



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 13, 2023 – 01:11 AM EDT

PDB ID : 4NMF
Title : Crystal structure of proline utilization A (PutA) from *Geobacter sulfurreducens* PCA inactivated by N-propargylglycine and complexed with menadione bisulfite
Authors : Singh, H.; Tanner, J.J.
Deposited on : 2013-11-14
Resolution : 1.95 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

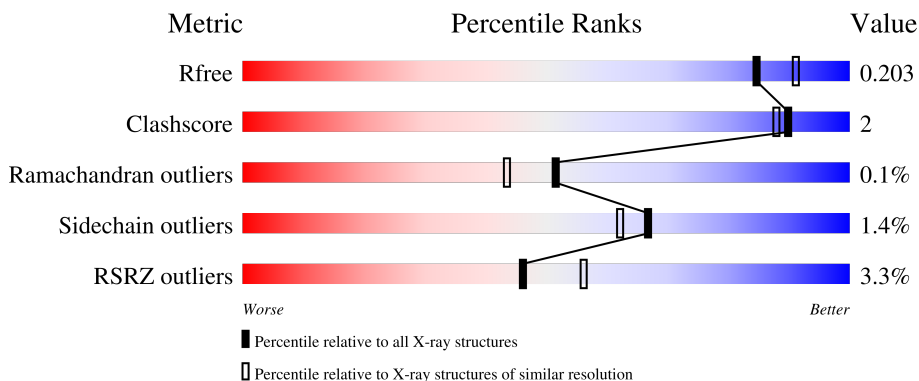
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

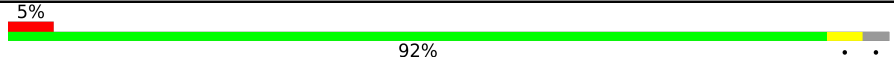
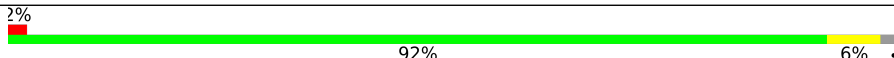
The reported resolution of this entry is 1.95 Å.

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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	1005	
1	B	1005	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 16394 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.

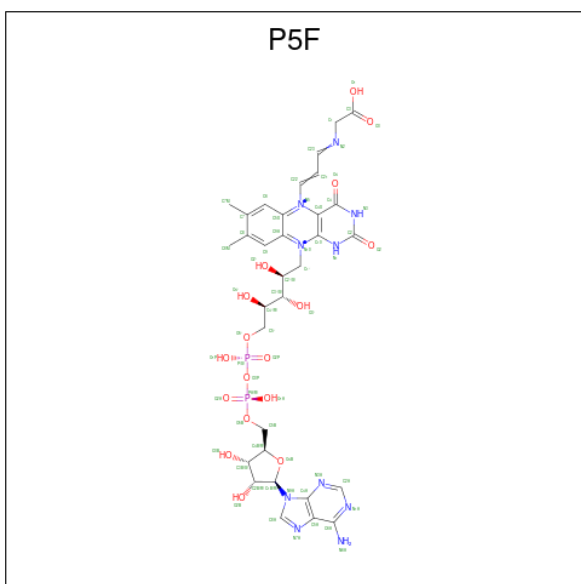
- Molecule 1 is a protein called Proline dehydrogenase and Delta-1-pyrroline-5-carboxylate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	973	Total	C	N	O	S	0	4	0
			7568	4837	1313	1379	39			
1	B	979	Total	C	N	O	S	0	2	0
			7624	4864	1323	1399	38			

There are 2 discrepancies between the modelled and reference sequences:

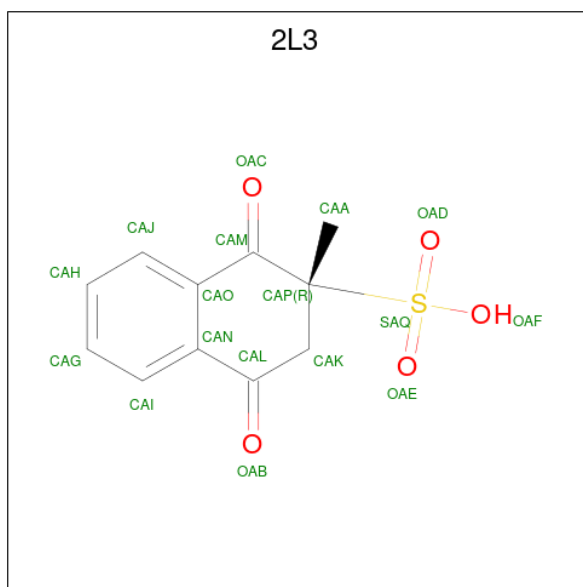
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q746X3
B	0	SER	-	expression tag	UNP Q746X3

- Molecule 2 is N-propargylglycine-modified flavin adenine dinucleotide (three-letter code: P5F) (formula: $C_{32}H_{40}N_{10}O_{17}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	56	30	9	15	2	0	0
2	B	1	56	30	9	15	2	0	0

- Molecule 3 is (2R)-2-methyl-1,4-dioxo-1,2,3,4-tetrahydronaphthalene-2-sulfonic acid (three-letter code: 2L3) (formula: C₁₁H₁₀O₅S).



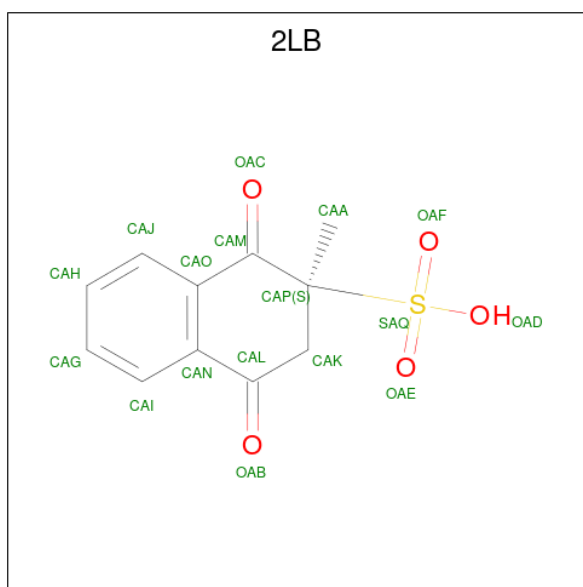
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	S		
3	A	1	17	11	5	1	0	0
3	B	1	17	11	5	1	0	1

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0

- Molecule 5 is (2S)-2-methyl-1,4-dioxo-1,2,3,4-tetrahydronaphthalene-2-sulfonic acid (three-letter code: 2LB) (formula: C₁₁H₁₀O₅S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	O	S	0	1
			17	11	5	1		

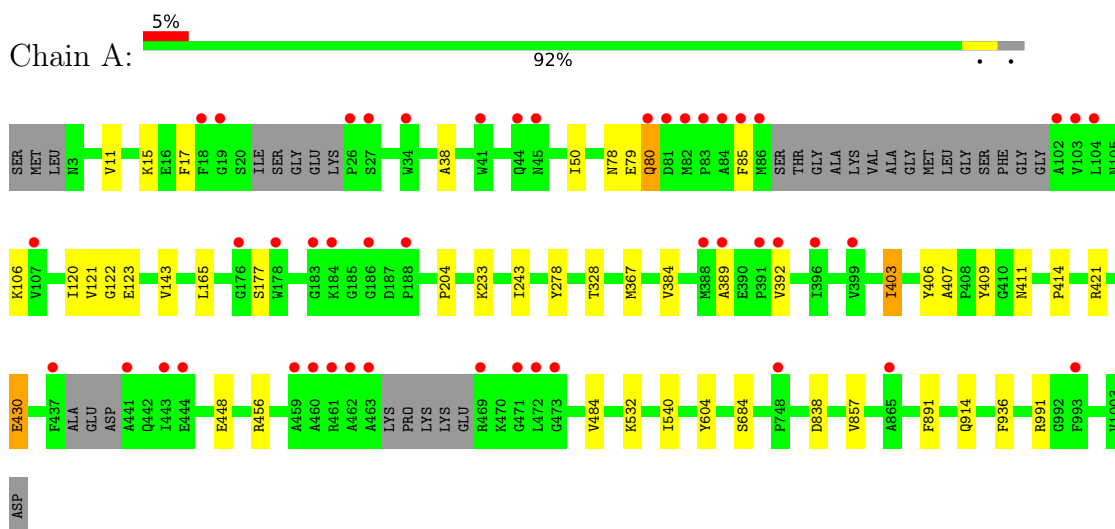
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	497	Total	O	0	0
			497	497		
6	B	510	Total	O	0	0
			510	510		

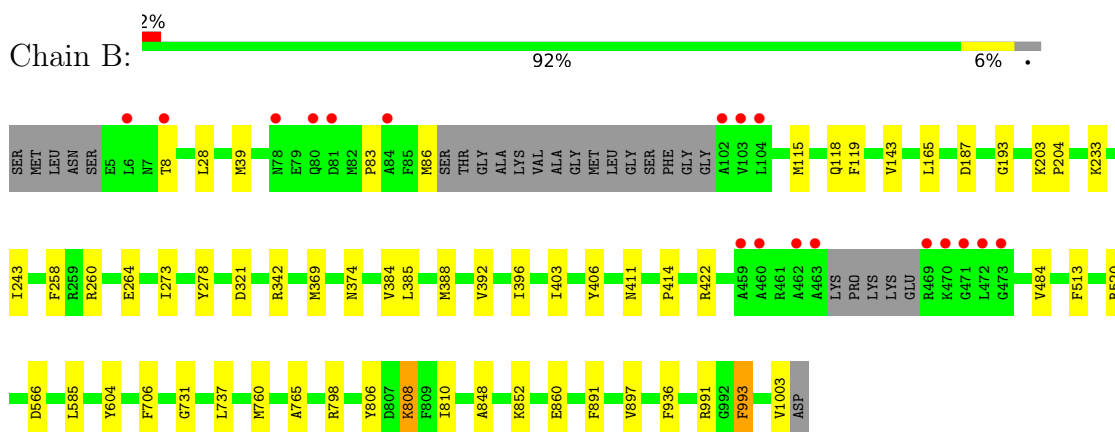
3 Residue-property plots [i](#)

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: Proline dehydrogenase and Delta-1-pyrroline-5-carboxylate dehydrogenase



- Molecule 1: Proline dehydrogenase and Delta-1-pyrroline-5-carboxylate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	95.59Å 151.74Å 176.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.60 – 1.95 47.60 – 1.95	Depositor EDS
% Data completeness (in resolution range)	98.2 (47.60-1.95) 98.2 (47.60-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 1.95Å)	Xtrriage
Refinement program	PHENIX 1.8_1069	Depositor
R, R_{free}	0.173 , 0.203 0.173 , 0.203	Depositor DCC
R_{free} test set	9211 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	24.3	Xtrriage
Anisotropy	0.486	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.3	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	0.96	EDS
Total number of atoms	16394	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.45% 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: EDO, 2L3, 2LB, P5F

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	0.36	0/7751	0.51	0/10500
1	B	0.37	0/7802	0.50	0/10570
All	All	0.36	0/15553	0.51	0/21070

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	7568	0	7425	24	0
1	B	7624	0	7485	32	0
2	A	56	0	35	1	0
2	B	56	0	35	3	0
3	A	17	0	9	1	0
3	B	17	0	9	1	0
4	A	20	0	30	0	0
4	B	12	0	18	0	0
5	B	17	0	9	3	0
6	A	497	0	0	1	0
6	B	510	0	0	1	0
All	All	16394	0	15055	56	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 2.

All (56) 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:422:ARG:NH1	3:B:1102[A]:2L3:OAF	2.25	0.69
1:B:28:LEU:HD22	1:B:39:MET:HG3	1.77	0.66
1:A:389:ALA:HB1	1:A:392:VAL:HB	1.77	0.66
1:A:406:TYR:HB2	2:A:1101:P5F:H18	1.81	0.62
1:B:321:ASP:HB2	6:B:1649:HOH:O	2.00	0.62
1:B:143:VAL:HG13	1:B:406:TYR:HA	1.83	0.61
1:A:120:ILE:HG22	1:A:122:GLY:H	1.66	0.59
1:B:203:LYS:HE3	5:B:1103[B]:2LB:H5	1.85	0.58
1:B:810:ILE:HD11	1:B:897:VAL:HG11	1.86	0.56
1:B:852:LYS:NZ	1:B:860:GLU:OE2	2.39	0.54
2:B:1101:P5F:C1'	5:B:1103[B]:2LB:H10	2.39	0.53
1:B:83:PRO:HD2	1:B:86:MET:HG3	1.90	0.53
1:B:187:ASP:O	1:B:193:GLY:HA2	2.10	0.52
1:A:78:ASN:O	1:A:80:GLN:N	2.43	0.51
1:B:406:TYR:HB2	2:B:1101:P5F:H18	1.92	0.51
1:A:384:VAL:HG22	1:A:403:ILE:HD11	1.92	0.51
1:B:808:LYS:NZ	1:B:808:LYS:HB2	2.25	0.50
1:A:121:VAL:HG11	1:A:407:ALA:HB1	1.93	0.49
1:A:448:GLU:OE2	1:A:456:ARG:NH1	2.45	0.49
1:B:731:GLY:O	1:B:760:MET:HA	2.14	0.48
1:A:17:PHE:CE2	1:A:367:MET:HE1	2.50	0.47
1:A:143:VAL:HG13	1:A:406:TYR:HA	1.96	0.47
1:A:50:ILE:HG12	1:B:993:PHE:HZ	1.81	0.45
1:A:914:GLN:NE2	6:A:1490:HOH:O	2.38	0.45
1:B:385:LEU:HD21	5:B:1103[B]:2LB:H9	2.00	0.44
1:A:484:VAL:HB	1:A:604:TYR:CG	2.53	0.44
1:B:806:TYR:O	1:B:810:ILE:HG12	2.18	0.43
1:A:120:ILE:HG23	1:A:409:TYR:HA	2.00	0.43
1:B:848:ALA:O	1:B:852:LYS:HG3	2.18	0.43
1:A:421:ARG:NE	3:A:1102:2L3:OAF	2.43	0.43
1:B:484:VAL:HB	1:B:604:TYR:CG	2.54	0.43
1:B:258:PHE:HB2	1:B:273:ILE:HD13	2.01	0.43
1:B:991:ARG:HA	1:B:991:ARG:HD3	1.92	0.42
1:B:411:ASN:HB3	1:B:414:PRO:HD2	2.01	0.42
1:B:342:ARG:HB2	1:B:369:MET:SD	2.59	0.42
1:A:11:VAL:O	1:A:15:LYS:HG3	2.19	0.42
2:B:1101:P5F:C4	2:B:1101:P5F:H35	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:765:ALA:HA	1:B:798:ARG:O	2.20	0.42
1:A:328:THR:HG21	1:B:1003:VAL:HG11	2.02	0.42
1:B:165:LEU:HD22	1:B:233:LYS:HG3	2.01	0.42
1:A:411:ASN:HB3	1:A:414:PRO:HD2	2.02	0.41
1:A:165:LEU:HD22	1:A:233:LYS:HG3	2.01	0.41
1:B:384:VAL:HG22	1:B:403:ILE:HD11	2.02	0.41
1:A:204:PRO:HG3	1:A:243:ILE:HG23	2.02	0.41
1:A:484:VAL:HB	1:A:604:TYR:CD1	2.55	0.41
1:B:392:VAL:O	1:B:396:ILE:HG13	2.20	0.41
1:A:38:ALA:HA	1:A:85:PHE:CZ	2.56	0.41
1:A:430:GLU:H	1:A:430:GLU:CD	2.24	0.41
1:A:540:ILE:HG13	1:A:684:SER:HB2	2.03	0.41
1:B:513:PHE:HB3	1:B:706:PHE:HB3	2.03	0.41
1:B:119:PHE:HE1	1:B:388:MET:HB3	1.87	0.41
1:A:991:ARG:HD3	1:A:991:ARG:HA	1.89	0.40
1:B:204:PRO:HG3	1:B:243:ILE:HG23	2.03	0.40
1:B:115:MET:O	1:B:118:GLN:HB3	2.21	0.40
1:B:737:LEU:HD23	1:B:737:LEU:HA	1.90	0.40
1:B:260:ARG:O	1:B:264:GLU:HG3	2.22	0.40

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	Percentiles
1	A	967/1005 (96%)	948 (98%)	17 (2%)	2 (0%)	47 38
1	B	975/1005 (97%)	953 (98%)	22 (2%)	0	100 100
All	All	1942/2010 (97%)	1901 (98%)	39 (2%)	2 (0%)	51 43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	79	GLU
1	A	80	GLN

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	768/821 (94%)	757 (99%)	11 (1%)	67	62
1	B	777/821 (95%)	767 (99%)	10 (1%)	69	65
All	All	1545/1642 (94%)	1524 (99%)	21 (1%)	67	62

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	LYS
1	A	123	GLU
1	A	177	SER
1	A	278	TYR
1	A	403	ILE
1	A	430	GLU
1	A	532	LYS
1	A	838	ASP
1	A	857	VAL
1	A	891	PHE
1	A	936	PHE
1	B	8	THR
1	B	278	TYR
1	B	374	ASN
1	B	520	ARG
1	B	566	ASP
1	B	585	LEU
1	B	808	LYS
1	B	891	PHE
1	B	936	PHE
1	B	993	PHE

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](#)

13 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	$\# Z > 2$
4	EDO	B	1106	-	3,3,3	0.57	0	2,2,2	0.47	0
2	P5F	B	1101	1	52,61,66	2.49	17 (32%)	59,93,99	2.30	11 (18%)
4	EDO	A	1104	-	3,3,3	0.59	0	2,2,2	0.33	0
4	EDO	A	1106	-	3,3,3	0.52	0	2,2,2	0.27	0
3	2L3	B	1102[A]	-	15,18,18	4.35	6 (40%)	18,29,29	1.55	4 (22%)
4	EDO	B	1105	-	3,3,3	0.56	0	2,2,2	0.43	0
4	EDO	A	1107	-	3,3,3	0.59	0	2,2,2	0.35	0
4	EDO	A	1103	-	3,3,3	0.51	0	2,2,2	0.43	0
4	EDO	B	1104	-	3,3,3	0.54	0	2,2,2	0.12	0
5	2LB	B	1103[B]	-	15,18,18	4.32	6 (40%)	18,29,29	1.63	6 (33%)
2	P5F	A	1101	1	52,61,66	2.50	18 (34%)	59,93,99	2.35	13 (22%)
3	2L3	A	1102	-	15,18,18	4.30	6 (40%)	18,29,29	1.50	3 (16%)
4	EDO	A	1105	-	3,3,3	0.45	0	2,2,2	0.32	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	1106	-	-	1/1/1/1	-
2	P5F	B	1101	1	-	6/31/53/58	0/6/6/6
4	EDO	A	1104	-	-	1/1/1/1	-
4	EDO	A	1106	-	-	1/1/1/1	-
3	2L3	B	1102[A]	-	-	2/2/28/28	0/2/2/2
4	EDO	B	1105	-	-	0/1/1/1	-
4	EDO	A	1107	-	-	1/1/1/1	-
4	EDO	A	1103	-	-	0/1/1/1	-
4	EDO	B	1104	-	-	0/1/1/1	-
5	2LB	B	1103[B]	-	-	2/2/28/28	0/2/2/2
2	P5F	A	1101	1	-	2/31/53/58	0/6/6/6
3	2L3	A	1102	-	-	0/2/28/28	0/2/2/2
4	EDO	A	1105	-	-	0/1/1/1	-

All (53) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1102[A]	2L3	OAC-CAM	9.81	1.37	1.21
5	B	1103[B]	2LB	OAC-CAM	9.72	1.37	1.21
3	A	1102	2L3	OAB-CAL	9.67	1.37	1.22
3	A	1102	2L3	OAC-CAM	9.62	1.37	1.21
5	B	1103[B]	2LB	OAB-CAL	9.59	1.37	1.22
3	B	1102[A]	2L3	OAB-CAL	9.55	1.37	1.22
2	B	1101	P5F	O4-C4	7.30	1.38	1.23
2	A	1101	P5F	O4-C4	7.20	1.37	1.23
2	B	1101	P5F	O2-C2	6.92	1.37	1.23
2	A	1101	P5F	O2-C2	6.70	1.37	1.23
3	B	1102[A]	2L3	CAI-CAN	5.82	1.49	1.39
5	B	1103[B]	2LB	CAI-CAN	5.70	1.49	1.39
2	B	1101	P5F	C4X-N5	5.50	1.45	1.37
3	A	1102	2L3	CAI-CAN	5.47	1.48	1.39
3	A	1102	2L3	CAJ-CAO	5.40	1.48	1.39
3	B	1102[A]	2L3	CAJ-CAO	5.34	1.48	1.39
2	A	1101	P5F	C4X-N5	5.17	1.44	1.37
5	B	1103[B]	2LB	CAJ-CAO	5.15	1.48	1.39
2	A	1101	P5F	C9A-N10	4.60	1.46	1.40
2	A	1101	P5F	C6-C5X	4.56	1.49	1.40
2	B	1101	P5F	C5X-N5	4.45	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1103[B]	2LB	CAO-CAN	-4.24	1.33	1.40
2	B	1101	P5F	C6-C5X	4.09	1.48	1.40
2	A	1101	P5F	C5X-N5	4.02	1.44	1.38
3	B	1102[A]	2L3	CAO-CAN	-3.93	1.34	1.40
2	B	1101	P5F	C2-N1	3.93	1.44	1.36
2	A	1101	P5F	C9-C9A	3.86	1.48	1.40
3	B	1102[A]	2L3	CAA-CAP	-3.83	1.49	1.53
2	B	1101	P5F	C9A-N10	3.74	1.45	1.40
3	A	1102	2L3	CAA-CAP	-3.72	1.49	1.53
3	A	1102	2L3	CAO-CAN	-3.64	1.34	1.40
2	A	1101	P5F	C2-N1	3.62	1.43	1.36
5	B	1103[B]	2LB	CAA-CAP	-3.54	1.49	1.53
2	B	1101	P5F	C9-C9A	3.48	1.47	1.40
2	A	1101	P5F	C10-N10	3.44	1.43	1.36
2	A	1101	P5F	C2B-C1B	-3.40	1.48	1.53
2	A	1101	P5F	C2A-N3A	3.27	1.37	1.32
2	B	1101	P5F	C2B-C1B	-3.20	1.48	1.53
2	B	1101	P5F	C9A-C5X	-3.19	1.35	1.42
2	B	1101	P5F	C10-N10	2.95	1.42	1.36
2	A	1101	P5F	C6A-N6A	2.94	1.44	1.34
2	B	1101	P5F	C2A-N3A	2.91	1.36	1.32
2	B	1101	P5F	C6A-N6A	2.87	1.44	1.34
2	A	1101	P5F	C9A-C5X	-2.84	1.36	1.42
2	B	1101	P5F	PA-O5B	-2.58	1.48	1.59
2	A	1101	P5F	PA-O5B	-2.43	1.49	1.59
2	A	1101	P5F	C2-N3	2.20	1.41	1.37
2	B	1101	P5F	C2-N3	2.12	1.41	1.37
2	A	1101	P5F	O3B-C3B	-2.09	1.38	1.43
2	B	1101	P5F	O3B-C3B	-2.09	1.38	1.43
2	B	1101	P5F	P-O2P	2.05	1.58	1.50
2	A	1101	P5F	P-O2P	2.05	1.58	1.50
2	A	1101	P5F	C9-C8	2.01	1.42	1.37

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	P5F	N1-C2-N3	8.52	120.85	115.90
2	B	1101	P5F	N1-C2-N3	8.48	120.82	115.90
2	A	1101	P5F	C4-N3-C2	-8.36	120.28	126.88
2	B	1101	P5F	C4-N3-C2	-7.22	121.18	126.88
2	A	1101	P5F	C1'-N10-C9A	6.67	123.54	118.29
2	B	1101	P5F	C10-N1-C2	-6.63	117.74	128.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1101	P5F	C1'-N10-C9A	6.18	123.16	118.29
2	A	1101	P5F	C10-N1-C2	-5.79	119.06	128.11
2	A	1101	P5F	N3A-C2A-N1A	-5.26	120.45	128.68
2	B	1101	P5F	N3A-C2A-N1A	-4.84	121.11	128.68
5	B	1103[B]	2LB	CAK-CAL-CAN	3.61	121.48	116.49
3	A	1102	2L3	OAE-SAQ-OAD	-3.57	103.82	117.56
3	B	1102[A]	2L3	CAK-CAL-CAN	3.24	120.97	116.49
5	B	1103[B]	2LB	OAE-SAQ-OAF	-2.93	106.27	117.56
3	B	1102[A]	2L3	OAE-SAQ-OAD	-2.89	106.42	117.56
2	B	1101	P5F	O5B-PA-O2A	-2.82	98.05	109.07
3	A	1102	2L3	OAF-SAQ-OAE	-2.80	102.97	112.91
5	B	1103[B]	2LB	OAD-SAQ-OAE	-2.68	103.40	112.91
2	A	1101	P5F	P-O3P-PA	-2.66	123.70	132.83
2	A	1101	P5F	C1B-N9A-C4A	-2.55	122.15	126.64
3	B	1102[A]	2L3	OAF-SAQ-OAD	-2.55	103.88	112.91
3	B	1102[A]	2L3	OAF-SAQ-OAE	-2.53	103.95	112.91
2	A	1101	P5F	C6-C5X-N5	-2.49	118.80	121.68
2	A	1101	P5F	C6-C5X-C9A	2.45	121.55	119.06
3	A	1102	2L3	CAK-CAL-CAN	2.43	119.85	116.49
2	B	1101	P5F	O1P-P-O5'	2.42	119.00	107.75
2	A	1101	P5F	O5B-PA-O2A	-2.42	99.60	109.07
2	A	1101	P5F	O5'-P-O2P	-2.36	99.85	109.07
2	A	1101	P5F	O2-C2-N1	-2.34	118.70	122.08
5	B	1103[B]	2LB	OAC-CAM-CAO	-2.34	118.84	122.31
2	A	1101	P5F	O1P-P-O5'	2.27	118.27	107.75
2	B	1101	P5F	O3'-C3'-C4'	-2.26	103.35	108.81
2	B	1101	P5F	O5'-C5'-C4'	-2.24	103.39	109.36
5	B	1103[B]	2LB	OAB-CAL-CAN	-2.21	119.14	121.71
5	B	1103[B]	2LB	OAD-SAQ-OAF	-2.15	105.27	112.91
2	B	1101	P5F	O5'-P-O2P	-2.14	100.72	109.07
2	B	1101	P5F	O1A-PA-O5B	2.06	117.30	107.75

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1101	P5F	C5'-O5'-P-O2P
3	B	1102[A]	2L3	CAK-CAP-SAQ-OAD
3	B	1102[A]	2L3	CAK-CAP-SAQ-OAE
4	B	1106	EDO	O1-C1-C2-O2
5	B	1103[B]	2LB	CAK-CAP-SAQ-OAF
5	B	1103[B]	2LB	CAK-CAP-SAQ-OAE

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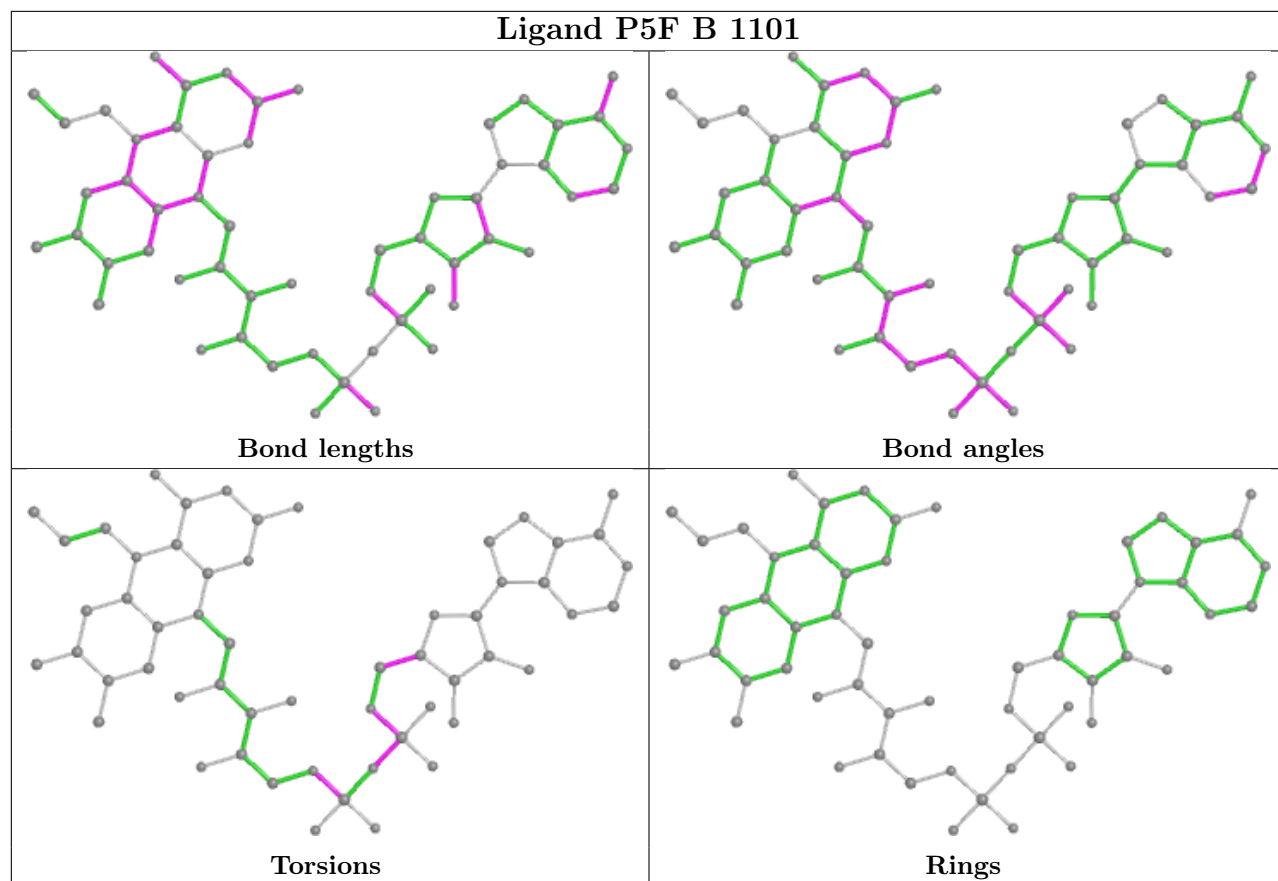
Mol	Chain	Res	Type	Atoms
4	A	1106	EDO	O1-C1-C2-O2
2	A	1101	P5F	C2'-C3'-C4'-O4'
2	B	1101	P5F	P-O3P-PA-O5B
2	B	1101	P5F	C5'-O5'-P-O3P
2	B	1101	P5F	C5'-O5'-P-O1P
4	A	1107	EDO	O1-C1-C2-O2
2	A	1101	P5F	O3'-C3'-C4'-O4'
2	B	1101	P5F	C3B-C4B-C5B-O5B
4	A	1104	EDO	O1-C1-C2-O2
2	B	1101	P5F	C5B-O5B-PA-O3P

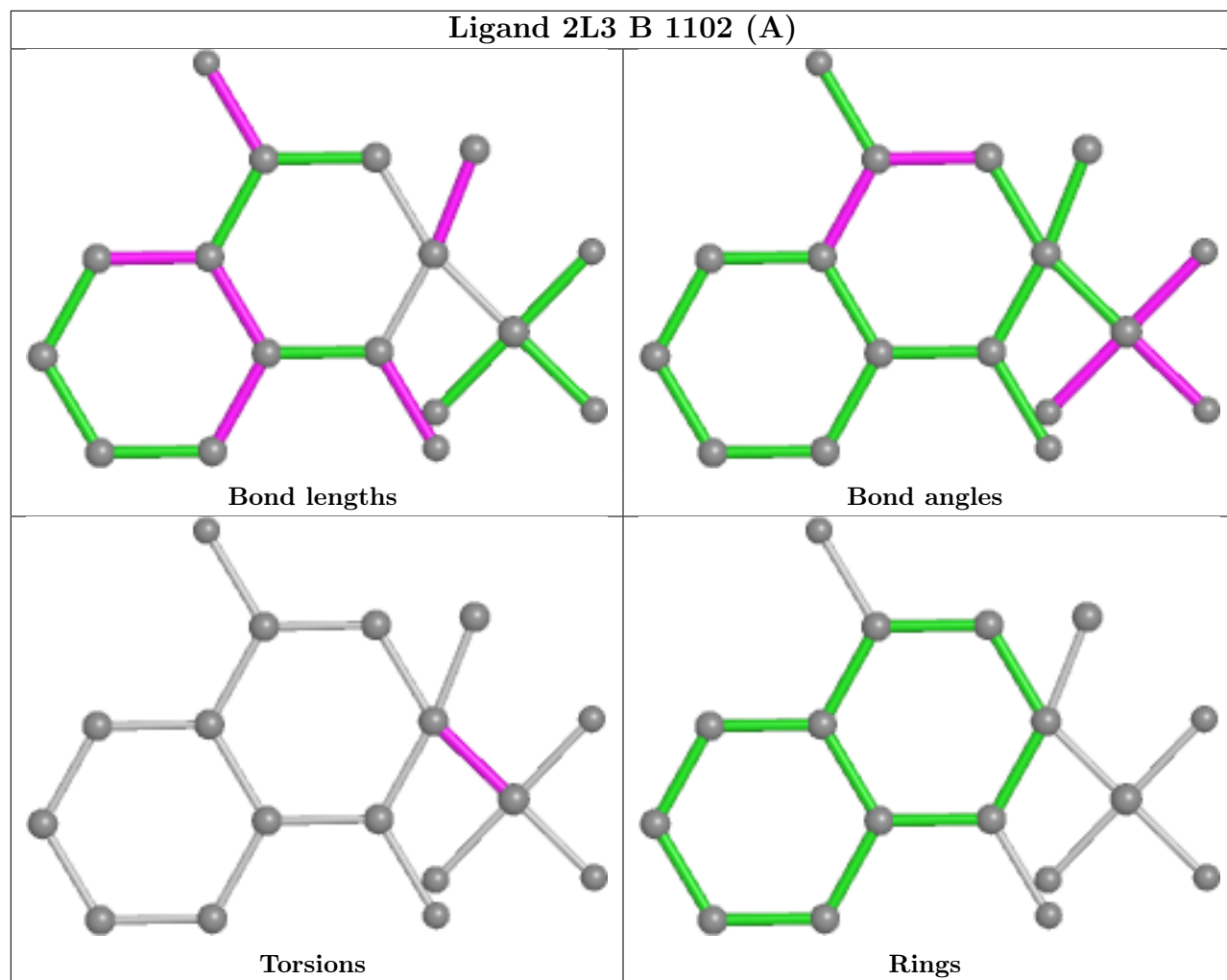
There are no ring outliers.

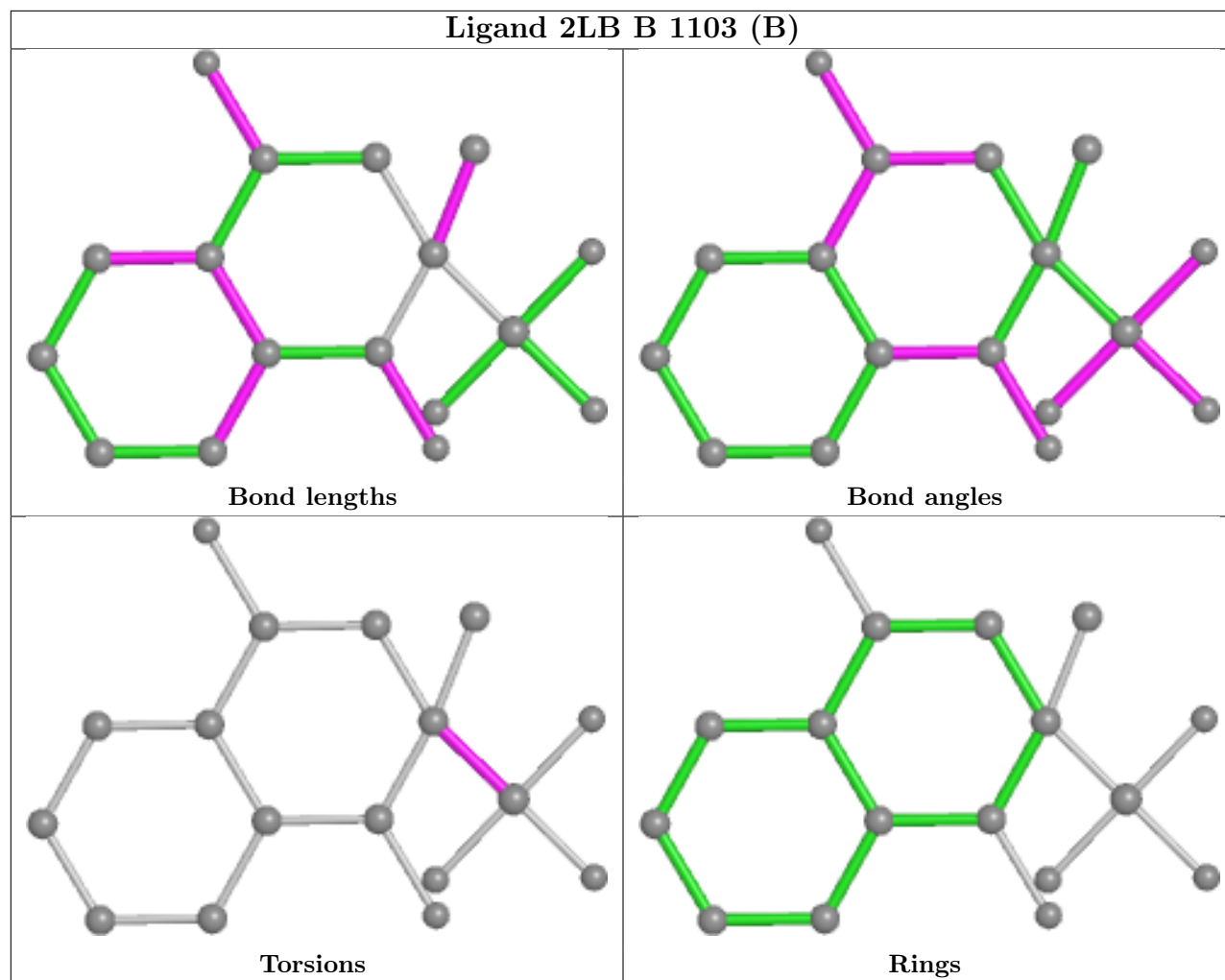
5 monomers are involved in 8 short contacts:

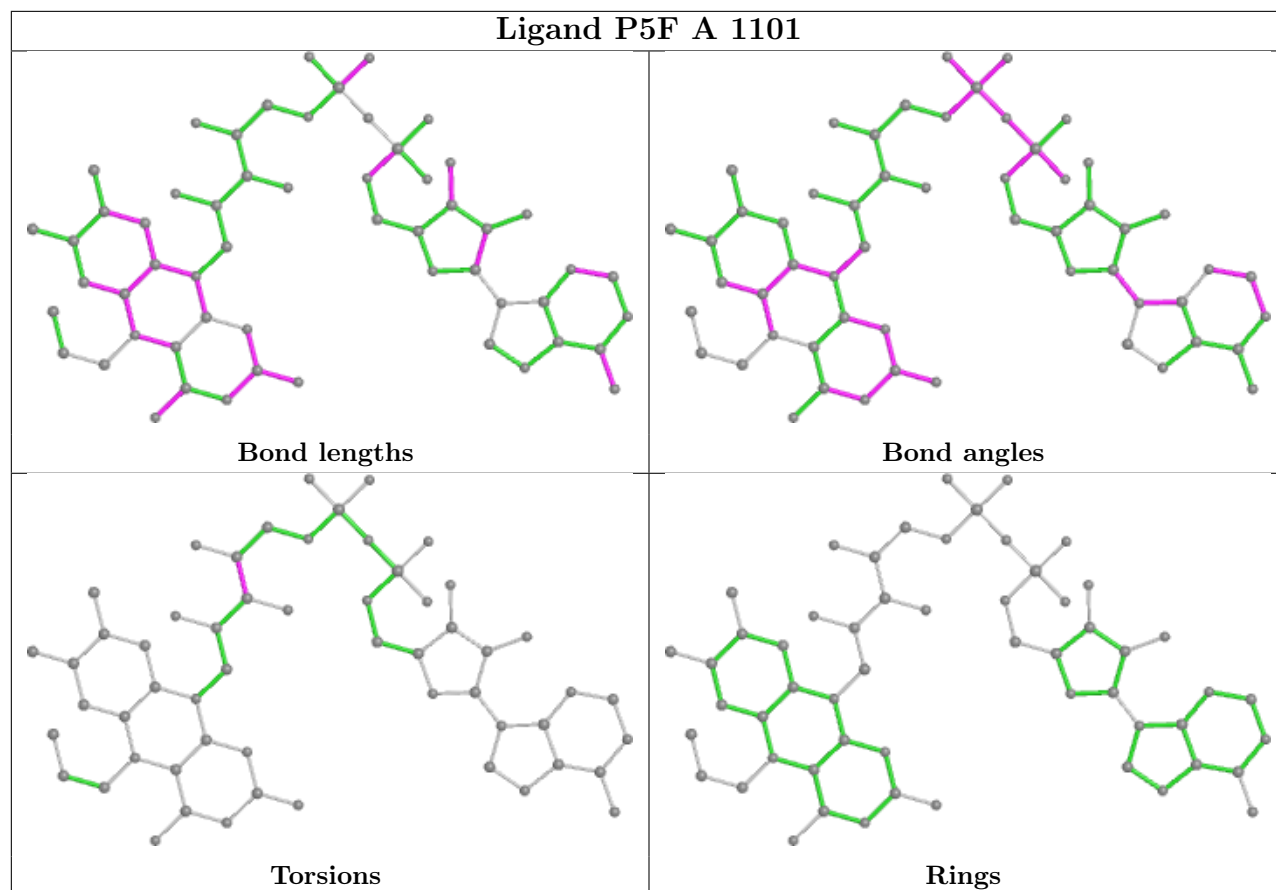
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1101	P5F	3	0
3	B	1102[A]	2L3	1	0
5	B	1103[B]	2LB	3	0
2	A	1101	P5F	1	0
3	A	1102	2L3	1	0

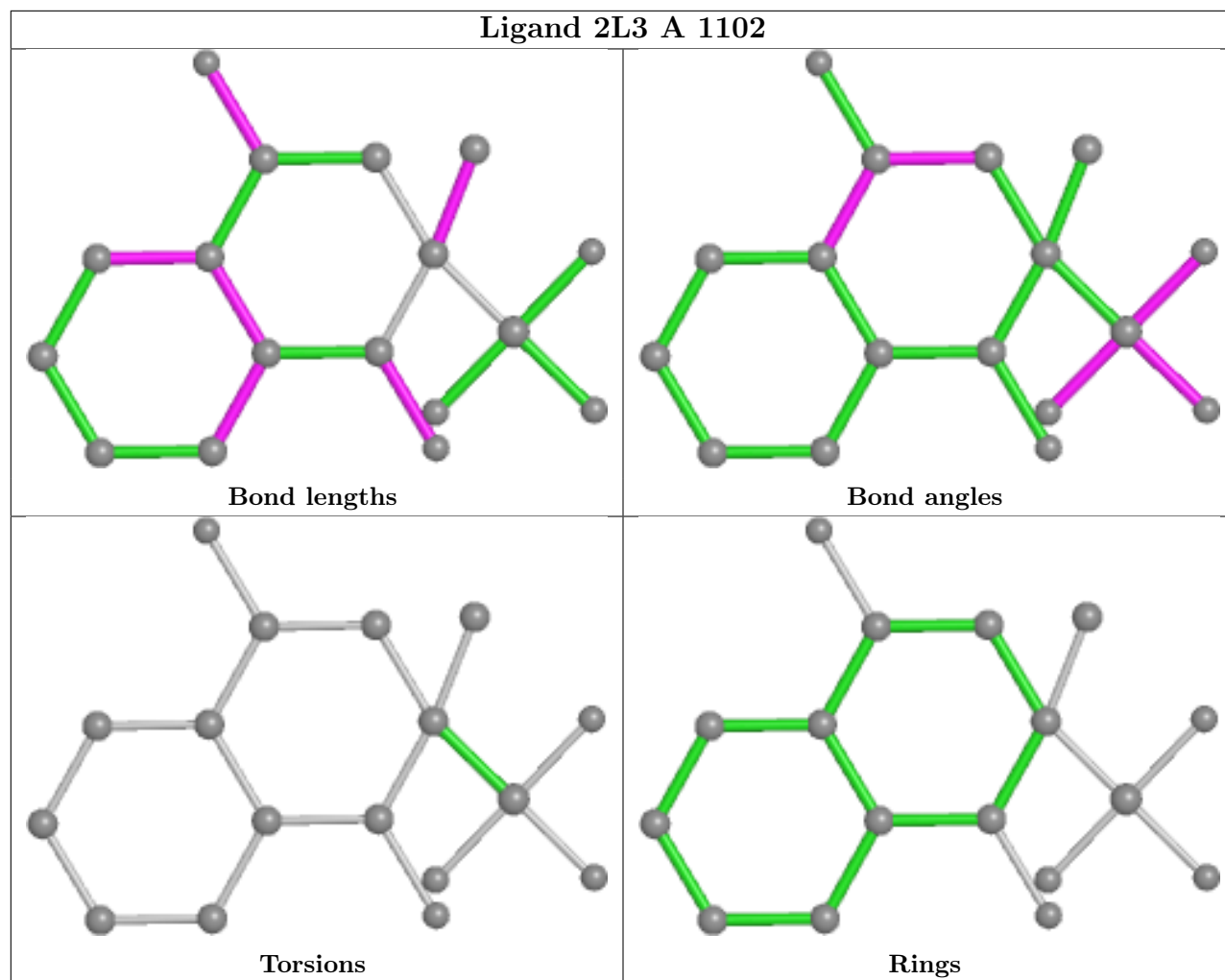
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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

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	973/1005 (96%)	0.10	47 (4%) 30 40	10, 19, 47, 64	0
1	B	979/1005 (97%)	-0.05	18 (1%) 68 76	9, 18, 37, 61	0
All	All	1952/2010 (97%)	0.02	65 (3%) 46 56	9, 19, 42, 64	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	85	PHE	5.7
1	A	103	VAL	5.7
1	A	392	VAL	5.3
1	A	102	ALA	5.1
1	A	389	ALA	4.7
1	A	86	MET	4.7
1	A	84	ALA	4.5
1	A	460	ALA	4.4
1	A	748	PRO	4.3
1	A	104	LEU	4.1
1	B	463	ALA	4.1
1	A	107	VAL	4.0
1	A	82	MET	4.0
1	B	103	VAL	3.8
1	A	473	GLY	3.8
1	B	471	GLY	3.6
1	A	178	TRP	3.6
1	B	462	ALA	3.6
1	A	462	ALA	3.5
1	B	80	GLN	3.4
1	A	441	ALA	3.3
1	A	19	GLY	3.3
1	A	463	ALA	3.2
1	B	104	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	81	ASP	3.1
1	A	83	PRO	3.1
1	B	459	ALA	3.1
1	A	471	GLY	3.1
1	A	388	MET	3.0
1	A	865	ALA	2.9
1	B	84	ALA	2.9
1	A	41	TRP	2.9
1	A	472	LEU	2.9
1	A	81	ASP	2.9
1	A	18	PHE	2.9
1	A	80	GLN	2.8
1	A	176	GLY	2.7
1	B	473	GLY	2.7
1	A	469	ARG	2.7
1	A	993[A]	PHE	2.6
1	B	6	LEU	2.6
1	B	469	ARG	2.6
1	B	472	LEU	2.5
1	B	460	ALA	2.4
1	A	396	ILE	2.4
1	A	44	GLN	2.3
1	B	78	ASN	2.3
1	A	459	ALA	2.3
1	A	391	PRO	2.3
1	A	186	GLY	2.3
1	A	461	ARG	2.3
1	A	184	LYS	2.2
1	A	34	TRP	2.2
1	A	443	ILE	2.2
1	A	26	PRO	2.2
1	B	102	ALA	2.2
1	A	399	VAL	2.2
1	A	437	PHE	2.2
1	B	8	THR	2.1
1	A	27	SER	2.1
1	A	444	GLU	2.1
1	B	470	LYS	2.1
1	A	188	PRO	2.1
1	A	45	ASN	2.1
1	A	183	GLY	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 monosaccharides in this entry.

6.4 Ligands [i](#)

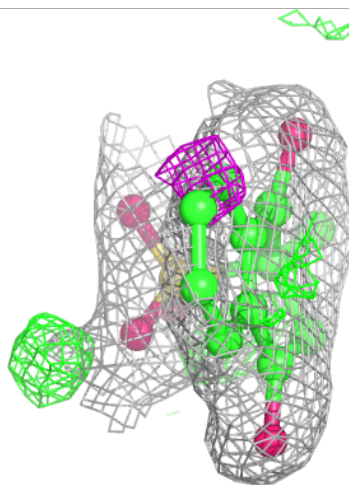
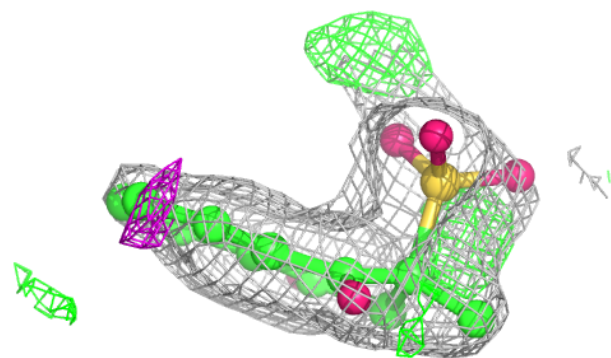
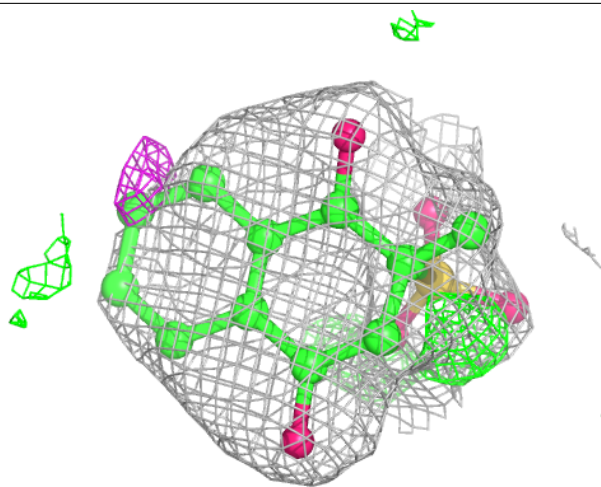
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	2L3	B	1102[A]	17/17	0.90	0.18	26,28,33,36	17
3	2L3	A	1102	17/17	0.91	0.16	29,34,44,46	0
5	2LB	B	1103[B]	17/17	0.91	0.17	23,27,33,37	17
4	EDO	A	1104	4/4	0.94	0.11	31,32,32,33	0
4	EDO	A	1107	4/4	0.95	0.23	19,35,43,47	0
4	EDO	B	1105	4/4	0.95	0.10	20,21,22,29	0
4	EDO	B	1106	4/4	0.95	0.19	18,28,29,30	0
4	EDO	A	1105	4/4	0.95	0.10	19,19,20,23	0
2	P5F	A	1101	56/61	0.96	0.11	8,15,27,30	0
2	P5F	B	1101	56/61	0.96	0.11	10,15,26,37	0
4	EDO	A	1103	4/4	0.97	0.09	15,21,27,27	0
4	EDO	B	1104	4/4	0.98	0.09	20,26,27,28	0
4	EDO	A	1106	4/4	0.98	0.09	14,21,25,29	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.

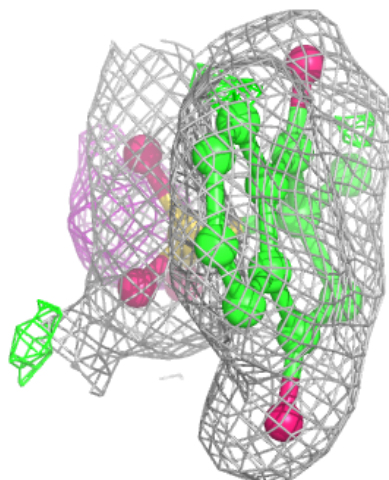
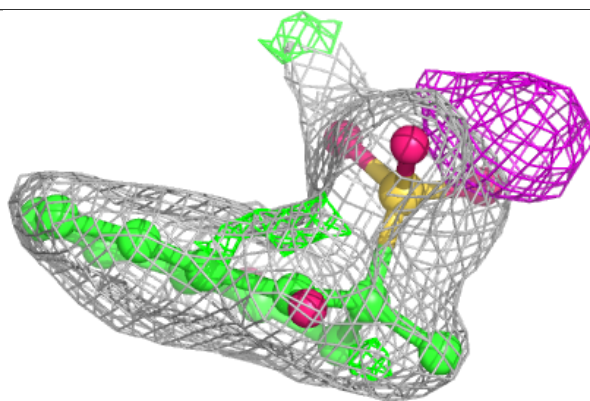
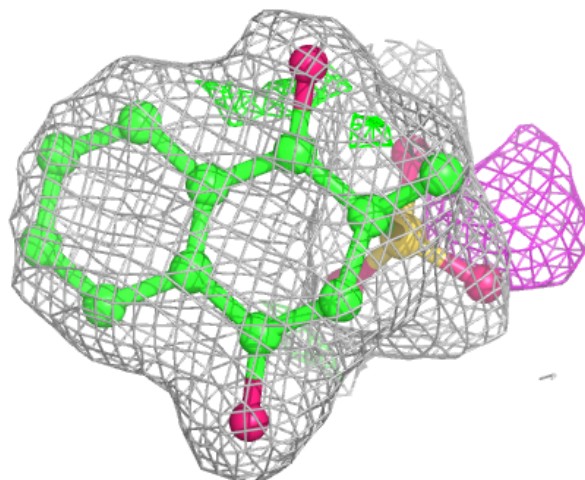
Electron density around 2L3 B 1102 (A):

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



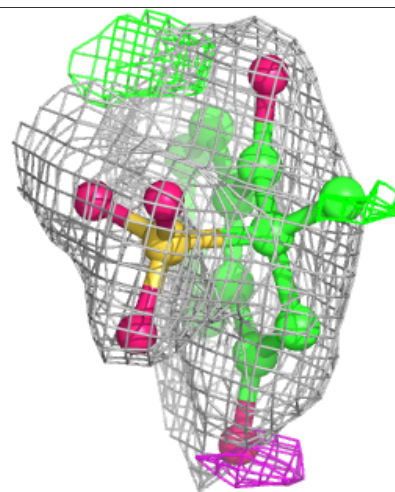
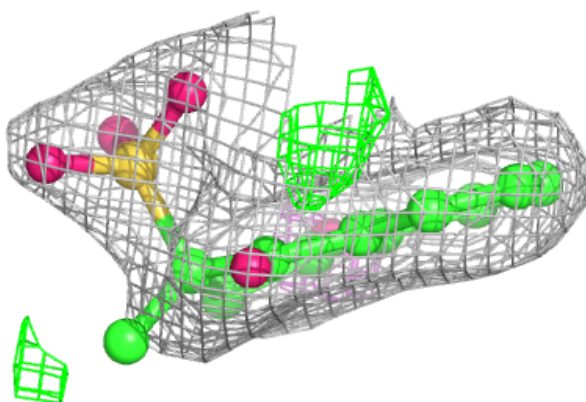
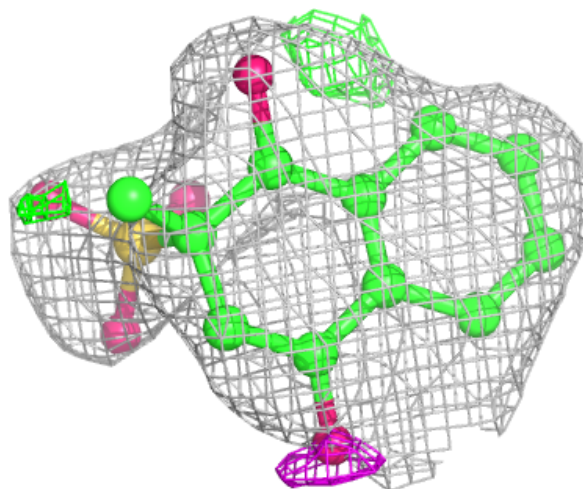
Electron density around 2L3 A 1102:

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



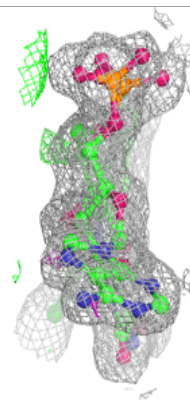
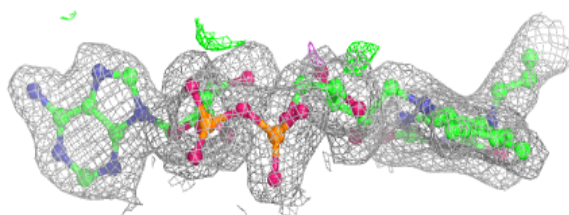
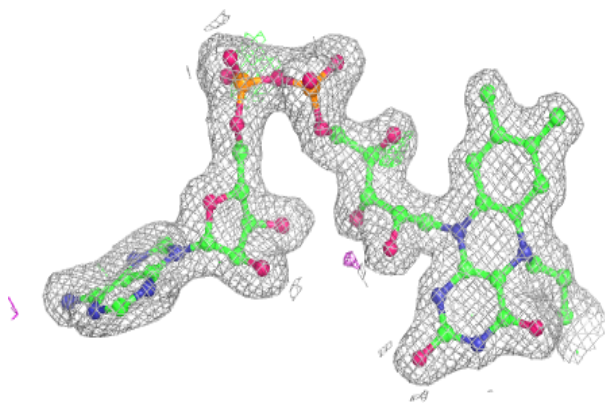
Electron density around 2LB B 1103 (B):

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

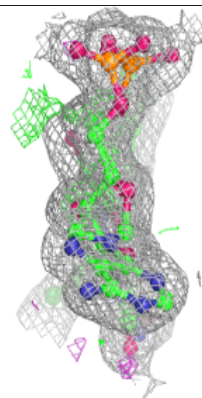
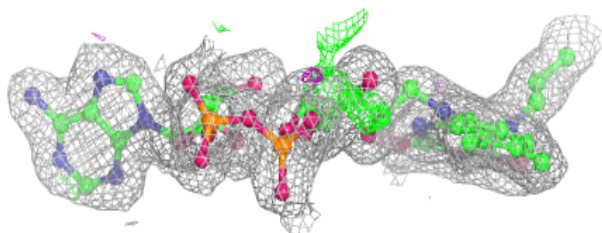
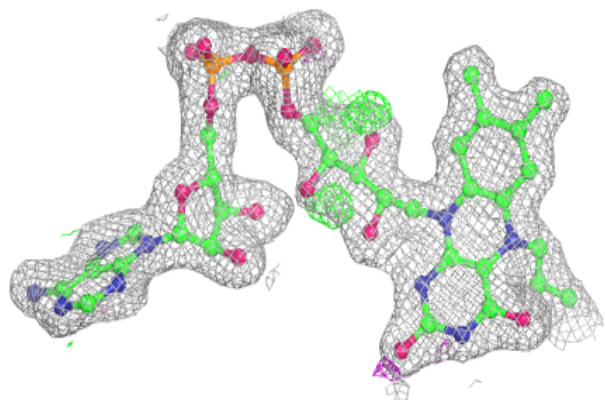


Electron density around P5F A 1101:

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

**Electron density around P5F 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.