



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:49 PM GMT

PDB ID : 1QRT
Title : GLUTAMINYL-TRNA SYNTHETASE MUTANT D235G COMPLEXED
WITH GLUTAMINE TRANSFER RNA
Authors : Arnez, J.G.; Steitz, T.A.
Deposited on : 1996-06-14
Resolution : 2.70 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

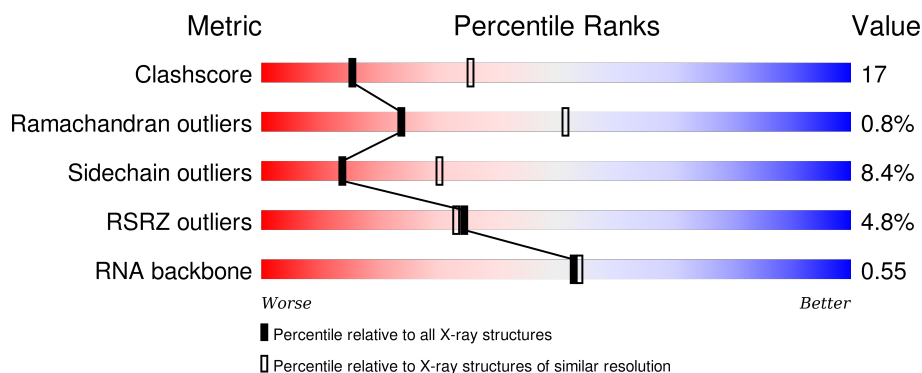
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)
RNA backbone	2183	1069 (3.10-2.30)

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	B	75	
2	A	553	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7158 atoms, of which 1207 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 RNA chain called TRNAGLN2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	74	Total	C	H	N	O	P	0	0	0
			1802	702	229	279	518	74			

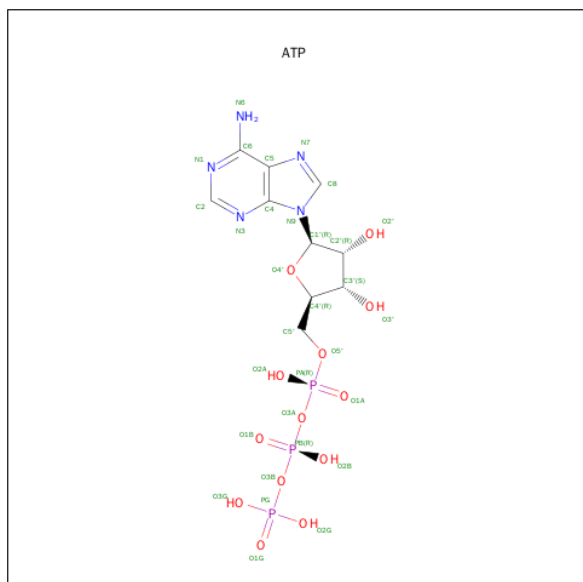
- Molecule 2 is a protein called PROTEIN (GLUTAMINYL-TRNA SYNTHETASE (E.C.6.1.1.18)).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	A	529	Total	C	H	N	O	S	0	0	0
			5250	2702	975	752	800	21			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	235	GLY	ASP	ENGINEERED	UNP P00962

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	P	0	0
			34	10	3	5	13	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	62	Total	O	0	0
			62	62		
4	B	10	Total	O	0	0
			10	10		

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	243.50 Å 93.70 Å 115.90 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.50 – 2.70 8.07 – 2.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.50-2.70) 88.7 (8.07-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.07 (at 2.59 Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.216 , (Not available) 0.216 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 66.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 35000 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7158	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: ATP

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	0.50	0/1756	0.84	5/2734 (0.2%)
2	A	0.47	0/4375	0.69	1/5922 (0.0%)
All	All	0.48	0/6131	0.74	6/8656 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	18	G	N9-C1'-C2'	5.84	121.60	114.00
1	B	74	C	N1-C1'-C2'	5.43	121.06	114.00
2	A	49	LEU	CA-CB-CG	5.38	127.68	115.30
1	B	73	G	N9-C1'-C2'	5.09	120.62	114.00
1	B	19	G	N9-C1'-C2'	5.06	120.58	114.00
1	B	48	C	N1-C1'-C2'	5.03	120.54	114.00

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	B	1573	229	800	69	0
2	A	4275	975	4171	131	1
3	A	31	3	12	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	62	0	0	4	0
4	B	10	0	0	0	0
All	All	5951	1207	4983	186	1

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 17.

All (186) 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:46:U:H4'	1:B:47:U:C6	1.72	1.25
1:B:58:A:H4'	1:B:59:A:OP1	1.51	1.09
1:B:7:A:H5'	2:A:318:GLN:HE22	1.23	1.00
1:B:36:G:H4'	1:B:37:A:OP2	1.62	0.99
1:B:46:U:H4'	1:B:47:U:C5	2.01	0.95
1:B:16:C:OP2	1:B:16:C:H3'	1.67	0.94
1:B:34:C:H4'	1:B:35:U:OP1	1.74	0.88
2:A:522:GLY:HA2	2:A:544:LEU:HD13	1.53	0.88
1:B:42:G:O2'	1:B:43:G:H5'	1.78	0.84
1:B:47:U:H2'	1:B:47:U:O2	1.77	0.83
2:A:39:LEU:HD13	2:A:81:ILE:HG12	1.59	0.82
1:B:7:A:C5'	2:A:318:GLN:HE22	1.92	0.80
1:B:40:C:O2'	1:B:41:C:H5'	1.83	0.79
1:B:41:C:H2'	1:B:42:G:C8	2.18	0.79
2:A:44:ALA:HB2	2:A:293:ILE:HD11	1.64	0.77
1:B:38:U:H4'	2:A:336:ASN:ND2	2.01	0.75
1:B:21:A:H62	1:B:47:U:H1'	1.51	0.74
2:A:407:LYS:HG2	2:A:408:GLU:H	1.54	0.73
2:A:46:SER:OG	3:A:999:ATP:H1'	1.89	0.72
2:A:471:LEU:HB2	2:A:497:VAL:HG13	1.71	0.72
1:B:46:U:C4'	1:B:47:U:C5	2.72	0.72
2:A:423:GLU:HB2	2:A:434:PHE:HE2	1.56	0.71
1:B:58:A:C4'	1:B:59:A:OP1	2.36	0.70
1:B:40:C:H2'	1:B:41:C:C6	2.27	0.69
1:B:46:U:H4'	1:B:47:U:H6	1.49	0.69
1:B:41:C:H2'	1:B:42:G:H8	1.58	0.69
2:A:30:ARG:NH1	2:A:215:HIS:CE1	2.60	0.68
1:B:40:C:H2'	1:B:41:C:H6	1.58	0.68
2:A:423:GLU:HB3	2:A:432:THR:HG23	1.76	0.66
2:A:401:LYS:O	2:A:454:GLY:HA2	1.97	0.64
2:A:30:ARG:NH1	2:A:215:HIS:HE1	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:545:ARG:HG2	2:A:547:THR:HG23	1.80	0.64
2:A:534:GLU:O	2:A:536:PRO:HD3	1.97	0.63
2:A:166:GLU:O	2:A:169:LYS:HB2	1.99	0.63
2:A:441:THR:O	2:A:442:LEU:HB3	1.99	0.62
1:B:7:A:H5'	2:A:318:GLN:NE2	2.06	0.61
1:B:45:A:O2'	1:B:47:U:H1'	2.01	0.61
2:A:324:MET:O	2:A:328:GLU:HG3	2.00	0.61
2:A:252:HIS:HD2	2:A:253:PRO:O	1.85	0.60
1:B:34:C:C4'	1:B:35:U:OP1	2.48	0.60
1:B:74:C:N3	2:A:168:GLY:HA2	2.17	0.60
2:A:349:VAL:HB	2:A:389:ARG:HD3	1.83	0.60
1:B:38:U:H4'	2:A:336:ASN:HD21	1.67	0.59
1:B:60:U:OP2	1:B:60:U:H6	1.85	0.59
1:B:35:U:C5	1:B:37:A:C2	2.91	0.59
2:A:363:VAL:HG23	2:A:381:PHE:HB2	1.83	0.59
1:B:15:G:H3'	1:B:16:C:C5	2.38	0.59
2:A:365:MET:SD	2:A:379:VAL:HG21	2.43	0.58
1:B:50:G:O2'	1:B:51:A:H5'	2.03	0.58
2:A:263:LEU:HB3	2:A:266:THR:CG2	2.33	0.58
1:B:36:G:C4'	1:B:37:A:OP2	2.46	0.58
2:A:98:TYR:O	2:A:101:ASP:HB2	2.04	0.58
2:A:511:VAL:HG23	2:A:514:LYS:HB2	1.85	0.57
2:A:331:ILE:HG13	2:A:335:LEU:HD22	1.87	0.57
2:A:114:ILE:HG21	2:A:154:LEU:HD13	1.85	0.57
2:A:512:ALA:HB1	2:A:528:SER:HB2	1.87	0.56
2:A:403:LEU:HD13	2:A:409:VAL:HG22	1.87	0.56
1:B:60:U:H5''	1:B:61:C:H5	1.71	0.56
1:B:7:A:C5'	2:A:318:GLN:NE2	2.65	0.56
1:B:16:C:H4'	1:B:18:G:OP2	2.04	0.56
2:A:60:GLN:HA	2:A:60:GLN:HE21	1.70	0.55
1:B:47:U:C2'	1:B:47:U:O2	2.50	0.55
1:B:60:U:H5''	1:B:61:C:C5	2.42	0.55
2:A:277:VAL:HG13	2:A:284:GLY:HA2	1.88	0.55
2:A:397:ASN:HD21	2:A:399:GLN:HG2	1.72	0.55
2:A:408:GLU:HB3	2:A:417:ILE:O	2.08	0.54
2:A:393:ARG:O	2:A:404:VAL:HA	2.07	0.54
1:B:36:G:C6	2:A:401:LYS:HG2	2.42	0.54
2:A:464:ALA:HB1	2:A:502:PHE:HB3	1.90	0.54
2:A:263:LEU:HB3	2:A:266:THR:HG23	1.89	0.54
1:B:9:C:H5'	1:B:10:G:OP2	2.09	0.53
2:A:71:VAL:HG23	2:A:72:LYS:HG2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:G:O2'	1:B:19:G:P	2.66	0.53
2:A:288:PRO:HG3	2:A:298:ARG:HG2	1.90	0.53
2:A:517:GLN:HG3	2:A:523:TYR:CE1	2.43	0.53
2:A:298:ARG:NH1	2:A:529:ARG:HD2	2.23	0.53
2:A:418:LYS:O	2:A:435:CYS:HB3	2.09	0.52
2:A:173:ARG:HD2	2:A:187:ASP:O	2.08	0.52
2:A:40:HIS:HA	2:A:292:THR:HA	1.91	0.52
2:A:346:ILE:HG12	2:A:469:ILE:CD1	2.38	0.52
1:B:34:C:O2'	2:A:412:ARG:NH2	2.41	0.52
2:A:335:LEU:HB3	2:A:523:TYR:CD1	2.43	0.52
2:A:63:LEU:HD23	2:A:96:VAL:HG22	1.91	0.52
2:A:363:VAL:HG11	2:A:413:ASN:O	2.10	0.52
2:A:32:PRO:HA	2:A:64:ARG:O	2.10	0.52
2:A:545:ARG:HG3	2:A:545:ARG:HH11	1.75	0.51
1:B:18:G:HO2'	1:B:19:G:P	2.33	0.51
1:B:36:G:C4	2:A:401:LYS:HD3	2.46	0.51
1:B:43:G:H2'	1:B:44:C:O4'	2.12	0.51
2:A:30:ARG:HH12	2:A:215:HIS:CE1	2.28	0.50
1:B:45:A:O2'	1:B:47:U:C1'	2.58	0.50
1:B:15:G:H3'	1:B:16:C:H5	1.76	0.50
2:A:35:PRO:HG2	2:A:73:GLU:HB3	1.94	0.50
2:A:368:HIS:HD2	2:A:370:ASN:H	1.58	0.50
2:A:545:ARG:NH1	2:A:545:ARG:HG3	2.26	0.50
1:B:46:U:C4'	1:B:47:U:C6	2.67	0.49
2:A:124:LEU:HD22	2:A:143:SER:HB2	1.93	0.49
2:A:39:LEU:CD1	2:A:81:ILE:HG12	2.38	0.49
2:A:232:GLU:H	2:A:232:GLU:CD	2.15	0.49
2:A:526:LEU:HD12	2:A:537:VAL:O	2.13	0.49
2:A:391:ASP:OD1	2:A:402:ARG:HD2	2.13	0.49
1:B:15:G:O2'	1:B:16:C:OP1	2.30	0.49
2:A:363:VAL:CG2	2:A:381:PHE:HB2	2.42	0.49
1:B:15:G:HO2'	1:B:16:C:P	2.36	0.49
2:A:467:VAL:HG23	2:A:536:PRO:HB2	1.95	0.49
2:A:394:GLU:HA	2:A:405:LEU:HB2	1.95	0.49
2:A:121:VAL:H	2:A:153:ASN:ND2	2.11	0.48
2:A:487:PHE:O	2:A:490:VAL:HG12	2.13	0.48
1:B:61:C:O2'	1:B:62:C:H5'	2.12	0.48
1:B:15:G:P	1:B:16:C:H41	2.35	0.48
2:A:410:ARG:HH21	2:A:442:LEU:HA	1.78	0.48
2:A:345:VAL:HA	4:A:732:HOH:O	2.13	0.48
2:A:322:ILE:N	2:A:322:ILE:HD12	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:C:HO2'	1:B:41:C:H5'	1.74	0.48
2:A:178:MET:O	2:A:186:ARG:HD2	2.15	0.47
2:A:467:VAL:HG21	2:A:538:PHE:CE1	2.50	0.47
2:A:433:ILE:HG22	2:A:435:CYS:SG	2.55	0.47
1:B:21:A:N6	1:B:47:U:O2'	2.48	0.47
2:A:463:HIS:O	2:A:505:PRO:HD3	2.15	0.47
1:B:18:G:O2'	1:B:57:G:N2	2.48	0.46
2:A:160:MET:HG2	2:A:193:ILE:HD11	1.97	0.46
2:A:301:TYR:HE2	2:A:327:LEU:HD22	1.81	0.46
2:A:388:ASP:HB3	2:A:391:ASP:OD2	2.16	0.46
1:B:9:C:H4'	1:B:10:G:OP2	2.16	0.46
2:A:425:ASP:HB3	2:A:431:THR:HG23	1.96	0.46
2:A:402:ARG:HG2	2:A:456:ILE:HG22	1.97	0.46
2:A:42:GLY:HA2	2:A:261:LEU:HD23	1.98	0.46
2:A:264:GLU:O	2:A:265:TYR:HB2	2.16	0.46
2:A:391:ASP:OD1	2:A:520:ARG:NH2	2.49	0.46
1:B:36:G:N3	2:A:401:LYS:HD3	2.31	0.45
1:B:9:C:C4'	1:B:10:G:OP2	2.64	0.45
1:B:58:A:H4'	1:B:60:U:H5	1.81	0.45
1:B:42:G:C2'	1:B:43:G:H5'	2.47	0.45
2:A:64:ARG:HD3	2:A:222:GLU:OE2	2.17	0.45
2:A:423:GLU:HB2	2:A:434:PHE:CE2	2.45	0.45
2:A:261:LEU:HD12	2:A:322:ILE:HD13	1.99	0.44
2:A:343:MET:HE3	2:A:458:TRP:H	1.82	0.44
1:B:8:U:H2'	1:B:13:A:H61	1.82	0.44
1:B:51:A:H2'	1:B:52:G:O4'	2.16	0.44
2:A:132:TYR:CD2	2:A:141:LYS:HG3	2.53	0.44
2:A:264:GLU:HG2	2:A:265:TYR:CD2	2.51	0.44
2:A:403:LEU:HA	2:A:409:VAL:HG11	1.99	0.44
2:A:231:LEU:HB2	4:A:768:HOH:O	2.16	0.44
2:A:30:ARG:HH11	2:A:215:HIS:CE1	2.35	0.44
2:A:506:SER:C	2:A:508:LYS:H	2.20	0.44
2:A:212:ASP:HB2	4:A:718:HOH:O	2.18	0.44
2:A:30:ARG:HH12	2:A:215:HIS:HE1	1.64	0.44
2:A:346:ILE:HG12	2:A:469:ILE:HD12	2.00	0.44
1:B:38:U:C4	1:B:39:U:C4	3.06	0.43
2:A:287:ASP:OD1	2:A:289:ARG:HD3	2.18	0.43
2:A:335:LEU:HA	2:A:335:LEU:HD12	1.82	0.43
2:A:113:LEU:HD12	2:A:113:LEU:HA	1.84	0.43
1:B:72:A:C4'	2:A:183:ILE:HG21	2.48	0.43
2:A:46:SER:HG	3:A:999:ATP:H1'	1.80	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:402:ARG:HD3	2:A:520:ARG:HH22	1.84	0.43
2:A:163:GLY:HA2	4:A:744:HOH:O	2.17	0.43
1:B:34:C:O2'	2:A:412:ARG:CZ	2.67	0.43
2:A:507:LEU:HA	2:A:507:LEU:HD13	1.81	0.43
2:A:345:VAL:CG1	2:A:348:PRO:HB3	2.48	0.43
2:A:75:ILE:O	2:A:78:VAL:HG22	2.19	0.43
1:B:32:U:C4	1:B:33:U:C4	3.07	0.42
2:A:51:PHE:O	2:A:55:GLN:HG2	2.18	0.42
1:B:8:U:OP2	2:A:318:GLN:NE2	2.53	0.42
2:A:349:VAL:HB	2:A:389:ARG:CD	2.49	0.42
2:A:505:PRO:O	2:A:508:LYS:HB2	2.19	0.42
2:A:415:TYR:CE1	2:A:442:LEU:HD22	2.55	0.42
2:A:124:LEU:HA	2:A:124:LEU:HD12	1.85	0.42
2:A:174:ALA:O	2:A:187:ASP:HA	2.20	0.41
2:A:13:GLN:O	2:A:17:GLU:HG3	2.20	0.41
1:B:44:C:H5'	1:B:44:C:H6	1.85	0.41
1:B:36:G:H5'	1:B:37:A:H8	1.84	0.41
1:B:38:U:H2'	1:B:39:U:H6	1.86	0.41
2:A:263:LEU:HD23	2:A:263:LEU:HA	1.86	0.41
2:A:158:GLU:OE1	2:A:161:ARG:NH1	2.54	0.41
2:A:377:ARG:HG2	2:A:377:ARG:HH11	1.85	0.41
1:B:9:C:C5'	1:B:10:G:OP2	2.68	0.41
2:A:286:ASP:O	2:A:298:ARG:HD3	2.21	0.41
2:A:506:SER:C	2:A:508:LYS:N	2.74	0.41
2:A:381:PHE:CG	2:A:382:SER:N	2.89	0.41
2:A:207:ILE:O	2:A:207:ILE:HG13	2.21	0.41
2:A:78:VAL:HG23	2:A:79:GLU:N	2.36	0.41
2:A:252:HIS:NE2	2:A:254:ARG:NH1	2.69	0.40
2:A:347:ASP:O	2:A:388:ASP:HA	2.21	0.40
2:A:289:ARG:HD2	2:A:472:TYR:HD1	1.86	0.40
1:B:19:G:C4	1:B:57:G:N2	2.90	0.40
2:A:19:LEU:HD12	2:A:19:LEU:HA	1.93	0.40

All (1) 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)
2:A:105:GLN:HE22	2:A:105:GLN:HE22[4_565]	1.26	0.34

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	
2	A	525/553 (95%)	485 (92%)	36 (7%)	4 (1%)	24	51

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	32	PRO
2	A	439	ALA
2	A	348	PRO
2	A	176	ILE

5.3.2 Protein sidechains [i](#)

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	462/480 (96%)	423 (92%)	39 (8%)	14	30

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	19	LEU
2	A	30	ARG
2	A	32	PRO
2	A	33	PRO
2	A	49	LEU
2	A	60	GLN
2	A	72	LYS

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Mol	Chain	Res	Type
2	A	95	ASN
2	A	106	LEU
2	A	113	LEU
2	A	124	LEU
2	A	154	LEU
2	A	156	LEU
2	A	186	ARG
2	A	190	LEU
2	A	192	ARG
2	A	231	LEU
2	A	263	LEU
2	A	266	THR
2	A	268	MET
2	A	271	ARG
2	A	318	GLN
2	A	326	SER
2	A	327	LEU
2	A	335	LEU
2	A	341	ARG
2	A	362	MET
2	A	395	GLU
2	A	410	ARG
2	A	412	ARG
2	A	422	VAL
2	A	425	ASP
2	A	432	THR
2	A	474	ARG
2	A	475	LEU
2	A	490	VAL
2	A	497	VAL
2	A	514	LYS
2	A	544	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	55	GLN
2	A	60	GLN
2	A	95	ASN
2	A	142	ASN
2	A	153	ASN
2	A	215	HIS

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Mol	Chain	Res	Type
2	A	274	ASN
2	A	318	GLN
2	A	336	ASN
2	A	368	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	B	73/75 (97%)	25 (34%)	16 (21%)

All (25) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	B	7	A
1	B	8	U
1	B	9	C
1	B	10	G
1	B	16	C
1	B	18	G
1	B	19	G
1	B	20	U
1	B	21	A
1	B	29	G
1	B	30	G
1	B	35	U
1	B	36	G
1	B	37	A
1	B	45	A
1	B	46	U
1	B	49	C
1	B	59	A
1	B	60	U
1	B	61	C
1	B	63	U
1	B	64	C
1	B	73	G
1	B	74	C
1	B	75	C

All (16) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	B	7	A
1	B	9	C
1	B	15	G
1	B	16	C
1	B	18	G
1	B	19	G
1	B	20	U
1	B	34	C
1	B	35	U
1	B	36	G
1	B	45	A
1	B	48	C
1	B	58	A
1	B	60	U
1	B	73	G
1	B	74	C

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](#)

1 ligand is 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	$\# Z > 2$
3	ATP	A	999	-	24,33,33	0.73	0	31,52,52	0.79	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
3	ATP	A	999	-	-	0/18/38/38	0/3/3/3

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 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	999	ATP	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	B	74/75 (98%)	0.61	6 (8%) 15 12	9, 52, 89, 105	0
2	A	529/553 (95%)	-0.28	23 (4%) 39 38	2, 19, 68, 81	0
All	All	603/628 (96%)	-0.17	29 (4%) 34 33	2, 21, 73, 105	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	46	U	6.4
1	B	47	U	6.2
2	A	440	ASP	4.8
2	A	533	ALA	4.4
2	A	355	ASN	3.8
1	B	45	A	3.8
2	A	393	ARG	3.7
2	A	425	ASP	3.6
1	B	20	U	3.6
2	A	534	GLU	3.5
2	A	441	THR	3.4
2	A	407	LYS	3.4
2	A	395	GLU	3.3
2	A	426	ALA	3.2
2	A	532	THR	3.2
2	A	429	ASN	3.1
2	A	454	GLY	3.0
1	B	30	G	2.8
2	A	398	LYS	2.5
2	A	409	VAL	2.5
2	A	442	LEU	2.4
1	B	44	C	2.4
2	A	404	VAL	2.3
2	A	397	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
2	A	463	HIS	2.2
2	A	399	GLN	2.2
2	A	127	GLU	2.1
2	A	427	GLU	2.0
2	A	357	GLN	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
3	ATP	A	999	31/31	0.93	0.16	0.57	0,28,59,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.