



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:37 AM UTC

PDB ID : 7MP2 / pdb_00007mp2
Title : CRYSTAL STRUCTURE OF NATIVE BOVINE ARRESTIN 1 IN COM-
PLEX WITH 1D-MYO-INOSITOL 1,5-BISDIPHOSPHATE TETRAK-
ISPHOSPHATE (1,5-PP IP4)
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Deposited on : 2021-05-04
Resolution : 3.00 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	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

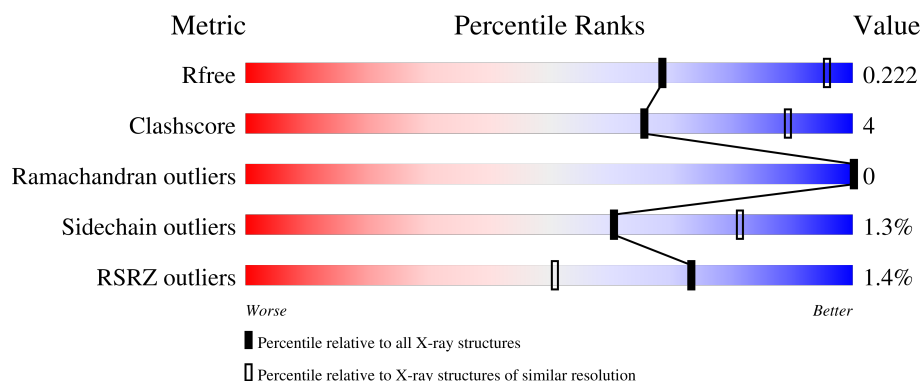
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	404	2% 81% 11% 9%
1	B	404	% 80% 10% 11%
1	C	404	% 79% 11% 10%
1	D	404	% 81% 10% 9%

2 Entry composition [i](#)

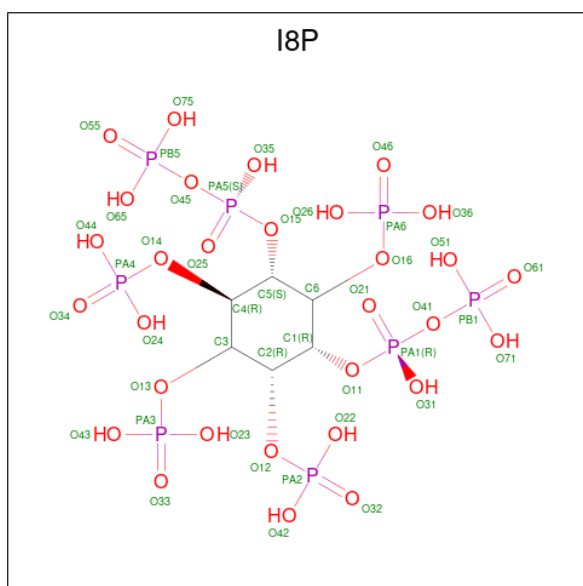
There are 3 unique types of molecules in this entry. The entry contains 11777 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 S-arrestin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	S	0	0	0
			2906	1859	489	549	9			
1	B	361	Total	C	N	O	S	0	0	0
			2850	1825	479	537	9			
1	C	364	Total	C	N	O	S	0	0	0
			2866	1834	484	540	8			
1	D	367	Total	C	N	O	S	0	0	0
			2882	1844	483	546	9			

- Molecule 2 is (1R,3S,4R,5S,6R)-2,4,5,6-tetrakis(phosphonooxy)cyclohexane-1,3-diyl bis[trihydrogen (diphosphate)] (CCD ID: I8P) (formula: $C_6H_{20}O_{30}P_8$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			44	6	30	8		

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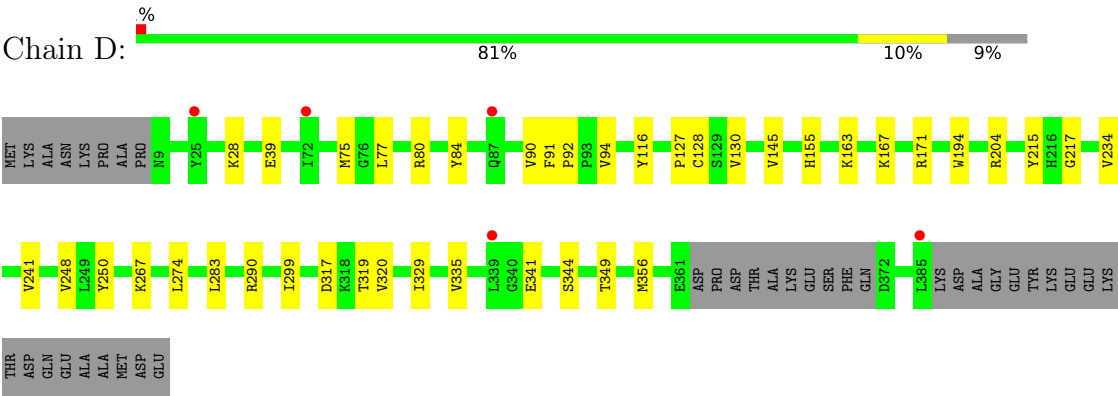
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	O	P	0	0
			44	6	30	8		
2	C	1	Total	C	O	P	0	0
			44	6	30	8		
2	C	1	Total	C	O	P	0	0
			44	6	30	8		
2	D	1	Total	C	O	P	0	0
			44	6	30	8		
2	D	1	Total	C	O	P	0	0
			44	6	30	8		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	O	0	0
			4	4		
3	C	2	Total	O	0	0
			2	2		
3	D	3	Total	O	0	0
			3	3		

● Molecule 1: S-arrestin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	168.90Å 187.09Å 90.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.16 – 3.00 49.16 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.4 (49.16-3.00) 97.4 (49.16-3.00)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.210 , 0.245 0.222 , 0.222	Depositor DCC
R_{free} test set	2844 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	90.4	Xtriage
Anisotropy	0.473	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 61.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11777	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% 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: I8P

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	0.99	0/2963	1.27	0/4017
1	B	1.00	0/2905	1.25	0/3938
1	C	0.99	0/2922	1.26	0/3960
1	D	1.01	0/2939	1.27	0/3988
All	All	1.00	0/11729	1.27	0/15903

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	2906	0	2963	24	0
1	B	2850	0	2897	25	0
1	C	2866	0	2923	25	0
1	D	2882	0	2924	22	0
2	A	44	0	6	0	0
2	B	44	0	6	0	0
2	C	88	0	12	1	0
2	D	88	0	12	2	0
3	A	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	2	0	0	0	0
3	D	3	0	0	0	0
All	All	11777	0	11743	90	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (90) 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:52:VAL:HG12	1:B:52:VAL:O	1.98	0.62
1:A:343:THR:H	1:A:346:GLU:HB2	1.65	0.62
1:D:90:VAL:HG22	1:D:116:TYR:HB2	1.82	0.61
1:B:71:ASP:HB3	1:B:74:VAL:HB	1.82	0.60
1:C:343:THR:H	1:C:346:GLU:HB2	1.67	0.59
1:B:80:ARG:CZ	1:C:342:LEU:HD21	2.34	0.58
1:D:128:CYS:O	1:D:130:VAL:HG23	2.06	0.56
1:C:188:PRO:HG2	1:C:209:LEU:HB2	1.87	0.56
1:B:128:CYS:O	1:B:130:VAL:HG23	2.06	0.56
1:A:128:CYS:O	1:A:130:VAL:HG23	2.06	0.55
1:D:248:VAL:HG22	1:D:319:THR:HG22	1.89	0.55
1:B:52:VAL:HG12	1:B:91:PHE:CE1	2.42	0.55
1:A:99:ALA:HB3	1:A:105:GLU:HG3	1.89	0.54
1:C:128:CYS:O	1:C:130:VAL:HG23	2.07	0.54
1:A:188:PRO:HG2	1:A:209:LEU:HB2	1.89	0.54
1:B:248:VAL:HG22	1:B:319:THR:HG22	1.89	0.53
1:C:264:ALA:HB2	1:C:278:LEU:HD21	1.91	0.53
1:C:248:VAL:HG22	1:C:319:THR:HG22	1.90	0.53
1:A:264:ALA:HB2	1:A:278:LEU:HD21	1.91	0.52
1:D:335:VAL:O	1:D:344:SER:HB2	2.09	0.52
1:C:127:PRO:HG2	1:C:145:VAL:HG11	1.91	0.52
1:B:127:PRO:HG2	1:B:145:VAL:HG11	1.91	0.52
1:B:52:VAL:O	1:B:52:VAL:CG1	2.57	0.51
1:A:127:PRO:HG2	1:A:145:VAL:HG11	1.92	0.51
1:D:127:PRO:HG2	1:D:145:VAL:HG11	1.92	0.51
1:D:194:TRP:CD2	1:D:349:THR:HG21	2.45	0.51
1:C:190:ALA:HB2	1:C:351:VAL:HG22	1.93	0.50
1:B:335:VAL:O	1:B:344:SER:HB2	2.12	0.50
1:B:100:THR:HG22	1:B:101:THR:N	2.27	0.50
1:D:90:VAL:HG22	1:D:116:TYR:CB	2.42	0.49
1:A:217:GLY:HA2	1:A:283:LEU:HD21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:THR:CG2	1:B:101:THR:N	2.76	0.49
1:B:217:GLY:HA2	1:B:283:LEU:HD21	1.94	0.49
1:C:184:MET:HE2	1:C:213:ILE:HB	1.95	0.48
1:C:217:GLY:HA2	1:C:283:LEU:HD21	1.95	0.48
1:B:194:TRP:CD2	1:B:349:THR:HG21	2.49	0.48
1:A:189:ARG:HE	1:D:75:MET:HE2	1.78	0.48
1:D:217:GLY:HA2	1:D:283:LEU:HD21	1.94	0.48
1:B:157:THR:O	1:B:161:GLU:HB2	2.14	0.48
1:A:184:MET:HE2	1:A:213:ILE:HB	1.96	0.47
1:B:194:TRP:CZ2	1:B:349:THR:HB	2.49	0.47
1:B:47:VAL:HB	1:B:52:VAL:HG21	1.98	0.46
1:C:190:ALA:CB	1:C:351:VAL:HG22	2.46	0.46
1:A:190:ALA:HB2	1:A:351:VAL:HG22	1.98	0.46
1:C:300:LYS:NZ	2:C:502:I8P:O34	2.49	0.45
1:A:91:PHE:HA	1:A:92:PRO:C	2.42	0.45
1:B:164:ILE:HD12	1:C:343:THR:CG2	2.47	0.45
1:B:84:TYR:HA	1:C:197:PHE:CE1	2.52	0.45
1:D:77:LEU:HD21	1:D:250:TYR:OH	2.17	0.44
1:C:91:PHE:HA	1:C:92:PRO:C	2.42	0.44
1:A:190:ALA:CB	1:A:351:VAL:HG22	2.47	0.44
1:A:342:LEU:HD21	1:D:80:ARG:CZ	2.48	0.44
1:A:215:TYR:HA	1:A:356:MET:O	2.18	0.44
1:B:85:PHE:HZ	1:B:87:GLN:HE21	1.64	0.44
1:C:215:TYR:HA	1:C:356:MET:O	2.18	0.44
1:C:335:VAL:HB	1:C:338:LEU:HD12	2.00	0.44
1:A:290:ARG:HD3	1:A:299:ILE:HG23	2.00	0.43
1:B:91:PHE:HA	1:B:92:PRO:C	2.43	0.43
1:B:215:TYR:HA	1:B:356:MET:O	2.18	0.43
1:D:91:PHE:HA	1:D:92:PRO:C	2.42	0.43
1:D:194:TRP:CZ2	1:D:349:THR:HB	2.53	0.43
1:C:317:ASP:O	1:C:320:VAL:HG22	2.19	0.43
1:A:317:ASP:O	1:A:320:VAL:HG22	2.19	0.42
1:A:240:LEU:HD12	1:A:240:LEU:N	2.33	0.42
1:C:361:GLU:N	1:C:361:GLU:OE2	2.53	0.42
1:B:317:ASP:O	1:B:320:VAL:HG22	2.19	0.42
1:D:163:LYS:HE2	2:D:502:I8P:O33	2.20	0.42
1:D:215:TYR:HA	1:D:356:MET:O	2.19	0.42
1:D:171:ARG:HD3	2:D:501:I8P:PB1	2.60	0.42
1:A:75:MET:HE2	1:A:75:MET:HB2	1.91	0.42
1:D:317:ASP:O	1:D:320:VAL:HG22	2.20	0.41
1:C:156:SER:HB2	1:C:162:ASP:OD1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:THR:HG22	1:A:164:ILE:HG12	2.02	0.41
1:C:154:THR:HG22	1:C:164:ILE:HG12	2.03	0.41
1:B:124:ASP:OD1	1:B:124:ASP:N	2.47	0.41
1:C:90:VAL:CG1	1:C:118:PHE:HB3	2.51	0.41
1:D:234:VAL:O	1:D:267:LYS:HA	2.21	0.41
1:B:100:THR:HG22	1:B:101:THR:O	2.20	0.41
1:B:241:VAL:HG22	1:B:329:ILE:CD1	2.51	0.41
1:C:110:LYS:HG3	1:C:375:PHE:HD2	1.86	0.41
1:C:180:ALA:HA	1:C:356:MET:SD	2.61	0.41
1:D:290:ARG:HD3	1:D:299:ILE:HG23	2.03	0.41
1:A:180:ALA:HA	1:A:356:MET:SD	2.61	0.40
1:A:331:VAL:O	1:A:348:ALA:HA	2.21	0.40
1:D:28:LYS:HE3	1:D:39:GLU:OE1	2.20	0.40
1:A:335:VAL:HB	1:A:338:LEU:HD12	2.02	0.40
1:D:241:VAL:HG22	1:D:329:ILE:CD1	2.50	0.40
1:C:331:VAL:O	1:C:348:ALA:HA	2.22	0.40
1:A:131:MET:HE1	1:A:298:LYS:HG3	2.03	0.40
1:A:197:PHE:CE1	1:D:84:TYR:HA	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	365/404 (90%)	345 (94%)	20 (6%)	0	100	100
1	B	353/404 (87%)	337 (96%)	16 (4%)	0	100	100
1	C	358/404 (89%)	338 (94%)	20 (6%)	0	100	100
1	D	363/404 (90%)	346 (95%)	17 (5%)	0	100	100
All	All	1439/1616 (89%)	1366 (95%)	73 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	329/359 (92%)	326 (99%)	3 (1%)	70	85
1	B	323/359 (90%)	318 (98%)	5 (2%)	57	80
1	C	324/359 (90%)	321 (99%)	3 (1%)	70	85
1	D	325/359 (90%)	319 (98%)	6 (2%)	51	77
All	All	1301/1436 (91%)	1284 (99%)	17 (1%)	61	81

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	274	LEU
1	A	341	GLU
1	A	349	THR
1	B	155	HIS
1	B	160	GLU
1	B	204	ARG
1	B	274	LEU
1	B	385	LEU
1	C	274	LEU
1	C	341	GLU
1	C	349	THR
1	D	94	VAL
1	D	155	HIS
1	D	167	LYS
1	D	204	ARG
1	D	274	LEU
1	D	341	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	357	HIS
1	A	374	ASN

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Mol	Chain	Res	Type
1	B	87	GLN
1	B	357	HIS
1	B	374	ASN
1	C	114	ASN
1	C	357	HIS
1	C	374	ASN
1	D	69	GLN
1	D	114	ASN
1	D	357	HIS

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](#)

6 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	$\# Z > 2$
2	I8P	D	501	-	42,44,44	0.74	2 (4%)	68,74,74	0.82	3 (4%)
2	I8P	A	501	-	42,44,44	0.76	1 (2%)	68,74,74	0.75	3 (4%)
2	I8P	B	501	-	42,44,44	0.74	1 (2%)	68,74,74	0.68	3 (4%)
2	I8P	C	502	-	42,44,44	0.74	2 (4%)	68,74,74	0.82	3 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	I8P	D	502	-	42,44,44	0.68	1 (2%)	68,74,74	0.66	3 (4%)
2	I8P	C	501	-	42,44,44	0.74	1 (2%)	68,74,74	0.77	2 (2%)

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
2	I8P	D	501	-	-	10/42/66/66	0/1/1/1
2	I8P	A	501	-	-	9/42/66/66	0/1/1/1
2	I8P	B	501	-	-	9/42/66/66	0/1/1/1
2	I8P	C	502	-	-	16/42/66/66	0/1/1/1
2	I8P	D	502	-	-	11/42/66/66	0/1/1/1
2	I8P	C	501	-	-	16/42/66/66	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	I8P	PA3-O13	3.55	1.65	1.59
2	C	501	I8P	PA3-O13	3.48	1.65	1.59
2	D	501	I8P	PA3-O13	3.40	1.65	1.59
2	B	501	I8P	PA3-O13	3.37	1.65	1.59
2	C	502	I8P	PA3-O13	3.22	1.65	1.59
2	D	502	I8P	PA3-O13	2.86	1.64	1.59
2	C	502	I8P	PA5-O15	2.16	1.65	1.59
2	D	501	I8P	PA5-O15	2.02	1.65	1.59

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	I8P	PA4-O14-C4	3.69	133.28	123.43
2	C	502	I8P	PA6-O16-C6	3.47	132.69	123.43
2	C	502	I8P	PA4-O14-C4	3.31	132.28	123.43
2	A	501	I8P	PA4-O14-C4	3.26	132.13	123.43
2	B	501	I8P	PA6-O16-C6	3.24	132.09	123.43
2	D	501	I8P	PA6-O16-C6	3.22	132.04	123.43
2	C	501	I8P	PA2-O12-C2	3.15	131.85	123.43
2	A	501	I8P	PA6-O16-C6	3.10	131.71	123.43
2	D	502	I8P	PA4-O14-C4	3.03	131.53	123.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	502	I8P	PA6-O16-C6	2.97	131.37	123.43
2	A	501	I8P	PA2-O12-C2	2.92	131.24	123.43
2	C	502	I8P	PA2-O12-C2	2.85	131.04	123.43
2	B	501	I8P	PA4-O14-C4	2.79	130.88	123.43
2	C	501	I8P	PA6-O16-C6	2.63	130.45	123.43
2	B	501	I8P	PA2-O12-C2	2.35	129.72	123.43
2	D	502	I8P	PA2-O12-C2	2.18	129.24	123.43
2	D	501	I8P	PA2-O12-C2	2.07	128.95	123.43

There are no chirality outliers.

All (71) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	I8P	C6-C5-O15-PA5
2	A	501	I8P	C5-O15-PA5-O45
2	A	501	I8P	PA1-O41-PB1-O71
2	B	501	I8P	C4-C5-O15-PA5
2	B	501	I8P	C6-C5-O15-PA5
2	B	501	I8P	C5-C6-O16-PA6
2	B	501	I8P	C5-O15-PA5-O45
2	B	501	I8P	PA1-O41-PB1-O51
2	C	501	I8P	C1-C2-O12-PA2
2	C	501	I8P	C3-C2-O12-PA2
2	C	501	I8P	C5-O15-PA5-O45
2	C	502	I8P	C4-C5-O15-PA5
2	C	502	I8P	C6-C5-O15-PA5
2	C	502	I8P	C1-O11-PA1-O41
2	D	501	I8P	C2-C3-O13-PA3
2	D	501	I8P	C4-C3-O13-PA3
2	D	501	I8P	C6-C5-O15-PA5
2	D	501	I8P	PB1-O41-PA1-O11
2	D	502	I8P	C1-O11-PA1-O21
2	D	502	I8P	C1-O11-PA1-O41
2	D	502	I8P	C5-O15-PA5-O45
2	D	502	I8P	PA1-O41-PB1-O71
2	C	502	I8P	C5-C4-O14-PA4
2	A	501	I8P	C6-O16-PA6-O46
2	C	501	I8P	PA1-O41-PB1-O61
2	C	502	I8P	PB1-O41-PA1-O21
2	C	502	I8P	C3-C4-O14-PA4
2	D	501	I8P	C4-C5-O15-PA5
2	C	502	I8P	PB1-O41-PA1-O11

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Mol	Chain	Res	Type	Atoms
2	C	501	I8P	PA5-O45-PB5-O55
2	D	502	I8P	PA5-O45-PB5-O55
2	C	501	I8P	PA1-O41-PB1-O71
2	D	501	I8P	PA5-O45-PB5-O65
2	B	501	I8P	C5-O15-PA5-O25
2	C	501	I8P	C5-O15-PA5-O25
2	A	501	I8P	C3-O13-PA3-O33
2	C	501	I8P	C3-O13-PA3-O33
2	B	501	I8P	PB1-O41-PA1-O31
2	C	502	I8P	C2-C3-O13-PA3
2	A	501	I8P	C2-O12-PA2-O22
2	C	502	I8P	C4-O14-PA4-O24
2	B	501	I8P	C1-C6-O16-PA6
2	C	502	I8P	C4-C3-O13-PA3
2	C	501	I8P	C2-O12-PA2-O32
2	C	501	I8P	C6-O16-PA6-O46
2	C	502	I8P	C2-O12-PA2-O32
2	D	502	I8P	C2-O12-PA2-O32
2	D	502	I8P	PB5-O45-PA5-O25
2	C	501	I8P	PA5-O45-PB5-O65
2	C	501	I8P	PA5-O45-PB5-O75
2	C	502	I8P	PA1-O41-PB1-O51
2	C	502	I8P	PA5-O45-PB5-O65
2	D	502	I8P	PA5-O45-PB5-O65
2	D	502	I8P	PA5-O45-PB5-O75
2	B	501	I8P	C5-O15-PA5-O35
2	C	501	I8P	C5-O15-PA5-O35
2	C	502	I8P	C1-O11-PA1-O21
2	C	502	I8P	C1-O11-PA1-O31
2	D	501	I8P	C1-O11-PA1-O21
2	D	502	I8P	C1-O11-PA1-O31
2	A	501	I8P	C3-O13-PA3-O23
2	A	501	I8P	C3-O13-PA3-O43
2	C	501	I8P	C2-O12-PA2-O22
2	C	501	I8P	C3-O13-PA3-O43
2	C	501	I8P	C6-O16-PA6-O26
2	C	502	I8P	C2-O12-PA2-O42
2	D	501	I8P	C2-O12-PA2-O22
2	D	501	I8P	C4-O14-PA4-O44
2	D	501	I8P	C2-O12-PA2-O32
2	D	502	I8P	C1-C6-O16-PA6
2	A	501	I8P	PB1-O41-PA1-O31

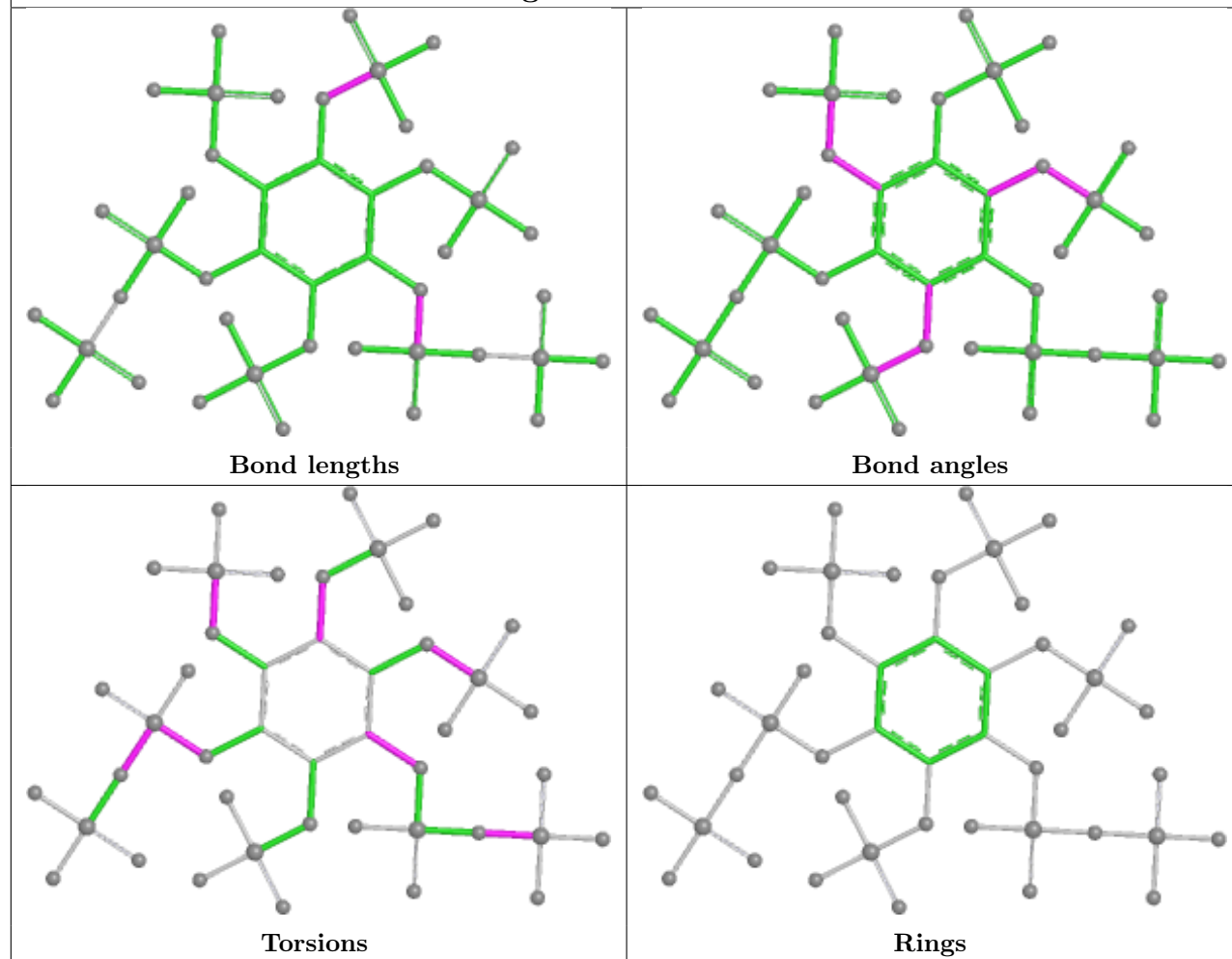
There are no ring outliers.

3 monomers are involved in 3 short contacts:

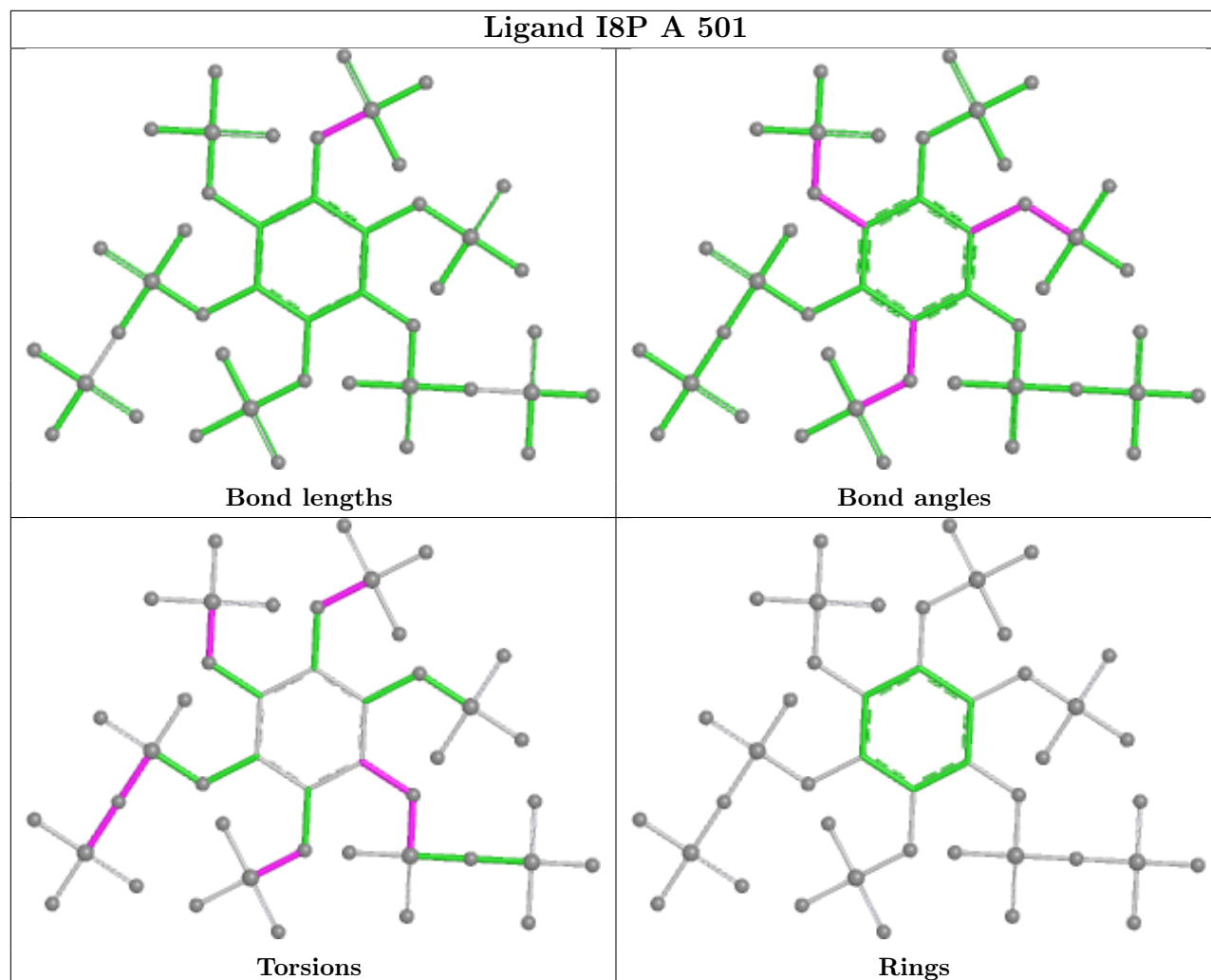
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	I8P	1	0
2	C	502	I8P	1	0
2	D	502	I8P	1	0

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.

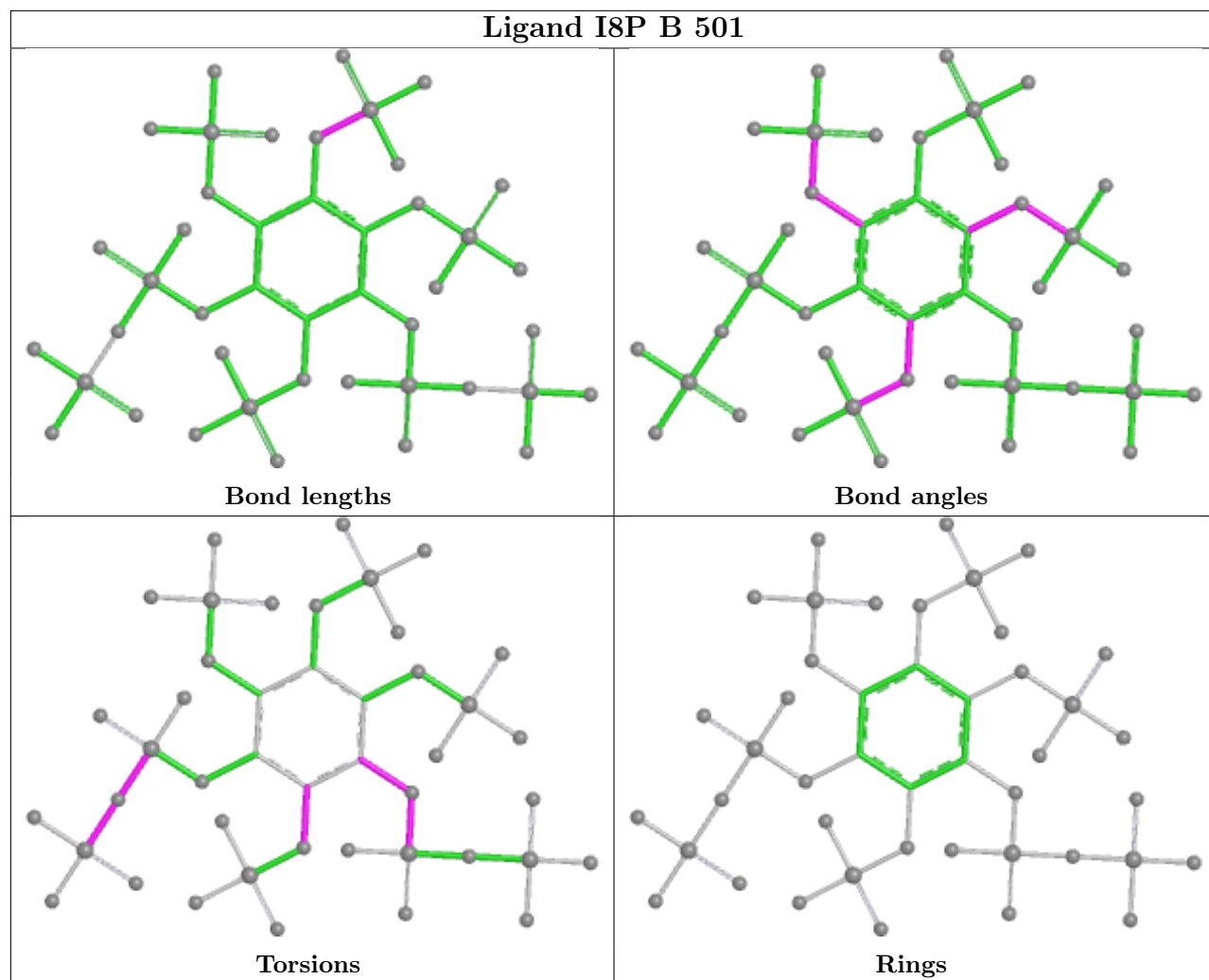
Ligand I8P D 501



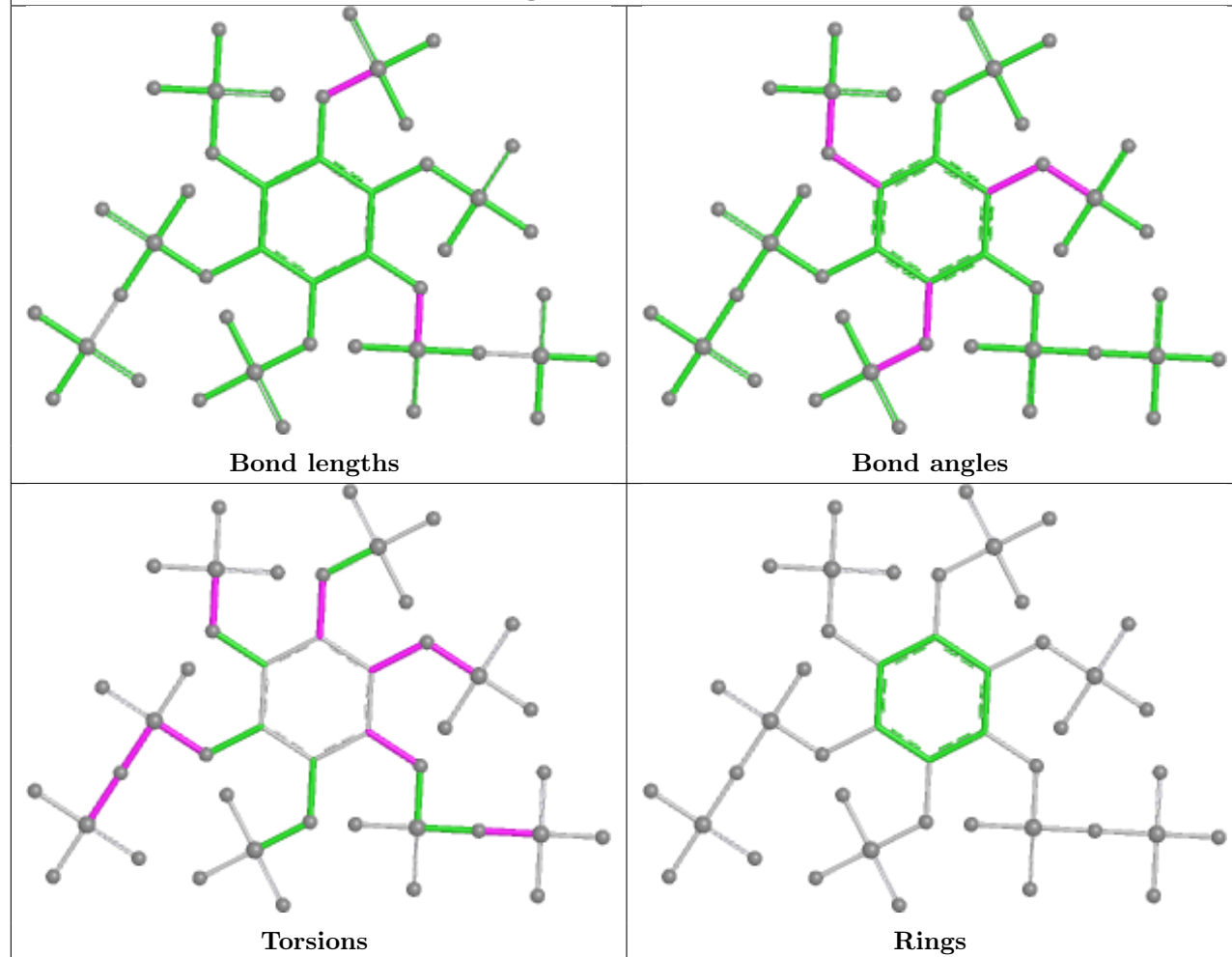
Ligand I8P A 501



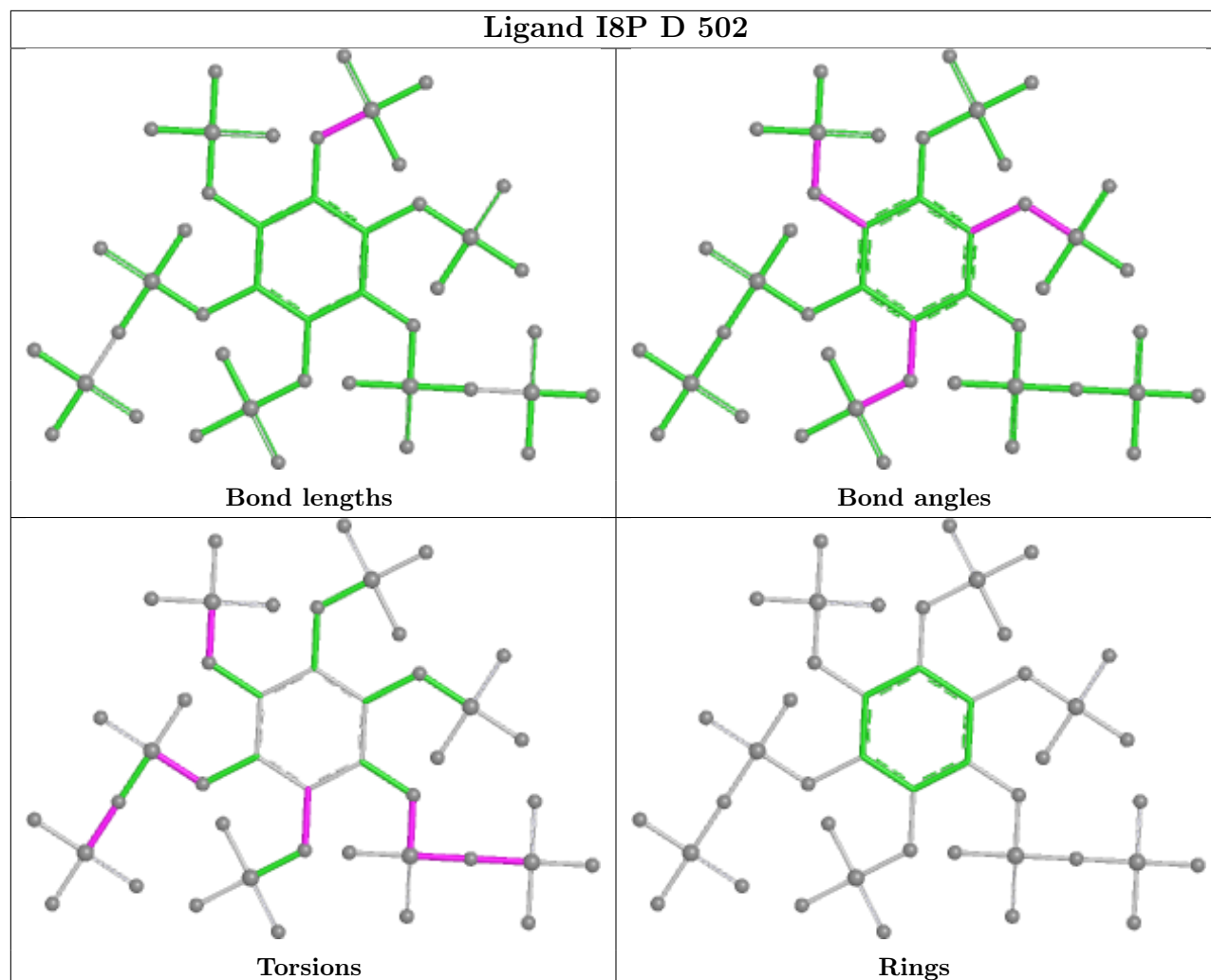
Ligand I8P B 501

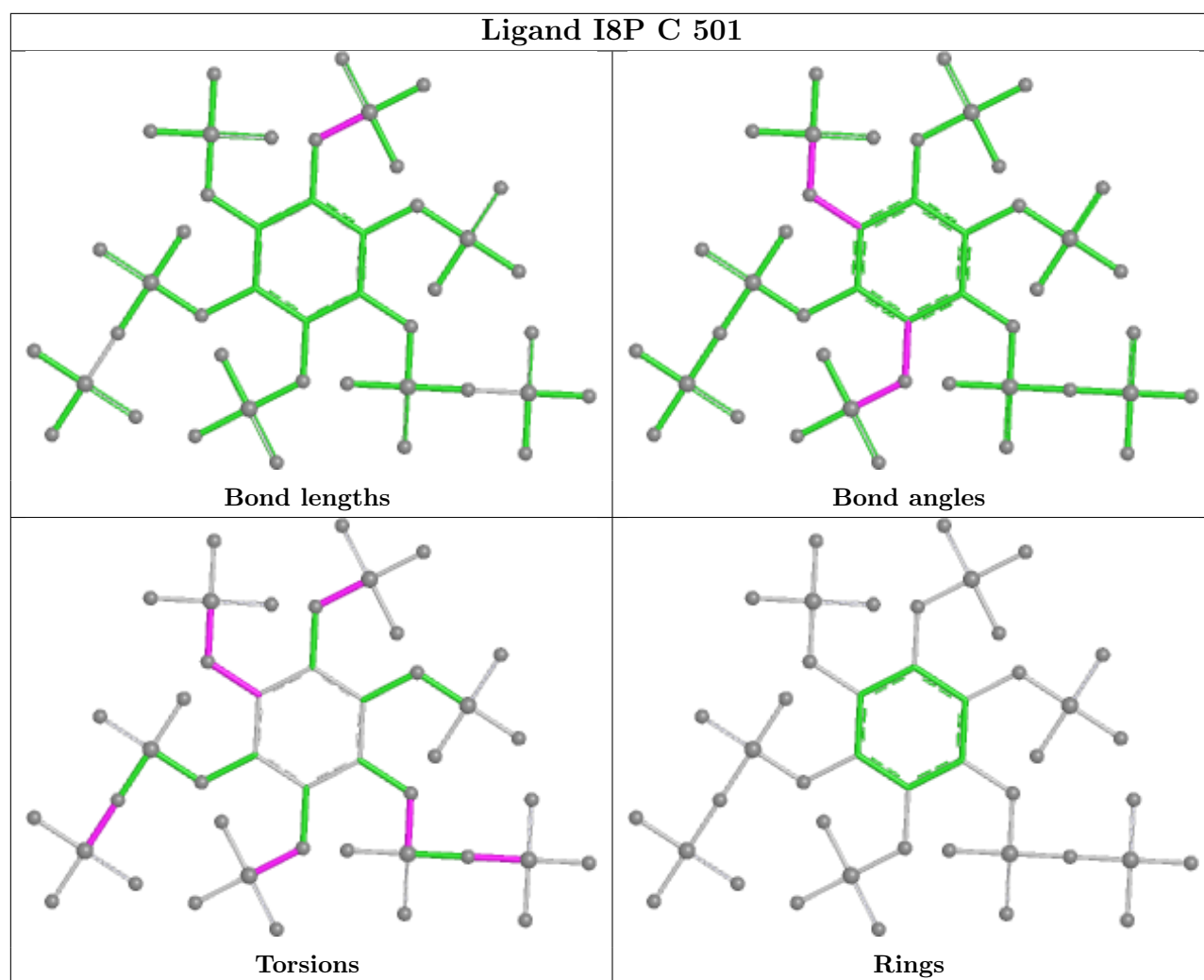


Ligand I8P C 502



Ligand I8P D 502





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	369/404 (91%)	-0.14	7 (1%)	66 43	72, 94, 148, 177	0
1	B	361/404 (89%)	0.07	5 (1%)	73 51	74, 112, 182, 211	0
1	C	364/404 (90%)	-0.20	4 (1%)	78 57	73, 95, 141, 177	0
1	D	367/404 (90%)	-0.02	5 (1%)	73 51	69, 108, 169, 203	0
All	All	1461/1616 (90%)	-0.07	21 (1%)	73 51	69, 101, 167, 211	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	342	LEU	3.4
1	C	345	SER	3.1
1	B	159	VAL	3.0
1	A	139	VAL	2.9
1	C	344	SER	2.6
1	A	133	GLN	2.6
1	B	338	LEU	2.5
1	A	339	LEU	2.4
1	A	344	SER	2.4
1	D	72	ILE	2.3
1	B	315	GLY	2.3
1	D	87	GLN	2.3
1	A	345	SER	2.2
1	B	386	LYS	2.2
1	D	25	TYR	2.2
1	A	72	ILE	2.1
1	D	385	LEU	2.1
1	D	339	LEU	2.1
1	B	380	PHE	2.1
1	C	139	VAL	2.0
1	C	342	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

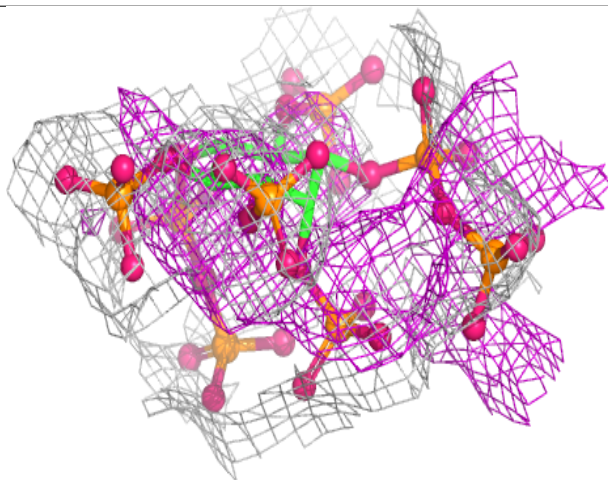
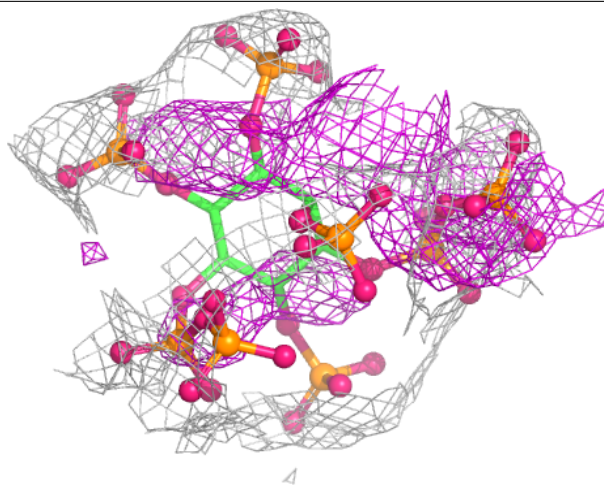
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	I8P	A	501	44/44	0.59	0.10	187,237,270,288	0
2	I8P	B	501	44/44	0.65	0.09	167,218,239,252	0
2	I8P	D	501	44/44	0.70	0.08	152,200,228,232	0
2	I8P	C	501	44/44	0.76	0.09	171,206,239,248	0
2	I8P	C	502	44/44	0.77	0.08	165,216,243,252	0
2	I8P	D	502	44/44	0.78	0.08	172,214,238,249	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.

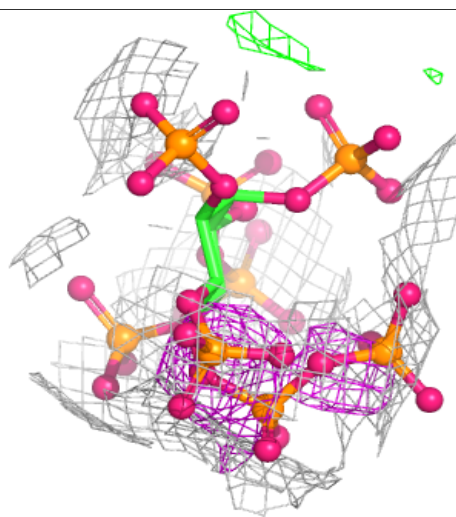
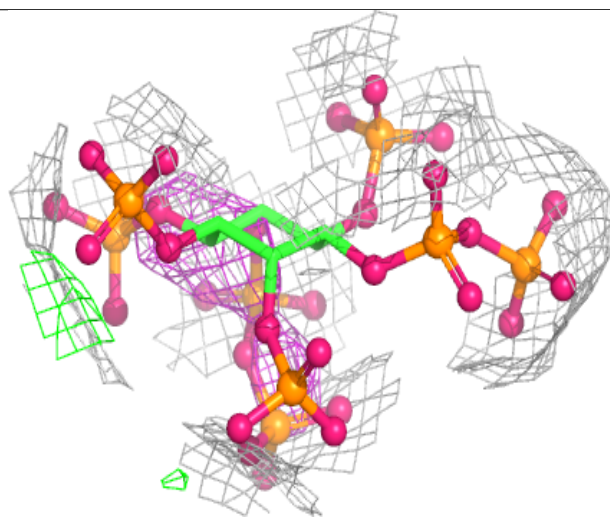
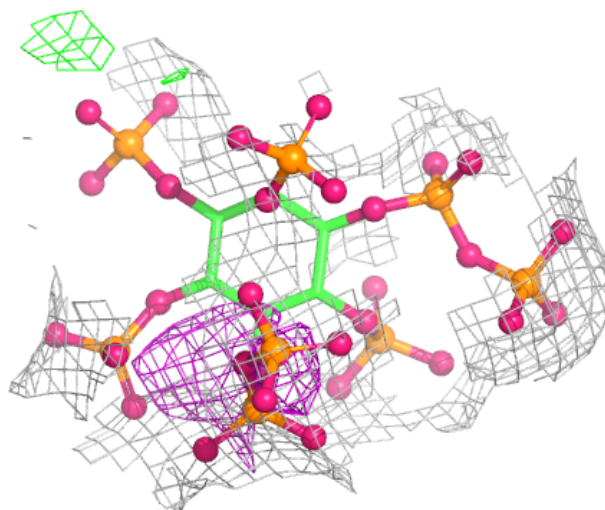
Electron density around I8P A 501:

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



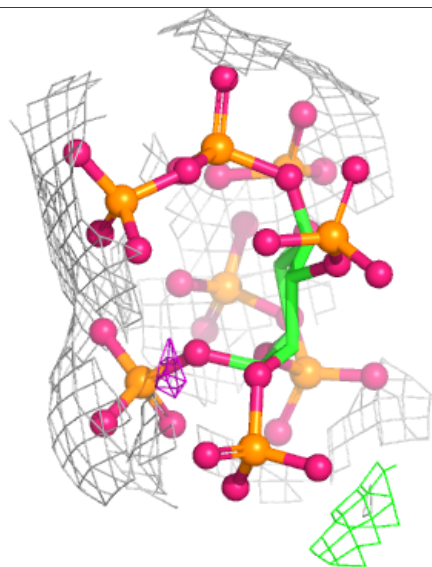
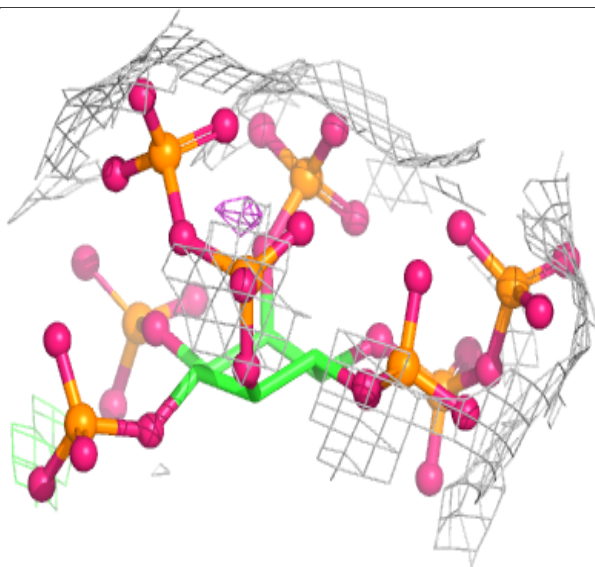
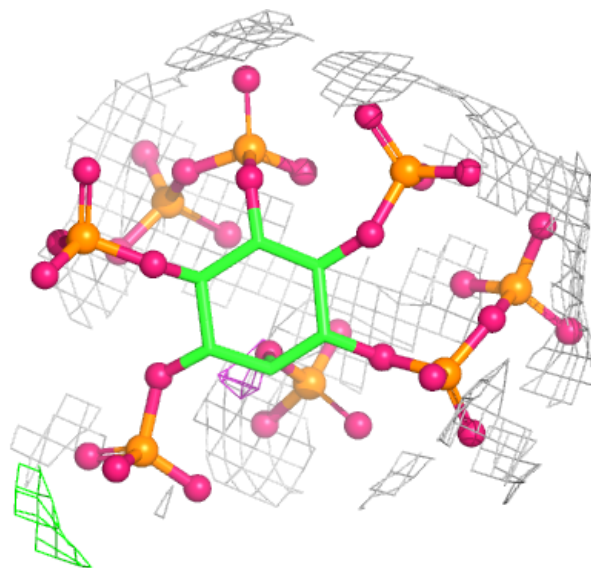
Electron density around I8P B 501:

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



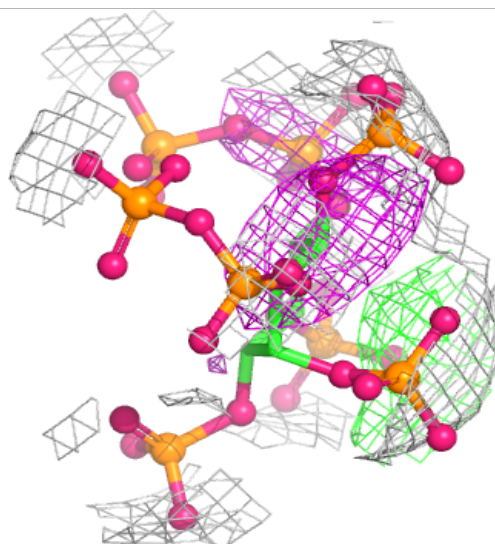
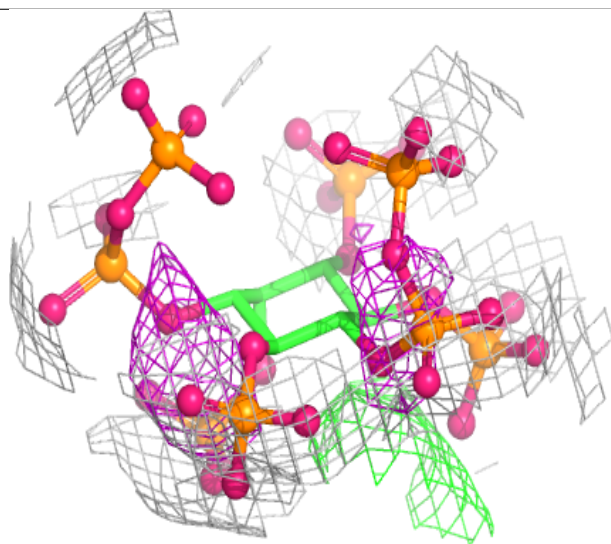
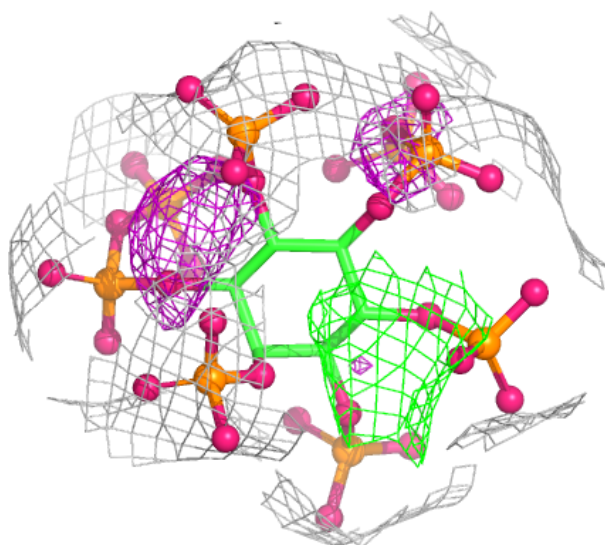
Electron density around I8P D 501:

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



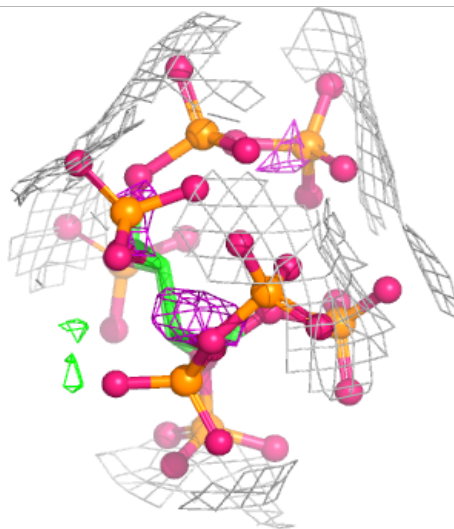
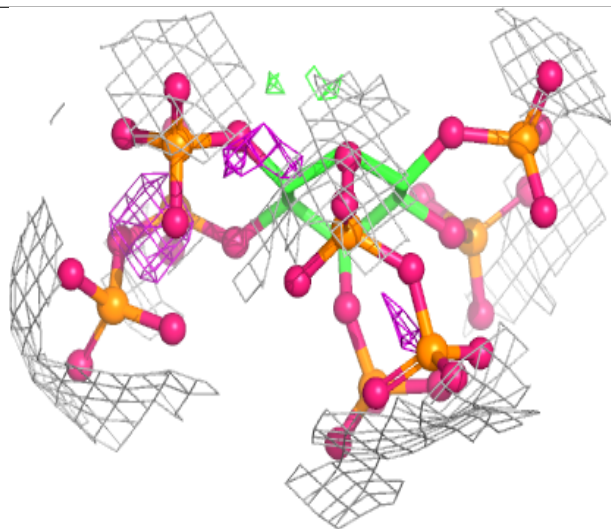
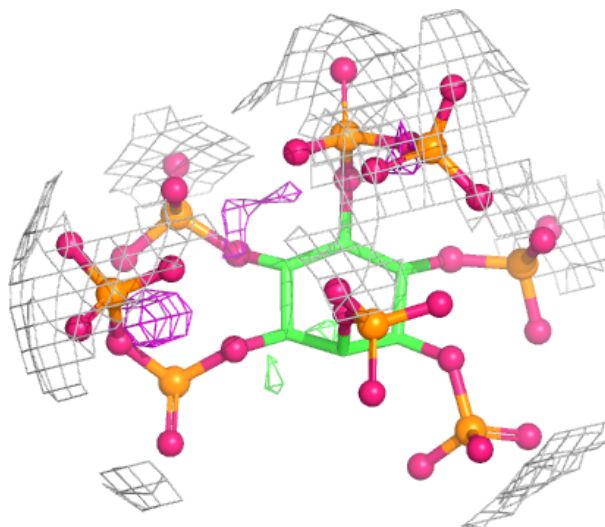
Electron density around I8P C 501:

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



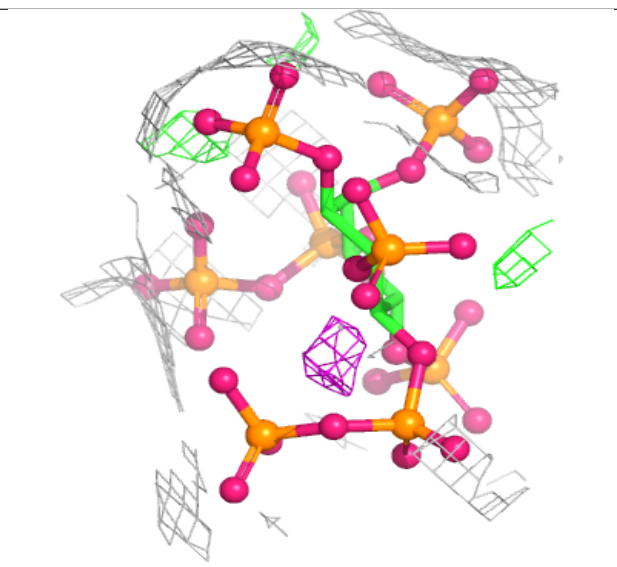
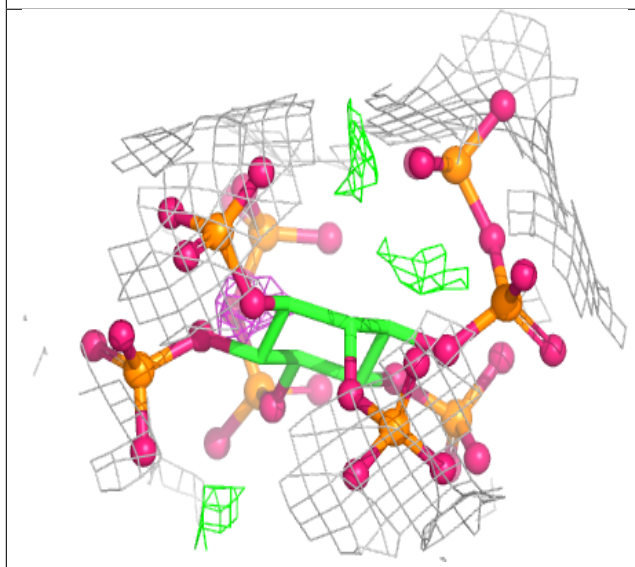
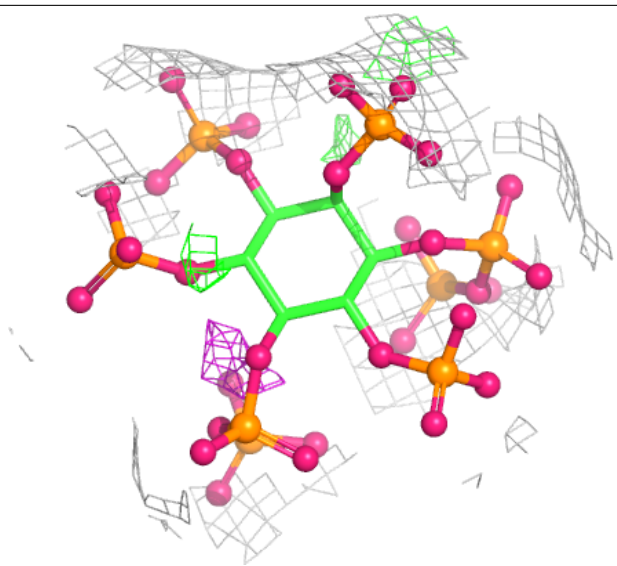
Electron density around I8P C 502:

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



Electron density around I8P D 502:

$2mF_o-DF_c$ (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.