

wwPDB X-ray Structure Validation Summary Report (i)

Nov 27, 2023 – 08:12 PM EST

PDB ID : 8GKM

Title: A1 Leu graft + Leu: Adenylation domain 1 core construct from teicoplanin

biosynthesis, leucine selection pocket graft; leucine bound

Authors: Hansen, M.H.; Cryle, M.J.

Deposited on : 2023-03-20

Resolution : 1.89 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.orgA user guide is available at

https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

MolProbity : 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS: 2.36

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

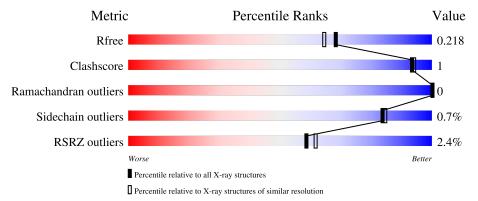
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

The reported resolution of this entry is 1.89 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 $(\# \mathrm{Entries})$	Similar resolution $(\# \text{Entries, resolution range}(\text{\AA}))$
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 <=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	399	92%		5%	-
1	В	399	93%			
2	С	69	84%	• 1	13%	_
2	D	69	90%		7%	,



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 13708 atoms, of which 6643 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 Non-ribosomal peptide synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	А	385	Total		Н		О	S	0	3	0
-	11	900	5733	1798	2869	525	530	11	U		
1	D	382	Total	С	Н	N	Ο	S	0	4	0
1	Б	362	5701	1788	2858	522	522	11	0	4	

There are 24 discrepancies between the modelled and reference sequences:

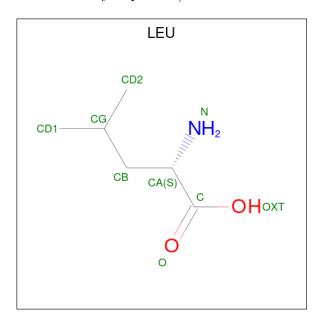
Chain	Residue	Modelled	Actual	Comment	Reference
A	8	MET	-	initiating methionine	UNP Q70AZ9
A	237	TYR	HIS	engineered mutation	UNP Q70AZ9
A	287	MET	LEU	engineered mutation	UNP Q70AZ9
A	295	MET	LEU	engineered mutation	UNP Q70AZ9
A	399	GLY	-	expression tag	UNP Q70AZ9
A	400	SER	-	expression tag	UNP Q70AZ9
A	401	HIS	-	expression tag	UNP Q70AZ9
A	402	HIS	-	expression tag	UNP Q70AZ9
A	403	HIS	-	expression tag	UNP Q70AZ9
A	404	HIS	-	expression tag	UNP Q70AZ9
A	405	HIS	-	expression tag	UNP Q70AZ9
A	406	HIS	-	expression tag	UNP Q70AZ9
В	8	MET	-	initiating methionine	UNP Q70AZ9
В	237	TYR	HIS	engineered mutation	UNP Q70AZ9
В	287	MET	LEU	engineered mutation	UNP Q70AZ9
В	295	MET	LEU	engineered mutation	UNP Q70AZ9
В	399	GLY	-	expression tag	UNP Q70AZ9
В	400	SER	-	expression tag	UNP Q70AZ9
В	401	HIS	-	expression tag	UNP Q70AZ9
В	402	HIS	-	expression tag	UNP Q70AZ9
В	403	HIS	-	expression tag	UNP Q70AZ9
В	404	HIS	- expression tag		UNP Q70AZ9
В	405	HIS	-	expression tag	UNP Q70AZ9
В	406	HIS	-	expression tag	UNP Q70AZ9



• Molecule 2 is a protein called MbtH-like short polypeptide.

Mol	Chain	Residues		A	Atom	S			ZeroOcc	AltConf	Trace
2	2 C	60	Total	С	Н	N	О	S	0	1	0
	00	894	297	423	81	92	1	0	1		
2	D	64	Total	С	Н	N	О	S	0	1	0
2	ע	04	955	315	456	85	98	1	U		U

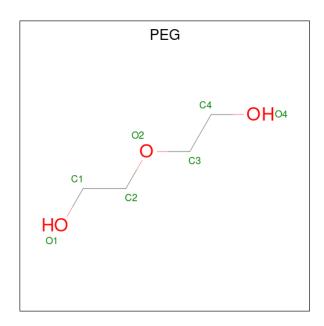
• Molecule 3 is LEUCINE (three-letter code: LEU) (formula: $C_6H_{13}NO_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
9	Λ	1	Total	С	Н	N	О	0	0	
$\begin{array}{ c c c c c } \hline 3 & A & \end{array}$	A	1	21	6	12	1	2	0		
9	D	1	Total	С	Н	N	О	0	0	
3	В	1	21	6	12	1	2	U	U	

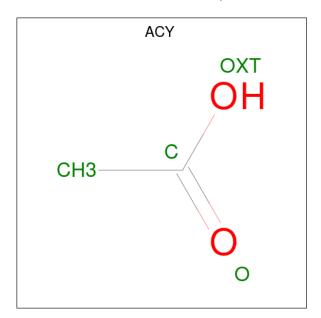
 $\bullet \ \ Molecule\ 4 \ is\ DI(HYDROXYETHYL)ETHER\ (three-letter\ code:\ PEG)\ (formula:\ C_4H_{10}O_3).$





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	С	Н	О	0	0
	11	_	17	4	10	3		

 \bullet Molecule 5 is ACETIC ACID (three-letter code: ACY) (formula: $\mathrm{C_2H_4O_2}).$



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	Λ	1	Total	С	Н	О	0	0
)	A	1	7	2	3	2	U	0

• Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Na 1 1	0	0

• Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	175	Total O 175 175	0	0
7	В	142	Total O 142 142	0	0
7	С	20	Total O 20 20	0	0
7	D	21	Total O 21 21	0	0

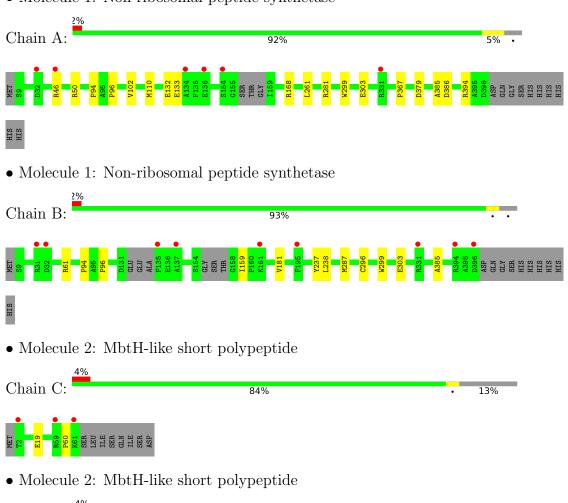


Chain D:

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: Non-ribosomal peptide synthetase



90%



4 Data and refinement statistics (i)

Property	Value	Source	
Space group	P 2 21 21	Depositor	
Cell constants	41.80Å 123.42Å 175.17Å	Donositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	43.79 - 1.89	Depositor	
Resolution (A)	43.79 - 1.88	EDS	
% Data completeness	99.5 (43.79-1.89)	Depositor	
(in resolution range)	99.3 (43.79-1.88)	EDS	
R_{merge}	0.04	Depositor	
R_{sym}	0.04	Depositor	
$< I/\sigma(I) > 1$	1.01 (at 1.88Å)	Xtriage	
Refinement program	PHENIX 1.20.1_4487	Depositor	
D D	0.189 , 0.220	Depositor	
R, R_{free}	0.187 , 0.218	DCC	
R_{free} test set	3732 reflections (5.01%)	wwPDB-VP	
Wilson B-factor (Å ²)	27.5	Xtriage	
Anisotropy	0.747	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.44, 44.7	EDS	
L-test for twinning ²	$ < L >=0.49, < L^2>=0.33$	Xtriage	
Estimated twinning fraction	No twinning to report.	Xtriage	
F_o, F_c correlation	0.96	EDS	
Total number of atoms	13708	wwPDB-VP	
Average B, all atoms (Å ²)	39.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 17.22% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: ACY, NA, 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		
MIOI		RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.29	0/2935	0.61	0/4002	
1	В	0.28	0/2919	0.61	0/3979	
2	С	0.27	0/495	0.51	0/678	
2	D	0.28	0/523	0.52	0/716	
All	All	0.29	0/6872	0.60	0/9375	

There are no bond length outliers.

There are no bond angle outliers.

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	2864	2869	2858	13	0
1	В	2843	2858	2834	7	0
2	С	471	423	416	1	0
2	D	499	456	448	2	0
3	A	9	12	10	0	0
3	В	9	12	10	0	0
4	A	7	10	10	0	0
5	A	4	3	3	0	0
6	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	175	0	0	2	0
7	В	142	0	0	0	0
7	С	20	0	0	0	0
7	D	21	0	0	1	0
All	All	7065	6643	6589	19	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 1.

The worst 5 of 19 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
2:D:45:GLN:NE2	7:D:101:HOH:O	1.93	0.99
1:A:386:ASP:N	7:A:601:HOH:O	2.17	0.76
1:A:46:ARG:NH2	1:A:132:GLU:OE1	2.20	0.75
1:A:303:GLU:OE2	1:B:303:GLU:OE2	2.10	0.67
1:A:50:ARG:NH2	1:A:133:GLU:OE2	2.36	0.56

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	Perce	entiles
1	A	384/399 (96%)	378 (98%)	6 (2%)	0	100	100
1	В	380/399~(95%)	374 (98%)	6 (2%)	0	100	100
2	С	59/69 (86%)	59 (100%)	0	0	100	100
2	D	63/69 (91%)	63 (100%)	0	0	100	100
All	All	886/936 (95%)	874 (99%)	12 (1%)	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	$286/295 \ (97\%)$	285 (100%)	1 (0%)	92	93	
1	В	285/295~(97%)	283 (99%)	2 (1%)	84	84	
2	\mathbf{C}	51/59 (86%)	51 (100%)	0	100	100	
2	D	55/59~(93%)	55 (100%)	0	100	100	
All	All	$677/708 \; (96\%)$	674 (100%)	3 (0%)	84	91	

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	299	TRP
1	В	61	ARG
1	В	299	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 1 is 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	Mol Type Chain Res		es Link	Bond lengths			Bond angles			
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
3	LEU	A	501	-	7,8,8	0.80	0	9,10,10	1.13	1 (11%)
3	LEU	В	501	-	7,8,8	0.77	0	9,10,10	1.06	2 (22%)
5	ACY	A	503	-	3,3,3	0.63	0	3,3,3	0.62	0
4	PEG	A	502	-	6,6,6	0.07	0	5,5,5	0.15	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	LEU	A	501	-	-	3/8/8/8	-
3	LEU	В	501	-	-	3/8/8/8	-
4	PEG	A	502	-	-	2/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
3	В	501	LEU	OXT-C-CA	2.02	120.27	113.38
3	A	501	LEU	OXT-C-O	-2.02	119.51	124.09
3	В	501	LEU	OXT-C-O	-2.00	119.54	124.09

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

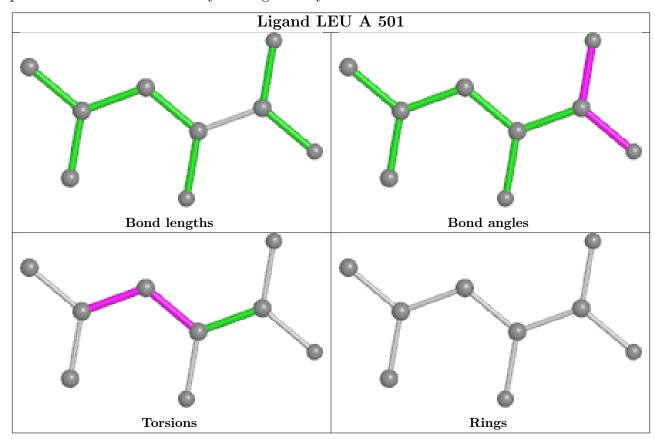
Mol	Chain	Res	Type	Atoms
3	A	501	LEU	N-CA-CB-CG
3	В	501	LEU	N-CA-CB-CG
3	В	501	LEU	C-CA-CB-CG
4	A	502	PEG	O2-C3-C4-O4
3	A	501	LEU	C-CA-CB-CG



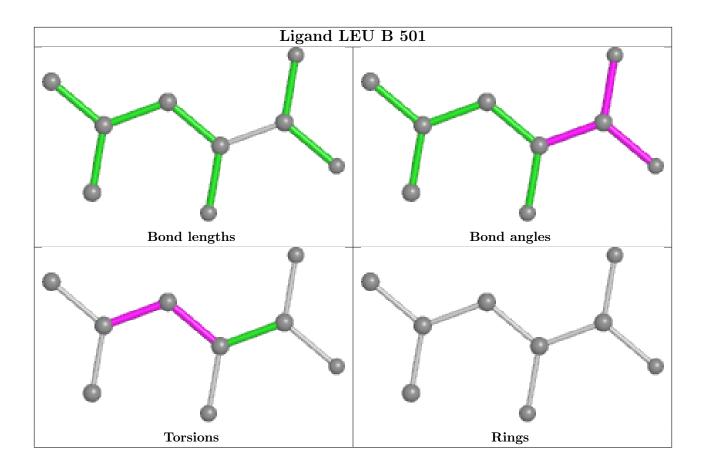
There are no ring outliers.

No monomer is involved in short contacts.

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 then 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, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ $>$	$\# \mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q < 0.9
1	A	385/399~(96%)	0.29	6 (1%) 72 74	23, 31, 48, 69	0
1	В	382/399~(95%)	0.27	9 (2%) 59 62	23, 34, 51, 78	0
2	С	60/69 (86%)	0.33	3 (5%) 28 32	30, 40, 66, 79	0
2	D	64/69 (92%)	0.38	3 (4%) 31 34	31, 37, 55, 68	0
All	All	891/936 (95%)	0.29	21 (2%) 59 62	23, 34, 53, 79	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	32	ASP	5.9
2	С	59	ARG	5.0
1	В	32	ASP	5.0
1	В	135	PHE	4.3
1	A	46	ARG	3.6

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 monosaccharides in this entry.

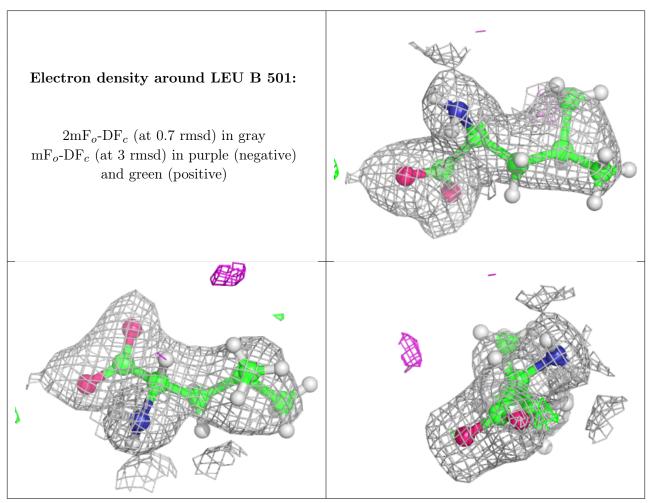
6.4 Ligands (i)

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, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

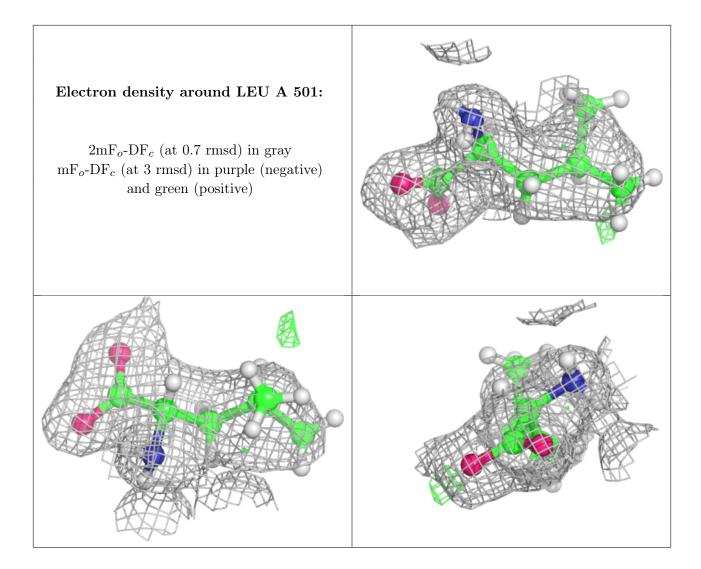


Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	PEG	A	502	7/7	0.74	0.15	43,55,66,67	0
3	LEU	В	501	9/9	0.83	0.15	35,43,51,51	0
3	LEU	A	501	9/9	0.91	0.13	31,40,46,46	0
6	NA	A	504	1/1	0.91	0.12	43,43,43,43	0
5	ACY	A	503	4/4	0.92	0.09	34,42,49,49	0

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.

