0.58
1:C:21:PRO:HG3	1:C:27:ILE:CD1	2.32	0.58
1:C:165:VAL:HG23	1:C:189:LEU:HD11	1.84	0.58
1:A:82:PRO:HB3	1:A:123:HIS:NE2	2.18	0.58
1:A:86:PHE:C	1:A:87:ASN:OD1	2.41	0.58
1:D:149:GLN:O	1:D:152:ARG:O	2.21	0.58
1:D:121:ILE:HG13	1:D:122:PRO:HD3	1.84	0.58
1:C:39:LYS:HB3	1:C:136:GLN:HG2	1.84	0.58
1:D:247:VAL:HG22	1:D:248:ALA:N	2.18	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:135:GLN:HE22	1:D:147:LEU:HD12	1.68	0.58
1:A:10:THR:HG22	1:A:11:VAL:N	2.17	0.58
1:B:201:GLN:HE21	1:B:207:TRP:HD1	1.52	0.58
1:D:87:ASN:N	1:D:87:ASN:HD22	2.01	0.58
1:A:10:THR:CG2	1:A:11:VAL:N	2.67	0.57
1:B:15:TRP:CE2	1:D:94:ILE:HD11	2.40	0.57
1:C:6:THR:HG22	1:C:135:GLN:HG2	1.85	0.57
1:A:222:ALA:O	1:A:223:GLU:HB2	2.05	0.57
1:A:100:ASP:CB	1:C:23:LYS:HZ3	2.17	0.57
1:D:9:ILE:HG21	1:D:69:ILE:HG21	1.86	0.57
1:C:235:LEU:HD12	1:C:235:LEU:O	2.04	0.57
1:A:117:LEU:HD22	1:A:143:LYS:HZ2	1.70	0.57
1:B:109:THR:CG2	1:D:28:HIS:HE1	2.17	0.57
1:C:115:PHE:O	1:C:120:MSE:HE3	2.04	0.57
1:A:121:ILE:HD11	1:A:150:TYR:CZ	2.40	0.57
1:A:106:ILE:HD11	1:C:21:PRO:HG2	1.86	0.57
1:A:222:ALA:O	1:A:223:GLU:CB	2.53	0.57
1:A:20:ILE:HG13	1:A:97:ASN:ND2	2.18	0.56
1:D:43:GLN:O	1:D:47:LEU:HB2	2.05	0.56
1:B:141:LEU:HD11	1:B:172:LEU:CD1	2.35	0.56
1:A:2:ARG:HD2	5:A:308:HOH:O	2.04	0.56
1:C:87:ASN:H	1:C:87:ASN:HD22	1.53	0.56
1:D:2:ARG:HH11	1:D:2:ARG:HG2	1.70	0.56
1:D:36:ILE:HD11	1:D:72:GLN:HG3	1.87	0.56
1:B:8:ASN:N	1:B:8:ASN:HD22	1.94	0.56
1:A:173:TRP:CE3	1:A:176:ILE:HD12	2.40	0.56
1:A:261:ALA:O	1:A:265:VAL:HG23	2.04	0.56
1:A:182:ASN:OD1	1:A:185:GLU:HG3	2.05	0.56
1:A:121:ILE:HD11	1:A:150:TYR:CD1	2.40	0.56
1:B:23:LYS:HD3	1:D:102:ASP:H	1.69	0.56
1:B:95:ILE:HG22	1:B:104:ALA:O	2.06	0.56
1:C:180:VAL:HG21	1:C:235:LEU:HD13	1.88	0.56
1:B:65:ASN:O	1:B:69:ILE:HG13	2.06	0.56
1:C:8:ASN:H	1:C:8:ASN:HD22	1.53	0.56
1:B:22:LYS:HB2	1:B:22:LYS:NZ	2.20	0.56
1:A:237:VAL:HG21	1:A:260:ALA:CB	2.36	0.56
1:B:104:ALA:C	1:B:105:ILE:HD12	2.26	0.55
1:D:141:LEU:HD11	1:D:172:LEU:CD1	2.35	0.55
1:D:98:SER:HB2	1:D:100:ASP:H	1.71	0.55
1:A:181:VAL:HG12	1:A:182:ASN:O	2.06	0.55
1:D:99:ALA:O	1:D:100:ASP:CB	2.54	0.55
1:B:210:GLN:HA	1:B:210:GLN:HE21	1.72	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:165:VAL:HG21	1:D:188:LEU:HD23	1.89	0.55
1:D:105:ILE:HG22	1:D:106:ILE:N	2.22	0.55
1:D:255:ALA:O	1:D:259:ARG:HD3	2.07	0.55
1:B:61:GLY:O	1:B:85:HIS:ND1	2.40	0.55
1:B:17:ILE:HD13	1:B:95:ILE:HD11	1.88	0.55
1:A:11:VAL:HG23	1:A:69:ILE:HD11	1.89	0.54
1:C:166:ASN:ND2	1:C:168:ASP:HB2	2.17	0.54
1:B:100:ASP:C	1:D:23:LYS:HZ2	2.10	0.54
1:A:8:ASN:H	1:A:8:ASN:HD22	1.55	0.54
1:B:63:ASP:HB3	1:B:89:HIS:HE2	1.72	0.54
1:D:115:PHE:CG	1:D:120:MSE:HE3	2.42	0.54
1:D:127:ALA:O	1:D:154:ARG:NH2	2.41	0.54
1:B:63:ASP:CB	1:B:89:HIS:CE1	2.87	0.54
1:A:168:ASP:C	1:A:170:CYS:N	2.61	0.54
1:A:23:LYS:O	1:C:103:ASN:O	2.26	0.54
1:D:245:ARG:O	1:D:246:GLY:C	2.46	0.54
1:B:245:ARG:NH2	1:B:250:ASP:OD2	2.40	0.54
1:A:221:PRO:O	1:A:222:ALA:HB2	2.08	0.54
1:D:182:ASN:N	1:D:182:ASN:ND2	2.54	0.54
1:B:105:ILE:O	1:D:26:SER:HA	2.08	0.53
1:A:222:ALA:O	1:A:223:GLU:HG3	2.09	0.53
1:D:150:TYR:CZ	1:D:154:ARG:HD3	2.43	0.53
1:A:184:SER:O	1:A:187:GLU:O	2.26	0.53
1:D:152:ARG:HG2	1:D:152:ARG:HH11	1.72	0.53
1:B:226:ASP:HB3	1:B:267:ARG:O	2.08	0.53
1:A:188:LEU:O	1:A:189:LEU:CB	2.50	0.53
1:B:100:ASP:HB3	1:D:23:LYS:HZ2	1.72	0.53
1:B:136:GLN:HA	1:B:161:ASN:O	2.08	0.53
1:B:270:THR:OG1	1:B:271:LEU:N	2.42	0.53
1:D:227:THR:HG21	5:D:323:HOH:O	2.09	0.53
1:A:121:ILE:HG23	1:A:122:PRO:HD3	1.88	0.53
1:B:211:GLU:H	1:B:211:GLU:CD	2.10	0.53
1:D:105:ILE:N	1:D:105:ILE:HD12	2.24	0.53
1:A:52:ILE:HD12	1:A:243:LEU:HD23	1.91	0.53
1:A:152:ARG:NH1	1:A:175:LEU:HA	2.23	0.52
1:A:2:ARG:HG2	1:A:53:GLU:HB3	1.91	0.52
1:A:87:ASN:N	1:A:87:ASN:OD1	2.43	0.52
1:C:6:THR:O	1:C:6:THR:CG2	2.56	0.52
1:B:66:GLY:N	1:B:89:HIS:ND1	2.57	0.52
1:A:6:THR:HG22	1:A:6:THR:O	2.08	0.52
1:B:20:ILE:HG21	1:D:20:ILE:HG22	1.92	0.52
1:A:65:ASN:O	1:A:69:ILE:HG13	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:117:LEU:C	1:A:119:GLU:N	2.62	0.52
1:B:65:ASN:HB2	1:B:89:HIS:CG	2.45	0.52
1:C:43:GLN:NE2	1:C:136:GLN:HE21	2.08	0.52
1:B:41:ALA:O	1:B:45:ILE:HG13	2.10	0.51
1:D:107:THR:HG22	1:D:108:THR:O	2.09	0.51
1:B:271:LEU:HA	1:B:274:PHE:CD2	2.45	0.51
1:D:182:ASN:ND2	1:D:185:GLU:OE1	2.44	0.51
1:A:98:SER:O	1:A:99:ALA:C	2.49	0.51
1:C:6:THR:CG2	1:C:139:PHE:HE2	2.22	0.51
1:A:20:ILE:HD11	5:A:313:HOH:O	2.11	0.51
1:B:23:LYS:CD	1:D:101:GLY:HA3	2.39	0.51
1:B:95:ILE:CG2	1:B:104:ALA:HB3	2.38	0.51
1:A:39:LYS:HE3	5:A:355:HOH:O	2.10	0.51
1:D:2:ARG:NH1	5:D:304:HOH:O	2.42	0.51
1:A:141:LEU:HD22	1:A:168:ASP:HB3	1.91	0.51
1:B:116:SER:HB2	1:B:119:GLU:OE2	2.10	0.51
1:B:124:MSE:HE2	1:B:156:MSE:SE	2.60	0.51
1:C:244:LEU:CD1	1:C:284:LEU:HD13	2.41	0.51
1:D:87:ASN:N	1:D:87:ASN:ND2	2.58	0.51
1:A:117:LEU:CD2	1:A:143:LYS:HD3	2.41	0.51
1:B:95:ILE:O	1:B:95:ILE:HG23	2.10	0.51
1:B:101:GLY:O	1:B:103:ASN:ND2	2.44	0.51
1:C:166:ASN:HB3	1:C:169:PHE:CD1	2.46	0.51
1:B:201:GLN:NE2	1:B:207:TRP:CD1	2.79	0.51
1:B:109:THR:HG22	1:D:28:HIS:CE1	2.44	0.51
1:A:117:LEU:HA	5:A:336:HOH:O	2.10	0.51
1:D:82:PRO:HD3	1:D:115:PHE:HZ	1.73	0.51
1:D:2:ARG:NH2	1:D:126:ASP:O	2.44	0.51
1:D:87:ASN:ND2	5:D:318:HOH:O	2.44	0.50
1:B:6:THR:OG1	1:B:135:GLN:HG2	2.10	0.50
1:D:173:TRP:CZ2	1:D:191:PRO:HG3	2.47	0.50
1:A:117:LEU:O	1:A:118:ASP:HB2	2.11	0.50
1:A:121:ILE:HB	5:A:336:HOH:O	2.10	0.50
1:C:79:MSE:CE	1:C:81:LEU:HD21	2.35	0.50
1:D:46:ILE:CD1	1:D:233:THR:HA	2.40	0.50
1:A:114:THR:O	1:A:114:THR:HG22	2.10	0.50
1:D:129:ALA:O	1:D:155:GLY:O	2.29	0.50
1:C:177:ASP:HA	1:C:195:LYS:HZ2	1.77	0.50
1:C:188:LEU:O	1:C:188:LEU:HD23	2.11	0.50
1:A:183:GLU:OE1	1:A:201:GLN:HG2	2.11	0.50
1:A:168:ASP:C	1:A:170:CYS:H	2.15	0.50
1:A:116:SER:C	1:A:117:LEU:O	2.49	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:180:VAL:HG21	1:D:235:LEU:CD2	2.41	0.50
1:C:32:VAL:O	1:C:32:VAL:HG22	2.12	0.50
1:A:108:THR:HG22	1:C:29:GLY:O	2.12	0.50
1:A:279:GLU:HG3	5:B:329:HOH:O	2.12	0.49
1:D:9:ILE:HG21	1:D:69:ILE:CG2	2.41	0.49
1:C:154:ARG:NH1	1:C:154:ARG:HG2	2.27	0.49
1:B:192:TYR:O	1:B:194:VAL:N	2.42	0.49
1:A:133:LEU:HD22	1:A:151:ALA:HB2	1.93	0.49
1:A:20:ILE:N	1:A:97:ASN:ND2	2.60	0.49
1:A:268:ARG:HG2	1:A:269:GLY:N	2.26	0.49
1:A:278:ARG:NH2	1:B:131:ASP:OD1	2.44	0.49
1:A:121:ILE:CG2	1:A:122:PRO:HD3	2.42	0.49
1:B:10:THR:HB	1:B:91:ASP:HA	1.94	0.49
1:B:11:VAL:CG1	1:B:91:ASP:HB3	2.43	0.49
1:B:135:GLN:HB2	1:B:148:PHE:CZ	2.47	0.49
1:D:267:ARG:HH11	1:D:267:ARG:HG2	1.78	0.49
1:D:177:ASP:HA	1:D:195:LYS:CD	2.39	0.49
1:B:245:ARG:O	1:B:247:VAL:HG13	2.13	0.49
1:D:52:ILE:HD12	1:D:52:ILE:N	2.28	0.49
1:C:63:ASP:HB3	1:C:66:GLY:H	1.77	0.49
1:C:87:ASN:HD22	1:C:87:ASN:N	2.09	0.49
1:A:210:GLN:O	1:A:213:GLN:HG2	2.12	0.48
1:D:178:ILE:HG12	1:D:196:THR:HB	1.94	0.48
1:A:10:THR:CG2	1:A:91:ASP:HA	2.24	0.48
1:D:98:SER:O	1:D:99:ALA:CB	2.61	0.48
1:C:79:MSE:HE2	1:C:81:LEU:HD11	1.96	0.48
1:D:36:ILE:N	1:D:36:ILE:HD12	2.28	0.48
1:D:71:GLN:NE2	1:D:72:GLN:NE2	2.60	0.48
1:C:225:LEU:HD22	1:C:225:LEU:H	1.77	0.48
1:B:13:GLU:OE2	1:B:31:LYS:HE3	2.13	0.48
1:A:20:ILE:N	1:A:97:ASN:HD21	2.10	0.48
1:D:141:LEU:CD1	1:D:172:LEU:HD11	2.42	0.48
1:B:141:LEU:HD22	1:B:168:ASP:HB3	1.95	0.48
1:D:182:ASN:ND2	1:D:182:ASN:H	2.11	0.48
1:D:9:ILE:HG12	1:D:41:ALA:HB2	1.96	0.48
1:B:225:LEU:H	1:B:225:LEU:HD23	1.76	0.48
1:D:9:ILE:HD11	1:D:56:LEU:HD21	1.96	0.48
1:B:267:ARG:HB2	1:B:273:ALA:HB1	1.95	0.48
1:D:50:CYS:SG	1:D:280:LEU:HD13	2.53	0.48
1:A:181:VAL:O	1:A:199:ILE:HA	2.14	0.48
1:A:138:ASN:O	1:A:139:PHE:CD1	2.67	0.48
1:D:183:GLU:OE1	1:D:201:GLN:HG2	2.14	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:117:LEU:HD22	1:A:143:LYS:HD3	1.96	0.48
1:A:121:ILE:HD11	1:A:150:TYR:CE2	2.49	0.48
1:B:226:ASP:OD2	1:B:269:GLY:N	2.44	0.48
1:A:10:THR:HG23	1:A:91:ASP:CA	2.23	0.47
1:A:181:VAL:HG13	1:A:185:GLU:CB	2.29	0.47
1:A:117:LEU:CA	5:A:336:HOH:O	2.62	0.47
1:B:46:ILE:HD13	1:B:233:THR:HA	1.96	0.47
1:C:152:ARG:NH1	1:C:175:LEU:HD23	2.28	0.47
1:B:6:THR:HG1	1:B:135:GLN:HG2	1.79	0.47
1:D:180:VAL:HG11	1:D:235:LEU:HD23	1.96	0.47
1:A:187:GLU:O	1:A:188:LEU:O	2.32	0.47
1:A:117:LEU:HD22	1:A:143:LYS:CD	2.45	0.47
1:D:36:ILE:HD11	1:D:68:TRP:CE2	2.49	0.47
1:C:105:ILE:HD12	1:C:106:ILE:N	2.29	0.47
1:C:121:ILE:HG12	1:C:122:PRO:HD3	1.95	0.47
1:A:65:ASN:HB2	1:A:89:HIS:CE1	2.47	0.47
1:B:225:LEU:HD12	1:B:268:ARG:NH2	2.29	0.47
1:A:33:SER:CB	1:A:268:ARG:NH2	2.78	0.47
1:A:11:VAL:O	1:A:91:ASP:HB2	2.14	0.47
1:B:22:LYS:NZ	1:B:22:LYS:CB	2.78	0.47
1:D:65:ASN:O	1:D:69:ILE:HG13	2.15	0.47
1:D:8:ASN:N	1:D:8:ASN:HD22	2.08	0.47
1:B:43:GLN:HE22	1:B:161:ASN:HD22	1.62	0.47
1:D:115:PHE:CB	1:D:120:MSE:HE3	2.42	0.47
1:C:152:ARG:NH2	1:C:174:PRO:O	2.47	0.47
1:B:94:ILE:CD1	1:B:94:ILE:N	2.76	0.47
1:D:165:VAL:HG22	1:D:166:ASN:N	2.30	0.47
1:A:65:ASN:HB3	1:A:89:HIS:HE1	1.77	0.47
1:B:213:GLN:HB2	5:B:307:HOH:O	2.14	0.47
1:D:175:LEU:HD22	1:D:175:LEU:N	2.30	0.47
1:A:204:ALA:O	1:A:216:PHE:HZ	1.98	0.47
1:B:20:ILE:CG2	1:D:20:ILE:CG2	2.93	0.46
1:B:238:MSE:HE3	1:B:249:PRO:HG3	1.96	0.46
1:D:134:LEU:HD13	1:D:159:VAL:HG13	1.98	0.46
1:A:55:ARG:HD3	1:A:79:MSE:HE2	1.97	0.46
1:C:135:GLN:HE22	1:C:144:THR:HA	1.79	0.46
1:B:63:ASP:N	1:B:89:HIS:CE1	2.68	0.46
1:D:9:ILE:HD11	1:D:56:LEU:CD2	2.45	0.46
1:D:183:GLU:H	1:D:183:GLU:HG2	1.55	0.46
1:D:119:GLU:HG2	5:D:305:HOH:O	2.16	0.46
1:A:188:LEU:HG	1:A:188:LEU:O	2.14	0.46
1:B:11:VAL:HG12	1:B:91:ASP:HB3	1.97	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:215:GLN:OE1	1:D:251:ALA:HB2	2.14	0.46
1:C:6:THR:HG23	1:C:139:PHE:CE2	2.39	0.46
1:B:22:LYS:O	1:B:23:LYS:O	2.34	0.46
1:D:267:ARG:HG2	1:D:267:ARG:NH1	2.31	0.46
1:B:87:ASN:OD1	1:B:88:GLN:N	2.38	0.46
1:C:-1:LEU:HD12	1:C:2:ARG:H	1.76	0.46
1:C:94:ILE:HG22	1:C:96:LEU:HD22	1.97	0.46
1:D:182:ASN:ND2	1:D:185:GLU:CD	2.69	0.46
1:D:226:ASP:HB3	1:D:267:ARG:O	2.15	0.46
1:B:70:ARG:CD	1:B:85:HIS:CE1	2.97	0.46
1:D:150:TYR:OH	1:D:154:ARG:NH1	2.49	0.46
1:C:121:ILE:N	1:C:122:PRO:CD	2.79	0.46
1:B:82:PRO:HG3	1:B:86:PHE:CZ	2.50	0.46
1:B:15:TRP:CZ2	1:D:94:ILE:HD11	2.52	0.45
1:D:219:ALA:O	1:D:221:PRO:HD3	2.16	0.45
1:D:2:ARG:HG2	1:D:2:ARG:NH1	2.31	0.45
1:D:90:SER:O	1:D:92:THR:HG23	2.17	0.45
1:A:121:ILE:HD11	1:A:150:TYR:CG	2.51	0.45
1:C:152:ARG:HH11	1:C:175:LEU:HD23	1.82	0.45
1:D:163:SER:HB2	1:D:185:GLU:OE2	2.17	0.45
1:D:281:ALA:O	1:D:285:THR:HG23	2.16	0.45
1:A:121:ILE:HG23	1:A:122:PRO:CD	2.46	0.45
1:A:124:MSE:HE1	1:A:147:LEU:HD12	1.99	0.45
1:D:145:ARG:HH11	1:D:145:ARG:HB2	1.77	0.45
1:A:160:PHE:CD2	1:A:176:ILE:HD13	2.52	0.45
1:D:271:LEU:HA	1:D:274:PHE:CD2	2.52	0.45
1:A:194:VAL:HG12	1:A:196:THR:H	1.82	0.45
1:A:132:ILE:HD11	1:A:243:LEU:HD11	1.98	0.45
1:C:237:VAL:HG21	1:C:260:ALA:CB	2.47	0.45
1:C:267:ARG:HB2	1:C:273:ALA:HB1	1.97	0.45
1:A:150:TYR:O	1:A:153:SER:HB3	2.17	0.45
1:C:152:ARG:NH2	5:C:357:HOH:O	2.31	0.45
1:B:39:LYS:CE	1:B:232:ASP:OD1	2.61	0.45
1:B:124:MSE:HE3	1:B:156:MSE:CE	2.43	0.45
1:A:212:GLY:N	5:A:320:HOH:O	2.46	0.45
1:B:23:LYS:HZ2	1:D:101:GLY:HA3	1.79	0.45
1:A:23:LYS:NZ	1:C:100:ASP:CB	2.79	0.45
1:A:232:ASP:HB3	5:A:355:HOH:O	2.17	0.45
1:D:276:GLY:O	1:D:277:SER:C	2.55	0.45
1:D:71:GLN:HE22	1:D:72:GLN:HE21	1.60	0.45
1:A:209:VAL:CG1	1:A:210:GLN:N	2.80	0.44
1:A:43:GLN:OE1	1:A:134:LEU:HD21	2.18	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:56:LEU:HD23	1:C:80:LEU:HD13	2.00	0.44
1:D:11:VAL:CG2	1:D:65:ASN:HB3	2.42	0.44
1:B:109:THR:OG1	1:B:113:ASP:CG	2.56	0.44
1:D:152:ARG:HG2	1:D:152:ARG:NH1	2.33	0.44
1:C:63:ASP:OD1	1:C:89:HIS:HA	2.17	0.44
1:D:43:GLN:HG3	1:D:232:ASP:HB3	1.99	0.44
1:C:166:ASN:HB3	1:C:169:PHE:CE1	2.52	0.44
1:A:282:ALA:HB3	5:A:311:HOH:O	2.16	0.44
1:B:211:GLU:N	5:B:307:HOH:O	2.51	0.44
1:D:221:PRO:O	1:D:222:ALA:HB2	2.17	0.44
1:D:197:LEU:HB3	1:D:209:VAL:HB	1.99	0.44
1:D:152:ARG:HE	1:D:175:LEU:HA	1.83	0.44
1:B:86:PHE:HB2	1:B:87:ASN:H	1.49	0.44
1:B:11:VAL:HG11	1:B:65:ASN:OD1	2.18	0.44
1:A:224:ALA:CB	1:A:265:VAL:CG1	2.96	0.44
1:C:132:ILE:HG12	1:C:157:THR:HB	2.00	0.44
1:B:128:VAL:HG12	1:B:129:ALA:O	2.17	0.44
1:D:76:GLU:C	1:D:78:LEU:H	2.20	0.44
1:D:145:ARG:HH11	1:D:145:ARG:CG	2.29	0.44
1:D:98:SER:O	1:D:99:ALA:HB2	2.17	0.44
1:B:55:ARG:HB2	1:B:55:ARG:HE	1.61	0.44
1:B:247:VAL:O	1:B:248:ALA:CB	2.61	0.44
1:A:141:LEU:O	1:A:141:LEU:HD12	2.18	0.44
1:A:121:ILE:HD11	1:A:150:TYR:CD2	2.53	0.44
1:B:237:VAL:HG21	1:B:260:ALA:CB	2.48	0.44
1:B:-1:LEU:C	1:B:-1:LEU:HD23	2.38	0.44
1:C:145:ARG:O	1:C:149:GLN:HG3	2.18	0.44
1:C:135:GLN:HB2	1:C:148:PHE:CZ	2.53	0.43
1:A:43:GLN:CG	1:A:232:ASP:HB2	2.47	0.43
1:D:237:VAL:HG21	1:D:260:ALA:HB1	1.99	0.43
1:A:106:ILE:CD1	1:C:21:PRO:HG2	2.48	0.43
1:D:154:ARG:HE	1:D:154:ARG:HB3	1.54	0.43
1:D:182:ASN:HD21	1:D:185:GLU:CD	2.21	0.43
1:C:31:LYS:NZ	1:C:31:LYS:HB2	2.33	0.43
1:B:10:THR:HG22	1:B:90:SER:HB2	2.00	0.43
1:A:194:VAL:O	1:A:210:GLN:CG	2.63	0.43
1:C:13:GLU:OE2	1:C:31:LYS:HD3	2.18	0.43
1:A:63:ASP:OD1	1:A:89:HIS:NE2	2.52	0.43
1:D:145:ARG:HH11	1:D:145:ARG:HB3	1.82	0.43
1:D:208:LEU:HB2	1:D:254:LEU:CD2	2.49	0.43
1:D:61:GLY:C	1:D:63:ASP:H	2.22	0.43
1:C:173:TRP:N	1:C:174:PRO:HD2	2.34	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:225:LEU:HD23	1:C:225:LEU:H	1.82	0.43
1:C:248:ALA:O	1:C:249:PRO:C	2.55	0.43
1:C:197:LEU:HB3	1:C:209:VAL:HB	2.01	0.43
1:C:17:ILE:HD13	1:C:96:LEU:HD12	2.00	0.43
1:B:11:VAL:O	1:B:11:VAL:HG13	2.19	0.43
1:B:23:LYS:HZ3	1:D:101:GLY:HA3	1.84	0.43
1:A:109:THR:O	1:A:110:ALA:HB3	2.18	0.43
1:A:21:PRO:HD3	1:A:27:ILE:HD11	2.00	0.43
1:D:18:PRO:HA	1:D:99:ALA:HB2	2.01	0.43
1:C:115:PHE:CB	1:C:120:MSE:HE3	2.49	0.43
1:B:225:LEU:HD21	1:B:266:SER:C	2.38	0.43
1:D:94:ILE:HG22	1:D:96:LEU:HD22	2.00	0.43
1:A:187:GLU:O	1:A:188:LEU:CG	2.67	0.42
1:D:198:VAL:HG21	1:D:238:MSE:SE	2.69	0.42
1:C:259:ARG:HB3	1:C:283:LEU:HD11	2.00	0.42
1:C:141:LEU:HD11	1:C:172:LEU:HD13	2.00	0.42
1:A:259:ARG:HD2	1:A:283:LEU:HB2	2.01	0.42
1:A:106:ILE:N	1:A:106:ILE:HD12	2.34	0.42
1:C:8:ASN:N	1:C:8:ASN:HD22	2.14	0.42
1:B:116:SER:O	1:B:120:MSE:HG3	2.19	0.42
1:C:230:ALA:HB2	1:C:265:VAL:CG1	2.49	0.42
1:D:105:ILE:CG2	1:D:106:ILE:N	2.83	0.42
1:B:49:ARG:HD3	1:B:76:GLU:OE1	2.19	0.42
1:A:267:ARG:HD2	1:A:273:ALA:O	2.20	0.42
1:A:185:GLU:O	1:A:189:LEU:HB2	2.19	0.42
1:B:22:LYS:HB2	1:B:22:LYS:HZ2	1.83	0.42
1:D:135:GLN:HE22	1:D:147:LEU:CD1	2.33	0.42
1:D:46:ILE:HD12	1:D:233:THR:HA	2.00	0.42
1:A:109:THR:O	1:A:111:ALA:N	2.47	0.42
1:A:188:LEU:HD12	1:A:188:LEU:C	2.40	0.42
1:A:209:VAL:HG12	1:A:210:GLN:N	2.33	0.42
1:D:63:ASP:OD1	1:D:64:SER:N	2.49	0.42
1:A:12:ASP:HB3	1:A:94:ILE:HD11	2.02	0.42
1:A:12:ASP:OD1	1:A:92:THR:CG2	2.67	0.42
1:B:43:GLN:O	1:B:47:LEU:HB2	2.20	0.42
1:B:198:VAL:HG21	1:B:238:MSE:SE	2.70	0.42
1:B:50:CYS:SG	1:B:280:LEU:HG	2.59	0.42
1:A:235:LEU:C	1:A:235:LEU:HD23	2.40	0.42
1:D:48:SER:OG	1:D:77:PRO:HG2	2.19	0.42
1:B:28:HIS:NE2	1:D:164:PRO:CG	2.82	0.42
1:B:20:ILE:HG21	1:D:20:ILE:CG2	2.48	0.42
1:D:35:ASP:OD1	1:D:36:ILE:N	2.51	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:121:ILE:N	1:D:122:PRO:HD2	2.34	0.42
1:C:107:THR:HG22	1:C:108:THR:O	2.20	0.42
1:B:228:THR:CG2	1:B:270:THR:HG22	2.40	0.41
1:D:9:ILE:HD13	1:D:73:ILE:CD1	2.50	0.41
1:B:20:ILE:H	1:B:20:ILE:HG13	1.58	0.41
1:A:121:ILE:CG2	1:A:122:PRO:CD	2.97	0.41
1:C:79:MSE:HE2	1:C:81:LEU:CD2	2.41	0.41
1:B:213:GLN:N	5:B:307:HOH:O	2.43	0.41
1:C:87:ASN:O	1:C:87:ASN:ND2	2.53	0.41
1:B:45:ILE:O	1:B:49:ARG:HG3	2.20	0.41
1:D:95:ILE:CG2	1:D:103:ASN:HD22	2.33	0.41
1:D:182:ASN:N	1:D:182:ASN:HD22	2.18	0.41
1:B:12:ASP:OD2	1:B:92:THR:CG2	2.68	0.41
1:A:121:ILE:HA	1:A:124:MSE:HE3	2.02	0.41
1:A:150:TYR:CZ	1:A:154:ARG:HD2	2.56	0.41
1:A:96:LEU:HD12	1:A:97:ASN:N	2.35	0.41
1:D:43:GLN:HE21	1:D:134:LEU:HD21	1.86	0.41
1:B:210:GLN:C	5:B:307:HOH:O	2.58	0.41
1:C:225:LEU:HD23	1:C:226:ASP:N	2.36	0.41
1:D:235:LEU:HD13	1:D:235:LEU:C	2.41	0.41
1:D:85:HIS:CD2	1:D:85:HIS:H	2.38	0.41
1:A:96:LEU:HA	1:A:103:ASN:OD1	2.21	0.41
1:C:235:LEU:C	1:C:235:LEU:HD12	2.41	0.41
1:B:55:ARG:HG2	1:B:81:LEU:HD12	2.02	0.41
1:D:263:ILE:O	1:D:267:ARG:HD3	2.21	0.41
1:C:245:ARG:HH22	1:C:250:ASP:CG	2.24	0.41
1:B:133:LEU:HD22	1:B:151:ALA:HB2	2.02	0.41
1:C:169:PHE:O	1:C:172:LEU:HB2	2.21	0.41
1:A:183:GLU:H	1:A:183:GLU:HG2	1.52	0.41
1:A:222:ALA:O	1:A:223:GLU:CG	2.69	0.41
1:B:238:MSE:HG3	1:B:254:LEU:HD12	2.02	0.41
1:B:-1:LEU:HD23	1:B:2:ARG:N	2.36	0.41
1:B:235:LEU:C	1:B:235:LEU:HD12	2.41	0.41
1:B:73:ILE:HG13	1:B:80:LEU:HD22	2.03	0.41
1:A:90:SER:O	1:A:91:ASP:HB3	2.21	0.41
1:D:180:VAL:HG21	1:D:235:LEU:HD22	2.02	0.41
1:A:6:THR:HG21	1:A:147:LEU:HG	2.02	0.40
1:B:210:GLN:HB2	1:B:211:GLU:OE1	2.21	0.40
1:D:135:GLN:HG3	5:D:300:HOH:O	2.20	0.40
1:A:207:TRP:CZ3	1:A:216:PHE:HB2	2.57	0.40
1:C:150:TYR:OH	1:C:154:ARG:HD2	2.21	0.40
1:D:276:GLY:C	1:D:278:ARG:N	2.70	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:283:LEU:HD12	1:A:284:LEU:N	2.36	0.40
1:C:21:PRO:HG3	1:C:27:ILE:HG12	2.03	0.40
1:D:129:ALA:O	1:D:155:GLY:C	2.60	0.40
1:C:61:GLY:C	1:C:63:ASP:H	2.24	0.40
1:A:212:GLY:CA	5:A:320:HOH:O	2.69	0.40
1:B:12:ASP:OD2	1:B:92:THR:HG22	2.22	0.40
1:A:173:TRP:CE3	1:A:173:TRP:HA	2.56	0.40
1:C:201:GLN:HB2	1:C:201:GLN:HE21	1.65	0.40
1:A:96:LEU:HD12	1:A:96:LEU:C	2.41	0.40
1:B:76:GLU:C	1:B:78:LEU:H	2.25	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	
1	A	283/299 (95%)	247 (87%)	27 (10%)	9 (3%)	6	1
1	B	276/299 (92%)	252 (91%)	19 (7%)	5 (2%)	13	4
1	C	283/299 (95%)	268 (95%)	12 (4%)	3 (1%)	21	11
1	D	283/299 (95%)	253 (89%)	18 (6%)	12 (4%)	4	1
All	All	1125/1196 (94%)	1020 (91%)	76 (7%)	29 (3%)	8	2

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	101	GLY
1	A	210	GLN
1	A	222	ALA
1	A	223	GLU
1	B	23	LYS
1	B	110	ALA
1	D	98	SER
1	D	99	ALA

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Mol	Chain	Res	Type
1	D	100	ASP
1	A	88	GLN
1	A	189	LEU
1	B	193	GLY
1	C	101	GLY
1	D	246	GLY
1	C	100	ASP
1	D	129	ALA
1	D	222	ALA
1	D	278	ARG
1	A	139	PHE
1	A	162	PRO
1	A	188	LEU
1	D	87	ASN
1	D	89	HIS
1	B	211	GLU
1	C	162	PRO
1	D	153	SER
1	B	249	PRO
1	D	247	VAL
1	D	18	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	223/228 (98%)	203 (91%)	20 (9%)	14	8
1	B	220/228 (96%)	199 (90%)	21 (10%)	12	7
1	C	223/228 (98%)	204 (92%)	19 (8%)	15	10
1	D	223/228 (98%)	201 (90%)	22 (10%)	11	7
All	All	889/912 (98%)	807 (91%)	82 (9%)	13	8

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR

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Mol	Chain	Res	Type
1	A	8	ASN
1	A	35	ASP
1	A	71	GLN
1	A	74	LYS
1	A	83	ASP
1	A	87	ASN
1	A	88	GLN
1	A	119	GLU
1	A	126	ASP
1	A	138	ASN
1	A	147	LEU
1	A	170	CYS
1	A	183	GLU
1	A	188	LEU
1	A	189	LEU
1	A	210	GLN
1	A	232	ASP
1	A	243	LEU
1	A	283	LEU
1	B	8	ASN
1	B	35	ASP
1	B	55	ARG
1	B	62	ASN
1	B	71	GLN
1	B	89	HIS
1	B	96	LEU
1	B	108	THR
1	B	109	THR
1	B	115	PHE
1	B	152	ARG
1	B	189	LEU
1	B	210	GLN
1	B	211	GLU
1	B	225	LEU
1	B	227	THR
1	B	235	LEU
1	B	243	LEU
1	B	280	LEU
1	B	283	LEU
1	B	285	THR
1	C	-1	LEU
1	C	6	THR

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Mol	Chain	Res	Type
1	C	8	ASN
1	C	47	LEU
1	C	63	ASP
1	C	72	GLN
1	C	87	ASN
1	C	100	ASP
1	C	102	ASP
1	C	105	ILE
1	C	115	PHE
1	C	121	ILE
1	C	142	ASP
1	C	172	LEU
1	C	183	GLU
1	C	213	GLN
1	C	225	LEU
1	C	227	THR
1	C	243	LEU
1	D	-1	LEU
1	D	8	ASN
1	D	34	GLN
1	D	47	LEU
1	D	85	HIS
1	D	87	ASN
1	D	89	HIS
1	D	90	SER
1	D	98	SER
1	D	102	ASP
1	D	108	THR
1	D	115	PHE
1	D	142	ASP
1	D	145	ARG
1	D	182	ASN
1	D	184	SER
1	D	195	LYS
1	D	223	GLU
1	D	227	THR
1	D	243	LEU
1	D	267	ARG
1	D	283	LEU

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

Mol	Chain	Res	Type
1	A	8	ASN
1	A	71	GLN
1	A	85	HIS
1	A	97	ASN
1	A	135	GLN
1	A	161	ASN
1	A	215	GLN
1	A	256	HIS
1	B	8	ASN
1	B	62	ASN
1	B	71	GLN
1	B	72	GLN
1	B	88	GLN
1	B	89	HIS
1	B	135	GLN
1	B	161	ASN
1	B	171	HIS
1	B	201	GLN
1	B	210	GLN
1	B	256	HIS
1	C	8	ASN
1	C	43	GLN
1	C	72	GLN
1	C	85	HIS
1	C	87	ASN
1	C	88	GLN
1	C	135	GLN
1	C	138	ASN
1	C	161	ASN
1	C	166	ASN
1	C	201	GLN
1	D	8	ASN
1	D	28	HIS
1	D	34	GLN
1	D	43	GLN
1	D	71	GLN
1	D	87	ASN
1	D	88	GLN
1	D	103	ASN
1	D	135	GLN
1	D	138	ASN
1	D	161	ASN
1	D	171	HIS

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Mol	Chain	Res	Type
1	D	182	ASN
1	D	210	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 ligands 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	$\# Z > 2$
2	SO4	B	298	-	4,4,4	0.26	0	6,6,6	0.07	0
2	SO4	B	299	-	4,4,4	0.31	0	6,6,6	0.07	0
2	SO4	B	300	-	4,4,4	0.23	0	6,6,6	0.10	0
2	SO4	C	1	-	4,4,4	0.19	0	6,6,6	0.15	0
2	SO4	C	298	-	4,4,4	0.22	0	6,6,6	0.17	0
3	GOL	C	501	-	5,5,5	0.23	0	5,5,5	0.36	0
2	SO4	D	298	-	4,4,4	0.37	0	6,6,6	0.13	0
4	RIB	D	311	-	10,10,10	1.51	3 (30%)	14,14,14	1.55	3 (21%)

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
2	SO4	B	298	-	-	0/0/0/0	0/0/0/0
2	SO4	B	299	-	-	0/0/0/0	0/0/0/0
2	SO4	B	300	-	-	0/0/0/0	0/0/0/0
2	SO4	C	1	-	-	0/0/0/0	0/0/0/0
2	SO4	C	298	-	-	0/0/0/0	0/0/0/0
3	GOL	C	501	-	-	0/4/4/4	0/0/0/0
2	SO4	D	298	-	-	0/0/0/0	0/0/0/0
4	RIB	D	311	-	-	0/2/18/18	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	311	RIB	O4'-C1'	2.99	1.45	1.42
4	D	311	RIB	C1'-C2'	2.24	1.54	1.52
4	D	311	RIB	O1'-C1'	2.10	1.47	1.39

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	311	RIB	C1'-C2'-C3'	2.95	106.14	102.33
4	D	311	RIB	O4'-C4'-C3'	2.58	110.39	105.17
4	D	311	RIB	O4'-C1'-C2'	2.55	107.16	104.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	A	285/299 (95%)	0.34	12 (4%) 35 38	25, 39, 58, 71	0
1	B	282/299 (94%)	0.31	10 (3%) 42 46	26, 39, 61, 83	0
1	C	285/299 (95%)	0.25	5 (1%) 65 71	27, 38, 52, 62	0
1	D	285/299 (95%)	0.45	11 (3%) 37 41	27, 43, 64, 80	0
All	All	1137/1196 (95%)	0.34	38 (3%) 44 48	25, 39, 61, 83	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	222	ALA	5.3
1	A	99	ALA	5.1
1	D	221	PRO	5.0
1	B	99	ALA	4.3
1	D	220	VAL	4.2
1	C	100	ASP	4.2
1	D	100	ASP	4.0
1	B	101	GLY	3.9
1	A	100	ASP	3.8
1	B	100	ASP	3.8
1	D	99	ALA	3.5
1	B	104	ALA	3.4
1	D	21	PRO	3.2
1	A	210	GLN	3.0
1	A	118	ASP	3.0
1	C	235	LEU	2.9
1	C	211	GLU	2.8
1	D	87	ASN	2.7
1	B	22	LYS	2.5
1	D	223	GLU	2.5
1	B	87	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	224	ALA	2.4
1	B	225	LEU	2.3
1	A	89	HIS	2.3
1	A	98	SER	2.3
1	A	126	ASP	2.3
1	D	23	LYS	2.2
1	C	152	ARG	2.2
1	A	224	ALA	2.2
1	A	222	ALA	2.2
1	C	128	VAL	2.1
1	A	86	PHE	2.1
1	B	283	LEU	2.1
1	A	102	ASP	2.1
1	B	105	ILE	2.1
1	B	211	GLU	2.1
1	A	41	ALA	2.0
1	D	268	ARG	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q < 0.9
2	SO4	C	1	5/5	0.22	5.47	74,74,76,77	0
3	GOL	C	501	6/6	0.26	2.16	41,43,44,48	0
4	RIB	D	311	10/10	0.22	1.68	53,55,58,59	0
2	SO4	B	298	5/5	0.16	1.14	72,72,73,74	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SO4	B	300	5/5	0.18	0.75	89,90,90,91	0
2	SO4	C	298	5/5	0.21	0.35	42,45,46,48	0
2	SO4	D	298	5/5	0.14	-0.29	62,62,64,64	0
2	SO4	B	299	5/5	0.13	-1.30	82,83,83,84	0

6.5 Other polymers

There are no such residues in this entry.