



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 03:53 PM GMT

PDB ID : 1F7U
Title : CRYSTAL STRUCTURE OF THE ARGINYL-TRNA SYNTHETASE COM-
PLEXED WITH THE TRNA(ARG) AND L-ARG
Authors : Delagoutte, B.; Moras, D.; Cavarelli, J.
Deposited on : 2000-06-28
Resolution : 2.20 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

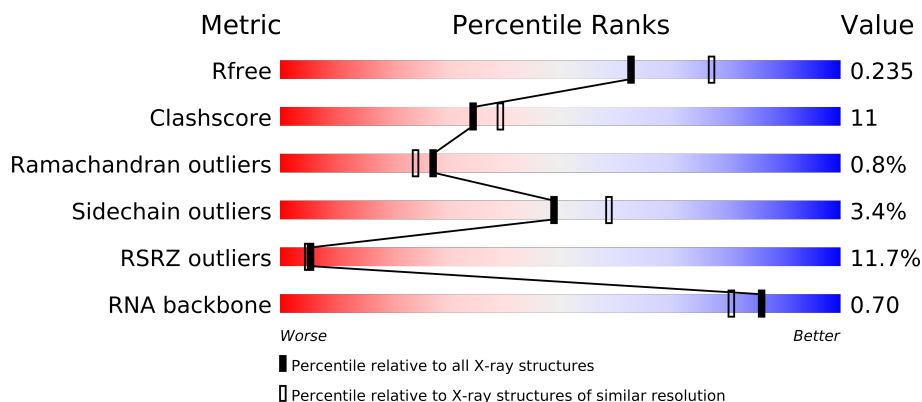
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.20 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}	66092	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)
RNA backbone	1838	1120 (3.00-1.40)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	B	76	
2	A	607	

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 7126 atoms, of which 0 are hydrogen and 0 are deuterium.

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 RNA chain called TRNA(ARG).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	76	Total	C	N	O	P	0	0	0
			1629	730	290	533	76			

- Molecule 2 is a protein called ARGINYL-TRNA SYNTHETASE.

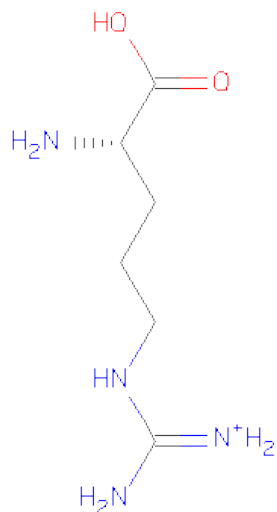
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	606	Total	C	N	O	S	0	0	0
			4892	3138	828	905	21			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 4 is ARGinine (three-letter code: ARG) (formula: C₆H₁₅N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			12	6	4	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	348	Total	O	0	0
			348	348		
5	B	240	Total	O	0	0
			240	240		

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 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.

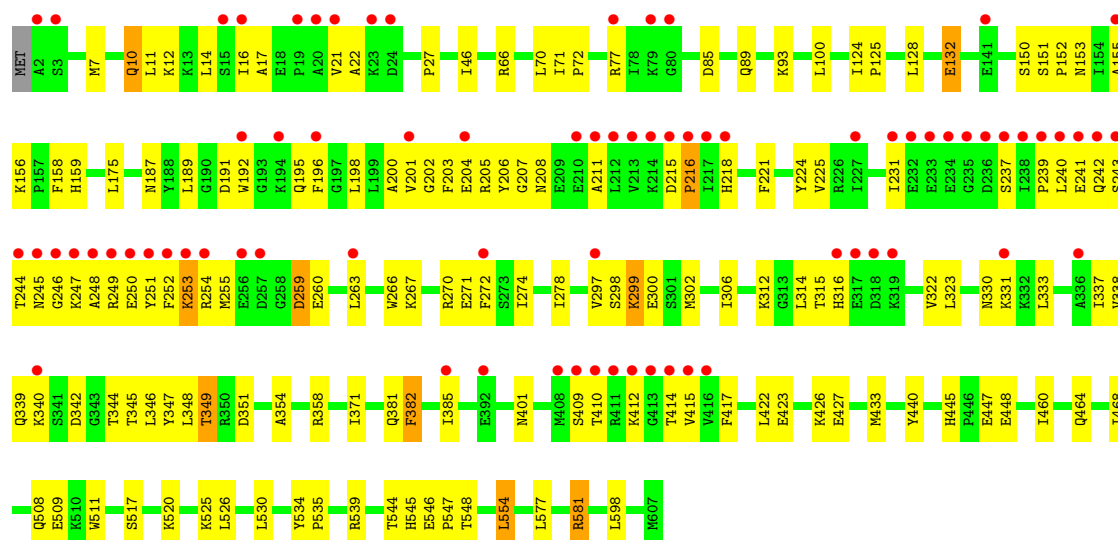
• Molecule 1: TRNA(ARG)

Chain B: 



• Molecule 2: ARGINYL-TRNA SYNTHETASE

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	129.65Å 107.47Å 71.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.98 – 2.20 24.82 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (14.98-2.20) 99.8 (24.82-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.71 (at 2.19Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.190 , 0.233 0.193 , 0.235	Depositor DCC
R_{free} test set	2562 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	20.7	Xtriage
Anisotropy	0.725	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 49.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 51242 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7126	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 5MU, H2U, SO4, 2MG, 5MC, 1MA, M2G, 1MG, PSU

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	B	0.30	0/1532	0.69	0/2387
2	A	0.30	0/4996	0.55	0/6736
All	All	0.30	0/6528	0.59	0/9123

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	2	0

There are no bond length outliers.

There are no bond angle outliers.

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	934	I	C3',C1'

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1629	0	834	18	0
2	A	4892	0	4906	119	0
3	B	5	0	0	0	0
4	A	12	0	12	1	0
5	A	348	0	0	3	0
5	B	240	0	0	2	0
All	All	7126	0	5752	133	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (133) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:299:LYS:H	2:A:299:LYS:HE3	1.11	1.15
2:A:331:LYS:H	2:A:331:LYS:HD2	1.24	1.00
2:A:323:LEU:HB3	2:A:337:ILE:HD13	1.53	0.90
2:A:299:LYS:N	2:A:299:LYS:HE3	1.94	0.83
2:A:331:LYS:HD2	2:A:331:LYS:N	1.97	0.79
2:A:299:LYS:H	2:A:299:LYS:CE	1.96	0.76
2:A:409:SER:HB2	2:A:414:THR:HG22	1.69	0.73
2:A:331:LYS:H	2:A:331:LYS:CD	2.02	0.72
2:A:322:VAL:HG12	2:A:339:GLN:HB3	1.71	0.72
2:A:520:LYS:O	2:A:525:LYS:HE3	1.91	0.71
2:A:153:ASN:HB3	2:A:195:GLN:OE1	1.93	0.69
2:A:246:GLY:O	2:A:249:ARG:HG2	1.93	0.69
2:A:241:GLU:HA	2:A:246:GLY:HA3	1.74	0.68
2:A:244:THR:HA	2:A:247:LYS:HE2	1.76	0.68
2:A:17:ALA:HB1	2:A:21:VAL:HG13	1.76	0.67
1:B:974:C:H2'	2:A:340:LYS:HG3	1.78	0.66
2:A:231:ILE:HD12	2:A:245:ASN:HD22	1.61	0.66
2:A:460:ILE:O	2:A:464:GLN:HG3	1.97	0.64
2:A:581:ARG:HD3	5:A:1043:HOH:O	1.98	0.63
1:B:942:G:O2'	1:B:943:A:H5'	2.00	0.62
2:A:300:GLU:CD	2:A:300:GLU:H	2.02	0.62
2:A:525:LYS:HE2	5:A:1072:HOH:O	2.00	0.62
2:A:322:VAL:CG1	2:A:339:GLN:HB3	2.30	0.60
2:A:322:VAL:HG13	2:A:338:VAL:HG23	1.84	0.60
2:A:155:ALA:C	2:A:156:LYS:HD2	2.21	0.60
2:A:316:HIS:NE2	2:A:323:LEU:HD11	2.17	0.60
2:A:198:LEU:HD23	2:A:224:TYR:CE1	2.38	0.59
2:A:348:LEU:CD1	2:A:382:PHE:HB3	2.32	0.59
2:A:205:ARG:HG2	2:A:206:TYR:CE1	2.36	0.59
2:A:371:ILE:O	2:A:401:ASN:HA	2.03	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:251:TYR:O	2:A:255:MET:HG3	2.03	0.59
1:B:928:C:O2'	1:B:929:U:H5'	2.03	0.58
2:A:445:HIS:HB3	2:A:448:GLU:OE1	2.03	0.58
2:A:196:PHE:CZ	2:A:266:TRP:HA	2.39	0.58
2:A:208:ASN:HB3	2:A:211:ALA:HB3	1.84	0.58
2:A:200:ALA:O	2:A:204:GLU:HG3	2.03	0.57
2:A:239:PRO:HB2	2:A:242:GLN:HB3	1.86	0.57
2:A:259:ASP:O	2:A:263:LEU:HD23	2.04	0.57
2:A:241:GLU:CA	2:A:246:GLY:HA3	2.34	0.56
2:A:196:PHE:HZ	2:A:266:TRP:HA	1.71	0.56
2:A:338:VAL:O	2:A:349:THR:HG21	2.06	0.56
2:A:70:LEU:C	2:A:70:LEU:HD12	2.25	0.56
2:A:346:LEU:H	2:A:349:THR:CG2	2.20	0.55
2:A:267:LYS:O	2:A:271:GLU:HB2	2.07	0.55
2:A:348:LEU:HD12	2:A:382:PHE:HB3	1.89	0.55
2:A:10:GLN:N	2:A:10:GLN:HE21	2.06	0.54
1:B:926:M2G:HM22	5:B:1244:HOH:O	2.08	0.53
2:A:17:ALA:HB1	2:A:21:VAL:CG1	2.38	0.53
2:A:221:PHE:O	2:A:225:VAL:HG23	2.09	0.53
1:B:974:C:O2	2:A:340:LYS:HG3	2.10	0.52
2:A:539:ARG:O	2:A:539:ARG:HD3	2.10	0.52
2:A:517:SER:O	2:A:520:LYS:HE3	2.11	0.51
2:A:156:LYS:N	2:A:156:LYS:HD2	2.25	0.51
2:A:156:LYS:HE3	2:A:224:TYR:CE2	2.45	0.51
2:A:345:THR:HB	2:A:349:THR:OG1	2.11	0.51
2:A:330:ASN:HB3	2:A:333:LEU:HG	1.93	0.51
2:A:577:LEU:O	2:A:581:ARG:HG3	2.12	0.50
2:A:546:GLU:HB3	2:A:548:THR:HG22	1.93	0.50
2:A:254:ARG:HE	2:A:259:ASP:CG	2.14	0.49
2:A:342:ASP:OD1	2:A:344:THR:HG23	2.12	0.49
2:A:423:GLU:O	2:A:427:GLU:HG3	2.13	0.49
2:A:315:THR:HB	2:A:322:VAL:CG2	2.43	0.49
1:B:964:G:O2'	1:B:965:G:H5'	2.14	0.48
2:A:534:TYR:HB3	2:A:535:PRO:HD3	1.96	0.48
2:A:302:MET:O	2:A:306:ILE:HG12	2.13	0.48
2:A:12:LYS:HA	2:A:27:PRO:HB2	1.96	0.48
2:A:14:LEU:HG	2:A:16:ILE:HG23	1.96	0.48
1:B:973:G:O2'	2:A:337:ILE:HG13	2.13	0.48
2:A:347:TYR:OH	4:A:800:ARG:HG2	2.13	0.48
2:A:252:PHE:HA	2:A:255:MET:HE3	1.95	0.48
2:A:323:LEU:HD12	2:A:323:LEU:C	2.34	0.48
2:A:241:GLU:N	2:A:246:GLY:HA3	2.29	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:252:PHE:HA	2:A:255:MET:CE	2.44	0.48
2:A:203:PHE:O	2:A:207:GLY:HA3	2.13	0.48
2:A:508:GLN:HA	2:A:511:TRP:CD2	2.50	0.47
1:B:901:PSU:H2'	1:B:902:U:C6	2.50	0.47
2:A:246:GLY:C	2:A:248:ALA:H	2.18	0.46
2:A:132:GLU:CD	2:A:132:GLU:H	2.18	0.46
2:A:409:SER:HB2	2:A:414:THR:CG2	2.44	0.46
2:A:152:PRO:HG3	2:A:158:PHE:CZ	2.51	0.46
2:A:85:ASP:O	2:A:89:GLN:HG3	2.16	0.46
2:A:70:LEU:O	2:A:70:LEU:HD12	2.15	0.46
2:A:191:ASP:OD2	2:A:270:ARG:NH1	2.49	0.46
2:A:215:ASP:OD2	2:A:218:HIS:HB2	2.15	0.46
2:A:274:ILE:O	2:A:278:ILE:HG13	2.16	0.46
2:A:433:MET:HG2	2:A:440:TYR:HD1	1.80	0.46
2:A:250:GLU:O	2:A:253:LYS:HG3	2.16	0.46
2:A:300:GLU:N	2:A:300:GLU:CD	2.68	0.45
2:A:348:LEU:HD11	2:A:382:PHE:HB3	1.99	0.45
1:B:919:H2U:H52	2:A:66:ARG:HD3	1.99	0.44
2:A:159:HIS:HA	2:A:417:PHE:HA	1.99	0.44
1:B:903:C:H2'	1:B:904:C:H6	1.83	0.44
2:A:151:SER:HA	2:A:187:ASN:OD1	2.18	0.44
2:A:205:ARG:HH21	2:A:244:THR:HG23	1.82	0.44
2:A:312:LYS:HB2	2:A:314:LEU:HG	1.99	0.44
2:A:381:GLN:O	2:A:385:ILE:HG13	2.18	0.43
2:A:298:SER:HB2	2:A:300:GLU:OE1	2.17	0.43
2:A:253:LYS:HD2	2:A:254:ARG:N	2.33	0.43
2:A:201:VAL:HG13	2:A:202:GLY:N	2.32	0.43
2:A:7:MET:O	2:A:11:LEU:HD22	2.18	0.43
2:A:547:PRO:HD2	5:A:1002:HOH:O	2.18	0.43
2:A:189:LEU:HD12	2:A:270:ARG:NH1	2.32	0.43
2:A:124:ILE:HB	2:A:125:PRO:HD3	2.01	0.43
2:A:17:ALA:O	2:A:22:ALA:HB2	2.19	0.42
2:A:253:LYS:C	2:A:253:LYS:HD2	2.39	0.42
2:A:175:LEU:HA	2:A:598:LEU:HD11	2.01	0.42
2:A:409:SER:O	2:A:415:VAL:HG13	2.19	0.42
2:A:346:LEU:H	2:A:349:THR:HG23	1.83	0.42
2:A:241:GLU:HA	2:A:246:GLY:CA	2.45	0.42
2:A:71:ILE:HA	2:A:72:PRO:HD3	1.85	0.42
2:A:426:LYS:CE	2:A:447:GLU:HG3	2.48	0.42
2:A:539:ARG:C	2:A:539:ARG:HD3	2.38	0.42
1:B:902:U:H2'	1:B:903:C:C6	2.54	0.42
1:B:903:C:H2'	1:B:904:C:C6	2.55	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:297:VAL:O	2:A:297:VAL:HG13	2.20	0.42
1:B:943:A:O2'	1:B:944:A:H5'	2.19	0.42
5:B:1192:HOH:O	2:A:468:ILE:HD11	2.20	0.41
1:B:912:C:H2'	1:B:913:C:O4'	2.19	0.41
1:B:966:C:H2'	1:B:967:G:C8	2.55	0.41
2:A:46:ILE:HG23	2:A:93:LYS:HD2	2.02	0.41
2:A:298:SER:O	2:A:302:MET:HG2	2.20	0.41
1:B:906:C:O2'	1:B:907:G:H5'	2.20	0.41
1:B:909:IMG:HM11	1:B:909:IMG:HN21	1.52	0.41
2:A:354:ALA:O	2:A:358:ARG:HG3	2.21	0.41
2:A:243:SER:O	2:A:247:LYS:HE2	2.21	0.41
2:A:205:ARG:NH2	2:A:244:THR:HG23	2.35	0.41
1:B:903:C:O2'	1:B:904:C:H5'	2.21	0.41
2:A:554:LEU:HD12	2:A:554:LEU:HA	1.89	0.41
2:A:299:LYS:N	2:A:299:LYS:CE	2.70	0.40
2:A:10:GLN:HE21	2:A:10:GLN:H	1.69	0.40
2:A:410:THR:C	2:A:412:LYS:H	2.24	0.40
2:A:544:THR:O	2:A:545:HIS:HB2	2.22	0.40
2:A:216:PRO:HG2	2:A:272:PHE:HB3	2.03	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
2	A	604/607 (100%)	576 (95%)	23 (4%)	5 (1%)	27 24

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	237	SER
2	A	259	ASP
2	A	240	LEU
2	A	150	SER

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Mol	Chain	Res	Type
2	A	216	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
2	A	525/526 (100%)	507 (97%)	18 (3%)	49 59

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	10	GLN
2	A	77	ARG
2	A	100	LEU
2	A	128	LEU
2	A	132	GLU
2	A	192	TRP
2	A	253	LYS
2	A	260	GLU
2	A	299	LYS
2	A	349	THR
2	A	351	ASP
2	A	382	PHE
2	A	422	LEU
2	A	509	GLU
2	A	526	LEU
2	A	530	LEU
2	A	554	LEU
2	A	581	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	10	GLN
2	A	62	ASN
2	A	89	GLN

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Mol	Chain	Res	Type
2	A	245	ASN
2	A	406	GLN
2	A	442	GLN
2	A	508	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	B	75/76 (98%)	2 (2%)	0

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	B	909	1MG
1	B	974	C

There are no RNA pucker outliers to report.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

12 non-standard protein/DNA/RNA residues 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
1	PSU	B	901	1	22,22,22	1.94	2 (9%)	29,33,33	2.48	3 (10%)
1	1MG	B	909	1	24,26,27	0.82	0	34,39,42	11.89	6 (17%)
1	2MG	B	910	1	24,26,27	0.86	1 (4%)	32,38,41	5.62	4 (12%)
1	H2U	B	916	1	19,21,22	0.73	1 (5%)	27,30,33	1.04	1 (3%)
1	H2U	B	919	1	19,21,22	0.77	1 (5%)	27,30,33	1.17	1 (3%)
1	M2G	B	926	1	25,27,28	0.86	0	34,40,43	5.05	6 (17%)
1	PSU	B	927	1	19,21,22	2.15	2 (10%)	23,30,33	2.79	3 (13%)
1	H2U	B	947	1	19,21,22	0.84	1 (5%)	27,30,33	1.25	1 (3%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	5MC	B	949	1	20,22,23	1.04	2 (10%)	26,32,35	1.39	2 (7%)
1	5MU	B	954	1	20,22,23	0.85	1 (5%)	25,32,35	1.15	3 (12%)
1	PSU	B	955	1	19,21,22	2.14	2 (10%)	23,30,33	2.71	3 (13%)
1	1MA	B	958	1	23,25,26	0.87	0	32,37,40	0.91	1 (3%)

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
1	PSU	B	901	1	-	0/10/26/26	0/2/2/2
1	1MG	B	909	1	-	0/8/25/26	0/1/3/3
1	2MG	B	910	1	-	0/10/27/28	0/1/3/3
1	H2U	B	916	1	-	0/8/38/39	0/2/2/2
1	H2U	B	919	1	-	0/8/38/39	0/2/2/2
1	M2G	B	926	1	-	0/12/29/30	0/1/3/3
1	PSU	B	927	1	-	0/8/25/26	0/2/2/2
1	H2U	B	947	1	-	0/8/38/39	0/2/2/2
1	5MC	B	949	1	-	0/6/25/26	0/2/2/2
1	5MU	B	954	1	-	0/6/25/26	0/2/2/2
1	PSU	B	955	1	-	0/8/25/26	0/2/2/2
1	1MA	B	958	1	-	0/8/25/26	0/1/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	927	PSU	C6-N1	-8.73	1.24	1.32
1	B	955	PSU	C6-N1	-8.63	1.24	1.32
1	B	901	PSU	C6-N1	-8.46	1.24	1.32
1	B	949	5MC	C5-C4	2.54	1.45	1.41
1	B	947	H2U	C2-N1	2.44	1.39	1.35
1	B	901	PSU	C6-C5	-2.22	1.34	1.38
1	B	955	PSU	C6-C5	-2.21	1.34	1.38
1	B	954	5MU	C6-C5	-2.18	1.34	1.40
1	B	927	PSU	C6-C5	-2.14	1.34	1.38
1	B	919	H2U	P-OP1	2.13	1.49	1.46
1	B	949	5MC	C2-N1	2.07	1.40	1.38
1	B	916	H2U	C2-N1	2.06	1.38	1.35
1	B	910	2MG	CM2-N2	2.03	1.49	1.45

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	909	1MG	C6-C5-N7	-67.32	130.66	134.24
1	B	910	2MG	C6-C5-N7	-29.80	130.13	134.14
1	B	926	M2G	C6-C5-N7	-27.28	130.47	134.14
1	B	909	1MG	C6-N1-C2	14.58	124.71	120.71
1	B	927	PSU	O4'-C1'-C5	12.49	125.09	109.55
1	B	901	PSU	O4'-C1'-C5	12.21	124.74	109.55
1	B	955	PSU	O4'-C1'-C5	12.07	124.56	109.55
1	B	910	2MG	CM2-N2-C2	-9.60	108.19	123.73
1	B	926	M2G	C6-N1-C2	6.93	125.17	120.28
1	B	926	M2G	CM2-N2-CM1	5.84	135.88	115.78
1	B	947	H2U	C5-C6-N1	5.73	117.13	110.71
1	B	949	5MC	C2-N3-C4	5.66	120.54	115.41
1	B	919	H2U	C5-C6-N1	5.25	116.59	110.71
1	B	916	H2U	C5-C6-N1	4.26	115.48	110.71
1	B	909	1MG	C5-C6-N1	-3.62	111.78	119.39
1	B	909	1MG	CM1-N1-C2	-3.53	117.40	119.51
1	B	954	5MU	C6-N1-C2	-3.50	121.41	122.41
1	B	955	PSU	C5-C6-N1	3.35	125.58	120.68
1	B	958	1MA	C2-N3-C4	-3.34	110.52	116.23
1	B	949	5MC	C6-N1-C2	3.25	120.22	118.62
1	B	926	M2G	CM1-N2-C2	-3.22	111.56	120.79
1	B	901	PSU	C5-C6-N1	3.19	125.35	120.68
1	B	927	PSU	C5-C6-N1	3.11	125.22	120.68
1	B	910	2MG	C6-N1-C2	3.10	125.21	119.87
1	B	926	M2G	CM2-N2-C2	-2.88	112.56	120.79
1	B	955	PSU	C5-C4-N3	-2.69	113.96	118.86
1	B	901	PSU	C5-C4-N3	-2.64	114.05	118.86
1	B	927	PSU	C5-C4-N3	-2.57	114.18	118.86
1	B	909	1MG	C2-N3-C4	-2.47	112.02	115.30
1	B	910	2MG	C2-N3-C4	-2.36	111.84	115.03
1	B	909	1MG	N2-C2-N1	-2.10	116.23	119.39
1	B	926	M2G	C2-N3-C4	-2.08	112.19	115.14
1	B	954	5MU	C5M-C5-C6	2.05	122.94	118.59
1	B	954	5MU	C5-C6-N1	2.04	123.57	121.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 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	$\# Z > 2$
4	ARG	A	800	-	11,11,11	0.55	0	13,13,13	0.42	0
3	SO4	B	900	-	4,4,4	0.18	0	6,6,6	0.07	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
4	ARG	A	800	-	-	0/11/11/11	0/0/0/0
3	SO4	B	900	-	-	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.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	B	76/76 (100%)	-0.38	4 (5%) 25 25	7, 17, 46, 75	0
2	A	606/607 (99%)	0.59	76 (12%) 5 4	5, 26, 83, 120	0
All	All	682/683 (99%)	0.48	80 (11%) 5 5	5, 24, 79, 120	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	238	ILE	13.2
2	A	410	THR	12.7
2	A	235	GLY	11.7
2	A	237	SER	9.6
2	A	251	TYR	9.3
2	A	236	ASP	8.8
2	A	409	SER	8.2
2	A	413	GLY	8.2
2	A	411	ARG	7.3
2	A	234	GLU	6.8
2	A	214	LYS	6.8
2	A	239	PRO	6.7
2	A	240	LEU	6.5
2	A	415	VAL	6.3
2	A	233	GLU	6.2
2	A	231	ILE	6.0
2	A	215	ASP	5.8
2	A	211	ALA	5.8
1	B	920	C	5.7
2	A	212	LEU	5.5
2	A	412	LYS	5.4
1	B	973	G	5.4
2	A	414	THR	5.3
2	A	249	ARG	5.3

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Mol	Chain	Res	Type	RSRZ
2	A	213	VAL	5.1
2	A	244	THR	5.0
2	A	241	GLU	4.6
2	A	210	GLU	4.6
2	A	242	GLN	4.5
2	A	243	SER	4.4
1	B	974	C	4.4
2	A	232	GLU	4.1
2	A	20	ALA	4.0
2	A	217	ILE	4.0
2	A	19	PRO	3.8
2	A	216	PRO	3.8
2	A	218	HIS	3.8
2	A	246	GLY	3.7
2	A	2	ALA	3.7
2	A	3	SER	3.6
1	B	901	PSU	3.6
2	A	317	GLU	3.5
2	A	318	ASP	3.4
2	A	254	ARG	3.3
2	A	80	GLY	3.2
2	A	263	LEU	3.1
2	A	408	MET	3.1
2	A	227	ILE	3.0
2	A	256	GLU	2.9
2	A	319	LYS	2.9
2	A	24	ASP	2.9
2	A	253	LYS	2.8
2	A	16	ILE	2.8
2	A	247	LYS	2.8
2	A	316	HIS	2.8
2	A	23	LYS	2.7
2	A	196	PHE	2.7
2	A	21	VAL	2.7
2	A	250	GLU	2.7
2	A	252	PHE	2.7
2	A	297	VAL	2.6
2	A	340	LYS	2.6
2	A	416	VAL	2.6
2	A	272	PHE	2.6
2	A	192	TRP	2.5
2	A	257	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
2	A	79	LYS	2.4
2	A	15	SER	2.4
2	A	204	GLU	2.3
2	A	155	ALA	2.3
2	A	336	ALA	2.3
2	A	141	GLU	2.2
2	A	194	LYS	2.2
2	A	248	ALA	2.2
2	A	392	GLU	2.1
2	A	331	LYS	2.1
2	A	77	ARG	2.1
2	A	385	ILE	2.0
2	A	245	ASN	2.0
2	A	201	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	PSU	B	901	21/21	0.31	7.29	56,58,68,69	0
1	PSU	B	955	20/21	0.12	0.59	15,19,22,24	0
1	1MG	B	909	24/25	0.09	0.53	8,12,17,21	0
1	H2U	B	947	20/21	0.22	0.35	34,44,57,57	0
1	2MG	B	910	24/25	0.08	0.12	4,11,17,18	0
1	1MA	B	958	23/24	0.09	0.02	8,11,15,15	0
1	M2G	B	926	25/26	0.10	-0.24	10,15,20,22	0
1	5MU	B	954	21/22	0.09	-0.38	8,14,17,21	0
1	H2U	B	919	20/21	0.12	-0.49	22,25,30,38	0
1	H2U	B	916	20/21	0.09	-0.94	11,14,18,19	0
1	5MC	B	949	21/22	0.08	-1.03	3,12,14,14	0
1	PSU	B	927	20/21	0.11	-1.72	14,21,25,30	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	B	900	5/5	0.13	0.77	30,33,35,36	0
4	ARG	A	800	12/12	0.14	0.75	23,25,30,32	0

6.5 Other polymers

There are no such residues in this entry.