



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 27, 2026 – 11:48 am BST

PDB ID : 9TWV / pdb_00009twv
Title : Complex of E. coli leucyl-tRNA synthetase AN3365-resistant mutant D345A bound to tRNA(leu)
Authors : Hoffmann, G.; Palencia, A.; Muller, B.
Deposited on : 2026-01-15
Resolution : 2.11 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	:	1.8.4, CSD as541be (2020)
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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

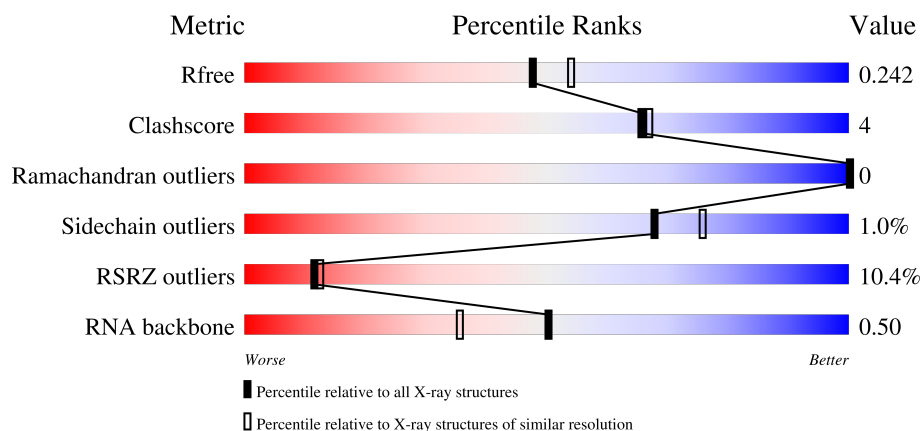
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 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	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)
RNA backbone	3983	1037 (2.46-1.78)

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	
2	B	87	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8702 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	3	0
			6479	4129	1097	1215	38			

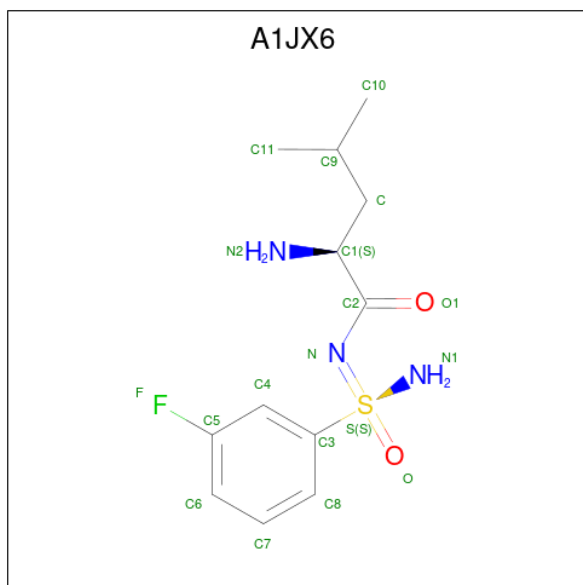
There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	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
A	345	ALA	ASP	engineered mutation	UNP P07813

- Molecule 2 is a RNA chain called tRNA(leu).

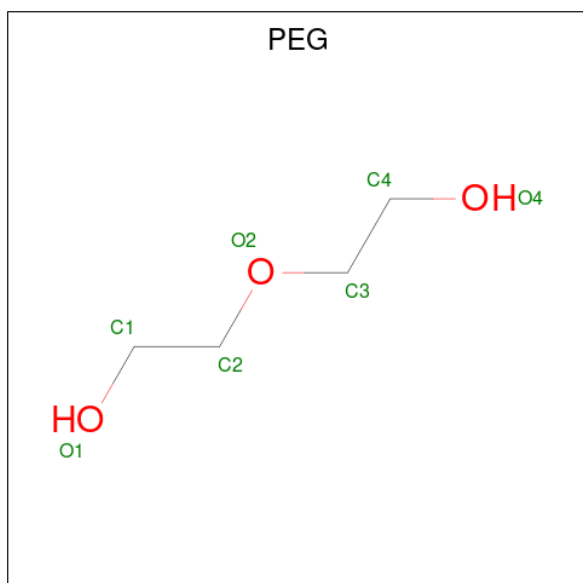
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	81	Total	C	N	O	P	0	0	0
			1731	770	309	571	81			

- Molecule 3 is (2 {S})-2-azanyl- {N}-[azanyl-(3-fluorophenyl)-oxidanylidene- λ^6 -sulfa nylidene]-4-methyl-pentanamide (CCD ID: A1JX6) (formula: $C_{12}H_{18}FN_3O_2S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	S	0	0
			19	12	1	3	2	1		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

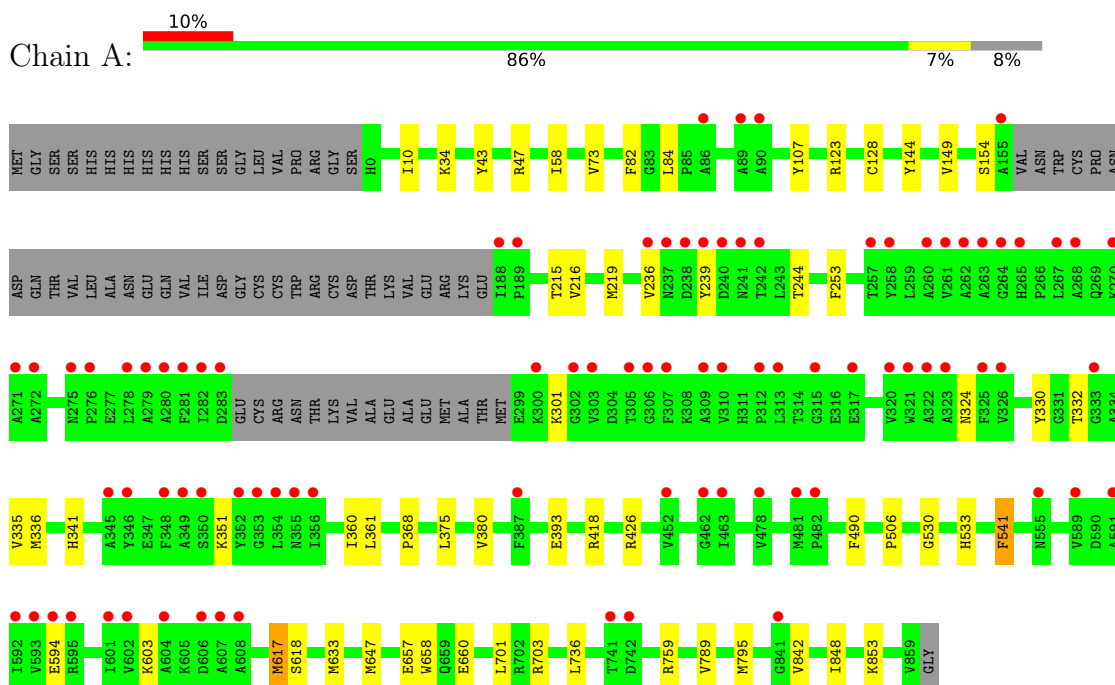
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	388	Total	O	0	0
			388	388		
6	B	77	Total	O	0	0
			77	77		

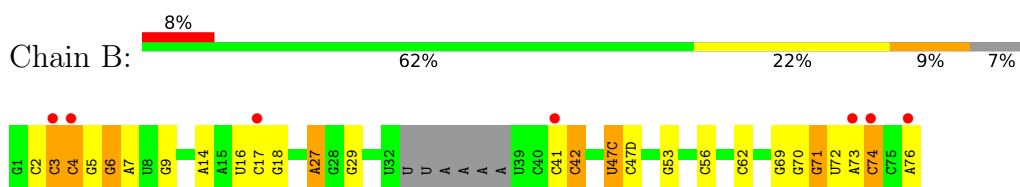
3 Residue-property plots [i](#)

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 2: tRNA(Leu)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.80Å 118.95Å 140.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.56 – 2.11 60.56 – 2.11	Depositor EDS
% Data completeness (in resolution range)	88.5 (60.56-2.11) 88.5 (60.56-2.11)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 2.10Å)	Xtriage
Refinement program	BUSTER 2.10.4	Depositor
R, R_{free}	0.215 , 0.252 0.206 , 0.242	Depositor DCC
R_{free} test set	3254 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	47.5	Xtriage
Anisotropy	0.007	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8702	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, A1JX6, PEG

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	0.71	1/6641 (0.0%)	0.99	7/9015 (0.1%)
2	B	0.57	0/1932	0.68	0/3007
All	All	0.68	1/8573 (0.0%)	0.92	7/12022 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	617	MET	SD-CE	-13.01	1.47	1.79

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	43	TYR	N-CA-C	-5.55	103.04	109.93
1	A	10	ILE	N-CA-CB	-5.48	107.15	111.64
1	A	73	VAL	N-CA-C	5.33	115.98	108.36
1	A	490	PHE	CA-C-N	5.28	128.26	120.82
1	A	490	PHE	C-N-CA	5.28	128.26	120.82
1	A	541	PHE	CA-CB-CG	5.24	119.04	113.80
1	A	530	GLY	N-CA-C	5.09	117.79	111.93

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	6479	0	6347	34	0
2	B	1731	0	877	28	0
3	A	19	0	0	0	0
4	A	7	0	10	3	0
5	B	1	0	0	0	0
6	A	388	0	0	1	0
6	B	77	0	0	0	0
All	All	8702	0	7234	58	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 4.

All (58) 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:5:G:H2'	2:B:6:G:H5''	1.43	0.98
1:A:330:TYR:HD1	1:A:418:ARG:HH12	1.25	0.85
1:A:84:LEU:HD21	1:A:426:ARG:HE	1.47	0.80
1:A:330:TYR:HD1	1:A:418:ARG:NH1	1.82	0.78
1:A:215:THR:HG22	1:A:219:MET:HE2	1.65	0.77
2:B:3:C:H42	2:B:70:G:H1	1.35	0.72
2:B:5:G:C2'	2:B:6:G:H5''	2.20	0.71
1:A:47:ARG:NH2	1:A:107:TYR:OH	2.27	0.68
2:B:5:G:H2'	2:B:6:G:C5'	2.23	0.64
1:A:236:VAL:HB	1:A:239:TYR:HB3	1.82	0.61
2:B:4:C:H42	2:B:69:G:H1	1.46	0.60
2:B:3:C:H42	2:B:70:G:H22	1.49	0.60
2:B:3:C:N4	2:B:70:G:H22	2.00	0.60
2:B:47(C):U:H5'	2:B:47(D):C:H5	1.67	0.59
1:A:703:ARG:HG3	1:A:795:MET:HA	1.85	0.58
1:A:633:MET:HE2	1:A:658:TRP:HZ2	1.69	0.58
1:A:759:ARG:HH12	4:A:902:PEG:H11	1.67	0.58
1:A:324:ASN:ND2	2:B:74:C:N4	2.54	0.56
2:B:5:G:C6	2:B:6:G:N7	2.76	0.54
2:B:5:G:N2	2:B:69:G:C2	2.76	0.53
2:B:71:G:N7	2:B:72:U:C4	2.78	0.52
2:B:3:C:H42	2:B:70:G:N2	2.08	0.51
2:B:47(C):U:H5'	2:B:47(D):C:C5	2.44	0.51
1:A:244:THR:HG22	1:A:332:THR:HG21	1.92	0.51
2:B:5:G:C4	2:B:6:G:C8	2.99	0.50
1:A:324:ASN:ND2	2:B:74:C:H42	2.09	0.50
2:B:41:C:H2'	2:B:41:C:O2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:27:A:OP2	2:B:27:A:H8	1.95	0.49
1:A:301:LYS:HA	1:A:351:LYS:NZ	2.28	0.48
1:A:123:ARG:HA	1:A:506:PRO:HG3	1.95	0.48
1:A:144:TYR:HA	1:A:149:VAL:HB	1.94	0.48
1:A:789:VAL:H	4:A:902:PEG:H22	1.78	0.48
2:B:5:G:C2	2:B:69:G:C2	3.01	0.48
2:B:5:G:C2	2:B:69:G:N2	2.81	0.48
1:A:330:TYR:CD1	1:A:418:ARG:NH1	2.73	0.48
1:A:533:HIS:HE1	6:A:1273:HOH:O	1.96	0.48
2:B:69:G:H2'	2:B:70:G:C8	2.49	0.47
1:A:84:LEU:HD21	1:A:426:ARG:NE	2.24	0.47
1:A:253:PHE:HD1	1:A:336:MET:SD	2.37	0.47
1:A:594:GLU:HB2	1:A:603:LYS:HG3	1.96	0.47
1:A:617:MET:HG3	1:A:657:GLU:CD	2.40	0.47
1:A:341:HIS:CE1	1:A:360:ILE:HB	2.50	0.46
2:B:5:G:C5	2:B:6:G:C8	3.04	0.46
1:A:617:MET:HE2	1:A:657:GLU:HG2	1.96	0.45
1:A:848:ILE:HG13	2:B:56:C:C2	2.51	0.45
2:B:53:G:N2	2:B:62:C:C2	2.85	0.45
2:B:29:G:N2	2:B:42:C:C2	2.85	0.44
1:A:58:ILE:HG13	1:A:647:MET:HE1	1.98	0.44
1:A:701:LEU:HD11	1:A:736:LEU:HD23	1.99	0.44
2:B:4:C:N4	2:B:69:G:H1	2.15	0.44
1:A:335:VAL:HG12	2:B:76:A:N7	2.34	0.43
1:A:368:PRO:HB3	1:A:375:LEU:HD22	2.01	0.42
1:A:216:VAL:HA	1:A:219:MET:HE3	2.02	0.42
2:B:5:G:C2	2:B:6:G:C8	3.08	0.41
1:A:361:LEU:HD12	1:A:380:VAL:HB	2.01	0.41
1:A:759:ARG:HH22	4:A:902:PEG:H32	1.85	0.41
1:A:853:LYS:HB3	1:A:853:LYS:HE2	1.94	0.40
1:A:82:PHE:CE2	1:A:128:CYS:HA	2.57	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	810/880 (92%)	788 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	683/740 (92%)	676 (99%)	7 (1%)	68	76

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LYS
1	A	154	SER
1	A	393	GLU
1	A	541	PHE
1	A	618	SER
1	A	660	GLU
1	A	842	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	449	GLN
1	A	762	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	79/87 (90%)	15 (18%)	3 (3%)

All (15) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	2	C
2	B	3	C
2	B	4	C
2	B	6	G
2	B	7	A
2	B	9	G
2	B	14	A
2	B	16	U
2	B	17	C
2	B	18	G
2	B	27	A
2	B	47(C)	U
2	B	71	G
2	B	73	A
2	B	74	C

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	6	G
2	B	17	C
2	B	42	C

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	A1JX6	A	901	-	15,19,19	0.55	0	20,27,27	1.16	1 (5%)
4	PEG	A	902	-	6,6,6	0.16	0	5,5,5	0.18	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	A1JX6	A	901	-	-	4/12/19/19	0/1/1/1
4	PEG	A	902	-	-	4/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	A1JX6	C3-S-N1	-4.52	101.99	108.40

There are no chirality outliers.

All (8) torsion outliers are listed below:

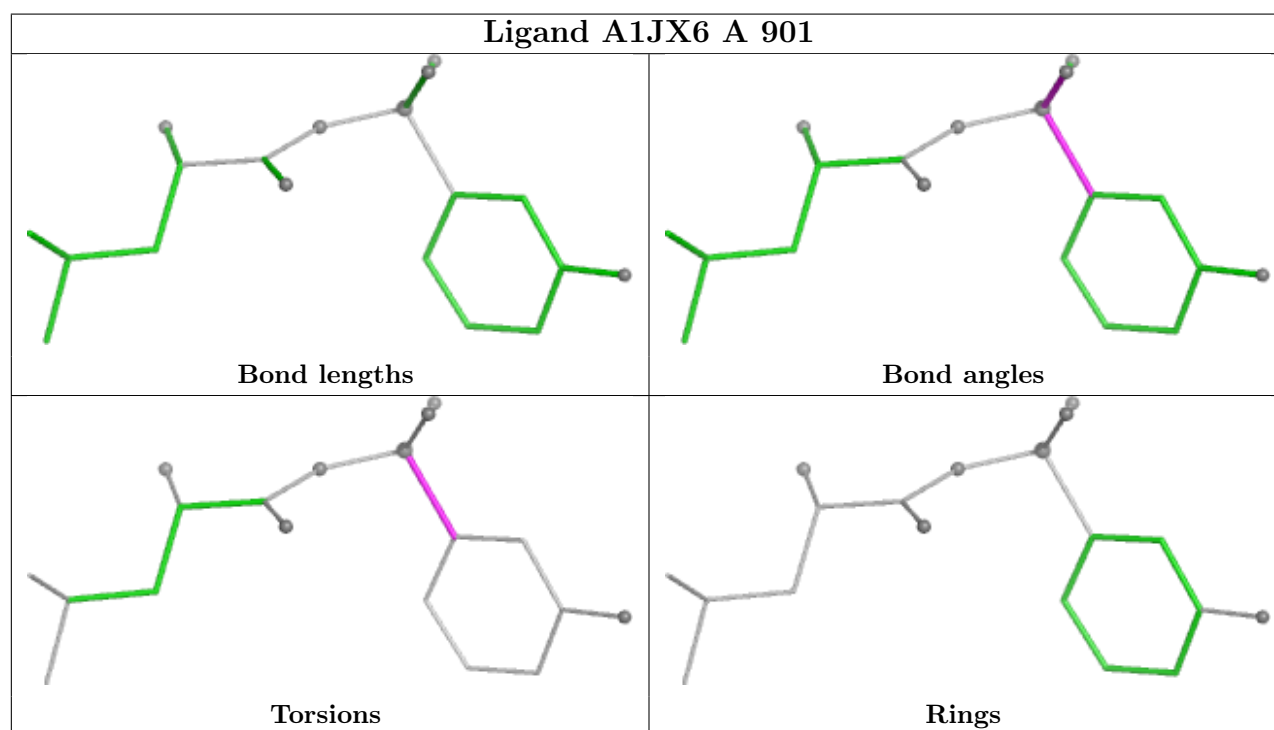
Mol	Chain	Res	Type	Atoms
4	A	902	PEG	O2-C3-C4-O4
3	A	901	A1JX6	C4-C3-S-N1
3	A	901	A1JX6	C8-C3-S-N1
4	A	902	PEG	C1-C2-O2-C3
4	A	902	PEG	O1-C1-C2-O2
4	A	902	PEG	C4-C3-O2-C2
3	A	901	A1JX6	C4-C3-S-O
3	A	901	A1JX6	C8-C3-S-O

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	902	PEG	3	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

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%)	0.56	86 (10%) 11 12	21, 50, 86, 119	3 (0%)
2	B	81/87 (93%)	0.40	7 (8%) 16 17	44, 82, 158, 165	0
All	All	894/967 (92%)	0.55	93 (10%) 11 12	21, 52, 106, 165	3 (0%)

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	281	PHE	5.3
1	A	278	LEU	5.2
1	A	282	ILE	5.1
1	A	268	ALA	5.0
1	A	741	THR	4.8
1	A	349	ALA	4.5
1	A	348	PHE	4.5
1	A	280	ALA	4.4
1	A	305	THR	4.3
1	A	608	ALA	4.2
1	A	606	ASP	4.1
1	A	310	VAL	4.1
1	A	607	ALA	4.0
1	A	279	ALA	3.9
1	A	263	ALA	3.7
1	A	307	PHE	3.7
1	A	262	ALA	3.7
1	A	352	TYR	3.7
1	A	591	ALA	3.7
1	A	267	LEU	3.7
1	A	593	VAL	3.7
1	A	481	MET	3.5
1	A	188	ILE	3.4
1	A	589	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	261	VAL	3.3
1	A	265	HIS	3.2
1	A	303	VAL	3.2
1	A	321	TRP	3.2
1	A	592	ILE	3.2
1	A	601	ILE	3.2
1	A	300	LYS	3.2
1	A	260	ALA	3.2
1	A	309	ALA	3.2
1	A	313	LEU	3.1
1	A	240	ASP	3.1
1	A	239	TYR	3.1
1	A	353	GLY	3.1
1	A	271	ALA	3.1
1	A	323	ALA	3.0
1	A	350	SER	3.0
1	A	325	PHE	3.0
1	A	841	GLY	3.0
2	B	76	A	2.9
1	A	595	ARG	2.9
1	A	354	LEU	2.9
1	A	346	TYR	2.9
1	A	333	GLY	2.8
1	A	283	ASP	2.8
1	A	326	VAL	2.8
1	A	594	GLU	2.7
1	A	742	ASP	2.7
1	A	355	ASN	2.7
1	A	322	ALA	2.7
1	A	242	THR	2.7
1	A	276	PRO	2.7
1	A	356	ILE	2.6
1	A	302	GLY	2.6
1	A	272	ALA	2.6
1	A	238	ASP	2.6
1	A	306	GLY	2.6
1	A	237	ASN	2.6
2	B	74	C	2.5
1	A	478	VAL	2.5
1	A	86	ALA	2.5
1	A	482	PRO	2.5
1	A	602	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	4	C	2.4
1	A	258	TYR	2.4
1	A	604	ALA	2.4
2	B	3	C	2.4
2	B	17	C	2.4
1	A	155	ALA	2.4
1	A	345	ALA	2.4
1	A	275	ASN	2.4
1	A	236	VAL	2.3
1	A	241	ASN	2.3
1	A	462	GLY	2.3
1	A	315	GLY	2.3
1	A	270	LYS	2.3
2	B	41	C	2.3
2	B	73	A	2.2
1	A	257	THR	2.2
1	A	312	PRO	2.1
1	A	387	PHE	2.1
1	A	89	ALA	2.1
1	A	189	PRO	2.1
1	A	90	ALA	2.1
1	A	317	GLU	2.1
1	A	463	ILE	2.1
1	A	264	GLY	2.0
1	A	320	VAL	2.0
1	A	452	VAL	2.0
1	A	555	ASN	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.

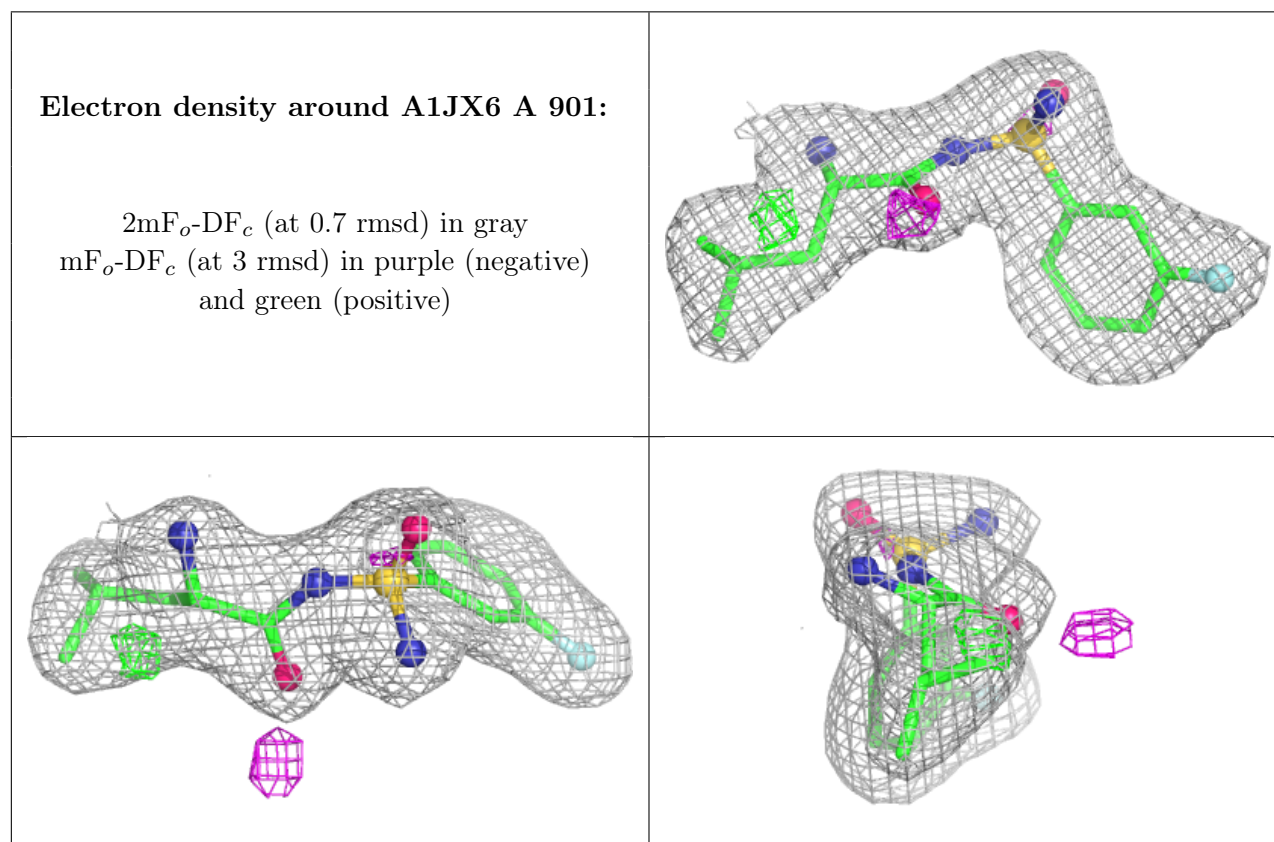
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
5	MG	B	101	1/1	0.73	0.23	91,91,91,91	0
4	PEG	A	902	7/7	0.80	0.18	72,73,74,74	0
3	A1JX6	A	901	19/19	0.95	0.09	38,41,43,44	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 ⓘ

There are no such residues in this entry.