



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 12:15 pm GMT

PDB ID : 5C1I
Title : m1A58 tRNA methyltransferase mutant - D170A
Authors : Ponchon, L.; Degut, C.; Folly-Klan, M.; Barraud, P.; Tisne, C.
Deposited on : 2015-06-14
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

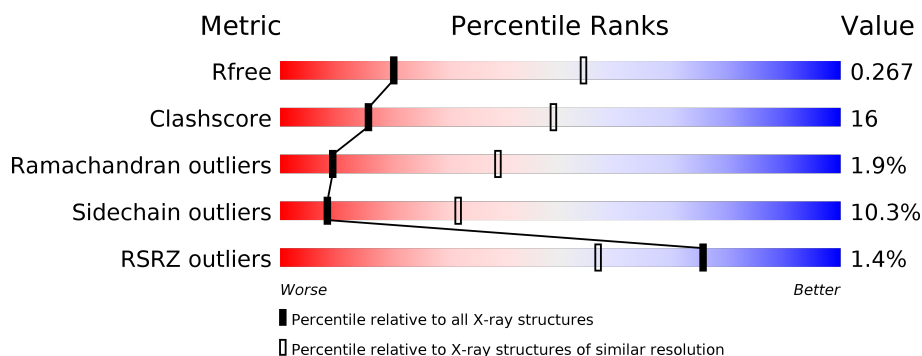
1 Overall quality at a glance

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}	100719	1001 (3.12-3.08)
Clashscore	112137	1099 (3.12-3.08)
Ramachandran outliers	110173	1057 (3.12-3.08)
Sidechain outliers	110143	1057 (3.12-3.08)
RSRZ outliers	101464	1006 (3.12-3.08)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	251	<div> <div>2%</div> <div>61% 24% • 12%</div> </div>
1	B	251	<div> <div>70% 26% • •</div> </div>
1	C	251	<div> <div>2%</div> <div>46% 30% 8% 16%</div> </div>
1	D	251	<div> <div>51% 24% • 22%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6987 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 tRNA (adenine(58)-N(1))-methyltransferase TrmI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	0	0
			1733	1121	310	298	4			
1	B	251	Total	C	N	O	S	0	0	0
			1986	1276	362	344	4			
1	C	210	Total	C	N	O	S	0	0	0
			1667	1076	298	289	4			
1	D	197	Total	C	N	O	S	0	0	0
			1564	1006	280	274	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	170	ALA	ASP	engineered mutation	UNP Q8GBB2
A	254	GLY	ALA	conflict	UNP Q8GBB2
B	170	ALA	ASP	engineered mutation	UNP Q8GBB2
B	254	GLY	ALA	conflict	UNP Q8GBB2
C	170	ALA	ASP	engineered mutation	UNP Q8GBB2
C	254	GLY	ALA	conflict	UNP Q8GBB2
D	170	ALA	ASP	engineered mutation	UNP Q8GBB2
D	254	GLY	ALA	conflict	UNP Q8GBB2

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



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

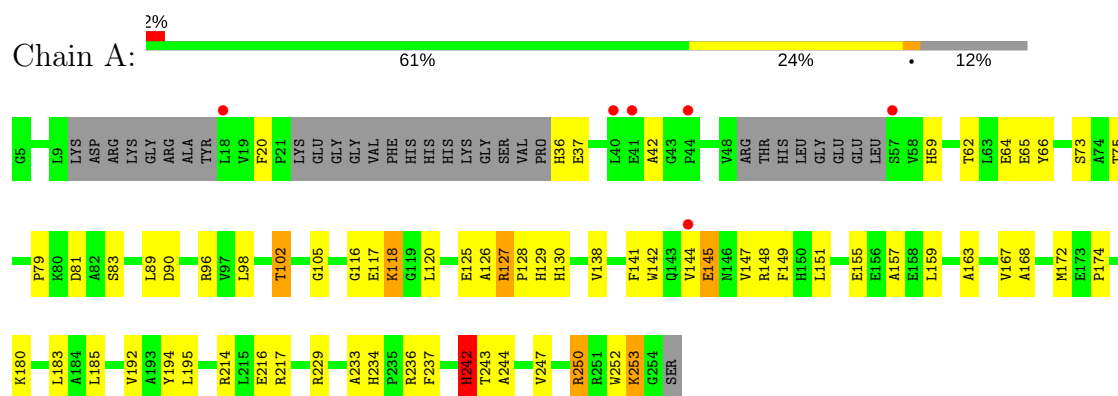
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total	O	0	0
			7	7		
3	B	7	Total	O	0	0
			7	7		
3	C	2	Total	O	0	0
			2	2		
3	D	6	Total	O	0	0
			6	6		

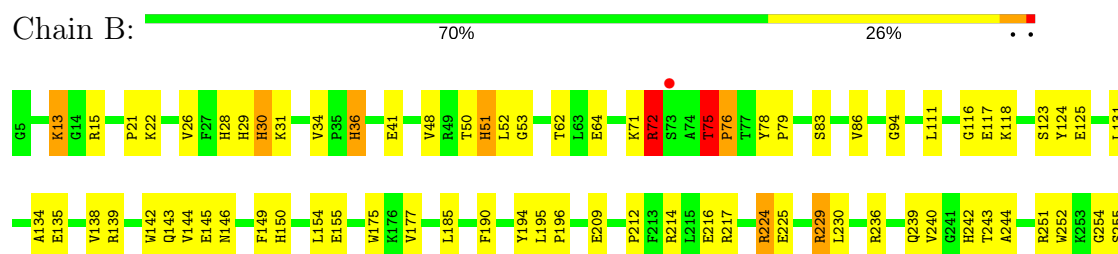
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

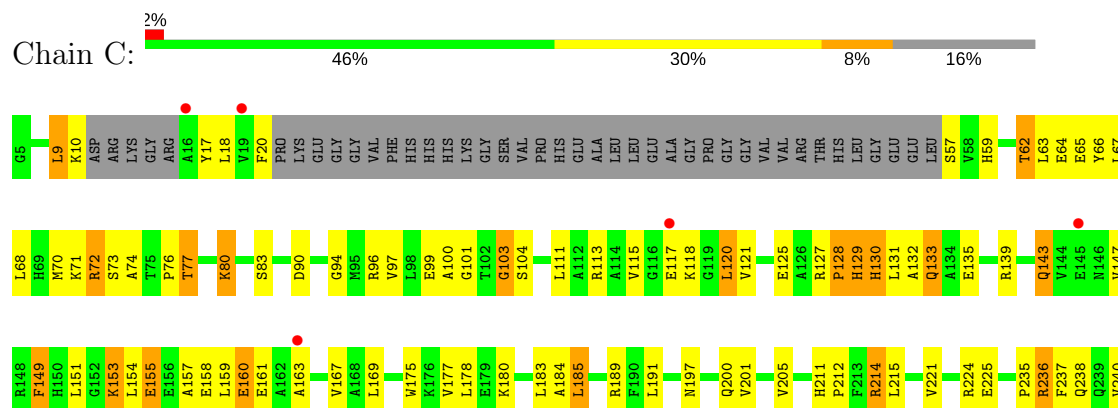
- Molecule 1: tRNA (adenine(58)-N(1))-methyltransferase TrmI

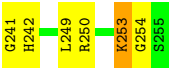


- Molecule 1: tRNA (adenine(58)-N(1))-methyltransferase TrmI

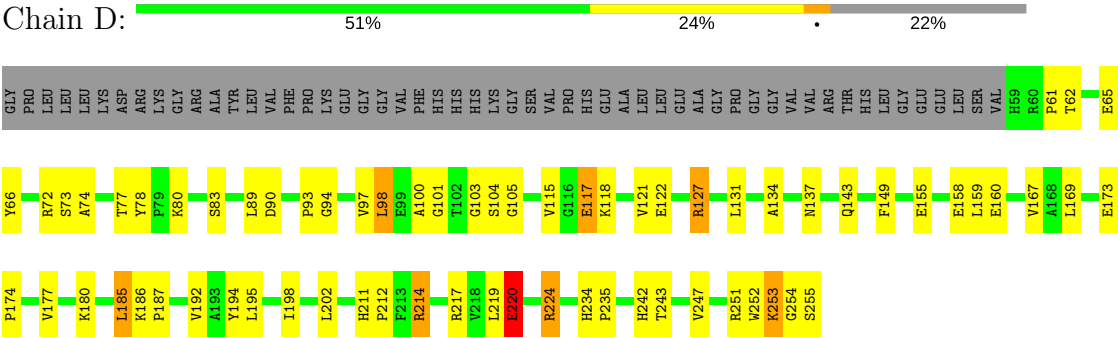


- Molecule 1: tRNA (adenine(58)-N(1))-methyltransferase TrmI





● Molecule 1: tRNA (adenine(58)-N(1))-methyltransferase TrmI



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	110.43Å 110.43Å 306.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	51.94 – 3.10 81.13 – 3.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (51.94-3.10) 100.0 (81.13-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 3.13Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.200 , 0.269 0.195 , 0.267	Depositor DCC
R_{free} test set	1077 reflections (5.14%)	DCC
Wilson B-factor (Å ²)	84.0	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 62.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6987	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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: 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	0.80	0/1775	0.87	1/2410 (0.0%)
1	B	0.82	1/2038 (0.0%)	0.93	4/2763 (0.1%)
1	C	0.79	0/1706	0.89	0/2312
1	D	0.81	1/1603 (0.1%)	0.87	1/2176 (0.0%)
All	All	0.81	2/7122 (0.0%)	0.89	6/9661 (0.1%)

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	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	72	ARG	CG-CD	6.27	1.67	1.51
1	D	220	GLU	CG-CD	5.14	1.59	1.51

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	72	ARG	NE-CZ-NH2	7.76	124.18	120.30
1	B	75	THR	N-CA-CB	5.90	121.50	110.30
1	A	250	ARG	NE-CZ-NH2	-5.57	117.52	120.30
1	B	72	ARG	CA-CB-CG	5.36	125.20	113.40
1	D	185	LEU	CA-CB-CG	5.24	127.34	115.30
1	B	224	ARG	NE-CZ-NH1	5.11	122.86	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	118	LYS	Peptide

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	1733	0	1756	53	0
1	B	1986	0	2006	73	0
1	C	1667	0	1693	80	0
1	D	1564	0	1573	42	0
2	B	10	0	0	1	0
2	D	5	0	0	0	0
3	A	7	0	0	2	0
3	B	7	0	0	5	0
3	C	2	0	0	1	0
3	D	6	0	0	4	0
All	All	6987	0	7028	229	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 16.

All (229) 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:B:72:ARG:HH21	1:B:76:PRO:HB2	1.34	0.92
1:A:96:ARG:NH1	1:A:163:ALA:O	2.02	0.92
1:A:157:ALA:O	1:A:180:LYS:NZ	2.02	0.92
1:D:253:LYS:O	1:D:255:SER:N	2.10	0.84
1:C:157:ALA:O	1:C:180:LYS:NZ	2.12	0.83
1:B:72:ARG:HH21	1:B:76:PRO:CB	1.95	0.79
1:C:80:LYS:NZ	1:C:221:VAL:O	2.15	0.79
1:C:154:LEU:O	1:C:157:ALA:N	2.19	0.75
1:B:72:ARG:HD2	1:B:76:PRO:HB2	1.69	0.75
1:D:74:ALA:HB3	1:D:103:GLY:HA2	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:104:SER:HB3	1:D:134:ALA:HB2	1.68	0.73
1:C:153:LYS:HZ3	1:C:154:LEU:H	1.37	0.71
1:B:229:ARG:NH1	3:B:401:HOH:O	2.08	0.71
1:A:217:ARG:NH2	1:C:225:GLU:OE2	2.24	0.70
1:B:72:ARG:HD2	1:B:76:PRO:CB	2.21	0.70
1:A:81:ASP:OD2	1:A:243:THR:OG1	2.08	0.70
1:D:167:VAL:HG23	1:D:185:LEU:HD13	1.73	0.69
1:A:253:LYS:NZ	3:A:302:HOH:O	2.25	0.69
1:B:118:LYS:HD3	1:D:117:GLU:HG2	1.73	0.69
1:C:215:LEU:HA	1:C:249:LEU:HD23	1.74	0.69
1:B:118:LYS:HD3	1:D:117:GLU:CG	2.24	0.68
1:C:167:VAL:HG23	1:C:185:LEU:HD12	1.76	0.67
1:C:175:TRP:HA	1:C:178:LEU:CD1	2.26	0.66
1:C:96:ARG:HG3	1:C:120:LEU:HD23	1.79	0.65
1:B:64:GLU:OE2	1:D:217:ARG:NH2	2.30	0.65
1:C:96:ARG:CG	1:C:120:LEU:HD23	2.28	0.64
1:A:90:ASP:HB2	1:C:62:THR:HG22	1.79	0.64
1:C:154:LEU:O	1:C:180:LYS:NZ	2.27	0.63
2:B:301:SO4:O3	3:B:402:HOH:O	2.14	0.62
1:B:72:ARG:CZ	1:B:76:PRO:HG2	2.30	0.62
1:D:62:THR:OG1	1:D:65:GLU:HG3	2.00	0.61
1:B:72:ARG:NH2	1:B:76:PRO:HG2	2.16	0.61
1:D:155:GLU:HG3	1:D:177:VAL:HG12	1.82	0.61
1:B:34:VAL:HG22	1:B:50:THR:HA	1.84	0.59
1:A:127:ARG:HH11	1:A:128:PRO:HD2	1.67	0.59
1:B:72:ARG:NH1	1:B:78:TYR:CE1	2.70	0.59
1:C:236:ARG:NH1	1:C:237:PHE:CZ	2.71	0.59
1:B:72:ARG:HD2	1:B:76:PRO:CG	2.33	0.58
1:C:127:ARG:HB3	1:C:130:HIS:CE1	2.38	0.58
1:A:89:LEU:HD11	1:A:168:ALA:HB2	1.85	0.58
1:B:52:LEU:H	1:B:53:GLY:HA2	1.68	0.58
1:C:139:ARG:O	1:C:143:GLN:NE2	2.38	0.56
1:C:131:LEU:O	1:C:135:GLU:HG3	2.05	0.56
1:C:201:VAL:O	1:C:205:VAL:HG23	2.05	0.56
1:A:229:ARG:NH2	1:A:236:ARG:HH22	2.03	0.55
1:C:153:LYS:HZ2	1:C:153:LYS:HB3	1.72	0.55
1:D:117:GLU:OE1	3:D:401:HOH:O	2.18	0.55
1:B:225:GLU:N	1:B:239:GLN:OE1	2.39	0.55
1:B:72:ARG:CD	1:B:76:PRO:HG2	2.36	0.55
1:C:236:ARG:NH1	1:C:237:PHE:CE2	2.75	0.55
1:B:21:PRO:HA	1:B:36:HIS:CD2	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:VAL:HG23	1:B:28:HIS:CE1	2.41	0.54
1:B:145:GLU:H	1:B:145:GLU:CD	2.11	0.54
1:D:78:TYR:CD2	1:D:243:THR:HG21	2.42	0.54
1:C:96:ARG:HH12	1:C:163:ALA:C	2.11	0.53
1:C:175:TRP:HA	1:C:178:LEU:HD13	1.88	0.53
1:C:161:GLU:HB2	1:C:183:LEU:HD11	1.90	0.53
1:A:167:VAL:HG23	1:A:185:LEU:HD13	1.90	0.53
1:C:161:GLU:HB2	1:C:183:LEU:CD1	2.39	0.52
1:D:174:PRO:HG3	1:D:195:LEU:HD21	1.91	0.52
1:B:29:HIS:O	1:B:30:HIS:HB2	2.09	0.52
1:C:63:LEU:HD22	1:C:113:ARG:HD2	1.92	0.52
1:B:22:LYS:H	1:B:36:HIS:HD2	1.57	0.51
1:C:101:GLY:C	1:C:125:GLU:HG2	2.30	0.51
1:B:72:ARG:NH2	1:B:76:PRO:CG	2.73	0.51
1:C:155:GLU:HA	1:C:180:LYS:CE	2.40	0.51
1:C:211:HIS:HB3	1:C:212:PRO:HD2	1.92	0.51
1:B:118:LYS:CD	1:D:117:GLU:HG2	2.41	0.51
1:B:131:LEU:HD11	1:B:149:PHE:HB3	1.92	0.51
1:B:72:ARG:NH2	1:B:76:PRO:CB	2.71	0.51
1:A:127:ARG:HG3	1:A:128:PRO:HD2	1.93	0.51
1:B:26:VAL:HG23	1:B:28:HIS:NE2	2.25	0.51
1:C:77:THR:OG1	3:C:301:HOH:O	2.18	0.50
1:B:62:THR:HA	1:D:90:ASP:OD2	2.11	0.50
1:D:194:TYR:O	1:D:195:LEU:HD23	2.11	0.50
1:A:242:HIS:ND1	1:A:242:HIS:C	2.65	0.50
1:B:139:ARG:HG3	1:B:139:ARG:NH1	2.26	0.50
1:C:94:GLY:HA2	1:C:118:LYS:HG2	1.94	0.50
1:B:214:ARG:HB2	1:B:252:TRP:HE1	1.76	0.50
1:A:144:VAL:HG12	1:A:145:GLU:N	2.27	0.49
1:A:217:ARG:HH21	1:C:225:GLU:CD	2.15	0.49
1:B:124:TYR:CZ	1:B:150:HIS:HD2	2.30	0.49
1:C:132:ALA:O	1:C:135:GLU:N	2.45	0.49
1:D:131:LEU:HD11	1:D:149:PHE:HB3	1.94	0.49
1:B:131:LEU:O	1:B:135:GLU:HG3	2.13	0.49
1:A:62:THR:HG22	1:C:90:ASP:HB2	1.94	0.49
1:D:234:HIS:HD2	1:D:235:PRO:O	1.96	0.49
1:B:251:ARG:HH12	1:B:255:SER:CB	2.26	0.48
1:B:155:GLU:OE1	1:B:177:VAL:HG12	2.13	0.48
1:C:99:GLU:HG2	1:C:100:ALA:N	2.27	0.48
1:B:139:ARG:HG3	1:B:139:ARG:HH11	1.78	0.48
1:C:111:LEU:O	1:C:115:VAL:HG22	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:211:HIS:HB3	1:D:212:PRO:HD2	1.95	0.48
1:A:229:ARG:NH2	1:A:236:ARG:NH2	2.62	0.48
1:B:236:ARG:NH2	3:B:401:HOH:O	2.41	0.48
1:A:127:ARG:HG3	1:A:128:PRO:CD	2.44	0.48
1:C:62:THR:OG1	1:C:65:GLU:OE1	2.31	0.48
1:C:18:LEU:HD21	1:C:66:TYR:CD1	2.49	0.48
1:D:185:LEU:HD23	1:D:251:ARG:HB2	1.96	0.48
1:A:42:ALA:O	1:C:253:LYS:NZ	2.41	0.48
1:A:217:ARG:NH1	1:C:64:GLU:OE2	2.47	0.47
1:C:97:VAL:CG1	1:C:121:VAL:HG22	2.44	0.47
1:C:155:GLU:HA	1:C:180:LYS:HD3	1.96	0.47
1:C:167:VAL:CG2	1:C:185:LEU:HD12	2.42	0.47
1:C:17:TYR:CD1	1:C:17:TYR:N	2.79	0.47
1:A:66:TYR:CE2	1:A:141:PHE:HB2	2.49	0.47
1:B:83:SER:OG	1:D:83:SER:OG	2.32	0.47
1:C:149:PHE:N	1:C:149:PHE:HD1	2.12	0.47
1:A:96:ARG:HG3	1:A:120:LEU:HB3	1.97	0.47
1:B:34:VAL:HG11	1:B:48:VAL:HG11	1.96	0.47
1:B:52:LEU:N	1:B:53:GLY:HA2	2.29	0.47
1:D:224:ARG:NH1	1:D:224:ARG:HG2	2.29	0.47
1:D:212:PRO:HA	1:D:252:TRP:CE3	2.50	0.47
1:A:242:HIS:ND1	1:A:242:HIS:O	2.48	0.47
1:C:153:LYS:HZ1	1:C:177:VAL:HG12	1.80	0.47
1:D:186:LYS:HB3	1:D:187:PRO:HD2	1.97	0.47
1:A:126:ALA:O	1:A:151:LEU:HD11	2.14	0.47
1:B:117:GLU:HG3	3:D:401:HOH:O	2.15	0.46
1:B:229:ARG:NH2	3:B:401:HOH:O	2.47	0.46
1:D:104:SER:HA	1:D:105:GLY:HA2	1.79	0.46
1:B:125:GLU:HG3	1:B:131:LEU:HD12	1.96	0.46
1:C:10:LYS:HE3	1:C:59:HIS:NE2	2.31	0.46
1:C:149:PHE:N	1:C:149:PHE:CD1	2.83	0.46
1:D:98:LEU:HD12	1:D:122:GLU:HB2	1.97	0.46
1:A:118:LYS:HE3	1:A:145:GLU:O	2.15	0.46
1:A:64:GLU:HG3	1:A:79:PRO:HG2	1.96	0.46
1:B:72:ARG:CD	1:B:76:PRO:CG	2.94	0.46
1:A:127:ARG:NH1	1:A:128:PRO:HD2	2.29	0.46
1:B:224:ARG:HB3	1:D:220:GLU:HG3	1.96	0.46
1:C:127:ARG:HH11	1:C:128:PRO:HD2	1.80	0.46
1:C:67:LEU:HA	1:C:70:MET:CE	2.46	0.46
1:B:229:ARG:CZ	3:B:401:HOH:O	2.60	0.46
1:C:127:ARG:NH2	1:C:129:HIS:ND1	2.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:ARG:HH11	1:A:127:ARG:HG3	1.81	0.45
1:A:105:GLY:HA3	1:A:138:VAL:CG2	2.46	0.45
1:A:144:VAL:HG12	1:A:145:GLU:H	1.80	0.45
1:A:194:TYR:CE2	1:A:244:ALA:HB3	2.52	0.45
1:D:219:LEU:C	1:D:219:LEU:HD12	2.36	0.45
1:A:155:GLU:HA	1:A:180:LYS:HE2	1.98	0.45
1:C:155:GLU:HA	1:C:180:LYS:HE2	1.99	0.45
1:C:71:LYS:HG2	1:C:133:GLN:NE2	2.31	0.45
1:C:161:GLU:CB	1:C:183:LEU:HD11	2.46	0.45
1:C:169:LEU:HD11	1:C:191:LEU:HD11	1.98	0.45
1:A:120:LEU:HD11	1:A:148:ARG:HG3	1.99	0.45
1:B:29:HIS:O	1:B:30:HIS:CB	2.64	0.45
1:D:97:VAL:HG13	1:D:121:VAL:HG13	1.98	0.45
1:B:86:VAL:HG22	1:B:111:LEU:HD23	1.98	0.45
1:B:94:GLY:HA2	1:B:118:LYS:HG3	1.99	0.45
1:A:73:SER:HG	1:A:130:HIS:CD2	2.34	0.45
1:C:253:LYS:N	1:C:254:GLY:CA	2.81	0.44
1:A:147:VAL:CG1	1:A:149:PHE:CE1	3.01	0.44
1:A:59:HIS:HB3	1:C:250:ARG:HH12	1.83	0.44
1:B:72:ARG:HH21	1:B:76:PRO:CG	2.29	0.44
1:C:73:SER:OG	1:C:103:GLY:O	2.32	0.44
1:D:219:LEU:O	1:D:219:LEU:HD12	2.17	0.44
1:A:216:GLU:HG3	1:A:250:ARG:HB2	2.00	0.44
1:C:153:LYS:HZ3	1:C:154:LEU:N	2.08	0.44
1:C:175:TRP:O	1:C:178:LEU:HD13	2.18	0.44
1:C:197:ASN:ND2	1:C:200:GLN:OE1	2.41	0.44
1:A:73:SER:OG	1:A:130:HIS:NE2	2.50	0.44
1:D:118:LYS:HE2	3:D:401:HOH:O	2.18	0.44
1:D:66:TYR:OH	1:D:137:ASN:O	2.20	0.44
1:B:230:LEU:HD21	1:D:202:LEU:HD22	2.00	0.43
1:B:212:PRO:HA	1:B:252:TRP:CE3	2.53	0.43
1:A:233:ALA:O	1:D:198:ILE:HG22	2.18	0.43
1:A:98:LEU:HB3	1:A:167:VAL:HG13	1.99	0.43
1:B:75:THR:HA	1:B:76:PRO:HD3	1.63	0.43
1:C:236:ARG:CZ	1:C:237:PHE:CZ	3.01	0.43
1:A:234:HIS:CE1	1:A:237:PHE:CE2	3.06	0.43
1:A:252:TRP:HB2	3:A:306:HOH:O	2.18	0.43
1:B:224:ARG:HH22	1:C:242:HIS:HB2	1.84	0.43
1:C:73:SER:HA	1:C:74:ALA:HA	1.77	0.43
1:A:118:LYS:HE3	1:A:145:GLU:C	2.39	0.43
1:A:147:VAL:HG11	1:A:149:PHE:CE1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:155:GLU:N	1:C:180:LYS:HE2	2.33	0.43
1:A:142:TRP:CE2	1:A:144:VAL:HB	2.54	0.43
1:C:185:LEU:HD21	1:C:189:ARG:C	2.39	0.43
1:A:194:TYR:O	1:A:195:LEU:HD23	2.19	0.43
1:C:147:VAL:CG1	1:C:149:PHE:CE1	3.01	0.43
1:D:94:GLY:N	1:D:115:VAL:O	2.46	0.43
1:A:229:ARG:HH21	1:A:236:ARG:HH22	1.67	0.43
1:C:160:GLU:HB2	1:C:163:ALA:HB2	2.00	0.43
1:C:9:LEU:CD1	1:C:9:LEU:N	2.81	0.43
1:D:100:ALA:O	1:D:169:LEU:HA	2.19	0.43
1:B:142:TRP:CE2	1:B:144:VAL:HB	2.54	0.43
1:B:216:GLU:C	1:B:217:ARG:HG3	2.39	0.43
1:D:224:ARG:HG2	1:D:224:ARG:HH11	1.83	0.43
1:B:34:VAL:HG11	1:B:48:VAL:CG1	2.49	0.42
1:D:89:LEU:HD23	1:D:192:VAL:HG23	2.00	0.42
1:A:118:LYS:HA	1:A:118:LYS:HD3	1.89	0.42
1:A:174:PRO:CG	1:A:195:LEU:HD21	2.49	0.42
1:D:214:ARG:NH2	3:D:402:HOH:O	2.52	0.42
1:C:129:HIS:CD2	1:C:130:HIS:HB3	2.54	0.42
1:B:64:GLU:CD	1:D:217:ARG:HH22	2.20	0.42
1:B:72:ARG:NH1	1:B:78:TYR:CZ	2.87	0.42
1:C:159:LEU:HD13	1:C:184:ALA:HB2	2.02	0.42
1:C:72:ARG:HB2	1:C:76:PRO:HG3	2.01	0.42
1:C:155:GLU:HA	1:C:180:LYS:CD	2.49	0.42
1:A:20:PHE:CD1	1:A:20:PHE:N	2.88	0.42
1:B:116:GLY:O	1:B:146:ASN:ND2	2.53	0.42
1:B:78:TYR:CD1	1:B:243:THR:HG21	2.55	0.42
1:A:192:VAL:HA	1:A:247:VAL:O	2.18	0.42
1:B:79:PRO:O	1:B:83:SER:HB3	2.20	0.42
1:B:71:LYS:HD3	1:B:71:LYS:HA	1.92	0.41
1:B:134:ALA:O	1:B:138:VAL:HG23	2.21	0.41
1:B:154:LEU:HD12	1:B:154:LEU:N	2.34	0.41
1:C:129:HIS:CD2	1:C:129:HIS:C	2.92	0.41
1:B:13:LYS:CD	1:B:15:ARG:HH21	2.34	0.41
1:B:72:ARG:HG3	1:B:76:PRO:HG3	2.02	0.41
1:B:13:LYS:HD2	1:B:15:ARG:HE	1.85	0.41
1:B:194:TYR:CE2	1:B:244:ALA:HB3	2.55	0.41
1:B:50:THR:O	1:B:51:HIS:ND1	2.53	0.41
1:C:214:ARG:HA	1:C:214:ARG:HD3	1.81	0.41
1:C:9:LEU:HD13	1:C:9:LEU:N	2.35	0.41
1:A:155:GLU:C	1:A:180:LYS:HZ1	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:93:PRO:HA	1:D:115:VAL:O	2.21	0.41
1:B:22:LYS:H	1:B:36:HIS:CD2	2.38	0.41
1:C:211:HIS:HB3	1:C:212:PRO:CD	2.50	0.41
1:C:96:ARG:NH1	1:C:163:ALA:HB1	2.36	0.41
1:A:98:LEU:HD23	1:A:167:VAL:HG13	2.03	0.41
1:D:127:ARG:HG2	1:D:127:ARG:H	1.76	0.41
1:B:242:HIS:CB	1:C:235:PRO:HB2	2.51	0.40
1:C:127:ARG:CZ	1:C:129:HIS:ND1	2.84	0.40
1:C:20:PHE:N	1:C:20:PHE:CD1	2.89	0.40
1:B:190:PHE:CD1	1:B:190:PHE:N	2.89	0.40
1:B:72:ARG:NH2	1:B:76:PRO:C	2.74	0.40
1:A:65:GLU:OE2	1:C:250:ARG:NH2	2.55	0.40
1:B:195:LEU:HA	1:B:196:PRO:HD2	1.93	0.40
1:D:192:VAL:HA	1:D:247:VAL:O	2.21	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	212/251 (84%)	195 (92%)	13 (6%)	4 (2%)	9	39
1	B	249/251 (99%)	237 (95%)	8 (3%)	4 (2%)	11	43
1	C	204/251 (81%)	188 (92%)	12 (6%)	4 (2%)	9	37
1	D	195/251 (78%)	181 (93%)	10 (5%)	4 (2%)	8	36
All	All	860/1004 (86%)	801 (93%)	43 (5%)	16 (2%)	9	39

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	102	THR

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Mol	Chain	Res	Type
1	B	30	HIS
1	D	254	GLY
1	A	116	GLY
1	B	76	PRO
1	A	117	GLU
1	B	254	GLY
1	C	241	GLY
1	D	61	PRO
1	A	242	HIS
1	C	240	VAL
1	D	101	GLY
1	B	240	VAL
1	D	173	GLU
1	C	103	GLY
1	C	128	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	175/201 (87%)	160 (91%)	15 (9%)	12	44
1	B	201/201 (100%)	188 (94%)	13 (6%)	20	55
1	C	169/201 (84%)	142 (84%)	27 (16%)	3	12
1	D	157/201 (78%)	140 (89%)	17 (11%)	7	30
All	All	702/804 (87%)	630 (90%)	72 (10%)	8	32

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	37	GLU
1	A	75	THR
1	A	83	SER
1	A	102	THR
1	A	125	GLU

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Mol	Chain	Res	Type
1	A	127	ARG
1	A	129	HIS
1	A	145	GLU
1	A	159	LEU
1	A	172	MET
1	A	183	LEU
1	A	214	ARG
1	A	242	HIS
1	A	253	LYS
1	B	13	LYS
1	B	31	LYS
1	B	36	HIS
1	B	41	GLU
1	B	51	HIS
1	B	72	ARG
1	B	75	THR
1	B	123	SER
1	B	143	GLN
1	B	175	TRP
1	B	185	LEU
1	B	209	GLU
1	B	229	ARG
1	C	9	LEU
1	C	57	SER
1	C	62	THR
1	C	68	LEU
1	C	72	ARG
1	C	77	THR
1	C	80	LYS
1	C	83	SER
1	C	104	SER
1	C	117	GLU
1	C	120	LEU
1	C	129	HIS
1	C	130	HIS
1	C	133	GLN
1	C	143	GLN
1	C	149	PHE
1	C	151	LEU
1	C	153	LYS
1	C	155	GLU
1	C	158	GLU

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Mol	Chain	Res	Type
1	C	160	GLU
1	C	185	LEU
1	C	214	ARG
1	C	224	ARG
1	C	236	ARG
1	C	238	GLN
1	C	253	LYS
1	D	72	ARG
1	D	73	SER
1	D	77	THR
1	D	80	LYS
1	D	98	LEU
1	D	117	GLU
1	D	127	ARG
1	D	143	GLN
1	D	158	GLU
1	D	159	LEU
1	D	160	GLU
1	D	180	LYS
1	D	214	ARG
1	D	220	GLU
1	D	224	ARG
1	D	242	HIS
1	D	253	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	GLN
1	B	36	HIS
1	B	150	HIS
1	C	143	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry

3 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	301	-	4,4,4	0.25	0	6,6,6	0.53	0
2	SO4	B	302	-	4,4,4	0.10	0	6,6,6	0.32	0
2	SO4	D	301	-	4,4,4	0.09	0	6,6,6	0.21	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	SO4	B	301	-	-	0/0/0/0	0/0/0/0
2	SO4	B	302	-	-	0/0/0/0	0/0/0/0
2	SO4	D	301	-	-	0/0/0/0	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	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	220/251 (87%)	0.18	6 (2%) 55 30	45, 71, 107, 125	0
1	B	251/251 (100%)	-0.16	1 (0%) 92 84	43, 65, 107, 135	0
1	C	210/251 (83%)	0.14	5 (2%) 59 37	49, 76, 109, 127	0
1	D	197/251 (78%)	-0.16	0 100 100	46, 69, 102, 118	0
All	All	878/1004 (87%)	-0.01	12 (1%) 75 57	43, 71, 107, 135	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	57	SER	3.3
1	A	44	PRO	2.6
1	A	18	LEU	2.4
1	C	16	ALA	2.3
1	C	163	ALA	2.3
1	A	40	LEU	2.3
1	C	19	VAL	2.3
1	C	117	GLU	2.3
1	C	145	GLU	2.2
1	A	41	GLU	2.2
1	A	144	VAL	2.1
1	B	73	SER	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	SO4	D	301	5/5	0.97	0.12	-1.54	90,94,95,98	0
2	SO4	B	302	5/5	0.97	0.15	-1.63	92,100,113,116	0
2	SO4	B	301	5/5	0.99	0.13	-2.17	74,77,83,88	0

6.5 Other polymers [i](#)

There are no such residues in this entry.