



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 04:08 PM UTC

PDB ID : 6SLI / pdb_00006sli
Title : Structure of the RagAB peptide transporter
Authors : Madej, M.; Ranson, N.A.; White, J.B.R.
Deposited on : 2019-08-19
Resolution : 3.38 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 : 4-5-2 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

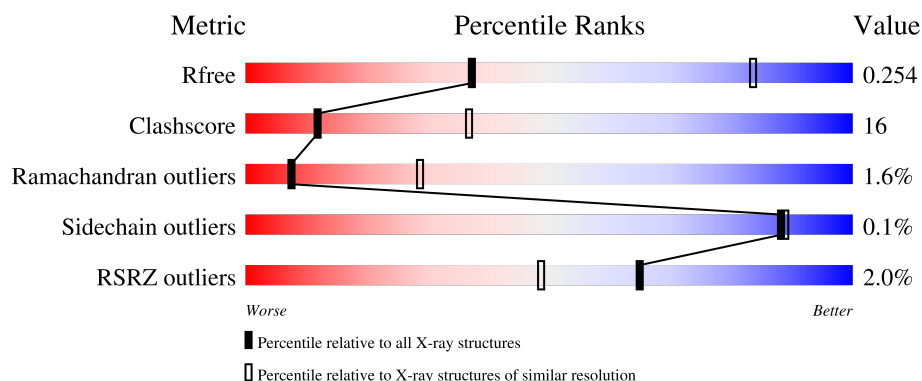
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.38 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	1047 (3.42-3.34)
Clashscore	190562	1067 (3.42-3.34)
Ramachandran outliers	187476	1056 (3.42-3.34)
Sidechain outliers	187428	1056 (3.42-3.34)
RSRZ outliers	180081	1047 (3.42-3.34)

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	488	75% 23% .
1	C	488	76% 22% .
1	F	488	78% 20% .
1	I	488	78% 21% .
2	B	997	55% 33% 10% .

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Mol	Chain	Length	Quality of chain
2	D	997	3% 54% 35% 10%
2	G	997	3% 57% 32% 10%
2	J	997	2% 58% 32% 10%
3	H	13	15% 69% 31%
3	P	13	8% 69% 31%
4	E	13	15% 38% 54% 8%
4	K	13	15% 69% 31%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 43926 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 Lipoprotein RagB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	482	Total	C	N	O	S	0	0	0
			3842	2440	656	737	9			
1	C	482	Total	C	N	O	S	0	0	0
			3842	2440	656	737	9			
1	F	482	Total	C	N	O	S	0	0	0
			3842	2440	656	737	9			
1	I	482	Total	C	N	O	S	0	0	0
			3842	2440	656	737	9			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	502	HIS	-	expression tag	UNP F5H948
A	503	HIS	-	expression tag	UNP F5H948
A	504	HIS	-	expression tag	UNP F5H948
A	505	HIS	-	expression tag	UNP F5H948
A	506	HIS	-	expression tag	UNP F5H948
A	507	HIS	-	expression tag	UNP F5H948
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
F	502	HIS	-	expression tag	UNP F5H948
F	503	HIS	-	expression tag	UNP F5H948
F	504	HIS	-	expression tag	UNP F5H948
F	505	HIS	-	expression tag	UNP F5H948
F	506	HIS	-	expression tag	UNP F5H948
F	507	HIS	-	expression tag	UNP F5H948
I	502	HIS	-	expression tag	UNP F5H948
I	503	HIS	-	expression tag	UNP F5H948
I	504	HIS	-	expression tag	UNP F5H948

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Chain	Residue	Modelled	Actual	Comment	Reference
I	505	HIS	-	expression tag	UNP F5H948
I	506	HIS	-	expression tag	UNP F5H948
I	507	HIS	-	expression tag	UNP F5H948

- Molecule 2 is a protein called RagA protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	900	Total	C	N	O	S	0	0	0
			7048	4466	1179	1371	32			
2	D	900	Total	C	N	O	S	0	0	0
			7048	4466	1179	1371	32			
2	G	900	Total	C	N	O	S	0	0	0
			7048	4466	1179	1371	32			
2	J	900	Total	C	N	O	S	0	0	0
			7048	4466	1179	1371	32			

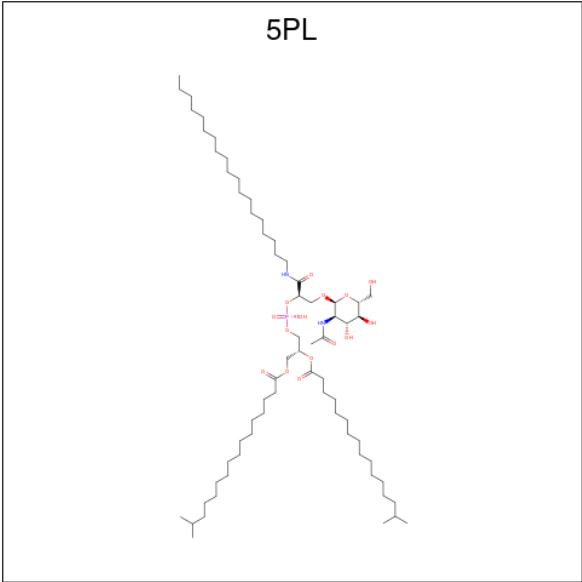
- Molecule 3 is a protein called ALA-SER-THR-THR-GLY-GLY-ASN-SER-GLN-ARG-GLY-SER-GLY.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	13	Total	C	N	O	0	0	0
			81	43	18	20			
3	H	13	Total	C	N	O	0	0	0
			81	43	18	20			

- Molecule 4 is a protein called ASTTGGSQRGGG.

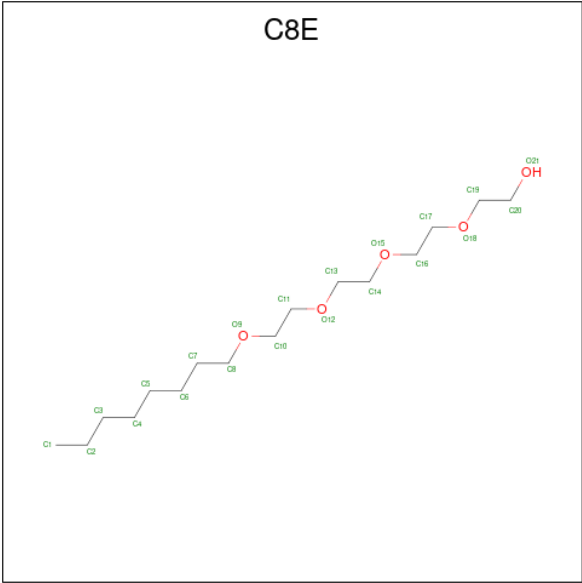
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	13	Total	C	N	O	0	0	0
			79	42	18	19			
4	K	13	Total	C	N	O	0	0	0
			79	42	18	19			

- Molecule 5 is (1R,4S,6R)-6-([2-(ACETYLAMINO)-2-DEOXY-ALPHA-D-GLUCOPYRANOSYL]OXY)METHYL)-4-HYDROXY-1-([(15-METHYLHEXADECANOYL)OXY]METHYL)-4-OXIDO-7-OXO-3,5-DIOXA-8-AZA-4-PHOSPHAHEPTACOS-1-YL 15-METHYLHEXADECANOATE (CCD ID: 5PL) (formula: C₆₇H₁₂₉N₂O₁₅P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			20	16	4		

- Molecule 6 is (HYDROXYETHYLOXY)TRI(ETHYLOXY)OCTANE (CCD ID: C8E) (formula: C₁₆H₃₄O₅).

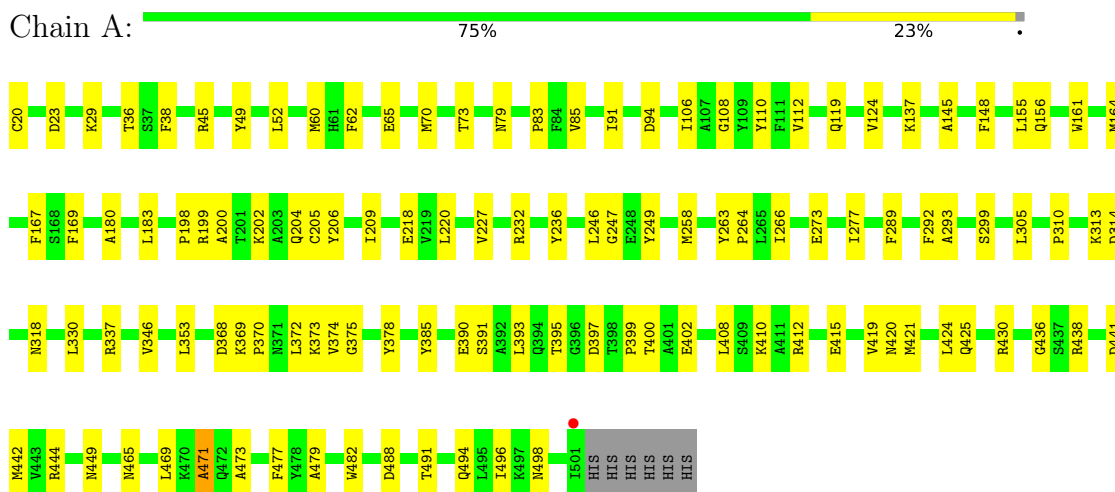


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			13	11	2		
6	B	1	Total	C	O	0	0
			13	11	2		

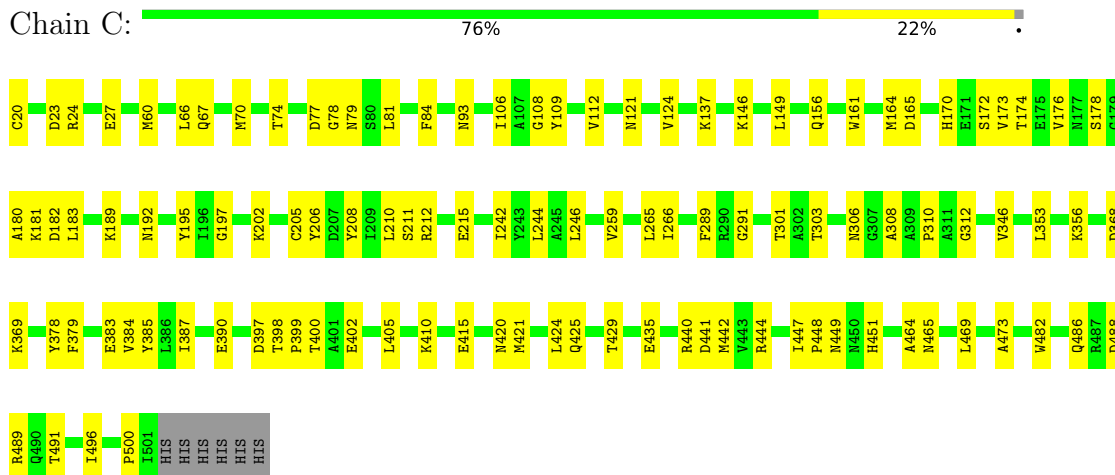
3 Residue-property plots

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.

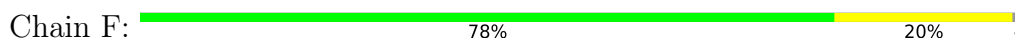
• Molecule 1: Lipoprotein RagB

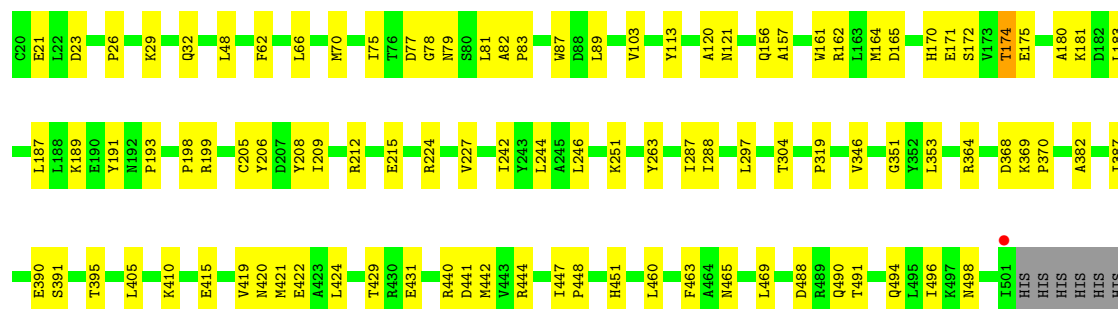


• Molecule 1: Lipoprotein RagB

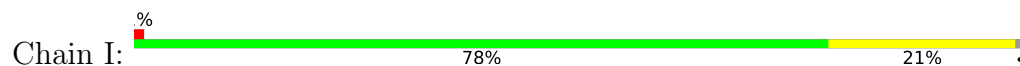


• Molecule 1: Lipoprotein RagB

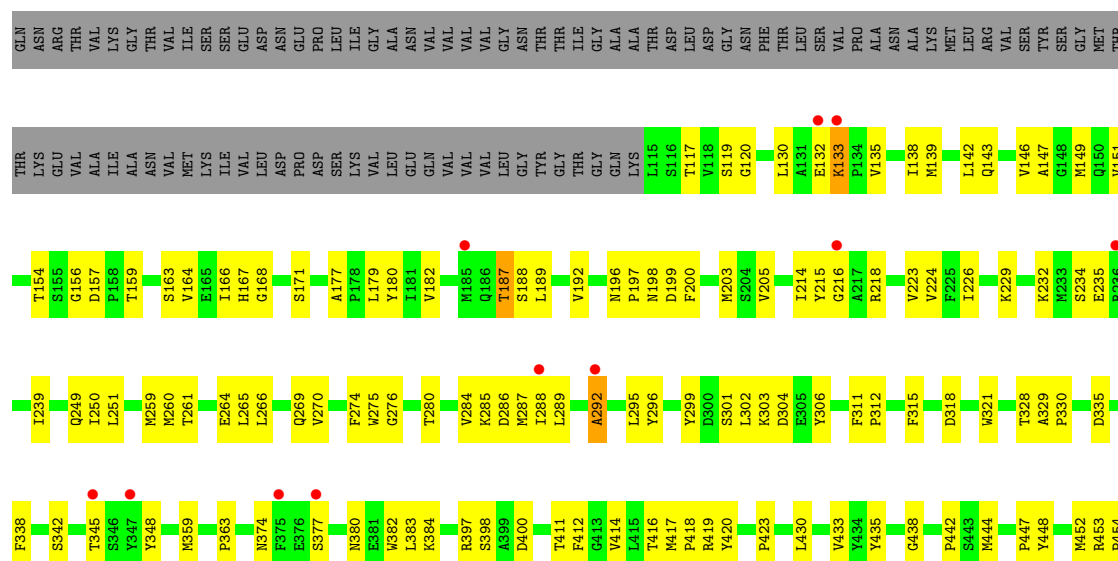


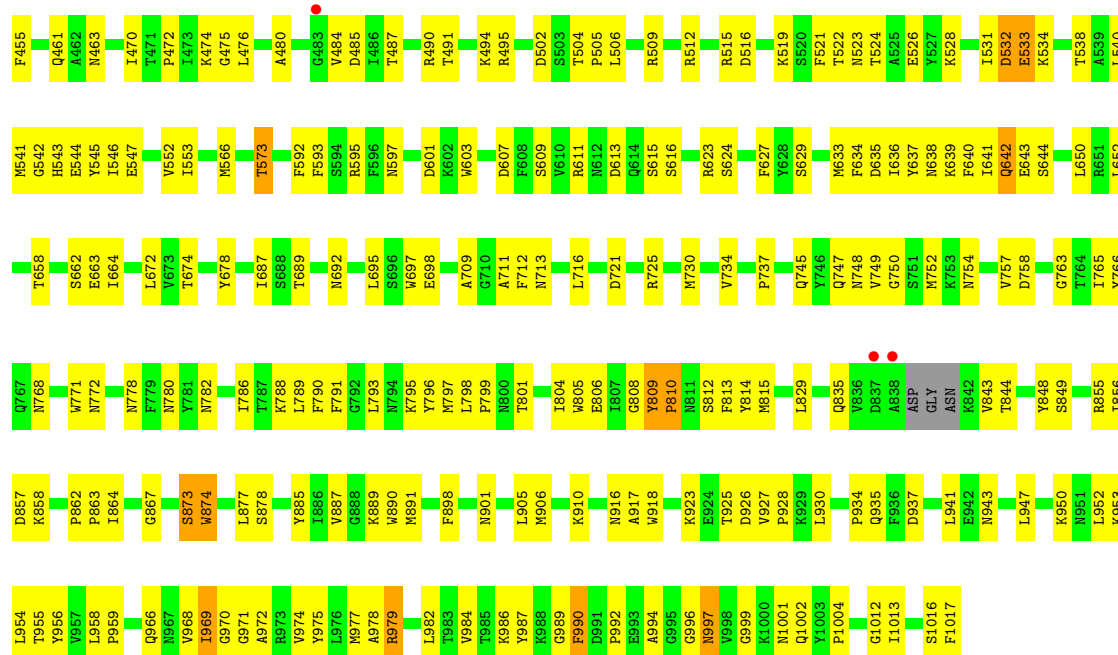


• Molecule 1: Lipoprotein RagB

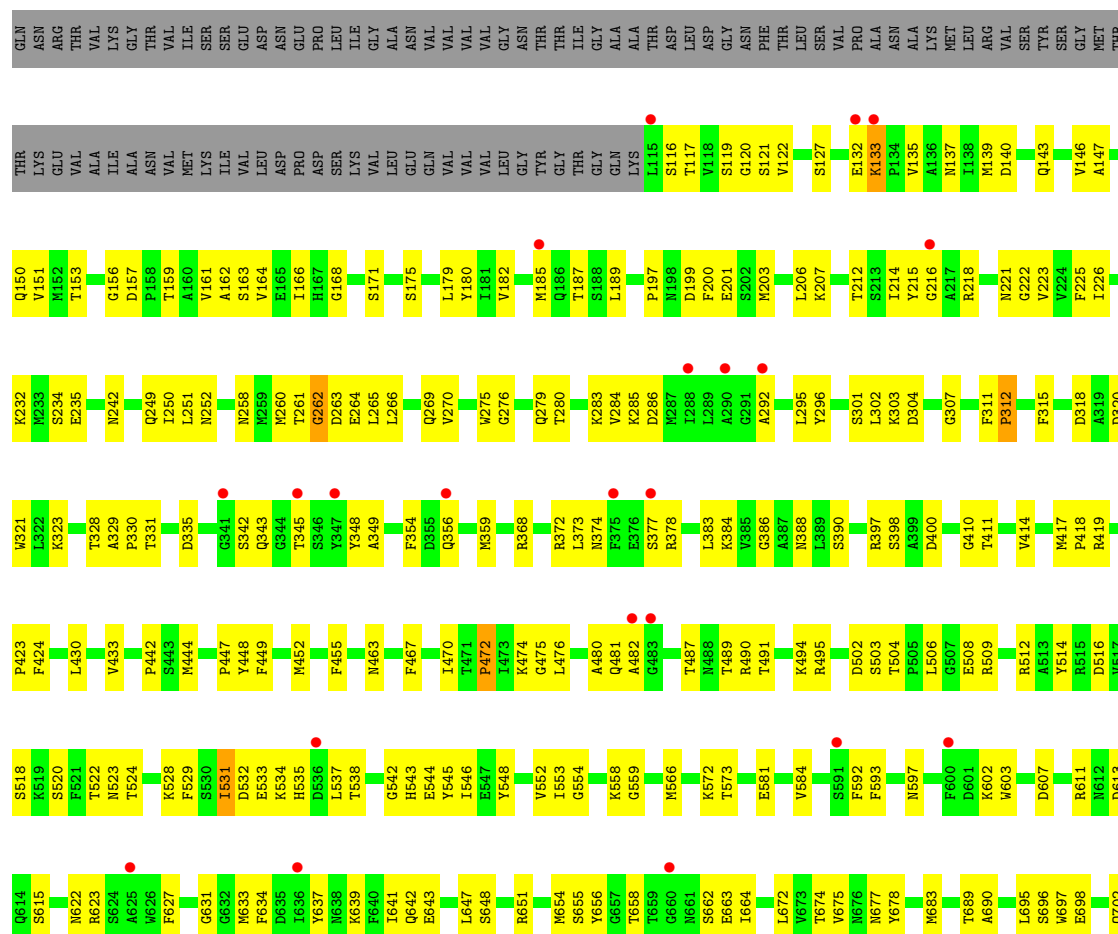


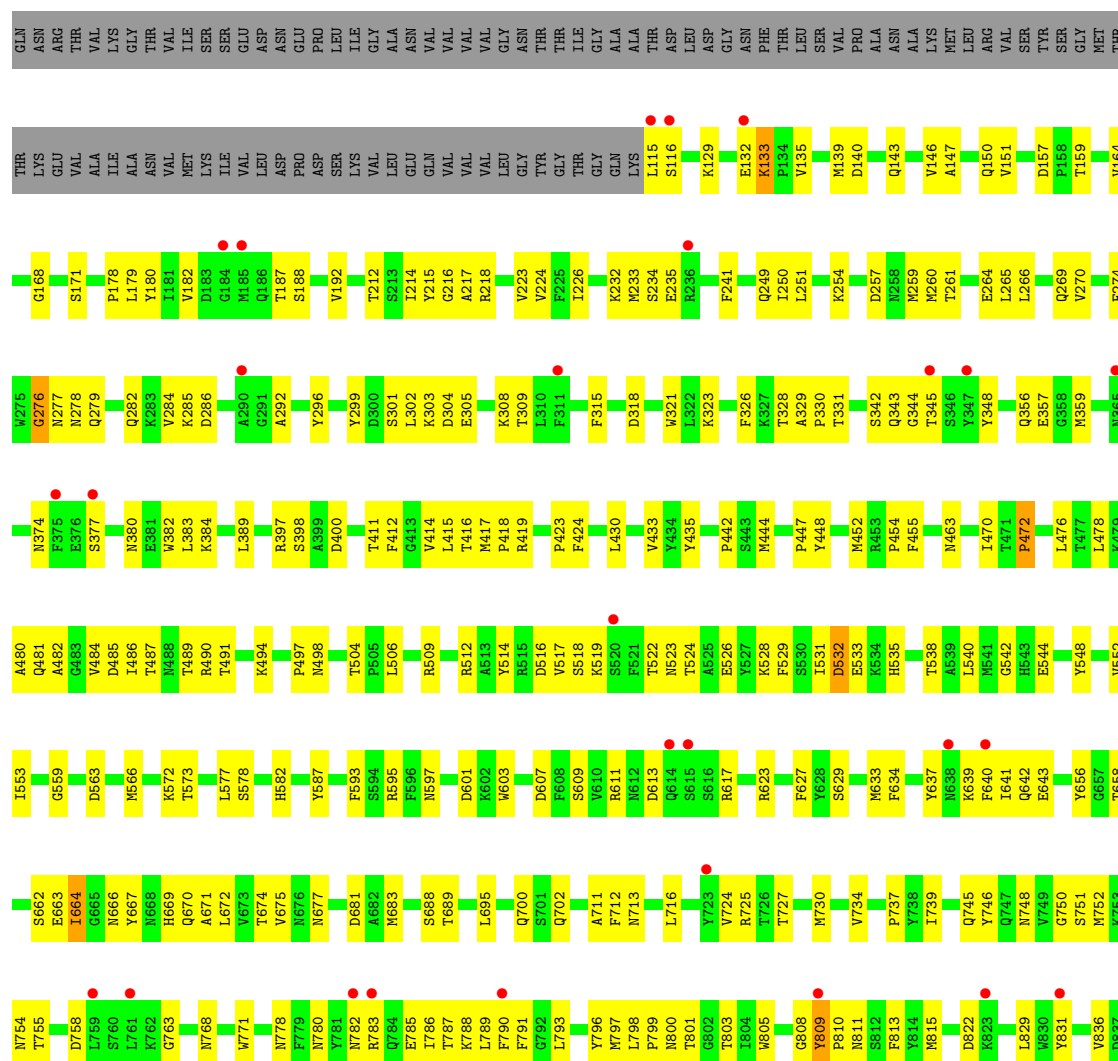
• Molecule 2: RagA protein

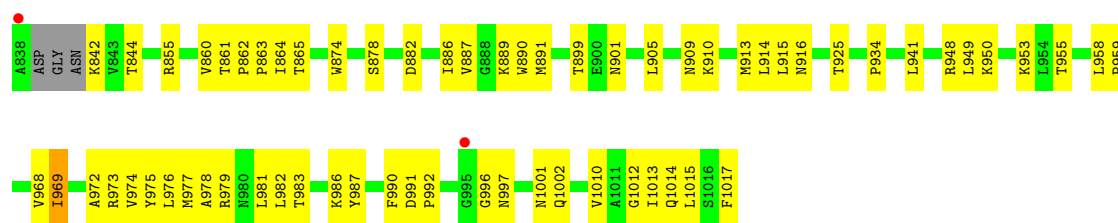




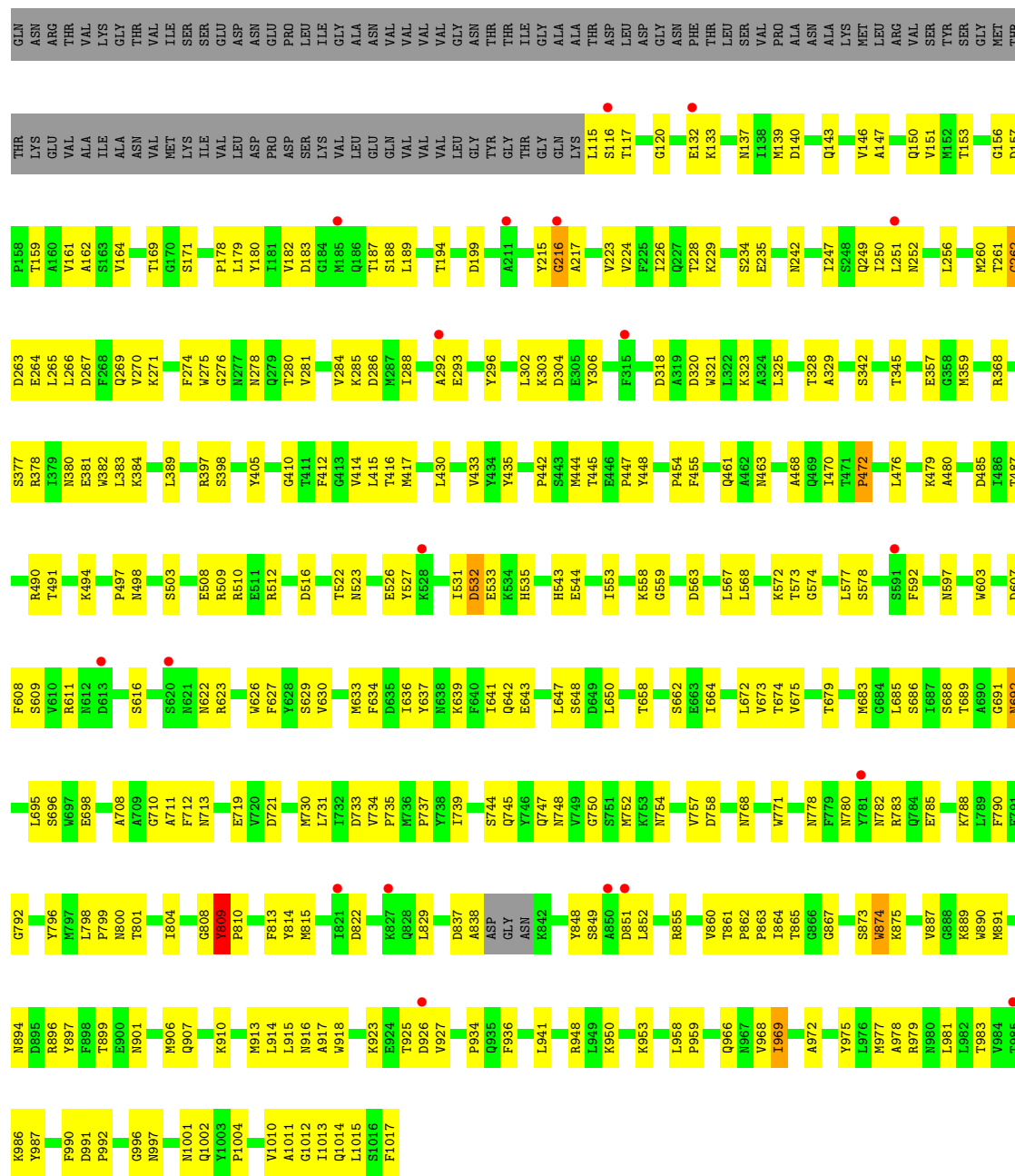
Molecule 2: RagA protein







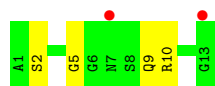
• Molecule 2: RagA protein



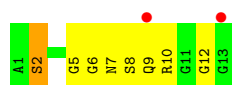
• Molecule 3: ALA-SER-THR-THR-GLY-GLY-ASN-SER-GLN-ARG-GLY-SER-GLY



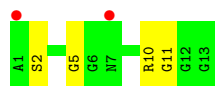
- Molecule 3: ALA-SER-THR-THR-GLY-GLY-ASN-SER-GLN-ARG-GLY-SER-GLY



- Molecule 4: ASTTGGNSQRGGG



- Molecule 4: ASTTGGNSQRGGG



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	190.56Å 376.87Å 369.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	83.92 – 3.38 83.92 – 3.38	Depositor EDS
% Data completeness (in resolution range)	99.7 (83.92-3.38) 99.9 (83.92-3.38)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 3.41Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472, PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.204 , 0.251 0.210 , 0.254	Depositor DCC
R_{free} test set	9224 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	83.0	Xtriage
Anisotropy	0.759	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 125.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	43926	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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: C8E, 5PL

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	0.60	0/3928	0.88	0/5331
1	C	0.59	0/3928	0.86	0/5331
1	F	0.54	0/3928	0.80	0/5331
1	I	0.41	0/3928	0.66	0/5331
2	B	0.59	0/7211	0.93	7/9753 (0.1%)
2	D	0.55	0/7211	0.88	1/9753 (0.0%)
2	G	0.46	0/7211	0.78	0/9753
2	J	0.45	0/7211	0.77	0/9753
3	H	0.46	0/80	0.83	0/105
3	P	0.57	0/80	1.04	0/105
4	E	0.40	0/78	0.82	0/102
4	K	0.51	0/78	0.88	0/102
All	All	0.53	0/44872	0.83	8/60750 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	B	0	1
2	D	0	1
2	J	0	1
All	All	0	4

There are no bond length outliers.

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	475	GLY	N-CA-C	6.90	121.61	112.77
2	B	573	THR	CA-C-N	-5.72	110.20	121.41
2	B	573	THR	C-N-CA	-5.72	110.20	121.41
2	B	990	PHE	CA-C-N	-5.37	112.78	122.28
2	B	990	PHE	C-N-CA	-5.37	112.78	122.28

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	94	ASP	Peptide
2	B	292	ALA	Peptide
2	D	292	ALA	Peptide
2	J	809	TYR	Peptide

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	3842	0	3739	78	0
1	C	3842	0	3740	76	0
1	F	3842	0	3740	82	0
1	I	3842	0	3740	72	0
2	B	7048	0	6792	301	1
2	D	7048	0	6792	287	1
2	G	7048	0	6792	259	0
2	J	7048	0	6792	256	0
3	H	81	0	75	5	0
3	P	81	0	75	4	0
4	E	79	0	73	10	0
4	K	79	0	73	3	0
5	B	20	0	21	4	0
6	B	26	0	42	1	0
All	All	43926	0	42486	1367	2

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 16.

The worst 5 of 1367 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:142:LEU:CD2	2:B:203:MET:HE3	1.74	1.16
1:I:465:ASN:OD1	2:J:276:GLY:N	1.88	1.05
2:B:142:LEU:HD21	2:B:203:MET:CE	1.87	1.04
2:B:754:ASN:HB3	2:B:782:ASN:HD21	1.25	0.98
2:B:250:ILE:HB	2:B:359:MET:HE1	1.48	0.96

All (2) 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)
2:D:342:SER:OG	2:D:342:SER:OG[3_655]	1.97	0.23
2:B:342:SER:OG	2:B:342:SER:OG[3_755]	2.09	0.11

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	480/488 (98%)	439 (92%)	38 (8%)	3 (1%)	21	49
1	C	480/488 (98%)	444 (92%)	33 (7%)	3 (1%)	21	49
1	F	480/488 (98%)	444 (92%)	33 (7%)	3 (1%)	21	49
1	I	480/488 (98%)	444 (92%)	33 (7%)	3 (1%)	21	49
2	B	896/997 (90%)	760 (85%)	118 (13%)	18 (2%)	6	24
2	D	896/997 (90%)	760 (85%)	117 (13%)	19 (2%)	5	24
2	G	896/997 (90%)	768 (86%)	111 (12%)	17 (2%)	6	25
2	J	896/997 (90%)	759 (85%)	118 (13%)	19 (2%)	5	24
3	H	11/13 (85%)	7 (64%)	4 (36%)	0	100	100
3	P	11/13 (85%)	7 (64%)	4 (36%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	E	11/13 (85%)	7 (64%)	3 (27%)	1 (9%)	0	3
4	K	11/13 (85%)	6 (54%)	4 (36%)	1 (9%)	0	3
All	All	5548/5992 (93%)	4845 (87%)	616 (11%)	87 (2%)	7	29

5 of 87 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	642	GLN
2	B	712	PHE
2	B	809	TYR
2	B	874	TRP
2	D	532	ASP

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	402/408 (98%)	402 (100%)	0	100	100
1	C	402/408 (98%)	402 (100%)	0	100	100
1	F	402/408 (98%)	402 (100%)	0	100	100
1	I	402/408 (98%)	402 (100%)	0	100	100
2	B	752/832 (90%)	750 (100%)	2 (0%)	86	84
2	D	752/832 (90%)	752 (100%)	0	100	100
2	G	752/832 (90%)	751 (100%)	1 (0%)	88	89
2	J	752/832 (90%)	752 (100%)	0	100	100
3	H	8/8 (100%)	8 (100%)	0	100	100
3	P	8/8 (100%)	7 (88%)	1 (12%)	4	18
4	E	7/7 (100%)	7 (100%)	0	100	100
4	K	7/7 (100%)	7 (100%)	0	100	100
All	All	4646/4990 (93%)	4642 (100%)	4 (0%)	88	89

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	187	THR
2	B	979	ARG
3	P	12	SER
2	G	886	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 61 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	960	ASN
2	J	754	ASN
2	G	380	ASN
2	J	748	ASN
2	J	951	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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
5	5PL	B	1101	1	19,19,85	1.82	5 (26%)	21,21,101	1.79	4 (19%)
6	C8E	B	1102	-	12,12,20	0.45	0	11,11,19	0.56	0
6	C8E	B	1103	-	12,12,20	0.44	0	11,11,19	0.71	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
5	5PL	B	1101	1	-	9/20/20/107	-
6	C8E	B	1102	-	-	5/10/10/18	-
6	C8E	B	1103	-	-	7/10/10/18	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1101	5PL	OCL-CBL	4.14	1.45	1.33
5	B	1101	5PL	OCK-CBH	-3.42	1.41	1.47
5	B	1101	5PL	OCK-CBU	3.13	1.43	1.34
5	B	1101	5PL	CBM-CBL	2.34	1.57	1.50
5	B	1101	5PL	CBI-CBH	2.31	1.55	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1101	5PL	OCK-CBU-CBV	4.36	120.91	111.48
5	B	1101	5PL	OCL-CBL-CBM	3.72	123.17	111.83
5	B	1101	5PL	CBH-OCK-CBU	-2.72	113.86	117.78
5	B	1101	5PL	OCK-CBH-CBI	2.57	112.13	106.21

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

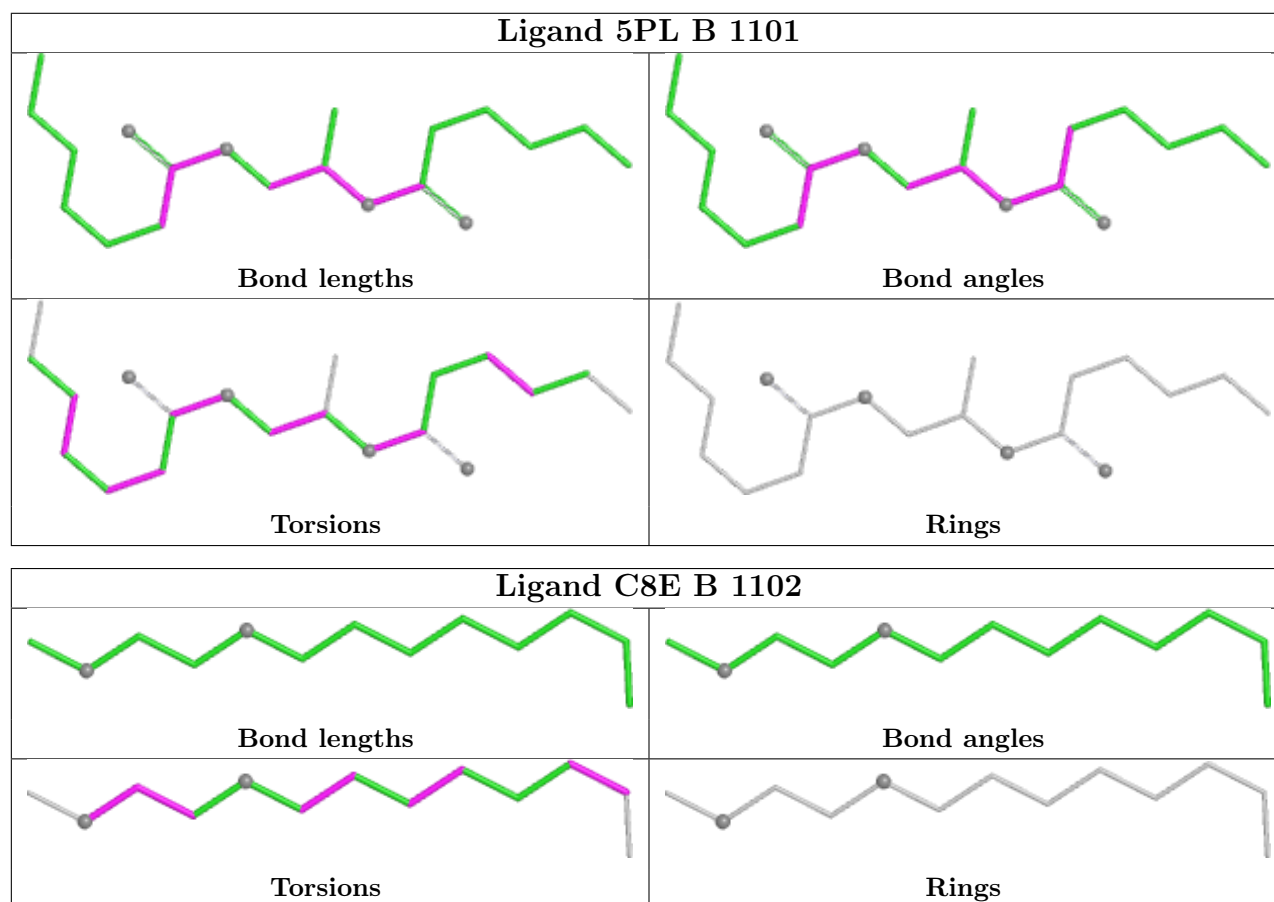
Mol	Chain	Res	Type	Atoms
5	B	1101	5PL	OCK-CBH-CBI-OCL
5	B	1101	5PL	CBV-CBU-OCK-CBH
5	B	1101	5PL	OCN-CBU-OCK-CBH
5	B	1101	5PL	CBM-CBL-OCL-CBI
5	B	1101	5PL	OCO-CBL-OCL-CBI

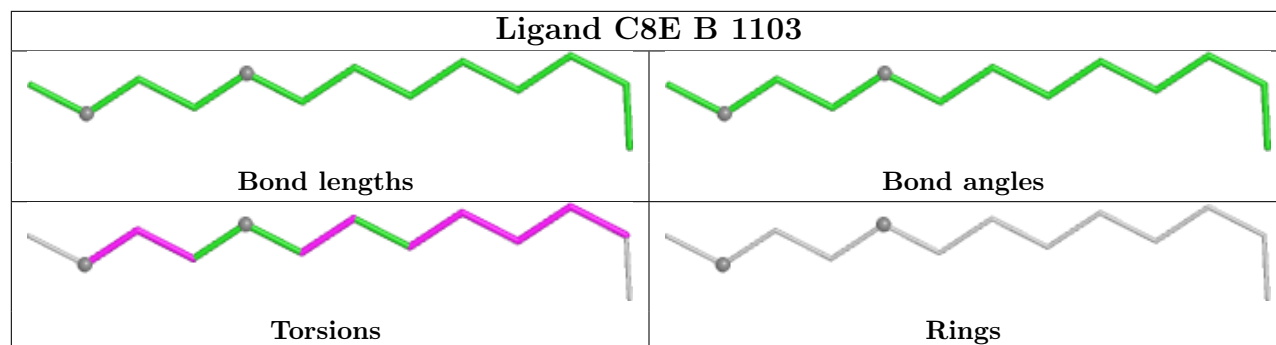
There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1101	5PL	4	0
6	B	1102	C8E	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	482/488 (98%)	-0.50	1 (0%) 91 87	40, 63, 94, 133	0
1	C	482/488 (98%)	-0.50	0 100 100	44, 67, 100, 132	0
1	F	482/488 (98%)	-0.41	1 (0%) 91 87	48, 76, 105, 146	0
1	I	482/488 (98%)	-0.04	4 (0%) 82 70	68, 115, 152, 186	0
2	B	900/997 (90%)	-0.10	14 (1%) 70 55	45, 84, 132, 214	0
2	D	900/997 (90%)	0.25	34 (3%) 44 31	56, 115, 180, 265	0
2	G	900/997 (90%)	0.19	29 (3%) 50 36	53, 115, 182, 253	0
2	J	900/997 (90%)	0.18	19 (2%) 63 47	56, 112, 170, 233	0
3	H	13/13 (100%)	1.01	2 (15%) 5 6	94, 107, 140, 141	0
3	P	13/13 (100%)	0.54	1 (7%) 19 16	74, 93, 108, 121	0
4	E	13/13 (100%)	0.68	2 (15%) 5 6	88, 100, 137, 152	0
4	K	13/13 (100%)	0.84	2 (15%) 5 6	102, 120, 133, 153	0
All	All	5580/5992 (93%)	-0.04	109 (1%) 65 49	40, 98, 161, 265	0

The worst 5 of 109 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	216	GLY	4.7
2	D	347	TYR	4.5
2	G	185	MET	4.3
2	J	216	GLY	4.1
2	D	660	GLY	3.9

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](#)

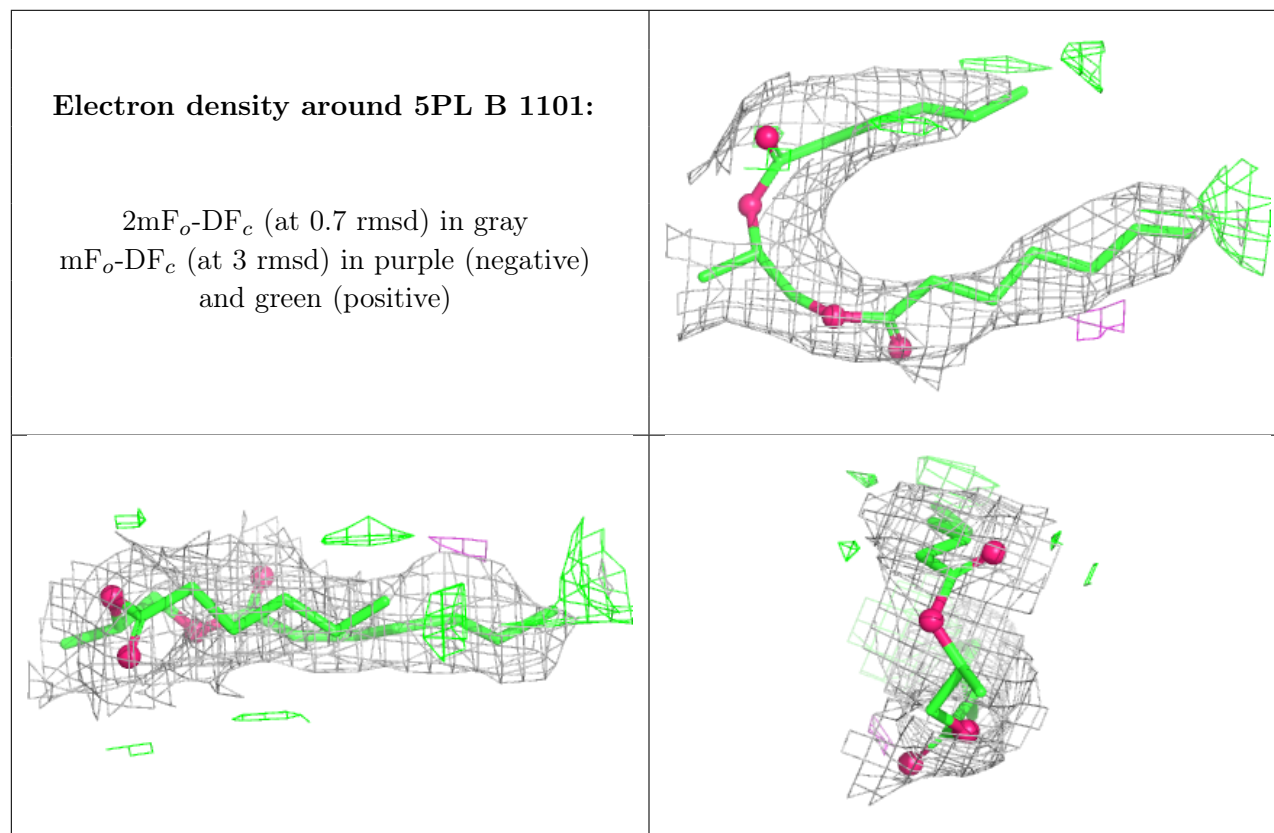
There are no oligosaccharides in this entry.

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.

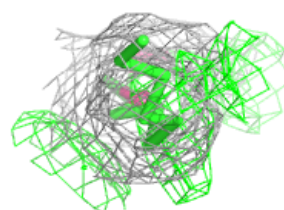
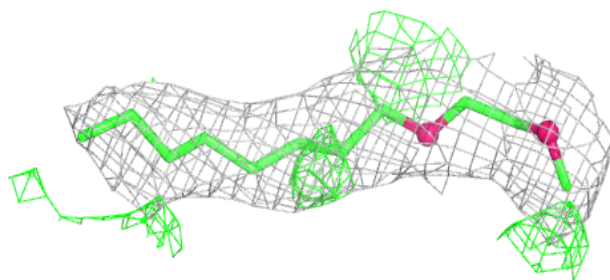
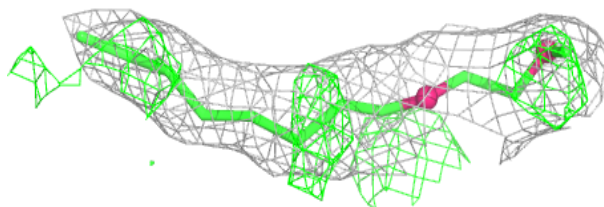
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	5PL	B	1101	20/85	0.81	0.20	87,90,99,100	0
6	C8E	B	1102	13/21	0.94	0.17	66,87,110,110	0
6	C8E	B	1103	13/21	0.95	0.14	36,53,89,92	0

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.

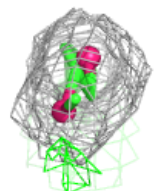
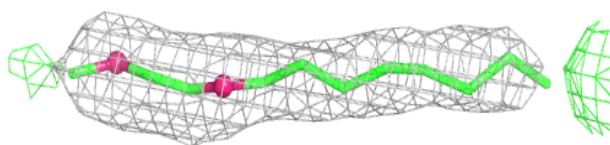
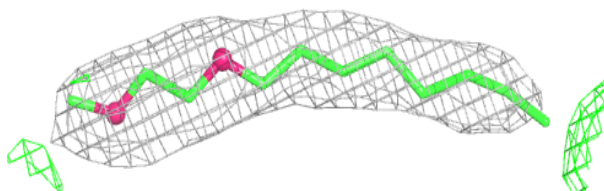


Electron density around C8E B 1102:

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

**Electron density around C8E 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.