



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:52 PM GMT

PDB ID : 4AS1  
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 : 2012-04-27  
Resolution : 2.02 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](mailto: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.

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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

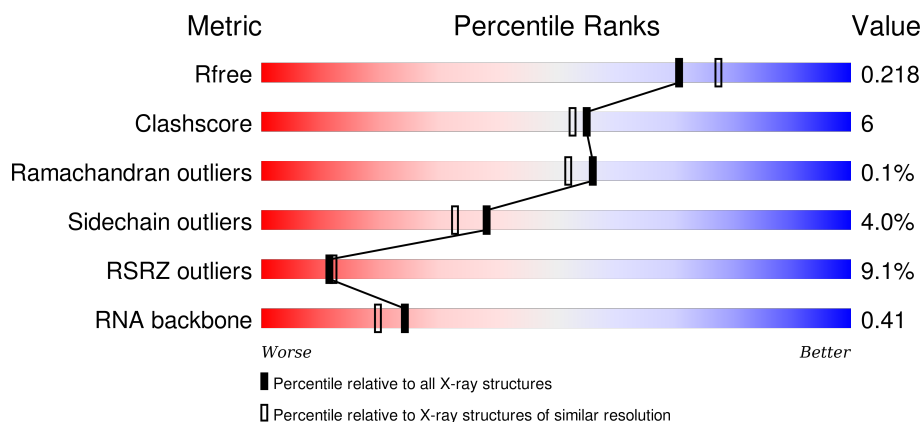
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.02 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	7858 (2.04-2.00)
Clashscore	102246	9060 (2.04-2.00)
Ramachandran outliers	100387	8952 (2.04-2.00)
Sidechain outliers	100360	8951 (2.04-2.00)
RSRZ outliers	91569	7873 (2.04-2.00)
RNA backbone	2183	1004 (2.74-1.30)

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  $\geq 3$ , 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	880	<div> <div>6%</div> <div> <div></div> <div>79%</div> <div>12%</div> <div>8%</div> </div> </div>
2	B	87	<div> <div>29%</div> <div> <div></div> <div>51%</div> <div>31%</div> <div>14%</div> <div>5%</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8665 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	812	Total	C	N	O	S	0	2	1
			6476	4119	1098	1219	40			

There are 20 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
A	0	HIS	-	EXPRESSION TAG	UNP P07813

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	83	Total	B	C	N	O	P	0	0
			1786	1	798	322	582	83		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Mg 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	365	Total 365	O 365	0	0
4	B	37	Total 37	O 37	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.68Å 77.11Å 91.14Å 90.00° 102.24° 90.00°	Depositor
Resolution (Å)	19.85 – 2.02 43.82 – 2.02	Depositor EDS
% Data completeness (in resolution range)	98.1 (19.85-2.02) 98.2 (43.82-2.02)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.26 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.6.0116	Depositor
R, $R_{free}$	0.208 , 0.249 0.211 , 0.218	Depositor DCC
$R_{free}$ test set	3933 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.6	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 53.3	EDS
Estimated twinning fraction	0.025 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 78347 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8665	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.80	7/6633 (0.1%)	0.82	8/8999 (0.1%)
2	B	0.46	1/1960 (0.1%)	0.90	5/3051 (0.2%)
All	All	0.74	8/8593 (0.1%)	0.84	13/12050 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	G	OP3-P	-10.10	1.49	1.61
1	A	575	TYR	CE2-CZ	-6.37	1.30	1.38
1	A	585	TRP	CD2-CE2	5.64	1.48	1.41
1	A	773	TRP	CD2-CE2	5.55	1.48	1.41
1	A	18	TRP	CD2-CE2	5.51	1.48	1.41
1	A	321	TRP	CD2-CE2	5.39	1.47	1.41
1	A	223	TRP	CD2-CE2	5.36	1.47	1.41
1	A	674	TRP	CD2-CE2	5.06	1.47	1.41

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	23	C	O5'-P-OP2	-8.58	97.98	105.70
2	B	14	A	C5'-C4'-C3'	-6.80	105.12	116.00
1	A	123	ARG	NE-CZ-NH1	6.56	123.58	120.30
2	B	47(D)	C	P-O3'-C3'	6.47	127.46	119.70
2	B	13	G	O3'-P-O5'	-6.26	92.10	104.00
1	A	344	ARG	NE-CZ-NH1	6.18	123.39	120.30
1	A	47	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	A	606	ASP	CB-CG-OD1	5.71	123.44	118.30
1	A	668	ARG	NE-CZ-NH2	-5.48	117.56	120.30
1	A	391	ASP	CB-CG-OD1	5.43	123.18	118.30
1	A	668	ARG	NE-CZ-NH1	5.05	122.83	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	67	ARG	NE-CZ-NH2	-5.01	117.79	120.30
2	B	49	G	O3'-P-O5'	-5.00	94.49	104.00

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	6476	0	6323	83	0
2	B	1786	0	904	16	0
3	B	1	0	0	0	0
4	A	365	0	0	19	0
4	B	37	0	0	1	0
All	All	8665	0	7227	97	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 6.

All (97) 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:570:LEU:C	1:A:617:MET:HE1	1.74	1.07
1:A:570:LEU:HB2	1:A:617:MET:HE2	1.40	1.02
1:A:477:THR:HG22	4:A:2236:HOH:O	1.63	0.97
1:A:633:MET:HE1	1:A:637:TYR:HE2	1.31	0.95
1:A:579:GLU:HG3	4:A:2279:HOH:O	1.68	0.92
1:A:439:ASP:HB3	1:A:441:THR:HG23	1.51	0.90
1:A:248:THR:HG23	2:B:76:N79:H5'1	1.55	0.86
1:A:633:MET:HE1	1:A:637:TYR:CE2	2.11	0.86
1:A:477:THR:CG2	4:A:2236:HOH:O	2.22	0.83
1:A:570:LEU:HB2	1:A:617:MET:CE	2.11	0.79
1:A:248:THR:CG2	2:B:76:N79:H5'1	2.13	0.78
1:A:570:LEU:C	1:A:617:MET:CE	2.53	0.77
1:A:571:ALA:N	1:A:617:MET:HE1	1.99	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:570:LEU:CB	1:A:617:MET:HE2	2.16	0.74
1:A:667:ASN:OD1	1:A:671:LYS:HE3	1.88	0.73
1:A:570:LEU:CB	1:A:617:MET:CE	2.68	0.72
2:B:27:A:H5''	2:B:27:A:H8	1.55	0.72
1:A:589:VAL:HG23	4:A:2292:HOH:O	1.91	0.70
1:A:826:ARG:NH1	1:A:847:VAL:HG21	2.07	0.70
1:A:798:ASP:HB2	4:A:2361:HOH:O	1.93	0.69
1:A:571:ALA:N	1:A:617:MET:CE	2.57	0.67
1:A:816:VAL:HG11	1:A:825:VAL:HG22	1.78	0.65
1:A:826:ARG:HH12	1:A:847:VAL:HG21	1.61	0.64
1:A:633:MET:CE	1:A:637:TYR:CE2	2.81	0.64
1:A:54:ARG:HG3	1:A:646:MET:HE2	1.78	0.64
2:B:71:G:C2'	2:B:72:U:O5'	2.48	0.61
2:B:73:A:H4'	2:B:74:C:OP2	2.00	0.61
1:A:507:GLN:NE2	4:A:2097:HOH:O	2.16	0.61
1:A:420:TRP:NE1	1:A:493[A]:PHE:CE1	2.68	0.60
1:A:749:LEU:HD12	4:A:2348:HOH:O	2.02	0.60
1:A:435:VAL:HG22	1:A:483:ALA:HB1	1.84	0.60
1:A:491:ASP:HB3	1:A:493[A]:PHE:CE1	2.37	0.59
1:A:570:LEU:O	1:A:617:MET:HE1	2.01	0.59
1:A:847:VAL:HG13	1:A:857:LEU:HD23	1.84	0.59
1:A:439:ASP:H	1:A:481:MET:HE1	1.67	0.59
2:B:71:G:H2'	2:B:72:U:O5'	2.03	0.59
2:B:27:A:C5'	2:B:27:A:H8	2.16	0.58
1:A:695:THR:HG22	1:A:697:ASN:H	1.69	0.58
1:A:435:VAL:CG2	1:A:483:ALA:HB1	2.34	0.58
1:A:194:LYS:NZ	4:A:2114:HOH:O	2.37	0.56
1:A:855:LEU:HD21	1:A:857:LEU:HD21	1.89	0.54
2:B:4:C:H2'	2:B:5:G:H5''	1.90	0.54
1:A:633:MET:CE	1:A:637:TYR:HE2	2.10	0.53
1:A:686:VAL:HG12	4:A:2328:HOH:O	2.08	0.53
1:A:553:MET:HE1	4:A:2109:HOH:O	2.06	0.53
2:B:24:A:N7	4:B:2022:HOH:O	2.33	0.53
1:A:469:ALA:O	1:A:471:PRO:HD3	2.08	0.53
1:A:633:MET:HE2	1:A:658:TRP:HZ2	1.73	0.53
1:A:620:MET:HE2	1:A:628:ILE:HG13	1.92	0.52
1:A:360:ILE:HD13	1:A:381:LEU:HD22	1.91	0.52
1:A:575:TYR:CE2	1:A:613:VAL:HG22	2.44	0.52
2:B:27:A:C5'	2:B:27:A:C8	2.92	0.51
2:B:16:U:O2'	2:B:17:C:O2	2.17	0.51
1:A:798:ASP:C	1:A:818:VAL:HG23	2.31	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:VAL:HG13	1:A:265:HIS:HB3	1.94	0.50
2:B:27:A:H5''	2:B:27:A:C8	2.41	0.49
1:A:596:ASP:HB3	1:A:602:VAL:HG11	1.94	0.49
1:A:690:ASN:H	1:A:745:GLN:HE22	1.61	0.48
1:A:360:ILE:CD1	1:A:381:LEU:HD22	2.44	0.48
1:A:287:ASN:O	1:A:288:THR:CB	2.62	0.48
1:A:804:VAL:HG11	1:A:835:VAL:HG21	1.95	0.47
1:A:822:GLU:HG2	1:A:849:TYR:CE2	2.49	0.47
1:A:262:ALA:CA	1:A:326:VAL:HG22	2.45	0.47
2:B:66:U:H2'	2:B:67:C:O4'	2.16	0.46
1:A:203:LEU:HD11	4:A:2144:HOH:O	2.16	0.46
1:A:477:THR:HB	4:A:2236:HOH:O	2.14	0.45
1:A:136:GLU:OE2	1:A:495:GLU:OE2	2.34	0.45
1:A:481:MET:HG3	1:A:482:PRO:HD2	1.98	0.45
2:B:17:C:H5''	2:B:18:G:OP1	2.17	0.45
1:A:265:HIS:ND1	1:A:266:PRO:HD2	2.32	0.45
1:A:416:ARG:NH1	4:A:2209:HOH:O	2.43	0.45
1:A:269:GLN:O	1:A:272:ALA:HB3	2.18	0.44
1:A:477:THR:CB	4:A:2236:HOH:O	2.57	0.44
1:A:797:GLU:O	1:A:818:VAL:HG21	2.18	0.44
1:A:526:ILE:O	1:A:526:ILE:HG23	2.17	0.44
1:A:196:THR:HG21	4:A:2208:HOH:O	2.18	0.44
1:A:265:HIS:CE1	1:A:266:PRO:HD2	2.53	0.44
1:A:832:GLU:HB3	1:A:835:VAL:HB	2.00	0.44
1:A:194:LYS:CE	4:A:2114:HOH:O	2.67	0.43
1:A:254:MET:SD	1:A:313:LEU:HD12	2.58	0.43
1:A:253:PHE:HD1	1:A:336:MET:SD	2.41	0.43
1:A:272:ALA:N	1:A:278:LEU:HD23	2.34	0.42
1:A:814:ILE:HD12	1:A:816:VAL:CG2	2.49	0.42
1:A:571:ALA:N	1:A:617:MET:HE3	2.35	0.42
1:A:611:GLU:HG2	4:A:2301:HOH:O	2.18	0.42
1:A:822:GLU:HG2	1:A:849:TYR:CD2	2.54	0.42
1:A:439:ASP:CB	1:A:441:THR:HG23	2.35	0.42
1:A:82:PHE:CE2	1:A:128:CYS:HA	2.55	0.41
1:A:262:ALA:HA	1:A:326:VAL:HG22	2.03	0.41
1:A:568:MET:SD	1:A:619:LYS:HG3	2.60	0.41
1:A:826:ARG:O	1:A:829:ALA:HB3	2.20	0.41
1:A:443:MET:HE3	4:A:2221:HOH:O	2.20	0.41
1:A:803:VAL:HG11	1:A:810:VAL:HG22	2.03	0.41
1:A:43:TYR:CD1	1:A:44:PRO:HD2	2.56	0.41
1:A:691:VAL:HG22	4:A:2329:HOH:O	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:30:G:C6	2:B:31:A:C6	3.09	0.40
2:B:5:G:H5''	2:B:5:G:H8	1.86	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	806/880 (92%)	786 (98%)	19 (2%)	1 (0%)	56 52

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	439	ASP

### 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	684/741 (92%)	656 (96%)	28 (4%)	37 32

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LYS

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Mol	Chain	Res	Type
1	A	36	TYR
1	A	84	LEU
1	A	94	ASN
1	A	261	VAL
1	A	286	ARG
1	A	314	THR
1	A	316	GLU
1	A	376	THR
1	A	381	LEU
1	A	477	THR
1	A	484	LEU
1	A	493[A]	PHE
1	A	493[B]	PHE
1	A	516	GLU
1	A	541	PHE
1	A	613	VAL
1	A	618	SER
1	A	622	LYS
1	A	625	ASN
1	A	646	MET
1	A	686	VAL
1	A	761	LEU
1	A	798	ASP
1	A	801	LEU
1	A	819	ASP
1	A	834	LEU
1	A	857	LEU

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

Mol	Chain	Res	Type
1	A	425	GLN
1	A	533	HIS
1	A	625	ASN
1	A	745	GLN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	80/87 (91%)	29 (36%)	2 (2%)

All (29) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	2	C
2	B	5	G
2	B	9	G
2	B	16	U
2	B	17	C
2	B	18	G
2	B	27	A
2	B	29	G
2	B	30	G
2	B	31	A
2	B	39	U
2	B	42	C
2	B	45	G
2	B	47(C)	U
2	B	47(D)	C
2	B	47(E)	G
2	B	47(F)	C
2	B	50	C
2	B	51	G
2	B	52	G
2	B	56	C
2	B	63	C
2	B	67	C
2	B	68	C
2	B	69	G
2	B	70	G
2	B	71	G
2	B	72	U
2	B	73	A

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	27	A
2	B	47(D)	C

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

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	1.98	3 (14%)	22,55,58	2.67	5 (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	N79	B	76	2	-	0/3/47/48	0/6/6/6

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	76	N79	C10-C11	2.80	1.53	1.50
2	B	76	N79	O1-C10	4.39	1.47	1.41
2	B	76	N79	C11-C12	6.49	1.44	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	76	N79	N3-C2-N1	-9.08	121.94	128.89
2	B	76	N79	C13-C11-C12	-4.40	119.33	122.17
2	B	76	N79	C5'-C4'-C3'	-3.41	102.02	114.31
2	B	76	N79	C16-C12-C11	3.62	120.32	117.46
2	B	76	N79	O4'-C1'-N9	3.80	116.05	108.10

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
2	B	76	N79	2	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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.

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	812/880 (92%)	0.42	56 (6%) 20 21	18, 38, 79, 115	0
2	B	82/87 (94%)	1.31	25 (30%) 1 1	31, 84, 139, 169	0
All	All	894/967 (92%)	0.50	81 (9%) 11 12	18, 40, 92, 169	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	818	VAL	7.3
2	B	2	C	5.2
2	B	31	A	5.1
2	B	36	A	5.1
1	A	821	THR	4.9
1	A	820	ALA	4.6
1	A	94	ASN	4.5
2	B	64	G	4.5
1	A	847	VAL	4.4
2	B	1	G	4.3
2	B	73	A	4.2
2	B	53	G	4.2
1	A	91	VAL	4.1
1	A	848	ILE	4.0
1	A	817	PRO	4.0
1	A	819	ASP	3.7
1	A	288	THR	3.5
1	A	62	ILE	3.5
2	B	3	C	3.4
2	B	51	G	3.4
1	A	823	GLU	3.3
2	B	30	G	3.3
1	A	528	ILE	3.2
1	A	493[A]	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	155	ALA	3.2
1	A	806	VAL	3.2
2	B	52	G	3.2
2	B	71	G	3.1
2	B	65	C	3.0
2	B	63	C	3.0
1	A	835	VAL	3.0
2	B	28	G	2.9
2	B	54	U	2.8
2	B	55	U	2.8
1	A	827	GLU	2.8
1	A	857	LEU	2.8
1	A	56	TYR	2.7
1	A	693	ALA	2.7
1	A	38	LEU	2.7
1	A	58	ILE	2.7
1	A	692	ASP	2.7
1	A	564	LEU	2.7
2	B	38	A	2.7
2	B	41	C	2.7
1	A	798	ASP	2.6
1	A	686	VAL	2.6
2	B	37	A	2.6
1	A	61	VAL	2.6
1	A	796	VAL	2.6
1	A	652	ALA	2.6
1	A	849	TYR	2.6
1	A	286	ARG	2.6
1	A	41	LEU	2.5
1	A	694	LEU	2.5
1	A	691	VAL	2.5
1	A	526	ILE	2.5
1	A	37	CYS	2.5
2	B	69	G	2.4
1	A	283	ASP	2.4
1	A	684	GLY	2.3
1	A	527	TYR	2.3
2	B	72	U	2.3
2	B	62	C	2.3
1	A	284	GLU	2.2
1	A	824	GLN	2.2
2	B	17	C	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	40	C	2.2
1	A	716	ILE	2.2
1	A	57	THR	2.2
1	A	77	ILE	2.2
1	A	287	ASN	2.2
1	A	748	ALA	2.2
1	A	39	SER	2.1
1	A	721	THR	2.1
1	A	764	PHE	2.1
1	A	578	GLY	2.1
1	A	829	ALA	2.1
1	A	78	GLY	2.1
1	A	59	GLY	2.0
1	A	40	MET	2.0
1	A	92	LYS	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<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
2	N79	B	76	31/32	0.96	0.10	-0.23	29,31,34,36	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<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	MG	B	1076	1/1	0.92	0.06	-	50,50,50,50	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.