



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:22 PM UTC

PDB ID : 2ZTC / pdb_00002ztc
Title : MtRuvA Form II
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Deposited on : 2008-10-01
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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

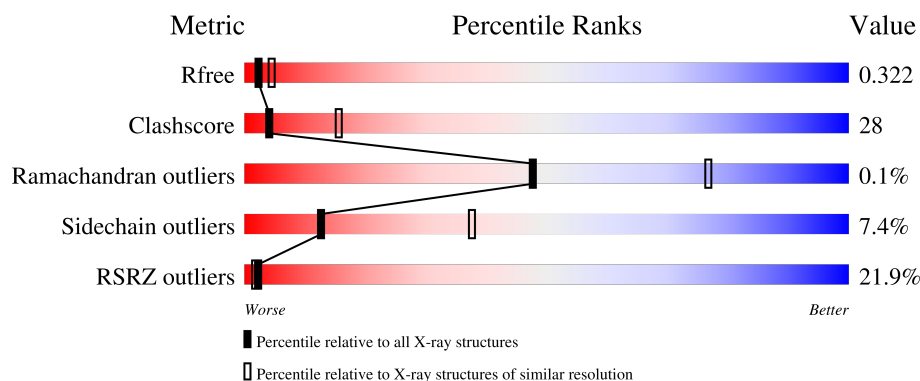
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	212	13% 52% 33% • 11%
1	B	212	30% 50% 35% • 12%
1	C	212	13% 49% 36% • 12%
1	D	212	21% 49% 36% • 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5372 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 Holliday junction ATP-dependent DNA helicase ruvA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	188	Total	C	N	O	S	0	0	0
			1329	831	240	253	5			
1	B	187	Total	C	N	O	S	0	0	0
			1324	828	239	252	5			
1	C	187	Total	C	N	O	S	0	0	0
			1324	828	239	252	5			
1	D	188	Total	C	N	O	S	0	0	0
			1323	828	237	253	5			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	expression tag	UNP P66744
A	-14	ALA	-	expression tag	UNP P66744
A	-13	SER	-	expression tag	UNP P66744
A	-12	MET	-	expression tag	UNP P66744
A	-11	THR	-	expression tag	UNP P66744
A	-10	GLY	-	expression tag	UNP P66744
A	-9	GLY	-	expression tag	UNP P66744
A	-8	GLN	-	expression tag	UNP P66744
A	-7	GLN	-	expression tag	UNP P66744
A	-6	MET	-	expression tag	UNP P66744
A	-5	GLY	-	expression tag	UNP P66744
A	-4	ARG	-	expression tag	UNP P66744
A	-3	GLY	-	expression tag	UNP P66744
A	-2	SER	-	expression tag	UNP P66744
A	-1	GLU	-	expression tag	UNP P66744
A	0	PHE	-	expression tag	UNP P66744
B	-15	MET	-	expression tag	UNP P66744
B	-14	ALA	-	expression tag	UNP P66744
B	-13	SER	-	expression tag	UNP P66744
B	-12	MET	-	expression tag	UNP P66744
B	-11	THR	-	expression tag	UNP P66744

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	GLY	-	expression tag	UNP P66744
B	-9	GLY	-	expression tag	UNP P66744
B	-8	GLN	-	expression tag	UNP P66744
B	-7	GLN	-	expression tag	UNP P66744
B	-6	MET	-	expression tag	UNP P66744
B	-5	GLY	-	expression tag	UNP P66744
B	-4	ARG	-	expression tag	UNP P66744
B	-3	GLY	-	expression tag	UNP P66744
B	-2	SER	-	expression tag	UNP P66744
B	-1	GLU	-	expression tag	UNP P66744
B	0	PHE	-	expression tag	UNP P66744
C	-15	MET	-	expression tag	UNP P66744
C	-14	ALA	-	expression tag	UNP P66744
C	-13	SER	-	expression tag	UNP P66744
C	-12	MET	-	expression tag	UNP P66744
C	-11	THR	-	expression tag	UNP P66744
C	-10	GLY	-	expression tag	UNP P66744
C	-9	GLY	-	expression tag	UNP P66744
C	-8	GLN	-	expression tag	UNP P66744
C	-7	GLN	-	expression tag	UNP P66744
C	-6	MET	-	expression tag	UNP P66744
C	-5	GLY	-	expression tag	UNP P66744
C	-4	ARG	-	expression tag	UNP P66744
C	-3	GLY	-	expression tag	UNP P66744
C	-2	SER	-	expression tag	UNP P66744
C	-1	GLU	-	expression tag	UNP P66744
C	0	PHE	-	expression tag	UNP P66744
D	-15	MET	-	expression tag	UNP P66744
D	-14	ALA	-	expression tag	UNP P66744
D	-13	SER	-	expression tag	UNP P66744
D	-12	MET	-	expression tag	UNP P66744
D	-11	THR	-	expression tag	UNP P66744
D	-10	GLY	-	expression tag	UNP P66744
D	-9	GLY	-	expression tag	UNP P66744
D	-8	GLN	-	expression tag	UNP P66744
D	-7	GLN	-	expression tag	UNP P66744
D	-6	MET	-	expression tag	UNP P66744
D	-5	GLY	-	expression tag	UNP P66744
D	-4	ARG	-	expression tag	UNP P66744
D	-3	GLY	-	expression tag	UNP P66744
D	-2	SER	-	expression tag	UNP P66744
D	-1	GLU	-	expression tag	UNP P66744

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Chain	Residue	Modelled	Actual	Comment	Reference
D	0	PHE	-	expression tag	UNP P66744

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		

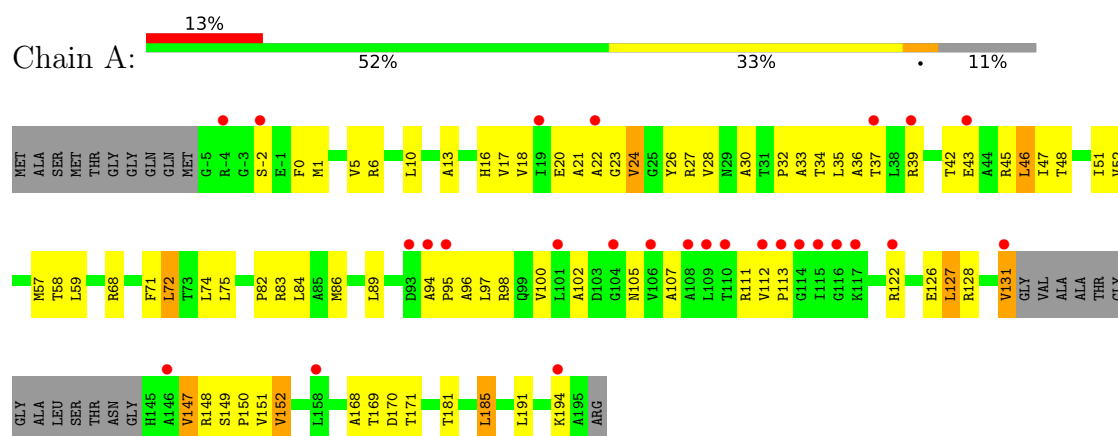
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	18	Total	O	0	0
			18	18		
3	B	5	Total	O	0	0
			5	5		
3	C	14	Total	O	0	0
			14	14		
3	D	11	Total	O	0	0
			11	11		

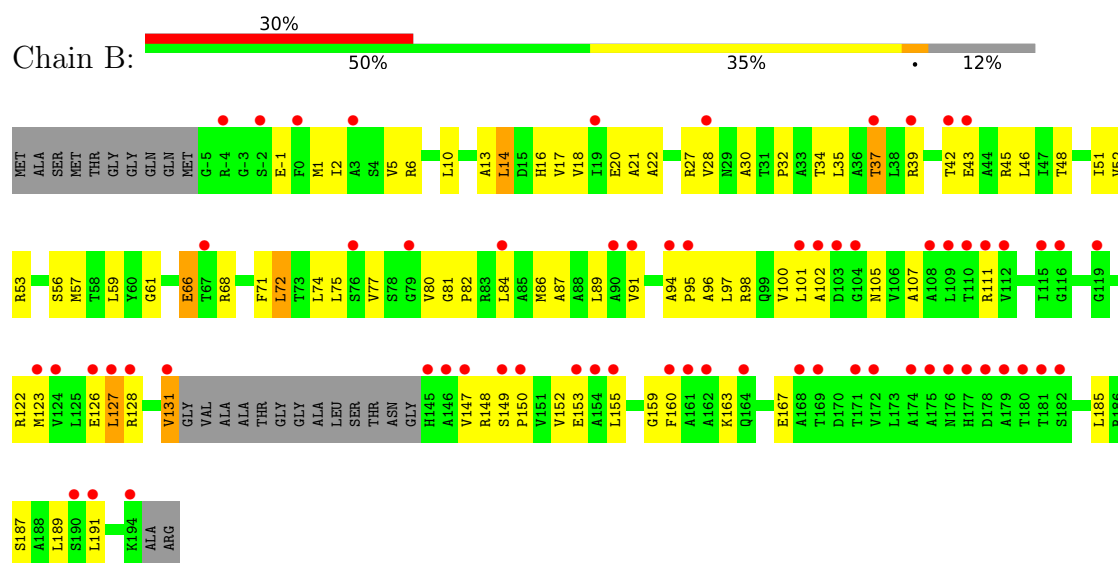
3 Residue-property plots [i](#)

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: Holliday junction ATP-dependent DNA helicase ruvA

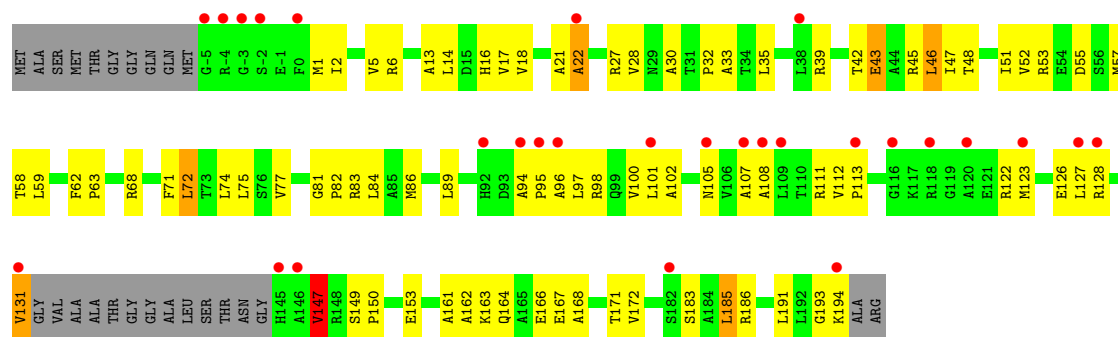


- Molecule 1: Holliday junction ATP-dependent DNA helicase ruvA

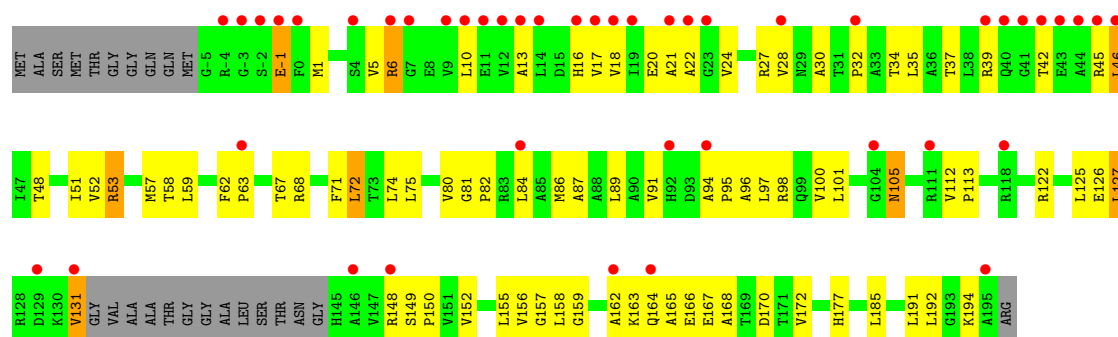


- Molecule 1: Holliday junction ATP-dependent DNA helicase ruvA





- Molecule 1: Holliday junction ATP-dependent DNA helicase ruvA



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	131.20Å 95.67Å 77.21Å 90.00° 121.80° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 30.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	91.8 (30.00-2.80) 97.1 (30.00-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.76Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.229 , 0.289 0.277 , 0.322	Depositor DCC
R_{free} test set	1018 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	57.9	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 79.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5372	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.17% 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: GOL

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	0.64	0/1341	1.02	4/1825 (0.2%)
1	B	0.56	0/1336	0.97	2/1818 (0.1%)
1	C	0.58	0/1336	1.01	3/1818 (0.2%)
1	D	0.58	0/1335	1.01	3/1818 (0.2%)
All	All	0.59	0/5348	1.00	12/7279 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	148	ARG	N-CA-C	8.29	120.40	111.36
1	C	172	VAL	N-CA-C	5.40	116.03	110.36
1	A	127	LEU	N-CA-C	5.34	119.02	112.03
1	A	-2	SER	N-CA-C	5.29	120.65	114.62
1	A	24	VAL	N-CA-C	5.25	115.82	108.89
1	C	147	VAL	N-CA-C	5.13	120.00	109.34
1	B	-1	GLU	N-CA-C	5.09	117.03	108.73
1	D	-1	GLU	N-CA-C	5.09	117.03	108.73
1	D	127	LEU	N-CA-C	5.08	118.69	112.03
1	B	127	LEU	N-CA-C	5.06	118.66	112.03
1	C	22	ALA	CB-CA-C	-5.05	110.36	117.23
1	A	147	VAL	N-CA-C	5.03	119.80	109.34

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1329	0	1349	72	0
1	B	1324	0	1344	95	0
1	C	1324	0	1344	84	0
1	D	1323	0	1338	85	0
2	A	6	0	8	2	0
2	B	6	0	8	1	0
2	C	6	0	8	2	0
2	D	6	0	8	2	0
3	A	18	0	0	3	0
3	B	5	0	0	0	0
3	C	14	0	0	3	0
3	D	11	0	0	1	0
All	All	5372	0	5407	296	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 28.

All (296) 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:6:ARG:HG3	1:D:45:ARG:HD3	1.15	1.15
1:D:6:ARG:HG2	1:D:45:ARG:HB2	1.39	1.03
1:B:98:ARG:HG2	1:B:131:VAL:HG11	1.44	1.00
1:C:98:ARG:HG2	1:C:131:VAL:HG11	1.47	0.96
1:A:98:ARG:HG2	1:A:131:VAL:HG11	1.48	0.96
1:B:6:ARG:HG3	1:B:45:ARG:HB2	1.48	0.94
1:C:6:ARG:HG3	1:C:45:ARG:HB2	1.49	0.94
1:B:52:VAL:HG12	1:B:57:MET:HG3	1.48	0.94
1:B:185:LEU:HD21	1:D:32:PRO:HG3	1.49	0.93
1:A:1:MET:HE1	1:B:57:MET:HG2	1.51	0.92
1:A:194:LYS:HD2	1:B:111:ARG:HE	1.34	0.92
1:D:6:ARG:CG	1:D:45:ARG:HB2	2.00	0.91
1:C:14:LEU:HD12	1:D:185:LEU:HD22	1.52	0.91
1:D:52:VAL:HG12	1:D:57:MET:HG3	1.53	0.90
1:D:98:ARG:HG2	1:D:131:VAL:HG11	1.51	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:6:ARG:HG3	1:D:45:ARG:CD	2.01	0.89
1:A:1:MET:CE	1:B:57:MET:HG2	2.03	0.88
1:A:6:ARG:HG3	1:A:45:ARG:HB2	1.56	0.85
1:C:52:VAL:HG12	1:C:57:MET:HG3	1.59	0.85
1:B:185:LEU:HD23	1:D:32:PRO:HB3	1.59	0.84
1:A:52:VAL:HG22	1:A:57:MET:HG3	1.60	0.83
1:D:80:VAL:HG13	1:D:84:LEU:HD22	1.60	0.82
1:A:6:ARG:HG3	1:A:45:ARG:HD3	1.62	0.81
1:D:6:ARG:HB3	1:D:6:ARG:CZ	2.10	0.80
1:D:101:LEU:HD13	1:D:127:LEU:CD1	2.10	0.80
1:A:0:PHE:HA	3:A:503:HOH:O	1.81	0.79
1:C:6:ARG:HG3	1:C:45:ARG:HD3	1.64	0.79
1:D:81:GLY:HA3	2:D:501:GOL:H2	1.63	0.79
1:A:194:LYS:HD2	1:B:111:ARG:NE	1.97	0.79
1:C:101:LEU:HD13	1:C:127:LEU:CD1	2.13	0.78
1:B:53:ARG:HB2	1:B:56:SER:OG	1.83	0.78
1:B:6:ARG:CG	1:B:45:ARG:HB2	2.14	0.78
1:B:81:GLY:HA3	2:B:500:GOL:H2	1.63	0.78
1:D:101:LEU:HD13	1:D:127:LEU:HD11	1.65	0.78
1:C:101:LEU:HD13	1:C:127:LEU:HD11	1.67	0.77
1:C:6:ARG:CG	1:C:45:ARG:HB2	2.14	0.76
1:C:83:ARG:HB2	2:C:503:GOL:O2	1.85	0.76
1:D:13:ALA:O	1:D:35:LEU:HD13	1.84	0.76
1:B:101:LEU:HD13	1:B:127:LEU:CD1	2.17	0.75
1:C:111:ARG:HD3	1:D:194:LYS:HB2	1.68	0.75
1:A:83:ARG:HB2	2:A:502:GOL:H2	1.69	0.74
1:B:75:LEU:HD22	1:B:82:PRO:HD3	1.69	0.74
1:B:101:LEU:HD13	1:B:127:LEU:HD11	1.70	0.74
1:D:6:ARG:CG	1:D:45:ARG:HD3	2.08	0.72
1:B:6:ARG:HG3	1:B:45:ARG:HD3	1.69	0.72
1:A:83:ARG:HB2	2:A:502:GOL:C2	2.20	0.71
1:D:162:ALA:O	1:D:166:GLU:HB2	1.89	0.71
1:A:28:VAL:HG12	1:A:59:LEU:HB3	1.74	0.70
1:D:28:VAL:HG12	1:D:59:LEU:HB3	1.73	0.70
1:B:80:VAL:HA	1:B:84:LEU:HD23	1.74	0.69
1:C:13:ALA:O	1:C:35:LEU:HD13	1.92	0.69
1:A:6:ARG:CG	1:A:45:ARG:HB2	2.23	0.68
1:B:28:VAL:HG12	1:B:59:LEU:HB3	1.75	0.67
1:C:57:MET:HG2	1:D:1:MET:HE1	1.77	0.66
1:B:28:VAL:HG12	1:B:59:LEU:CB	2.27	0.65
1:B:34:THR:O	1:B:37:THR:HG23	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:VAL:HG13	1:C:2:ILE:HG23	1.79	0.65
1:B:1:MET:HE1	1:D:57:MET:HG2	1.79	0.64
1:C:14:LEU:CD1	1:D:185:LEU:HD22	2.26	0.64
1:A:168:ALA:O	1:A:171:THR:HB	1.98	0.63
1:C:113:PRO:HB3	1:D:159:GLY:O	1.99	0.63
1:B:185:LEU:CD2	1:D:32:PRO:HB3	2.27	0.63
1:B:1:MET:CE	1:D:57:MET:HG2	2.29	0.62
1:D:68:ARG:HD2	1:D:72:LEU:CD2	2.29	0.62
1:B:68:ARG:HD2	1:B:72:LEU:CD2	2.29	0.61
1:D:163:LYS:O	1:D:167:GLU:HG3	2.01	0.61
1:C:74:LEU:CD1	1:C:89:LEU:HD11	2.31	0.61
1:A:75:LEU:HD22	1:A:82:PRO:HD3	1.82	0.61
1:B:163:LYS:O	1:B:167:GLU:HG3	2.01	0.61
1:A:94:ALA:HB3	1:A:95:PRO:HD3	1.82	0.61
1:B:160:PHE:CD1	1:D:113:PRO:HG3	2.35	0.61
1:C:28:VAL:HG12	1:C:59:LEU:HB3	1.82	0.61
1:C:74:LEU:HD11	1:C:89:LEU:HD11	1.82	0.61
1:C:51:ILE:HG23	1:C:53:ARG:NH1	2.16	0.60
1:C:57:MET:HG2	1:D:1:MET:CE	2.30	0.60
1:C:94:ALA:HB3	1:C:95:PRO:HD3	1.84	0.60
1:A:34:THR:O	1:A:37:THR:HG22	2.01	0.60
1:C:33:ALA:HB2	3:C:515:HOH:O	2.01	0.60
1:A:194:LYS:HB2	1:B:111:ARG:NE	2.17	0.60
1:B:46:LEU:HD23	1:B:61:GLY:HA3	1.83	0.60
1:C:68:ARG:HD2	1:C:72:LEU:CD2	2.32	0.60
1:D:75:LEU:HD22	1:D:82:PRO:HD3	1.84	0.59
1:B:94:ALA:HB3	1:B:95:PRO:HD3	1.84	0.59
1:D:28:VAL:HG12	1:D:59:LEU:CB	2.32	0.59
1:A:28:VAL:HG12	1:A:59:LEU:CB	2.32	0.59
1:B:32:PRO:HA	1:B:35:LEU:HD12	1.84	0.59
1:B:147:VAL:C	1:B:150:PRO:HD2	2.28	0.58
1:C:28:VAL:HG12	1:C:59:LEU:CB	2.33	0.58
1:A:74:LEU:CD1	1:A:89:LEU:HD11	2.33	0.58
1:B:80:VAL:HG22	1:B:84:LEU:CD2	2.34	0.57
1:D:81:GLY:HA3	2:D:501:GOL:C2	2.35	0.57
1:A:74:LEU:HD11	1:A:89:LEU:HD11	1.86	0.57
1:D:34:THR:O	1:D:37:THR:HG22	2.05	0.57
1:A:1:MET:HE3	1:B:57:MET:HG2	1.85	0.56
1:A:68:ARG:HD2	1:A:72:LEU:CD2	2.34	0.56
1:D:94:ALA:HB3	1:D:95:PRO:HD3	1.87	0.56
1:D:96:ALA:O	1:D:100:VAL:HG23	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:6:ARG:HG3	1:C:45:ARG:CB	2.30	0.56
1:D:17:VAL:HG22	1:D:18:VAL:N	2.20	0.56
1:B:149:SER:O	1:B:153:GLU:HB2	2.05	0.56
1:A:149:SER:HB3	1:A:150:PRO:HD3	1.88	0.56
1:B:96:ALA:O	1:B:100:VAL:HG23	2.06	0.56
1:D:192:LEU:C	1:D:194:LYS:H	2.12	0.55
1:B:89:LEU:HD23	1:B:97:LEU:HD23	1.89	0.55
1:D:30:ALA:HA	1:D:86:MET:HE1	1.89	0.55
1:B:6:ARG:HG3	1:B:45:ARG:CB	2.30	0.55
1:D:155:LEU:HD12	1:D:165:ALA:HB1	1.90	0.54
1:B:75:LEU:CD2	1:B:82:PRO:HD3	2.38	0.54
1:A:147:VAL:O	1:A:151:VAL:HG23	2.09	0.53
1:A:57:MET:HG2	1:C:1:MET:CE	2.39	0.53
1:C:96:ALA:O	1:C:100:VAL:HG23	2.08	0.53
1:B:66:GLU:CD	1:B:66:GLU:H	2.16	0.53
1:C:71:PHE:HA	1:C:74:LEU:HD12	1.90	0.52
1:D:71:PHE:HA	1:D:74:LEU:HD12	1.91	0.52
1:C:107:ALA:O	1:C:111:ARG:HG3	2.09	0.52
1:D:51:ILE:HD13	1:D:82:PRO:HG3	1.90	0.52
1:B:5:VAL:HG23	1:B:48:THR:HG21	1.91	0.52
1:D:125:LEU:HB2	3:D:507:HOH:O	2.08	0.52
1:D:168:ALA:O	1:D:172:VAL:HG23	2.09	0.52
1:C:147:VAL:C	1:C:150:PRO:HD2	2.35	0.52
1:B:71:PHE:HA	1:B:74:LEU:HD12	1.90	0.52
1:C:113:PRO:CB	1:D:159:GLY:O	2.58	0.51
1:B:17:VAL:HG22	1:B:18:VAL:N	2.25	0.51
1:A:13:ALA:O	1:A:35:LEU:HD13	2.10	0.51
1:A:6:ARG:HG3	1:A:45:ARG:CB	2.37	0.51
1:C:168:ALA:O	1:C:171:THR:HB	2.11	0.51
1:C:55:ASP:OD1	1:D:53:ARG:HA	2.10	0.51
1:C:51:ILE:HG23	1:C:53:ARG:CZ	2.41	0.51
1:B:30:ALA:HA	1:B:86:MET:HE1	1.92	0.51
1:B:46:LEU:O	1:B:48:THR:HG23	2.11	0.51
1:B:98:ARG:HD3	1:B:131:VAL:HG12	1.93	0.51
1:B:107:ALA:O	1:B:111:ARG:HG3	2.10	0.51
1:C:149:SER:HB3	1:C:150:PRO:HD3	1.93	0.50
1:A:71:PHE:HA	1:A:74:LEU:HD12	1.93	0.50
1:D:89:LEU:HD23	1:D:97:LEU:HD23	1.92	0.50
1:D:149:SER:N	1:D:150:PRO:CD	2.74	0.50
1:D:6:ARG:HG3	1:D:45:ARG:HB2	1.90	0.50
1:B:80:VAL:HG22	1:B:84:LEU:HD23	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:ILE:HD13	1:C:82:PRO:HG3	1.93	0.50
1:A:96:ALA:O	1:A:100:VAL:HG23	2.12	0.50
1:B:149:SER:N	1:B:150:PRO:CD	2.75	0.50
1:B:98:ARG:HG2	1:B:131:VAL:CG1	2.30	0.50
1:C:81:GLY:HA3	2:C:503:GOL:H2	1.94	0.50
1:C:163:LYS:O	1:C:167:GLU:HG3	2.11	0.50
1:C:183:SER:HA	1:C:186:ARG:CD	2.41	0.50
1:B:149:SER:HA	1:B:152:VAL:CG2	2.42	0.49
1:C:98:ARG:CG	1:C:131:VAL:HG11	2.32	0.49
1:C:183:SER:HA	1:C:186:ARG:HD2	1.94	0.49
1:A:16:HIS:CE1	1:A:27:ARG:HH21	2.31	0.49
1:C:32:PRO:HG3	1:D:185:LEU:HD21	1.93	0.49
1:B:51:ILE:HD13	1:B:82:PRO:HG3	1.95	0.49
1:A:194:LYS:HG2	3:A:506:HOH:O	2.13	0.49
1:D:127:LEU:HD12	1:D:127:LEU:O	2.12	0.49
1:A:57:MET:HG2	1:C:1:MET:HE1	1.94	0.49
1:A:194:LYS:HB2	1:B:111:ARG:CZ	2.43	0.49
1:C:5:VAL:HG23	1:C:48:THR:HG21	1.94	0.49
1:A:17:VAL:HG22	1:A:18:VAL:N	2.28	0.48
1:B:160:PHE:CE1	1:D:113:PRO:HG3	2.48	0.48
1:B:189:LEU:C	1:B:191:LEU:N	2.70	0.48
1:C:30:ALA:HA	1:C:86:MET:HE1	1.94	0.48
1:D:46:LEU:O	1:D:48:THR:HG23	2.13	0.48
1:A:32:PRO:HG3	1:C:185:LEU:HD13	1.96	0.48
1:A:107:ALA:O	1:A:111:ARG:HG3	2.14	0.48
1:C:6:ARG:HG3	1:C:45:ARG:CD	2.40	0.47
1:C:21:ALA:O	1:C:22:ALA:HB3	2.14	0.47
1:A:6:ARG:HG3	1:A:45:ARG:CD	2.41	0.47
1:C:17:VAL:HG22	1:C:18:VAL:N	2.29	0.47
1:A:98:ARG:HD3	1:A:131:VAL:HG12	1.97	0.47
1:A:181:THR:HG22	3:A:513:HOH:O	2.14	0.47
1:B:80:VAL:HA	1:B:84:LEU:CD2	2.44	0.47
1:A:30:ALA:HA	1:A:86:MET:HE1	1.97	0.47
1:C:161:ALA:O	1:C:162:ALA:C	2.57	0.47
1:A:194:LYS:HB2	1:B:111:ARG:CD	2.45	0.46
1:B:13:ALA:O	1:B:35:LEU:HD13	2.15	0.46
1:B:185:LEU:CD2	1:D:32:PRO:HG3	2.34	0.46
1:B:185:LEU:HD21	1:D:32:PRO:CG	2.33	0.46
1:B:127:LEU:HD12	1:B:127:LEU:O	2.15	0.46
1:B:187:SER:O	1:B:191:LEU:HB2	2.15	0.46
1:C:55:ASP:C	1:D:52:VAL:HG23	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:GLY:HA3	1:C:21:ALA:HB1	1.97	0.46
1:C:185:LEU:HD23	1:C:185:LEU:HA	1.51	0.46
1:B:80:VAL:CA	1:B:84:LEU:HD23	2.46	0.45
1:C:46:LEU:O	1:C:48:THR:HG23	2.16	0.45
1:D:167:GLU:O	1:D:170:ASP:HB2	2.16	0.45
1:C:58:THR:HG22	1:C:59:LEU:N	2.31	0.45
1:A:1:MET:HB2	1:B:57:MET:O	2.16	0.45
1:B:66:GLU:CD	1:B:66:GLU:N	2.75	0.45
1:B:98:ARG:HD3	1:B:131:VAL:CG1	2.47	0.45
1:C:98:ARG:HD3	1:C:131:VAL:HG12	1.98	0.45
1:A:17:VAL:HG12	1:A:28:VAL:HG23	1.99	0.45
1:A:39:ARG:O	1:A:42:THR:HB	2.17	0.45
1:C:33:ALA:CB	3:C:515:HOH:O	2.63	0.45
1:D:127:LEU:HD12	1:D:127:LEU:C	2.41	0.45
1:C:112:VAL:HA	1:C:113:PRO:HD2	1.82	0.45
1:A:105:ASN:C	1:A:105:ASN:OD1	2.60	0.45
1:B:53:ARG:HB2	1:B:56:SER:HG	1.78	0.45
1:B:80:VAL:HG13	1:B:84:LEU:HD23	1.99	0.45
1:C:42:THR:HG22	1:C:43:GLU:O	2.16	0.45
1:A:17:VAL:CG1	1:A:28:VAL:HG23	2.47	0.45
1:B:155:LEU:O	1:B:160:PHE:HD2	2.00	0.45
3:C:517:HOH:O	1:D:158:LEU:HD21	2.17	0.45
1:D:16:HIS:CE1	1:D:27:ARG:HH21	2.35	0.45
1:B:10:LEU:HD11	1:B:20:GLU:HB2	1.99	0.44
1:B:39:ARG:O	1:B:42:THR:HB	2.17	0.44
1:B:149:SER:HA	1:B:152:VAL:HG22	1.99	0.44
1:C:75:LEU:HD21	1:C:82:PRO:HA	1.99	0.44
1:C:127:LEU:O	1:C:127:LEU:HD12	2.16	0.44
1:C:186:ARG:H	1:C:186:ARG:HG3	1.55	0.44
1:B:16:HIS:CE1	1:B:27:ARG:HH21	2.35	0.44
1:B:97:LEU:HD12	1:B:97:LEU:O	2.18	0.44
1:C:149:SER:O	1:C:153:GLU:HB2	2.16	0.44
1:D:5:VAL:HG23	1:D:48:THR:HG21	1.99	0.44
1:B:87:ALA:O	1:B:91:VAL:HG23	2.17	0.44
1:C:183:SER:HA	1:C:186:ARG:NE	2.32	0.44
1:A:32:PRO:HA	1:A:35:LEU:HD12	2.00	0.44
1:B:98:ARG:CG	1:B:131:VAL:HG11	2.30	0.44
1:B:148:ARG:O	1:B:152:VAL:HG22	2.17	0.44
1:C:62:PHE:HB3	1:C:63:PRO:HD2	1.99	0.44
1:A:46:LEU:O	1:A:48:THR:HG23	2.17	0.44
1:D:105:ASN:OD1	1:D:105:ASN:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:152:VAL:O	1:D:156:VAL:HG23	2.17	0.44
1:A:169:THR:O	1:A:170:ASP:C	2.60	0.44
1:D:98:ARG:HD3	1:D:131:VAL:HG12	2.00	0.44
1:C:97:LEU:HD12	1:C:97:LEU:O	2.18	0.43
1:A:6:ARG:NH1	1:B:20:GLU:OE2	2.50	0.43
1:A:97:LEU:HD12	1:A:97:LEU:O	2.18	0.43
1:A:21:ALA:O	1:A:22:ALA:HB3	2.18	0.43
1:A:51:ILE:HD13	1:A:82:PRO:HG3	1.98	0.43
1:C:101:LEU:HD22	1:C:127:LEU:HD11	2.00	0.43
1:D:10:LEU:HD11	1:D:20:GLU:HB2	2.00	0.43
1:B:2:ILE:HG23	1:D:24:VAL:HG13	2.00	0.43
1:B:127:LEU:HD12	1:B:127:LEU:C	2.44	0.43
1:C:16:HIS:CE1	1:C:27:ARG:HH21	2.36	0.43
1:A:102:ALA:HA	1:A:128:ARG:HG3	2.00	0.43
1:D:6:ARG:HG2	1:D:45:ARG:CB	2.29	0.43
1:C:127:LEU:HD12	1:C:127:LEU:C	2.44	0.43
1:A:16:HIS:ND1	1:A:27:ARG:NH2	2.67	0.43
1:A:149:SER:O	1:A:150:PRO:C	2.61	0.43
1:A:149:SER:HB3	1:A:150:PRO:CD	2.49	0.43
1:B:77:VAL:HG11	1:B:123:MET:HG2	2.00	0.43
1:D:87:ALA:O	1:D:91:VAL:HG23	2.19	0.42
1:A:32:PRO:CG	1:C:185:LEU:HD13	2.49	0.42
1:D:58:THR:HG22	1:D:59:LEU:N	2.35	0.42
1:D:67:THR:HG23	1:D:94:ALA:HB2	2.01	0.42
1:A:5:VAL:HG23	1:A:48:THR:HG21	2.00	0.42
1:A:112:VAL:HA	1:A:113:PRO:HD2	1.83	0.42
1:C:17:VAL:HG12	1:C:28:VAL:HG23	2.00	0.42
1:A:185:LEU:HD12	1:B:14:LEU:HD13	2.01	0.42
1:C:39:ARG:O	1:C:42:THR:HB	2.19	0.42
1:B:16:HIS:ND1	1:B:27:ARG:NH2	2.68	0.42
1:B:80:VAL:HG22	1:B:84:LEU:HD21	2.00	0.42
1:C:17:VAL:CG1	1:C:28:VAL:HG23	2.50	0.42
1:D:112:VAL:HA	1:D:113:PRO:HD2	1.88	0.42
1:A:148:ARG:O	1:A:152:VAL:HG13	2.20	0.42
1:C:102:ALA:HA	1:C:128:ARG:HG3	2.01	0.42
1:A:33:ALA:O	1:A:36:ALA:HB3	2.20	0.42
1:A:185:LEU:HA	1:A:185:LEU:HD23	1.81	0.42
1:D:101:LEU:HD13	1:D:127:LEU:HD13	1.99	0.42
1:B:148:ARG:C	1:B:150:PRO:HD2	2.44	0.42
1:D:97:LEU:O	1:D:97:LEU:HD12	2.19	0.42
1:B:21:ALA:O	1:B:22:ALA:HB3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:ALA:O	1:C:111:ARG:HB2	2.20	0.41
1:C:166:GLU:O	1:C:167:GLU:C	2.61	0.41
1:B:68:ARG:O	1:B:71:PHE:HB3	2.20	0.41
1:C:16:HIS:ND1	1:C:27:ARG:NH2	2.68	0.41
1:D:157:GLY:C	1:D:159:GLY:H	2.28	0.41
1:B:102:ALA:HA	1:B:128:ARG:HG3	2.01	0.41
1:B:160:PHE:CE1	1:D:113:PRO:CG	3.03	0.41
1:D:98:ARG:CG	1:D:131:VAL:HG11	2.36	0.41
1:C:77:VAL:HG11	1:C:123:MET:HG2	2.02	0.41
1:D:39:ARG:NH1	1:D:42:THR:OG1	2.51	0.41
1:A:10:LEU:HD11	1:A:20:GLU:HB2	2.01	0.41
1:A:58:THR:HG22	1:A:59:LEU:N	2.36	0.41
1:B:6:ARG:HG3	1:B:45:ARG:CD	2.44	0.41
1:C:68:ARG:O	1:C:71:PHE:HB3	2.20	0.41
1:B:149:SER:N	1:B:150:PRO:HD2	2.36	0.41
1:C:98:ARG:HG2	1:C:131:VAL:CG1	2.34	0.41
1:A:98:ARG:HD3	1:A:131:VAL:CG1	2.51	0.41
1:B:17:VAL:CG1	1:B:28:VAL:HG23	2.50	0.41
1:B:42:THR:HG22	1:B:43:GLU:O	2.21	0.41
1:B:159:GLY:O	1:D:113:PRO:HB3	2.21	0.41
1:C:105:ASN:OD1	1:C:105:ASN:C	2.64	0.41
1:D:17:VAL:CG1	1:D:28:VAL:HG23	2.51	0.41
1:D:17:VAL:CG2	1:D:18:VAL:N	2.84	0.41
1:D:21:ALA:O	1:D:22:ALA:HB3	2.21	0.41
1:D:68:ARG:HD2	1:D:72:LEU:HD21	2.01	0.41
1:A:24:VAL:HB	1:A:26:TYR:CE1	2.56	0.41
1:B:68:ARG:HD2	1:B:72:LEU:HD21	2.01	0.41
1:A:47:ILE:N	1:A:47:ILE:CD1	2.84	0.40
1:C:47:ILE:N	1:C:47:ILE:CD1	2.84	0.40
1:D:163:LYS:O	1:D:164:GLN:C	2.64	0.40
1:C:193:GLY:O	1:C:194:LYS:HB2	2.22	0.40
1:D:62:PHE:HB3	1:D:63:PRO:HD2	2.03	0.40
1:C:27:ARG:NH1	1:D:-1:GLU:CB	2.85	0.40
1:B:105:ASN:OD1	1:B:105:ASN:C	2.64	0.40
1:D:101:LEU:CD1	1:D:127:LEU:HD11	2.44	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	184/212 (87%)	163 (89%)	21 (11%)	0	100	100
1	B	183/212 (86%)	163 (89%)	20 (11%)	0	100	100
1	C	183/212 (86%)	159 (87%)	24 (13%)	0	100	100
1	D	184/212 (87%)	161 (88%)	22 (12%)	1 (0%)	24	55
All	All	734/848 (87%)	646 (88%)	87 (12%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	105	ASN

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	129/154 (84%)	118 (92%)	11 (8%)	10	31
1	B	129/154 (84%)	122 (95%)	7 (5%)	20	51
1	C	129/154 (84%)	118 (92%)	11 (8%)	10	31
1	D	128/154 (83%)	119 (93%)	9 (7%)	14	40
All	All	515/616 (84%)	477 (93%)	38 (7%)	13	37

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	GLU
1	A	46	LEU
1	A	72	LEU
1	A	84	LEU
1	A	122	ARG
1	A	126	GLU
1	A	127	LEU
1	A	131	VAL
1	A	152	VAL
1	A	185	LEU
1	A	191	LEU
1	B	14	LEU
1	B	37	THR
1	B	66	GLU
1	B	72	LEU
1	B	122	ARG
1	B	126	GLU
1	B	131	VAL
1	C	43	GLU
1	C	46	LEU
1	C	72	LEU
1	C	84	LEU
1	C	122	ARG
1	C	126	GLU
1	C	131	VAL
1	C	147	VAL
1	C	164	GLN
1	C	185	LEU
1	C	191	LEU
1	D	6	ARG
1	D	46	LEU
1	D	53	ARG
1	D	72	LEU
1	D	122	ARG
1	D	126	GLU
1	D	131	VAL
1	D	177	HIS
1	D	191	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

4 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	GOL	C	503	-	5,5,5	0.40	0	5,5,5	0.50	0
2	GOL	B	500	-	5,5,5	0.28	0	5,5,5	0.62	0
2	GOL	A	502	-	5,5,5	0.27	0	5,5,5	0.74	0
2	GOL	D	501	-	5,5,5	0.34	0	5,5,5	0.61	0

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	GOL	C	503	-	-	2/4/4/4	-
2	GOL	B	500	-	-	2/4/4/4	-
2	GOL	A	502	-	-	4/4/4/4	-
2	GOL	D	501	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	502	GOL	O1-C1-C2-C3
2	A	502	GOL	C1-C2-C3-O3
2	B	500	GOL	O1-C1-C2-O2
2	B	500	GOL	O1-C1-C2-C3
2	C	503	GOL	O1-C1-C2-C3
2	D	501	GOL	C1-C2-C3-O3
2	D	501	GOL	O2-C2-C3-O3
2	A	502	GOL	O1-C1-C2-O2
2	A	502	GOL	O2-C2-C3-O3
2	C	503	GOL	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	503	GOL	2	0
2	B	500	GOL	1	0
2	A	502	GOL	2	0
2	D	501	GOL	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	188/212 (88%)	1.07	27 (14%) 6 5	50, 76, 123, 184	0
1	B	187/212 (88%)	1.68	64 (34%) 1 1	62, 96, 166, 227	0
1	C	187/212 (88%)	1.09	28 (14%) 5 4	48, 77, 129, 184	0
1	D	188/212 (88%)	1.43	45 (23%) 2 1	56, 86, 131, 230	0
All	All	750/848 (88%)	1.32	164 (21%) 2 2	48, 83, 142, 230	0

All (164) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	39	ARG	4.6
1	A	110	THR	4.5
1	D	9	VAL	4.4
1	D	43	GLU	4.1
1	A	146	ALA	3.9
1	D	22	ALA	3.9
1	B	150	PRO	3.9
1	B	102	ALA	3.9
1	D	12	VAL	3.8
1	C	146	ALA	3.6
1	D	195	ALA	3.6
1	B	116	GLY	3.6
1	A	37	THR	3.6
1	A	109	LEU	3.6
1	D	44	ALA	3.5
1	D	94	ALA	3.5
1	D	7	GLY	3.5
1	D	-2	SER	3.5
1	B	178	ASP	3.4
1	A	22	ALA	3.4
1	D	131	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	127	LEU	3.3
1	B	155	LEU	3.2
1	B	194	LYS	3.2
1	B	149	SER	3.2
1	D	21	ALA	3.2
1	A	113	PRO	3.2
1	C	194	LYS	3.1
1	B	191	LEU	3.1
1	C	-5	GLY	3.1
1	D	-3	GLY	3.1
1	D	11	GLU	3.1
1	D	84	LEU	3.1
1	D	28	VAL	3.1
1	B	128	ARG	3.1
1	B	181	THR	3.1
1	B	123	MET	3.0
1	D	41	GLY	3.0
1	D	148	ARG	3.0
1	B	168	ALA	3.0
1	B	177	HIS	3.0
1	D	14	LEU	3.0
1	D	42	THR	3.0
1	B	0	PHE	3.0
1	B	169	THR	3.0
1	C	94	ALA	3.0
1	A	194	LYS	2.9
1	B	131	VAL	2.9
1	B	172	VAL	2.9
1	C	0	PHE	2.9
1	A	112	VAL	2.9
1	B	111	ARG	2.9
1	C	-2	SER	2.9
1	B	37	THR	2.9
1	C	95	PRO	2.8
1	C	128	ARG	2.8
1	D	10	LEU	2.8
1	B	147	VAL	2.8
1	A	95	PRO	2.8
1	B	43	GLU	2.8
1	B	153	GLU	2.8
1	C	127	LEU	2.8
1	D	4	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	116	GLY	2.8
1	D	164	GLN	2.7
1	B	3	ALA	2.7
1	B	180	THR	2.7
1	D	17	VAL	2.7
1	A	101	LEU	2.7
1	B	176	ASN	2.7
1	B	119	GLY	2.7
1	D	111	ARG	2.7
1	B	19	ILE	2.7
1	B	115	ILE	2.7
1	C	107	ALA	2.7
1	D	-1	GLU	2.7
1	B	171	THR	2.6
1	A	-2	SER	2.6
1	B	179	ALA	2.6
1	D	13	ALA	2.6
1	D	23	GLY	2.6
1	B	101	LEU	2.6
1	B	109	LEU	2.6
1	A	115	ILE	2.6
1	D	129	ASP	2.6
1	C	145	HIS	2.6
1	A	116	GLY	2.6
1	B	42	THR	2.6
1	D	40	GLN	2.6
1	B	164	GLN	2.5
1	B	146	ALA	2.5
1	D	0	PHE	2.5
1	B	145	HIS	2.5
1	B	160	PHE	2.5
1	C	120	ALA	2.5
1	C	-4	ARG	2.5
1	D	92	HIS	2.5
1	C	182	SER	2.5
1	C	109	LEU	2.5
1	A	114	GLY	2.5
1	B	67	THR	2.5
1	C	118	ARG	2.4
1	C	123	MET	2.4
1	B	104	GLY	2.4
1	B	108	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	92	HIS	2.4
1	B	-2	SER	2.4
1	D	162	ALA	2.4
1	B	84	LEU	2.4
1	A	108	ALA	2.4
1	D	146	ALA	2.4
1	C	101	LEU	2.4
1	B	103	ASP	2.3
1	B	76	SER	2.3
1	B	182	SER	2.3
1	A	19	ILE	2.3
1	B	39	ARG	2.3
1	C	113	PRO	2.3
1	D	63	PRO	2.3
1	C	131	VAL	2.3
1	A	94	ALA	2.3
1	B	94	ALA	2.3
1	B	174	ALA	2.3
1	C	38	LEU	2.3
1	B	190	SER	2.3
1	A	131	VAL	2.3
1	B	95	PRO	2.3
1	B	126	GLU	2.3
1	B	162	ALA	2.3
1	B	161	ALA	2.2
1	D	46	LEU	2.2
1	A	104	GLY	2.2
1	B	154	ALA	2.2
1	B	175	ALA	2.2
1	C	22	ALA	2.2
1	B	-4	ARG	2.2
1	A	39	ARG	2.2
1	D	19	ILE	2.2
1	D	16	HIS	2.2
1	B	110	THR	2.2
1	D	104	GLY	2.1
1	A	158	LEU	2.1
1	A	122	ARG	2.1
1	B	91	VAL	2.1
1	B	79	GLY	2.1
1	B	90	ALA	2.1
1	C	108	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	-4	ARG	2.1
1	D	6	ARG	2.1
1	D	45	ARG	2.1
1	D	118	ARG	2.1
1	A	106	VAL	2.1
1	B	112	VAL	2.1
1	C	-3	GLY	2.1
1	A	43	GLU	2.1
1	A	93	ASP	2.1
1	C	96	ALA	2.1
1	B	124	VAL	2.0
1	D	32	PRO	2.0
1	B	28	VAL	2.0
1	D	18	VAL	2.0
1	A	-4	ARG	2.0
1	C	105	ASN	2.0
1	A	117	LYS	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($\text{\AA}^2$)	Q<0.9
2	GOL	A	502	6/6	0.80	0.14	46,63,86,89	0
2	GOL	D	501	6/6	0.80	0.16	69,89,103,107	0
2	GOL	C	503	6/6	0.81	0.14	43,80,99,110	0
2	GOL	B	500	6/6	0.86	0.11	41,59,86,120	0

6.5 Other polymers [i](#)

There are no such residues in this entry.