



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 02:21 PM GMT

PDB ID : 3CK5
Title : Crystal structure of a racemase from *Streptomyces coelicolor* A3(2) with bound magnesium
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Deposited on : 2008-03-14
Resolution : 2.30 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>

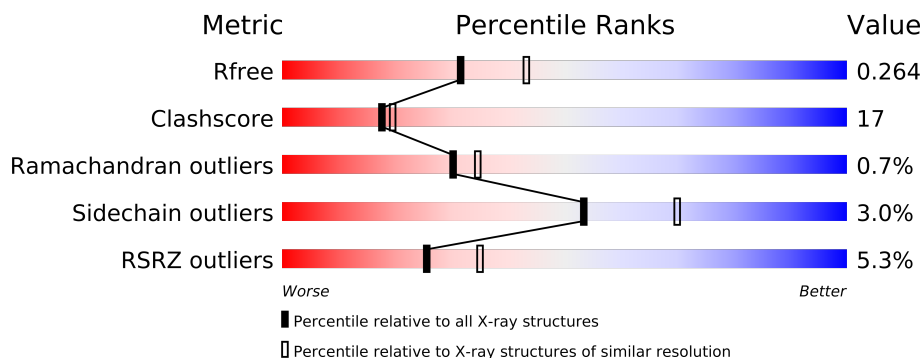
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.30 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	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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	371	
1	B	371	
1	C	371	
1	D	371	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11164 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 Putative racemase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	S	Se	0	0	0
			2749	1738	492	506	2	11			
1	B	357	Total	C	N	O	S	Se	0	0	0
			2749	1738	492	506	2	11			
1	C	357	Total	C	N	O	S	Se	0	0	0
			2749	1738	492	506	2	11			
1	D	334	Total	C	N	O	S	Se	0	0	0
			2577	1631	462	474	2	8			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	EXPRESSION TAG	UNP Q9RKF7
A	2	SER	-	EXPRESSION TAG	UNP Q9RKF7
A	3	LEU	-	EXPRESSION TAG	UNP Q9RKF7
A	364	GLU	-	EXPRESSION TAG	UNP Q9RKF7
A	365	GLY	-	EXPRESSION TAG	UNP Q9RKF7
A	366	HIS	-	EXPRESSION TAG	UNP Q9RKF7
A	367	HIS	-	EXPRESSION TAG	UNP Q9RKF7
A	368	HIS	-	EXPRESSION TAG	UNP Q9RKF7
A	369	HIS	-	EXPRESSION TAG	UNP Q9RKF7
A	370	HIS	-	EXPRESSION TAG	UNP Q9RKF7
A	371	HIS	-	EXPRESSION TAG	UNP Q9RKF7
B	1	MSE	-	EXPRESSION TAG	UNP Q9RKF7
B	2	SER	-	EXPRESSION TAG	UNP Q9RKF7
B	3	LEU	-	EXPRESSION TAG	UNP Q9RKF7
B	364	GLU	-	EXPRESSION TAG	UNP Q9RKF7
B	365	GLY	-	EXPRESSION TAG	UNP Q9RKF7
B	366	HIS	-	EXPRESSION TAG	UNP Q9RKF7
B	367	HIS	-	EXPRESSION TAG	UNP Q9RKF7
B	368	HIS	-	EXPRESSION TAG	UNP Q9RKF7
B	369	HIS	-	EXPRESSION TAG	UNP Q9RKF7
B	370	HIS	-	EXPRESSION TAG	UNP Q9RKF7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	371	HIS	-	EXPRESSION TAG	UNP Q9RKF7
C	1	MSE	-	EXPRESSION TAG	UNP Q9RKF7
C	2	SER	-	EXPRESSION TAG	UNP Q9RKF7
C	3	LEU	-	EXPRESSION TAG	UNP Q9RKF7
C	364	GLU	-	EXPRESSION TAG	UNP Q9RKF7
C	365	GLY	-	EXPRESSION TAG	UNP Q9RKF7
C	366	HIS	-	EXPRESSION TAG	UNP Q9RKF7
C	367	HIS	-	EXPRESSION TAG	UNP Q9RKF7
C	368	HIS	-	EXPRESSION TAG	UNP Q9RKF7
C	369	HIS	-	EXPRESSION TAG	UNP Q9RKF7
C	370	HIS	-	EXPRESSION TAG	UNP Q9RKF7
C	371	HIS	-	EXPRESSION TAG	UNP Q9RKF7
D	1	MSE	-	EXPRESSION TAG	UNP Q9RKF7
D	2	SER	-	EXPRESSION TAG	UNP Q9RKF7
D	3	LEU	-	EXPRESSION TAG	UNP Q9RKF7
D	364	GLU	-	EXPRESSION TAG	UNP Q9RKF7
D	365	GLY	-	EXPRESSION TAG	UNP Q9RKF7
D	366	HIS	-	EXPRESSION TAG	UNP Q9RKF7
D	367	HIS	-	EXPRESSION TAG	UNP Q9RKF7
D	368	HIS	-	EXPRESSION TAG	UNP Q9RKF7
D	369	HIS	-	EXPRESSION TAG	UNP Q9RKF7
D	370	HIS	-	EXPRESSION TAG	UNP Q9RKF7
D	371	HIS	-	EXPRESSION TAG	UNP Q9RKF7

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	99	Total O 99 99	0	0

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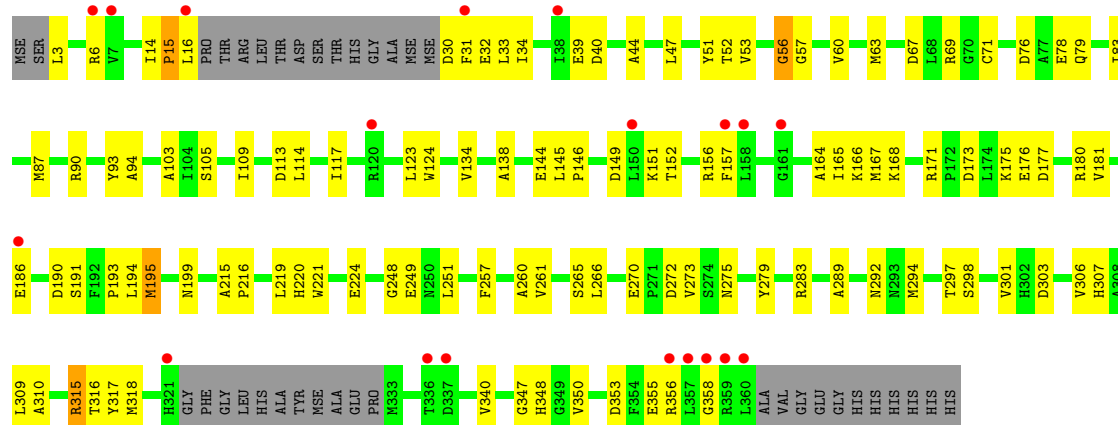
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	113	Total 113	O 113	0	0
3	C	73	Total 73	O 73	0	0
3	D	51	Total 51	O 51	0	0



● Molecule 1: Putative racemase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	177.60Å 177.60Å 112.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.26 – 2.30 49.26 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.9 (49.26-2.30) 95.0 (49.26-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.97 (at 2.32Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.264 0.224 , 0.264	Depositor DCC
R_{free} test set	1876 reflections (2.49%)	DCC
Wilson B-factor (Å ²)	34.0	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 24.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 79252 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11164	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.48% 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: MG

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	0/2802	0.64	0/3794
1	B	0.39	0/2802	0.66	0/3794
1	C	0.41	0/2802	0.67	0/3794
1	D	0.40	0/2628	0.67	0/3561
All	All	0.40	0/11034	0.66	0/14943

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	93	TYR	Sidechain

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	2749	0	2711	82	0
1	B	2749	0	2711	87	0
1	C	2749	0	2711	112	0
1	D	2577	0	2543	98	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	99	0	0	2	0
3	B	113	0	0	2	0
3	C	73	0	0	0	0
3	D	51	0	0	2	0
All	All	11164	0	10676	372	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 17.

All (372) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:60:VAL:HA	1:D:63:MSE:HE2	1.40	1.03
1:A:60:VAL:HA	1:A:63:MSE:HE3	1.39	1.03
1:C:19:ARG:HH22	1:C:29:MSE:HB2	1.25	0.98
1:A:348:HIS:HD2	1:A:350:VAL:H	1.11	0.96
1:C:15:PRO:HA	1:C:30:ASP:HB3	1.48	0.93
1:B:70:GLY:HA3	1:B:90:ARG:HH12	1.34	0.91
1:A:20:LEU:HB3	1:A:141:ILE:HB	1.52	0.90
1:B:60:VAL:HA	1:B:63:MSE:CE	2.02	0.90
1:C:348:HIS:HD2	1:C:350:VAL:H	1.16	0.89
1:B:28:MSE:HE2	1:B:53:VAL:HG11	1.54	0.88
1:B:60:VAL:HA	1:B:63:MSE:HE2	1.54	0.88
1:C:83:ILE:HG22	1:C:87:MSE:HE2	1.55	0.87
1:D:289:ALA:HA	1:D:294:MSE:HE3	1.57	0.86
1:C:270:GLU:HG3	1:C:297:THR:HG23	1.58	0.86
1:A:23:SER:OG	1:A:201:LYS:HB2	1.77	0.85
1:A:289:ALA:HA	1:A:294:MSE:HE3	1.56	0.84
1:C:60:VAL:HA	1:C:63:MSE:HE3	1.61	0.82
1:D:292:ASN:HB2	1:D:294:MSE:HE2	1.60	0.82
1:C:14:ILE:O	1:C:30:ASP:HB2	1.80	0.81
1:B:289:ALA:HA	1:B:294:MSE:HE3	1.62	0.81
1:A:63:MSE:HE1	1:A:99:HIS:HB3	1.62	0.81
1:B:348:HIS:HD2	1:B:350:VAL:H	1.29	0.80

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:195:MSE:HE2	1:D:220:HIS:HB3	1.62	0.80
1:B:71:CYS:SG	1:B:90:ARG:HD2	2.22	0.80
1:C:289:ALA:HA	1:C:294:MSE:HE3	1.65	0.79
1:D:40:ASP:OD2	1:D:44:ALA:HB3	1.82	0.79
1:D:272:ASP:HB3	1:D:275:ASN:ND2	1.98	0.79
1:B:47:LEU:O	1:B:348:HIS:HE1	1.65	0.78
1:C:29:MSE:HE2	1:C:30:ASP:HB3	1.65	0.78
1:D:113:ASP:O	1:D:117:ILE:HG12	1.84	0.76
1:A:19:ARG:NE	1:A:29:MSE:HB3	2.00	0.76
1:C:15:PRO:HA	1:C:30:ASP:CB	2.15	0.75
1:C:348:HIS:CD2	1:C:350:VAL:H	2.03	0.75
1:B:150:LEU:HG	1:B:184:LEU:HD11	1.68	0.75
1:C:131:ASP:HB3	1:C:315:ARG:NH2	2.01	0.75
1:A:329:MSE:HE3	1:A:350:VAL:HG13	1.66	0.75
1:C:131:ASP:HB3	1:C:315:ARG:HH22	1.51	0.74
1:C:60:VAL:HA	1:C:63:MSE:CE	2.16	0.74
1:B:348:HIS:CD2	1:B:350:VAL:H	2.06	0.73
1:D:279:TYR:O	1:D:283:ARG:HG3	1.89	0.73
1:A:34:ILE:N	1:A:34:ILE:HD12	2.03	0.73
1:C:21:THR:H	1:C:141:ILE:HG22	1.55	0.72
1:B:7:VAL:HB	1:B:69:ARG:HH12	1.55	0.72
1:D:348:HIS:HD2	1:D:350:VAL:H	1.38	0.71
1:D:195:MSE:HE2	1:D:220:HIS:CB	2.21	0.71
1:D:69:ARG:HH11	1:D:69:ARG:HG3	1.55	0.71
1:C:113:ASP:O	1:C:117:ILE:HG13	1.91	0.70
1:B:13:ARG:HH11	1:B:13:ARG:HG3	1.57	0.70
1:C:19:ARG:NH2	1:C:29:MSE:HB2	2.05	0.70
1:B:21:THR:N	1:B:141:ILE:HG22	2.06	0.70
1:A:60:VAL:CA	1:A:63:MSE:HE3	2.20	0.69
1:C:19:ARG:HH22	1:C:29:MSE:CB	2.02	0.69
1:A:348:HIS:CD2	1:A:350:VAL:H	2.03	0.69
1:A:6:ARG:NH1	1:A:8:ARG:HD3	2.08	0.69
1:C:70:GLY:HA3	1:C:90:ARG:HH12	1.58	0.69
1:C:6:ARG:HB3	1:C:39:GLU:HB3	1.75	0.68
1:B:23:SER:OG	1:B:201:LYS:HB2	1.93	0.68
1:C:19:ARG:HH12	1:C:29:MSE:HB3	1.59	0.68
1:C:40:ASP:OD2	1:C:44:ALA:HB3	1.94	0.68
1:B:16:LEU:HB2	1:B:29:MSE:O	1.94	0.68
1:A:131:ASP:HB3	1:A:315:ARG:HH22	1.58	0.67
1:D:194:LEU:CA	1:D:195:MSE:HE3	2.24	0.67
1:B:16:LEU:HD23	1:B:28:MSE:HB3	1.77	0.67
1:A:69:ARG:HH11	1:A:73:LEU:HD11	1.60	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:16:LEU:HD11	1:B:31:PHE:CD2	2.31	0.66
1:C:289:ALA:HA	1:C:294:MSE:CE	2.25	0.65
1:D:79:GLN:O	1:D:83:ILE:HG12	1.96	0.65
1:D:315:ARG:C	1:D:315:ARG:HD2	2.17	0.64
1:B:167:MSE:SE	1:B:184:LEU:HD22	2.48	0.64
1:C:67:ASP:O	1:C:90:ARG:HD3	1.97	0.64
1:A:289:ALA:HA	1:A:294:MSE:CE	2.28	0.64
1:A:53:VAL:O	1:A:54:ASN:HB2	1.98	0.64
1:D:164:ALA:O	1:D:165:ILE:HD13	1.98	0.64
1:B:131:ASP:HB3	1:B:315:ARG:HH22	1.62	0.64
1:C:18:THR:HG22	1:C:19:ARG:H	1.63	0.64
1:A:55:HIS:HB3	1:B:67:ASP:OD1	1.98	0.64
1:B:7:VAL:HB	1:B:69:ARG:NH1	2.12	0.63
1:D:78:GLU:HG3	3:D:409:HOH:O	1.97	0.63
1:B:105:SER:O	1:B:109:ILE:HG13	1.98	0.63
1:D:144:GLU:HA	1:D:171:ARG:HH12	1.63	0.63
1:C:20:LEU:HB3	1:C:141:ILE:CG2	2.29	0.62
1:A:292:ASN:HB3	1:A:294:MSE:HE2	1.80	0.62
1:B:70:GLY:CA	1:B:90:ARG:HH12	2.11	0.62
1:B:18:THR:HG23	1:B:19:ARG:HG3	1.81	0.62
1:B:28:MSE:HE2	1:B:53:VAL:CG1	2.30	0.62
1:D:63:MSE:HE3	1:D:103:ALA:HB2	1.82	0.62
1:D:194:LEU:C	1:D:195:MSE:HE3	2.20	0.62
1:A:63:MSE:HE1	1:A:99:HIS:CB	2.30	0.61
1:A:328:TYR:HE1	1:A:356:ARG:HG2	1.65	0.61
1:D:194:LEU:N	1:D:195:MSE:HE3	2.16	0.61
1:C:70:GLY:HA3	1:C:90:ARG:NH1	2.16	0.61
1:A:309:LEU:O	1:A:315:ARG:HG2	2.01	0.61
1:A:19:ARG:CZ	1:A:29:MSE:HB3	2.30	0.60
1:D:168:LYS:NZ	1:D:199:ASN:HD21	1.99	0.60
1:C:21:THR:HG21	1:C:144:GLU:OE1	2.01	0.60
1:D:60:VAL:HA	1:D:63:MSE:CE	2.25	0.60
1:C:20:LEU:HB3	1:C:141:ILE:HG21	1.83	0.59
1:A:168:LYS:NZ	1:A:199:ASN:HD21	2.00	0.59
1:C:150:LEU:HG	1:C:184:LEU:HD21	1.83	0.59
1:D:14:ILE:O	1:D:30:ASP:HB2	2.03	0.58
1:B:315:ARG:HH11	1:B:315:ARG:HG2	1.66	0.58
1:C:34:ILE:HD12	1:C:34:ILE:N	2.19	0.58
1:D:109:ILE:HD11	1:D:273:VAL:HG22	1.84	0.58
1:B:289:ALA:HA	1:B:294:MSE:CE	2.31	0.58
1:D:69:ARG:HG3	1:D:69:ARG:NH1	2.19	0.58
1:B:18:THR:HG23	1:B:19:ARG:CG	2.34	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:18:THR:HG22	1:C:19:ARG:N	2.18	0.58
1:C:166:LYS:HA	1:C:195:MSE:O	2.03	0.57
1:C:21:THR:HG21	1:C:144:GLU:CD	2.25	0.57
1:C:63:MSE:HE1	1:C:99:HIS:HB3	1.85	0.57
1:C:168:LYS:NZ	1:C:199:ASN:HD21	2.03	0.57
1:C:272:ASP:OD2	1:C:299:HIS:HD2	1.88	0.57
1:B:315:ARG:C	1:B:315:ARG:HD2	2.25	0.56
1:C:54:ASN:HB3	1:C:55:HIS:ND1	2.21	0.56
1:D:138:ALA:HB3	1:D:165:ILE:HD12	1.87	0.56
1:C:353:ASP:OD1	1:C:355:GLU:HB2	2.05	0.56
1:C:91:LEU:HD13	1:C:100:ALA:HB1	1.87	0.56
1:A:328:TYR:CE1	1:A:356:ARG:HG2	2.41	0.56
1:B:6:ARG:HB2	1:B:39:GLU:HB3	1.87	0.56
1:B:177:ASP:O	1:B:181:VAL:HG23	2.06	0.56
1:D:215:ALA:N	1:D:216:PRO:CD	2.69	0.56
1:B:70:GLY:HA3	1:B:90:ARG:NH1	2.14	0.56
1:D:152:THR:HG22	1:D:156:ARG:NH1	2.20	0.56
1:B:328:TYR:HE1	1:B:356:ARG:HG2	1.70	0.56
1:C:152:THR:HG22	1:C:156:ARG:NH1	2.21	0.56
1:A:292:ASN:CB	1:A:294:MSE:HE2	2.35	0.55
1:B:292:ASN:HB2	1:B:294:MSE:HE2	1.89	0.55
1:D:309:LEU:O	1:D:315:ARG:HG2	2.06	0.55
1:B:16:LEU:HD13	1:B:30:ASP:HA	1.89	0.55
1:C:109:ILE:HD11	1:C:273:VAL:HG22	1.89	0.55
1:C:20:LEU:HD13	1:C:141:ILE:HB	1.89	0.55
1:C:29:MSE:HE2	1:C:30:ASP:CB	2.34	0.55
1:C:315:ARG:O	1:C:315:ARG:HD2	2.05	0.55
1:D:32:GLU:HG2	1:D:57:GLY:HA3	1.87	0.55
1:C:16:LEU:HD13	1:C:20:LEU:HD12	1.89	0.55
1:C:284:LYS:O	1:C:288:LEU:HD13	2.07	0.55
1:C:321:HIS:CG	1:C:321:HIS:O	2.60	0.55
1:A:173:ASP:OD1	1:A:175:LYS:HB3	2.06	0.55
1:D:93:TYR:O	1:D:94:ALA:HB3	2.07	0.54
1:A:115:LYS:HB2	3:A:420:HOH:O	2.06	0.54
1:A:19:ARG:HE	1:A:29:MSE:HA	1.73	0.54
1:C:173:ASP:OD2	1:C:175:LYS:HB3	2.07	0.54
1:A:11:LEU:HG	1:A:362:VAL:HG13	1.89	0.54
1:C:29:MSE:HG3	1:C:30:ASP:N	2.22	0.54
1:D:6:ARG:HB2	1:D:39:GLU:HB3	1.89	0.54
1:B:3:LEU:O	1:B:40:ASP:HB2	2.08	0.54
1:A:211:ALA:CB	1:A:241:SER:HB2	2.37	0.54
1:A:47:LEU:O	1:A:348:HIS:HE1	1.91	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:166:LYS:HA	1:B:195:MSE:O	2.08	0.54
1:D:14:ILE:HG23	1:D:15:PRO:HD2	1.90	0.54
1:C:118:ARG:HH11	1:C:118:ARG:HG2	1.73	0.54
1:C:16:LEU:HG	1:C:30:ASP:HA	1.89	0.53
1:D:173:ASP:HB3	1:D:176:GLU:OE1	2.08	0.53
1:A:87:MSE:HE3	1:A:104:ILE:HG23	1.91	0.53
1:B:14:ILE:HD12	1:B:31:PHE:CE1	2.44	0.53
1:A:302:HIS:HB2	1:A:318:MSE:HE3	1.91	0.53
1:B:83:ILE:HG22	1:B:87:MSE:HE2	1.91	0.53
1:D:193:PRO:C	1:D:195:MSE:HE3	2.28	0.53
1:B:69:ARG:HH11	1:B:69:ARG:HG3	1.74	0.53
1:A:188:LEU:HD13	1:A:192:PHE:CD2	2.44	0.53
1:A:33:LEU:C	1:A:34:ILE:HD12	2.30	0.52
1:D:67:ASP:O	1:D:90:ARG:HD3	2.10	0.52
1:D:145:LEU:HD12	1:D:146:PRO:HD2	1.92	0.52
1:A:13:ARG:HB2	1:A:13:ARG:HH11	1.75	0.52
1:B:56:GLY:O	1:B:60:VAL:HG23	2.09	0.52
1:C:270:GLU:HG3	1:C:297:THR:CG2	2.36	0.52
1:B:13:ARG:HG3	1:B:13:ARG:NH1	2.25	0.52
1:C:163:ARG:NH2	1:C:337:ASP:HA	2.25	0.52
1:A:302:HIS:HA	1:A:305:THR:HB	1.91	0.51
1:D:356:ARG:C	1:D:358:GLY:H	2.14	0.51
1:D:39:GLU:HA	1:D:44:ALA:O	2.10	0.51
1:D:353:ASP:OD1	1:D:355:GLU:HB3	2.11	0.51
1:A:6:ARG:NH1	1:A:8:ARG:CD	2.73	0.51
1:D:181:VAL:HG11	1:D:219:LEU:HD21	1.93	0.51
1:B:137:TYR:CZ	1:B:319:GLU:HB2	2.46	0.51
1:A:309:LEU:HD12	1:A:318:MSE:SE	2.61	0.50
1:B:163:ARG:NH2	1:B:337:ASP:HA	2.25	0.50
1:C:21:THR:N	1:C:141:ILE:HG22	2.23	0.50
1:B:14:ILE:HD12	1:B:31:PHE:HE1	1.77	0.50
1:D:272:ASP:HB3	1:D:275:ASN:HD21	1.76	0.50
1:C:272:ASP:HB3	1:C:275:ASN:ND2	2.26	0.50
1:A:67:ASP:OD1	1:B:55:HIS:HB3	2.11	0.50
1:C:345:ARG:NH1	1:C:349:GLY:O	2.45	0.50
1:C:19:ARG:NH1	1:C:29:MSE:HB3	2.26	0.50
1:C:152:THR:HG22	1:C:156:ARG:HH12	1.76	0.50
1:B:361:ALA:HB1	3:B:519:HOH:O	2.10	0.50
1:C:19:ARG:NH2	1:C:29:MSE:CB	2.69	0.50
1:D:303:ASP:O	1:D:306:VAL:HG12	2.11	0.50
1:D:34:ILE:HD13	1:D:60:VAL:HB	1.94	0.49
1:C:13:ARG:HG3	1:C:13:ARG:HH11	1.76	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:257:PHE:O	1:D:261:VAL:HG23	2.12	0.49
1:D:83:ILE:O	1:D:87:MSE:HG3	2.12	0.49
1:C:171:ARG:HH12	1:C:180:ARG:NE	2.09	0.49
1:B:309:LEU:O	1:B:315:ARG:HG2	2.13	0.49
1:D:292:ASN:CB	1:D:294:MSE:HE2	2.38	0.49
1:C:63:MSE:HE1	1:C:99:HIS:CB	2.43	0.49
1:B:143:LEU:HD11	1:B:171:ARG:HG2	1.95	0.48
1:C:154:ALA:O	1:C:157:PHE:HB2	2.12	0.48
1:C:51:TYR:HE1	1:D:93:TYR:HH	1.58	0.48
1:C:171:ARG:NH1	1:C:180:ARG:NE	2.61	0.48
1:C:157:PHE:O	1:C:162:PHE:HB2	2.13	0.48
1:B:163:ARG:HH21	1:B:337:ASP:HA	1.78	0.48
1:B:190:ASP:O	1:B:191:SER:HB2	2.13	0.48
1:D:33:LEU:HG	1:D:34:ILE:N	2.27	0.48
1:C:171:ARG:HH12	1:C:180:ARG:CZ	2.26	0.48
1:A:6:ARG:HH12	1:A:8:ARG:HD3	1.77	0.48
1:C:51:TYR:HE2	1:C:301:VAL:HG13	1.79	0.48
1:D:193:PRO:C	1:D:195:MSE:CE	2.82	0.48
1:C:138:ALA:HB3	1:C:165:ILE:CD1	2.44	0.48
1:A:109:ILE:HD13	1:A:304:LEU:HD13	1.96	0.48
1:D:51:TYR:O	1:D:52:THR:HG23	2.14	0.48
1:A:19:ARG:NE	1:A:29:MSE:CB	2.76	0.48
1:A:151:LYS:O	1:A:154:ALA:HB3	2.14	0.48
1:A:20:LEU:HB3	1:A:141:ILE:CB	2.36	0.47
1:D:14:ILE:CG2	1:D:15:PRO:HD2	2.44	0.47
1:C:93:TYR:OH	1:D:51:TYR:HE1	1.97	0.47
1:B:87:MSE:SE	1:B:107:VAL:HG11	2.64	0.47
1:C:12:TYR:O	1:C:32:GLU:HA	2.14	0.47
1:C:165:ILE:HG12	1:C:192:PHE:HE2	1.79	0.47
1:A:32:GLU:HB2	1:A:52:THR:OG1	2.13	0.47
1:C:124:TRP:HB2	1:C:310:ALA:HB3	1.96	0.47
1:C:123:LEU:HA	1:C:126:LEU:HB3	1.96	0.47
1:B:31:PHE:HB2	1:B:52:THR:O	2.14	0.47
1:C:149:ASP:O	1:C:152:THR:HB	2.15	0.47
1:D:270:GLU:HG3	1:D:297:THR:HG23	1.97	0.47
1:D:248:GLY:HA2	1:D:251:LEU:HG	1.96	0.47
1:A:124:TRP:CZ2	1:A:125:LYS:HE3	2.50	0.47
1:A:53:VAL:HG23	1:A:53:VAL:O	2.15	0.47
1:D:134:VAL:O	1:D:340:VAL:HG22	2.15	0.47
1:A:69:ARG:NH1	1:A:73:LEU:HD11	2.27	0.47
1:D:138:ALA:HB3	1:D:165:ILE:CD1	2.45	0.47
1:D:298:SER:OG	1:D:318:MSE:HG3	2.14	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:34:ILE:HD11	1:D:57:GLY:HA2	1.97	0.46
1:B:211:ALA:CB	1:B:241:SER:HB2	2.45	0.46
1:B:4:ILE:CG1	1:B:77:ALA:HB2	2.45	0.46
1:B:60:VAL:CA	1:B:63:MSE:HE2	2.38	0.46
1:D:171:ARG:HG2	1:D:177:ASP:OD1	2.14	0.46
1:D:173:ASP:OD1	1:D:175:LYS:HB3	2.15	0.46
1:C:13:ARG:NH1	1:C:362:VAL:HG21	2.30	0.46
1:D:105:SER:O	1:D:109:ILE:HG12	2.14	0.46
1:C:230:ASP:OD2	1:C:233:GLY:HA3	2.16	0.46
1:A:34:ILE:CD1	1:A:34:ILE:N	2.73	0.46
1:C:55:HIS:HE2	1:D:90:ARG:NH1	2.14	0.46
1:C:180:ARG:HH11	1:C:180:ARG:HG2	1.80	0.46
1:C:326:HIS:O	1:C:328:TYR:N	2.48	0.46
1:D:3:LEU:O	1:D:40:ASP:HA	2.15	0.46
1:A:113:ASP:O	1:A:117:ILE:HG13	2.16	0.46
1:A:93:TYR:O	1:A:94:ALA:HB3	2.16	0.46
1:A:303:ASP:HA	1:A:333:MSE:HE2	1.98	0.46
1:A:23:SER:CB	1:A:201:LYS:HB2	2.46	0.45
1:C:71:CYS:SG	1:C:90:ARG:NE	2.89	0.45
1:B:147:VAL:HG12	1:B:151:LYS:HE3	1.98	0.45
1:B:13:ARG:CG	1:B:13:ARG:NH1	2.79	0.45
1:B:51:TYR:CD2	1:B:51:TYR:N	2.82	0.45
1:B:328:TYR:CE1	1:B:356:ARG:HG2	2.50	0.45
1:C:37:ARG:NH1	1:C:354:PHE:CD1	2.84	0.45
1:A:147:VAL:O	1:A:150:LEU:HB3	2.16	0.45
1:C:55:HIS:HD2	1:D:90:ARG:HD3	1.82	0.45
1:B:127:PHE:O	1:B:283:ARG:NH2	2.49	0.45
1:D:168:LYS:HZ3	1:D:199:ASN:HD21	1.64	0.45
1:C:329:MSE:HE3	1:C:350:VAL:HG13	1.98	0.45
1:D:194:LEU:C	1:D:195:MSE:CE	2.84	0.45
1:A:315:ARG:HD2	1:A:315:ARG:C	2.36	0.45
1:C:20:LEU:C	1:C:21:THR:HG23	2.37	0.45
1:D:71:CYS:SG	1:D:90:ARG:NE	2.90	0.45
1:A:71:CYS:SG	1:A:90:ARG:NE	2.90	0.45
1:D:76:ASP:OD1	1:D:78:GLU:HB2	2.16	0.45
1:B:4:ILE:HG12	1:B:77:ALA:HB2	1.99	0.45
1:B:141:ILE:N	1:B:141:ILE:HD12	2.32	0.45
1:D:166:LYS:HA	1:D:195:MSE:O	2.17	0.45
1:C:117:ILE:HA	1:C:346:PRO:HB3	1.99	0.45
1:D:124:TRP:HB2	1:D:310:ALA:HB3	1.99	0.45
1:C:22:ASP:HA	1:C:143:LEU:HB2	1.98	0.44
1:B:151:LYS:O	1:B:154:ALA:HB3	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:166:LYS:HG3	1:C:195:MSE:O	2.18	0.44
1:A:211:ALA:HB1	1:A:241:SER:HB2	2.00	0.44
1:D:358:GLY:HA2	3:D:402:HOH:O	2.17	0.44
1:D:16:LEU:HD12	1:D:30:ASP:N	2.32	0.44
1:A:31:PHE:HB2	1:A:52:THR:O	2.18	0.44
1:C:326:HIS:C	1:C:328:TYR:H	2.21	0.44
1:C:301:VAL:O	1:C:301:VAL:HG23	2.18	0.44
1:B:11:LEU:HD22	1:B:58:ALA:HA	2.00	0.44
1:A:266:LEU:HD12	1:A:266:LEU:C	2.38	0.44
1:B:215:ALA:N	1:B:216:PRO:CD	2.81	0.44
1:D:32:GLU:HB3	1:D:52:THR:OG1	2.18	0.43
1:D:195:MSE:CE	1:D:220:HIS:CB	2.93	0.43
1:D:168:LYS:HD2	1:D:199:ASN:ND2	2.33	0.43
1:B:38:ILE:O	1:B:45:THR:HA	2.18	0.43
1:D:151:LYS:HE2	1:D:151:LYS:HB3	1.80	0.43
1:D:266:LEU:C	1:D:266:LEU:HD12	2.39	0.43
1:C:5:GLU:OE1	1:C:41:SER:HA	2.18	0.43
1:A:234:ASN:CB	1:A:265:SER:HB2	2.49	0.43
1:C:157:PHE:CD2	1:C:165:ILE:HD12	2.54	0.43
1:B:63:MSE:HE3	1:B:103:ALA:HB2	2.01	0.43
1:C:165:ILE:HG12	1:C:192:PHE:CE2	2.52	0.43
1:C:142:ASP:HB3	1:C:145:LEU:HD22	2.01	0.43
1:C:16:LEU:HD13	1:C:20:LEU:CD1	2.48	0.43
1:C:60:VAL:HA	1:C:63:MSE:HE2	2.00	0.43
1:B:32:GLU:O	1:B:51:TYR:HA	2.19	0.43
1:A:215:ALA:N	1:A:216:PRO:CD	2.82	0.43
1:B:47:LEU:O	1:B:348:HIS:CE1	2.56	0.43
1:D:47:LEU:O	1:D:348:HIS:HE1	2.01	0.43
1:B:321:HIS:O	1:B:321:HIS:CG	2.72	0.43
1:D:152:THR:HG22	1:D:156:ARG:HH11	1.82	0.43
1:D:34:ILE:CD1	1:D:60:VAL:HB	2.48	0.43
1:A:215:ALA:HB3	1:A:216:PRO:HD3	2.00	0.43
1:A:168:LYS:HD2	1:A:199:ASN:ND2	2.34	0.42
1:C:54:ASN:HB3	1:C:55:HIS:CE1	2.54	0.42
1:B:301:VAL:HG23	1:B:301:VAL:O	2.19	0.42
1:B:314:HIS:N	1:B:314:HIS:ND1	2.63	0.42
1:B:150:LEU:HG	1:B:184:LEU:CD1	2.45	0.42
1:A:6:ARG:HB2	1:A:39:GLU:HB3	2.01	0.42
1:B:306:VAL:HG13	1:B:307:HIS:N	2.34	0.42
1:C:302:HIS:HB3	1:C:319:GLU:O	2.19	0.42
1:B:16:LEU:HD13	1:B:31:PHE:H	1.84	0.42
1:D:40:ASP:OD2	1:D:114:LEU:HD11	2.19	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:171:ARG:HH12	1:B:180:ARG:NH1	2.17	0.42
1:A:32:GLU:O	1:A:51:TYR:HA	2.18	0.42
1:B:279:TYR:O	1:B:283:ARG:HG3	2.19	0.42
1:C:4:ILE:HG13	1:C:77:ALA:HB2	2.01	0.42
1:A:18:THR:HG22	1:A:19:ARG:N	2.34	0.42
1:D:224:GLU:OE2	1:D:249:GLU:HG2	2.19	0.42
1:D:31:PHE:HE1	1:D:51:TYR:CD1	2.38	0.42
1:B:53:VAL:O	1:B:54:ASN:HB2	2.19	0.42
1:C:123:LEU:CD2	1:C:126:LEU:HD23	2.50	0.42
1:B:325:LEU:HG	1:B:325:LEU:O	2.19	0.42
1:D:315:ARG:HD2	1:D:315:ARG:O	2.19	0.42
1:D:167:MSE:HE1	1:D:180:ARG:HB3	2.01	0.42
1:A:22:ASP:OD1	1:A:25:HIS:HD2	2.03	0.42
1:A:309:LEU:CD1	1:A:318:MSE:HB2	2.50	0.42
1:A:221:TRP:CD1	1:A:221:TRP:C	2.92	0.42
1:B:21:THR:H	1:B:141:ILE:HG22	1.84	0.42
1:A:25:HIS:HE1	3:B:477:HOH:O	2.00	0.42
1:D:34:ILE:HD13	1:D:60:VAL:CG1	2.50	0.42
1:A:166:LYS:HA	1:A:195:MSE:O	2.19	0.42
1:B:109:ILE:HG21	1:B:304:LEU:HD13	2.02	0.41
1:C:76:ASP:HB3	1:C:79:GLN:HE21	1.85	0.41
1:A:196:VAL:HG13	1:A:219:LEU:HD13	2.02	0.41
1:A:156:ARG:HG3	1:A:156:ARG:HH11	1.85	0.41
1:D:32:GLU:CG	1:D:57:GLY:HA3	2.49	0.41
1:A:20:LEU:CD1	1:A:141:ILE:HD12	2.50	0.41
1:B:60:VAL:HA	1:B:63:MSE:HE3	1.94	0.41
1:C:163:ARG:HH21	1:C:337:ASP:HA	1.85	0.41
1:A:298:SER:OG	1:A:318:MSE:HG3	2.20	0.41
1:C:165:ILE:HG22	1:C:166:LYS:N	2.35	0.41
1:C:192:PHE:HA	1:C:193:PRO:HD3	1.92	0.41
1:D:53:VAL:HG23	1:D:53:VAL:O	2.20	0.41
1:D:260:ALA:HA	1:D:265:SER:OG	2.20	0.41
1:D:32:GLU:CD	1:D:57:GLY:HA3	2.40	0.41
1:B:220:HIS:O	1:B:221:TRP:HB3	2.20	0.41
1:A:152:THR:O	1:A:155:ASP:HB2	2.21	0.41
1:D:316:THR:O	1:D:317:TYR:C	2.59	0.41
1:C:80:ILE:HG12	1:C:111:LEU:HB3	2.02	0.41
1:B:315:ARG:O	1:B:315:ARG:HD2	2.20	0.41
1:B:137:TYR:OH	1:B:319:GLU:HB2	2.21	0.41
1:D:272:ASP:CB	1:D:275:ASN:ND2	2.79	0.41
1:C:121:THR:O	1:C:346:PRO:HA	2.21	0.41
1:A:6:ARG:NH1	1:A:8:ARG:NE	2.69	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:6:ARG:NE	1:C:8:ARG:HE	2.18	0.41
1:A:11:LEU:HG	1:A:362:VAL:CG1	2.49	0.41
1:A:83:ILE:O	1:A:87:MSE:HG3	2.21	0.41
1:D:306:VAL:HG13	1:D:307:HIS:N	2.34	0.41
1:C:257:PHE:O	1:C:261:VAL:HG23	2.20	0.41
1:B:51:TYR:N	1:B:51:TYR:HD2	2.19	0.41
1:D:157:PHE:HB3	1:D:165:ILE:HD11	2.03	0.40
1:C:266:LEU:HD12	1:C:266:LEU:C	2.41	0.40
1:A:13:ARG:NE	3:A:445:HOH:O	2.54	0.40
1:C:83:ILE:HG21	1:C:111:LEU:HD11	2.03	0.40
1:D:123:LEU:HG	1:D:347:GLY:C	2.42	0.40
1:D:194:LEU:N	1:D:195:MSE:CE	2.83	0.40
1:A:29:MSE:O	1:A:30:ASP:HB3	2.22	0.40
1:C:67:ASP:CG	1:D:56:GLY:H	2.24	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	353/371 (95%)	336 (95%)	17 (5%)	0	100	100
1	B	353/371 (95%)	332 (94%)	19 (5%)	2 (1%)	33	39
1	C	353/371 (95%)	328 (93%)	21 (6%)	4 (1%)	21	21
1	D	328/371 (88%)	289 (88%)	35 (11%)	4 (1%)	19	19
All	All	1387/1484 (94%)	1285 (93%)	92 (7%)	10 (1%)	30	34

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	191	SER
1	C	327	ALA
1	C	361	ALA
1	D	191	SER

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Mol	Chain	Res	Type
1	D	190	ASP
1	C	27	ALA
1	B	301	VAL
1	C	344	ASP
1	D	301	VAL
1	D	56	GLY

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	283/281 (101%)	272 (96%)	11 (4%)	43	57
1	B	283/281 (101%)	276 (98%)	7 (2%)	60	77
1	C	283/281 (101%)	274 (97%)	9 (3%)	51	67
1	D	265/281 (94%)	259 (98%)	6 (2%)	63	80
All	All	1114/1124 (99%)	1081 (97%)	33 (3%)	53	70

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	19	ARG
1	A	29	MSE
1	A	195	MSE
1	A	221	TRP
1	A	229	ASP
1	A	288	LEU
1	A	292	ASN
1	A	315	ARG
1	A	326	HIS
1	A	362	VAL
1	B	28	MSE
1	B	30	ASP
1	B	51	TYR
1	B	195	MSE
1	B	221	TRP

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Mol	Chain	Res	Type
1	B	276	ILE
1	B	315	ARG
1	C	25	HIS
1	C	29	MSE
1	C	195	MSE
1	C	221	TRP
1	C	229	ASP
1	C	292	ASN
1	C	315	ARG
1	C	321	HIS
1	C	325	LEU
1	D	15	PRO
1	D	149	ASP
1	D	186	GLU
1	D	195	MSE
1	D	221	TRP
1	D	315	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	HIS
1	A	54	ASN
1	A	199	ASN
1	A	299	HIS
1	A	348	HIS
1	B	153	GLN
1	B	199	ASN
1	B	321	HIS
1	B	348	HIS
1	C	54	ASN
1	C	79	GLN
1	C	92	HIS
1	C	199	ASN
1	C	299	HIS
1	C	348	HIS
1	D	54	ASN
1	D	79	GLN
1	D	153	GLN
1	D	199	ASN
1	D	250	ASN
1	D	348	HIS

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 ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	357/371 (96%)	0.06	15 (4%) 35 45	18, 29, 55, 77	0
1	B	357/371 (96%)	0.11	4 (1%) 77 85	19, 32, 53, 68	0
1	C	357/371 (96%)	0.68	37 (10%) 7 11	24, 42, 63, 81	0
1	D	334/371 (90%)	0.33	19 (5%) 23 32	25, 41, 56, 65	0
All	All	1405/1484 (94%)	0.29	75 (5%) 25 35	18, 36, 57, 81	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	25	HIS	9.5
1	C	17	PRO	6.6
1	C	20	LEU	6.5
1	C	362	VAL	6.0
1	A	20	LEU	5.4
1	C	361	ALA	5.0
1	A	24	THR	5.0
1	C	26	GLY	4.6
1	D	360	LEU	4.5
1	C	18	THR	4.4
1	D	31	PHE	4.3
1	C	27	ALA	3.9
1	A	26	GLY	3.8
1	A	362	VAL	3.6
1	C	358	GLY	3.5
1	C	29	MSE	3.2
1	D	6	ARG	3.2
1	A	18	THR	3.2
1	D	359	ARG	3.2
1	D	161	GLY	3.2
1	B	16	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	38	ILE	3.2
1	A	25	HIS	3.2
1	B	19	ARG	3.2
1	C	359	ARG	3.1
1	A	359	ARG	3.1
1	A	17	PRO	2.9
1	C	155	ASP	2.9
1	C	152	THR	2.8
1	C	165	ILE	2.8
1	C	43	GLY	2.7
1	C	16	LEU	2.7
1	A	321	HIS	2.7
1	A	29	MSE	2.7
1	C	3	LEU	2.6
1	D	357	LEU	2.6
1	C	23	SER	2.6
1	D	120	ARG	2.6
1	C	186	GLU	2.6
1	D	186	GLU	2.6
1	C	150	LEU	2.6
1	A	21	THR	2.5
1	C	145	LEU	2.5
1	D	321	HIS	2.5
1	C	19	ARG	2.5
1	C	360	LEU	2.4
1	A	31	PHE	2.4
1	C	118	ARG	2.4
1	C	159	ALA	2.4
1	D	16	LEU	2.4
1	C	330	ALA	2.3
1	D	356	ARG	2.3
1	C	327	ALA	2.3
1	D	7	VAL	2.2
1	B	17	PRO	2.2
1	A	360	LEU	2.2
1	B	6	ARG	2.2
1	C	355	GLU	2.2
1	C	188	LEU	2.2
1	C	153	GLN	2.1
1	D	358	GLY	2.1
1	C	22	ASP	2.1
1	C	28	MSE	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	337	ASP	2.1
1	C	21	THR	2.1
1	A	19	ARG	2.1
1	C	357	LEU	2.1
1	D	158	LEU	2.1
1	C	6	ARG	2.0
1	C	120	ARG	2.0
1	D	157	PHE	2.0
1	A	22	ASP	2.0
1	D	150	LEU	2.0
1	C	7	VAL	2.0
1	D	336	THR	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	MG	A	400	1/1	0.10	-0.55	31,31,31,31	0
2	MG	C	400	1/1	0.10	-1.11	42,42,42,42	0
2	MG	B	400	1/1	0.10	-1.29	31,31,31,31	0
2	MG	D	400	1/1	0.07	-2.46	40,40,40,40	0

6.5 Other polymers ⓘ

There are no such residues in this entry.