



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 13, 2024 – 12:53 pm GMT

PDB ID : 6SLJ
Title : Structure of the RagAB peptide transporter
Authors : Madej, M.; Ranson, N.A.; White, J.B.R.
Deposited on : 2019-08-20
Resolution : 3.04 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.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 : 4.02b-467
Mogul : 1.8.4, 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 : 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

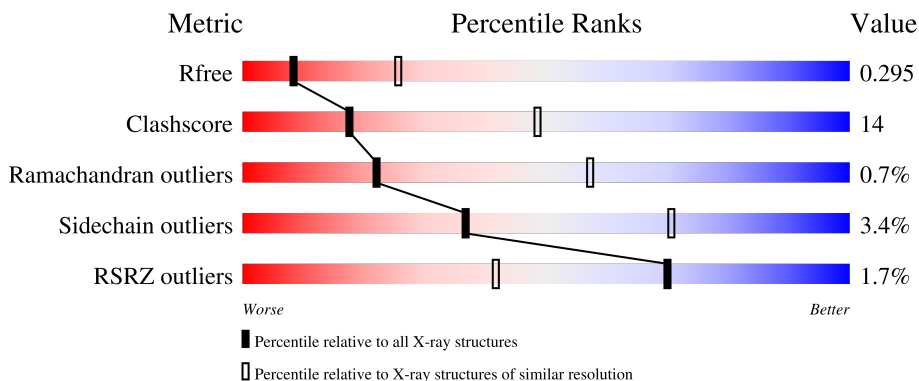
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.04 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}	130704	2752 (3.08-3.00)
Clashscore	141614	3096 (3.08-3.00)
Ramachandran outliers	138981	2986 (3.08-3.00)
Sidechain outliers	138945	2988 (3.08-3.00)
RSRZ outliers	127900	2636 (3.08-3.00)

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	997	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 58%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 31%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="margin-left: 20px;">2% 58% 31% • 10%</p>
1	B	997	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 61%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 28%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 10%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="margin-left: 20px;">3% 61% 28% • 10%</p>
2	C	488	<div style="display: flex; align-items: center;"> <div style="width: 0%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 72%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 26%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="margin-left: 20px;">% 72% 26% ••</p>
2	D	488	<div style="display: flex; align-items: center;"> <div style="width: 0%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 73%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 25%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="margin-left: 20px;">73% 25% ••</p>
3	P	13	<div style="display: flex; align-items: center;"> <div style="width: 69%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: orange; margin-right: 5px;"></div> </div> <p style="margin-left: 20px;">69% 15% 15%</p>

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Mol	Chain	Length	Quality of chain
4	Q	10	 50% 30% 20%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	C8E	A	1102	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 22178 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 RagA protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	900	Total 7069	C 4482	N 1183	O 1372	S 32	0	3	0
1	A	898	Total 7035	C 4459	N 1177	O 1367	S 32	0	0	0

- Molecule 2 is a protein called Lipoprotein RagB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	482	Total 3842	C 2440	N 656	O 737	S 9	0	0	0
2	D	482	Total 3841	C 2440	N 656	O 736	S 9	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	502	HIS	-	expression tag	UNP F5H948
C	503	HIS	-	expression tag	UNP F5H948
C	504	HIS	-	expression tag	UNP F5H948
C	505	HIS	-	expression tag	UNP F5H948
C	506	HIS	-	expression tag	UNP F5H948
C	507	HIS	-	expression tag	UNP F5H948
D	502	HIS	-	expression tag	UNP F5H948
D	503	HIS	-	expression tag	UNP F5H948
D	504	HIS	-	expression tag	UNP F5H948
D	505	HIS	-	expression tag	UNP F5H948
D	506	HIS	-	expression tag	UNP F5H948
D	507	HIS	-	expression tag	UNP F5H948

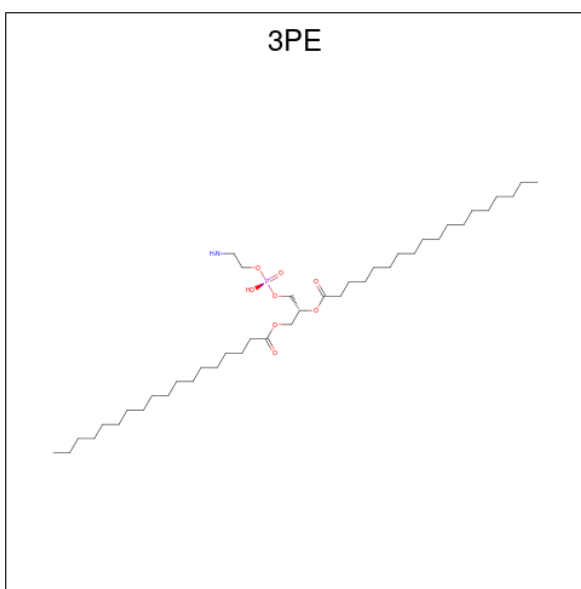
- Molecule 3 is a protein called ALA-SER-THR-THR-GLY-ALA-ASN-SER-GLN-ARG-GLY-SER-GLY.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	P	13	82	44	18	20	0	0	0

- Molecule 4 is a protein called ALA-SER-THR-THR-GLY-ALA-ASN-SER-GLN-ARG.

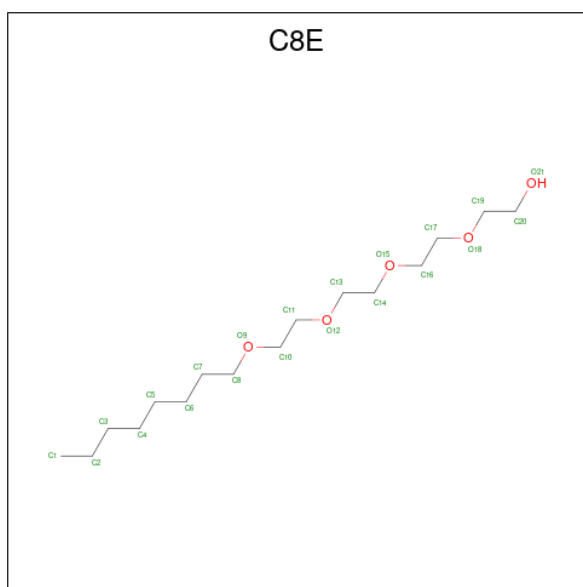
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	Q	10	68	37	15	16	0	0	0

- Molecule 5 is 1,2-Distearoyl-sn-glycerophosphoethanolamine (three-letter code: 3PE) (formula: $C_{41}H_{82}NO_8P$).



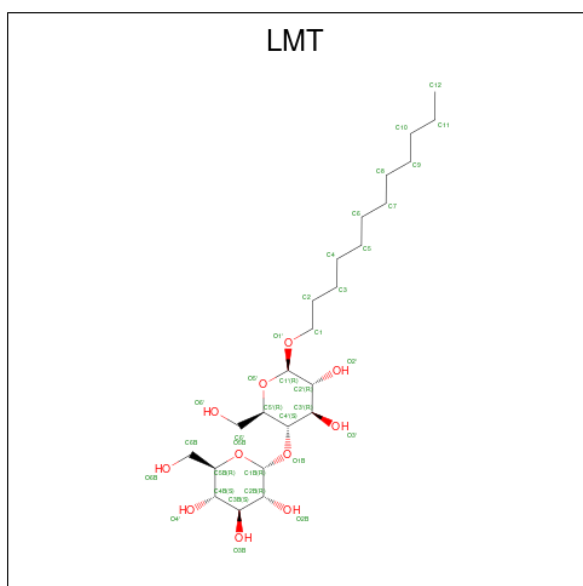
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	B	1	34	30	4	0	0
5	A	1	34	30	4	0	0

- Molecule 6 is (HYDROXYETHYLOXY)TRI(ETHYLOXY)OCTANE (three-letter code: C8E) (formula: $C_{16}H_{34}O_5$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	C O	0	0
			21	16 5		
6	A	1	Total	C O	0	0
			21	16 5		

- Molecule 7 is DODECYL-BETA-D-MALTOSE (three-letter code: LMT) (formula: $C_{24}H_{46}O_{11}$).



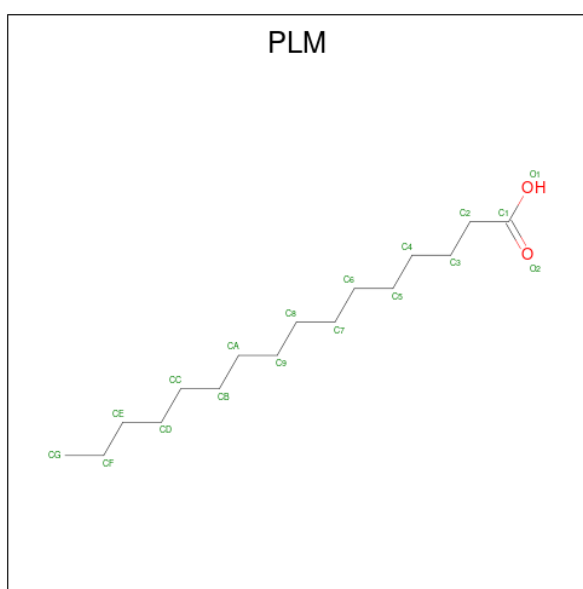
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	C O	0	0
			28	17 11		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			13	12	1		
7	A	1	Total	C	O	0	0
			13	12	1		
7	C	1	Total	C	O	0	0
			34	23	11		
7	C	1	Total	C	O	0	0
			13	12	1		

- Molecule 8 is PALMITIC ACID (three-letter code: PLM) (formula: C₁₆H₃₂O₂).

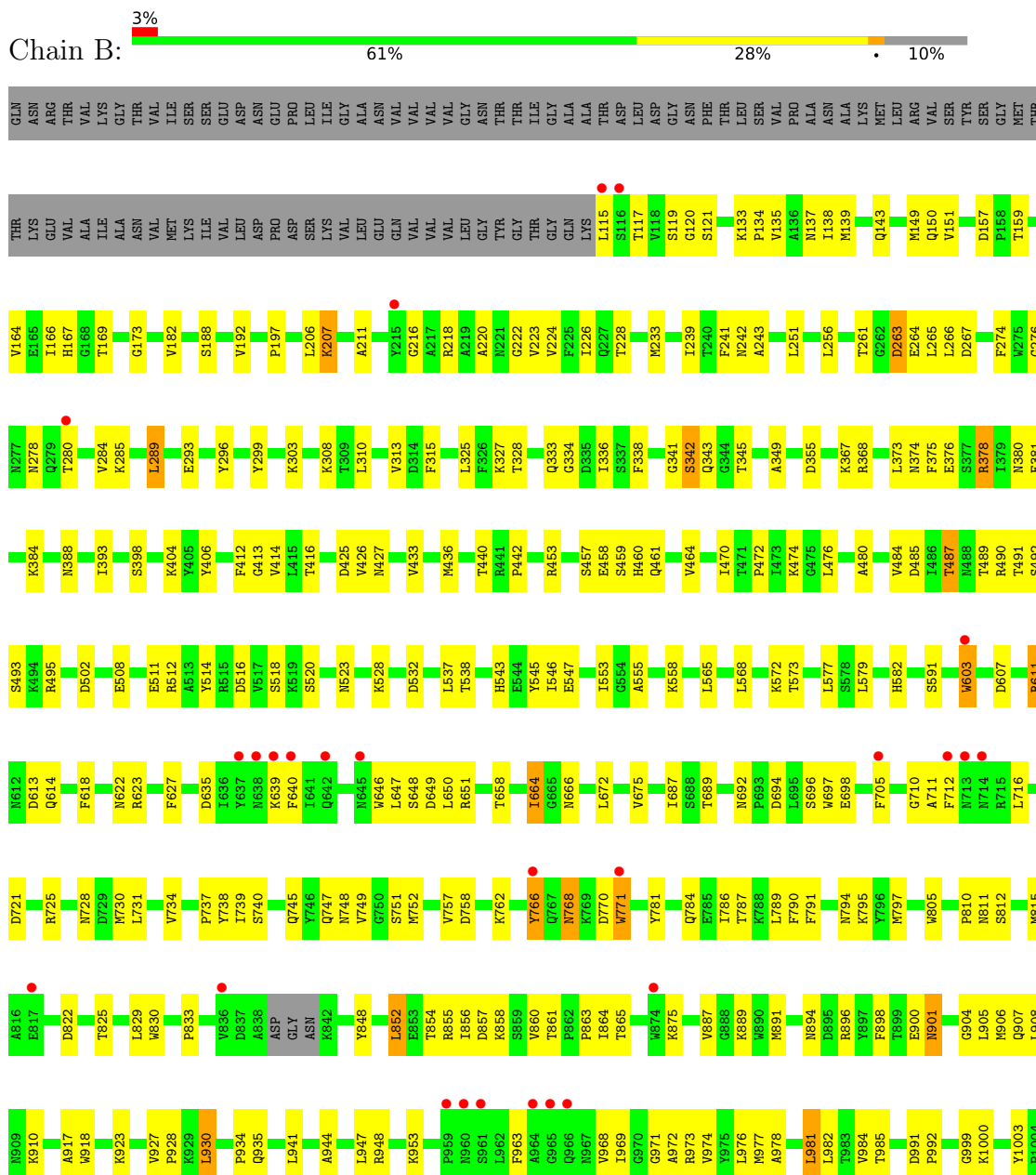


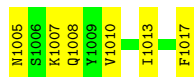
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	C	1	Total	C	O	0	0
			14	13	1		
8	D	1	Total	C	O	0	0
			16	15	1		

3 Residue-property plots

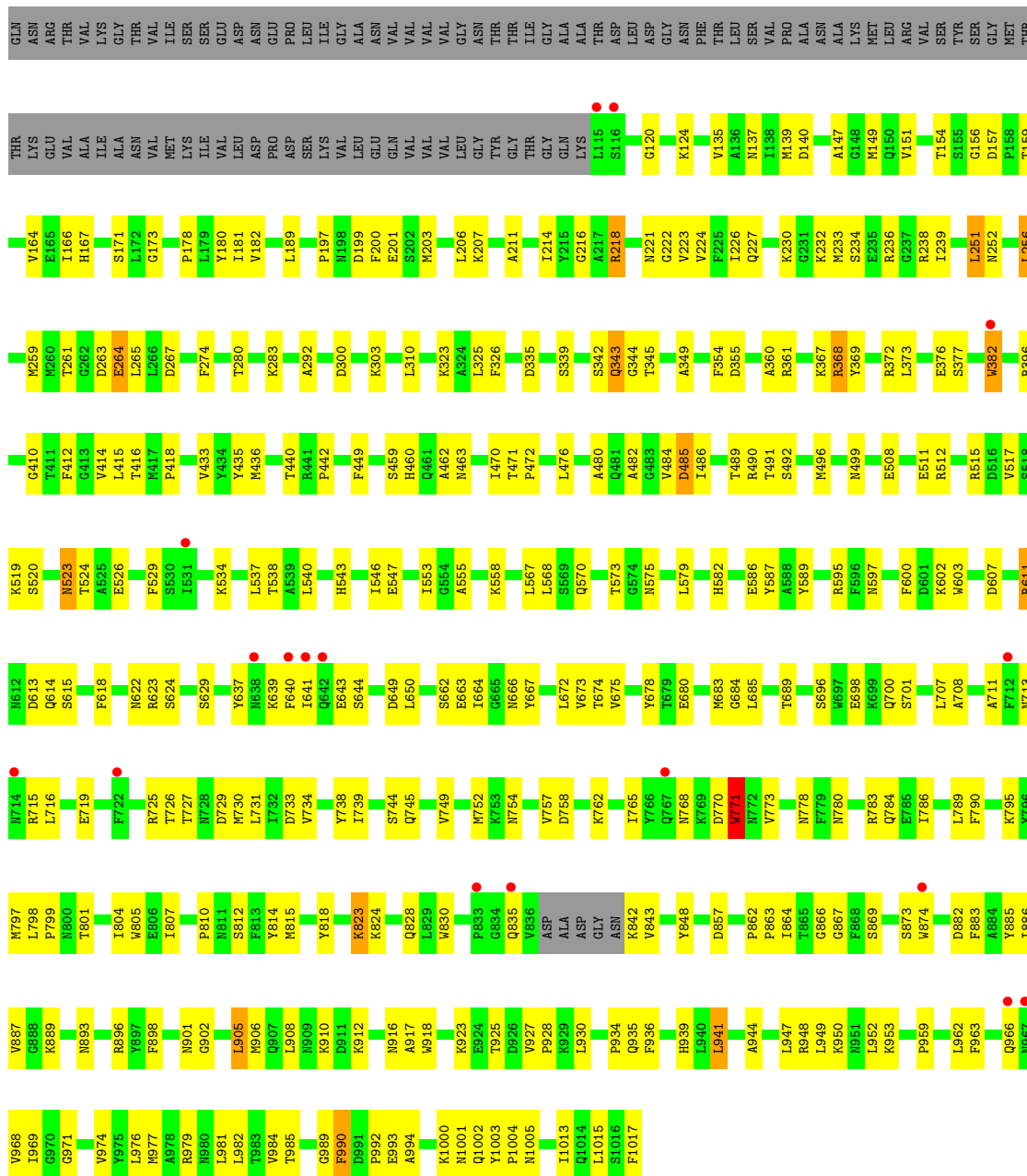
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



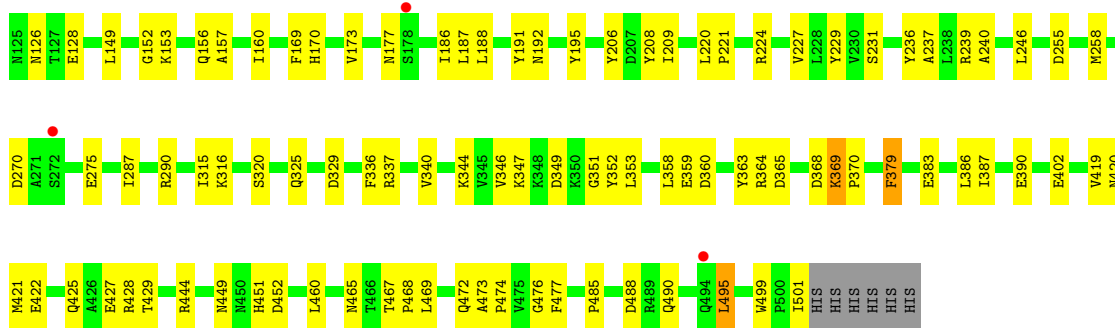


• Molecule 1: RagA protein

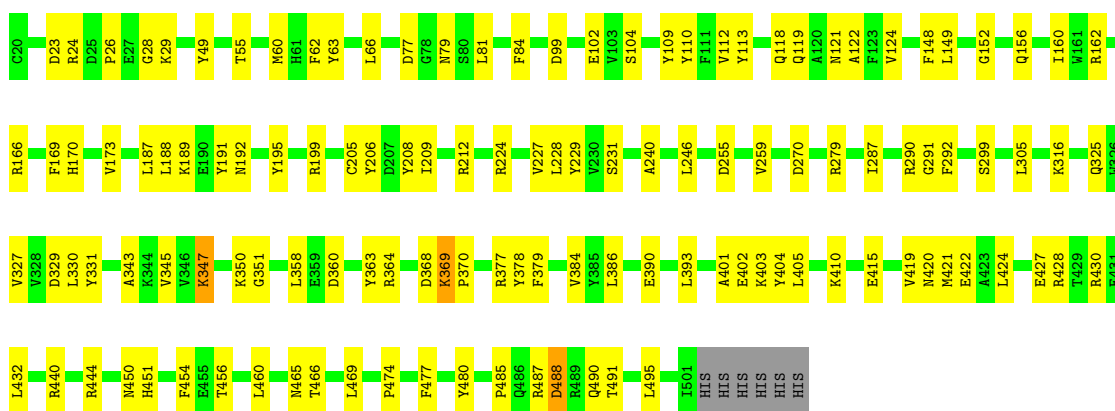


• Molecule 2: Lipoprotein RagB





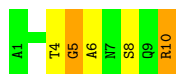
• Molecule 2: Lipoprotein RagB



• Molecule 3: ALA-SER-THR-THR-GLY-ALA-ASN-SER-GLN-ARG-GLY-SER-GLY



• Molecule 4: ALA-SER-THR-THR-GLY-ALA-ASN-SER-GLN-ARG



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	130.60Å 142.02Å 241.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.01 – 3.04 75.22 – 3.04	Depositor EDS
% Data completeness (in resolution range)	99.8 (71.01-3.04) 99.9 (75.22-3.04)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 3.01Å)	Xtrriage
Refinement program	PHENIX 1.15.2_3472, PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.222 , 0.295 0.221 , 0.295	Depositor DCC
R_{free} test set	4354 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	43.6	Xtrriage
Anisotropy	0.938	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	22178	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.90% 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: LMT, 3PE, C8E, PLM

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	0.52	0/7198	0.76	7/9735 (0.1%)
1	B	0.52	0/7242	0.75	7/9794 (0.1%)
2	C	0.50	0/3928	0.68	1/5331 (0.0%)
2	D	0.50	0/3927	0.69	0/5330
3	P	0.45	0/81	0.90	0/107
4	Q	0.49	0/67	0.72	0/89
All	All	0.51	0/22443	0.73	15/30386 (0.0%)

There are no bond length outliers.

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	981	LEU	CA-CB-CG	8.05	133.81	115.30
1	A	256	LEU	CA-CB-CG	7.46	132.45	115.30
1	A	251	LEU	CB-CG-CD2	6.55	122.14	111.00
2	C	22	LEU	CA-CB-CG	5.98	129.06	115.30
1	B	852	LEU	CA-CB-CG	5.93	128.95	115.30

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	7035	0	6783	235	1
1	B	7069	0	6820	210	1
2	C	3842	0	3739	98	0
2	D	3841	0	3739	89	0
3	P	82	0	77	8	0
4	Q	68	0	66	9	0
5	A	34	0	49	6	0
5	B	34	0	49	2	0
6	A	21	0	34	6	0
6	B	21	0	34	0	0
7	A	26	0	45	2	0
7	B	28	0	28	1	0
7	C	47	0	64	7	0
8	C	14	0	22	3	0
8	D	16	0	26	0	0
All	All	22178	0	21575	601	1

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 14.

The worst 5 of 601 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:B:440:THR:HG23	2:D:316:LYS:HB2	1.50	0.93
1:A:218:ARG:HD3	1:A:663:GLU:HG2	1.48	0.93
1:A:768:ASN:HB3	1:A:771:TRP:HB2	1.52	0.91
1:B:672:LEU:H	1:B:689:THR:HG22	1.37	0.89
3:P:10:ARG:H	3:P:10:ARG:HD3	1.36	0.89

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:694:ASP:OD1	1:A:280:THR:OG1[3_645]	2.05	0.15

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	894/997 (90%)	817 (91%)	71 (8%)	6 (1%)	22	57
1	B	899/997 (90%)	821 (91%)	70 (8%)	8 (1%)	17	52
2	C	480/488 (98%)	453 (94%)	27 (6%)	0	100	100
2	D	480/488 (98%)	449 (94%)	30 (6%)	1 (0%)	47	80
3	P	11/13 (85%)	7 (64%)	2 (18%)	2 (18%)	0	0
4	Q	8/10 (80%)	5 (62%)	2 (25%)	1 (12%)	0	1
All	All	2772/2993 (93%)	2552 (92%)	202 (7%)	18 (1%)	22	60

5 of 18 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	343	GLN
1	B	771	TRP
1	A	343	GLN
1	A	771	TRP
1	A	1001	ASN

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	751/832 (90%)	718 (96%)	33 (4%)	28	63
1	B	755/832 (91%)	723 (96%)	32 (4%)	30	64
2	C	402/408 (98%)	395 (98%)	7 (2%)	60	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	D	402/408 (98%)	396 (98%)	6 (2%)	65 86
3	P	8/8 (100%)	7 (88%)	1 (12%)	4 18
4	Q	7/7 (100%)	5 (71%)	2 (29%)	0 1
All	All	2325/2495 (93%)	2244 (96%)	81 (4%)	37 69

5 of 81 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	823	LYS
2	C	495	LEU
1	A	857	ASP
2	C	64	GLN
2	D	379	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	156	GLN
2	C	156	GLN
2	C	54	ASN
1	A	780	ASN
2	C	118	GLN

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

11 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	C8E	B	1102	-	20,20,20	0.61	0	19,19,19	0.61	0
7	LMT	B	1103	-	29,29,36	1.18	2 (6%)	40,40,47	0.96	2 (5%)
5	3PE	A	1101	2	33,33,50	1.16	4 (12%)	35,35,55	1.40	5 (14%)
7	LMT	A	1104	-	12,12,36	0.16	0	11,11,47	0.57	0
7	LMT	A	1103	-	12,12,36	0.33	0	11,11,47	0.79	0
7	LMT	C	602	-	35,35,36	1.25	6 (17%)	46,46,47	1.13	2 (4%)
6	C8E	A	1102	-	20,20,20	0.55	0	19,19,19	0.52	0
8	PLM	C	601	2	13,13,17	0.77	0	12,12,17	0.58	0
5	3PE	B	1101	2	33,33,50	1.12	3 (9%)	35,35,55	1.94	5 (14%)
8	PLM	D	601	2	15,15,17	0.86	0	14,14,17	0.67	0
7	LMT	C	603	-	12,12,36	0.12	0	11,11,47	0.74	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
6	C8E	B	1102	-	-	9/18/18/18	-
7	LMT	B	1103	-	-	8/14/54/61	0/2/2/2
5	3PE	A	1101	2	-	12/34/34/54	-
7	LMT	A	1104	-	-	8/10/10/61	-
7	LMT	A	1103	-	-	7/10/10/61	-
7	LMT	C	602	-	-	11/20/60/61	0/2/2/2
6	C8E	A	1102	-	-	9/18/18/18	-
8	PLM	C	601	2	-	7/10/11/15	-
5	3PE	B	1101	2	-	17/34/34/54	-
8	PLM	D	601	2	-	5/12/13/15	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	LMT	C	603	-	-	4/10/10/61	-

The worst 5 of 15 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1101	3PE	O21-C2	-3.78	1.40	1.47
7	C	602	LMT	O3'-C3'	-3.11	1.35	1.43
5	B	1101	3PE	O21-C21	3.11	1.43	1.34
7	B	1103	LMT	O3'-C3'	-2.70	1.36	1.43
5	A	1101	3PE	O21-C21	2.51	1.41	1.34

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1101	3PE	C2-O21-C21	6.64	126.42	117.88
5	B	1101	3PE	O21-C21-C22	6.56	125.64	111.50
5	A	1101	3PE	O21-C21-C22	4.79	121.82	111.50
7	C	602	LMT	C3'-C4'-C5'	-3.69	102.46	110.93
5	A	1101	3PE	O31-C31-C32	3.07	121.55	111.91

There are no chirality outliers.

5 of 97 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1101	3PE	O22-C21-O21-C2
5	A	1101	3PE	C22-C21-O21-C2
7	B	1103	LMT	C2'-C1'-O1'-C1
7	B	1103	LMT	O5'-C1'-O1'-C1
5	A	1101	3PE	O22-C21-O21-C2

There are no ring outliers.

8 monomers are involved in 20 short contacts:

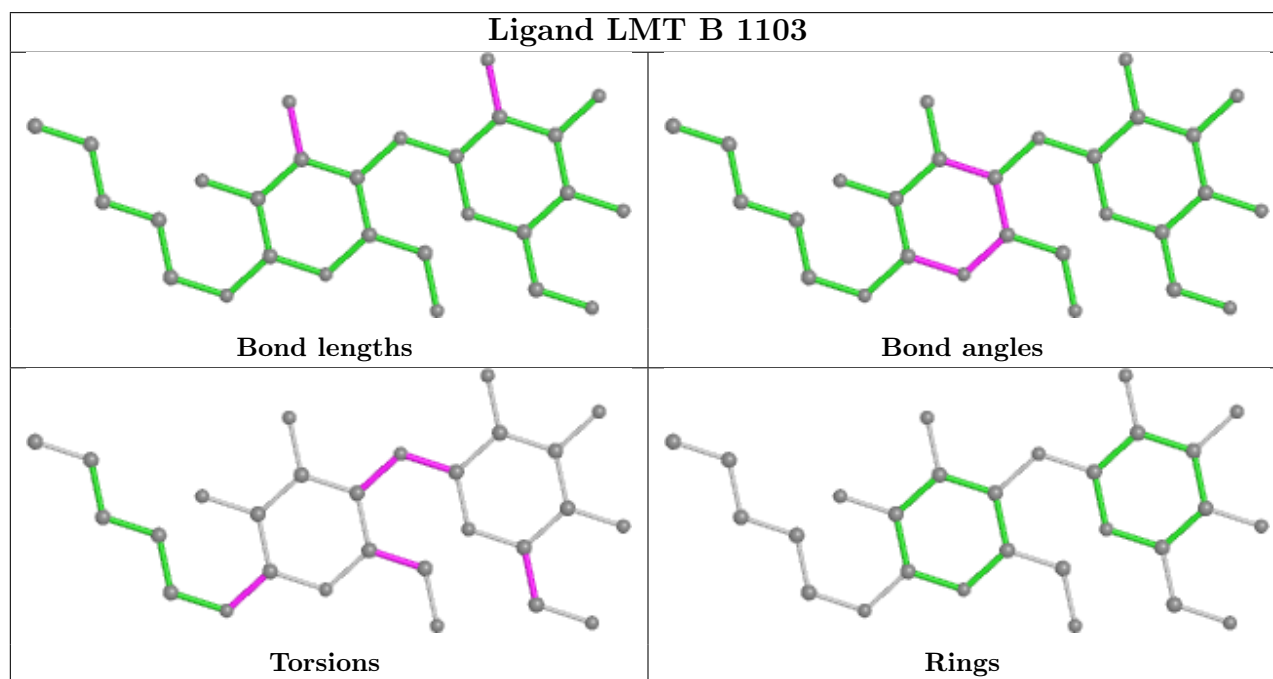
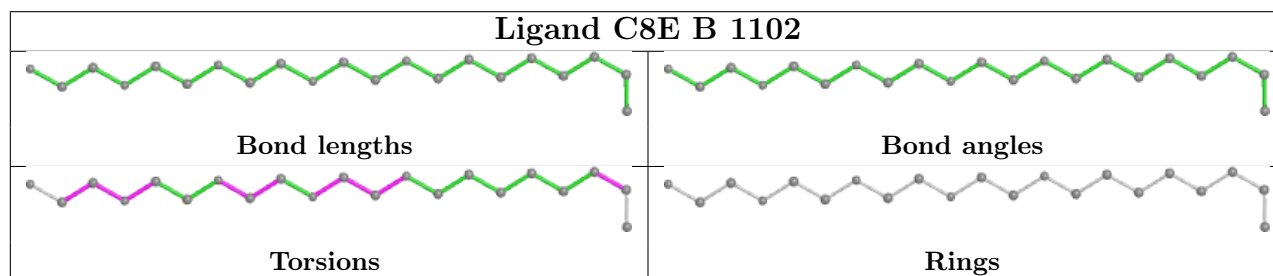
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	1103	LMT	1	0
5	A	1101	3PE	6	0
7	A	1104	LMT	2	0
7	C	602	LMT	2	0
6	A	1102	C8E	6	0
8	C	601	PLM	3	0
5	B	1101	3PE	2	0

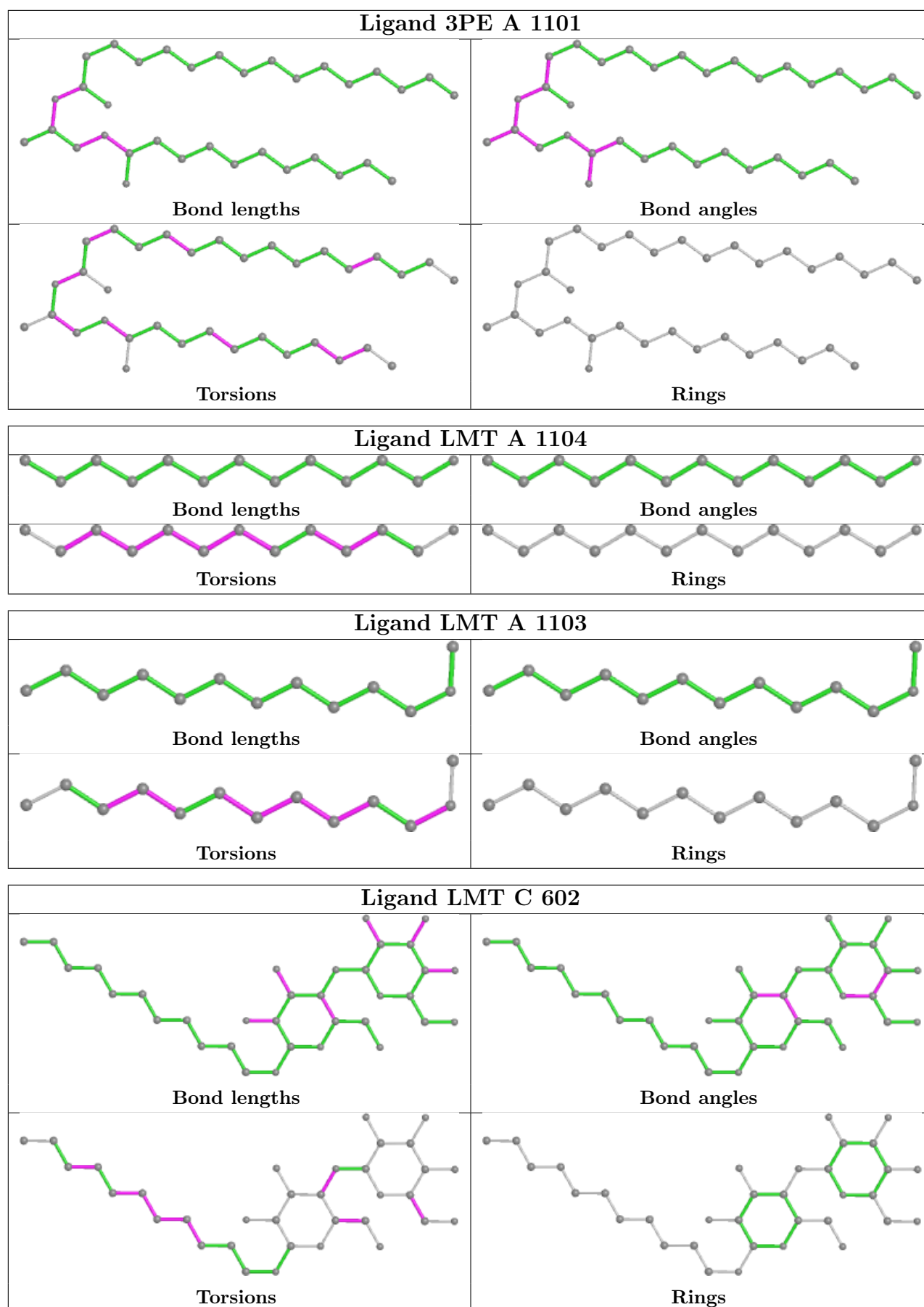
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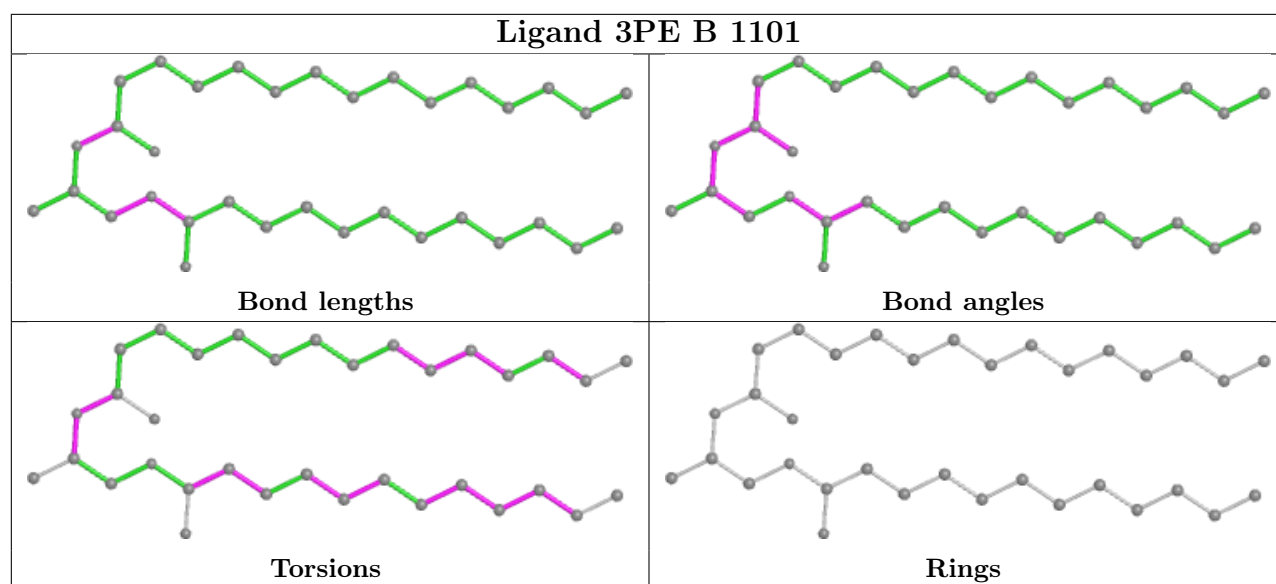
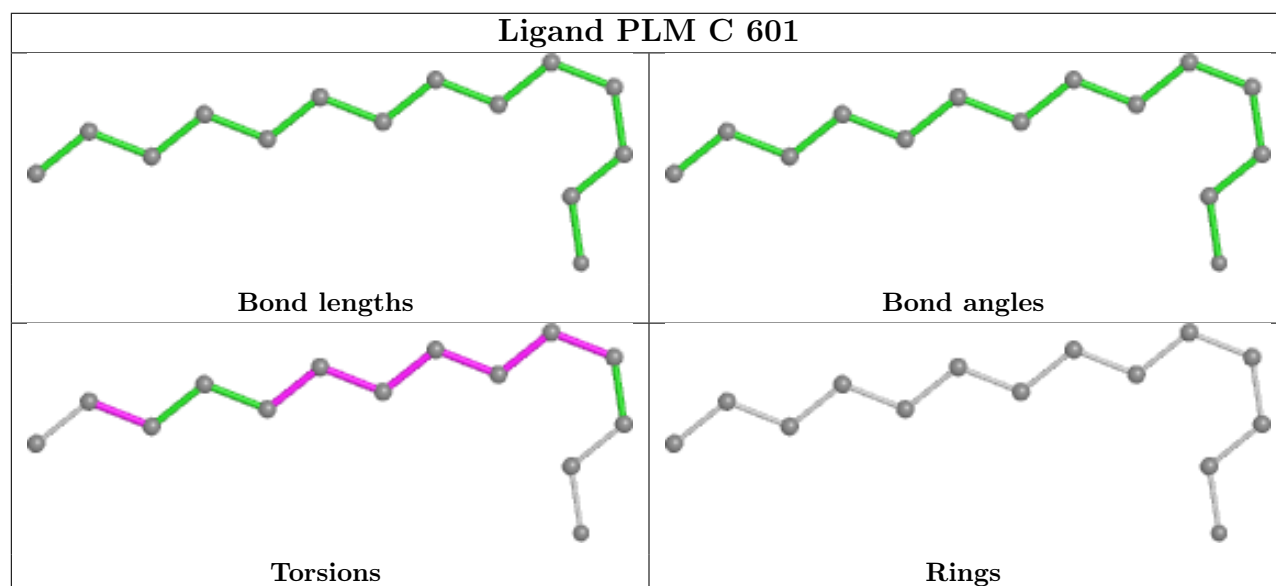
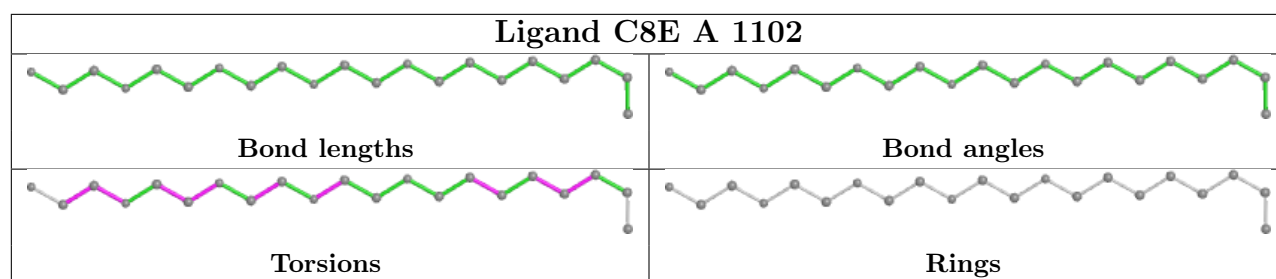
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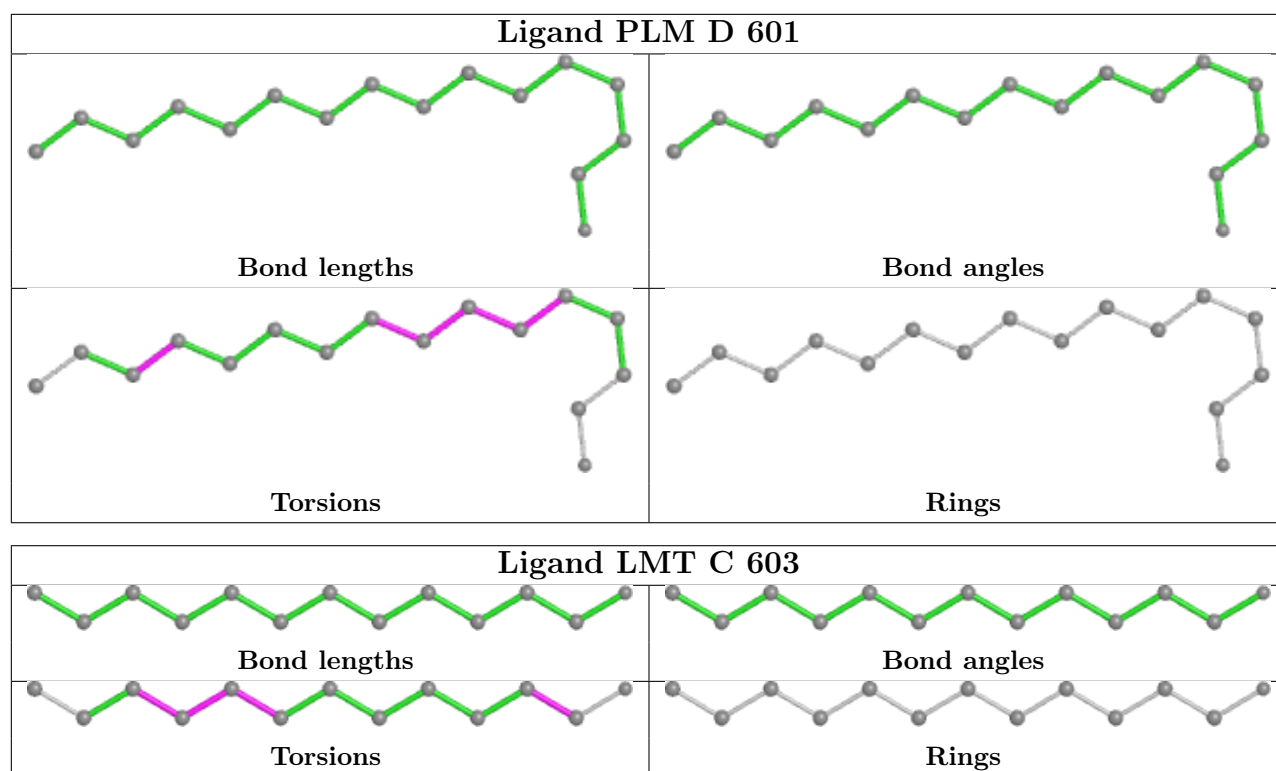
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	603	LMT	5	0

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	898/997 (90%)	-0.15	17 (1%) 66 38	19, 40, 69, 94	0
1	B	900/997 (90%)	-0.07	26 (2%) 51 23	21, 40, 76, 108	0
2	C	482/488 (98%)	-0.06	3 (0%) 89 72	21, 38, 58, 73	0
2	D	482/488 (98%)	-0.16	0 100 100	21, 35, 54, 72	0
3	P	13/13 (100%)	0.71	0 100 100	50, 62, 96, 117	0
4	Q	10/10 (100%)	0.26	0 100 100	54, 61, 66, 68	0
All	All	2785/2993 (93%)	-0.11	46 (1%) 70 42	19, 39, 67, 117	0

The worst 5 of 46 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	645	ASN	4.7
1	B	642	GLN	4.6
1	A	116	SER	4.4
1	B	714	ASN	4.2
1	B	959	PRO	4.0

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](#)

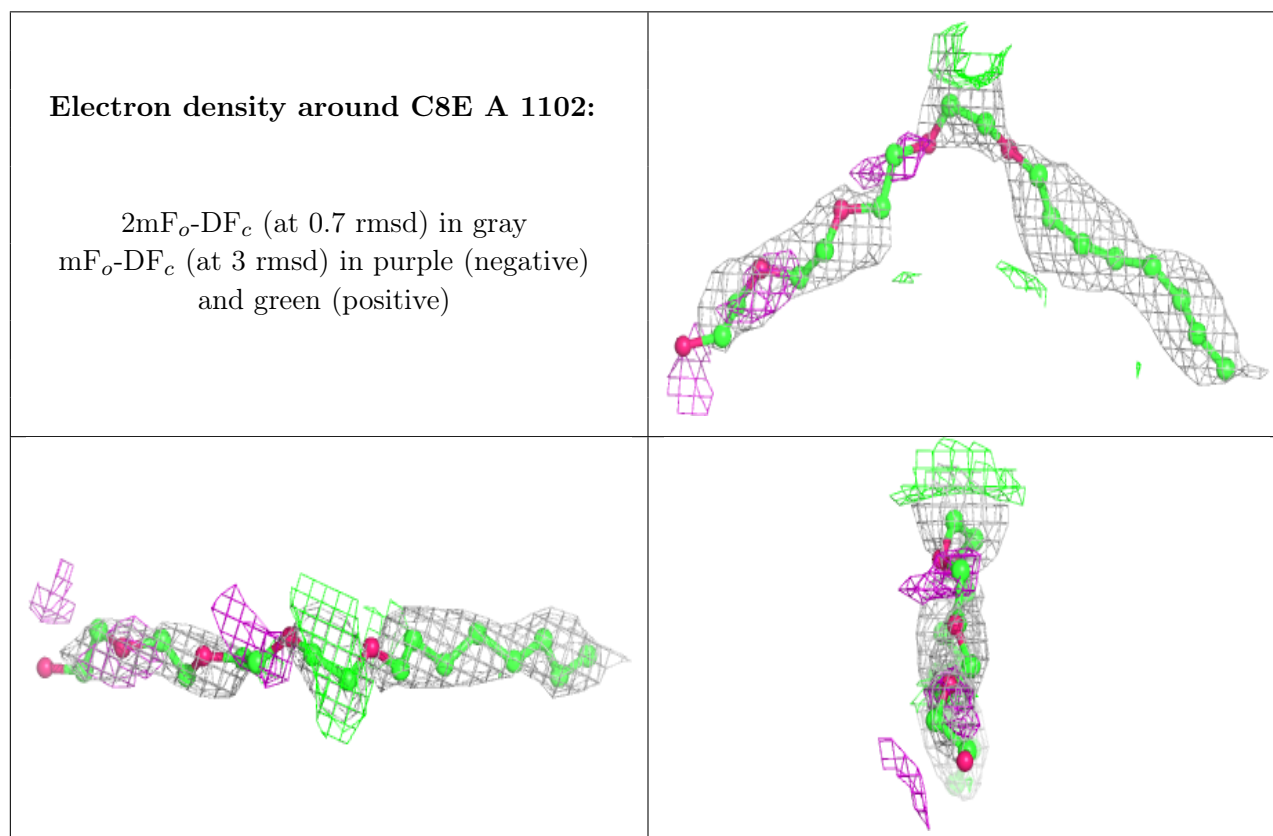
There are no monosaccharides in this entry.

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, 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.

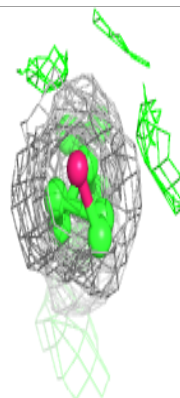
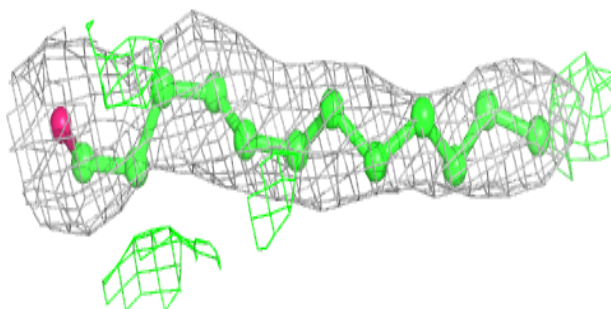
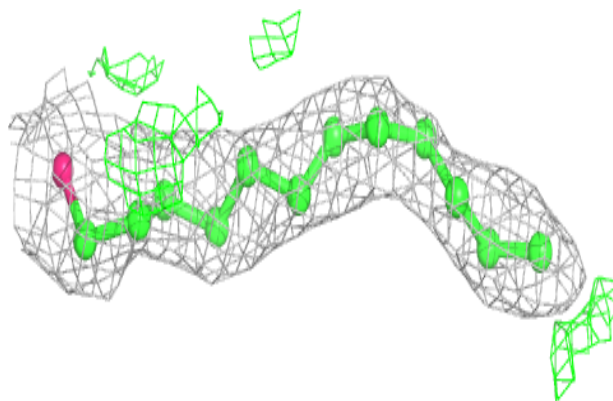
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	C8E	A	1102	21/21	0.79	0.45	19,34,60,62	0
7	LMT	A	1104	13/35	0.83	0.29	7,19,62,63	0
7	LMT	A	1103	13/35	0.86	0.23	37,42,58,63	0
6	C8E	B	1102	21/21	0.87	0.43	24,36,44,49	0
7	LMT	B	1103	28/35	0.88	0.23	26,54,70,88	0
7	LMT	C	602	34/35	0.90	0.27	33,48,65,71	0
5	3PE	B	1101	34/51	0.92	0.41	7,29,39,41	0
5	3PE	A	1101	34/51	0.92	0.35	24,38,56,64	0
7	LMT	C	603	13/35	0.92	0.30	24,34,46,49	0
8	PLM	C	601	14/18	0.92	0.25	7,20,43,51	0
8	PLM	D	601	16/18	0.92	0.48	17,37,47,47	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.

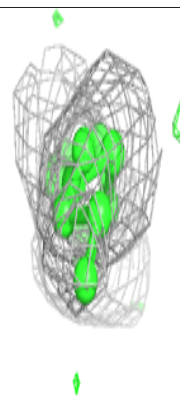
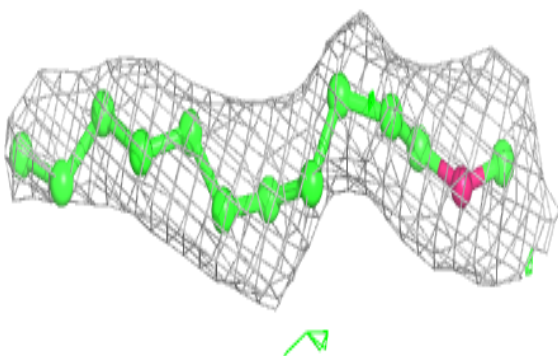
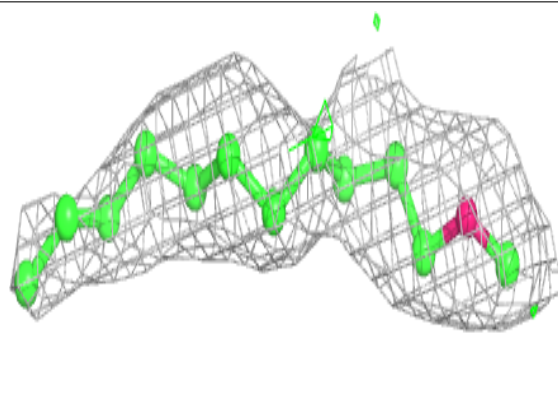


Electron density around LMT A 1104:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

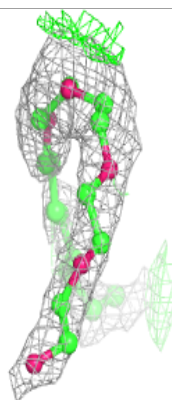
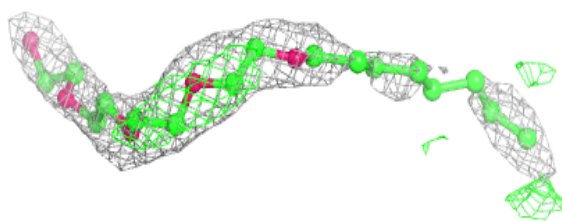
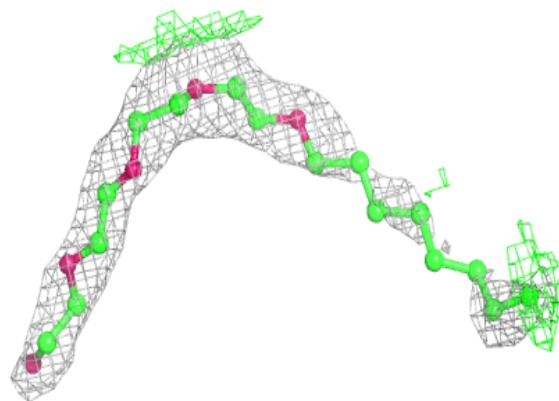
**Electron density around LMT A 1103:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

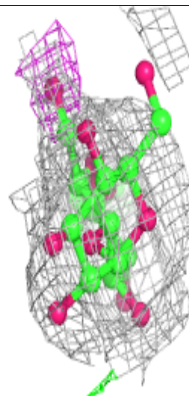
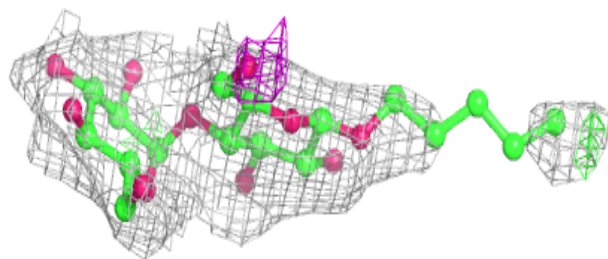
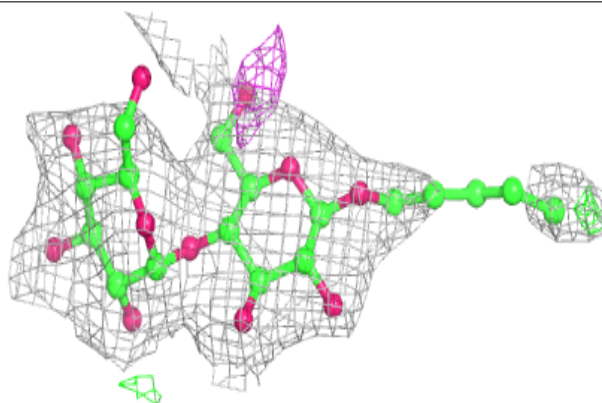


Electron density around C8E B 1102:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

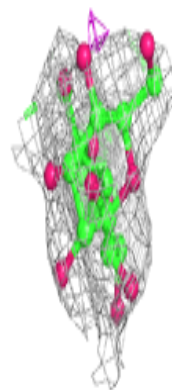
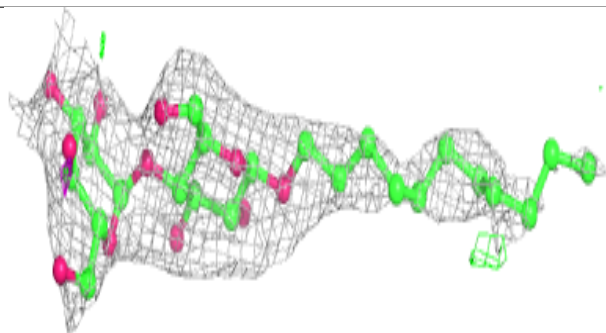
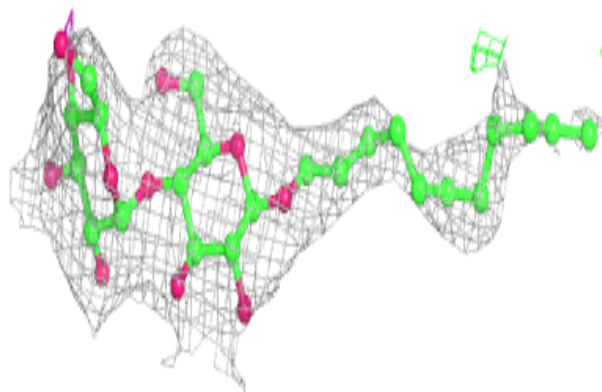
**Electron density around LMT B 1103:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

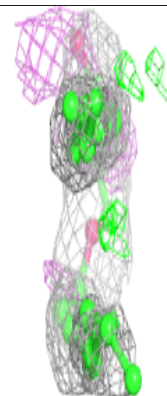
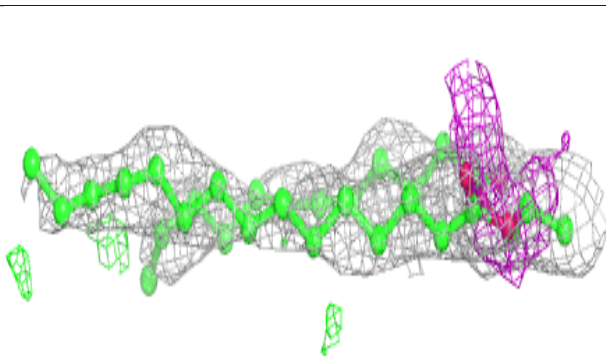
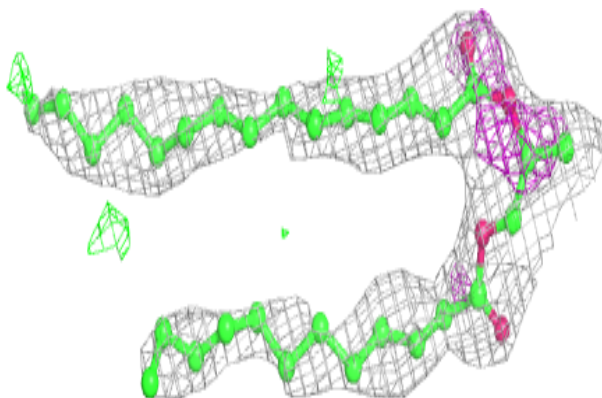


Electron density around LMT C 602:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

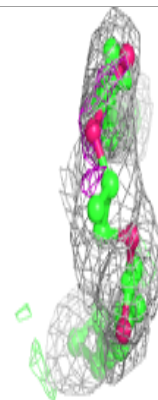
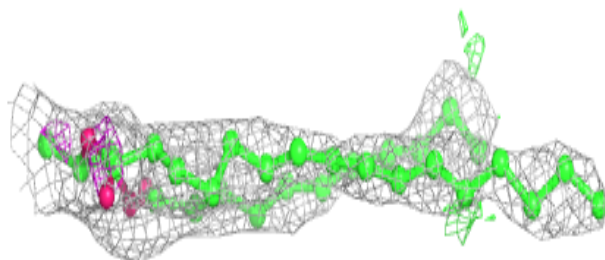
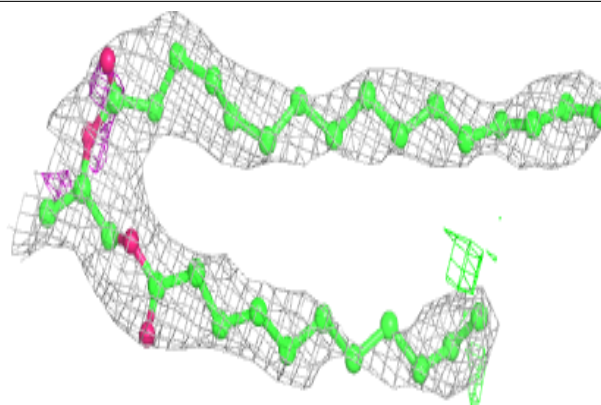
**Electron density around 3PE B 1101:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

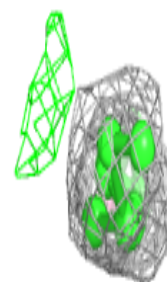
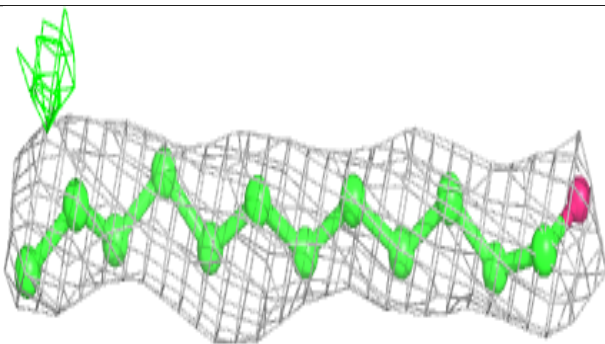
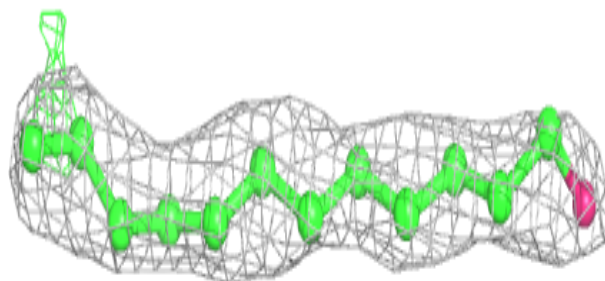


Electron density around 3PE A 1101:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

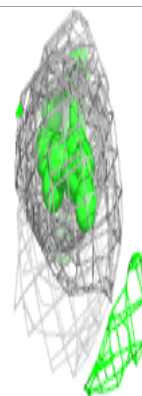
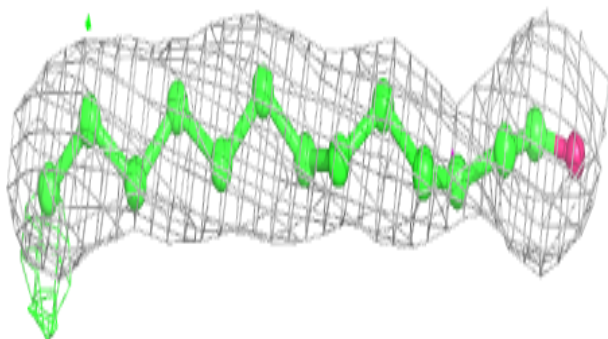
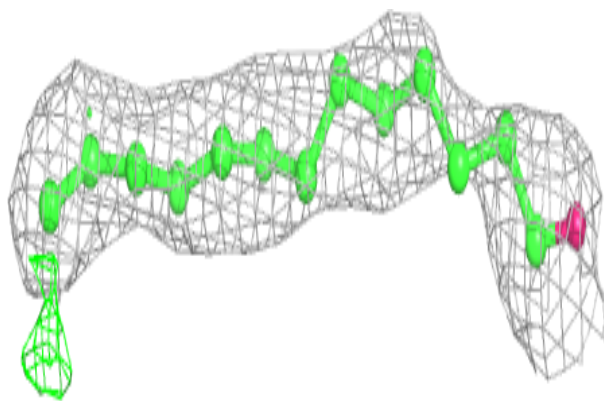
**Electron density around LMT C 603:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
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and green (positive)

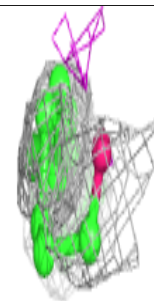
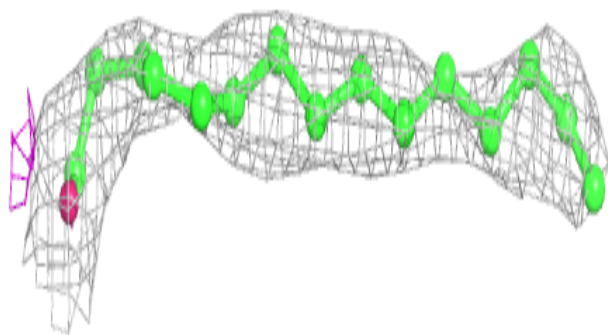
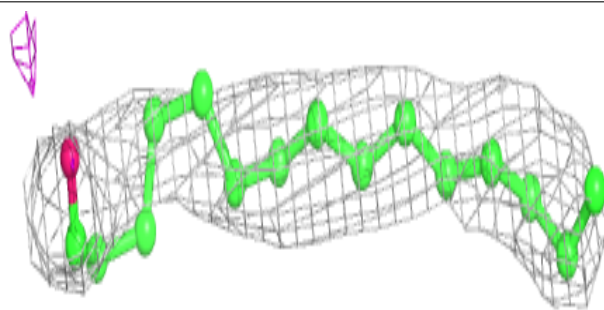


Electron density around PLM C 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
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and green (positive)

**Electron density around PLM D 601:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.