



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 6, 2016 – 07:37 AM EST

PDB ID : 5E9W
Title : Crystal structure of mRNA cap guanine-N7 methyltransferase obtained by limited proteolysis
Authors : PETIT, P.; COWLING, V.H.
Deposited on : 2015-10-15
Resolution : 2.28 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : rb-20028442

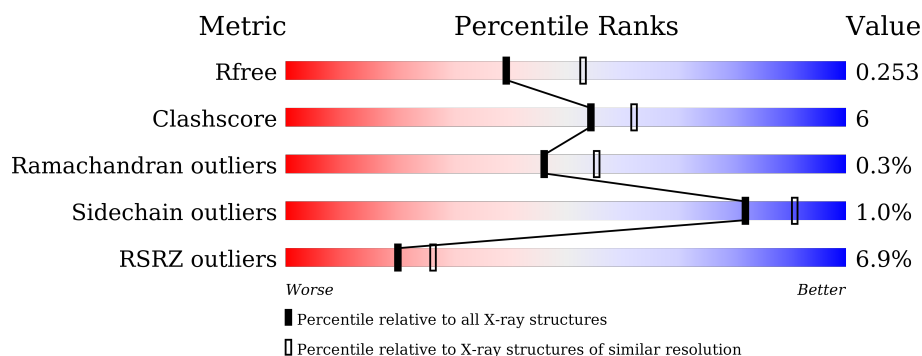
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.28 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}	91344	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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	312	<div> <div>8%</div> <div> <div></div> <div>73%</div> <div>13%</div> <div>•</div> <div>13%</div> </div> </div>
1	B	312	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>13%</div> <div></div> <div>13%</div> </div> </div>
1	C	312	<div> <div>8%</div> <div> <div></div> <div>73%</div> <div>14%</div> <div></div> <div>13%</div> </div> </div>
1	D	312	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>11%</div> <div>•</div> <div>12%</div> </div> </div>

2 Entry composition [i](#)

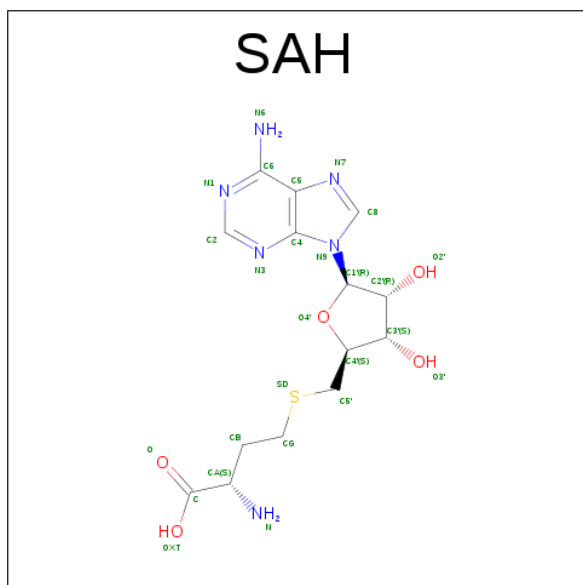
There are 3 unique types of molecules in this entry. The entry contains 9400 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 mRNA cap guanine-N7 methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2237	1441	364	414	18			
1	B	273	Total	C	N	O	S	0	0	0
			2252	1450	371	413	18			
1	C	273	Total	C	N	O	S	0	0	0
			2247	1447	368	414	18			
1	D	274	Total	C	N	O	S	0	0	0
			2253	1451	370	414	18			

- Molecule 2 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
2	B	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
2	D	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

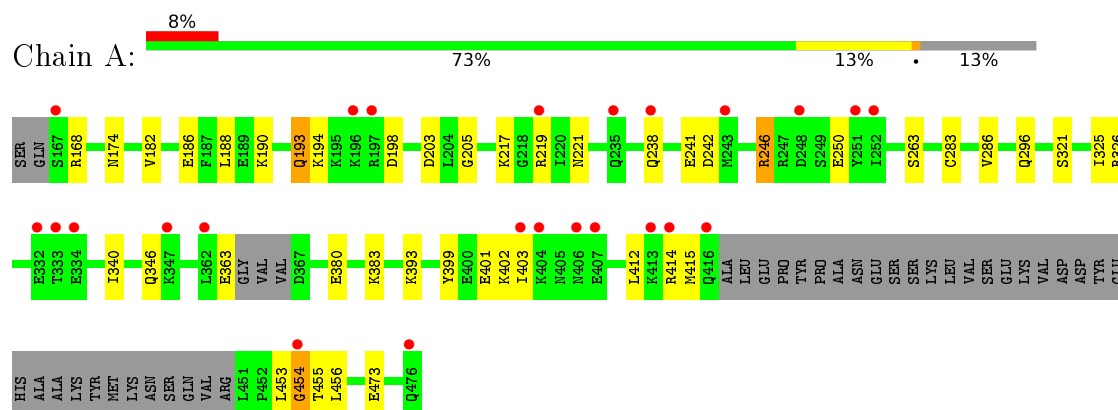
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	56	Total	O	0	0
			56	56		
3	B	94	Total	O	0	0
			94	94		
3	C	63	Total	O	0	0
			63	63		
3	D	94	Total	O	0	0
			94	94		

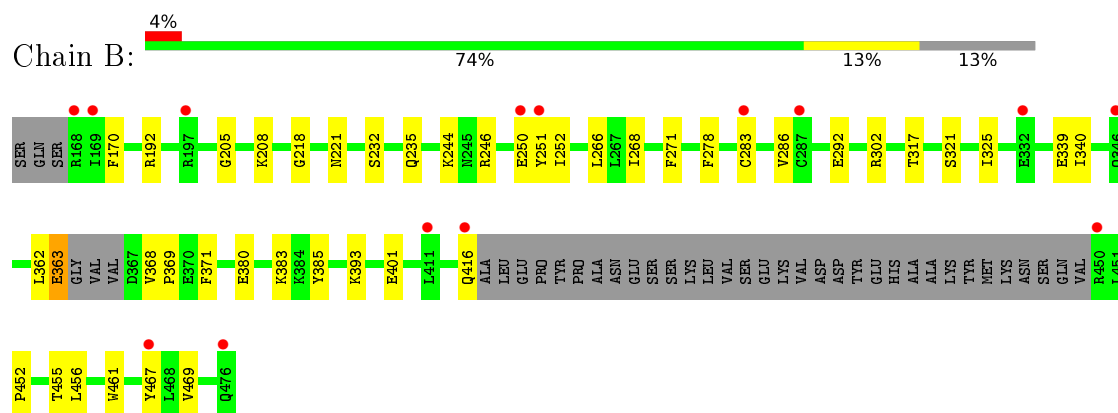
3 Residue-property plots [i](#)

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

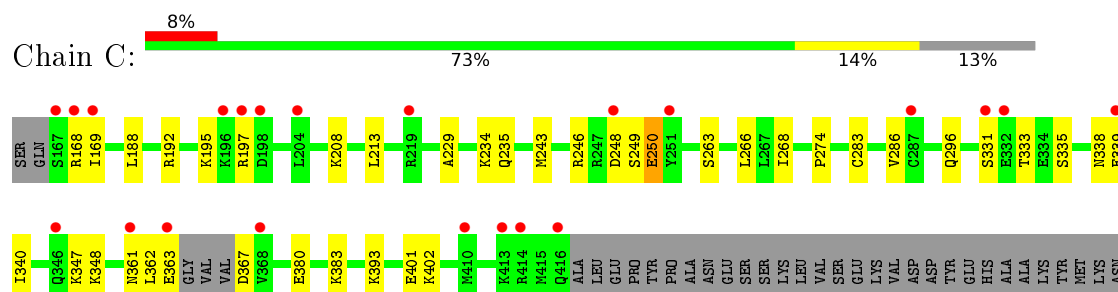
- Molecule 1: mRNA cap guanine-N7 methyltransferase



- Molecule 1: mRNA cap guanine-N7 methyltransferase

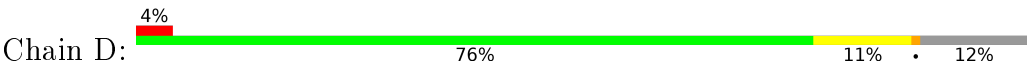


- Molecule 1: mRNA cap guanine-N7 methyltransferase





● Molecule 1: mRNA cap guanine-N7 methyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.46Å 99.16Å 167.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.58 – 2.28 29.58 – 2.28	Depositor EDS
% Data completeness (in resolution range)	96.3 (29.58-2.28) 96.4 (29.58-2.28)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.55 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1702)	Depositor
R, R_{free}	0.220 , 0.255 0.219 , 0.253	Depositor DCC
R_{free} test set	2888 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	21.9	Xtriage
Anisotropy	0.918	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9400	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 44.42 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.5293e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: SAH

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	0.49	0/2283	0.63	1/3062 (0.0%)
1	B	0.49	0/2298	0.62	0/3079
1	C	0.49	0/2293	0.63	1/3073 (0.0%)
1	D	0.48	0/2299	0.59	0/3081
All	All	0.49	0/9173	0.62	2/12295 (0.0%)

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	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	454	GLY	N-CA-C	8.33	133.93	113.10
1	C	451	LEU	CA-CB-CG	5.86	128.77	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	346	GLN	Peptide
1	A	453	LEU	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	2237	0	2192	30	4
1	B	2252	0	2222	27	4
1	C	2247	0	2214	38	2
1	D	2253	0	2220	21	2
2	A	26	0	19	0	0
2	B	26	0	19	0	0
2	C	26	0	19	0	0
2	D	26	0	19	0	0
3	A	56	0	0	3	0
3	B	94	0	0	2	0
3	C	63	0	0	4	0
3	D	94	0	0	3	0
All	All	9400	0	8924	115	6

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 6.

All (115) 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:246:ARG:HH11	1:B:246:ARG:HG2	1.29	0.96
1:A:198:ASP:OD2	1:A:221:ASN:ND2	2.06	0.89
1:D:192:ARG:NH2	1:D:218:GLY:O	2.12	0.82
1:C:168:ARG:HB2	1:C:363:GLU:OE1	1.81	0.81
1:C:333:THR:HG22	1:C:335:SER:H	1.52	0.75
1:C:476:GLN:O	3:C:601:HOH:O	2.06	0.74
1:C:195:LYS:HE2	1:C:197:ARG:HE	1.55	0.72
1:A:380:GLU:OE2	1:A:383:LYS:NZ	2.26	0.69
1:D:380:GLU:OE2	1:D:383:LYS:NZ	2.25	0.69
1:B:302:ARG:NH2	3:B:603:HOH:O	2.24	0.69
1:B:208:LYS:HD3	1:B:232:SER:HA	1.73	0.69
1:D:415:MET:O	1:D:416:GLN:HG2	1.93	0.68
1:C:361:ASN:ND2	1:C:367:ASP:OD1	2.26	0.68
1:B:246:ARG:NH1	1:B:246:ARG:HG2	1.99	0.68
1:C:451:LEU:HD12	1:C:452:PRO:HD2	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:339:GLU:OE2	1:C:459:SER:OG	2.09	0.65
1:C:380:GLU:OE2	1:C:383:LYS:NZ	2.29	0.65
1:A:186:GLU:OE2	1:A:190:LYS:NZ	2.30	0.64
1:B:467:TYR:OH	3:B:601:HOH:O	2.07	0.64
1:B:244:LYS:HG3	1:B:252:ILE:HG21	1.80	0.63
1:A:414:ARG:HG3	1:A:414:ARG:HH21	1.63	0.63
1:C:192:ARG:NH1	3:C:603:HOH:O	2.17	0.62
1:A:238:GLN:NE2	1:A:242:ASP:OD1	2.33	0.61
1:D:169:ILE:HD11	1:D:340:ILE:HG23	1.81	0.60
1:A:194:LYS:NZ	1:A:473:GLU:OE2	2.34	0.60
1:C:456:LEU:HB2	1:C:461:TRP:NE1	2.15	0.60
1:C:234:LYS:NZ	3:C:605:HOH:O	2.31	0.59
1:D:399:TYR:O	1:D:403:ILE:HG13	2.02	0.59
1:B:192:ARG:NH2	1:B:218:GLY:O	2.31	0.59
1:D:195:LYS:HE3	1:D:279:ASP:HB3	1.84	0.59
1:C:456:LEU:HB2	1:C:461:TRP:CD1	2.37	0.59
1:B:283:CYS:SG	1:B:286:VAL:HB	2.44	0.58
1:A:412:LEU:HD21	1:A:455:THR:HG21	1.86	0.57
1:B:244:LYS:CG	1:B:252:ILE:HG21	2.35	0.57
1:A:283:CYS:SG	1:A:286:VAL:HB	2.45	0.56
1:A:193:GLN:HG2	1:C:229:ALA:HB1	1.88	0.56
1:C:451:LEU:CD1	1:C:452:PRO:HD2	2.36	0.56
1:D:393:LYS:HE2	1:D:401:GLU:OE1	2.06	0.56
1:B:380:GLU:OE2	1:B:383:LYS:NZ	2.39	0.56
1:B:221:ASN:ND2	1:B:251:TYR:OH	2.38	0.55
1:B:170:PHE:HZ	1:B:416:GLN:HE21	1.55	0.55
1:A:174:ASN:ND2	1:A:415:MET:O	2.40	0.54
1:C:347:LYS:NZ	3:C:602:HOH:O	2.15	0.54
1:B:339:GLU:HG3	1:B:340:ILE:HG13	1.89	0.54
1:C:451:LEU:HD12	1:C:452:PRO:CD	2.38	0.53
1:A:198:ASP:CB	1:A:221:ASN:HD22	2.21	0.52
1:C:401:GLU:HG3	1:C:402:LYS:HD2	1.92	0.51
1:D:326:ARG:NH1	3:D:602:HOH:O	2.29	0.51
1:A:203:ASP:OD2	3:A:601:HOH:O	2.19	0.51
1:A:393:LYS:HE2	1:A:401:GLU:OE1	2.11	0.51
1:D:239:ARG:NH2	3:D:609:HOH:O	2.45	0.50
1:B:393:LYS:HE2	1:B:401:GLU:OE1	2.11	0.50
1:B:455:THR:HG22	1:B:456:LEU:HD12	1.93	0.50
1:D:191:VAL:HG13	1:D:195:LYS:HE2	1.94	0.50
1:A:340:ILE:O	1:A:363:GLU:HG2	2.12	0.50
1:C:393:LYS:HE2	1:C:401:GLU:OE1	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:GLY:HA3	1:B:286:VAL:HG21	1.94	0.49
1:A:454:GLY:O	1:A:456:LEU:N	2.42	0.49
1:D:269:ASP:O	3:D:601:HOH:O	2.19	0.49
1:D:266:LEU:HD11	1:D:302:ARG:HD3	1.94	0.49
1:C:213:LEU:HG	1:C:243:MET:HE1	1.96	0.48
1:A:241:GLU:HA	1:A:241:GLU:OE1	2.13	0.48
1:D:182:VAL:HG12	1:D:402:LYS:HE3	1.96	0.48
1:B:362:LEU:O	1:B:363:GLU:HB2	2.13	0.48
1:B:302:ARG:HB2	1:B:385:TYR:CZ	2.48	0.47
1:B:321:SER:O	1:B:325:ILE:HG12	2.13	0.47
1:D:275:GLN:N	1:D:275:GLN:OE1	2.41	0.47
1:C:168:ARG:CB	1:C:363:GLU:OE1	2.58	0.47
1:D:337:GLY:HA2	1:D:341:TYR:CZ	2.50	0.47
1:C:169:ILE:HG23	1:C:363:GLU:CD	2.36	0.46
1:B:452:PRO:HG2	1:B:461:TRP:CE2	2.51	0.46
1:C:283:CYS:SG	1:C:286:VAL:HB	2.55	0.46
1:D:283:CYS:SG	1:D:286:VAL:HB	2.55	0.46
1:D:266:LEU:HG	1:D:268:ILE:HG12	1.98	0.46
1:D:244:LYS:HB2	1:D:252:ILE:HG21	1.97	0.46
1:B:266:LEU:HG	1:B:268:ILE:HG12	1.98	0.45
1:C:333:THR:HG22	1:C:335:SER:N	2.26	0.45
1:B:455:THR:C	1:B:456:LEU:HD12	2.36	0.45
1:C:249:SER:O	1:C:250:GLU:O	2.33	0.45
1:A:174:ASN:HB3	3:A:605:HOH:O	2.17	0.45
1:A:205:GLY:HA3	1:A:286:VAL:HG21	1.99	0.45
1:C:266:LEU:HG	1:C:268:ILE:HG22	1.99	0.45
1:D:202:LEU:HB2	1:D:278:PHE:CE1	2.52	0.45
1:A:217:LYS:HA	1:A:217:LYS:HE3	2.00	0.44
1:A:246:ARG:HD2	1:A:246:ARG:HA	1.71	0.44
1:A:182:VAL:HG11	1:A:402:LYS:HG3	1.98	0.44
1:C:331:SER:O	1:C:348:LYS:NZ	2.51	0.44
1:B:271:PHE:CZ	1:B:278:PHE:HZ	2.36	0.43
1:B:292:GLU:HB2	1:B:371:PHE:CE2	2.53	0.43
1:A:198:ASP:HB3	1:A:221:ASN:HD22	1.83	0.43
1:A:263:SER:HB2	1:A:296:GLN:HB3	2.00	0.43
1:C:268:ILE:HA	1:C:268:ILE:HD12	1.81	0.43
1:C:208:LYS:HD3	1:C:235:GLN:OE1	2.19	0.43
1:D:395:PHE:CE1	1:D:467:TYR:HB3	2.53	0.43
1:C:263:SER:HB2	1:C:296:GLN:HB3	2.00	0.42
1:A:399:TYR:O	1:A:403:ILE:HG12	2.19	0.42
1:C:456:LEU:HB2	1:C:461:TRP:CE2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:452:PRO:HG2	1:C:461:TRP:CZ2	2.55	0.42
1:A:414:ARG:NH2	1:A:414:ARG:HG3	2.31	0.42
1:C:168:ARG:N	1:C:168:ARG:HD3	2.34	0.42
1:C:451:LEU:HD12	1:C:452:PRO:N	2.34	0.42
1:B:208:LYS:HE2	1:B:235:GLN:OE1	2.20	0.42
1:C:268:ILE:HD11	1:C:274:PRO:O	2.19	0.42
1:B:317:THR:HG22	1:B:469:VAL:HG12	2.01	0.41
1:A:188:LEU:HA	1:A:188:LEU:HD23	1.94	0.41
1:C:338:ASN:HB2	1:C:339:GLU:OE1	2.20	0.41
1:C:362:LEU:HA	1:C:362:LEU:HD12	1.94	0.41
1:B:368:VAL:HA	1:B:369:PRO:HD3	1.92	0.41
1:C:188:LEU:HD23	1:C:188:LEU:HA	1.97	0.41
1:C:339:GLU:HG2	1:C:340:ILE:HG13	2.02	0.41
1:A:198:ASP:CG	1:A:221:ASN:HD22	2.23	0.41
1:A:321:SER:O	1:A:325:ILE:HG12	2.20	0.41
1:D:343:VAL:HG22	1:D:360:PHE:HD1	1.86	0.41
1:A:326:ARG:NH2	3:A:602:HOH:O	2.29	0.40
1:A:198:ASP:CG	1:A:221:ASN:ND2	2.73	0.40

All (6) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:ARG:NH1	1:B:416:GLN:NE2[4_466]	0.40	1.80
1:C:246:ARG:NH1	1:D:416:GLN:NE2[4_366]	1.05	1.15
1:C:246:ARG:NH1	1:D:416:GLN:OE1[4_366]	1.51	0.69
1:A:246:ARG:CZ	1:B:416:GLN:NE2[4_466]	1.67	0.53
1:A:246:ARG:NH1	1:B:416:GLN:OE1[4_466]	2.08	0.12
1:A:246:ARG:CZ	1:B:416:GLN:CD[4_466]	2.16	0.04

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	267/312 (86%)	258 (97%)	8 (3%)	1 (0%)	39	47
1	B	267/312 (86%)	260 (97%)	6 (2%)	1 (0%)	39	47
1	C	267/312 (86%)	258 (97%)	8 (3%)	1 (0%)	39	47
1	D	268/312 (86%)	258 (96%)	10 (4%)	0	100	100
All	All	1069/1248 (86%)	1034 (97%)	32 (3%)	3 (0%)	46	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	250	GLU
1	A	250	GLU
1	B	250	GLU

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	244/285 (86%)	240 (98%)	4 (2%)	70	83
1	B	246/285 (86%)	245 (100%)	1 (0%)	93	97
1	C	246/285 (86%)	244 (99%)	2 (1%)	86	93
1	D	246/285 (86%)	243 (99%)	3 (1%)	78	88
All	All	982/1140 (86%)	972 (99%)	10 (1%)	82	91

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	168	ARG
1	A	193	GLN
1	A	219	ARG
1	A	246	ARG
1	B	363	GLU
1	C	248	ASP
1	C	451	LEU
1	D	169	ILE

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Mol	Chain	Res	Type
1	D	174	ASN
1	D	244	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	221	ASN
1	B	416	GLN

5.3.3 RNA [i](#)

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	SAH	A	500	-	22,28,28	0.98	2 (9%)	18,40,40	3.08	4 (22%)
2	SAH	B	501	-	22,28,28	1.05	2 (9%)	18,40,40	3.02	4 (22%)
2	SAH	C	501	-	22,28,28	0.97	2 (9%)	18,40,40	3.06	4 (22%)
2	SAH	D	501	-	22,28,28	1.05	2 (9%)	18,40,40	3.20	4 (22%)

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	SAH	A	500	-	-	0/7/31/31	0/3/3/3
2	SAH	B	501	-	-	0/7/31/31	0/3/3/3
2	SAH	C	501	-	-	0/7/31/31	0/3/3/3
2	SAH	D	501	-	-	0/7/31/31	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	SAH	C2-N1	2.03	1.37	1.33
2	A	500	SAH	C2-N1	2.06	1.37	1.33
2	D	501	SAH	C2-N1	2.17	1.38	1.33
2	B	501	SAH	C2-N1	2.27	1.38	1.33
2	D	501	SAH	C2-N3	2.76	1.37	1.32
2	C	501	SAH	C2-N3	3.15	1.37	1.32
2	A	500	SAH	C2-N3	3.25	1.37	1.32
2	B	501	SAH	C2-N3	3.41	1.38	1.32

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	SAH	N3-C2-N1	-12.18	119.31	128.87
2	A	500	SAH	N3-C2-N1	-11.56	119.79	128.87
2	B	501	SAH	N3-C2-N1	-11.38	119.93	128.87
2	C	501	SAH	N3-C2-N1	-11.16	120.10	128.87
2	C	501	SAH	C1'-N9-C4	-3.89	122.46	126.81
2	C	501	SAH	C5'-SD-CG	-3.29	92.42	102.42
2	A	500	SAH	C1'-N9-C4	-3.29	123.13	126.81
2	D	501	SAH	C1'-N9-C4	-3.03	123.42	126.81
2	A	500	SAH	C5'-SD-CG	-3.01	93.28	102.42
2	B	501	SAH	C1'-N9-C4	-2.93	123.54	126.81
2	B	501	SAH	C5'-SD-CG	-2.40	95.14	102.42
2	D	501	SAH	C5'-SD-CG	-2.32	95.36	102.42
2	C	501	SAH	O4'-C1'-N9	2.16	112.18	108.11
2	B	501	SAH	O4'-C1'-N9	2.80	113.39	108.11
2	A	500	SAH	O4'-C1'-N9	2.92	113.63	108.11
2	D	501	SAH	O4'-C1'-N9	3.36	114.45	108.11

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 [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	273/312 (87%)	0.37	24 (8%) 12 17	13, 30, 65, 93	0
1	B	273/312 (87%)	0.18	14 (5%) 32 39	10, 24, 54, 82	0
1	C	273/312 (87%)	0.33	24 (8%) 12 17	13, 28, 66, 84	0
1	D	274/312 (87%)	0.23	13 (4%) 35 43	10, 25, 60, 104	0
All	All	1093/1248 (87%)	0.28	75 (6%) 20 26	10, 26, 63, 104	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	167	SER	6.8
1	D	167	SER	5.9
1	A	251	TYR	5.3
1	B	416	GLN	5.3
1	C	451	LEU	4.9
1	A	167	SER	4.7
1	A	407	GLU	4.7
1	D	416	GLN	4.6
1	D	197	ARG	4.5
1	D	168	ARG	4.4
1	D	196	LYS	4.2
1	B	250	GLU	4.1
1	B	168	ARG	4.0
1	B	450	ARG	4.0
1	A	414	ARG	3.8
1	C	168	ARG	3.8
1	D	195	LYS	3.6
1	C	332	GLU	3.5
1	B	332	GLU	3.3
1	D	332	GLU	3.3
1	B	476	GLN	3.3

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Mol	Chain	Res	Type	RSRZ
1	C	196	LYS	3.2
1	A	332	GLU	3.2
1	C	197	ARG	3.1
1	B	251	TYR	3.1
1	A	413	LYS	3.0
1	A	404	LYS	3.0
1	C	287	CYS	3.0
1	C	339	GLU	3.0
1	D	219	ARG	3.0
1	C	363	GLU	2.9
1	A	347	LYS	2.9
1	B	287	CYS	2.9
1	C	219	ARG	2.8
1	D	346	GLN	2.8
1	C	169	ILE	2.7
1	B	197	ARG	2.7
1	A	476	GLN	2.7
1	B	346	GLN	2.7
1	C	410	MET	2.7
1	D	169	ILE	2.6
1	B	467	TYR	2.6
1	A	219	ARG	2.6
1	A	454	GLY	2.5
1	D	170	PHE	2.5
1	C	416	GLN	2.5
1	A	196	LYS	2.4
1	A	197	ARG	2.4
1	A	334	GLU	2.4
1	D	333	THR	2.4
1	C	476	GLN	2.4
1	C	198	ASP	2.3
1	A	243	MET	2.3
1	A	333	THR	2.3
1	B	169	ILE	2.3
1	B	411	LEU	2.3
1	C	251	TYR	2.3
1	C	413	LYS	2.3
1	A	416	GLN	2.2
1	C	331	SER	2.2
1	C	361	ASN	2.2
1	C	248	ASP	2.2
1	A	238	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	204	LEU	2.2
1	C	346	GLN	2.2
1	C	414	ARG	2.1
1	C	368	VAL	2.1
1	A	248	ASP	2.1
1	B	283	CYS	2.1
1	A	362	LEU	2.0
1	A	235	GLN	2.0
1	D	287	CYS	2.0
1	A	406	ASN	2.0
1	A	252	ILE	2.0
1	A	403	ILE	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(\AA^2)	Q<0.9
2	SAH	B	501	26/26	0.97	0.13	-0.11	9,14,15,15	0
2	SAH	C	501	26/26	0.95	0.12	-0.40	12,16,22,25	0
2	SAH	D	501	26/26	0.97	0.11	-0.64	10,12,14,24	0
2	SAH	A	500	26/26	0.97	0.10	-0.85	16,19,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.