



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:47 AM UTC

PDB ID : 4ARC / pdb_00004arc
Title : Ternary complex of E. coli leucyl-tRNA synthetase, tRNA(Leu) and leucine in the editing conformation
Authors : Palencia, A.; Crepin, T.; Vu, M.T.; Lincecum Jr, T.L.; Martinis, S.A.; Cusack, S.
Deposited on : 2012-04-23
Resolution : 2.00 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

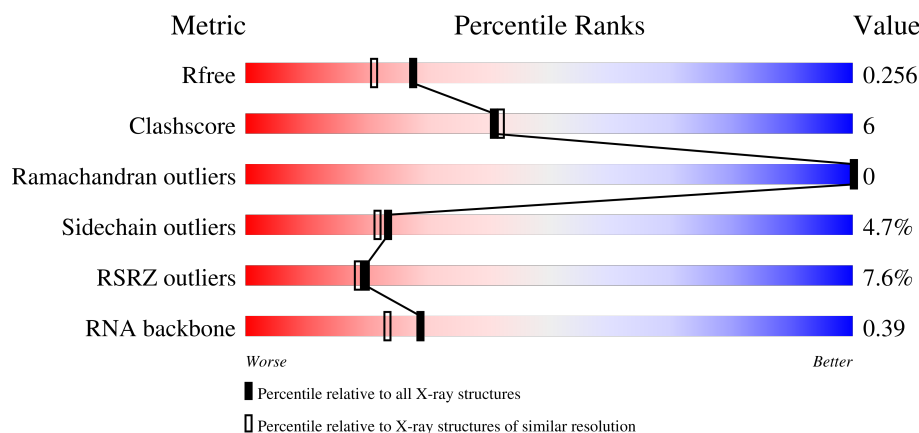
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)
RNA backbone	3983	1000 (2.36-1.64)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	8% 79% 12% 8%
2	B	87	% 36% 33% 13% 18%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8406 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	813	Total	C	N	O	S	0	0	0
			6465	4118	1091	1218	38			

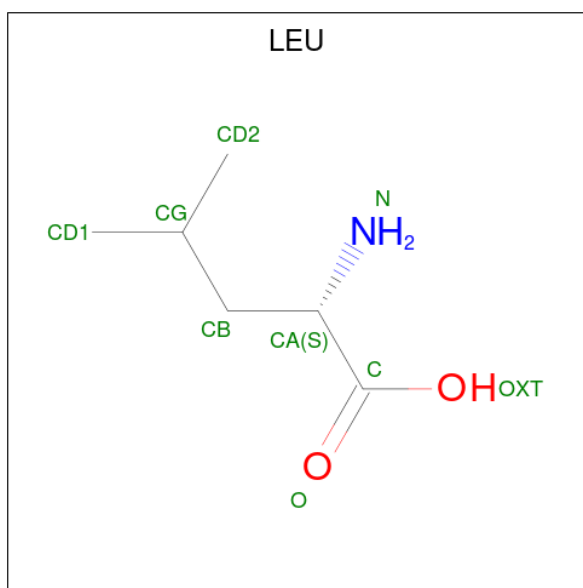
There are 20 discrepancies between the modelled and reference sequences:

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	71	Total	C	N	O	P	0	0	0
			1516	675	271	499	71			

- Molecule 3 is LEUCINE (CCD ID: LEU) (formula: $C_6H_{13}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	6	1	2		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	385	Total	O	0	0
			385	385		
5	B	30	Total	O	0	0
			30	30		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.08Å 119.37Å 141.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.57 – 2.00 45.57 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.3 (45.57-2.00) 99.3 (45.57-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.6.0116	Depositor
R, R_{free}	0.210 , 0.257 0.208 , 0.256	Depositor DCC
R_{free} test set	4411 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8406	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.44% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	1.11	4/6619 (0.1%)	1.00	3/8986 (0.0%)
2	B	0.50	0/1691	1.11	12/2630 (0.5%)
All	All	1.02	4/8310 (0.0%)	1.03	15/11616 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	30	GLU	C-O	-5.61	1.16	1.24
1	A	668	ARG	C-O	-5.34	1.17	1.24
1	A	612	LEU	C-O	-5.27	1.17	1.24
1	A	651	PRO	C-O	5.04	1.29	1.23

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	58	A	O3'-P-O5'	-8.14	91.78	104.00
2	B	76	A	C1'-C2'-O2'	6.52	118.17	108.40
2	B	48	U	O3'-P-O5'	-5.87	95.19	104.00
2	B	56	C	C3'-C2'-O2'	5.74	119.32	110.70
1	A	43	TYR	N-CA-C	-5.71	102.85	109.93
2	B	55	U	O3'-P-O5'	5.59	112.38	104.00
1	A	50	MET	N-CA-C	5.57	117.35	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	47(G)	G	C1'-C2'-O2'	5.51	116.67	108.40
2	B	16	U	C4'-C3'-O3'	5.47	117.61	109.40
2	B	50	C	O3'-P-O5'	-5.44	95.84	104.00
1	A	515	SER	N-CA-C	5.27	117.03	111.28
2	B	25	C	C2'-C3'-O3'	-5.26	105.81	113.70
2	B	12	G	O3'-P-O5'	-5.21	96.19	104.00
2	B	44	C	C3'-C2'-O2'	5.18	118.47	110.70
2	B	20(A)	A	C4'-C3'-O3'	-5.03	105.45	113.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	94	ASN	Peptide

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	A	6465	0	6331	82	0
2	B	1516	0	770	18	0
3	A	9	0	10	0	0
4	B	1	0	0	0	0
5	A	385	0	0	19	0
5	B	30	0	0	1	0
All	All	8406	0	7111	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)
2:B:26:A:H2'	2:B:27:A:H5'	1.41	0.98
1:A:512:MET:HE2	1:A:548:MET:SD	2.07	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:PHE:HB2	5:A:2157:HOH:O	1.71	0.90
2:B:26:A:C2'	2:B:27:A:H5'	2.03	0.88
1:A:575:TYR:CZ	1:A:613:VAL:HG11	2.08	0.88
1:A:223:TRP:CE3	1:A:535:ILE:HD12	2.16	0.81
1:A:468:LYS:NZ	1:A:487:THR:HG23	1.95	0.81
1:A:512:MET:CE	1:A:548:MET:SD	2.69	0.80
1:A:695:THR:HG22	5:A:2317:HOH:O	1.82	0.80
1:A:95:THR:HG21	5:A:2088:HOH:O	1.82	0.79
1:A:468:LYS:HZ3	1:A:487:THR:HG23	1.48	0.79
1:A:243:LEU:HB3	5:A:2157:HOH:O	1.92	0.70
1:A:492:THR:HG23	5:A:2082:HOH:O	1.92	0.70
1:A:575:TYR:CE1	1:A:613:VAL:HG11	2.25	0.70
1:A:569:VAL:HG12	1:A:620:MET:HE3	1.76	0.68
1:A:840:ASP:HB3	5:A:2383:HOH:O	1.94	0.68
2:B:52:G:N2	2:B:63:C:C2	2.62	0.68
1:A:593:VAL:HG12	1:A:594:GLU:N	2.10	0.65
1:A:695:THR:CG2	5:A:2317:HOH:O	2.40	0.65
1:A:244:THR:HG22	1:A:332:THR:HG21	1.76	0.65
1:A:50:MET:HE1	1:A:646:MET:SD	2.37	0.63
1:A:422:VAL:HG11	1:A:494:MET:HE1	1.82	0.62
1:A:593:VAL:HG11	1:A:601:ILE:HG12	1.81	0.62
1:A:593:VAL:HG13	1:A:601:ILE:HG23	1.80	0.62
1:A:841:GLY:O	1:A:842:VAL:HG13	1.99	0.61
1:A:142:GLU:HG2	5:A:2111:HOH:O	1.99	0.61
1:A:267:LEU:HD21	1:A:305:THR:HG21	1.84	0.60
1:A:393:GLU:CD	1:A:393:GLU:H	2.11	0.58
1:A:98:ALA:HB3	1:A:99:PRO:CD	2.34	0.58
1:A:593:VAL:CG1	1:A:594:GLU:N	2.67	0.58
1:A:349:ALA:HA	5:A:2189:HOH:O	2.03	0.58
1:A:492:THR:CG2	5:A:2082:HOH:O	2.50	0.57
1:A:156:VAL:HG21	5:A:2123:HOH:O	2.06	0.55
1:A:393:GLU:CD	1:A:393:GLU:N	2.66	0.54
1:A:142:GLU:CG	5:A:2111:HOH:O	2.55	0.54
1:A:191:TRP:CH2	1:A:434:MET:HE2	2.42	0.54
2:B:8:U:O2'	2:B:20(A):A:N1	2.37	0.54
1:A:352:TYR:HB2	5:A:2189:HOH:O	2.07	0.54
1:A:25:GLU:OE1	5:A:2029:HOH:O	2.18	0.54
1:A:391:ASP:HB3	5:A:2205:HOH:O	2.09	0.53
2:B:43:U:O2'	5:B:2022:HOH:O	2.19	0.52
1:A:463:ILE:HD12	1:A:463:ILE:N	2.25	0.52
1:A:595:ARG:HD2	1:A:599:GLY:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:28:G:H2'	2:B:29:G:O4'	2.11	0.51
1:A:498:TRP:CE3	1:A:512:MET:HE3	2.46	0.50
1:A:695:THR:HG22	1:A:697:ASN:H	1.76	0.50
1:A:267:LEU:C	1:A:267:LEU:HD23	2.36	0.50
1:A:468:LYS:HZ2	1:A:487:THR:HG23	1.74	0.50
1:A:828:ARG:HH11	1:A:828:ARG:HG3	1.77	0.49
2:B:47(G):G:H2'	2:B:47(H):C:C6	2.48	0.49
1:A:847:VAL:HB	5:A:2373:HOH:O	2.12	0.49
1:A:98:ALA:HB3	1:A:99:PRO:HD3	1.95	0.49
1:A:223:TRP:CD2	1:A:535:ILE:HD12	2.48	0.49
1:A:695:THR:HG22	1:A:696:GLU:N	2.29	0.48
1:A:828:ARG:HH11	1:A:828:ARG:CG	2.27	0.48
1:A:81:ALA:H	1:A:127:THR:HG1	1.60	0.47
2:B:5:G:H3'	2:B:6:G:H5''	1.95	0.47
1:A:512:MET:HE1	1:A:548:MET:SD	2.53	0.47
1:A:215:THR:O	1:A:219:MET:HE3	2.15	0.47
1:A:219:MET:HE1	1:A:531:ILE:HD11	1.97	0.47
1:A:572:ASP:O	1:A:588:PRO:HG3	2.15	0.47
1:A:828:ARG:HG3	1:A:828:ARG:NH1	2.30	0.47
1:A:360:ILE:HD13	1:A:381:LEU:HD22	1.96	0.46
1:A:587:SER:HA	1:A:625:ASN:O	2.16	0.46
1:A:575:TYR:CE1	1:A:613:VAL:CG1	2.98	0.45
2:B:53:G:C2	2:B:62:C:C2	3.04	0.45
1:A:344:ARG:NH2	2:B:76:A:N3	2.64	0.45
1:A:284:GLU:HA	5:A:2177:HOH:O	2.17	0.45
2:B:52:G:N2	2:B:63:C:O2	2.49	0.45
1:A:354:LEU:HD12	5:A:2189:HOH:O	2.16	0.45
1:A:668:ARG:HD3	5:A:2300:HOH:O	2.17	0.45
1:A:575:TYR:CZ	1:A:613:VAL:CG1	2.93	0.45
1:A:82:PHE:CE2	1:A:128:CYS:HA	2.52	0.45
1:A:50:MET:CE	1:A:646:MET:SD	3.04	0.44
2:B:51:G:C4	2:B:64:G:N2	2.86	0.44
1:A:849:TYR:CZ	1:A:851:PRO:HA	2.53	0.44
2:B:39:U:H2'	2:B:40:C:C6	2.54	0.43
1:A:368:PRO:HB3	1:A:375:LEU:HD22	2.00	0.42
2:B:10:G:C6	2:B:26:A:C2	3.07	0.42
1:A:426:ARG:HA	1:A:426:ARG:HD2	1.83	0.42
1:A:47:ARG:NH2	1:A:107:TYR:OH	2.52	0.42
1:A:327:LEU:O	1:A:333:GLY:HA2	2.20	0.42
1:A:258:TYR:CZ	1:A:337:ALA:HB1	2.55	0.42
1:A:154:SER:O	1:A:187:GLU:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:53:G:O2'	2:B:54:U:H5'	2.20	0.42
1:A:310:VAL:O	1:A:310:VAL:HG12	2.20	0.41
1:A:468:LYS:HZ2	1:A:487:THR:CG2	2.33	0.41
1:A:738:LYS:HE2	2:B:42:C:OP2	2.21	0.41
1:A:575:TYR:O	1:A:613:VAL:HG12	2.21	0.41
1:A:853:LYS:HE2	2:B:19:G:O2'	2.21	0.41
1:A:480:GLY:O	1:A:481:MET:C	2.63	0.41
2:B:9:G:H2'	2:B:9:G:N3	2.36	0.41
1:A:125:LEU:HD23	1:A:125:LEU:HA	1.93	0.40
1:A:847:VAL:HG11	1:A:855:LEU:HD11	2.03	0.40
1:A:305:THR:HG22	1:A:307:PHE:H	1.86	0.40
1:A:305:THR:HB	1:A:320:VAL:O	2.21	0.40
1:A:360:ILE:CD1	1:A:381:LEU:HD22	2.52	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	807/880 (92%)	783 (97%)	24 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	683/741 (92%)	651 (95%)	32 (5%)	23	22

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	VAL
1	A	95	THR
1	A	154	SER
1	A	188	ILE
1	A	234	PHE
1	A	284	GLU
1	A	310	VAL
1	A	329	GLU
1	A	335	VAL
1	A	336	MET
1	A	350	SER
1	A	381	LEU
1	A	435	VAL
1	A	450	LEU
1	A	484	LEU
1	A	492	THR
1	A	541	PHE
1	A	603	LYS
1	A	605	LYS
1	A	613	VAL
1	A	618	SER
1	A	622	LYS
1	A	623	SER
1	A	798	ASP
1	A	826	ARG
1	A	828	ARG
1	A	837	LYS
1	A	842	VAL
1	A	843	THR
1	A	845	ARG
1	A	848	ILE
1	A	859	VAL

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

Mol	Chain	Res	Type
1	A	2	GLN

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Mol	Chain	Res	Type
1	A	235	ASN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	68/87 (78%)	14 (20%)	3 (4%)

All (14) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	6	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	31	A
2	B	42	C
2	B	47(C)	U
2	B	47(D)	C
2	B	50	C
2	B	52	G
2	B	68	C

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	16	U
2	B	29	G
2	B	47(C)	U

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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	LEU	A	1001	-	6,8,8	0.77	0	5,10,10	1.41	1 (20%)

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	LEU	A	1001	-	-	0/8/8/8	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1001	LEU	OXT-C-O	-2.13	119.24	124.08

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

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	A	813/880 (92%)	0.37	66 (8%) 18 17	20, 38, 72, 100	0
2	B	71/87 (81%)	0.53	1 (1%) 73 73	31, 65, 104, 144	0
All	All	884/967 (91%)	0.38	67 (7%) 20 18	20, 41, 79, 144	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	281	PHE	5.3
1	A	282	ILE	5.0
1	A	279	ALA	4.9
1	A	305	THR	4.8
1	A	271	ALA	4.0
1	A	283	ASP	3.7
1	A	156	VAL	3.6
1	A	327	LEU	3.5
1	A	602	VAL	3.5
1	A	348	PHE	3.4
1	A	349	ALA	3.4
1	A	607	ALA	3.4
1	A	272	ALA	3.3
1	A	325	PHE	3.3
1	A	591	ALA	3.3
1	A	592	ILE	3.3
1	A	278	LEU	3.3
1	A	346	TYR	3.2
1	A	320	VAL	3.2
1	A	262	ALA	3.2
1	A	321	TRP	3.1
1	A	307	PHE	3.0
1	A	280	ALA	2.9
1	A	590	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	239	TYR	2.9
1	A	261	VAL	2.8
1	A	263	ALA	2.8
1	A	267	LEU	2.8
2	B	5	G	2.8
1	A	355	ASN	2.7
1	A	604	ALA	2.7
1	A	481	MET	2.6
1	A	236	VAL	2.6
1	A	601	ILE	2.6
1	A	94	ASN	2.6
1	A	310	VAL	2.5
1	A	333	GLY	2.5
1	A	480	GLY	2.5
1	A	401	ASP	2.4
1	A	589	VAL	2.4
1	A	593	VAL	2.4
1	A	330	TYR	2.4
1	A	237	ASN	2.4
1	A	350	SER	2.4
1	A	345	ASP	2.4
1	A	603	LYS	2.4
1	A	265	HIS	2.4
1	A	337	ALA	2.3
1	A	354	LEU	2.3
1	A	598	LYS	2.3
1	A	302	GLY	2.3
1	A	600	ARG	2.3
1	A	356	ILE	2.3
1	A	326	VAL	2.3
1	A	352	TYR	2.3
1	A	808	GLY	2.2
1	A	595	ARG	2.2
1	A	318	ILE	2.2
1	A	377	GLU	2.2
1	A	398	ALA	2.1
1	A	270	LYS	2.1
1	A	487	THR	2.1
1	A	188	ILE	2.1
1	A	482	PRO	2.0
1	A	353	GLY	2.0
1	A	260	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	257	THR	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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
3	LEU	A	1001	9/9	0.86	0.11	27,33,41,42	0
4	MG	B	1077	1/1	0.90	0.20	62,62,62,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.