



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 07:06 AM BST

PDB ID : 5YFZ
Title : Plasmodium vivax SHMT bound with PLP-glycine and S-GS626
Authors : Chitnumsub, P.; Jaruwat, A.; Leartsakulpanich, U.; Schwertz, G.; Diederich, F.
Deposited on : 2017-09-22
Resolution : 2.16 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 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)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
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) : 2.13.1

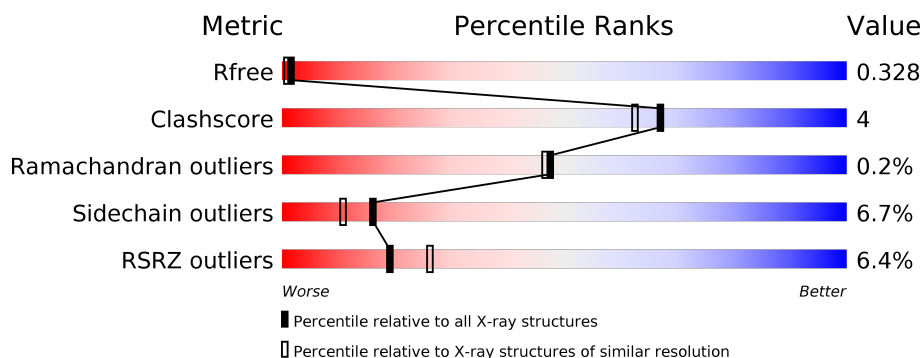
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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 on 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	442	<div> <div>8%</div> <div> <div></div> <div>85%</div> <div>14%</div> </div> <div>.</div> </div>
1	B	442	<div> <div>6%</div> <div> <div></div> <div>85%</div> <div>14%</div> </div> <div>.</div> </div>
1	C	442	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>14%</div> </div> <div>.</div> </div>

2 Entry composition [i](#)

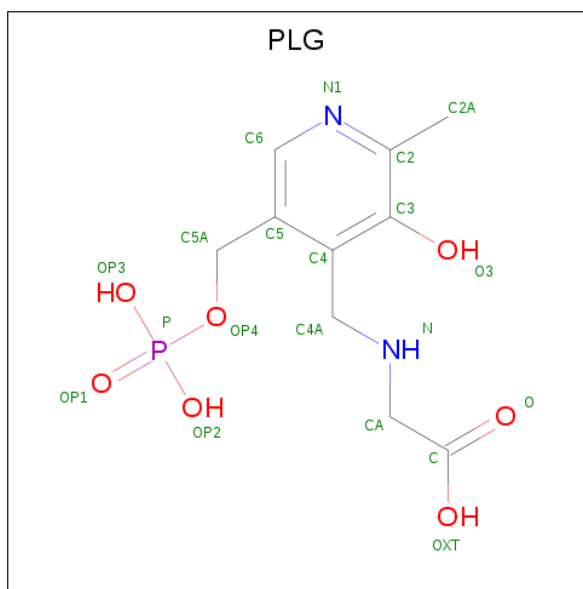
There are 5 unique types of molecules in this entry. The entry contains 10924 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Serine hydroxymethyltransferase.

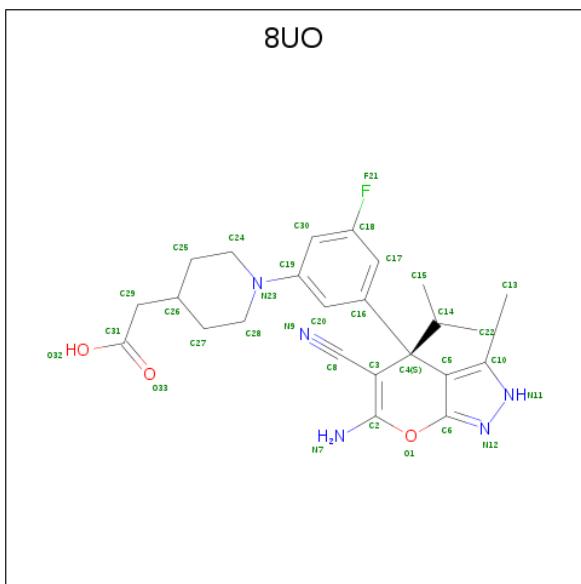
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	0	0	0
			3458	2186	600	655	17			
1	B	442	Total	C	N	O	S	0	0	0
			3458	2186	600	655	17			
1	C	442	Total	C	N	O	S	0	0	0
			3458	2186	600	655	17			

- Molecule 2 is N-GLYCINE-[3-HYDROXY-2-METHYL-5-PHOSPHONOXYMETHYL-PYRIDIN-4-YL-METHANE] (three-letter code: PLG) (formula: C₁₀H₁₅N₂O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			20	10	2	7	1		
2	B	1	Total	C	N	O	P	0	0
			20	10	2	7	1		
2	C	1	Total	C	N	O	P	0	0
			20	10	2	7	1		

- Molecule 3 is 2-[1-[3-[(4 {S})-6-azanyl-5-cyano-3-methyl-4-propan-2-yl-2 {H}-pyrano[2,3-c]pyrazol-4-yl]-5-fluoranyl-phenyl]piperidin-4-yl]ethanoic acid (three-letter code: 8UO) (formula: C₂₄H₂₈FN₅O₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			33	24	1	5	3		
3	B	1	Total	C	F	N	O	0	0
			33	24	1	5	3		
3	C	1	Total	C	F	N	O	0	0
			33	24	1	5	3		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	C	1	Total	Cl	0	0
			1	1		

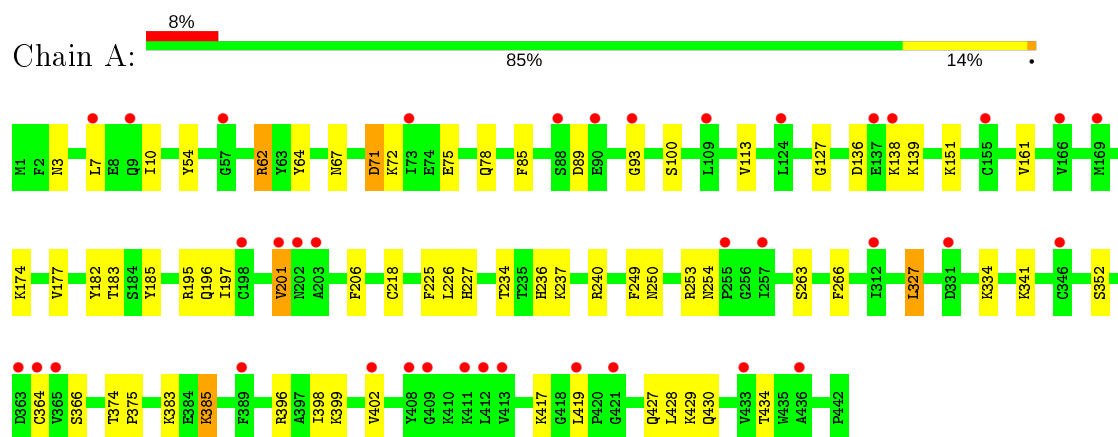
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	124	Total	O	0	0
			124	124		
5	B	131	Total	O	0	0
			131	131		
5	C	134	Total	O	0	0
			134	134		

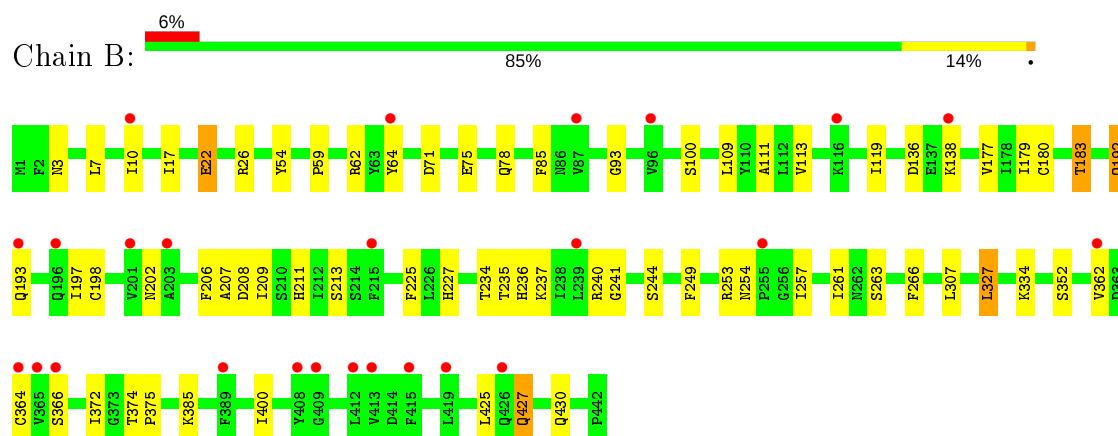
3 Residue-property plots

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.

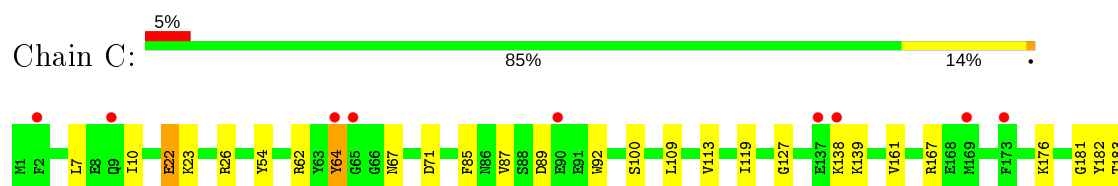
• Molecule 1: Serine hydroxymethyltransferase

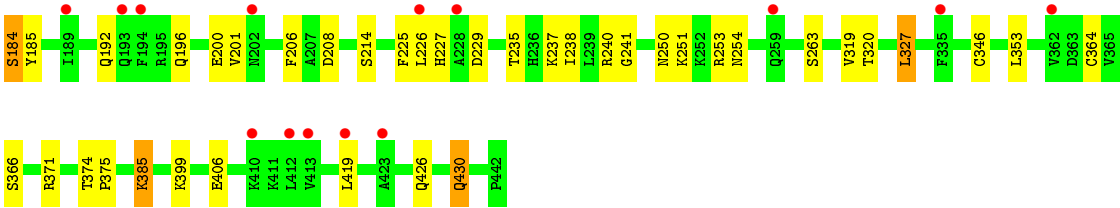


• Molecule 1: Serine hydroxymethyltransferase



• Molecule 1: Serine hydroxymethyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.88Å 58.80Å 235.73Å 90.00° 89.99° 90.00°	Depositor
Resolution (Å)	30.00 – 2.16 29.47 – 2.16	Depositor EDS
% Data completeness (in resolution range)	91.4 (30.00-2.16) 91.4 (29.47-2.16)	Depositor EDS
R_{merge}	0.02	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.98 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.263 , 0.333 0.264 , 0.328	Depositor DCC
R_{free} test set	6717 reflections (9.76%)	wwPDB-VP
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 14.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.024 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.025 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.470 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.469 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.023 for -h,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	10924	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% 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: PLG, 8UO, CL

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	0/3521	0.72	2/4754 (0.0%)
1	B	0.54	0/3521	0.71	2/4754 (0.0%)
1	C	0.56	0/3521	0.72	1/4754 (0.0%)
All	All	0.54	0/10563	0.72	5/14262 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	240	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	A	327	LEU	CA-CB-CG	5.39	127.70	115.30
1	A	240	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	B	240	ARG	NE-CZ-NH2	-5.04	117.78	120.30
1	C	327	LEU	CA-CB-CG	5.03	126.87	115.30

There are no chirality outliers.

There are no planarity outliers.

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	3458	0	3477	25	0
1	B	3458	0	3477	32	0
1	C	3458	0	3477	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	20	0	11	1	0
2	B	20	0	11	2	0
2	C	20	0	11	0	0
3	A	33	0	0	0	0
3	B	33	0	0	0	0
3	C	33	0	0	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	A	124	0	0	1	0
5	B	131	0	0	2	0
5	C	134	0	0	0	0
All	All	10924	0	10464	82	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 (82) 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:22:GLU:OE2	1:B:26:ARG:NH2	2.03	0.92
1:B:192:GLN:HG3	1:B:227:HIS:CE1	2.08	0.89
1:B:193:GLN:O	1:B:197:ILE:HG12	1.89	0.72
1:A:218:CYS:SG	5:A:705:HOH:O	2.15	0.70
1:A:85:PHE:HB3	1:A:225:PHE:CE2	2.34	0.62
1:C:67:ASN:O	1:C:71:ASP:OD1	2.17	0.62
1:B:85:PHE:HB3	1:B:225:PHE:CE2	2.39	0.58
1:B:109:LEU:HD22	1:B:119:ILE:HG21	1.86	0.57
1:A:7:LEU:HA	1:A:10:ILE:HG22	1.86	0.56
1:C:109:LEU:HD22	1:C:119:ILE:HG21	1.89	0.55
1:C:85:PHE:HB3	1:C:225:PHE:CE2	2.42	0.55
1:A:263:SER:HA	1:A:266:PHE:O	2.07	0.54
1:C:184:SER:HB3	1:C:371:ARG:HD3	1.90	0.53
1:B:183:THR:HB	2:B:501:PLG:O3	2.08	0.53
1:B:109:LEU:HD21	1:B:179:ILE:HD11	1.91	0.53
1:A:383:LYS:HB2	1:A:385:LYS:HE2	1.91	0.52
1:B:59:PRO:HA	1:B:71:ASP:OD1	2.09	0.52
1:C:229:ASP:OD1	1:C:253:ARG:HD2	2.09	0.52
1:A:195:ARG:HB2	1:A:227:HIS:HB3	1.91	0.52
1:C:161:VAL:HG23	1:C:185:TYR:CZ	2.45	0.51
1:B:93:GLY:HA3	1:B:249:PHE:CE1	2.46	0.51
1:C:7:LEU:HA	1:C:10:ILE:HG22	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:LEU:HB2	1:A:227:HIS:CD2	2.46	0.51
1:B:192:GLN:CG	1:B:227:HIS:CE1	2.90	0.51
1:C:89:ASP:O	1:C:251:LYS:NZ	2.44	0.51
1:C:385:LYS:HE3	1:C:385:LYS:H	1.77	0.50
1:A:174:LYS:HD3	1:A:201:VAL:HG12	1.92	0.50
1:C:22:GLU:OE2	1:C:26:ARG:NH2	2.35	0.50
1:B:427:GLN:HA	1:B:427:GLN:HE21	1.76	0.50
1:C:235:THR:O	1:C:241:GLY:N	2.45	0.49
1:B:235:THR:O	1:B:241:GLY:N	2.43	0.49
1:A:62:ARG:HG2	1:A:67:ASN:ND2	2.28	0.49
1:A:93:GLY:HA3	1:A:249:PHE:CE2	2.48	0.48
1:C:226:LEU:HB2	1:C:227:HIS:CD2	2.49	0.48
1:B:7:LEU:HA	1:B:10:ILE:CG2	2.43	0.48
1:B:3:ASN:ND2	1:B:10:ILE:HD13	2.29	0.48
1:A:113:VAL:CG1	1:A:177:VAL:HG23	2.43	0.48
1:C:319:VAL:HG12	1:C:320:THR:HG23	1.97	0.47
1:A:234:THR:HB	1:A:236:HIS:CE1	2.50	0.47
1:A:398:ILE:O	1:A:402:VAL:HG23	2.15	0.47
1:A:161:VAL:HG23	1:A:185:TYR:CZ	2.49	0.47
1:C:214:SER:HB2	1:C:238:ILE:HG12	1.97	0.47
1:A:72:LYS:HB3	1:B:17:ILE:HD13	1.97	0.47
1:A:127:GLY:O	1:A:182:TYR:HB3	2.15	0.46
1:A:7:LEU:HA	1:A:10:ILE:CG2	2.44	0.46
1:C:426:GLN:O	1:C:430:GLN:HB2	2.16	0.46
1:B:307:LEU:HD21	1:B:372:ILE:HD12	1.98	0.46
1:A:195:ARG:HD2	1:A:227:HIS:O	2.16	0.46
1:C:346:CYS:SG	1:C:353:LEU:HD21	2.55	0.46
1:B:249:PHE:HB2	1:B:257:ILE:HD13	1.98	0.45
1:A:174:LYS:HD3	1:A:201:VAL:CG1	2.46	0.45
1:A:67:ASN:O	1:A:71:ASP:OD1	2.34	0.45
1:C:64:TYR:N	1:C:64:TYR:CD1	2.82	0.45
1:B:111:ALA:CB	1:B:261:ILE:HD12	2.46	0.45
1:C:167:ARG:NH1	1:C:200:GLU:OE1	2.49	0.45
1:C:181:GLY:HA3	1:C:208:ASP:O	2.17	0.45
1:A:396:ARG:CZ	1:A:428:LEU:HD13	2.47	0.44
1:A:3:ASN:ND2	1:A:10:ILE:HD13	2.33	0.44
1:B:263:SER:HA	1:B:266:PHE:O	2.17	0.44
1:B:327:LEU:HD12	1:B:327:LEU:C	2.37	0.43
1:B:7:LEU:HA	1:B:10:ILE:HG22	2.01	0.43
1:C:385:LYS:CE	1:C:385:LYS:H	2.31	0.43
1:C:113:VAL:HA	1:C:176:LYS:HE3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:THR:N	1:A:375:PRO:CD	2.81	0.43
1:B:75:GLU:HA	1:B:78:GLN:OE1	2.19	0.43
1:A:75:GLU:HA	1:A:78:GLN:OE1	2.18	0.43
1:B:209:ILE:HD12	1:B:213:SER:HA	2.00	0.43
1:B:234:THR:HB	1:B:236:HIS:CE1	2.54	0.43
2:A:501:PLG:H4A1	5:B:650:HOH:O	2.18	0.42
1:B:183:THR:HA	1:B:211:HIS:CD2	2.55	0.42
1:A:197:ILE:O	1:A:201:VAL:HG23	2.20	0.42
1:B:253:ARG:NH2	5:B:608:HOH:O	2.53	0.42
1:B:400:ILE:HG21	1:B:425:LEU:HD21	2.01	0.41
1:C:192:GLN:HG3	1:C:227:HIS:CE1	2.56	0.41
1:B:198:CYS:O	1:B:202:ASN:N	2.52	0.41
1:C:374:THR:N	1:C:375:PRO:CD	2.84	0.41
1:B:208:ASP:OD2	2:B:501:PLG:N1	2.54	0.41
1:B:374:THR:N	1:B:375:PRO:CD	2.84	0.41
1:C:127:GLY:O	1:C:182:TYR:HB3	2.21	0.40
1:B:180:CYS:O	1:B:207:ALA:HA	2.21	0.40
1:C:87:VAL:HB	1:C:92:TRP:CD1	2.57	0.40
1:B:113:VAL:CG1	1:B:177:VAL:HG23	2.51	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	440/442 (100%)	423 (96%)	16 (4%)	1 (0%)	47	46
1	B	440/442 (100%)	427 (97%)	12 (3%)	1 (0%)	47	46
1	C	440/442 (100%)	426 (97%)	13 (3%)	1 (0%)	47	46
All	All	1320/1326 (100%)	1276 (97%)	41 (3%)	3 (0%)	47	46

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	237	LYS
1	B	237	LYS
1	C	237	LYS

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	381/381 (100%)	350 (92%)	31 (8%)	11	7
1	B	381/381 (100%)	360 (94%)	21 (6%)	21	17
1	C	381/381 (100%)	356 (93%)	25 (7%)	16	11
All	All	1143/1143 (100%)	1066 (93%)	77 (7%)	16	11

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	TYR
1	A	62	ARG
1	A	64	TYR
1	A	71	ASP
1	A	89	ASP
1	A	100	SER
1	A	136	ASP
1	A	138	LYS
1	A	139	LYS
1	A	151	LYS
1	A	183	THR
1	A	196	GLN
1	A	201	VAL
1	A	206	PHE
1	A	250	ASN
1	A	253	ARG
1	A	254	ASN
1	A	327	LEU
1	A	334	LYS
1	A	341	LYS

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Mol	Chain	Res	Type
1	A	352	SER
1	A	364	CYS
1	A	366	SER
1	A	385	LYS
1	A	399	LYS
1	A	417	LYS
1	A	419	LEU
1	A	427	GLN
1	A	429	LYS
1	A	430	GLN
1	A	434	THR
1	B	22	GLU
1	B	54	TYR
1	B	62	ARG
1	B	64	TYR
1	B	100	SER
1	B	136	ASP
1	B	138	LYS
1	B	183	THR
1	B	192	GLN
1	B	206	PHE
1	B	244	SER
1	B	254	ASN
1	B	327	LEU
1	B	334	LYS
1	B	352	SER
1	B	362	VAL
1	B	364	CYS
1	B	366	SER
1	B	385	LYS
1	B	427	GLN
1	B	430	GLN
1	C	22	GLU
1	C	23	LYS
1	C	54	TYR
1	C	62	ARG
1	C	64	TYR
1	C	100	SER
1	C	138	LYS
1	C	139	LYS
1	C	183	THR
1	C	184	SER

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Mol	Chain	Res	Type
1	C	196	GLN
1	C	201	VAL
1	C	206	PHE
1	C	240	ARG
1	C	250	ASN
1	C	254	ASN
1	C	263	SER
1	C	327	LEU
1	C	364	CYS
1	C	366	SER
1	C	385	LYS
1	C	399	LYS
1	C	406	GLU
1	C	419	LEU
1	C	430	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	GLN
1	A	196	GLN
1	A	227	HIS
1	A	299	GLN
1	A	356	ASN
1	A	427	GLN
1	B	219	ASN
1	B	227	HIS
1	B	299	GLN
1	B	356	ASN
1	B	427	GLN
1	C	227	HIS
1	C	250	ASN
1	C	299	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	PLG	A	501	-	17,20,20	2.57	3 (17%)	23,28,28	1.95	6 (26%)
2	PLG	C	501	-	17,20,20	3.05	4 (23%)	23,28,28	2.41	10 (43%)
3	8UO	A	502	-	28,36,36	2.76	4 (14%)	37,54,54	2.44	14 (37%)
2	PLG	B	501	-	17,20,20	2.68	3 (17%)	23,28,28	2.06	7 (30%)
3	8UO	C	502	-	28,36,36	2.79	3 (10%)	37,54,54	2.29	11 (29%)
3	8UO	B	502	-	28,36,36	2.73	2 (7%)	37,54,54	2.26	14 (37%)

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	PLG	A	501	-	-	6/10/12/12	0/1/1/1
2	PLG	C	501	-	-	6/10/12/12	0/1/1/1
3	8UO	A	502	-	-	2/18/52/52	0/4/4/4
2	PLG	B	501	-	-	6/10/12/12	0/1/1/1
3	8UO	C	502	-	-	0/18/52/52	0/4/4/4
3	8UO	B	502	-	-	2/18/52/52	0/4/4/4

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	502	8UO	C8-C3	-11.30	1.27	1.42
3	A	502	8UO	C8-C3	-11.09	1.27	1.42
3	B	502	8UO	C8-C3	-10.77	1.27	1.42
3	B	502	8UO	O1-C2	8.41	1.46	1.36
3	C	502	8UO	O1-C2	8.23	1.46	1.36
2	C	501	PLG	C3-C2	8.22	1.49	1.40
3	A	502	8UO	O1-C2	7.94	1.45	1.36
2	B	501	PLG	C3-C2	7.79	1.48	1.40
2	A	501	PLG	C3-C2	7.12	1.48	1.40
2	C	501	PLG	C5-C4	6.83	1.50	1.40
2	C	501	PLG	C3-C4	5.68	1.48	1.40
2	A	501	PLG	C5-C4	5.41	1.48	1.40
2	B	501	PLG	C5-C4	5.21	1.47	1.40
2	A	501	PLG	C3-C4	5.06	1.47	1.40
2	B	501	PLG	C3-C4	4.84	1.47	1.40
3	A	502	8UO	C20-C16	2.20	1.43	1.39
3	A	502	8UO	C24-N23	2.17	1.50	1.46
3	C	502	8UO	N12-N11	2.08	1.41	1.37
2	C	501	PLG	C4A-C4	2.03	1.54	1.51

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	8UO	C28-N23-C24	6.81	126.54	111.52
3	C	502	8UO	C28-N23-C24	6.70	126.32	111.52
3	A	502	8UO	O1-C2-N7	6.23	114.69	110.22
3	B	502	8UO	C28-N23-C24	6.21	125.23	111.52
2	C	501	PLG	OP4-C5A-C5	6.10	120.97	109.35
3	C	502	8UO	O1-C2-N7	6.07	114.58	110.22
3	B	502	8UO	O1-C2-N7	5.06	113.85	110.22
2	B	501	PLG	OP4-C5A-C5	4.82	118.54	109.35
3	C	502	8UO	C22-C14-C4	4.49	117.21	112.40
3	A	502	8UO	C3-C2-N7	-4.27	123.20	127.47
2	A	501	PLG	OP4-C5A-C5	4.20	117.35	109.35
3	C	502	8UO	C3-C2-N7	-4.03	123.44	127.47
2	A	501	PLG	CA-N-C4A	3.98	122.57	112.54
2	C	501	PLG	CA-N-C4A	3.89	122.36	112.54
3	A	502	8UO	C30-C19-N23	-3.72	117.33	121.33
3	A	502	8UO	C20-C19-N23	3.71	125.32	121.33
2	C	501	PLG	C4A-C4-C5	3.62	123.73	119.71
3	B	502	8UO	C3-C2-N7	-3.59	123.88	127.47
2	C	501	PLG	C3-C4-C5	-3.57	115.29	118.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	8UO	C30-C19-N23	-3.51	117.55	121.33
2	B	501	PLG	CA-N-C4A	3.51	121.38	112.54
2	B	501	PLG	OP3-P-OP4	-3.44	97.58	106.73
3	B	502	8UO	C5-C4-C3	3.43	111.31	104.34
2	B	501	PLG	OP3-P-OP2	3.40	120.62	107.64
3	A	502	8UO	C22-C14-C4	3.22	115.85	112.40
3	B	502	8UO	C22-C14-C4	3.17	115.80	112.40
3	A	502	8UO	C14-C4-C3	-3.11	105.32	110.32
3	B	502	8UO	C14-C4-C3	-3.11	105.33	110.32
3	B	502	8UO	C20-C19-N23	3.07	124.64	121.33
2	C	501	PLG	OP2-P-OP4	-3.06	98.58	106.73
3	C	502	8UO	C30-C19-N23	-3.03	118.07	121.33
3	A	502	8UO	C15-C14-C4	3.00	115.62	112.40
2	C	501	PLG	C4-C4A-N	3.00	117.19	111.58
2	C	501	PLG	OP3-P-OP4	-2.99	98.79	106.73
3	B	502	8UO	C8-C3-C2	2.97	121.58	117.64
3	A	502	8UO	C5-C4-C3	2.92	110.29	104.34
3	C	502	8UO	C16-C17-C18	2.89	121.43	118.30
3	B	502	8UO	C16-C17-C18	2.81	121.34	118.30
2	A	501	PLG	OP3-P-OP2	2.80	118.35	107.64
2	A	501	PLG	C6-N1-C2	2.73	124.22	119.17
3	A	502	8UO	C16-C17-C18	2.69	121.22	118.30
3	C	502	8UO	C14-C4-C3	-2.68	106.01	110.32
3	C	502	8UO	C5-C4-C3	2.68	109.79	104.34
3	B	502	8UO	C19-C30-C18	2.66	121.10	118.06
2	B	501	PLG	C6-N1-C2	2.59	123.96	119.17
2	A	501	PLG	OP3-P-OP4	-2.55	99.96	106.73
3	C	502	8UO	C6-C5-C10	-2.51	101.36	104.01
3	A	502	8UO	C19-C30-C18	2.45	120.85	118.06
2	A	501	PLG	OP2-P-OP4	-2.41	100.31	106.73
2	C	501	PLG	C5A-C5-C6	-2.40	115.42	119.37
2	B	501	PLG	C4A-C4-C3	2.39	122.61	120.04
3	A	502	8UO	C13-C10-N11	2.36	124.84	119.65
3	B	502	8UO	C15-C14-C4	2.34	114.91	112.40
3	A	502	8UO	C8-C3-C2	2.34	120.74	117.64
3	B	502	8UO	C17-C18-C30	-2.32	120.59	123.52
3	A	502	8UO	C20-C16-C4	2.26	124.55	120.67
3	C	502	8UO	C17-C18-C30	-2.24	120.69	123.52
3	B	502	8UO	C13-C10-N11	2.23	124.55	119.65
2	C	501	PLG	C6-N1-C2	2.21	123.26	119.17
2	C	501	PLG	OP3-P-OP2	2.14	115.82	107.64
2	B	501	PLG	C4-C4A-N	2.02	115.35	111.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	502	8UO	C13-C10-N11	2.01	124.06	119.65

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	PLG	C5-C4-C4A-N
2	A	501	PLG	C5A-OP4-P-OP2
2	A	501	PLG	C5A-OP4-P-OP3
2	C	501	PLG	C5A-OP4-P-OP2
2	C	501	PLG	C5A-OP4-P-OP3
2	B	501	PLG	C5-C4-C4A-N
2	B	501	PLG	C5A-OP4-P-OP1
2	B	501	PLG	C5A-OP4-P-OP2
2	B	501	PLG	C5A-OP4-P-OP3
3	B	502	8UO	C25-C26-C29-C31
3	B	502	8UO	C27-C26-C29-C31
2	A	501	PLG	C4-C4A-N-CA
2	C	501	PLG	C5-C4-C4A-N
2	B	501	PLG	C4-C4A-N-CA
3	A	502	8UO	C25-C26-C29-C31
3	A	502	8UO	C27-C26-C29-C31
2	C	501	PLG	C-CA-N-C4A
2	A	501	PLG	C3-C4-C4A-N
2	C	501	PLG	C3-C4-C4A-N
2	B	501	PLG	C3-C4-C4A-N
2	C	501	PLG	C5A-OP4-P-OP1
2	A	501	PLG	C-CA-N-C4A

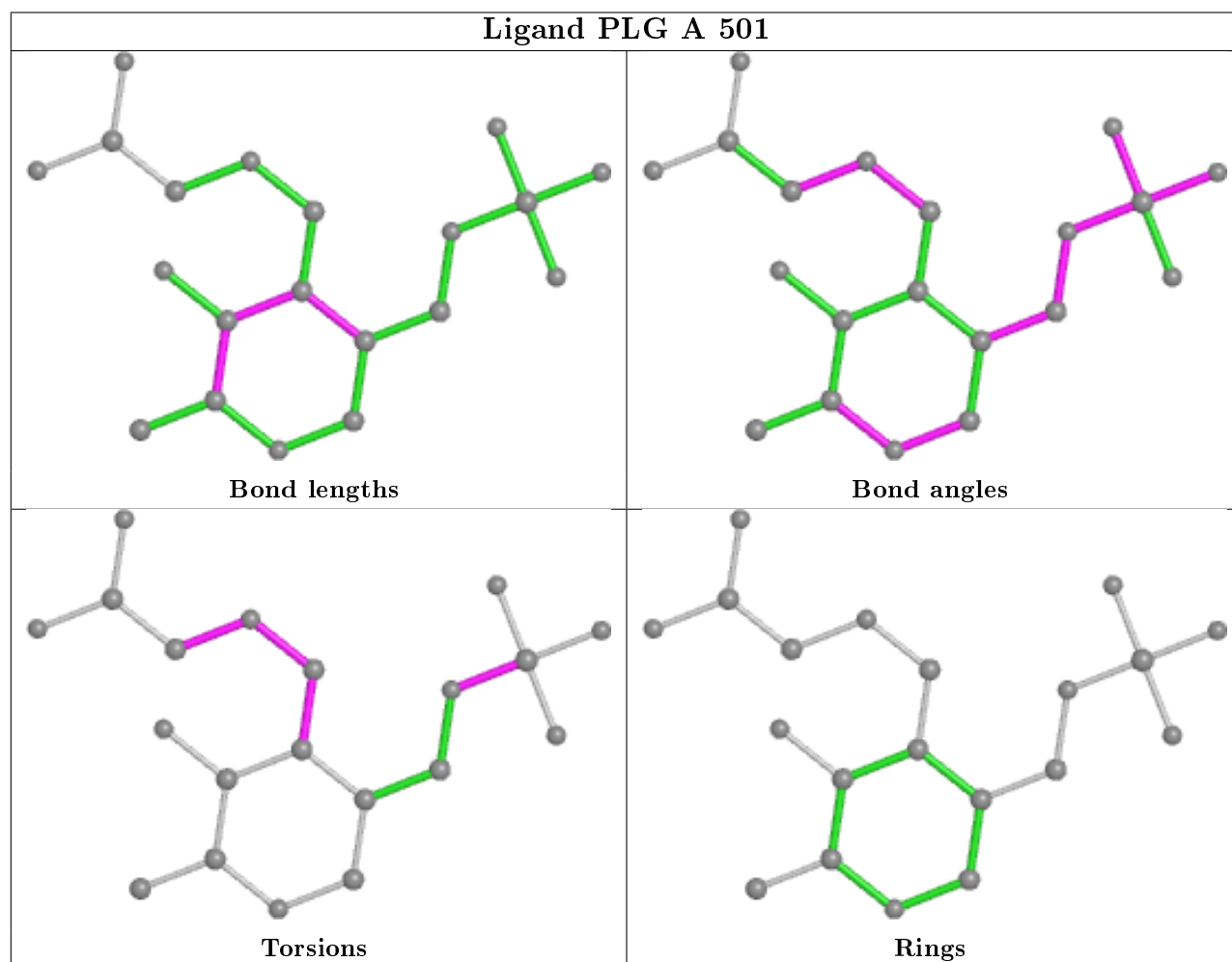
There are no ring outliers.

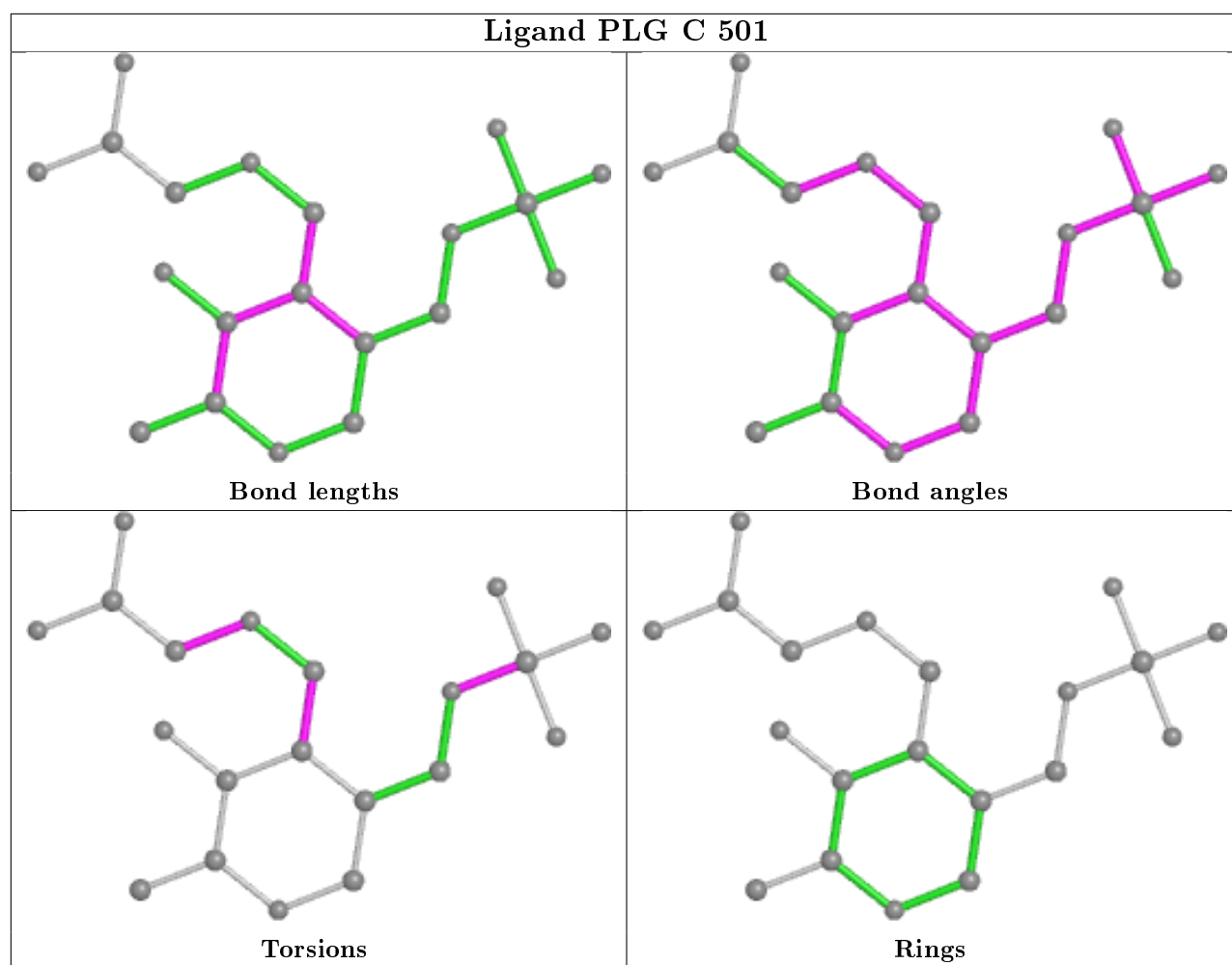
2 monomers are involved in 3 short contacts:

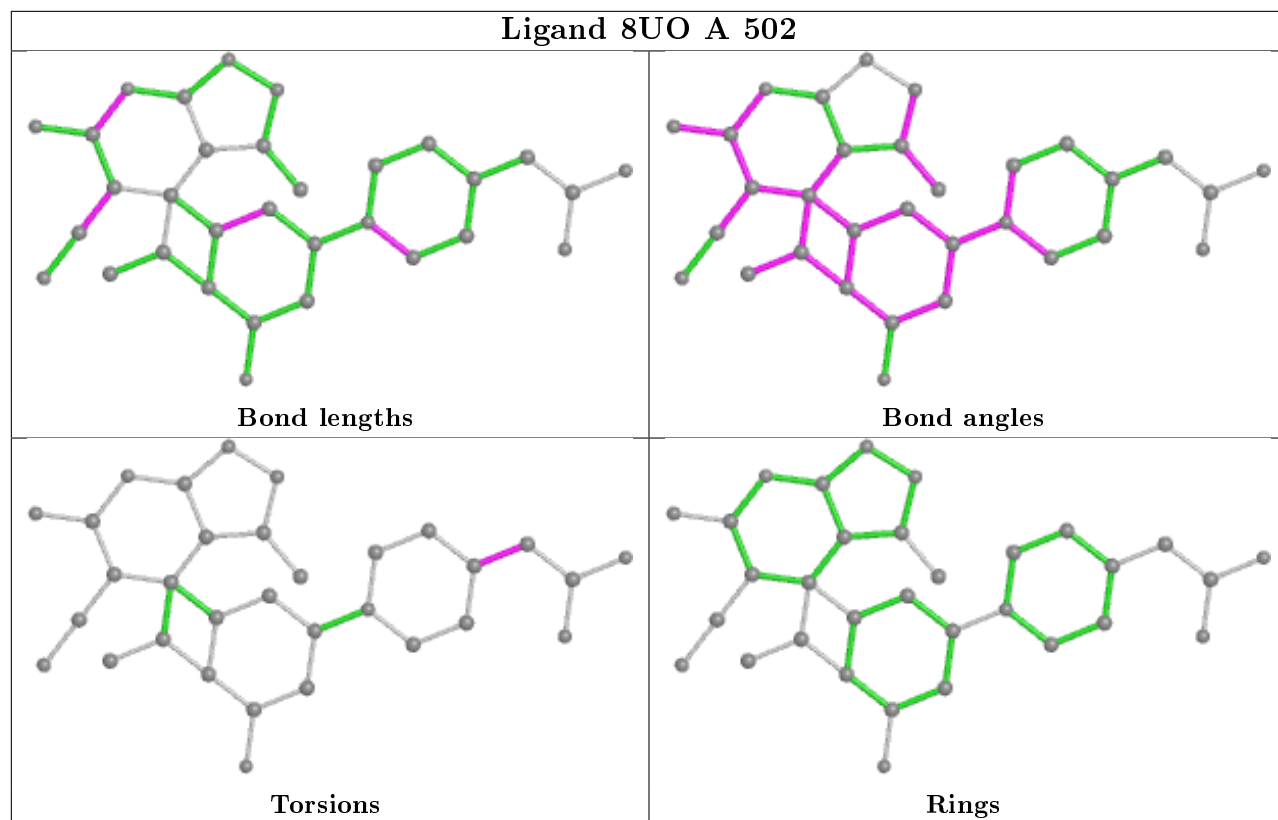
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	PLG	1	0
2	B	501	PLG	2	0

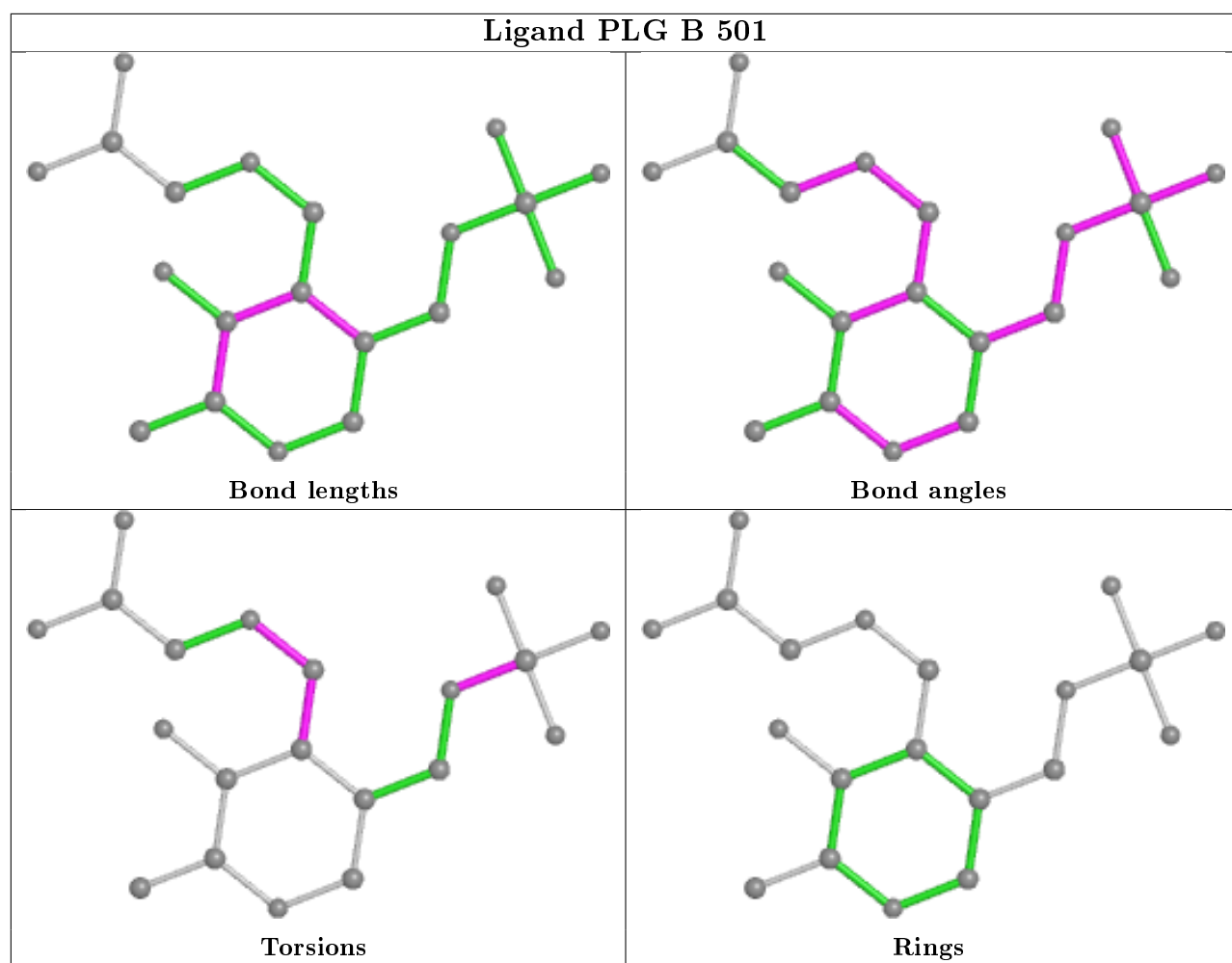
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

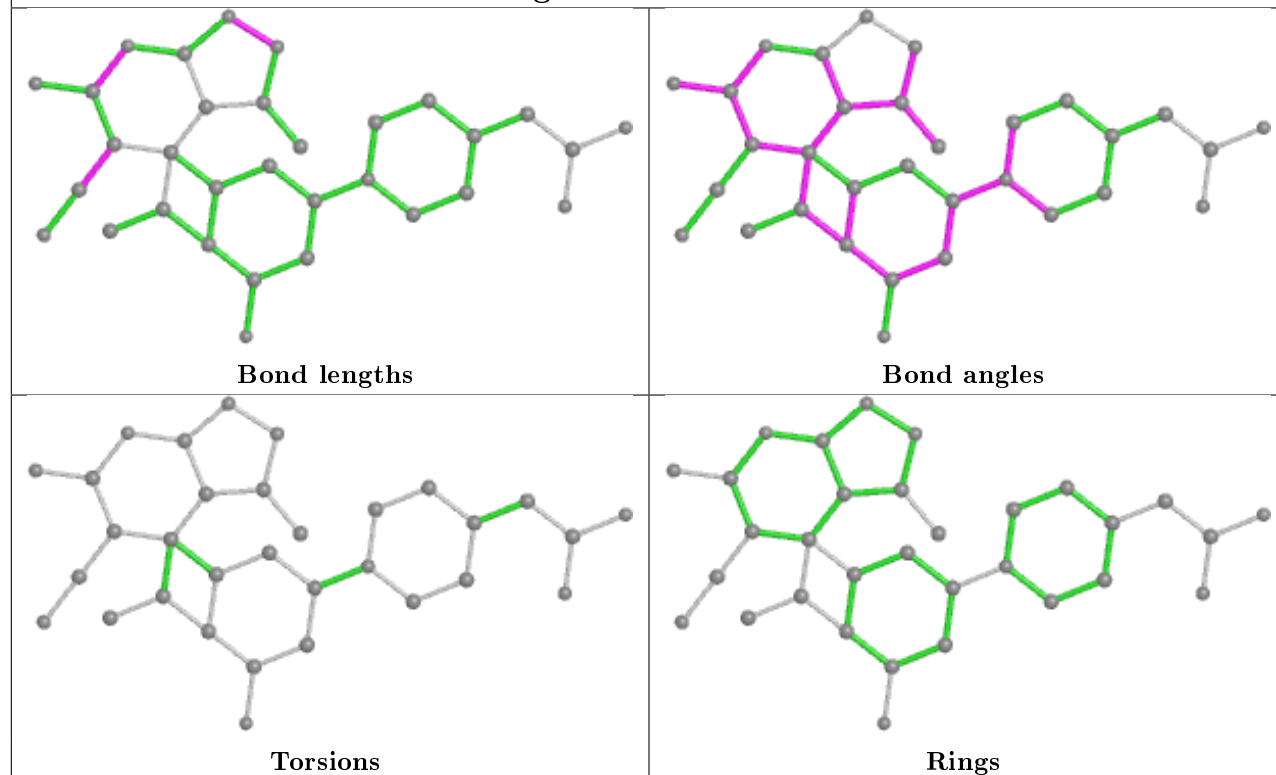




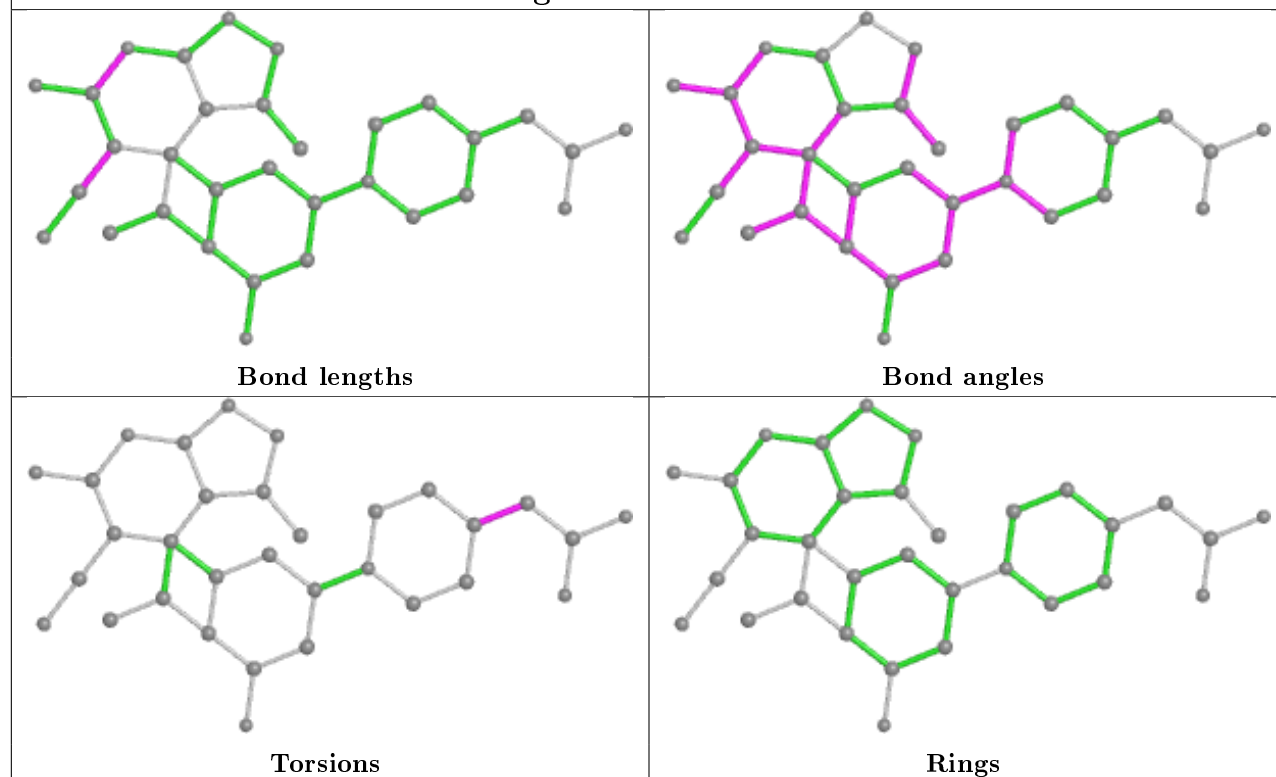




Ligand 8UO C 502



Ligand 8UO B 502



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	442/442 (100%)	0.84	37 (8%)	11 15	13, 28, 49, 78	0
1	B	442/442 (100%)	0.77	25 (5%)	23 32	13, 27, 48, 74	0
1	C	442/442 (100%)	0.75	23 (5%)	27 35	12, 26, 49, 71	0
All	All	1326/1326 (100%)	0.79	85 (6%)	19 26	12, 27, 49, 78	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	413	VAL	6.7
1	A	412	LEU	4.4
1	C	64	TYR	3.6
1	A	255	PRO	3.6
1	B	138	LYS	3.4
1	A	402	VAL	3.4
1	A	166	VAL	3.4
1	B	408	TYR	3.3
1	A	138	LYS	3.3
1	B	365	VAL	3.2
1	B	415	PHE	3.2
1	C	423	ALA	3.1
1	A	421	GLY	3.0
1	B	64	TYR	3.0
1	A	363	ASP	2.9
1	C	169	MET	2.9
1	A	413	VAL	2.9
1	A	198	CYS	2.9
1	A	202	ASN	2.8
1	A	312	ILE	2.8
1	B	412	LEU	2.8
1	B	10	ILE	2.8
1	A	90	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	194	PHE	2.8
1	C	138	LYS	2.8
1	A	88	SER	2.8
1	A	408	TYR	2.7
1	A	365	VAL	2.7
1	C	189	ILE	2.7
1	C	259	GLN	2.7
1	A	137	GLU	2.6
1	B	413	VAL	2.6
1	C	362	VAL	2.6
1	B	116	LYS	2.6
1	B	419	LEU	2.6
1	C	202	ASN	2.5
1	C	2	PHE	2.5
1	A	364	CYS	2.5
1	A	409	GLY	2.5
1	A	7	LEU	2.5
1	B	215	PHE	2.5
1	A	155	CYS	2.4
1	A	9	GLN	2.4
1	C	9	GLN	2.4
1	A	419	LEU	2.4
1	A	203	ALA	2.4
1	C	137	GLU	2.4
1	A	436	ALA	2.4
1	C	419	LEU	2.3
1	A	124	LEU	2.3
1	B	366	SER	2.3
1	B	364	CYS	2.3
1	A	257	ILE	2.3
1	A	169	MET	2.3
1	A	109	LEU	2.3
1	C	193	GLN	2.3
1	A	433	VAL	2.3
1	B	362	VAL	2.2
1	B	203	ALA	2.2
1	C	410	LYS	2.2
1	A	331	ASP	2.2
1	C	335	PHE	2.2
1	B	426	GLN	2.2
1	A	93	GLY	2.2
1	B	201	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	346	CYS	2.1
1	A	411	LYS	2.1
1	C	412	LEU	2.1
1	A	73	ILE	2.1
1	B	196	GLN	2.1
1	B	96	VAL	2.1
1	A	389	PHE	2.1
1	B	193	GLN	2.1
1	A	201	VAL	2.1
1	C	173	PHE	2.1
1	B	255	PRO	2.0
1	C	226	LEU	2.0
1	C	228	ALA	2.0
1	B	409	GLY	2.0
1	C	65	GLY	2.0
1	A	57	GLY	2.0
1	B	389	PHE	2.0
1	C	90	GLU	2.0
1	B	239	LEU	2.0
1	B	87	VAL	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	8UO	B	502	33/33	0.75	0.27	34,38,52,54	0
3	8UO	A	502	33/33	0.82	0.19	31,35,51,52	0
4	CL	A	503	1/1	0.83	0.09	41,41,41,41	0

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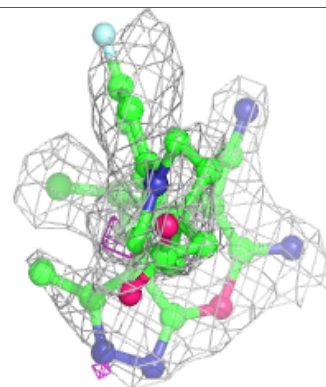
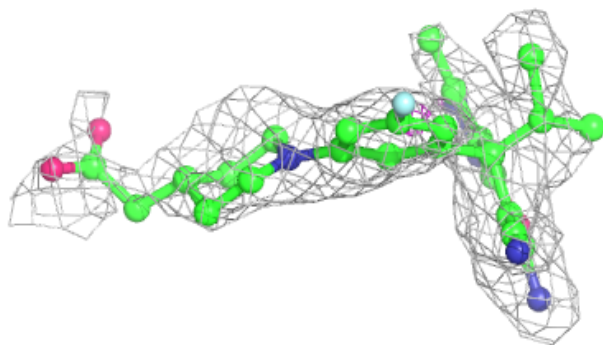
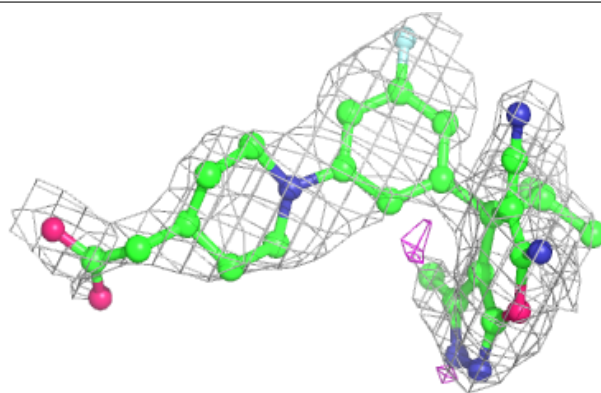
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	8UO	C	502	33/33	0.84	0.19	30,34,48,50	0
4	CL	C	503	1/1	0.88	0.13	15,15,15,15	1
2	PLG	C	501	20/20	0.90	0.15	15,19,20,20	0
2	PLG	A	501	20/20	0.91	0.15	17,21,27,30	0
2	PLG	B	501	20/20	0.91	0.16	16,21,23,23	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.

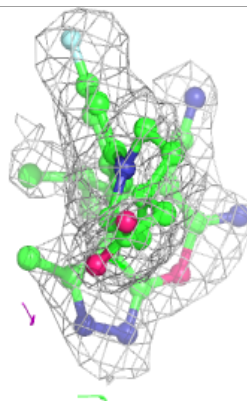
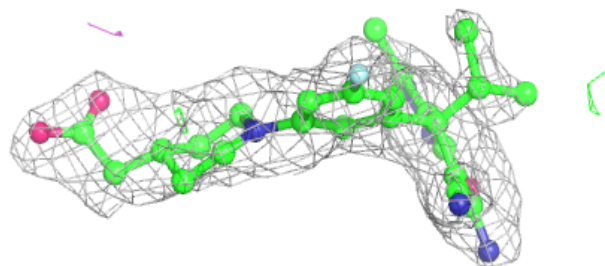
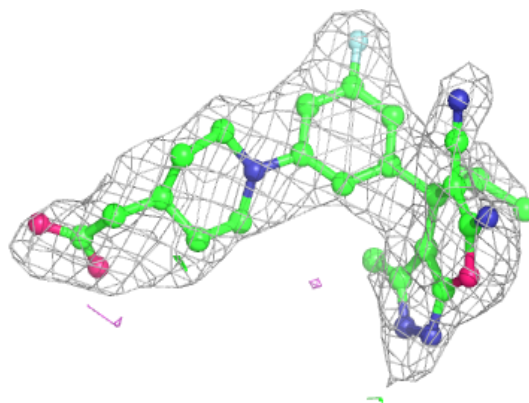
Electron density around 8UO B 502:

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

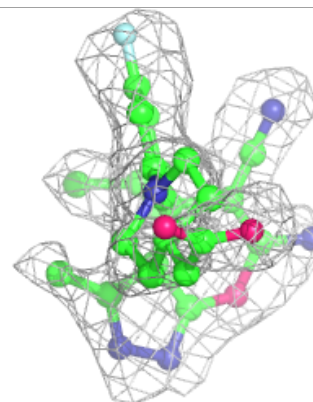
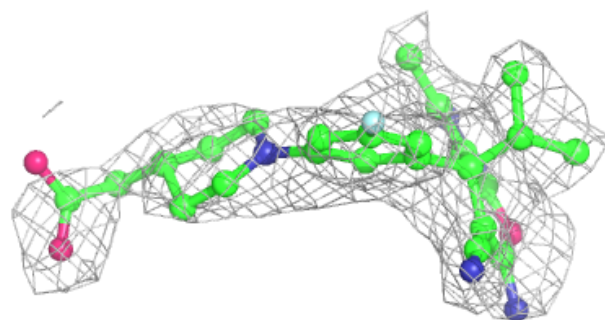
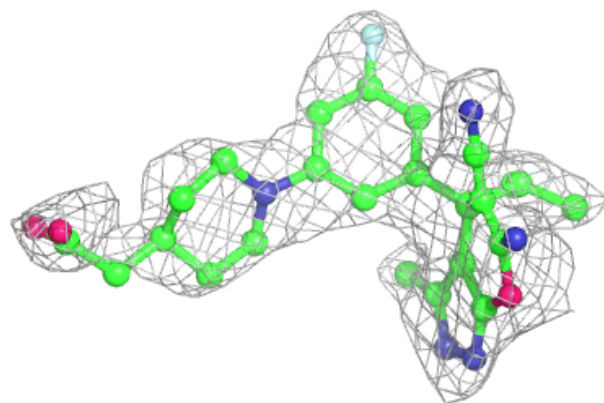


Electron density around 8UO A 502:

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

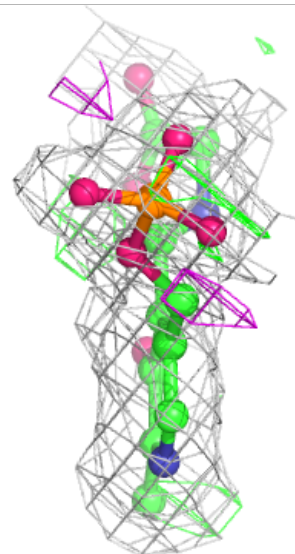
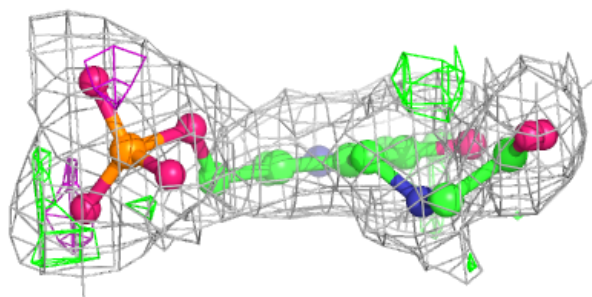
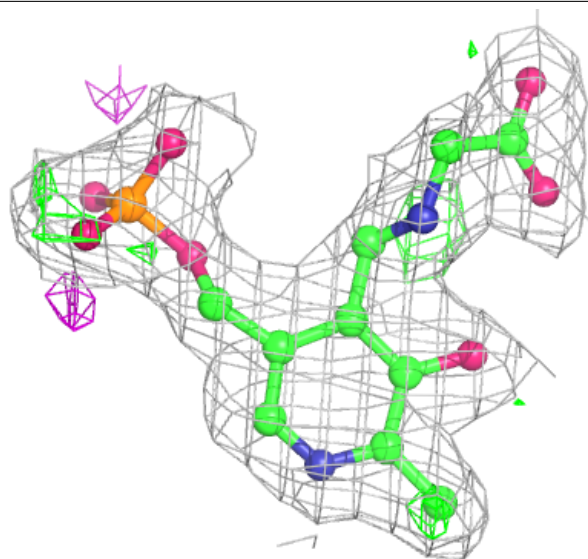
**Electron density around 8UO C 502:**

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



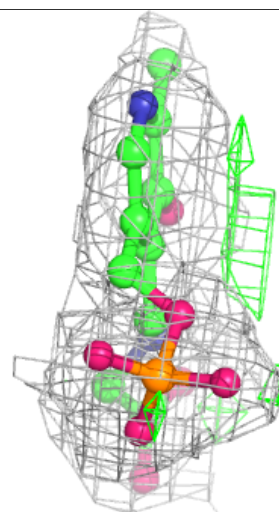
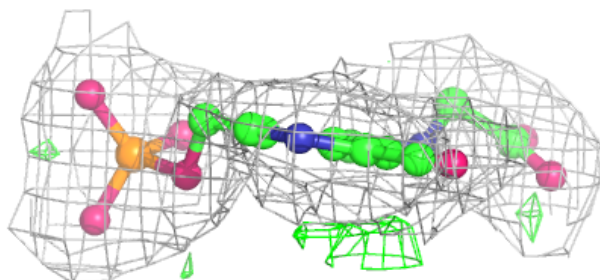
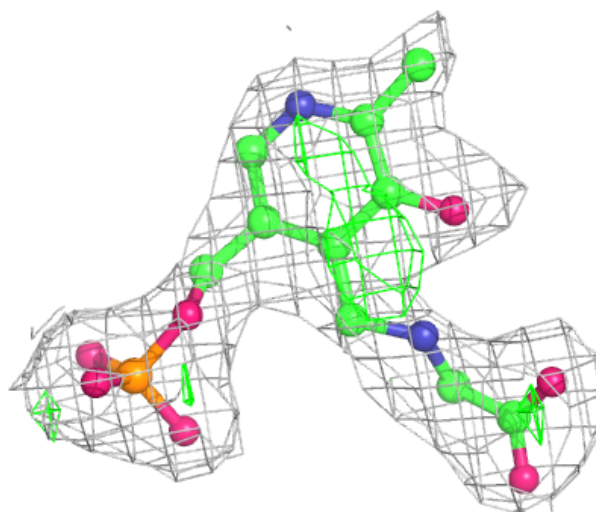
Electron density around PLG C 501:

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



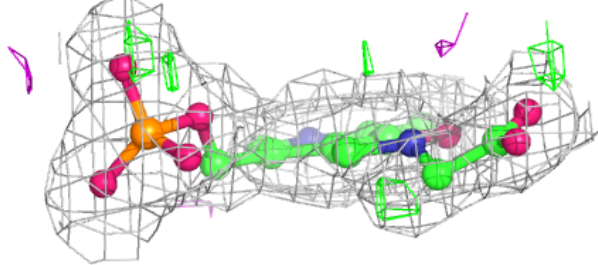
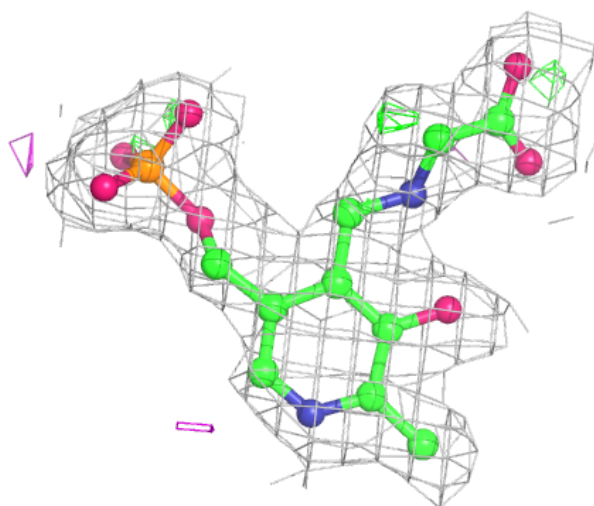
Electron density around PLG A 501:

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



Electron density around PLG B 501:

$2mF_o-DF_c$ (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.