



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 11:40 AM EDT

PDB ID : 4AQ7 / pdb_00004aq7
Title : Ternary complex of E. coli leucyl-tRNA synthetase, tRNA(leu) and leucyl-adenylate analogue in the aminoacylation conformation
Authors : Palencia, A.; Crepin, T.; Vu, M.T.; Lincecum Jr, T.L.; Martinis, S.A.; Cusack, S.
Deposited on : 2012-04-13
Resolution : 2.50 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

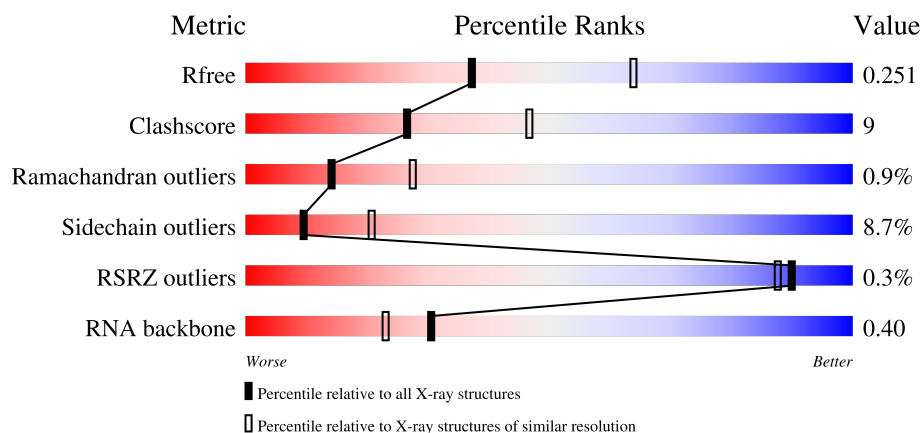
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)
RNA backbone	3983	1003 (2.78-2.22)

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	 73% 22% • •
1	D	880	 72% 23% • •
2	B	87	 49% 36% 5% • 9%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	E	87	% 45% 34% 8% 11%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 17212 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	860	Total	C	N	O	S	0	0	0
			6835	4339	1159	1292	45			
1	D	860	Total	C	N	O	S	0	0	0
			6835	4339	1159	1292	45			

There are 40 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
D	-19	MET	-	expression tag	UNP P07813
D	-18	GLY	-	expression tag	UNP P07813
D	-17	SER	-	expression tag	UNP P07813
D	-16	SER	-	expression tag	UNP P07813
D	-15	HIS	-	expression tag	UNP P07813

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	-14	HIS	-	expression tag	UNP P07813
D	-13	HIS	-	expression tag	UNP P07813
D	-12	HIS	-	expression tag	UNP P07813
D	-11	HIS	-	expression tag	UNP P07813
D	-10	HIS	-	expression tag	UNP P07813
D	-9	SER	-	expression tag	UNP P07813
D	-8	SER	-	expression tag	UNP P07813
D	-7	GLY	-	expression tag	UNP P07813
D	-6	LEU	-	expression tag	UNP P07813
D	-5	VAL	-	expression tag	UNP P07813
D	-4	PRO	-	expression tag	UNP P07813
D	-3	ARG	-	expression tag	UNP P07813
D	-2	GLY	-	expression tag	UNP P07813
D	-1	SER	-	expression tag	UNP P07813
D	0	HIS	-	expression tag	UNP P07813

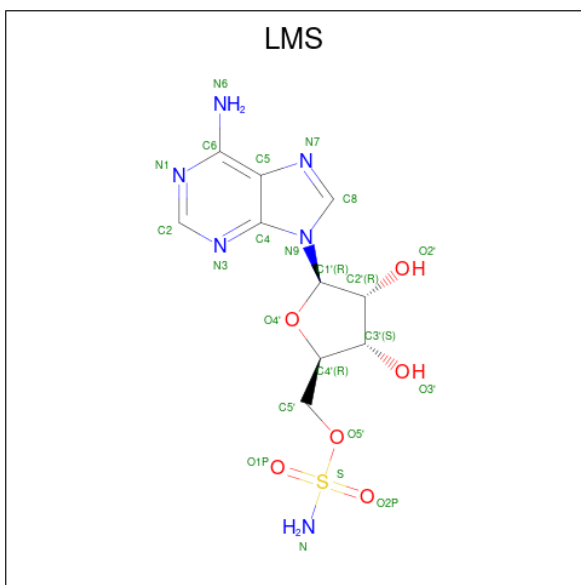
- Molecule 2 is a RNA chain called E. COLI TRNALEU UAA ISOACCEPTOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	79	Total	C	N	O	P	0	0	0
			1675	743	303	550	79			
2	E	77	Total	C	N	O	P	0	0	0
			1632	724	295	536	77			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

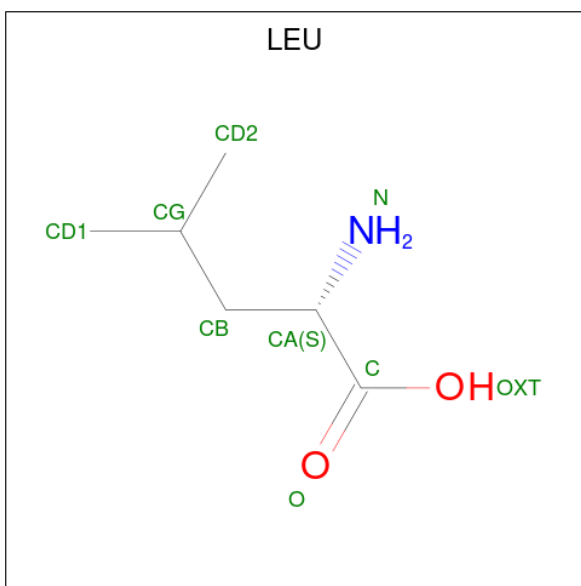
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		

- Molecule 4 is [(2R,3S,4R,5R)-5-(6-AMINO-9H-PURIN-9-YL)-3,4-DIHYDROXYTETRAHYDRO-2-FURANYL]METHYL SULFAMATE (CCD ID: LMS) (formula: C₁₀H₁₄N₆O₆S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			23	10	6	6	1		
4	D	1	Total	C	N	O	S	0	0
			23	10	6	6	1		

- Molecule 5 is LEUCINE (CCD ID: LEU) (formula: $C_6H_{13}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			8	6	1	1		
5	D	1	Total	C	N	O	0	0
			8	6	1	1		

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total 1	Mg 1	0	0
6	E	1	Total 1	Mg 1	0	0

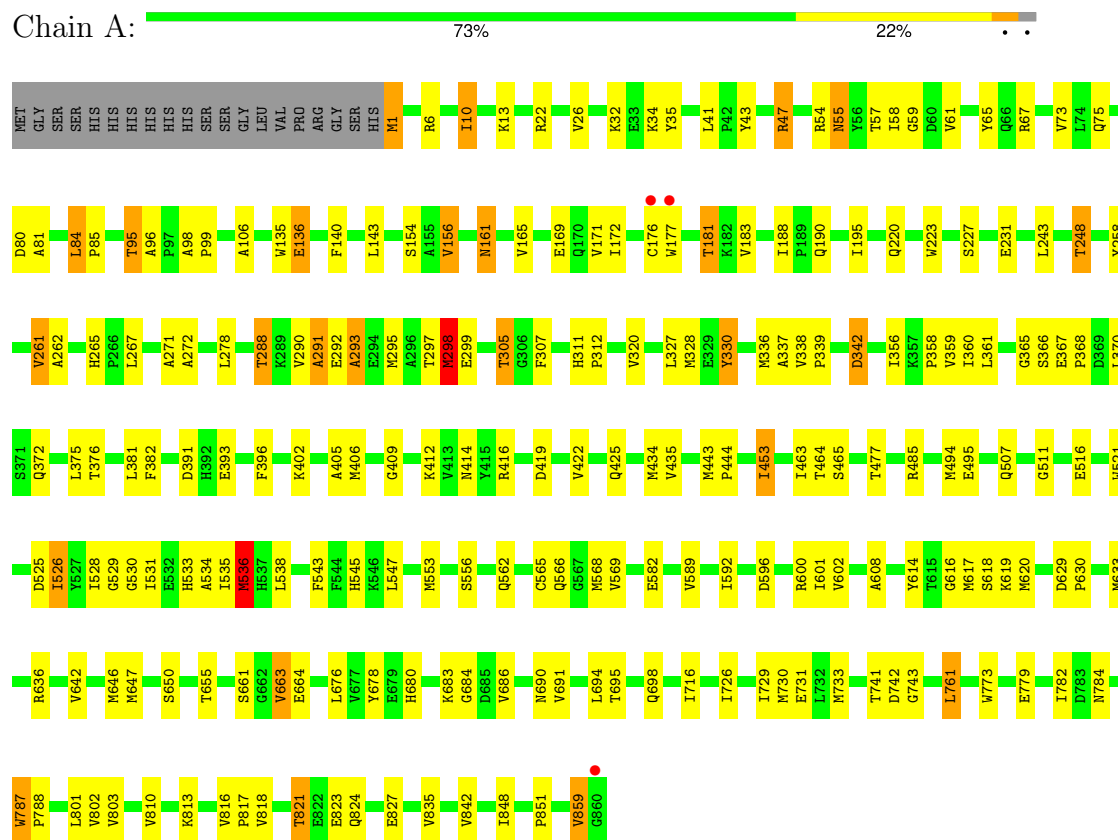
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	104	Total 104	O 104	0	0
7	B	15	Total 15	O 15	0	0
7	D	47	Total 47	O 47	0	0
7	E	3	Total 3	O 3	0	0

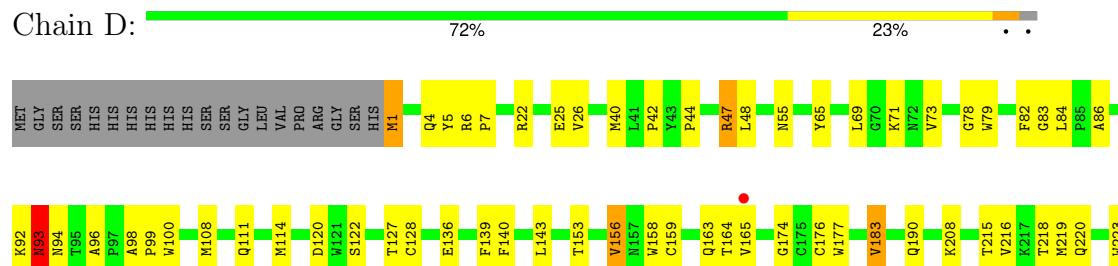
3 Residue-property plots [i](#)

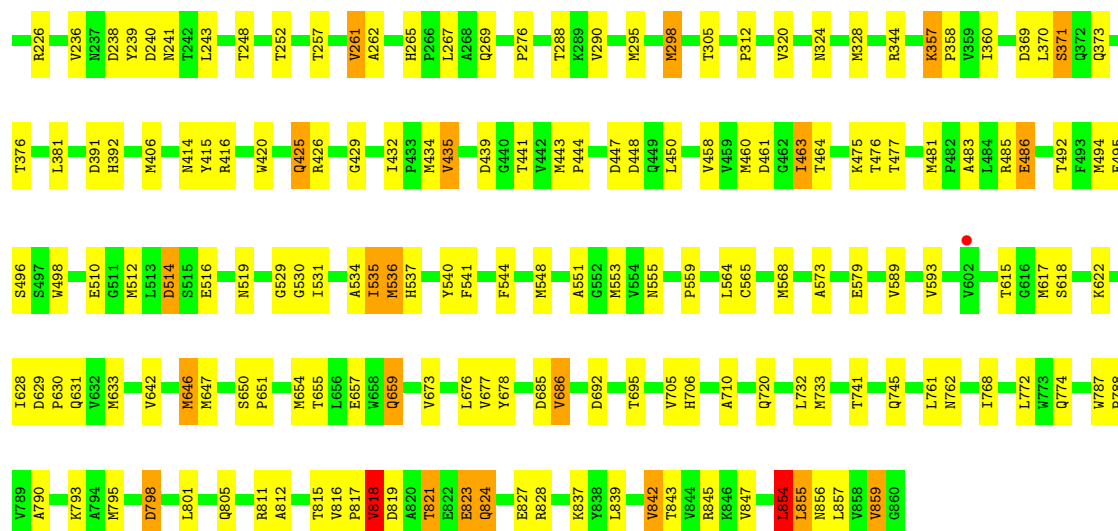
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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

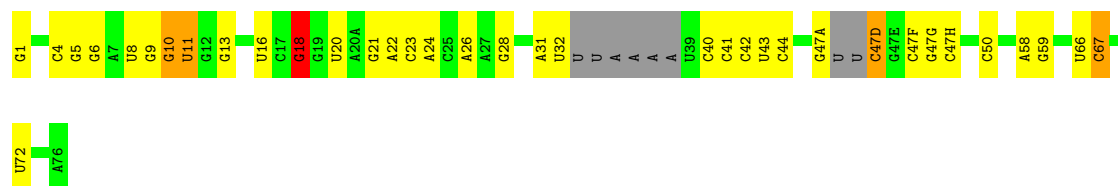


• Molecule 1: LEUCINE-TRNA LIGASE

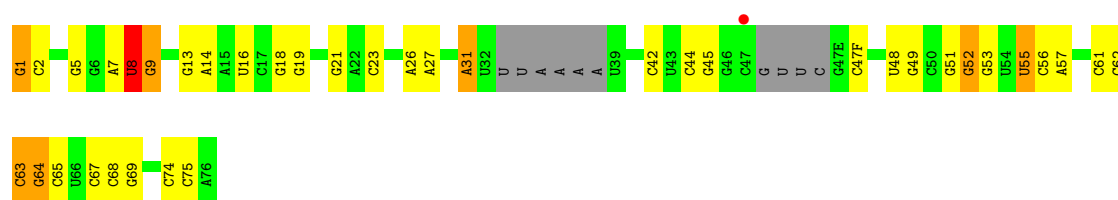




● Molecule 2: E. COLI TRNALEU UAA ISOACCEPTOR



● Molecule 2: E. COLI TRNALEU UAA ISOACCEPTOR



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	158.66Å 69.20Å 228.84Å 90.00° 104.35° 90.00°	Depositor
Resolution (Å)	39.69 – 2.50 39.69 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.3 (39.69-2.50) 90.3 (39.69-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.6.0116	Depositor
R, R_{free}	0.188 , 0.250 0.189 , 0.251	Depositor DCC
R_{free} test set	3826 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	50.5	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17212	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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, LMS, ZN

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	1.18	11/6996 (0.2%)	1.15	13/9497 (0.1%)
1	D	0.99	8/6996 (0.1%)	1.05	10/9497 (0.1%)
2	B	0.60	1/1869 (0.1%)	1.24	20/2908 (0.7%)
2	E	0.50	1/1821 (0.1%)	1.07	11/2833 (0.4%)
All	All	1.00	21/17682 (0.1%)	1.11	54/24735 (0.2%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	75	GLN	CD-OE1	8.07	1.38	1.23
2	B	1	G	OP3-P	7.35	1.63	1.48
1	A	507	GLN	CD-OE1	6.59	1.36	1.23
1	A	58	ILE	CA-CB	-6.15	1.47	1.54
1	A	414	ASN	CG-OD1	5.87	1.34	1.23
1	A	680	HIS	CG-ND1	-5.83	1.31	1.38
1	A	372	GLN	CD-OE1	5.61	1.34	1.23
1	A	220	GLN	C-O	-5.61	1.17	1.24
1	D	190	GLN	CD-OE1	5.57	1.34	1.23
1	D	93	ASN	CG-OD1	5.55	1.34	1.23
1	D	414	ASN	CG-OD1	5.54	1.34	1.23
2	E	1	G	OP3-P	5.49	1.59	1.48
1	A	26	VAL	C-O	-5.34	1.18	1.23
1	D	392	HIS	CG-CD2	5.29	1.41	1.35
1	A	106	ALA	N-CA	5.27	1.52	1.46
1	D	659	GLN	CD-OE1	5.25	1.33	1.23
1	D	706	HIS	CG-ND1	-5.22	1.32	1.38
1	A	680	HIS	CD2-NE2	-5.20	1.32	1.37
1	A	161	ASN	CG-OD1	5.17	1.33	1.23
1	D	373	GLN	CD-NE2	-5.09	1.22	1.33
1	D	218	THR	C-O	-5.06	1.18	1.24

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	47(D)	C	P-O3'-C3'	10.76	132.62	119.70
2	B	40	C	O3'-P-O5'	-8.48	91.28	104.00
2	B	18	G	O3'-P-O5'	-7.57	92.65	104.00
2	B	22	A	C2'-C3'-O3'	-7.49	102.46	113.70
1	A	521	TRP	N-CA-C	7.08	119.08	111.36
1	A	684	GLY	N-CA-C	7.06	120.32	111.85
1	A	43	TYR	N-CA-C	-6.98	101.00	109.72
1	A	311	HIS	CA-C-N	-6.97	112.72	120.45
1	A	311	HIS	C-N-CA	-6.97	112.72	120.45
2	B	13	G	C3'-C2'-O2'	6.76	120.84	110.70
2	E	8	U	O3'-P-O5'	-6.47	94.29	104.00
1	A	81	ALA	N-CA-C	6.44	121.14	113.28
2	B	50	C	C3'-C2'-O2'	6.43	120.35	110.70
2	B	24	A	C3'-C2'-O2'	-6.37	101.14	110.70
2	E	19	G	O3'-P-O5'	-6.28	94.58	104.00
2	B	4	C	O3'-P-O5'	-6.21	94.69	104.00
1	A	716	ILE	CB-CA-C	-6.01	104.19	112.24
2	B	58	A	O3'-P-O5'	-5.96	95.06	104.00
1	D	854	LEU	N-CA-C	5.92	118.11	108.76
2	E	69	G	O3'-P-O5'	-5.88	95.18	104.00
2	B	21	G	P-O3'-C3'	5.77	128.85	120.20
2	E	64	G	C3'-C2'-O2'	5.76	119.34	110.70
2	B	67	C	C2'-C3'-O3'	-5.75	105.07	113.70
2	E	52	G	C3'-C2'-O2'	5.75	119.33	110.70
2	B	72	U	O3'-P-O5'	-5.70	95.45	104.00
2	B	6	G	C2'-C3'-O3'	-5.57	105.35	113.70
2	B	21	G	C4'-C3'-O3'	-5.56	104.66	113.00
2	B	8	U	O3'-P-O5'	-5.53	95.71	104.00
1	A	453	ILE	N-CA-C	5.52	116.17	108.89
1	A	835	VAL	N-CA-C	-5.50	105.49	110.82
1	D	710	ALA	N-CA-C	-5.48	105.39	111.36
1	A	65	TYR	N-CA-C	-5.45	105.26	111.14
1	A	172	ILE	N-CA-C	5.41	116.46	108.45
2	E	42	C	C3'-C2'-O2'	5.41	118.81	110.70
2	E	44	C	O3'-P-O5'	-5.40	95.91	104.00
1	A	787	TRP	CA-C-N	5.37	125.28	119.76
1	A	787	TRP	C-N-CA	5.37	125.28	119.76
1	D	220	GLN	N-CA-C	5.33	117.08	111.28
1	D	762	ASN	CA-C-N	-5.31	114.20	119.56
1	D	762	ASN	C-N-CA	-5.31	114.20	119.56
2	B	23	C	O3'-P-O5'	-5.23	96.15	104.00
2	E	55	U	C1'-C2'-O2'	5.23	116.25	108.40

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	58	A	C3'-C2'-O2'	-5.23	106.76	114.60
2	B	5	G	C1'-C2'-O2'	5.20	116.19	108.40
2	B	59	G	P-O3'-C3'	5.19	127.99	120.20
2	E	23	C	O3'-P-O5'	-5.15	96.27	104.00
1	D	96	ALA	CA-C-N	5.12	124.73	119.05
1	D	96	ALA	C-N-CA	5.12	124.73	119.05
2	B	18	G	C2'-C3'-O3'	-5.11	101.83	109.50
2	E	63	C	C3'-C2'-O2'	5.10	118.36	110.70
1	D	248	THR	N-CA-C	-5.10	106.84	113.16
1	D	686	VAL	N-CA-C	5.05	117.10	108.86
1	D	216	VAL	CB-CA-C	-5.05	105.51	111.97
2	E	31	A	C4'-C3'-O3'	5.04	120.56	113.00

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	A	6835	0	6687	150	0
1	D	6835	0	6687	132	0
2	B	1675	0	847	12	0
2	E	1632	0	825	14	0
3	A	1	0	0	0	0
3	D	1	0	0	0	0
4	A	23	0	13	0	0
4	D	23	0	13	0	0
5	A	8	0	10	1	0
5	D	8	0	10	0	0
6	B	1	0	0	0	0
6	E	1	0	0	0	0
7	A	104	0	0	12	0
7	B	15	0	0	1	0
7	D	47	0	0	6	0
7	E	3	0	0	1	0
All	All	17212	0	15092	299	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 9.

All (299) 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:288:THR:HG23	1:A:295:MET:HE3	1.23	1.19
1:A:1:MET:HE1	1:A:678:TYR:HA	1.15	1.08
1:D:139:PHE:HE1	1:D:553:MET:HE1	1.34	0.91
1:D:223:TRP:CE3	1:D:535:ILE:HD13	2.04	0.91
1:D:463:ILE:HG22	1:D:464:THR:HG23	1.54	0.90
1:D:548:MET:HA	1:D:553:MET:HE3	1.55	0.86
1:A:1:MET:HE1	1:A:678:TYR:CA	2.04	0.84
1:D:267:LEU:O	1:D:267:LEU:HD23	1.79	0.83
1:A:360:ILE:HD13	1:A:381:LEU:HD23	1.60	0.82
1:D:262:ALA:HB3	1:D:328:MET:HE2	1.60	0.82
1:D:223:TRP:CD2	1:D:535:ILE:HD13	2.14	0.82
1:D:139:PHE:CE1	1:D:553:MET:HE1	2.15	0.81
1:A:633:MET:HE3	1:A:663:VAL:HG11	1.64	0.80
1:D:842:VAL:HG21	1:D:859:VAL:HG22	1.63	0.80
1:D:447:ASP:HA	1:D:450:LEU:HD12	1.64	0.79
1:A:726:ILE:HG22	1:A:730:MET:HE3	1.66	0.78
1:A:729:ILE:HD13	1:A:761:LEU:HD13	1.66	0.78
1:A:288:THR:HG23	1:A:295:MET:CE	2.11	0.77
1:A:288:THR:CG2	1:A:295:MET:HE3	2.11	0.75
1:A:1:MET:HE3	1:A:678:TYR:HD2	1.51	0.75
1:A:694:LEU:HD22	1:A:698:GLN:HB3	1.69	0.74
1:D:262:ALA:HB3	1:D:328:MET:CE	2.16	0.74
1:A:154:SER:HB3	1:A:190:GLN:HE21	1.51	0.74
1:D:98:ALA:HB3	1:D:99:PRO:HD3	1.69	0.73
1:D:535:ILE:HG22	1:D:536:MET:N	2.04	0.73
1:A:1:MET:HE3	1:A:678:TYR:CD2	2.24	0.72
1:D:288:THR:HG22	1:D:298:MET:HE1	1.71	0.72
1:D:223:TRP:CE3	1:D:535:ILE:CD1	2.73	0.71
1:A:813:LYS:CE	2:B:20:U:O2	2.39	0.71
1:A:422:VAL:HG11	1:A:494:MET:HE1	1.72	0.71
1:D:573:ALA:HB3	1:D:615:THR:OG1	1.91	0.69
1:A:267:LEU:HD21	1:A:305:THR:HG21	1.75	0.69
1:A:265:HIS:HB2	1:A:328:MET:HE1	1.74	0.69
1:A:1:MET:CE	1:A:678:TYR:HA	2.09	0.69
1:D:226:ARG:HG3	1:D:415:TYR:CE1	2.28	0.68
1:A:55:ASN:HD22	1:A:55:ASN:C	2.01	0.67
1:D:265:HIS:HB2	1:D:328:MET:HE1	1.77	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:THR:HB	1:A:320:VAL:O	1.95	0.66
1:A:530:GLY:HA3	1:A:568:MET:HE1	1.76	0.66
1:A:161:ASN:OD1	1:A:181:THR:HG21	1.95	0.65
1:A:272:ALA:CA	1:A:278:LEU:HD23	2.27	0.65
1:A:32:LYS:NZ	7:A:2013:HOH:O	2.29	0.65
1:A:813:LYS:HE2	2:B:20:U:O2	1.97	0.64
1:D:842:VAL:HG21	1:D:859:VAL:CG2	2.27	0.64
1:D:71:LYS:O	1:D:73:VAL:HG23	1.98	0.64
1:D:677:VAL:HG11	1:D:772:LEU:HD22	1.79	0.64
1:A:366:SER:HB3	7:A:2058:HOH:O	1.97	0.63
1:A:589:VAL:HG12	1:A:589:VAL:O	1.98	0.63
1:A:816:VAL:O	1:A:817:PRO:C	2.42	0.63
1:D:790:ALA:HB1	1:D:795:MET:HE1	1.79	0.63
1:A:136:GLU:OE2	1:A:495:GLU:OE2	2.16	0.62
1:D:486:GLU:OE1	1:D:486:GLU:HA	1.98	0.62
1:A:536:MET:HA	1:A:536:MET:HE2	1.82	0.62
1:A:358:PRO:HB3	1:A:370:LEU:HD11	1.81	0.62
1:D:633:MET:HE2	1:D:642:VAL:HG22	1.83	0.61
1:A:98:ALA:HB3	1:A:99:PRO:HD3	1.83	0.61
1:D:530:GLY:CA	1:D:568:MET:HE1	2.31	0.61
1:D:790:ALA:HB1	1:D:795:MET:CE	2.31	0.61
1:A:842:VAL:HG21	1:A:859:VAL:HG11	1.82	0.61
1:A:223:TRP:CD2	1:A:535:ILE:HG21	2.35	0.61
1:A:305:THR:HG22	1:A:307:PHE:H	1.65	0.61
1:A:596:ASP:HB2	1:A:600:ARG:O	2.01	0.61
1:D:1:MET:N	7:D:2001:HOH:O	2.32	0.61
1:A:55:ASN:HB2	1:A:566:GLN:NE2	2.17	0.60
1:D:111:GLN:NE2	1:D:629:ASP:OD1	2.34	0.60
1:A:176:CYS:SG	1:A:177:TRP:N	2.75	0.59
1:A:741:THR:N	1:A:742:ASP:OD1	2.35	0.59
1:A:813:LYS:NZ	2:B:20:U:O2	2.35	0.59
1:D:174:GLY:O	1:D:183:VAL:HG21	2.02	0.59
1:D:498:TRP:CE3	1:D:512:MET:HE3	2.38	0.59
1:A:265:HIS:CB	1:A:328:MET:HE1	2.32	0.58
1:A:629:ASP:OD1	1:A:630:PRO:HD2	2.04	0.58
1:D:530:GLY:HA3	1:D:568:MET:HE1	1.85	0.58
1:D:426:ARG:NH2	2:E:74:C:O2	2.35	0.58
1:A:272:ALA:N	1:A:278:LEU:HD23	2.19	0.57
1:A:156:VAL:HG13	1:A:165:VAL:HG13	1.85	0.57
1:D:811:ARG:O	1:D:812:ALA:HB2	2.04	0.57
1:D:805:GLN:OE1	1:D:854:LEU:HD11	2.05	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:817:PRO:O	1:D:818:VAL:C	2.48	0.57
1:D:305:THR:HG23	1:D:320:VAL:O	2.05	0.57
1:A:690:ASN:O	1:A:691:VAL:C	2.47	0.57
1:D:847:VAL:HG22	1:D:857:LEU:HD23	1.87	0.57
1:D:741:THR:HG22	1:D:741:THR:O	2.05	0.56
1:A:801:LEU:C	1:A:801:LEU:HD23	2.30	0.56
1:D:774:GLN:HG2	7:D:2046:HOH:O	2.05	0.56
1:D:114:MET:HE1	1:D:631:GLN:NE2	2.20	0.55
1:D:288:THR:CG2	1:D:298:MET:HE1	2.36	0.55
1:D:165:VAL:HG21	1:D:425:GLN:HG3	1.87	0.55
1:D:241:ASN:C	7:D:2024:HOH:O	2.48	0.55
1:D:288:THR:HG21	1:D:324:ASN:HB2	1.88	0.55
1:A:359:VAL:O	1:A:381:LEU:HD23	2.06	0.55
2:B:18:G:H5"	7:B:2005:HOH:O	2.07	0.55
1:D:435:VAL:HG13	1:D:483:ALA:HB1	1.89	0.55
1:D:531:ILE:HG22	1:D:568:MET:HE3	1.89	0.55
1:D:252:THR:HG22	1:D:360:ILE:HD11	1.88	0.55
1:A:366:SER:OG	1:A:367:GLU:N	2.39	0.55
1:A:98:ALA:HB3	1:A:99:PRO:CD	2.38	0.54
1:A:803:VAL:CG1	1:A:810:VAL:HG22	2.37	0.54
2:E:53:G:N2	2:E:62:C:C2	2.75	0.54
1:A:543:PHE:CZ	1:A:547:LEU:HD11	2.42	0.54
1:A:361:LEU:HD13	1:A:365:GLY:O	2.08	0.54
1:A:84:LEU:HD22	7:A:2023:HOH:O	2.08	0.54
1:D:548:MET:CA	1:D:553:MET:HE3	2.35	0.54
1:A:47:ARG:NH1	1:D:579:GLU:OE2	2.38	0.53
1:D:551:ALA:HB3	1:D:553:MET:HE2	1.90	0.53
1:A:262:ALA:HB3	1:A:328:MET:CE	2.38	0.53
1:A:368:PRO:HB3	1:A:375:LEU:HD22	1.90	0.53
1:A:526:ILE:HD12	1:A:562:GLN:HG2	1.90	0.53
1:D:548:MET:HA	1:D:553:MET:CE	2.32	0.53
1:D:120:ASP:OD1	1:D:122:SER:OG	2.26	0.53
1:D:5:TYR:CE1	1:D:768:ILE:HD12	2.44	0.53
2:E:53:G:C2	2:E:62:C:N3	2.77	0.53
1:A:272:ALA:HA	1:A:278:LEU:HD23	1.91	0.52
1:A:464:THR:HG22	1:A:465:SER:N	2.24	0.52
1:D:267:LEU:HD23	1:D:267:LEU:C	2.33	0.52
1:D:536:MET:HE2	1:D:540:TYR:OH	2.08	0.52
1:D:156:VAL:HG13	1:D:165:VAL:HG13	1.91	0.52
1:D:676:LEU:CD2	1:D:733:MET:HE1	2.39	0.52
1:A:358:PRO:HB3	1:A:370:LEU:CD1	2.38	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:82:PHE:CG	1:D:429:GLY:HA2	2.45	0.52
1:D:140:PHE:CE2	1:D:434:MET:HE3	2.46	0.51
1:A:360:ILE:HD13	1:A:381:LEU:CD2	2.36	0.51
1:A:553:MET:HE1	7:A:2030:HOH:O	2.10	0.51
1:D:369:ASP:OD1	1:D:371:SER:OG	2.28	0.51
1:A:608:ALA:HB1	1:D:93:ASN:OD1	2.11	0.51
2:B:47(D):C:O2	2:B:47(D):C:O4'	2.27	0.51
1:D:4:GLN:HB2	7:D:2002:HOH:O	2.11	0.51
1:A:140:PHE:CE1	1:A:434:MET:HE3	2.46	0.51
1:A:171:VAL:HG22	1:A:183:VAL:HG21	1.92	0.51
1:A:261:VAL:HG13	1:A:265:HIS:HB3	1.91	0.51
1:A:265:HIS:HB2	1:A:328:MET:CE	2.40	0.50
1:D:541:PHE:CZ	1:D:559:PRO:HB2	2.46	0.50
1:A:22:ARG:HD3	7:A:2009:HOH:O	2.10	0.50
1:A:135:TRP:CZ3	1:A:511:GLY:HA2	2.47	0.50
1:A:223:TRP:CZ2	1:A:536:MET:HE3	2.46	0.50
1:A:305:THR:CG2	1:A:307:PHE:HD2	2.24	0.50
1:A:227:SER:OG	1:A:248:THR:CG2	2.59	0.50
1:A:361:LEU:HD23	1:A:382:PHE:CD1	2.46	0.50
1:A:267:LEU:CD2	1:A:305:THR:HG21	2.40	0.50
1:D:215:THR:HG22	2:E:5:G:H5'	1.94	0.50
1:D:514:ASP:C	1:D:514:ASP:OD1	2.53	0.49
1:A:258:TYR:CZ	1:A:337:ALA:CB	2.95	0.49
2:B:10:G:C2	2:B:11:U:C6	3.00	0.49
1:D:1:MET:HE1	1:D:678:TYR:HA	1.93	0.49
1:D:536:MET:O	1:D:537:HIS:C	2.54	0.49
1:D:705:VAL:HG22	1:D:732:LEU:HD11	1.95	0.49
1:D:790:ALA:CB	1:D:795:MET:HE1	2.43	0.49
1:D:140:PHE:HE2	1:D:434:MET:HE3	1.77	0.49
1:D:646:MET:C	1:D:647:MET:HE2	2.38	0.49
1:A:67:ARG:NH1	1:A:784:ASN:OD1	2.46	0.49
1:D:564:LEU:HD22	1:D:720:GLN:HG2	1.94	0.49
1:A:298:MET:HG3	7:A:2056:HOH:O	2.13	0.48
1:A:553:MET:CE	7:A:2030:HOH:O	2.61	0.48
2:E:67:C:H2'	2:E:68:C:O4'	2.13	0.48
1:A:339:PRO:HB2	1:A:356:ILE:HG23	1.95	0.48
2:B:47(G):G:O2'	2:B:47(H):C:H5'	2.13	0.48
1:D:818:VAL:HG13	1:D:819:ASP:H	1.78	0.48
1:A:140:PHE:HE1	1:A:434:MET:HE3	1.79	0.48
1:A:223:TRP:CG	1:A:535:ILE:HG21	2.48	0.48
1:A:842:VAL:HB	1:A:859:VAL:HG13	1.95	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:LYS:O	1:A:405:ALA:HB3	2.13	0.48
1:D:443:MET:HB2	1:D:444:PRO:HD2	1.94	0.48
2:E:55:U:H2'	2:E:57:A:OP2	2.13	0.48
1:A:773:TRP:HB2	1:A:782:ILE:HD12	1.94	0.48
1:D:257:THR:HG22	1:D:357:LYS:HD3	1.95	0.48
1:D:855:LEU:C	1:D:855:LEU:HD12	2.39	0.48
1:A:265:HIS:CD2	1:A:328:MET:HE1	2.49	0.48
2:B:66:U:H2'	2:B:67:C:O4'	2.13	0.48
1:D:589:VAL:O	1:D:589:VAL:CG1	2.62	0.48
1:D:651:PRO:HG3	2:E:14:A:OP1	2.14	0.47
1:D:65:TYR:CZ	1:D:69:LEU:HD11	2.49	0.47
1:D:673:VAL:HG22	1:D:733:MET:HE2	1.95	0.47
2:E:9:G:N3	2:E:9:G:H2'	2.29	0.47
1:D:798:ASP:OD1	1:D:798:ASP:N	2.34	0.47
1:A:135:TRP:CH2	1:A:511:GLY:HA2	2.50	0.47
1:A:633:MET:HE2	1:A:642:VAL:HG22	1.95	0.47
2:B:43:U:H2'	2:B:44:C:O4'	2.15	0.47
1:D:823:GLU:OE1	1:D:823:GLU:HA	2.14	0.47
1:D:358:PRO:HB3	1:D:370:LEU:HD11	1.96	0.47
1:D:855:LEU:HD12	1:D:856:ASN:N	2.30	0.47
1:D:40:MET:HE3	1:D:529:GLY:HA3	1.97	0.47
1:A:530:GLY:CA	1:A:568:MET:HE1	2.44	0.46
1:A:57:THR:O	1:A:61:VAL:HG23	2.14	0.46
1:D:176:CYS:SG	1:D:177:TRP:N	2.87	0.46
1:A:842:VAL:HG21	1:A:859:VAL:CG1	2.46	0.46
2:E:26:A:H2'	2:E:27:A:O4'	2.14	0.46
1:A:55:ASN:HB2	1:A:566:GLN:HE22	1.80	0.46
1:A:569:VAL:HG12	1:A:620:MET:HE3	1.97	0.46
1:D:174:GLY:O	1:D:183:VAL:CG2	2.63	0.46
1:A:647:MET:N	1:A:647:MET:HE2	2.30	0.46
1:A:271:ALA:HB2	1:A:307:PHE:CE2	2.50	0.46
1:A:55:ASN:C	1:A:55:ASN:ND2	2.71	0.46
1:D:676:LEU:HD22	1:D:733:MET:HE1	1.98	0.46
1:A:34:LYS:NZ	1:A:525:ASP:OD2	2.45	0.45
1:D:439:ASP:HB3	1:D:441:THR:H	1.81	0.45
1:A:59:GLY:HA2	1:A:528:ILE:HD13	1.96	0.45
1:D:463:ILE:HG22	1:D:464:THR:N	2.31	0.45
1:D:22:ARG:NH1	1:D:25:GLU:OE2	2.49	0.45
2:B:31:A:H2'	2:B:32:U:O5'	2.16	0.45
1:D:79:TRP:O	1:D:127:THR:HG23	2.16	0.45
1:A:136:GLU:OE1	1:A:495:GLU:HG3	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:312:PRO:O	1:D:406:MET:HE1	2.17	0.45
1:A:95:THR:HG22	1:A:96:ALA:H	1.82	0.45
1:D:821:THR:HG23	1:D:824:GLN:OE1	2.17	0.45
1:D:114:MET:HE1	1:D:631:GLN:CD	2.41	0.45
1:D:842:VAL:CG2	1:D:859:VAL:HG22	2.42	0.45
1:A:231:GLU:O	1:A:409:GLY:HA2	2.18	0.45
1:A:614:TYR:CZ	1:A:616:GLY:HA2	2.52	0.44
1:A:290:VAL:O	1:A:291:ALA:HB2	2.18	0.44
1:D:136:GLU:OE1	1:D:495:GLU:HG3	2.18	0.44
1:D:519:ASN:OD1	1:D:555:ASN:HB2	2.18	0.44
1:D:241:ASN:HB2	7:D:2024:HOH:O	2.17	0.44
1:D:47:ARG:HE	1:D:47:ARG:HB3	1.67	0.44
1:D:439:ASP:CB	1:D:441:THR:H	2.31	0.44
1:A:41:LEU:HG	1:A:80:ASP:HB2	2.00	0.43
1:A:545:HIS:CE1	1:A:556:SER:HG	2.36	0.43
1:A:743:GLY:HA3	7:A:2097:HOH:O	2.18	0.43
1:D:236:VAL:HB	1:D:239:TYR:HB3	2.00	0.43
1:A:536:MET:HB2	7:A:2075:HOH:O	2.19	0.43
1:A:589:VAL:O	1:A:589:VAL:CG1	2.65	0.43
1:A:617:MET:HE1	7:A:2077:HOH:O	2.18	0.43
1:D:83:GLY:HA2	1:D:492:THR:HG21	1.99	0.43
1:D:787:TRP:CD2	1:D:788:PRO:HD2	2.53	0.43
1:A:821:THR:HG23	1:A:824:GLN:HG3	1.99	0.43
1:A:823:GLU:N	1:A:823:GLU:CD	2.76	0.43
1:D:447:ASP:HA	1:D:450:LEU:CD1	2.40	0.43
1:A:361:LEU:HD23	1:A:382:PHE:CE1	2.54	0.43
1:A:646:MET:C	1:A:647:MET:HE2	2.43	0.43
1:D:391:ASP:C	1:D:391:ASP:OD1	2.61	0.43
1:A:312:PRO:O	1:A:406:MET:HE1	2.19	0.43
1:A:396:PHE:CD1	1:A:396:PHE:C	2.96	0.43
2:E:1:G:C6	2:E:2:C:C4	3.06	0.43
1:A:531:ILE:HG22	1:A:568:MET:HE3	2.01	0.43
1:D:420:TRP:CZ2	1:D:494:MET:HE2	2.53	0.43
1:A:391:ASP:OD1	1:A:391:ASP:C	2.62	0.43
1:A:529:GLY:O	1:A:565:CYS:HA	2.18	0.43
1:A:726:ILE:CG2	1:A:730:MET:HE3	2.44	0.43
1:A:733:MET:HE3	7:A:2096:HOH:O	2.18	0.43
1:A:817:PRO:O	1:A:818:VAL:C	2.62	0.43
1:D:158:TRP:CZ2	1:D:163:GLN:HA	2.54	0.43
1:A:6:ARG:O	1:A:10:ILE:HD12	2.18	0.42
1:A:342:ASP:OD1	1:A:342:ASP:C	2.62	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:361:LEU:CD2	1:A:382:PHE:CD1	3.02	0.42
1:D:136:GLU:OE2	1:D:495:GLU:OE2	2.36	0.42
1:A:359:VAL:O	1:A:381:LEU:CD2	2.68	0.42
1:A:823:GLU:CD	1:A:823:GLU:H	2.26	0.42
1:A:292:GLU:O	1:A:293:ALA:C	2.62	0.42
1:D:265:HIS:N	1:D:328:MET:HE3	2.34	0.42
1:A:801:LEU:HD23	1:A:802:VAL:N	2.35	0.42
1:D:261:VAL:HG22	1:D:265:HIS:CG	2.55	0.42
1:D:633:MET:CE	1:D:642:VAL:HG13	2.50	0.42
1:A:443:MET:HB2	1:A:444:PRO:CD	2.50	0.42
2:B:31:A:C2'	2:B:32:U:O5'	2.68	0.42
1:A:195:ILE:HG22	1:A:419:ASP:HA	2.01	0.42
1:A:327:LEU:CD1	1:A:330:TYR:HE2	2.33	0.42
1:A:336:MET:HE3	1:A:338:VAL:HG23	2.02	0.42
1:D:828:ARG:O	1:D:828:ARG:HD3	2.20	0.42
1:A:729:ILE:CD1	1:A:761:LEU:HD13	2.42	0.42
1:D:434:MET:HE1	1:D:444:PRO:HB3	2.02	0.42
1:A:80:ASP:OD2	5:A:902:LEU:N	2.52	0.41
1:D:6:ARG:HA	1:D:7:PRO:HD3	1.91	0.41
1:A:494:MET:HB2	1:A:494:MET:HE2	1.76	0.41
1:D:86:ALA:HA	1:D:100:TRP:CZ3	2.55	0.41
1:A:485:ARG:HB2	7:A:2066:HOH:O	2.18	0.41
2:E:57:A:P	7:E:2003:HOH:O	2.78	0.41
1:D:44:PRO:HA	1:D:108:MET:SD	2.61	0.41
1:D:82:PHE:CE2	1:D:128:CYS:HA	2.56	0.41
1:D:219:MET:HE3	1:D:534:ALA:HB1	2.01	0.41
1:D:290:VAL:N	7:D:2018:HOH:O	2.29	0.41
1:D:476:THR:OG1	1:D:477:THR:N	2.54	0.41
1:A:328:MET:HE2	1:A:328:MET:HA	2.03	0.41
1:A:601:ILE:HD12	1:A:614:TYR:CE1	2.55	0.41
1:A:729:ILE:CD1	1:A:761:LEU:CD1	2.98	0.41
1:D:26:VAL:HG11	1:D:73:VAL:O	2.21	0.41
1:D:531:ILE:O	1:D:531:ILE:HG13	2.21	0.41
1:D:854:LEU:HD23	1:D:854:LEU:C	2.46	0.41
1:A:533:HIS:O	1:A:534:ALA:C	2.63	0.41
2:B:10:G:C6	2:B:26:A:C2	3.09	0.41
1:A:35:TYR:O	1:A:73:VAL:HA	2.21	0.41
1:A:787:TRP:CD2	1:A:788:PRO:HD2	2.56	0.41
1:D:498:TRP:CD2	1:D:512:MET:HE3	2.56	0.41
1:D:617:MET:HE3	1:D:657:GLU:OE1	2.21	0.41
1:D:647:MET:HE2	1:D:647:MET:N	2.36	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:7:A:H3'	2:E:8:U:H5'	2.03	0.41
2:E:7:A:C3'	2:E:8:U:H5'	2.51	0.41
1:A:54:ARG:HB2	1:A:646:MET:HE2	2.03	0.41
1:A:98:ALA:N	1:A:99:PRO:HD2	2.36	0.41
1:A:538:LEU:HD23	1:A:538:LEU:HA	1.87	0.41
1:A:647:MET:HE2	1:A:647:MET:HA	2.03	0.41
1:A:676:LEU:CD2	1:A:733:MET:HE1	2.51	0.41
1:D:84:LEU:HD23	1:D:426:ARG:HD3	2.03	0.41
2:E:8:U:H2'	2:E:21:G:N2	2.36	0.41
1:A:608:ALA:HB2	1:D:93:ASN:HD21	1.86	0.40
1:D:629:ASP:OD1	1:D:630:PRO:HD2	2.21	0.40
1:A:258:TYR:CE1	1:A:337:ALA:CB	3.04	0.40
1:A:618:SER:O	1:A:619:LYS:C	2.62	0.40
1:D:42:PRO:HD2	1:D:78:GLY:O	2.20	0.40
1:A:84:LEU:N	1:A:85:PRO:CD	2.84	0.40
1:D:25:GLU:HA	1:D:120:ASP:CG	2.46	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	
1	A	858/880 (98%)	808 (94%)	44 (5%)	6 (1%)	18	34
1	D	858/880 (98%)	802 (94%)	47 (6%)	9 (1%)	12	24
All	All	1716/1760 (98%)	1610 (94%)	91 (5%)	15 (1%)	14	27

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	463	ILE
1	D	463	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	293	ALA
1	A	536	MET
1	D	240	ASP
1	D	460	MET
1	D	818	VAL
1	A	291	ALA
1	D	93	ASN
1	D	536	MET
1	A	298	MET
1	D	94	ASN
1	D	276	PRO
1	A	851	PRO
1	D	535	ILE

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	
1	A	724/741 (98%)	671 (93%)	53 (7%)	13	27
1	D	724/741 (98%)	651 (90%)	73 (10%)	7	15
All	All	1448/1482 (98%)	1322 (91%)	126 (9%)	9	21

All (126) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	10	ILE
1	A	13	LYS
1	A	47	ARG
1	A	55	ASN
1	A	84	LEU
1	A	95	THR
1	A	136	GLU
1	A	143	LEU
1	A	156	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	169	GLU
1	A	181	THR
1	A	188	ILE
1	A	243	LEU
1	A	248	THR
1	A	261	VAL
1	A	288	THR
1	A	297	THR
1	A	298	MET
1	A	299	GLU
1	A	305	THR
1	A	330	TYR
1	A	342	ASP
1	A	376	THR
1	A	393	GLU
1	A	412	LYS
1	A	416	ARG
1	A	425	GLN
1	A	435	VAL
1	A	453	ILE
1	A	477	THR
1	A	516	GLU
1	A	526	ILE
1	A	536	MET
1	A	582	GLU
1	A	592	ILE
1	A	602	VAL
1	A	636	ARG
1	A	650	SER
1	A	655	THR
1	A	661	SER
1	A	663	VAL
1	A	664	GLU
1	A	683	LYS
1	A	686	VAL
1	A	695	THR
1	A	731	GLU
1	A	761	LEU
1	A	779	GLU
1	A	821	THR
1	A	827	GLU
1	A	848	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	859	VAL
1	D	1	MET
1	D	47	ARG
1	D	48	LEU
1	D	55	ASN
1	D	92	LYS
1	D	143	LEU
1	D	153	THR
1	D	156	VAL
1	D	159	CYS
1	D	164	THR
1	D	183	VAL
1	D	208	LYS
1	D	238	ASP
1	D	243	LEU
1	D	261	VAL
1	D	269	GLN
1	D	295	MET
1	D	298	MET
1	D	344	ARG
1	D	357	LYS
1	D	371	SER
1	D	376	THR
1	D	381	LEU
1	D	416	ARG
1	D	425	GLN
1	D	432	ILE
1	D	435	VAL
1	D	448	ASP
1	D	458	VAL
1	D	461	ASP
1	D	475	LYS
1	D	481	MET
1	D	485	ARG
1	D	486	GLU
1	D	496	SER
1	D	510	GLU
1	D	514	ASP
1	D	516	GLU
1	D	544	PHE
1	D	565	CYS
1	D	593	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	618	SER
1	D	622	LYS
1	D	628	ILE
1	D	646	MET
1	D	650	SER
1	D	654	MET
1	D	655	THR
1	D	659	GLN
1	D	685	ASP
1	D	686	VAL
1	D	692	ASP
1	D	695	THR
1	D	745	GLN
1	D	761	LEU
1	D	793	LYS
1	D	798	ASP
1	D	801	LEU
1	D	815	THR
1	D	816	VAL
1	D	818	VAL
1	D	821	THR
1	D	823	GLU
1	D	824	GLN
1	D	827	GLU
1	D	837	LYS
1	D	839	LEU
1	D	842	VAL
1	D	843	THR
1	D	845	ARG
1	D	854	LEU
1	D	855	LEU
1	D	859	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	211	HIS
1	A	269	GLN
1	A	762	ASN
1	A	774	GLN
1	A	805	GLN
1	A	824	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	372	GLN
1	D	383	ASN
1	D	533	HIS
1	D	631	GLN
1	D	784	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	75/87 (86%)	10 (13%)	0
2	E	73/87 (83%)	17 (23%)	1 (1%)
All	All	148/174 (85%)	27 (18%)	1 (0%)

All (27) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	9	G
2	B	10	G
2	B	11	U
2	B	16	U
2	B	18	G
2	B	28	G
2	B	41	C
2	B	42	C
2	B	47(A)	G
2	B	47(F)	C
2	E	8	U
2	E	9	G
2	E	13	G
2	E	16	U
2	E	18	G
2	E	31	A
2	E	45	G
2	E	47(F)	C
2	E	48	U
2	E	49	G
2	E	51	G
2	E	52	G
2	E	56	C
2	E	61	C
2	E	63	C

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	E	65	C
2	E	75	C

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	E	64	G

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 [i](#)

Of 8 ligands modelled in this entry, 4 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	$\# Z > 2$
4	LMS	D	901	5	25,25,25	1.94	5 (20%)	38,38,38	2.44	11 (28%)
4	LMS	A	901	5	25,25,25	2.16	6 (24%)	38,38,38	2.34	10 (26%)
5	LEU	D	902	4	5,7,8	0.45	0	6,8,10	1.10	1 (16%)
5	LEU	A	902	4	5,7,8	0.97	0	6,8,10	0.44	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	LMS	D	901	5	-	1/10/26/26	0/3/3/3
4	LMS	A	901	5	-	0/10/26/26	0/3/3/3
5	LEU	D	902	4	-	0/5/6/8	-
5	LEU	A	902	4	-	0/5/6/8	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	LMS	O2P-S	6.42	1.49	1.42
4	D	901	LMS	O1P-S	5.75	1.48	1.42
4	D	901	LMS	S-N	-4.32	1.53	1.58
4	A	901	LMS	C4-N9	-4.23	1.28	1.37
4	A	901	LMS	C5-N7	-4.14	1.31	1.39
4	D	901	LMS	O2P-S	3.81	1.46	1.42
4	A	901	LMS	O1P-S	3.19	1.45	1.42
4	D	901	LMS	O5'-S	-3.07	1.53	1.57
4	A	901	LMS	S-N	-3.04	1.55	1.58
4	D	901	LMS	C8-N7	2.40	1.36	1.31
4	A	901	LMS	O5'-S	-2.21	1.54	1.57

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	901	LMS	O2P-S-O1P	-8.71	111.07	119.94
4	A	901	LMS	O2P-S-O1P	-6.69	113.13	119.94
4	A	901	LMS	C5'-O5'-S	5.88	124.81	117.29
4	A	901	LMS	N3-C2-N1	-5.06	120.93	128.58
4	D	901	LMS	C5-C4-N3	-4.99	119.84	126.72
4	A	901	LMS	C5-C4-N3	-4.84	120.05	126.72
4	D	901	LMS	O1P-S-N	-4.71	102.32	109.13
4	D	901	LMS	N3-C2-N1	-3.71	122.97	128.58
4	D	901	LMS	N9-C8-N7	-3.50	108.97	113.94
4	A	901	LMS	O2P-S-N	3.47	114.15	109.13
4	A	901	LMS	C2-N3-C4	3.44	120.23	111.83
4	D	901	LMS	O2P-S-N	3.11	113.63	109.13
4	D	901	LMS	C2-N3-C4	3.08	119.36	111.83
4	D	901	LMS	C4-C5-N7	-2.99	107.17	110.58
4	D	901	LMS	N3-C4-N9	2.93	132.15	127.17
4	D	901	LMS	C4-N9-C8	2.79	108.67	105.74
4	A	901	LMS	N3-C4-N9	2.59	131.57	127.17
5	D	902	LEU	CB-CA-C	2.43	114.74	110.99
4	D	901	LMS	C5-N7-C8	2.39	107.21	103.45
4	A	901	LMS	C5-C4-N9	2.32	108.34	105.81

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	LMS	N9-C8-N7	-2.29	110.68	113.94
4	A	901	LMS	O3'-C3'-C2'	-2.17	104.86	111.82

There are no chirality outliers.

All (1) torsion outliers are listed below:

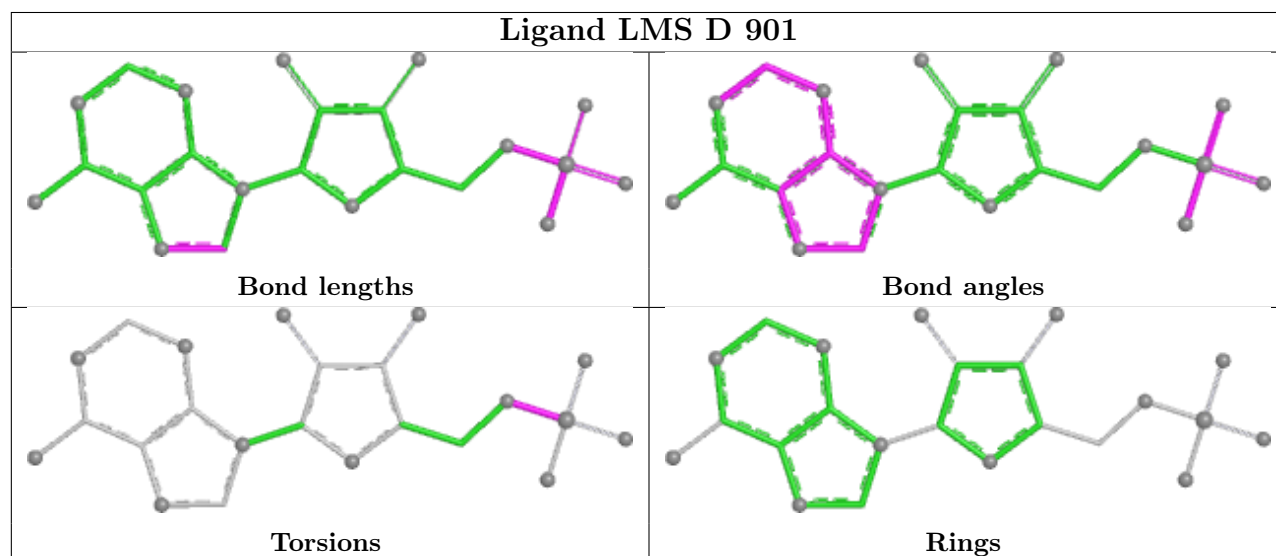
Mol	Chain	Res	Type	Atoms
4	D	901	LMS	C5'-O5'-S-N

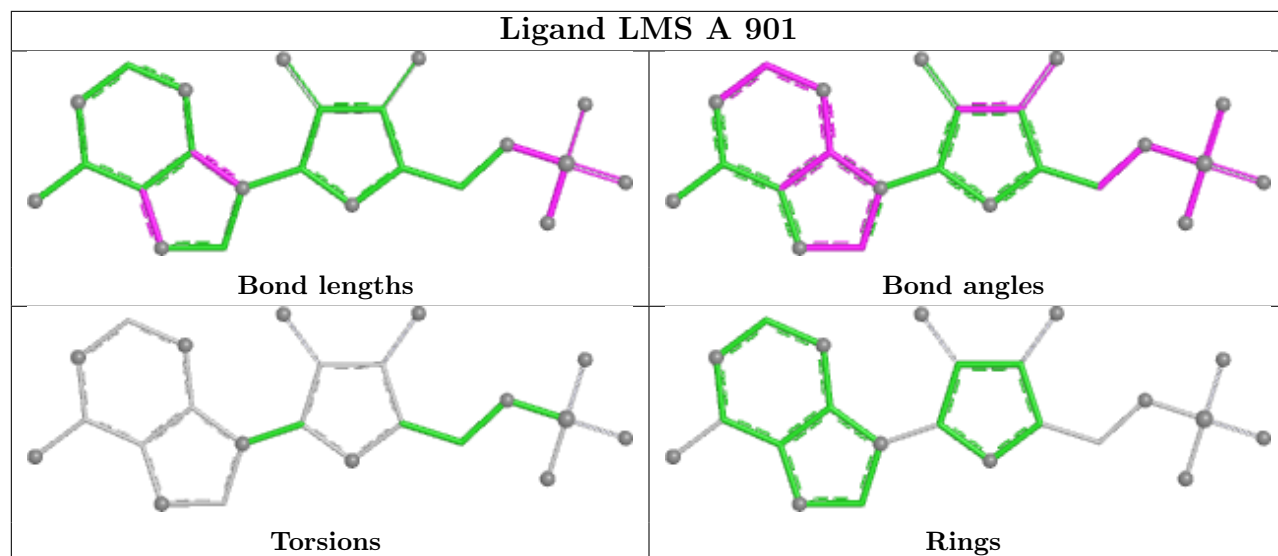
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	902	LEU	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	860/880 (97%)	-0.32	3 (0%) 90 87	20, 43, 80, 132	0
1	D	860/880 (97%)	-0.07	2 (0%) 91 89	33, 58, 102, 165	0
2	B	79/87 (90%)	-0.70	0 100 100	24, 45, 98, 122	0
2	E	77/87 (88%)	-0.21	1 (1%) 75 71	40, 72, 111, 141	0
All	All	1876/1934 (97%)	-0.22	6 (0%) 90 87	20, 51, 99, 165	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	177	TRP	2.8
1	A	860	GLY	2.7
1	D	165	VAL	2.2
1	A	176	CYS	2.2
1	D	602	VAL	2.1
2	E	47	C	2.1

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.

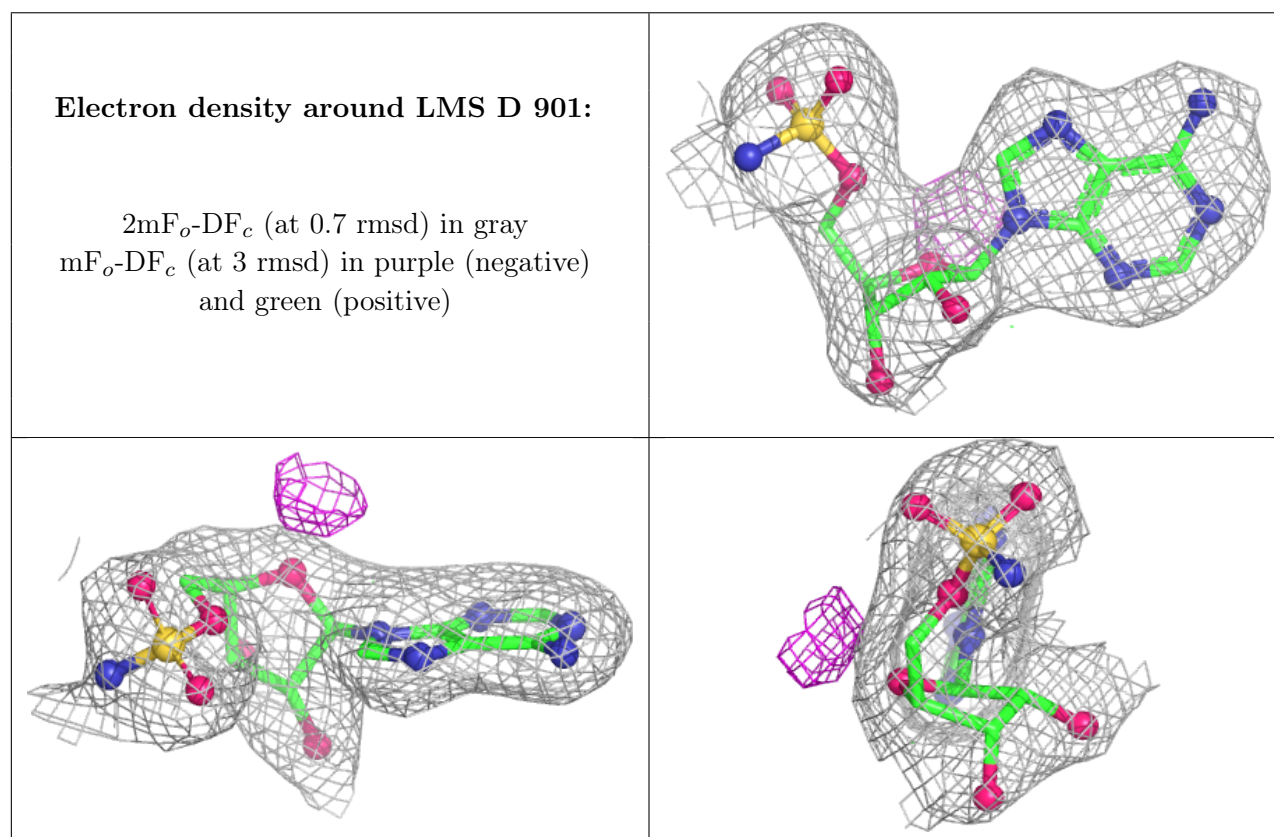
6.4 Ligands [i](#)

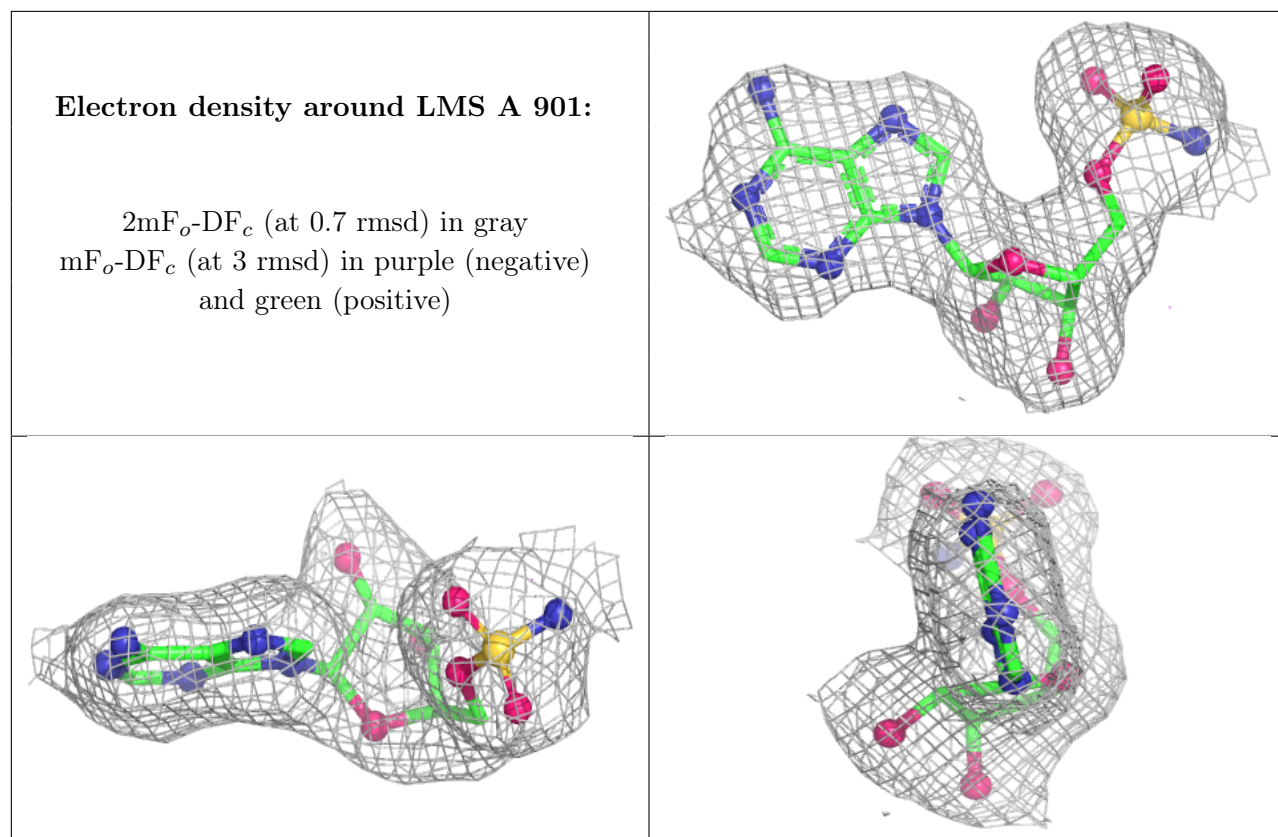
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

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
6	MG	B	1077	1/1	0.87	0.28	55,55,55,55	0
6	MG	E	1077	1/1	0.90	0.24	67,67,67,67	0
3	ZN	A	900	1/1	0.92	0.08	104,104,104,104	0
5	LEU	D	902	8/9	0.95	0.09	31,37,40,43	0
3	ZN	D	900	1/1	0.96	0.05	105,105,105,105	0
5	LEU	A	902	8/9	0.97	0.06	23,25,27,27	0
4	LMS	D	901	23/23	0.97	0.06	32,34,40,42	0
4	LMS	A	901	23/23	0.99	0.04	20,22,25,28	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers [i](#)

There are no such residues in this entry.