



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2024 – 01:11 AM EDT

PDB ID : 2IS3
Title : Crystal Structure of Escherichia coli RNase T
Authors : Zuo, Y.; Wang, Y.; Malhotra, A.
Deposited on : 2006-10-16
Resolution : 3.10 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

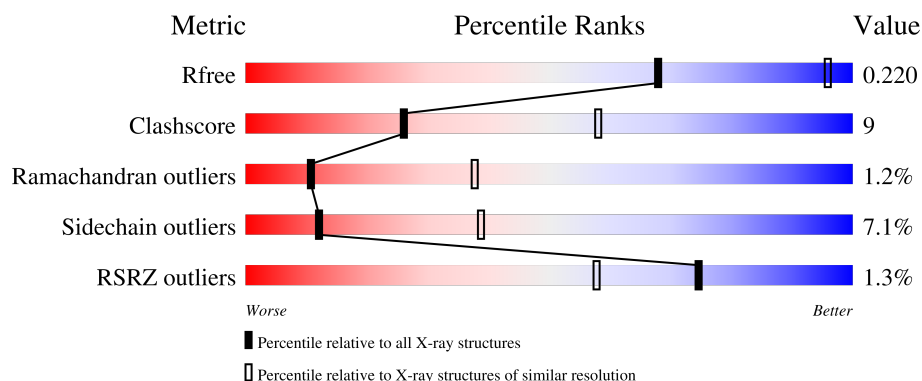
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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}	164625	1351 (3.10-3.10)
Clashscore	180529	1454 (3.10-3.10)
Ramachandran outliers	177936	1391 (3.10-3.10)
Sidechain outliers	177891	1391 (3.10-3.10)
RSRZ outliers	164620	1351 (3.10-3.10)

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	215	<div> <div>72%</div> <div>20%</div> <div>6%</div> </div>
1	B	215	<div> <div>60%</div> <div>18%</div> <div>20%</div> </div>
1	C	215	<div> <div>73%</div> <div>20%</div> <div>6%</div> </div>
1	D	215	<div> <div>72%</div> <div>21%</div> <div>6%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6104 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 Ribonuclease T.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	Se	0	0	0
			1568	998	272	288	4	6			
1	B	172	Total	C	N	O	S	Se	0	0	0
			1335	851	232	242	4	6			
1	C	203	Total	C	N	O	S	Se	0	0	0
			1568	998	272	288	4	6			
1	D	203	Total	C	N	O	S	Se	0	0	0
			1568	998	272	288	4	6			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP P30014
A	46	MSE	MET	modified residue	UNP P30014
A	53	MSE	MET	modified residue	UNP P30014
A	117	MSE	MET	modified residue	UNP P30014
A	129	MSE	MET	modified residue	UNP P30014
A	130	MSE	MET	modified residue	UNP P30014
A	173	MSE	MET	modified residue	UNP P30014
B	1	MSE	MET	modified residue	UNP P30014
B	46	MSE	MET	modified residue	UNP P30014
B	53	MSE	MET	modified residue	UNP P30014
B	117	MSE	MET	modified residue	UNP P30014
B	129	MSE	MET	modified residue	UNP P30014
B	130	MSE	MET	modified residue	UNP P30014
B	173	MSE	MET	modified residue	UNP P30014
C	1	MSE	MET	modified residue	UNP P30014
C	46	MSE	MET	modified residue	UNP P30014
C	53	MSE	MET	modified residue	UNP P30014
C	117	MSE	MET	modified residue	UNP P30014
C	129	MSE	MET	modified residue	UNP P30014
C	130	MSE	MET	modified residue	UNP P30014
C	173	MSE	MET	modified residue	UNP P30014

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Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MSE	MET	modified residue	UNP P30014
D	46	MSE	MET	modified residue	UNP P30014
D	53	MSE	MET	modified residue	UNP P30014
D	117	MSE	MET	modified residue	UNP P30014
D	129	MSE	MET	modified residue	UNP P30014
D	130	MSE	MET	modified residue	UNP P30014
D	173	MSE	MET	modified residue	UNP P30014

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



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

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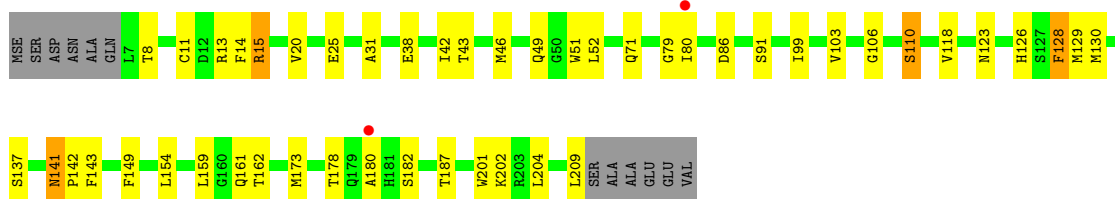
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

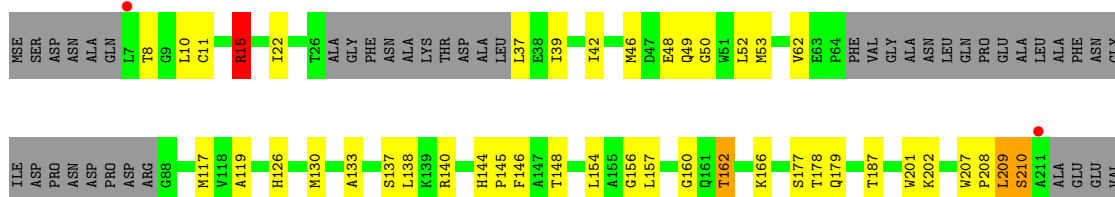
3 Residue-property plots [i](#)

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.

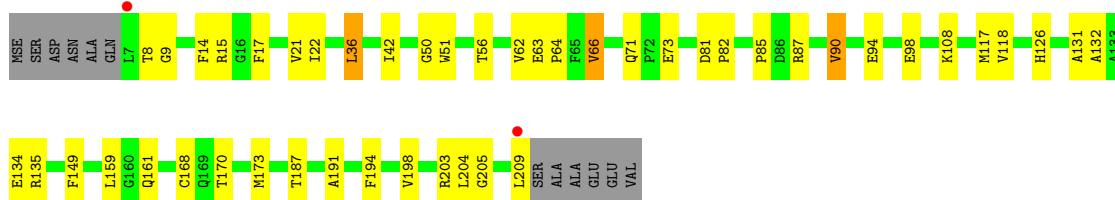
• Molecule 1: Ribonuclease T



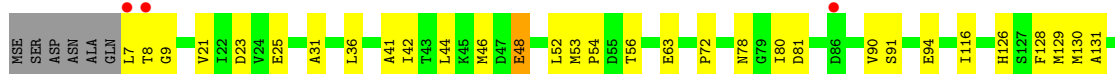
• Molecule 1: Ribonuclease T

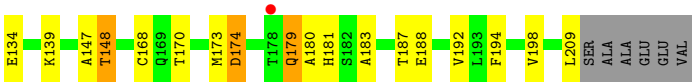


• Molecule 1: Ribonuclease T



• Molecule 1: Ribonuclease T





4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	213.14Å 213.14Å 149.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.83 – 3.10 29.83 – 3.10	Depositor EDS
% Data completeness (in resolution range)	88.3 (29.83-3.10) 88.2 (29.83-3.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 3.11Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.196 , 0.231 0.187 , 0.220	Depositor DCC
R_{free} test set	1418 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	85.2	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 74.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6104	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.45% 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 [i](#)

5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.66	1/1600 (0.1%)	0.63	0/2162
1	B	0.70	1/1359 (0.1%)	0.74	1/1829 (0.1%)
1	C	0.60	0/1600	0.63	0/2162
1	D	0.69	0/1600	0.69	0/2162
All	All	0.66	2/6159 (0.0%)	0.67	1/8315 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	11	CYS	CB-SG	-5.73	1.72	1.81
1	A	11	CYS	CB-SG	-5.57	1.72	1.81

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	15	ARG	NE-CZ-NH1	6.93	123.76	120.30

There are no chirality outliers.

There are no planarity outliers.

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	1568	0	1527	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1335	0	1308	34	0
1	C	1568	0	1527	30	0
1	D	1568	0	1527	26	0
2	A	15	0	0	1	0
2	B	20	0	0	1	0
2	C	15	0	0	0	0
2	D	15	0	0	1	0
All	All	6104	0	5889	112	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 9.

All (112) 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:8:THR:HG22	1:D:9:GLY:N	1.81	0.95
1:D:8:THR:HG22	1:D:9:GLY:H	1.41	0.85
1:D:8:THR:CG2	1:D:9:GLY:H	1.93	0.81
1:C:159:LEU:HD22	1:C:170:THR:HG22	1.63	0.80
1:D:8:THR:CG2	1:D:9:GLY:N	2.46	0.79
1:C:118:VAL:HG22	1:C:149:PHE:HB2	1.64	0.78
1:A:141:ASN:HD22	1:A:141:ASN:C	1.85	0.77
1:B:37:LEU:HG	1:B:62:VAL:HG23	1.64	0.77
1:B:15:ARG:CG	1:B:15:ARG:HH11	1.99	0.74
1:B:15:ARG:HH11	1:B:15:ARG:HG2	1.54	0.70
1:A:141:ASN:ND2	1:A:143:PHE:H	1.90	0.70
1:A:51:TRP:CZ3	1:A:202:LYS:HD3	2.33	0.63
1:A:15:ARG:HD2	1:B:162:THR:HB	1.79	0.63
1:A:141:ASN:HD22	1:A:142:PRO:N	1.96	0.62
1:A:118:VAL:HG22	1:A:149:PHE:HB2	1.81	0.62
1:B:42:ILE:HD12	1:B:187:THR:HG22	1.81	0.62
1:D:48:GLU:H	1:D:48:GLU:CD	2.03	0.62
1:C:71:GLN:HB3	1:C:73:GLU:OE1	2.02	0.60
1:A:201:TRP:CZ3	1:B:201:TRP:CE3	2.90	0.60
1:C:8:THR:HG22	1:C:9:GLY:N	2.16	0.60
1:A:79:GLY:O	1:A:80:ILE:HG13	2.02	0.59
1:B:117:MSE:HE3	1:B:119:ALA:HB2	1.84	0.59
1:B:8:THR:HG21	1:B:48:GLU:O	2.02	0.59
1:C:131:ALA:O	1:C:134:GLU:HB3	2.03	0.58
1:B:117:MSE:HE3	1:B:119:ALA:CB	2.33	0.58
1:B:133:ALA:HA	1:B:138:LEU:HD12	1.84	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:TRP:HA	1:B:208:PRO:O	2.03	0.57
1:A:141:ASN:C	1:A:141:ASN:ND2	2.55	0.57
1:B:207:TRP:HA	1:B:208:PRO:C	2.25	0.57
1:D:25:GLU:HG2	1:D:183:ALA:HB2	1.87	0.57
1:C:90:VAL:HG13	1:C:94:GLU:HB3	1.89	0.54
1:A:14:PHE:C	1:A:15:ARG:HG3	2.28	0.53
1:A:46:MSE:HE3	1:A:52:LEU:HD13	1.91	0.53
1:D:78:ASN:ND2	1:D:80:ILE:HD12	2.23	0.53
1:D:173:MSE:HB2	2:D:218:SO4:O4	2.09	0.52
1:D:81:ASP:OD1	1:D:81:ASP:C	2.48	0.52
1:C:22:ILE:HD12	1:C:117:MSE:HG3	1.91	0.52
1:B:8:THR:O	1:B:8:THR:HG22	2.10	0.52
1:A:126:HIS:O	1:A:130:MSE:HG2	2.11	0.51
1:C:8:THR:HG23	1:C:50:GLY:CA	2.41	0.50
1:A:49:GLN:HB2	1:A:51:TRP:CD1	2.46	0.50
1:D:179:GLN:O	1:D:181:HIS:N	2.43	0.50
1:B:15:ARG:HG2	1:B:15:ARG:NH1	2.26	0.50
1:D:42:ILE:HD12	1:D:187:THR:HG22	1.93	0.50
1:D:188:GLU:O	1:D:192:VAL:HG23	2.10	0.50
1:C:117:MSE:CE	1:C:126:HIS:HB2	2.42	0.49
1:A:20:VAL:HG22	1:A:43:THR:HG22	1.95	0.48
1:A:42:ILE:HD12	1:A:187:THR:HG22	1.95	0.48
1:C:51:TRP:CH2	1:C:203:ARG:HG3	2.49	0.48
1:B:37:LEU:HG	1:B:62:VAL:CG2	2.39	0.48
1:A:129:MSE:HG2	1:A:129:MSE:O	2.14	0.47
1:B:126:HIS:O	1:B:130:MSE:HG2	2.15	0.47
1:C:51:TRP:HH2	1:C:203:ARG:HG3	1.78	0.47
1:B:15:ARG:CG	1:B:15:ARG:NH1	2.68	0.47
1:C:42:ILE:HA	1:C:56:THR:O	2.15	0.46
1:A:99:ILE:O	1:A:103:VAL:HG23	2.15	0.46
1:B:22:ILE:HD12	1:B:117:MSE:HG3	1.97	0.46
1:A:154:LEU:HD23	1:B:154:LEU:HD23	1.98	0.46
1:B:209:LEU:O	1:B:210:SER:HB3	2.15	0.46
1:D:23:ASP:HB2	1:D:187:THR:OG1	2.15	0.46
1:D:194:PHE:O	1:D:198:VAL:HG23	2.16	0.46
1:C:21:VAL:HG21	1:C:191:ALA:HA	1.98	0.46
1:B:8:THR:HG23	1:B:50:GLY:CA	2.46	0.45
1:C:132:ALA:HA	1:C:135:ARG:NH1	2.31	0.45
1:A:31:ALA:HA	1:A:128:PHE:CE2	2.51	0.45
1:C:42:ILE:HD12	1:C:187:THR:HG22	1.98	0.45
1:C:117:MSE:HE1	1:C:126:HIS:HB2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:168:CYS:O	1:D:173:MSE:HG2	2.17	0.45
1:C:203:ARG:C	1:C:205:GLY:H	2.20	0.44
1:D:46:MSE:HE3	1:D:52:LEU:HD13	2.00	0.44
1:A:141:ASN:HD22	1:A:143:PHE:H	1.62	0.44
1:D:131:ALA:O	1:D:134:GLU:HB3	2.18	0.44
1:A:173:MSE:HB2	2:A:217:SO4:O2	2.18	0.44
1:D:126:HIS:O	1:D:130:MSE:HG2	2.18	0.44
1:C:21:VAL:HA	1:C:118:VAL:O	2.18	0.43
1:D:31:ALA:HA	1:D:128:PHE:CE2	2.53	0.43
1:C:63:GLU:HA	1:C:64:PRO:HD2	1.75	0.43
1:D:63:GLU:HG2	1:D:91:SER:HA	2.01	0.43
1:A:46:MSE:HA	1:A:51:TRP:O	2.19	0.43
1:D:179:GLN:H	1:D:179:GLN:HG2	1.57	0.43
1:C:36:LEU:CD2	1:C:62:VAL:HG11	2.49	0.42
1:C:161:GLN:HA	1:C:161:GLN:OE1	2.19	0.42
1:C:8:THR:HG22	1:C:9:GLY:H	1.84	0.42
1:B:49:GLN:O	1:B:202:LYS:NZ	2.50	0.42
1:C:8:THR:CG2	1:C:9:GLY:N	2.82	0.42
1:D:21:VAL:O	1:D:41:ALA:HA	2.19	0.42
1:B:144:HIS:HA	1:B:145:PRO:HD3	1.83	0.42
1:A:201:TRP:CE3	1:B:201:TRP:CE3	3.08	0.41
1:B:53:MSE:HE2	1:B:53:MSE:HB3	2.01	0.41
1:D:147:ALA:C	1:D:148:THR:HG22	2.41	0.41
1:D:44:LEU:HD23	1:D:54:PRO:HA	2.02	0.41
1:A:106:GLY:O	1:A:110:SER:HB3	2.20	0.41
1:B:46:MSE:HE3	1:B:52:LEU:HD13	2.01	0.41
1:C:168:CYS:O	1:C:173:MSE:HG2	2.20	0.41
1:A:79:GLY:C	1:A:80:ILE:HG13	2.40	0.41
1:C:118:VAL:HG22	1:C:149:PHE:CB	2.43	0.41
1:D:78:ASN:HD21	1:D:80:ILE:HD12	1.84	0.41
1:B:15:ARG:HH11	1:B:15:ARG:HG3	1.81	0.41
1:B:177:SER:O	1:B:178:THR:C	2.56	0.41
1:C:194:PHE:O	1:C:198:VAL:HG23	2.21	0.41
1:C:36:LEU:HD23	1:C:62:VAL:HG11	2.03	0.40
1:C:81:ASP:HA	1:C:82:PRO:HD3	1.90	0.40
1:C:85:PRO:C	1:C:87:ARG:H	2.25	0.40
1:A:13:ARG:NH2	1:B:157:LEU:O	2.47	0.40
1:A:25:GLU:HB2	1:A:38:GLU:HB3	2.03	0.40
1:A:161:GLN:HE22	1:B:15:ARG:NE	2.20	0.40
1:D:174:ASP:N	1:D:174:ASP:OD1	2.54	0.40
1:B:10:LEU:HA	1:B:10:LEU:HD23	1.83	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:PHE:O	1:C:17:PHE:HB2	2.22	0.40
1:A:201:TRP:CE3	1:B:201:TRP:CZ3	3.10	0.40
1:B:140:ARG:NH2	2:B:217:SO4:O1	2.54	0.40
1:B:156:GLY:O	1:B:160:GLY:HA2	2.21	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	201/215 (94%)	184 (92%)	16 (8%)	1 (0%)	25	58
1	B	166/215 (77%)	152 (92%)	12 (7%)	2 (1%)	11	38
1	C	201/215 (94%)	178 (89%)	20 (10%)	3 (2%)	8	33
1	D	201/215 (94%)	184 (92%)	14 (7%)	3 (2%)	8	33
All	All	769/860 (89%)	698 (91%)	62 (8%)	9 (1%)	11	38

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	180	ALA
1	B	210	SER
1	C	15	ARG
1	C	66	VAL
1	D	180	ALA
1	D	48	GLU
1	B	146	PHE
1	C	204	LEU
1	D	72	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	161/163 (99%)	145 (90%)	16 (10%)	6	24
1	B	138/163 (85%)	130 (94%)	8 (6%)	17	45
1	C	161/163 (99%)	155 (96%)	6 (4%)	29	59
1	D	161/163 (99%)	147 (91%)	14 (9%)	8	31
All	All	621/652 (95%)	577 (93%)	44 (7%)	12	39

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	THR
1	A	15	ARG
1	A	71	GLN
1	A	86	ASP
1	A	91	SER
1	A	110	SER
1	A	123	ASN
1	A	128	PHE
1	A	137	SER
1	A	141	ASN
1	A	159	LEU
1	A	162	THR
1	A	178	THR
1	A	182	SER
1	A	204	LEU
1	A	209	LEU
1	B	15	ARG
1	B	39	ILE
1	B	137	SER
1	B	148	THR
1	B	162	THR
1	B	166	LYS
1	B	179	GLN
1	B	209	LEU

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Mol	Chain	Res	Type
1	C	36	LEU
1	C	66	VAL
1	C	90	VAL
1	C	98	GLU
1	C	108	LYS
1	C	209	LEU
1	D	7	LEU
1	D	36	LEU
1	D	53	MSE
1	D	56	THR
1	D	90	VAL
1	D	94	GLU
1	D	116	ILE
1	D	129	MSE
1	D	139	LYS
1	D	148	THR
1	D	170	THR
1	D	174	ASP
1	D	179	GLN
1	D	209	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	59	HIS
1	A	141	ASN
1	A	161	GLN
1	B	97	HIS
1	C	123	ASN
1	D	78	ASN
1	D	123	ASN
1	D	179	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

13 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	218	-	4,4,4	0.20	0	6,6,6	0.37	0
2	SO4	D	218	-	4,4,4	0.20	0	6,6,6	0.29	0
2	SO4	A	217	-	4,4,4	0.24	0	6,6,6	0.22	0
2	SO4	A	218	-	4,4,4	0.29	0	6,6,6	0.30	0
2	SO4	C	217	-	4,4,4	0.25	0	6,6,6	0.13	0
2	SO4	C	218	-	4,4,4	0.29	0	6,6,6	0.33	0
2	SO4	D	216	-	4,4,4	0.21	0	6,6,6	0.40	0
2	SO4	B	217	-	4,4,4	0.21	0	6,6,6	0.26	0
2	SO4	D	217	-	4,4,4	0.31	0	6,6,6	0.32	0
2	SO4	B	216	-	4,4,4	0.31	0	6,6,6	0.37	0
2	SO4	B	219	-	4,4,4	0.30	0	6,6,6	0.25	0
2	SO4	C	216	-	4,4,4	0.34	0	6,6,6	0.35	0
2	SO4	A	216	-	4,4,4	0.29	0	6,6,6	0.33	0

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.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	218	SO4	1	0
2	A	217	SO4	1	0
2	B	217	SO4	1	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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	197/215 (91%)	-0.18	2 (1%) 79 64	64, 76, 89, 100	0
1	B	166/215 (77%)	-0.47	2 (1%) 76 60	64, 74, 85, 90	0
1	C	197/215 (91%)	-0.42	2 (1%) 79 64	65, 76, 89, 92	0
1	D	197/215 (91%)	0.02	4 (2%) 64 45	65, 75, 94, 104	0
All	All	757/860 (88%)	-0.25	10 (1%) 74 58	64, 75, 89, 104	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	180	ALA	3.5
1	C	7	LEU	3.4
1	D	7	LEU	3.4
1	B	211	ALA	3.2
1	D	8	THR	3.2
1	C	209	LEU	3.0
1	D	86	ASP	2.8
1	A	80	ILE	2.5
1	D	178	THR	2.2
1	B	7	LEU	2.2

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 monosaccharides 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	SO4	C	218	5/5	0.77	0.24	138,139,139,139	0
2	SO4	A	218	5/5	0.80	0.31	133,134,134,134	0
2	SO4	B	217	5/5	0.81	0.23	135,135,135,136	0
2	SO4	C	217	5/5	0.84	0.16	168,168,168,168	0
2	SO4	A	217	5/5	0.85	0.27	143,143,143,143	0
2	SO4	A	216	5/5	0.88	0.16	105,105,106,107	0
2	SO4	D	216	5/5	0.90	0.21	106,107,107,108	0
2	SO4	D	217	5/5	0.90	0.28	104,104,105,105	0
2	SO4	D	218	5/5	0.91	0.18	115,115,115,116	0
2	SO4	B	219	5/5	0.93	0.17	110,111,111,111	0
2	SO4	B	216	5/5	0.95	0.19	94,95,95,95	0
2	SO4	B	218	5/5	0.95	0.14	121,122,122,122	0
2	SO4	C	216	5/5	0.97	0.14	62,63,64,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.