



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:23 PM UTC

PDB ID : 3CK5 / pdb_00003ck5
Title : Crystal structure of a racemase from *Streptomyces coelicolor* A3(2) with bound magnesium
Authors : Rao, K.N.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2008-03-14
Resolution : 2.30 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

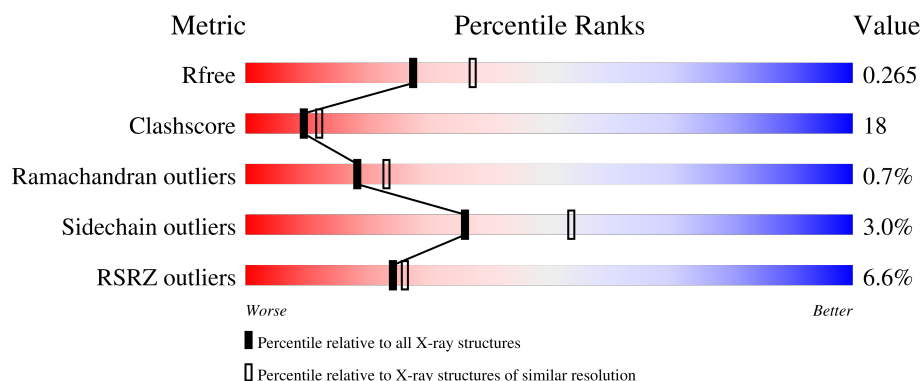
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	A	371	5% 70% 23% • •
1	B	371	4% 68% 25% • •
1	C	371	8% 59% 35% • •
1	D	371	6% 58% 30% • 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11164 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 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 (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	D	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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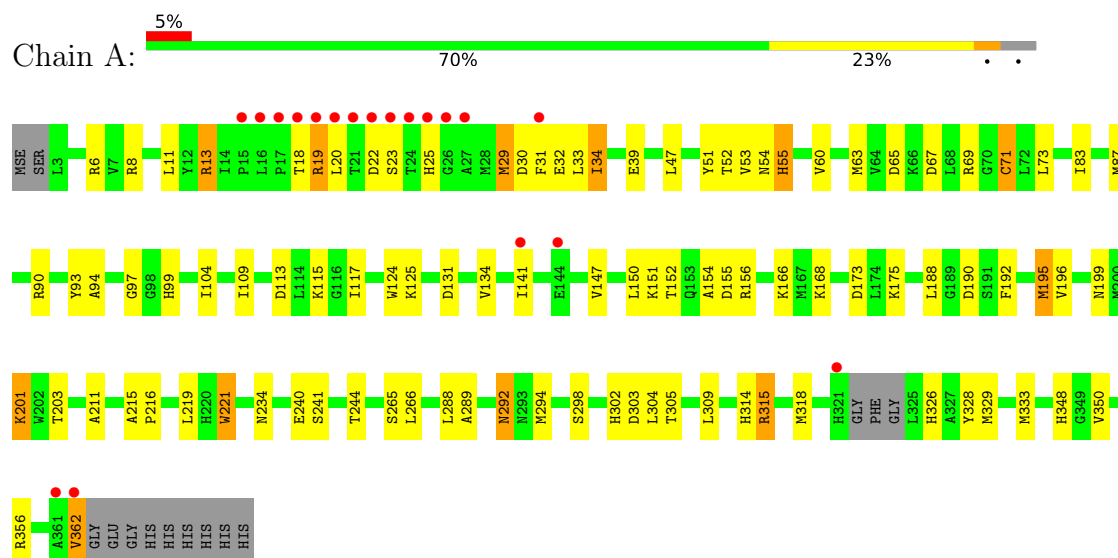
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	113	Total 113	O 113	0	0
3	C	73	Total 73	O 73	0	0
3	D	51	Total 51	O 51	0	0

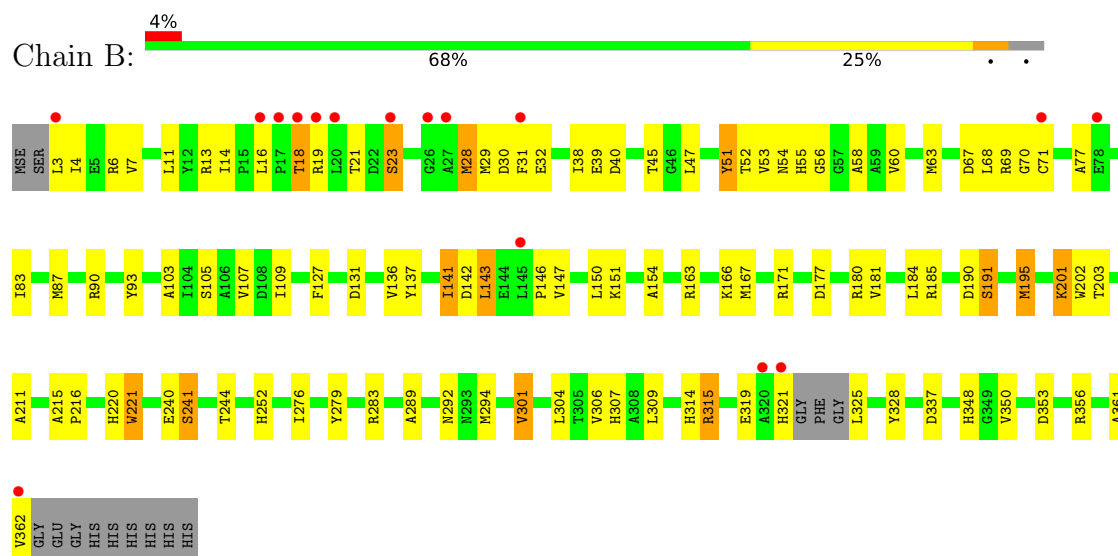
3 Residue-property plots

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

• Molecule 1: Putative racemase

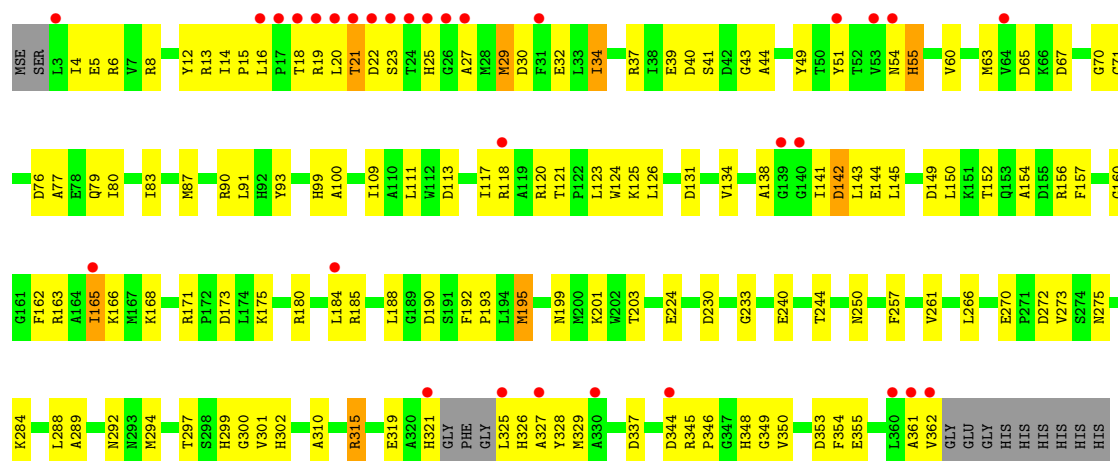


• Molecule 1: Putative racemase

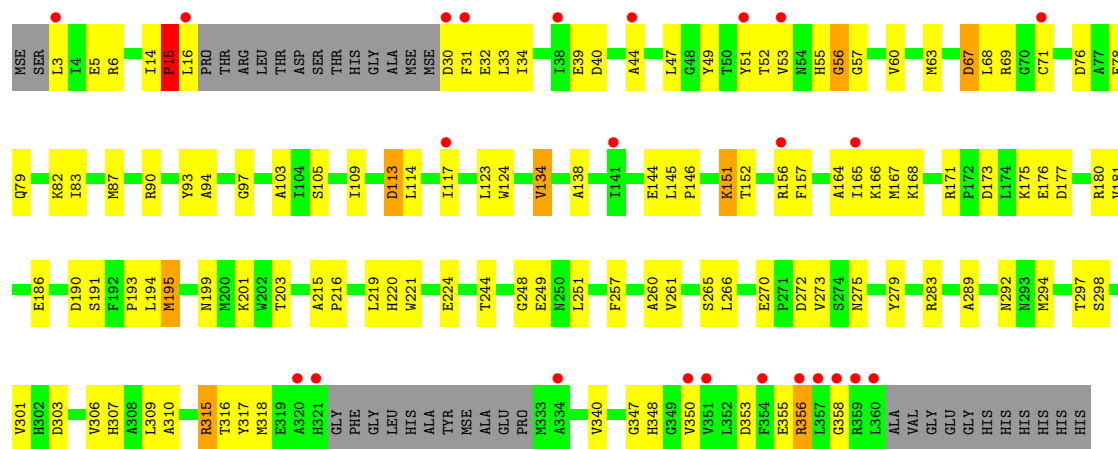


• Molecule 1: Putative racemase





• Molecule 1: Putative racemase



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 , 0.265	Depositor DCC
R_{free} test set	1876 reflections (2.37%)	wwPDB-VP
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 , 32.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	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.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.51	0/2802	0.97	11/3794 (0.3%)
1	B	0.51	0/2802	1.00	14/3794 (0.4%)
1	C	0.50	0/2802	1.04	19/3794 (0.5%)
1	D	0.47	0/2628	1.01	11/3561 (0.3%)
All	All	0.50	0/11034	1.01	55/14943 (0.4%)

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.

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	23	SER	N-CA-C	8.38	121.54	111.82
1	A	240	GLU	N-CA-C	7.86	123.93	113.72
1	A	203	THR	N-CA-C	-7.68	99.46	110.59
1	D	15	PRO	N-CA-C	7.62	124.49	111.77
1	C	201	LYS	N-CA-C	7.46	119.19	111.14
1	C	244	THR	N-CA-C	-7.37	99.37	110.28
1	A	190	ASP	N-CA-C	7.32	119.89	111.11
1	B	142	ASP	N-CA-C	7.16	121.97	111.87
1	B	240	GLU	N-CA-C	7.13	122.27	113.50
1	A	201	LYS	N-CA-C	7.05	120.00	111.82
1	C	134	VAL	N-CA-C	7.02	115.54	107.89
1	D	55	HIS	N-CA-C	6.85	117.32	108.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	203	THR	N-CA-C	-6.71	101.87	110.53
1	C	300	GLY	N-CA-C	6.59	120.99	112.68
1	B	201	LYS	N-CA-C	6.38	118.31	111.36
1	B	203	THR	N-CA-C	-6.34	101.40	110.59
1	A	65	ASP	N-CA-C	6.30	118.14	111.28
1	B	143	LEU	N-CA-C	6.29	117.80	111.07
1	C	23	SER	N-CA-C	6.27	121.72	111.37
1	C	240	GLU	N-CA-C	6.22	121.15	113.50
1	D	244	THR	N-CA-C	-6.12	100.77	109.96
1	C	160	GLY	N-CA-C	-6.07	107.99	114.67
1	C	203	THR	N-CA-C	-6.06	101.99	110.35
1	D	134	VAL	N-CA-C	6.04	115.18	108.05
1	D	201	LYS	N-CA-C	6.03	118.63	111.33
1	B	244	THR	N-CA-C	-5.97	101.01	109.96
1	A	55	HIS	N-CA-C	5.91	117.57	109.18
1	B	136	VAL	N-CA-C	5.89	117.19	108.65
1	A	244	THR	N-CA-C	-5.86	101.16	109.96
1	C	142	ASP	N-CA-C	5.85	117.92	110.61
1	B	18	THR	N-CA-C	-5.83	102.96	110.19
1	B	241	SER	N-CA-C	5.79	120.35	113.28
1	C	188	LEU	N-CA-C	5.75	120.36	113.23
1	B	252	HIS	N-CA-C	5.66	120.18	112.88
1	A	97	GLY	N-CA-C	-5.63	105.10	111.85
1	D	113	ASP	N-CA-C	-5.58	105.10	111.07
1	C	55	HIS	N-CA-C	5.56	117.86	109.41
1	C	165	ILE	N-CA-C	5.54	117.38	108.85
1	C	185	ARG	N-CA-C	-5.43	105.44	111.36
1	C	43	GLY	N-CA-C	5.41	121.52	115.08
1	D	49	TYR	N-CA-C	5.41	117.48	108.99
1	A	134	VAL	N-CA-C	5.41	113.78	107.89
1	C	21	THR	N-CA-C	5.41	117.06	108.79
1	C	190	ASP	N-CA-C	5.39	118.65	111.75
1	B	202	TRP	N-CA-C	5.38	118.49	110.52
1	D	356	ARG	N-CA-C	-5.36	105.55	111.71
1	A	71	CYS	N-CA-C	-5.28	106.89	113.38
1	C	125	LYS	N-CA-C	-5.24	106.57	113.12
1	B	185	ARG	N-CA-C	-5.23	105.47	111.07
1	C	49	TYR	N-CA-C	5.19	115.38	108.38
1	D	97	GLY	N-CA-C	-5.14	105.72	111.63
1	B	141	ILE	N-CA-C	-5.14	98.65	109.34
1	C	65	ASP	N-CA-C	5.09	116.52	111.07
1	D	67	ASP	N-CA-C	5.06	117.92	111.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(^o)	Ideal(^o)
1	A	314	HIS	N-CA-C	5.04	119.06	113.02

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	93	TYR	Sidechain

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2749	0	2711	82	0
1	B	2749	0	2711	92	0
1	C	2749	0	2711	115	0
1	D	2577	0	2543	100	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	383	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 18.

All (383) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:A:19:ARG:NE	1:A:29:MSE:HB3	2.00	0.76
1:D:113:ASP:O	1:D:117:ILE:HG12	1.84	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:195:MSE:HE2	1:D:220:HIS:CB	2.21	0.71
1:D:69:ARG:HG3	1:D:69:ARG:HH11	1.55	0.71
1:C:113:ASP:O	1:C:117:ILE:HG13	1.91	0.70
1:B:13:ARG:HG3	1:B:13:ARG:HH11	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:D:315:ARG:HD2	1:D:315:ARG:C	2.17	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:D:194:LEU:C	1:D:195:MSE:HE3	2.20	0.66
1:A:69:ARG:HH11	1:A:73:LEU:HD11	1.60	0.66
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: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:53:VAL:O	1:A:54:ASN:HB2	1.98	0.64
1:A:289:ALA:HA	1:A:294:MSE:CE	2.28	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:A:55:HIS:HB3	1:B:67:ASP:OD1	1.98	0.64
1:C:18:THR:HG22	1:C:19:ARG:H	1.63	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:18:THR:HG23	1:B:19:ARG:HG3	1.81	0.62
1:B:70:GLY:CA	1:B:90:ARG:HH12	2.11	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:21:THR:HG21	1:C:144:GLU:CD	2.25	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: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:B:315:ARG:HD2	1:B:315:ARG:C	2.25	0.61
1:D:194:LEU:N	1:D:195:MSE:HE3	2.16	0.61
1:A:309:LEU:O	1:A:315:ARG:HG2	2.01	0.61
1:C:70:GLY:HA3	1:C:90:ARG:NH1	2.16	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:HG2	1:B:315:ARG:HH11	1.66	0.58
1:C:34:ILE:N	1:C:34:ILE:HD12	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
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:63:MSE:HE1	1:C:99:HIS:HB3	1.85	0.57
1:D:193:PRO:C	1:D:195:MSE:HE3	2.28	0.57
1:A:33:LEU:C	1:A:34:ILE:HD12	2.30	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:C:54:ASN:HB3	1:C:55:HIS:ND1	2.21	0.56
1:C:353:ASP:OD1	1:C:355:GLU:HB2	2.05	0.56
1:D:138:ALA:HB3	1:D:165:ILE:HD12	1.87	0.56
1:C:91:LEU:HD13	1:C:100:ALA:HB1	1.87	0.56
1:D:356:ARG:C	1:D:358:GLY:H	2.14	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:B:328:TYR:HE1	1:B:356:ARG:HG2	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:152:THR:HG22	1:D:156:ARG:NH1	2.20	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:HD2	1:C:315:ARG:O	2.05	0.55
1:C:16:LEU:HD13	1:C:20:LEU:HD12	1.89	0.55
1:D:32:GLU:HG2	1:D:57:GLY:HA3	1.87	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:A:211:ALA:CB	1:A:241:SER:HB2	2.37	0.54
1:B:3:LEU:O	1:B:40:ASP:HB2	2.08	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:A:47:LEU:O	1:A:348:HIS:HE1	1.91	0.54
1:B:166:LYS:HA	1:B:195:MSE:O	2.08	0.54
1:C:118:ARG:HG2	1:C:118:ARG:HH11	1.73	0.54
1:D:14:ILE:HG23	1:D:15:PRO:HD2	1.90	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:A:302:HIS:HB2	1:A:318:MSE:HE3	1.91	0.53
1:B:14:ILE:HD12	1:B:31:PHE:CE1	2.44	0.53
1:B:83:ILE:HG22	1:B:87:MSE:HE2	1.91	0.53
1:A:188:LEU:HD13	1:A:192:PHE:CD2	2.44	0.53
1:B:69:ARG:HG3	1:B:69:ARG:HH11	1.74	0.53
1:D:193:PRO:C	1:D:195:MSE:CE	2.82	0.53
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:HH11	1:A:13:ARG:HB2	1.75	0.52
1:B:56:GLY:O	1:B:60:VAL:HG23	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:270:GLU:HG3	1:C:297:THR:CG2	2.36	0.52
1:A:302:HIS:HA	1:A:305:THR:HB	1.91	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:D:194:LEU:C	1:D:195:MSE:CE	2.84	0.51
1:A:6:ARG:NH1	1:A:8:ARG:CD	2.73	0.51
1:B:137:TYR:CZ	1:B:319:GLU:HB2	2.46	0.51
1:D:181:VAL:HG11	1:D:219:LEU:HD21	1.93	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:A:67:ASP:OD1	1:B:55:HIS:HB3	2.11	0.50
1:B:14:ILE:HD12	1:B:31:PHE:HE1	1.77	0.50
1:C:272:ASP:HB3	1:C:275:ASN:ND2	2.26	0.50
1:D:272:ASP:HB3	1:D:275:ASN:HD21	1.76	0.50
1:C:345:ARG:NH1	1:C:349:GLY:O	2.45	0.50
1:B:361:ALA:HB1	3:B:519:HOH:O	2.10	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: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:A:315:ARG:C	1:A:315:ARG:HD2	2.36	0.49
1:C:13:ARG:HG3	1:C:13:ARG:HH11	1.76	0.49
1:C:20:LEU:C	1:C:21:THR:HG23	2.37	0.49
1:D:34:ILE:HD13	1:D:60:VAL:HB	1.94	0.49
1:D:257:PHE:O	1:D:261:VAL:HG23	2.12	0.49
1:C:171:ARG:HH12	1:C:180:ARG:NE	2.09	0.49
1:D:83:ILE:O	1:D:87:MSE:HG3	2.12	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:171:ARG:NH1	1:C:180:ARG:NE	2.61	0.48
1:A:266:LEU:C	1:A:266:LEU:HD12	2.38	0.48
1:B:163:ARG:HH21	1:B:337:ASP:HA	1.78	0.48
1:C:157:PHE:O	1:C:162:PHE:HB2	2.13	0.48
1:B:190:ASP:O	1:B:191:SER:HB2	2.13	0.48
1:C:171:ARG:HH12	1:C:180:ARG:CZ	2.26	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:326:HIS:C	1:C:328:TYR:H	2.21	0.48
1:D:33:LEU:HG	1:D:34:ILE:N	2.27	0.48
1:D:266:LEU:C	1:D:266:LEU:HD12	2.39	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:A:109:ILE:HD13	1:A:304:LEU:HD13	1.96	0.48
1:C:138:ALA:HB3	1:C:165:ILE:CD1	2.44	0.48
1:A:19:ARG:NE	1:A:29:MSE:CB	2.76	0.48
1:D:51:TYR:O	1:D:52:THR:HG23	2.14	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: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:93:TYR:OH	1:D:51:TYR:HE1	1.97	0.47
1:A:32:GLU:HB2	1:A:52:THR:OG1	2.13	0.47
1:C:123:LEU:HA	1:C:126:LEU:HB3	1.96	0.47
1:C:124:TRP:HB2	1:C:310:ALA:HB3	1.96	0.47
1:C:165:ILE:HG12	1:C:192:PHE:HE2	1.79	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:A:124:TRP:CZ2	1:A:125:LYS:HE3	2.50	0.47
1:D:248:GLY:HA2	1:D:251:LEU:HG	1.96	0.47
1:A:53:VAL:O	1:A:53:VAL:HG23	2.15	0.47
1:A:221:TRP:CD1	1:A:221:TRP:C	2.92	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
1:B:211:ALA:CB	1:B:241:SER:HB2	2.45	0.46
1:D:34:ILE:HD11	1:D:57:GLY:HA2	1.97	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:C:13:ARG:NH1	1:C:362:VAL:HG21	2.30	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:230:ASP:OD2	1:C:233:GLY:HA3	2.16	0.46
1:D:105:SER:O	1:D:109:ILE:HG12	2.14	0.46
1:C:266:LEU:HD12	1:C:266:LEU:C	2.41	0.46
1:D:32:GLU:CD	1:D:57:GLY:HA3	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:ILE:N	1:A:34:ILE:CD1	2.73	0.46
1:C:55:HIS:HE2	1:D:90:ARG:NH1	2.14	0.46
1:C:180:ARG:HG2	1:C:180:ARG:HH11	1.80	0.46
1:C:326:HIS:O	1:C:328:TYR:N	2.48	0.46
1:D:316:THR:O	1:D:317:TYR:C	2.59	0.46
1:A:93:TYR:O	1:A:94:ALA:HB3	2.16	0.46
1:A:113:ASP:O	1:A:117:ILE:HG13	2.16	0.46
1:A:303:ASP:HA	1:A:333:MSE:HE2	1.98	0.46
1:C:67:ASP:CG	1:D:56:GLY:H	2.24	0.46
1:D:3:LEU:O	1:D:40:ASP:HA	2.15	0.46
1:A:23:SER:CB	1:A:201:LYS:HB2	2.46	0.45
1:B:147:VAL:HG12	1:B:151:LYS:HE3	1.98	0.45
1:C:71:CYS:SG	1:C:90:ARG:NE	2.89	0.45
1:B:13:ARG:NH1	1:B:13:ARG:CG	2.79	0.45
1:B:51:TYR:N	1:B:51:TYR:CD2	2.82	0.45
1:A:147:VAL:O	1:A:150:LEU:HB3	2.16	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:B:127:PHE:O	1:B:283:ARG:NH2	2.49	0.45
1:C:55:HIS:HD2	1:D:90:ARG:HD3	1.82	0.45
1:C:329:MSE:HE3	1:C:350:VAL:HG13	1.98	0.45
1:D:123:LEU:HG	1:D:347:GLY:C	2.42	0.45
1:A:71:CYS:SG	1:A:90:ARG:NE	2.90	0.45
1:D:71:CYS:SG	1:D:90:ARG:NE	2.90	0.45
1:B:4:ILE:HG12	1:B:77:ALA:HB2	1.99	0.45
1:D:76:ASP:OD1	1:D:78:GLU:HB2	2.16	0.45
1:B:40:ASP:OD1	1:B:40:ASP:C	2.59	0.45
1:B:141:ILE:N	1:B:141:ILE:HD12	2.32	0.45
1:C:22:ASP:OD1	1:C:22:ASP:C	2.60	0.45
1:C:54:ASN:C	1:C:55:HIS:ND1	2.75	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:D:166:LYS:HA	1:D:195:MSE:O	2.17	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
1:A:211:ALA:HB1	1:A:241:SER:HB2	2.00	0.44
1:C:166:LYS:HG3	1:C:195:MSE:O	2.18	0.44
1:C:224:GLU:CD	1:C:250:ASN:HD22	2.26	0.44
1:D:358:GLY:HA2	3:D:402:HOH:O	2.17	0.44
1:A:31:PHE:HB2	1:A:52:THR:O	2.18	0.44
1:D:16:LEU:HD12	1:D:30:ASP:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:301:VAL:HG23	1:C:301:VAL:O	2.18	0.44
1:B:11:LEU:HD22	1:B:58:ALA:HA	2.00	0.44
1:B:215:ALA:N	1:B:216:PRO:CD	2.81	0.44
1:B:38:ILE:O	1:B:45:THR:HA	2.18	0.43
1:D:32:GLU:HB3	1:D:52:THR:OG1	2.18	0.43
1:D:151:LYS:HE2	1:D:151:LYS:HB3	1.80	0.43
1:D:168:LYS:HD2	1:D:199:ASN:ND2	2.33	0.43
1:D:195:MSE:CE	1:D:220:HIS:CB	2.93	0.43
1:A:234:ASN:CB	1:A:265:SER:HB2	2.49	0.43
1:C:5:GLU:OE1	1:C:41:SER:HA	2.18	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:142:ASP:HB3	1:C:145:LEU:HD22	2.01	0.43
1:C:165:ILE:HG12	1:C:192:PHE:CE2	2.52	0.43
1:B:32:GLU:O	1:B:51:TYR:HA	2.19	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: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:B:321:HIS:O	1:B:321:HIS:CG	2.72	0.43
1:D:47:LEU:O	1:D:348:HIS:HE1	2.01	0.43
1:D:152:THR:HG22	1:D:156:ARG:HH11	1.82	0.43
1:A:215:ALA:HB3	1:A:216:PRO:HD3	2.00	0.43
1:D:34:ILE:CD1	1:D:60:VAL:HB	2.48	0.43
1:A:168:LYS:HD2	1:A:199:ASN:ND2	2.34	0.42
1:B:301:VAL:HG23	1:B:301:VAL:O	2.19	0.42
1:B:314:HIS:ND1	1:B:314:HIS:N	2.63	0.42
1:C:54:ASN:HB3	1:C:55:HIS:CE1	2.54	0.42
1:A:6:ARG:HB2	1:A:39:GLU:HB3	2.01	0.42
1:B:150:LEU:HG	1:B:184:LEU:CD1	2.45	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:A:32:GLU:O	1:A:51:TYR:HA	2.18	0.42
1:B:16:LEU:HD13	1:B:31:PHE:H	1.84	0.42
1:B:171:ARG:HH12	1:B:180:ARG:NH1	2.17	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:D:40:ASP:OD2	1:D:114:LEU:HD11	2.19	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:B:53:VAL:O	1:B:54:ASN:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:123:LEU:CD2	1:C:126:LEU:HD23	2.50	0.42
1:D:31:PHE:HE1	1:D:51:TYR:CD1	2.38	0.42
1:A:22:ASP:OD1	1:A:25:HIS:HD2	2.03	0.42
1:D:167:MSE:HE1	1:D:180:ARG:HB3	2.01	0.42
1:D:315:ARG:HD2	1:D:315:ARG:O	2.19	0.42
1:A:309:LEU:CD1	1:A:318:MSE:HB2	2.50	0.42
1:B:325:LEU:O	1:B:325:LEU:HG	2.19	0.42
1:D:221:TRP:CD1	1:D:221:TRP:C	2.98	0.42
1:A:25:HIS:HE1	3:B:477:HOH:O	2.00	0.42
1:B:21:THR:H	1:B:141:ILE:HG22	1.84	0.42
1:A:166:LYS:HA	1:A:195:MSE:O	2.19	0.42
1:D:34:ILE:HD13	1:D:60:VAL:CG1	2.50	0.42
1:A:196:VAL:HG13	1:A:219:LEU:HD13	2.02	0.41
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:156:ARG:HG3	1:A:156:ARG:HH11	1.85	0.41
1:D:68:LEU:O	1:D:69:ARG:C	2.63	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:D:32:GLU:CG	1:D:57:GLY:HA3	2.49	0.41
1:D:168:LYS:HZ2	1:D:199:ASN:HD21	1.66	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:260:ALA:HA	1:D:265:SER:OG	2.20	0.41
1:D:356:ARG:C	1:D:358:GLY:N	2.79	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:B:221:TRP:CD1	1:B:221:TRP:C	2.98	0.41
1:B:137:TYR:OH	1:B:319:GLU:HB2	2.21	0.41
1:B:315:ARG:HD2	1:B:315:ARG:O	2.20	0.41
1:C:80:ILE:HG12	1:C:111:LEU:HB3	2.02	0.41
1:D:53:VAL:O	1:D:53:VAL:HG23	2.20	0.41
1:A:6:ARG:NH1	1:A:8:ARG:NE	2.69	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:B:146:PRO:O	1:B:147:VAL:C	2.63	0.41
1:C:6:ARG:NE	1:C:8:ARG:HE	2.18	0.41
1:C:121:THR:O	1:C:346:PRO:HA	2.21	0.41
1:C:257:PHE:O	1:C:261:VAL:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:272:ASP:CB	1:D:275:ASN:ND2	2.79	0.41
1:D:306:VAL:HG13	1:D:307:HIS:N	2.34	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:A:13:ARG:NE	3:A:445:HOH:O	2.54	0.40
1:B:68:LEU:O	1:B:69:ARG:C	2.64	0.40
1:B:353:ASP:CG	1:B:356:ARG:NH1	2.79	0.40
1:C:83:ILE:HG21	1:C:111:LEU:HD11	2.03	0.40
1:A:29:MSE:O	1:A:30:ASP:HB3	2.22	0.40
1:C:16:LEU:C	1:C:18:THR:H	2.29	0.40
1:D:194:LEU:N	1:D:195:MSE:CE	2.83	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 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	
1	A	353/371 (95%)	336 (95%)	17 (5%)	0	100	100
1	B	353/371 (95%)	332 (94%)	19 (5%)	2 (1%)	21	27
1	C	353/371 (95%)	328 (93%)	21 (6%)	4 (1%)	11	13
1	D	328/371 (88%)	289 (88%)	35 (11%)	4 (1%)	10	12
All	All	1387/1484 (94%)	1285 (93%)	92 (7%)	10 (1%)	18	23

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%)	28	43
1	B	283/281 (101%)	276 (98%)	7 (2%)	42	60
1	C	283/281 (101%)	275 (97%)	8 (3%)	38	56
1	D	265/281 (94%)	258 (97%)	7 (3%)	40	59
All	All	1114/1124 (99%)	1081 (97%)	33 (3%)	36	53

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	34	ILE
1	A	195	MSE
1	A	221	TRP
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	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	B	362	VAL
1	C	25	HIS
1	C	29	MSE
1	C	34	ILE
1	C	120	ARG
1	C	195	MSE
1	C	292	ASN
1	C	315	ARG
1	C	325	LEU
1	D	5	GLU
1	D	15	PRO
1	D	82	LYS
1	D	151	LYS
1	D	186	GLU
1	D	195	MSE
1	D	315	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	HIS
1	A	54	ASN
1	A	79	GLN
1	A	153	GLN
1	A	199	ASN
1	A	292	ASN
1	A	299	HIS
1	A	348	HIS
1	B	153	GLN
1	B	199	ASN
1	B	234	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	220	HIS
1	C	299	HIS
1	C	348	HIS

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Mol	Chain	Res	Type
1	D	54	ASN
1	D	79	GLN
1	D	153	GLN
1	D	187	HIS
1	D	199	ASN
1	D	348	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides 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.

No monomer is involved in short contacts.

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	346/371 (93%)	0.22	19 (5%)	30 32	18, 29, 54, 69	0
1	B	346/371 (93%)	0.27	16 (4%)	37 39	19, 33, 52, 63	0
1	C	346/371 (93%)	0.73	31 (8%)	15 16	24, 43, 63, 75	0
1	D	326/371 (87%)	0.68	24 (7%)	20 22	25, 41, 55, 65	0
All	All	1364/1484 (91%)	0.47	90 (6%)	24 26	18, 36, 57, 75	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	25	HIS	6.4
1	D	16	LEU	5.1
1	D	360	LEU	5.1
1	A	24	THR	5.0
1	C	362	VAL	4.9
1	C	17	PRO	4.8
1	C	3	LEU	4.8
1	D	31	PHE	4.4
1	B	16	LEU	4.4
1	A	321	HIS	4.4
1	C	26	GLY	4.3
1	C	27	ALA	4.2
1	C	23	SER	4.1
1	C	20	LEU	3.9
1	B	17	PRO	3.9
1	C	24	THR	3.9
1	D	321	HIS	3.7
1	B	31	PHE	3.6
1	A	25	HIS	3.6
1	A	31	PHE	3.5
1	C	18	THR	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	23	SER	3.4
1	C	325	LEU	3.4
1	A	17	PRO	3.3
1	D	53	VAL	3.3
1	A	20	LEU	3.3
1	D	359	ARG	3.3
1	A	18	THR	3.3
1	A	26	GLY	3.2
1	D	358	GLY	3.2
1	C	16	LEU	3.0
1	C	165	ILE	2.9
1	D	141	ILE	2.9
1	A	362	VAL	2.9
1	B	362	VAL	2.9
1	B	18	THR	2.9
1	C	51	TYR	2.8
1	A	27	ALA	2.8
1	A	22	ASP	2.8
1	C	54	ASN	2.8
1	D	3	LEU	2.8
1	C	330	ALA	2.8
1	D	165	ILE	2.8
1	A	16	LEU	2.8
1	D	356	ARG	2.7
1	D	320	ALA	2.6
1	B	26	GLY	2.6
1	B	19	ARG	2.6
1	C	361	ALA	2.6
1	C	64	VAL	2.6
1	A	15	PRO	2.6
1	A	21	THR	2.6
1	C	22	ASP	2.5
1	D	357	LEU	2.5
1	C	31	PHE	2.5
1	C	360	LEU	2.5
1	C	321	HIS	2.4
1	D	334	ALA	2.4
1	D	30	ASP	2.3
1	D	51	TYR	2.3
1	C	327	ALA	2.3
1	D	44	ALA	2.3
1	D	117	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	350	VAL	2.3
1	A	19	ARG	2.3
1	C	184	LEU	2.2
1	C	19	ARG	2.2
1	A	361	ALA	2.2
1	C	53	VAL	2.2
1	D	351	VAL	2.2
1	D	354	PHE	2.2
1	B	23	SER	2.2
1	C	140	GLY	2.2
1	B	27	ALA	2.2
1	C	139	GLY	2.1
1	B	20	LEU	2.1
1	D	38	ILE	2.1
1	B	71	CYS	2.1
1	C	21	THR	2.1
1	A	141	ILE	2.1
1	B	321	HIS	2.1
1	B	3	LEU	2.1
1	B	145	LEU	2.1
1	D	156	ARG	2.1
1	A	144	GLU	2.0
1	C	344	ASP	2.0
1	B	78	GLU	2.0
1	B	320	ALA	2.0
1	D	71	CYS	2.0
1	C	118	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	B-factors(Å ²)	Q<0.9
2	MG	A	400	1/1	0.95	0.05	31,31,31,31	0
2	MG	C	400	1/1	0.96	0.05	42,42,42,42	0
2	MG	B	400	1/1	0.98	0.03	31,31,31,31	0
2	MG	D	400	1/1	0.98	0.04	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.