Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 02:51 PM GMT

PDB ID	:	4ARI
Title	:	Ternary complex of E. coli leucyl-tRNA synthetase, tRNA(leu) and the benzoxaborole AN2679 in the editing conformation
Authors	:	Palencia, A.; Crepin, T.; Vu, M.T.; Lincecum Jr, T.L.; Martinis, S.A.; Cusack, S.
Deposited on	:	2012-04-24
Resolution	:	2.08 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

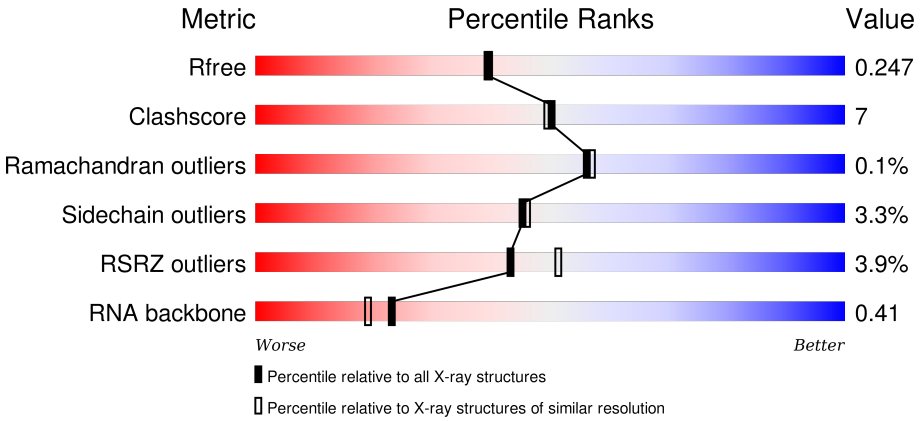
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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	4546 (2.10-2.06)
Clashscore	102246	5101 (2.10-2.06)
Ramachandran outliers	100387	5048 (2.10-2.06)
Sidechain outliers	100360	5049 (2.10-2.06)
RSRZ outliers	91569	4556 (2.10-2.06)
RNA backbone	2183	1003 (2.76-1.40)

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. 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 <=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	880	2% 81% 12% • 7%
2	B	87	16% 51% 29% 13% 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	1861	-	-	-	X
3	GOL	A	1862	-	-	-	X
3	GOL	A	1863	-	-	X	-
3	GOL	A	1864	-	-	-	X

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 8596 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 LEUCINE–TRNA LIGASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	821	Total	C	N	O	S	0	1	0
			6526	4153	1104	1229	40			

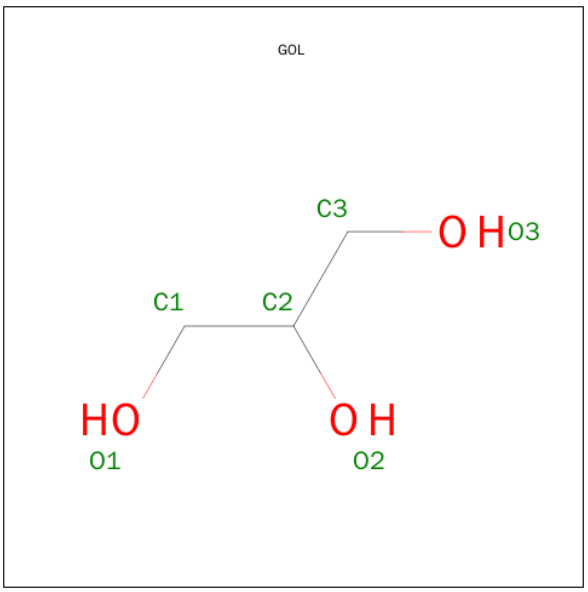
There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP P07813
A	-18	GLY	-	EXPRESSION TAG	UNP P07813
A	-17	SER	-	EXPRESSION TAG	UNP P07813
A	-16	SER	-	EXPRESSION TAG	UNP P07813
A	-15	HIS	-	EXPRESSION TAG	UNP P07813
A	-14	HIS	-	EXPRESSION TAG	UNP P07813
A	-13	HIS	-	EXPRESSION TAG	UNP P07813
A	-12	HIS	-	EXPRESSION TAG	UNP P07813
A	-11	HIS	-	EXPRESSION TAG	UNP P07813
A	-10	HIS	-	EXPRESSION TAG	UNP P07813
A	-9	SER	-	EXPRESSION TAG	UNP P07813
A	-8	SER	-	EXPRESSION TAG	UNP P07813
A	-7	GLY	-	EXPRESSION TAG	UNP P07813
A	-6	LEU	-	EXPRESSION TAG	UNP P07813
A	-5	VAL	-	EXPRESSION TAG	UNP P07813
A	-4	PRO	-	EXPRESSION TAG	UNP P07813
A	-3	ARG	-	EXPRESSION TAG	UNP P07813
A	-2	GLY	-	EXPRESSION TAG	UNP P07813
A	-1	SER	-	EXPRESSION TAG	UNP P07813

- Molecule 2 is a RNA chain called TRNA-LEU5 (UAA ISOACEPTOR).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	80	Total	B	C	N	O	P	0	0	0
			1720	1	768	307	564	80			

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Mg 1 1	0	0
4	A	1	Total Mg 1 1	0	0

- Molecule 5 is water.

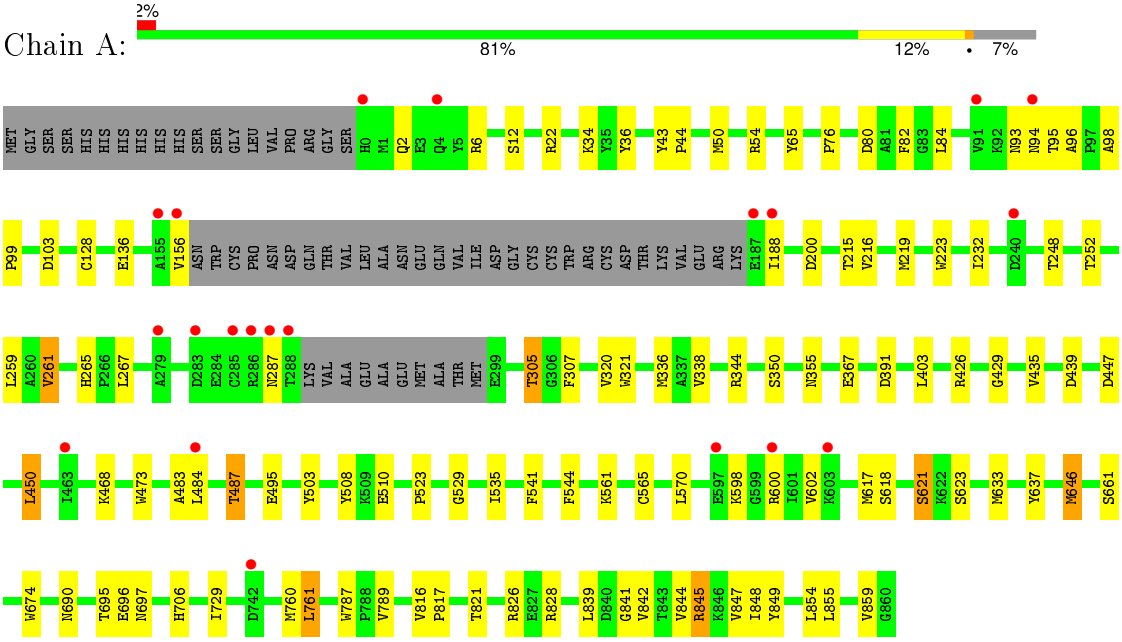
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	301	Total O 301 301	0	0
5	B	23	Total O 23 23	0	0

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.

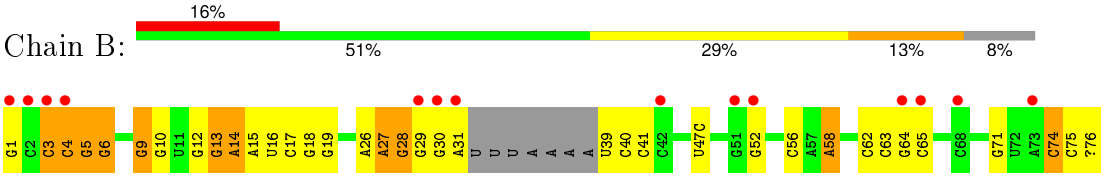
- Molecule 1: LEUCINE–TRNA LIGASE

Chain A:



- Molecule 2: TRNA-LEU5 (UAA ISOACEPTOR)

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	76.18Å 118.94Å 141.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	90.91 – 2.08 45.46 – 2.08	Depositor EDS
% Data completeness (in resolution range)	96.7 (90.91-2.08) 96.8 (45.46-2.08)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.74 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.6.0116	Depositor
R, R_{free}	0.199 , 0.244 0.202 , 0.247	Depositor DCC
R_{free} test set	3759 reflections (5.28%)	DCC
Wilson B-factor (Å^2)	28.5	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 47.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Outliers	0 of 74893 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8596	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.68% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, MG, N79

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	0.68	3/6683 (0.0%)	0.73	4/9069 (0.0%)
2	B	0.37	0/1885	0.85	2/2934 (0.1%)
All	All	0.62	3/8568 (0.0%)	0.76	6/12003 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	674	TRP	CD2-CE2	5.53	1.48	1.41
1	A	321	TRP	CD2-CE2	5.24	1.47	1.41
1	A	473	TRP	CD2-CE2	5.04	1.47	1.41

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	344	ARG	NE-CZ-NH2	-6.43	117.08	120.30
1	A	80	ASP	CB-CG-OD1	5.93	123.64	118.30
1	A	103	ASP	CB-CG-OD1	5.62	123.36	118.30
2	B	15	A	P-O3'-C3'	-5.30	113.34	119.70
2	B	12	G	P-O3'-C3'	5.19	125.93	119.70
1	A	344	ARG	NE-CZ-NH1	5.13	122.87	120.30

There are no chirality outliers.

There are no planarity outliers.

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	6526	0	6389	93	0
2	B	1720	0	871	17	0
3	A	24	0	32	10	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	301	0	0	14	0
5	B	23	0	0	0	0
All	All	8596	0	7292	109	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 7.

All (109) 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:A:355:ASN:HB2	5:A:2135:HOH:O	1.52	1.07
1:A:305:THR:HG23	1:A:307:PHE:H	1.35	0.89
1:A:633:MET:HE1	1:A:637:TYR:HE2	1.36	0.88
1:A:695:THR:CG2	5:A:2243:HOH:O	2.27	0.83
1:A:215:THR:HG22	1:A:219:MET:CE	2.12	0.79
1:A:729:ILE:HD13	1:A:761:LEU:HD13	1.65	0.77
1:A:598:LYS:HE3	5:A:2213:HOH:O	1.87	0.75
1:A:215:THR:O	1:A:219:MET:HE2	1.90	0.72
1:A:633:MET:HE1	1:A:637:TYR:CE2	2.23	0.72
1:A:468:LYS:NZ	1:A:487:THR:CG2	2.54	0.70
1:A:468:LYS:HZ3	1:A:487:THR:CG2	2.05	0.69
1:A:215:THR:HG22	1:A:219:MET:HE2	1.76	0.67
1:A:468:LYS:NZ	1:A:487:THR:HG23	2.10	0.66
1:A:468:LYS:HZ3	1:A:487:THR:HG23	1.61	0.66
1:A:215:THR:HG22	1:A:219:MET:HE1	1.75	0.66
1:A:633:MET:CE	1:A:637:TYR:HE2	2.07	0.65
2:B:27:A:H2'	2:B:28:G:O4'	1.97	0.65
1:A:729:ILE:HD13	1:A:761:LEU:CD1	2.27	0.65
1:A:94:ASN:HA	5:A:2058:HOH:O	1.97	0.64
5:A:2224:HOH:O	2:B:13:G:OP1	2.15	0.64
1:A:842:VAL:HG12	1:A:859:VAL:HB	1.81	0.63
1:A:261:VAL:HG22	1:A:265:HIS:CG	2.34	0.62
1:A:2:GLN:OE1	1:A:6:ARG:NH2	2.33	0.61
1:A:695:THR:HG22	5:A:2243:HOH:O	1.94	0.61
1:A:223:TRP:CE3	1:A:535:ILE:HD12	2.35	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:355:ASN:CB	5:A:2135:HOH:O	2.26	0.60
1:A:50:MET:HE1	1:A:646:MET:SD	2.40	0.60
1:A:729:ILE:HG21	1:A:761:LEU:HD13	1.82	0.60
1:A:816:VAL:HG22	1:A:817:PRO:HD2	1.82	0.60
1:A:633:MET:CE	1:A:637:TYR:CE2	2.85	0.59
1:A:729:ILE:HG21	1:A:761:LEU:CD1	2.32	0.59
1:A:841:GLY:O	1:A:842:VAL:HG23	2.02	0.59
1:A:849:TYR:HH	3:A:1862:GOL:HO3	1.32	0.59
1:A:391:ASP:HB3	5:A:2150:HOH:O	2.02	0.59
2:B:27:A:H2'	2:B:28:G:C1'	2.32	0.59
1:A:93:ASN:HB2	1:A:95:THR:HG22	1.85	0.58
1:A:367:GLU:OE2	5:A:2140:HOH:O	2.17	0.58
1:A:789:VAL:H	3:A:1863:GOL:H12	1.70	0.57
1:A:848:ILE:HG13	2:B:56:C:C2	2.39	0.57
2:B:39:U:H2'	2:B:40:C:C6	2.40	0.57
1:A:844:VAL:HA	1:A:859:VAL:HG12	1.87	0.56
1:A:336:MET:HG3	1:A:336:MET:O	2.06	0.55
1:A:447:ASP:HA	1:A:450:LEU:HD22	1.89	0.54
1:A:136:GLU:OE2	1:A:495:GLU:OE2	2.26	0.54
1:A:261:VAL:CG2	1:A:265:HIS:CG	2.91	0.54
1:A:50:MET:CE	1:A:646:MET:SD	2.96	0.53
1:A:523:PRO:HG2	1:A:561:LYS:HZ2	1.73	0.53
1:A:845:ARG:HG3	5:A:2301:HOH:O	2.08	0.52
1:A:847:VAL:HB	5:A:2294:HOH:O	2.08	0.52
1:A:305:THR:HG22	1:A:320:VAL:O	2.09	0.52
2:B:3:C:H2'	2:B:4:C:O4'	2.09	0.52
1:A:695:THR:HG22	1:A:697:ASN:H	1.74	0.52
1:A:695:THR:HG22	1:A:696:GLU:N	2.25	0.52
1:A:305:THR:CG2	1:A:307:PHE:H	2.16	0.52
1:A:435:VAL:CG1	1:A:483:ALA:HB1	2.40	0.52
1:A:529:GLY:O	1:A:565:CYS:HA	2.11	0.51
1:A:621:SER:HG	1:A:623:SER:HG	1.59	0.51
1:A:248:THR:HG23	2:B:76:N79:H5'1	1.94	0.49
1:A:600:ARG:HB2	5:A:2213:HOH:O	2.13	0.49
2:B:74:C:H4'	2:B:75:C:OP2	2.13	0.49
2:B:5:G:H3'	2:B:6:G:H5"	1.94	0.49
1:A:84:LEU:HD21	1:A:426:ARG:NH1	2.28	0.48
1:A:706:HIS:CE1	3:A:1864:GOL:O2	2.68	0.47
1:A:82:PHE:CE2	1:A:128:CYS:HA	2.50	0.47
1:A:849:TYR:OH	3:A:1862:GOL:O3	2.10	0.47
1:A:259:LEU:CD2	1:A:336:MET:HA	2.45	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:847:VAL:HG11	1:A:855:LEU:HD11	1.95	0.47
1:A:95:THR:OG1	1:A:96:ALA:N	2.46	0.47
2:B:13:G:H2'	2:B:14:A:H5"	1.97	0.46
3:A:1863:GOL:H11	3:A:1864:GOL:C3	2.46	0.46
1:A:259:LEU:HD22	1:A:336:MET:HA	1.98	0.46
1:A:468:LYS:HZ2	1:A:487:THR:CG2	2.28	0.45
1:A:261:VAL:HG22	1:A:265:HIS:CD2	2.52	0.45
2:B:52:G:N2	2:B:63:C:C2	2.85	0.45
1:A:65:TYR:OH	5:A:2043:HOH:O	2.14	0.44
1:A:261:VAL:HG22	1:A:265:HIS:CB	2.47	0.44
1:A:54:ARG:HG3	1:A:646:MET:HE2	1.99	0.44
1:A:789:VAL:H	3:A:1863:GOL:C1	2.31	0.44
1:A:200:ASP:N	3:A:1861:GOL:O2	2.41	0.44
1:A:468:LYS:HZ3	1:A:487:THR:HG22	1.80	0.44
1:A:43:TYR:CD1	1:A:44:PRO:HD2	2.53	0.44
1:A:789:VAL:N	3:A:1863:GOL:H12	2.33	0.43
2:B:63:C:H2'	2:B:64:G:O4'	2.18	0.43
2:B:5:G:H5"	2:B:5:G:C8	2.53	0.43
1:A:839:LEU:HD12	5:A:2298:HOH:O	2.18	0.43
1:A:600:ARG:HH11	1:A:602:VAL:HG12	1.83	0.43
1:A:646:MET:HE2	1:A:646:MET:HB3	1.84	0.43
1:A:706:HIS:HE1	3:A:1864:GOL:O2	2.02	0.43
1:A:570:LEU:HB3	1:A:617[A]:MET:HG3	1.99	0.43
2:B:1:G:H2'	2:B:1:G:N3	2.34	0.43
1:A:821:THR:HA	3:A:1862:GOL:H12	2.00	0.42
2:B:10:G:C6	2:B:26:A:C2	3.08	0.42
1:A:695:THR:CG2	1:A:696:GLU:N	2.83	0.42
1:A:261:VAL:CG2	1:A:265:HIS:CD2	3.03	0.42
1:A:76:PRO:CG	1:A:503:TYR:CD2	3.03	0.41
1:A:570:LEU:HB2	1:A:617[B]:MET:HE2	2.03	0.41
2:B:9:G:N3	2:B:9:G:H2'	2.36	0.41
1:A:439:ASP:OD1	1:A:439:ASP:C	2.58	0.41
1:A:508:TYR:CZ	1:A:510:GLU:HB2	2.56	0.41
2:B:18:G:H21	2:B:58:A:H5'	1.85	0.41
1:A:760:MET:HA	1:A:787:TRP:CH2	2.56	0.41
1:A:216:VAL:HA	1:A:219:MET:HE3	2.02	0.41
1:A:468:LYS:NZ	1:A:487:THR:HG22	2.35	0.41
1:A:82:PHE:CG	1:A:429:GLY:HA2	2.56	0.41
1:A:267:LEU:HD23	1:A:267:LEU:C	2.41	0.41
1:A:98:ALA:HB3	1:A:99:PRO:CD	2.51	0.41
1:A:252:THR:HB	1:A:338:VAL:HG11	2.03	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:ILE:HG12	1:A:403:LEU:HD13	2.03	0.40
1:A:690:ASN:C	1:A:690:ASN:OD1	2.60	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	816/880 (93%)	797 (98%)	18 (2%)	1 (0%)	56 57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	287	ASN

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
1	A	689/741 (93%)	666 (97%)	23 (3%)	45 46

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	22	ARG
1	A	34	LYS
1	A	36	TYR
1	A	156	VAL
1	A	188	ILE
1	A	261	VAL
1	A	305	THR
1	A	350	SER
1	A	450	LEU
1	A	484	LEU
1	A	487	THR
1	A	541	PHE
1	A	544	PHE
1	A	618	SER
1	A	621	SER
1	A	646	MET
1	A	661	SER
1	A	761	LEU
1	A	826	ARG
1	A	828	ARG
1	A	845	ARG
1	A	854	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	706	HIS

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	77/87 (88%)	21 (27%)	3 (3%)

All (21) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	3	C
2	B	4	C
2	B	5	G
2	B	6	G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	9	G
2	B	13	G
2	B	14	A
2	B	16	U
2	B	17	C
2	B	19	G
2	B	27	A
2	B	28	G
2	B	29	G
2	B	30	G
2	B	31	A
2	B	41	C
2	B	58	A
2	B	62	C
2	B	65	C
2	B	71	G
2	B	74	C

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	5	G
2	B	27	A
2	B	47(C)	U

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

1 non-standard protein/DNA/RNA residue 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	$\#\|Z\| > 2$
2	N79	B	76	2	21,36,37	1.75	2 (9%)	22,55,58	3.02	7 (31%)

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	N79	B	76	2	-	0/3/47/48	0/6/6/6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	76	N79	O1-C10	3.16	1.46	1.41
2	B	76	N79	C11-C12	6.38	1.44	1.40

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	76	N79	N3-C2-N1	-9.63	121.52	128.89
2	B	76	N79	C13-C11-C12	-4.12	119.52	122.17
2	B	76	N79	C4-C5-N7	-3.74	106.04	109.48
2	B	76	N79	C5'-C4'-C3'	-3.32	102.35	114.31
2	B	76	N79	C1'-N9-C4	-2.41	123.30	126.94
2	B	76	N79	O4'-C1'-N9	4.03	116.54	108.10
2	B	76	N79	C16-C12-C11	5.11	121.50	117.46

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	76	N79	1	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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	GOL	A	1861	-	5,5,5	0.52	0	5,5,5	1.06	1 (20%)
3	GOL	A	1862	-	5,5,5	0.37	0	5,5,5	0.49	0
3	GOL	A	1863	-	5,5,5	0.26	0	5,5,5	0.74	0
3	GOL	A	1864	-	5,5,5	0.49	0	5,5,5	0.95	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	GOL	A	1861	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1862	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1863	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1864	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1861	GOL	O1-C1-C2	2.07	120.20	110.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1861	GOL	1	0
3	A	1862	GOL	3	0
3	A	1863	GOL	4	0
3	A	1864	GOL	3	0

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, 95^{th} 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	821/880 (93%)	0.04	21 (2%) 59 65	14, 28, 55, 81	0
2	B	79/87 (90%)	0.63	14 (17%) 2 2	23, 47, 92, 110	0
All	All	900/967 (93%)	0.10	35 (3%) 43 51	14, 29, 64, 110	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	156	VAL	5.4
1	A	463	ILE	4.8
1	A	91	VAL	4.0
1	A	94	ASN	4.0
2	B	2	C	3.8
2	B	30	G	3.6
2	B	64	G	3.3
2	B	31	A	3.2
2	B	65	C	3.2
1	A	0	HIS	3.2
1	A	597	GLU	3.1
1	A	288	THR	3.1
1	A	155	ALA	3.0
1	A	600	ARG	2.9
1	A	283	ASP	2.9
1	A	188	ILE	2.9
2	B	29	G	2.8
1	A	279	ALA	2.8
2	B	73	A	2.8
1	A	286	ARG	2.8
1	A	742	ASP	2.7
2	B	4	C	2.6
2	B	42	C	2.6
2	B	3	C	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	187	GLU	2.4
2	B	1	G	2.4
1	A	484	LEU	2.4
1	A	285	CYS	2.3
2	B	51	G	2.3
2	B	52	G	2.2
2	B	68	C	2.2
1	A	603	LYS	2.1
1	A	4	GLN	2.1
1	A	240	ASP	2.0
1	A	287	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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, 95^{th} 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	N79	B	76	31/32	0.97	0.11	0.23	21,25,28,31	0

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, 95^{th} 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
3	GOL	A	1861	6/6	0.86	0.21	7.59	34,37,39,40	0
3	GOL	A	1862	6/6	0.94	0.17	6.31	30,32,32,35	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
3	GOL	A	1864	6/6	0.77	0.20	3.10	34,43,46,48	0
4	MG	A	1865	1/1	0.95	0.04	-2.14	51,51,51,51	0
3	GOL	A	1863	6/6	0.90	0.21	-	28,36,42,44	0
4	MG	B	1076	1/1	0.91	0.04	-	44,44,44,44	0

6.5 Other polymers (i)

There are no such residues in this entry.