



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 14, 2026 – 02:09 PM EDT

PDB ID : 9Y6D / pdb_00009y6d
Title : CRYSTAL STRUCTURE OF THE A149T VARIANT OF SERINE HYDROXYMETHYLTRANSFERASE 8 FROM SOYBEAN CULTIVAR FORREST IN COMPLEX WITH PLP-GLYCINE
Authors : Beamer, L.J.; Samarakoon, V.; Owuocha, L.F.
Deposited on : 2025-09-08
Resolution : 2.25 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

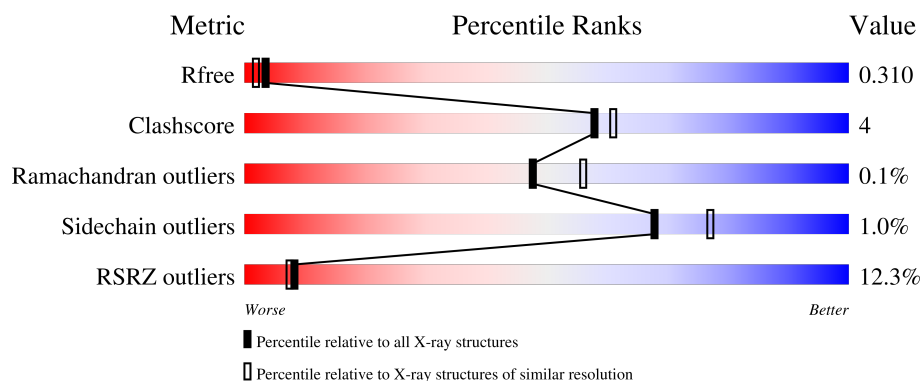
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	496	<div> <div>12%</div> <div> <div></div> <div>83%</div> <div>11%</div> <div>6%</div> </div> </div>
1	B	496	<div> <div>7%</div> <div> <div></div> <div>86%</div> <div>8%</div> <div>6%</div> </div> </div>
1	C	496	<div> <div>10%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>6%</div> </div> </div>
1	D	496	<div> <div>8%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>7%</div> </div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	E	496	<div><div></div><div></div><div></div><div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17890 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	467	Total	C	N	O	S	0	0	0
			3450	2184	595	654	17			
1	B	467	Total	C	N	O	S	0	0	0
			3502	2225	597	663	17			
1	C	466	Total	C	N	O	S	0	0	0
			3496	2223	599	657	17			
1	D	462	Total	C	N	O	S	0	0	0
			3481	2211	595	658	17			
1	E	449	Total	C	N	O	S	0	0	0
			3253	2052	559	626	16			

There are 130 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	initiating methionine	UNP K4FW35
A	-23	GLY	-	expression tag	UNP K4FW35
A	-22	MET	-	expression tag	UNP K4FW35
A	-21	HIS	-	expression tag	UNP K4FW35
A	-20	HIS	-	expression tag	UNP K4FW35
A	-19	HIS	-	expression tag	UNP K4FW35
A	-18	HIS	-	expression tag	UNP K4FW35
A	-17	HIS	-	expression tag	UNP K4FW35
A	-16	HIS	-	expression tag	UNP K4FW35
A	-15	SER	-	expression tag	UNP K4FW35
A	-14	SER	-	expression tag	UNP K4FW35
A	-13	GLY	-	expression tag	UNP K4FW35
A	-12	VAL	-	expression tag	UNP K4FW35
A	-11	ASP	-	expression tag	UNP K4FW35
A	-10	LEU	-	expression tag	UNP K4FW35
A	-9	GLY	-	expression tag	UNP K4FW35
A	-8	THR	-	expression tag	UNP K4FW35
A	-7	GLU	-	expression tag	UNP K4FW35
A	-6	ASN	-	expression tag	UNP K4FW35

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	LEU	-	expression tag	UNP K4FW35
A	-4	TYR	-	expression tag	UNP K4FW35
A	-3	PHE	-	expression tag	UNP K4FW35
A	-2	GLN	-	expression tag	UNP K4FW35
A	-1	SER	-	expression tag	UNP K4FW35
A	0	ASN	-	expression tag	UNP K4FW35
A	149	THR	ALA	engineered mutation	UNP K4FW35
B	-24	MET	-	initiating methionine	UNP K4FW35
B	-23	GLY	-	expression tag	UNP K4FW35
B	-22	MET	-	expression tag	UNP K4FW35
B	-21	HIS	-	expression tag	UNP K4FW35
B	-20	HIS	-	expression tag	UNP K4FW35
B	-19	HIS	-	expression tag	UNP K4FW35
B	-18	HIS	-	expression tag	UNP K4FW35
B	-17	HIS	-	expression tag	UNP K4FW35
B	-16	HIS	-	expression tag	UNP K4FW35
B	-15	SER	-	expression tag	UNP K4FW35
B	-14	SER	-	expression tag	UNP K4FW35
B	-13	GLY	-	expression tag	UNP K4FW35
B	-12	VAL	-	expression tag	UNP K4FW35
B	-11	ASP	-	expression tag	UNP K4FW35
B	-10	LEU	-	expression tag	UNP K4FW35
B	-9	GLY	-	expression tag	UNP K4FW35
B	-8	THR	-	expression tag	UNP K4FW35
B	-7	GLU	-	expression tag	UNP K4FW35
B	-6	ASN	-	expression tag	UNP K4FW35
B	-5	LEU	-	expression tag	UNP K4FW35
B	-4	TYR	-	expression tag	UNP K4FW35
B	-3	PHE	-	expression tag	UNP K4FW35
B	-2	GLN	-	expression tag	UNP K4FW35
B	-1	SER	-	expression tag	UNP K4FW35
B	0	ASN	-	expression tag	UNP K4FW35
B	149	THR	ALA	engineered mutation	UNP K4FW35
C	-24	MET	-	initiating methionine	UNP K4FW35
C	-23	GLY	-	expression tag	UNP K4FW35
C	-22	MET	-	expression tag	UNP K4FW35
C	-21	HIS	-	expression tag	UNP K4FW35
C	-20	HIS	-	expression tag	UNP K4FW35
C	-19	HIS	-	expression tag	UNP K4FW35
C	-18	HIS	-	expression tag	UNP K4FW35
C	-17	HIS	-	expression tag	UNP K4FW35
C	-16	HIS	-	expression tag	UNP K4FW35

Continued on next page...

Continued from previous page...

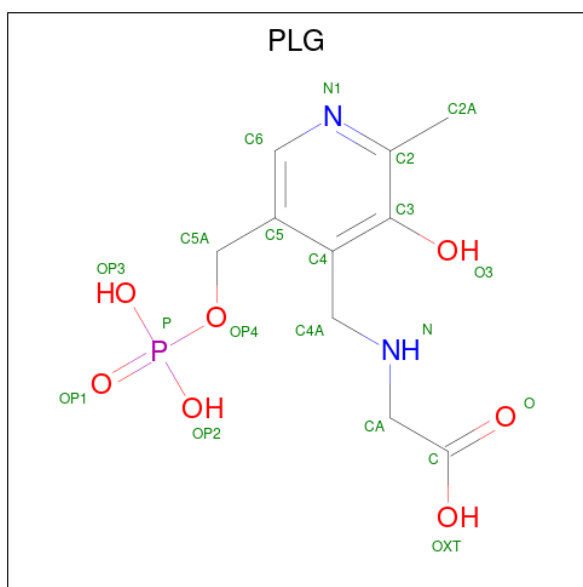
Chain	Residue	Modelled	Actual	Comment	Reference
C	-15	SER	-	expression tag	UNP K4FW35
C	-14	SER	-	expression tag	UNP K4FW35
C	-13	GLY	-	expression tag	UNP K4FW35
C	-12	VAL	-	expression tag	UNP K4FW35
C	-11	ASP	-	expression tag	UNP K4FW35
C	-10	LEU	-	expression tag	UNP K4FW35
C	-9	GLY	-	expression tag	UNP K4FW35
C	-8	THR	-	expression tag	UNP K4FW35
C	-7	GLU	-	expression tag	UNP K4FW35
C	-6	ASN	-	expression tag	UNP K4FW35
C	-5	LEU	-	expression tag	UNP K4FW35
C	-4	TYR	-	expression tag	UNP K4FW35
C	-3	PHE	-	expression tag	UNP K4FW35
C	-2	GLN	-	expression tag	UNP K4FW35
C	-1	SER	-	expression tag	UNP K4FW35
C	0	ASN	-	expression tag	UNP K4FW35
C	149	THR	ALA	engineered mutation	UNP K4FW35
D	-24	MET	-	initiating methionine	UNP K4FW35
D	-23	GLY	-	expression tag	UNP K4FW35
D	-22	MET	-	expression tag	UNP K4FW35
D	-21	HIS	-	expression tag	UNP K4FW35
D	-20	HIS	-	expression tag	UNP K4FW35
D	-19	HIS	-	expression tag	UNP K4FW35
D	-18	HIS	-	expression tag	UNP K4FW35
D	-17	HIS	-	expression tag	UNP K4FW35
D	-16	HIS	-	expression tag	UNP K4FW35
D	-15	SER	-	expression tag	UNP K4FW35
D	-14	SER	-	expression tag	UNP K4FW35
D	-13	GLY	-	expression tag	UNP K4FW35
D	-12	VAL	-	expression tag	UNP K4FW35
D	-11	ASP	-	expression tag	UNP K4FW35
D	-10	LEU	-	expression tag	UNP K4FW35
D	-9	GLY	-	expression tag	UNP K4FW35
D	-8	THR	-	expression tag	UNP K4FW35
D	-7	GLU	-	expression tag	UNP K4FW35
D	-6	ASN	-	expression tag	UNP K4FW35
D	-5	LEU	-	expression tag	UNP K4FW35
D	-4	TYR	-	expression tag	UNP K4FW35
D	-3	PHE	-	expression tag	UNP K4FW35
D	-2	GLN	-	expression tag	UNP K4FW35
D	-1	SER	-	expression tag	UNP K4FW35
D	0	ASN	-	expression tag	UNP K4FW35

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	149	THR	ALA	engineered mutation	UNP K4FW35
E	-24	MET	-	initiating methionine	UNP K4FW35
E	-23	GLY	-	expression tag	UNP K4FW35
E	-22	MET	-	expression tag	UNP K4FW35
E	-21	HIS	-	expression tag	UNP K4FW35
E	-20	HIS	-	expression tag	UNP K4FW35
E	-19	HIS	-	expression tag	UNP K4FW35
E	-18	HIS	-	expression tag	UNP K4FW35
E	-17	HIS	-	expression tag	UNP K4FW35
E	-16	HIS	-	expression tag	UNP K4FW35
E	-15	SER	-	expression tag	UNP K4FW35
E	-14	SER	-	expression tag	UNP K4FW35
E	-13	GLY	-	expression tag	UNP K4FW35
E	-12	VAL	-	expression tag	UNP K4FW35
E	-11	ASP	-	expression tag	UNP K4FW35
E	-10	LEU	-	expression tag	UNP K4FW35
E	-9	GLY	-	expression tag	UNP K4FW35
E	-8	THR	-	expression tag	UNP K4FW35
E	-7	GLU	-	expression tag	UNP K4FW35
E	-6	ASN	-	expression tag	UNP K4FW35
E	-5	LEU	-	expression tag	UNP K4FW35
E	-4	TYR	-	expression tag	UNP K4FW35
E	-3	PHE	-	expression tag	UNP K4FW35
E	-2	GLN	-	expression tag	UNP K4FW35
E	-1	SER	-	expression tag	UNP K4FW35
E	0	ASN	-	expression tag	UNP K4FW35
E	149	THR	ALA	engineered mutation	UNP K4FW35

- Molecule 2 is N-GLYCINE-[3-HYDROXY-2-METHYL-5-PHOSPHONOOXYMETHYL-PYRIDIN-4-YL-METHANE] (CCD ID: PLG) (formula: C₁₀H₁₅N₂O₇P) (labeled as "Ligand of Interest" by depositor).



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		
2	D	1	Total	C	N	O	P	0	0
			20	10	2	7	1		
2	E	1	Total	C	N	O	P	0	0
			20	10	2	7	1		

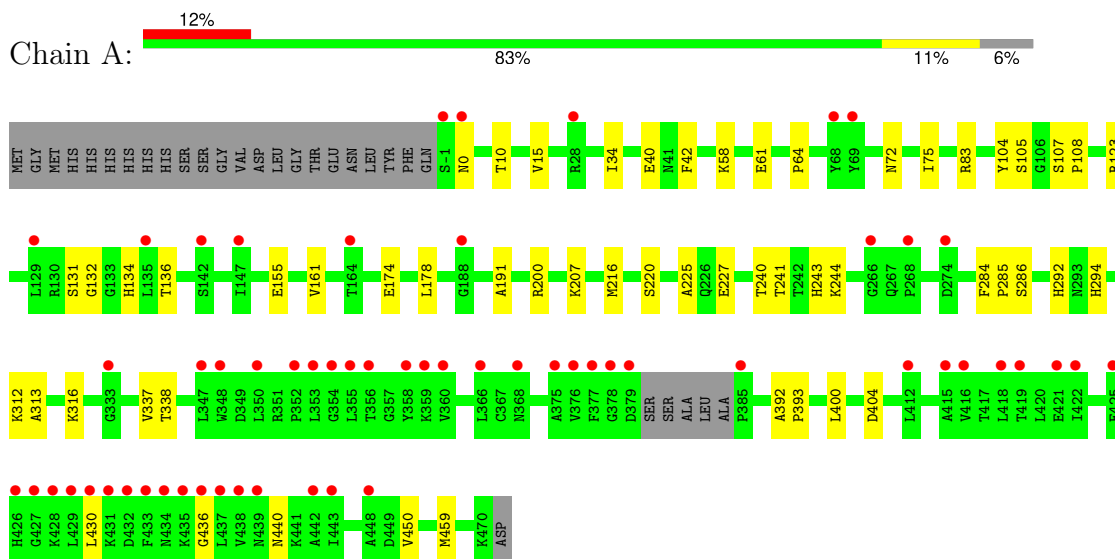
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	158	Total	O	0	0
			158	158		
3	B	129	Total	O	0	0
			129	129		
3	C	141	Total	O	0	0
			141	141		
3	D	126	Total	O	0	0
			126	126		
3	E	54	Total	O	0	0
			54	54		

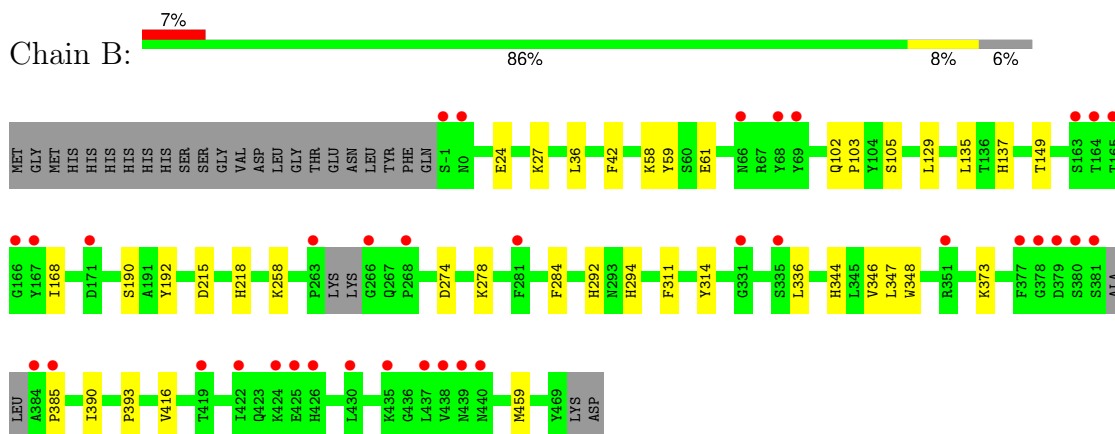
3 Residue-property plots [i](#)

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.

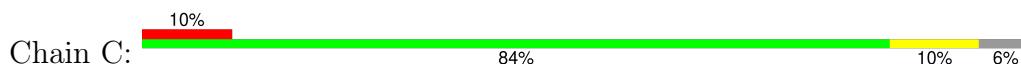
- Molecule 1: Serine hydroxymethyltransferase

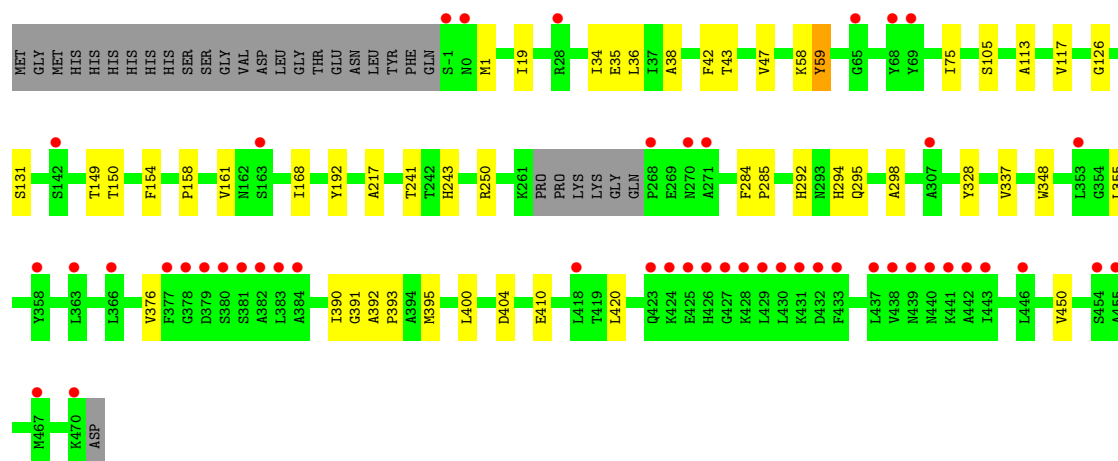


- Molecule 1: Serine hydroxymethyltransferase

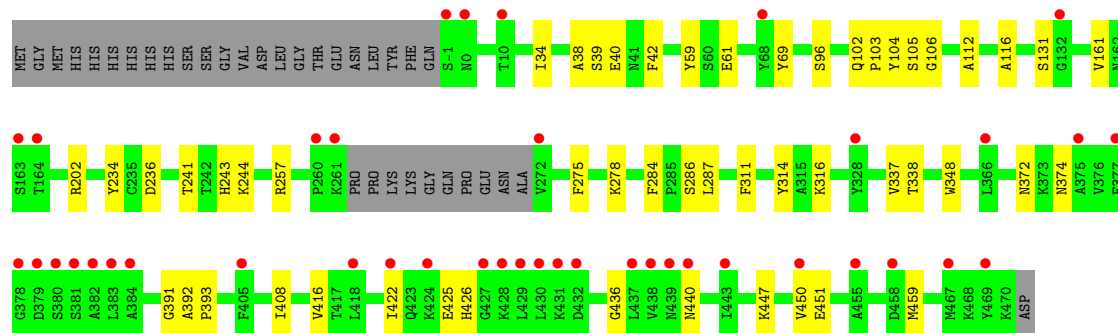
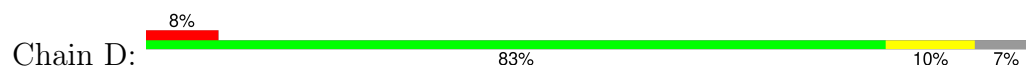


- Molecule 1: Serine hydroxymethyltransferase

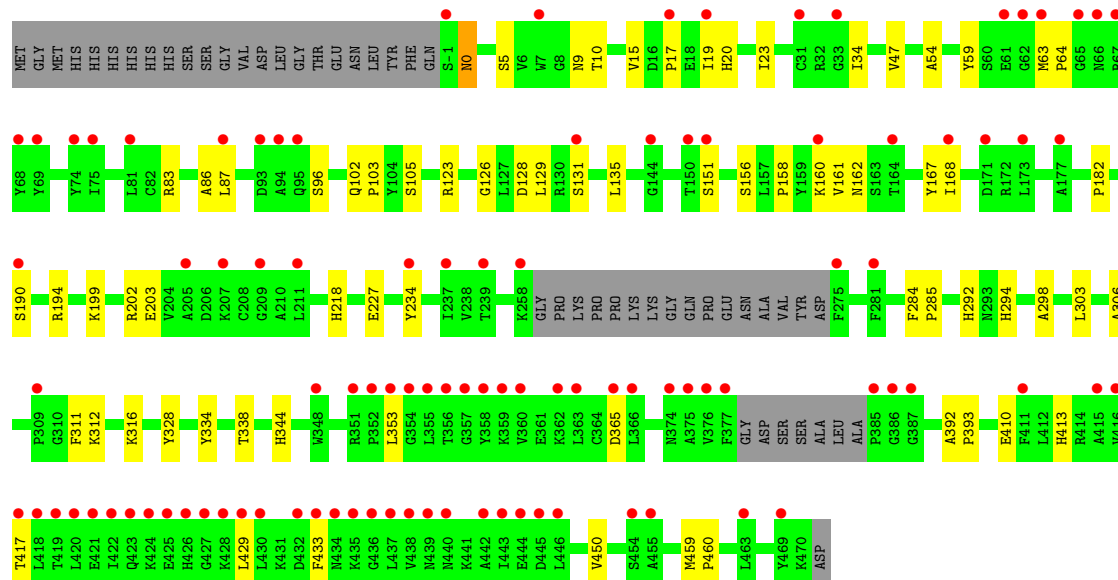
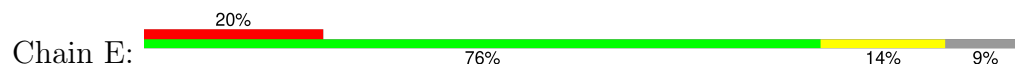




• Molecule 1: Serine hydroxymethyltransferase



• Molecule 1: Serine hydroxymethyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	175.31Å 175.31Å 186.33Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.86 – 2.25 48.86 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.86-2.25) 99.3 (48.86-2.25)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.01Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.265 , 0.311 0.266 , 0.310	Depositor DCC
R_{free} test set	10690 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.420	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	17890	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 28.15 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.9397e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PLG

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	0.35	0/3530	0.52	0/4799
1	B	0.33	0/3582	0.50	0/4866
1	C	0.32	0/3575	0.48	0/4853
1	D	0.32	0/3559	0.47	0/4830
1	E	0.30	0/3326	0.49	0/4530
All	All	0.32	0/17572	0.49	0/23878

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	3450	0	3204	33	0
1	B	3502	0	3325	22	0
1	C	3496	0	3328	29	0
1	D	3481	0	3323	30	0
1	E	3253	0	2939	40	0
2	A	20	0	12	1	0
2	B	20	0	12	1	0
2	C	20	0	12	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	20	0	12	2	0
2	E	20	0	12	0	0
3	A	158	0	0	0	0
3	B	129	0	0	1	0
3	C	141	0	0	4	0
3	D	126	0	0	0	0
3	E	54	0	0	0	0
All	All	17890	0	16179	150	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 (150) 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:C:34:ILE:HA	1:C:395:MET:HE2	1.54	0.87
1:A:285:PRO:HB2	1:B:135:LEU:HD21	1.61	0.81
1:A:178:LEU:HD11	1:A:207:LYS:HE3	1.67	0.77
1:A:58:LYS:HG2	1:A:75:ILE:HG13	1.74	0.70
1:E:47:VAL:HG13	1:E:298:ALA:HB1	1.78	0.66
1:A:61:GLU:HB3	1:A:284:PHE:CZ	2.31	0.66
1:D:131:SER:HB3	1:D:161:VAL:HG13	1.78	0.65
1:E:128:ASP:HB2	1:E:158:PRO:HB2	1.78	0.65
1:D:42:PHE:CD1	1:D:459:MET:HE1	2.31	0.65
1:A:42:PHE:CD1	1:A:459:MET:HE1	2.33	0.64
1:E:429:LEU:O	1:E:433:PHE:N	2.28	0.64
1:E:413:HIS:O	1:E:417:THR:OG1	2.14	0.62
1:A:34:ILE:HG13	1:A:450:VAL:HG13	1.81	0.61
1:D:34:ILE:HD13	1:D:408:ILE:HG12	1.83	0.61
1:B:137:HIS:HE1	1:B:215:ASP:OD2	1.84	0.61
1:D:34:ILE:HG13	1:D:450:VAL:HG13	1.82	0.59
1:A:174:GLU:OE1	1:A:200:ARG:NH2	2.34	0.59
1:E:131:SER:HG	1:E:161:VAL:HG22	1.68	0.58
1:C:131:SER:HB3	1:C:161:VAL:HG13	1.85	0.58
1:A:313:ALA:HA	1:A:316:LYS:HE2	1.86	0.57
1:A:40:GLU:OE2	1:B:58:LYS:HD2	2.05	0.56
1:E:128:ASP:OD2	1:E:129:LEU:N	2.39	0.56
1:E:328:TYR:OH	1:E:410:GLU:OE1	2.24	0.56
1:D:236:ASP:HA	1:D:257:ARG:HD2	1.88	0.55
1:E:292:HIS:HB3	1:E:294:HIS:CE1	2.41	0.55
1:D:337:VAL:HG12	1:D:338:THR:HG23	1.89	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:241:THR:HB	1:D:243:HIS:CE1	2.42	0.54
1:C:328:TYR:OH	1:C:410:GLU:OE1	2.24	0.53
1:B:348:TRP:CH2	1:B:416:VAL:HG21	2.44	0.53
1:E:17:PRO:HA	1:E:20:HIS:HB3	1.91	0.53
1:C:34:ILE:HG13	1:C:450:VAL:HG13	1.91	0.53
1:D:348:TRP:CH2	1:D:416:VAL:HG21	2.45	0.52
1:C:35:GLU:HG3	1:C:42:PHE:HZ	1.75	0.51
1:E:10:THR:HB	1:E:15:VAL:HG23	1.92	0.51
1:D:425:GLU:OE1	1:D:426:HIS:NE2	2.43	0.51
1:E:131:SER:HG	1:E:160:LYS:HZ2	1.56	0.51
1:E:202:ARG:HB2	1:E:234:TYR:HB3	1.93	0.50
1:E:128:ASP:OD1	1:E:160:LYS:HD3	2.11	0.50
1:C:58:LYS:HD2	1:D:40:GLU:OE2	2.11	0.50
1:E:131:SER:HB3	1:E:161:VAL:HG13	1.94	0.50
1:A:225:ALA:O	1:A:312:LYS:NZ	2.44	0.50
1:E:63:MET:HG3	1:E:64:PRO:HD2	1.93	0.50
1:B:168:ILE:HG13	1:B:192:TYR:CZ	2.47	0.50
1:E:151:SER:HB3	1:E:156:SER:OG	2.13	0.49
1:A:313:ALA:HA	1:A:316:LYS:CE	2.42	0.49
1:C:149:THR:HG23	1:D:286:SER:O	2.12	0.49
1:B:129:LEU:HD11	1:B:135:LEU:HA	1.94	0.49
1:B:42:PHE:CD1	1:B:459:MET:HE1	2.48	0.49
1:E:0:ASN:N	1:E:0:ASN:OD1	2.47	0.48
1:E:227:GLU:CD	1:E:316:LYS:HG2	2.39	0.48
2:D:501:PLG:O3	2:D:501:PLG:N	2.45	0.48
1:B:292:HIS:HB3	1:B:294:HIS:CE1	2.48	0.48
1:A:132:GLY:HA2	1:A:191:ALA:HB3	1.95	0.47
1:E:162:ASN:N	1:E:167:TYR:O	2.40	0.47
1:E:161:VAL:HA	1:E:168:ILE:HA	1.97	0.47
1:C:241:THR:HB	1:C:243:HIS:CE1	2.49	0.47
1:E:306:ALA:HA	1:E:311:PHE:CG	2.50	0.47
1:E:34:ILE:HG13	1:E:450:VAL:HG13	1.97	0.47
1:E:83:ARG:O	1:E:87:LEU:HD22	2.15	0.47
1:D:447:LYS:O	1:D:451:GLU:HG3	2.14	0.46
1:E:284:PHE:CD1	1:E:285:PRO:HA	2.50	0.46
1:D:422:ILE:HD11	1:D:440:ASN:ND2	2.30	0.46
1:B:344:HIS:HB2	1:B:393:PRO:HD3	1.98	0.46
1:D:39:SER:HB2	1:D:244:LYS:HE3	1.98	0.46
1:C:348:TRP:HE3	1:C:390:ILE:HD11	1.81	0.45
2:B:501:PLG:O3	2:B:501:PLG:N	2.49	0.45
1:A:123:ARG:HA	1:A:155:GLU:O	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:36:LEU:HB3	1:C:390:ILE:HG23	1.98	0.45
1:D:392:ALA:N	1:D:393:PRO:CD	2.79	0.45
1:E:334:TYR:HE1	1:E:353:LEU:HD21	1.81	0.45
1:B:61:GLU:HB3	1:B:284:PHE:CZ	2.52	0.45
3:C:610:HOH:O	2:D:501:PLG:H4A1	2.17	0.45
1:D:112:ALA:HB2	1:D:287:LEU:HD23	1.99	0.45
1:B:190:SER:HA	1:B:218:HIS:CD2	2.52	0.45
1:C:47:VAL:HG13	1:C:298:ALA:HB1	1.99	0.45
1:E:5:SER:HA	1:E:9:ASN:HB2	1.99	0.44
1:A:292:HIS:HB3	1:A:294:HIS:CE1	2.52	0.44
1:B:36:LEU:HB3	1:B:390:ILE:HG23	1.99	0.44
1:C:400:LEU:HD23	1:C:404:ASP:HB3	1.99	0.44
1:E:218:HIS:HB3	1:E:344:HIS:CE1	2.51	0.44
1:A:64:PRO:HD2	1:A:83:ARG:HH12	1.81	0.44
1:A:436:GLY:O	1:A:440:ASN:HB2	2.18	0.44
1:E:459:MET:SD	1:E:460:PRO:HD2	2.57	0.44
1:C:126:GLY:O	1:C:158:PRO:HA	2.17	0.44
1:D:236:ASP:HB3	1:D:275:PHE:CE2	2.52	0.44
1:D:311:PHE:O	1:D:314:TYR:HB3	2.17	0.44
1:A:216:MET:HG2	1:A:240:THR:HB	2.00	0.44
1:C:19:ILE:HD13	1:C:19:ILE:HA	1.89	0.44
1:C:284:PHE:CD1	1:C:285:PRO:HA	2.53	0.44
1:A:337:VAL:HG12	1:A:338:THR:HG23	1.98	0.44
1:D:61:GLU:HB3	1:D:284:PHE:CZ	2.53	0.44
1:D:316:LYS:HB3	1:D:316:LYS:HE3	1.74	0.44
1:D:106:GLY:HA3	1:D:241:THR:HG22	1.99	0.44
1:A:392:ALA:N	1:A:393:PRO:HD3	2.33	0.43
1:B:336:LEU:HD13	1:B:346:VAL:HB	1.99	0.43
1:B:373:LYS:HG3	1:B:385:PRO:HG2	2.00	0.43
1:E:19:ILE:O	1:E:23:ILE:HG13	2.19	0.43
1:A:241:THR:HG21	1:A:244:LYS:NZ	2.33	0.43
2:A:501:PLG:N	2:A:501:PLG:O3	2.51	0.43
1:C:1:MET:HE3	1:D:314:TYR:HA	2.00	0.43
1:A:216:MET:HG3	1:A:220:SER:HB3	2.00	0.43
1:D:38:ALA:HA	1:D:391:GLY:HA3	2.01	0.43
1:E:190:SER:HA	1:E:218:HIS:CD2	2.53	0.43
1:A:241:THR:HB	1:A:243:HIS:CE1	2.54	0.43
1:C:168:ILE:HG13	1:C:192:TYR:CZ	2.54	0.43
1:E:131:SER:OG	1:E:161:VAL:HG22	2.18	0.43
1:C:58:LYS:HG2	1:C:75:ILE:HG13	2.01	0.43
1:C:355:LEU:HD21	1:C:420:LEU:HD23	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:HIS:CE1	1:A:136:THR:HG23	2.54	0.42
1:E:123:ARG:HB2	1:E:182:PRO:HA	2.00	0.42
1:A:392:ALA:N	1:A:393:PRO:CD	2.83	0.42
1:D:422:ILE:HD12	1:D:436:GLY:HA3	2.01	0.42
1:B:459:MET:N	3:B:603:HOH:O	2.34	0.42
1:A:0:ASN:HD22	1:A:0:ASN:N	2.17	0.42
1:B:274:ASP:OD2	1:B:278:LYS:NZ	2.52	0.42
1:C:392:ALA:N	1:C:393:PRO:CD	2.83	0.42
1:E:129:LEU:HD21	1:E:135:LEU:HD23	2.02	0.42
1:E:199:LYS:O	1:E:203:GLU:HG3	2.20	0.42
3:C:738:HOH:O	1:D:287:LEU:HD11	2.19	0.42
1:A:107:SER:HB2	1:A:108:PRO:HD3	2.01	0.42
1:A:227:GLU:CD	1:A:316:LYS:HG2	2.44	0.42
1:E:86:ALA:HA	1:E:303:LEU:HD22	2.02	0.42
1:A:131:SER:OG	1:A:161:VAL:HG22	2.20	0.42
1:C:150:THR:O	1:C:154:PHE:HB2	2.20	0.42
1:C:295:GLN:NE2	3:C:603:HOH:O	2.34	0.42
1:E:126:GLY:O	1:E:158:PRO:HA	2.19	0.41
1:B:347:LEU:HD12	1:B:347:LEU:HA	1.89	0.41
1:A:104:TYR:O	1:B:292:HIS:HE1	2.04	0.41
1:B:311:PHE:O	1:B:314:TYR:HB3	2.20	0.41
1:E:160:LYS:HA	1:E:160:LYS:HD2	1.94	0.41
1:B:24:GLU:HA	1:B:27:LYS:HD3	2.02	0.41
1:D:372:ASN:HD21	1:D:374:ASN:HD22	1.67	0.41
1:A:58:LYS:HE2	1:A:72:ASN:OD1	2.21	0.41
1:C:243:HIS:ND1	1:C:250:ARG:HA	2.36	0.41
1:E:392:ALA:N	1:E:393:PRO:CD	2.84	0.41
1:A:286:SER:O	1:B:149:THR:HG23	2.21	0.41
1:C:337:VAL:HG21	1:C:376:VAL:HG23	2.02	0.41
1:D:102:GLN:N	1:D:103:PRO:CD	2.84	0.41
1:A:10:THR:HG22	1:A:15:VAL:HG23	2.03	0.40
1:A:400:LEU:HD23	1:A:404:ASP:HB3	2.02	0.40
1:C:38:ALA:HA	1:C:391:GLY:HA3	2.03	0.40
1:C:113:ALA:O	1:C:117:VAL:HG22	2.20	0.40
1:D:116:ALA:O	1:D:278:LYS:HD2	2.20	0.40
1:E:102:GLN:N	1:E:103:PRO:CD	2.85	0.40
1:B:102:GLN:N	1:B:103:PRO:CD	2.84	0.40
1:D:202:ARG:HB2	1:D:234:TYR:HB3	2.02	0.40
1:C:59:TYR:HA	3:C:632:HOH:O	2.21	0.40
1:C:292:HIS:HE1	1:D:104:TYR:O	2.05	0.40
1:E:194:ARG:NH2	1:E:338:THR:O	2.50	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:292:HIS:HB3	1:C:294:HIS:CE1	2.56	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	463/496 (93%)	442 (96%)	20 (4%)	1 (0%)	43	50
1	B	461/496 (93%)	446 (97%)	15 (3%)	0	100	100
1	C	462/496 (93%)	446 (96%)	15 (3%)	1 (0%)	43	50
1	D	458/496 (92%)	445 (97%)	13 (3%)	0	100	100
1	E	443/496 (89%)	420 (95%)	22 (5%)	1 (0%)	43	50
All	All	2287/2480 (92%)	2199 (96%)	85 (4%)	3 (0%)	48	56

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	430	LEU
1	E	54	ALA
1	C	217	ALA

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	334/408 (82%)	333 (100%)	1 (0%)	86	90
1	B	351/408 (86%)	348 (99%)	3 (1%)	70	79
1	C	348/408 (85%)	345 (99%)	3 (1%)	70	79
1	D	350/408 (86%)	346 (99%)	4 (1%)	65	75
1	E	305/408 (75%)	299 (98%)	6 (2%)	48	59
All	All	1688/2040 (83%)	1671 (99%)	17 (1%)	68	77

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

Mol	Chain	Res	Type
1	A	105	SER
1	B	59	TYR
1	B	105	SER
1	B	258	LYS
1	C	43	THR
1	C	59	TYR
1	C	105	SER
1	D	59	TYR
1	D	69	TYR
1	D	96	SER
1	D	105	SER
1	E	0	ASN
1	E	59	TYR
1	E	96	SER
1	E	105	SER
1	E	312	LYS
1	E	365	ASP

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	0	ASN
1	A	407	GLN
1	B	88	GLN
1	B	137	HIS
1	B	293	ASN
1	C	9	ASN
1	C	20	HIS
1	C	293	ASN
1	D	374	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	407	GLN
1	E	88	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

5 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	PLG	A	501	-	20,20,20	1.38	2 (10%)	26,28,28	1.26	1 (3%)
2	PLG	B	501	-	20,20,20	1.32	3 (15%)	26,28,28	1.39	5 (19%)
2	PLG	E	501	-	20,20,20	1.33	2 (10%)	26,28,28	1.25	3 (11%)
2	PLG	D	501	-	20,20,20	1.42	2 (10%)	26,28,28	1.45	6 (23%)
2	PLG	C	501	-	20,20,20	1.43	4 (20%)	26,28,28	2.00	7 (26%)

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	-	-	5/12/12/12	0/1/1/1
2	PLG	B	501	-	-	5/12/12/12	0/1/1/1
2	PLG	E	501	-	-	6/12/12/12	0/1/1/1
2	PLG	D	501	-	-	4/12/12/12	0/1/1/1
2	PLG	C	501	-	-	7/12/12/12	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	PLG	C5-C4	-3.44	1.35	1.40
2	A	501	PLG	C5-C4	-2.99	1.36	1.40
2	C	501	PLG	C5-C4	-2.65	1.36	1.40
2	C	501	PLG	C3-C2	-2.60	1.38	1.41
2	B	501	PLG	C3-C2	-2.40	1.38	1.41
2	E	501	PLG	C5-C4	-2.29	1.37	1.40
2	B	501	PLG	C5-C4	-2.26	1.37	1.40
2	C	501	PLG	C2A-C2	2.22	1.53	1.50
2	E	501	PLG	C3-C2	-2.17	1.38	1.41
2	D	501	PLG	C2A-C2	2.17	1.53	1.50
2	C	501	PLG	C3-C4	-2.14	1.36	1.40
2	A	501	PLG	C3-C2	-2.12	1.38	1.41
2	B	501	PLG	C2A-C2	2.09	1.53	1.50

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	PLG	C4A-N-CA	-7.25	104.28	112.72
2	E	501	PLG	C4A-N-CA	-3.14	109.06	112.72
2	C	501	PLG	C5-C6-N1	-3.02	118.92	123.83
2	D	501	PLG	C6-C5-C4	2.99	120.32	118.06
2	B	501	PLG	OP4-C5A-C5	2.95	114.89	109.36
2	D	501	PLG	OP4-P-OP1	2.84	114.12	106.44
2	A	501	PLG	OXT-C-CA	2.74	123.23	112.81
2	B	501	PLG	C6-C5-C4	2.73	120.12	118.06
2	D	501	PLG	OP4-C5A-C5	2.53	114.09	109.36
2	B	501	PLG	OP4-P-OP1	2.49	113.17	106.44
2	C	501	PLG	OP4-P-OP1	2.48	113.14	106.44
2	B	501	PLG	C5-C6-N1	-2.45	119.85	123.83
2	B	501	PLG	OXT-C-CA	2.38	121.84	112.81
2	C	501	PLG	OP4-C5A-C5	2.34	113.75	109.36
2	E	501	PLG	OP4-C5A-C5	2.28	113.62	109.36
2	C	501	PLG	C6-C5-C4	2.15	119.68	118.06

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	PLG	C4A-C4-C3	2.11	122.79	119.98
2	D	501	PLG	OXT-C-CA	2.11	120.83	112.81
2	E	501	PLG	C5-C6-N1	-2.11	120.41	123.83
2	C	501	PLG	C4A-C4-C5	-2.08	117.49	119.75
2	C	501	PLG	C3-C4-C5	2.01	120.55	118.73
2	D	501	PLG	C5-C6-N1	-2.01	120.56	123.83

There are no chirality outliers.

All (27) torsion outliers are listed below:

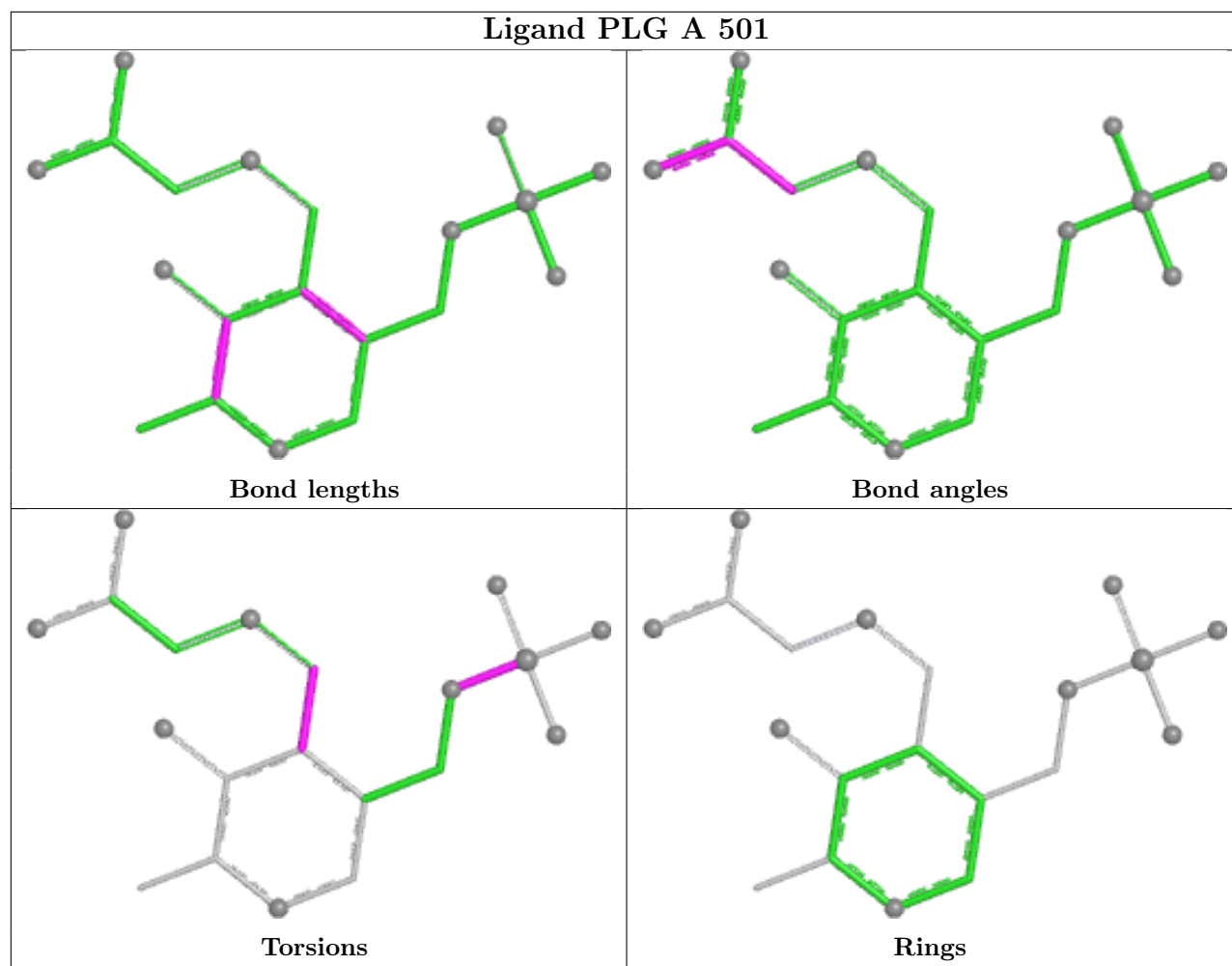
Mol	Chain	Res	Type	Atoms
2	A	501	PLG	C3-C4-C4A-N
2	A	501	PLG	C5-C4-C4A-N
2	A	501	PLG	C5A-OP4-P-OP2
2	A	501	PLG	C5A-OP4-P-OP3
2	B	501	PLG	C5-C4-C4A-N
2	B	501	PLG	C5A-OP4-P-OP1
2	B	501	PLG	C5A-OP4-P-OP3
2	C	501	PLG	C5A-OP4-P-OP1
2	D	501	PLG	C5-C4-C4A-N
2	E	501	PLG	C5A-OP4-P-OP2
2	E	501	PLG	C5A-OP4-P-OP3
2	C	501	PLG	C5-C4-C4A-N
2	E	501	PLG	C5-C4-C4A-N
2	B	501	PLG	C3-C4-C4A-N
2	C	501	PLG	C4-C4A-N-CA
2	E	501	PLG	C4-C4A-N-CA
2	A	501	PLG	C5A-OP4-P-OP1
2	D	501	PLG	C5A-OP4-P-OP1
2	E	501	PLG	C5A-OP4-P-OP1
2	C	501	PLG	C3-C4-C4A-N
2	E	501	PLG	C3-C4-C4A-N
2	B	501	PLG	C5A-OP4-P-OP2
2	C	501	PLG	C5A-OP4-P-OP3
2	D	501	PLG	C5A-OP4-P-OP3
2	D	501	PLG	C3-C4-C4A-N
2	C	501	PLG	OXT-C-CA-N
2	C	501	PLG	O-C-CA-N

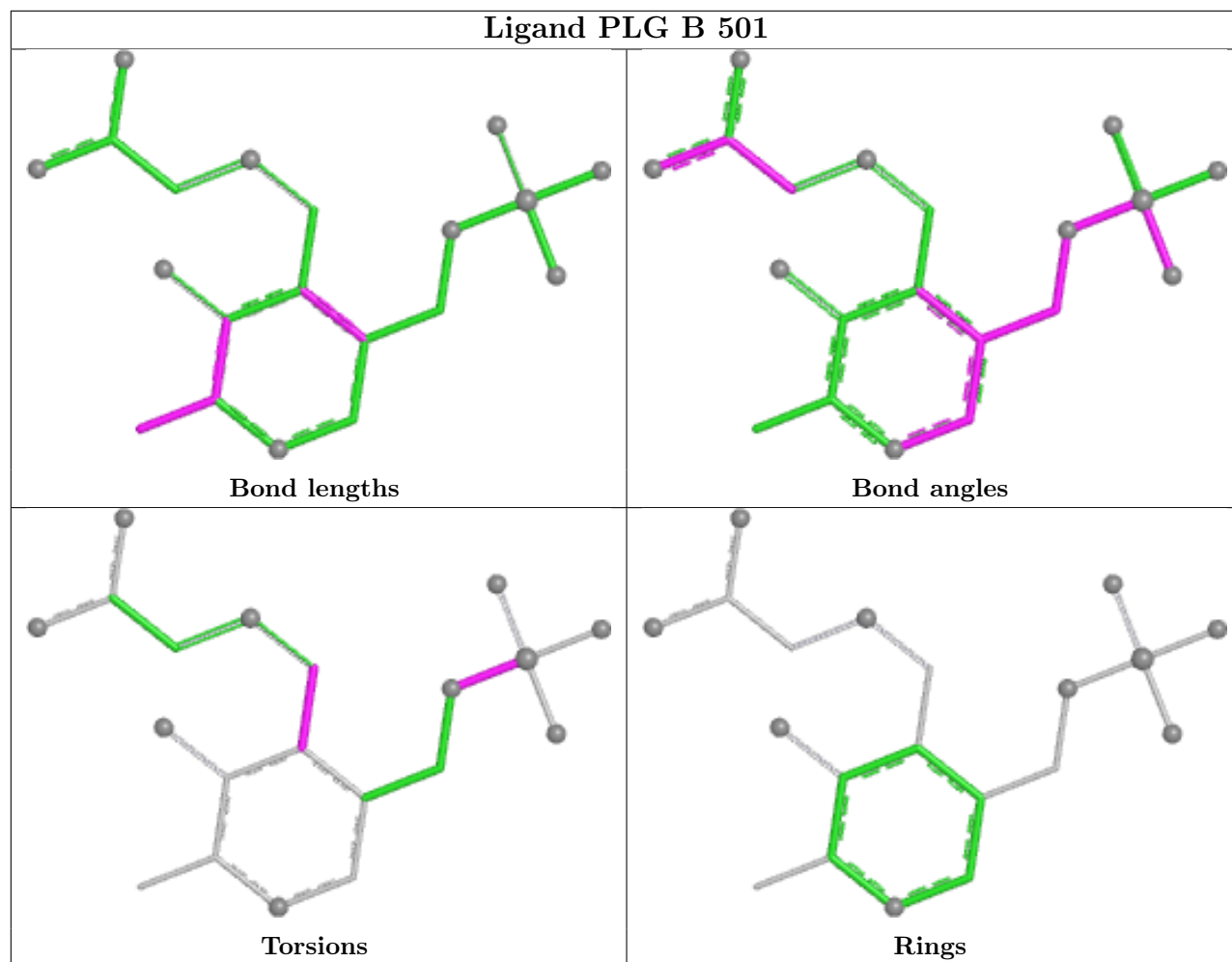
There are no ring outliers.

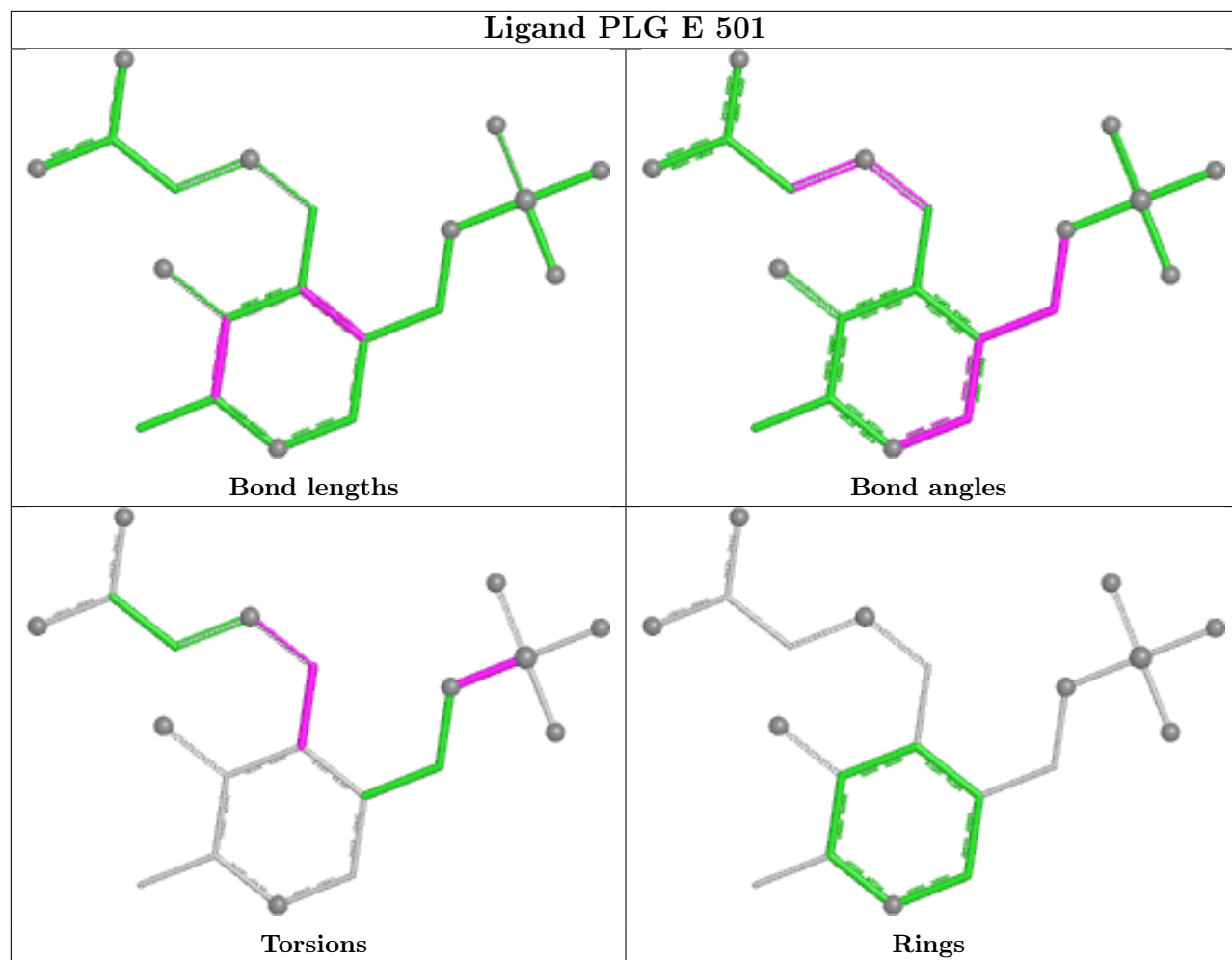
3 monomers are involved in 4 short contacts:

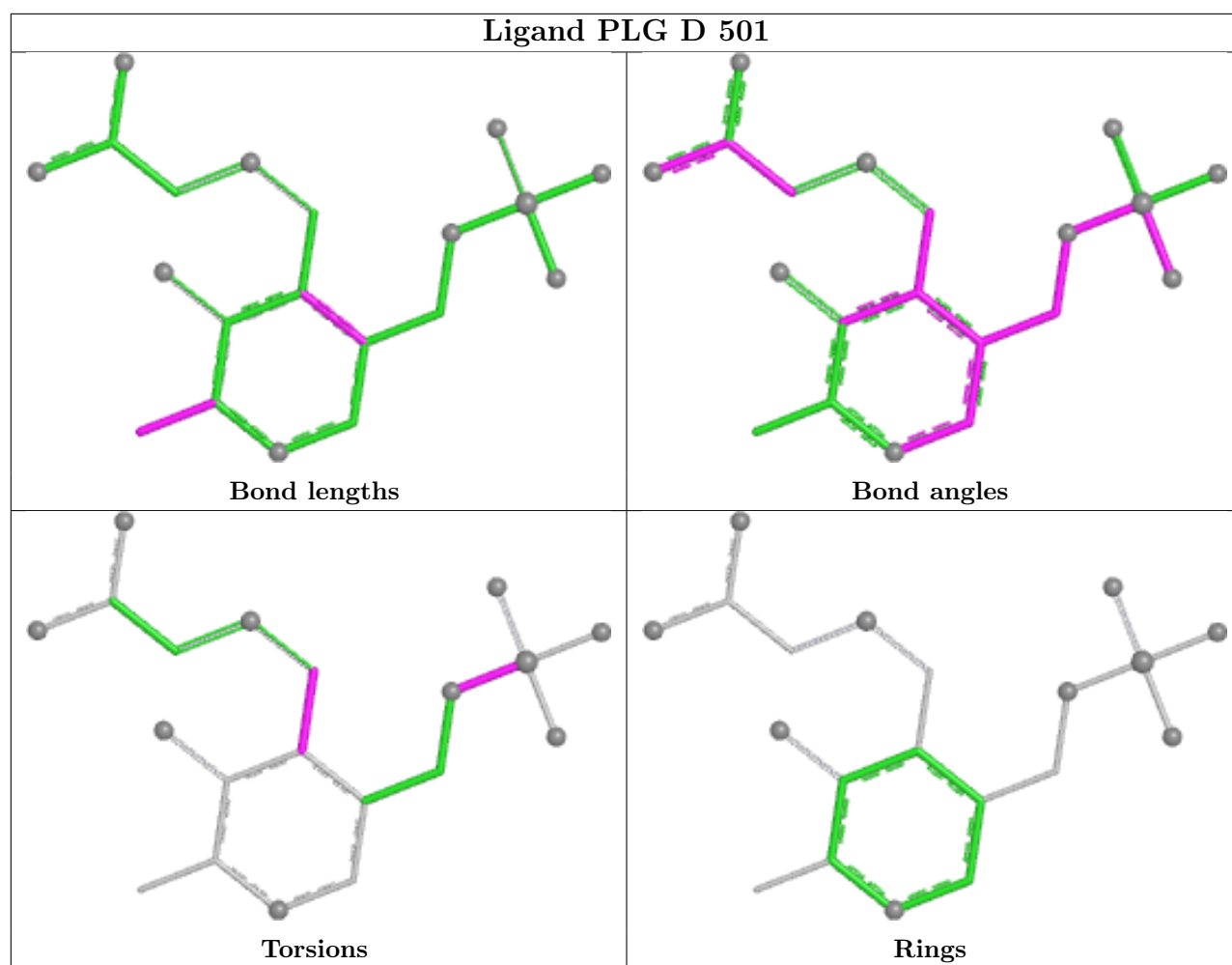
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	PLG	1	0
2	B	501	PLG	1	0
2	D	501	PLG	2	0

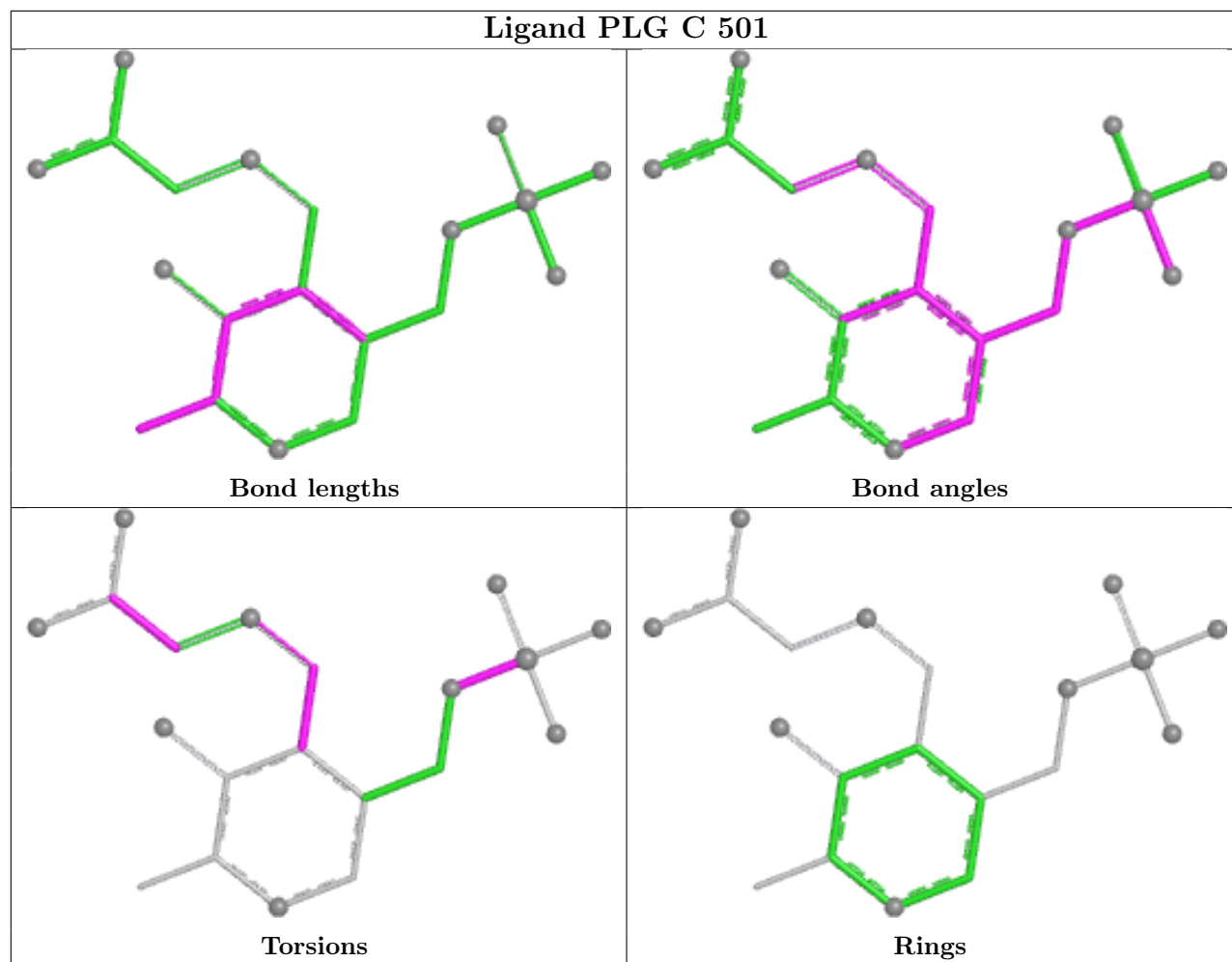
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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











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	467/496 (94%)	0.86	59 (12%) 8 7	20, 30, 65, 102	0
1	B	467/496 (94%)	0.77	36 (7%) 19 17	20, 33, 56, 76	0
1	C	466/496 (93%)	0.79	48 (10%) 12 11	22, 33, 67, 89	0
1	D	462/496 (93%)	0.77	41 (8%) 15 14	21, 34, 63, 89	0
1	E	449/496 (90%)	1.37	100 (22%) 2 2	30, 45, 78, 112	0
All	All	2311/2480 (93%)	0.91	284 (12%) 8 7	20, 35, 67, 112	0

All (284) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	433	PHE	6.9
1	E	385	PRO	6.0
1	A	437	LEU	5.5
1	A	429	LEU	5.4
1	E	377	PHE	5.4
1	A	432	ASP	5.3
1	A	377	PHE	5.3
1	B	163	SER	5.3
1	B	439	ASN	5.3
1	A	379	ASP	5.2
1	C	268	PRO	5.1
1	E	426	HIS	5.1
1	C	437	LEU	4.8
1	E	439	ASN	4.8
1	A	442	ALA	4.7
1	C	430	LEU	4.7
1	A	436	GLY	4.7
1	A	378	GLY	4.6
1	E	427	GLY	4.6
1	D	377	PHE	4.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	434	ASN	4.5
1	E	355	LEU	4.4
1	E	433	PHE	4.4
1	C	378	GLY	4.3
1	D	163	SER	4.3
1	A	435	LYS	4.3
1	E	66	ASN	4.2
1	E	281	PHE	4.2
1	A	355	LEU	4.2
1	E	422	ILE	4.2
1	D	68	TYR	4.1
1	C	383	LEU	4.1
1	C	439	ASN	4.1
1	D	430	LEU	4.1
1	E	436	GLY	4.1
1	E	275	PHE	4.0
1	E	375	ALA	4.0
1	A	430	LEU	3.9
1	E	420	LEU	3.9
1	B	266	GLY	3.9
1	C	432	ASP	3.9
1	C	381	SER	3.9
1	A	443	ILE	3.8
1	E	437	LEU	3.8
1	E	432	ASP	3.8
1	C	163	SER	3.8
1	E	376	VAL	3.8
1	E	440	ASN	3.8
1	C	429	LEU	3.8
1	E	363	LEU	3.8
1	B	380	SER	3.8
1	A	129	LEU	3.8
1	A	354	GLY	3.8
1	B	164	THR	3.7
1	E	356	THR	3.7
1	B	268	PRO	3.7
1	A	188	GLY	3.7
1	A	366	LEU	3.7
1	C	425	GLU	3.7
1	E	445	ASP	3.7
1	C	384	ALA	3.7
1	B	381	SER	3.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	438	VAL	3.6
1	E	430	LEU	3.6
1	A	431	LYS	3.6
1	D	380	SER	3.6
1	E	68	TYR	3.6
1	E	438	VAL	3.6
1	D	427	GLY	3.5
1	C	-1	SER	3.5
1	A	419	THR	3.5
1	C	427	GLY	3.5
1	A	-1	SER	3.4
1	D	405	PHE	3.4
1	E	424	LYS	3.4
1	E	360	VAL	3.4
1	D	0	ASN	3.4
1	E	353	LEU	3.4
1	E	429	LEU	3.4
1	C	438	VAL	3.4
1	D	450	VAL	3.3
1	B	331	GLY	3.3
1	B	384	ALA	3.3
1	D	-1	SER	3.3
1	A	439	ASN	3.3
1	E	444	GLU	3.3
1	C	431	LYS	3.3
1	E	95	GLN	3.3
1	E	415	ALA	3.2
1	E	386	GLY	3.2
1	E	418	LEU	3.2
1	E	168	ILE	3.2
1	E	428	LYS	3.2
1	A	352	PRO	3.2
1	E	423	GLN	3.2
1	C	68	TYR	3.2
1	D	261	LYS	3.2
1	A	422	ILE	3.2
1	A	427	GLY	3.1
1	C	382	ALA	3.1
1	D	328	TYR	3.1
1	E	-1	SER	3.1
1	E	358	TYR	3.1
1	D	438	VAL	3.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	385	PRO	3.1
1	A	418	LEU	3.1
1	B	430	LEU	3.1
1	E	87	LEU	3.1
1	B	379	ASP	3.1
1	C	442	ALA	3.1
1	A	353	LEU	3.0
1	B	165	THR	3.0
1	C	377	PHE	3.0
1	E	411	PHE	3.0
1	C	423	GLN	3.0
1	C	433	PHE	2.9
1	A	358	TYR	2.9
1	D	379	ASP	2.9
1	A	348	TRP	2.9
1	D	429	LEU	2.9
1	E	33	GLY	2.9
1	A	416	VAL	2.9
1	C	270	ASN	2.9
1	E	94	ALA	2.8
1	C	428	LYS	2.8
1	E	258	LYS	2.8
1	B	0	ASN	2.8
1	D	439	ASN	2.8
1	A	266	GLY	2.8
1	D	132	GLY	2.8
1	E	209	GLY	2.8
1	A	347	LEU	2.8
1	A	438	VAL	2.8
1	E	207	LYS	2.8
1	D	458	ASP	2.8
1	B	-1	SER	2.8
1	E	74	TYR	2.8
1	B	263	PRO	2.8
1	C	418	LEU	2.8
1	D	383	LEU	2.8
1	A	428	LYS	2.8
1	E	435	LYS	2.8
1	A	28	ARG	2.8
1	C	426	HIS	2.7
1	A	426	HIS	2.7
1	A	434	ASN	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	424	LYS	2.7
1	A	69	TYR	2.7
1	C	0	ASN	2.7
1	D	381	SER	2.7
1	D	467	MET	2.7
1	E	366	LEU	2.7
1	D	384	ALA	2.7
1	A	376	VAL	2.7
1	D	469	TYR	2.6
1	A	0	ASN	2.6
1	E	171	ASP	2.6
1	E	425	GLU	2.6
1	A	415	ALA	2.6
1	B	440	ASN	2.6
1	D	432	ASP	2.6
1	A	142	SER	2.6
1	C	454	SER	2.6
1	E	443	ILE	2.6
1	D	428	LYS	2.6
1	E	362	LYS	2.6
1	B	69	TYR	2.6
1	C	69	TYR	2.6
1	D	366	LEU	2.6
1	D	437	LEU	2.6
1	A	368	ASN	2.6
1	B	281	PHE	2.6
1	D	455	ALA	2.5
1	B	66	ASN	2.5
1	D	164	THR	2.5
1	E	160	LYS	2.5
1	D	260	PRO	2.5
1	E	61	GLU	2.5
1	A	375	ALA	2.5
1	E	177	ALA	2.5
1	E	205	ALA	2.5
1	A	333	GLY	2.5
1	E	421	GLU	2.5
1	D	424	LYS	2.5
1	E	93	ASP	2.5
1	E	65	GLY	2.5
1	E	354	GLY	2.5
1	E	69	TYR	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	274	ASP	2.4
1	B	171	ASP	2.4
1	C	455	ALA	2.4
1	A	356	THR	2.4
1	E	150	THR	2.4
1	C	443	ILE	2.4
1	C	379	ASP	2.4
1	D	382	ALA	2.4
1	D	10	THR	2.4
1	E	19	ILE	2.4
1	A	68	TYR	2.3
1	A	359	LYS	2.3
1	B	424	LYS	2.3
1	C	380	SER	2.3
1	E	63	MET	2.3
1	E	151	SER	2.3
1	C	440	ASN	2.3
1	C	271	ALA	2.3
1	E	442	ALA	2.3
1	B	419	THR	2.3
1	A	421	GLU	2.3
1	A	360	VAL	2.3
1	E	359	LYS	2.3
1	E	357	GLY	2.3
1	E	348	TRP	2.3
1	A	164	THR	2.3
1	C	441	LYS	2.3
1	C	470	LYS	2.3
1	C	65	GLY	2.2
1	A	425	GLU	2.2
1	E	164	THR	2.2
1	B	422	ILE	2.2
1	D	443	ILE	2.2
1	A	135	LEU	2.2
1	C	353	LEU	2.2
1	C	363	LEU	2.2
1	E	446	LEU	2.2
1	E	352	PRO	2.2
1	E	144	GLY	2.2
1	B	425	GLU	2.2
1	B	68	TYR	2.2
1	C	142	SER	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	446	LEU	2.2
1	E	416	VAL	2.2
1	D	378	GLY	2.2
1	E	62	GLY	2.2
1	E	190	SER	2.2
1	E	234	TYR	2.2
1	C	28	ARG	2.2
1	D	440	ASN	2.2
1	B	166	GLY	2.2
1	E	31	CYS	2.2
1	E	131	SER	2.2
1	E	454	SER	2.2
1	D	272	VAL	2.1
1	B	377	PHE	2.1
1	B	435	LYS	2.1
1	B	335	SER	2.1
1	E	239	THR	2.1
1	B	167	TYR	2.1
1	A	350	LEU	2.1
1	E	7	TRP	2.1
1	E	417	THR	2.1
1	D	422	ILE	2.1
1	E	81	LEU	2.1
1	E	365	ASP	2.1
1	E	374	ASN	2.1
1	E	469	TYR	2.1
1	B	426	HIS	2.1
1	A	268	PRO	2.1
1	E	67	ARG	2.1
1	A	448	ALA	2.1
1	C	307	ALA	2.1
1	D	431	LYS	2.1
1	E	455	ALA	2.1
1	C	467	MET	2.1
1	B	437	LEU	2.1
1	C	366	LEU	2.1
1	E	173	LEU	2.1
1	E	17	PRO	2.1
1	E	309	PRO	2.1
1	B	378	GLY	2.1
1	E	419	THR	2.0
1	D	375	ALA	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	412	LEU	2.0
1	D	418	LEU	2.0
1	E	211	LEU	2.0
1	A	147	ILE	2.0
1	E	75	ILE	2.0
1	B	385	PRO	2.0
1	C	358	TYR	2.0
1	E	351	ARG	2.0
1	E	387	GLY	2.0
1	E	463	LEU	2.0
1	B	351	ARG	2.0
1	E	237	ILE	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](#)

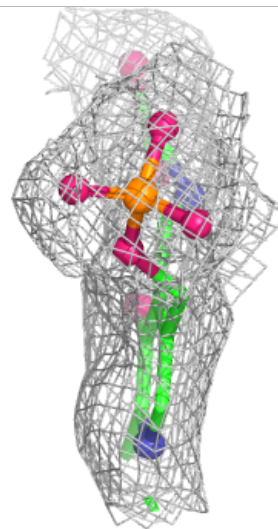
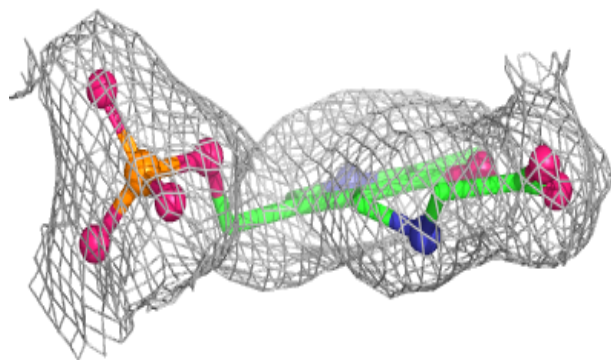
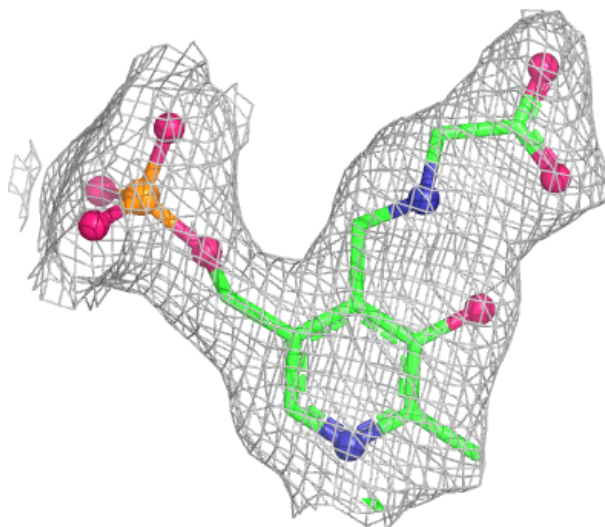
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PLG	E	501	20/20	0.93	0.10	30,35,38,38	0
2	PLG	D	501	20/20	0.95	0.10	21,27,32,33	0
2	PLG	B	501	20/20	0.95	0.10	23,26,29,32	0
2	PLG	A	501	20/20	0.96	0.08	20,26,31,32	0
2	PLG	C	501	20/20	0.97	0.07	24,28,30,31	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.

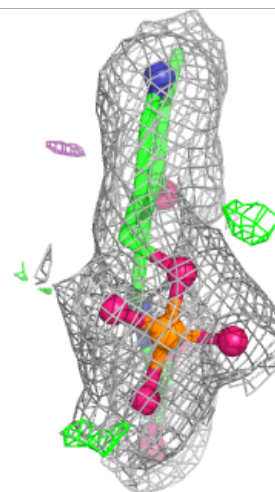
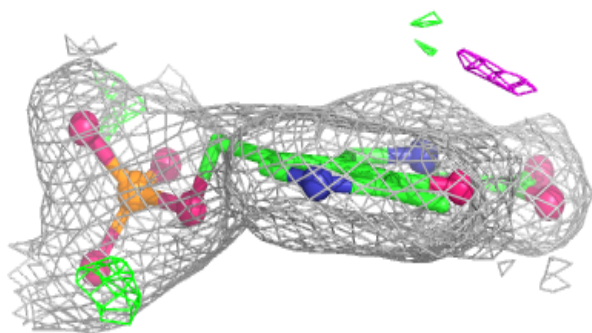
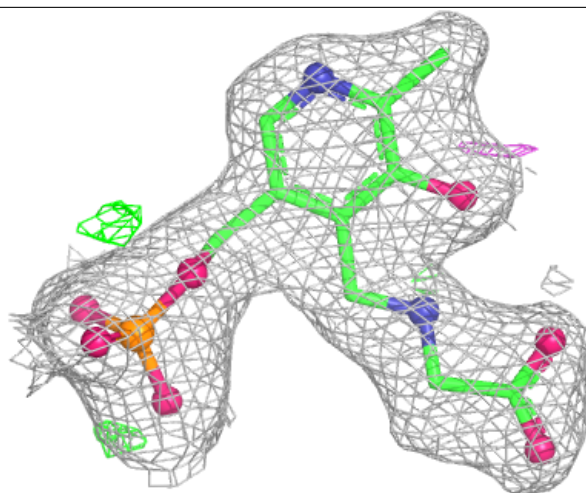
Electron density around PLG E 501:

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



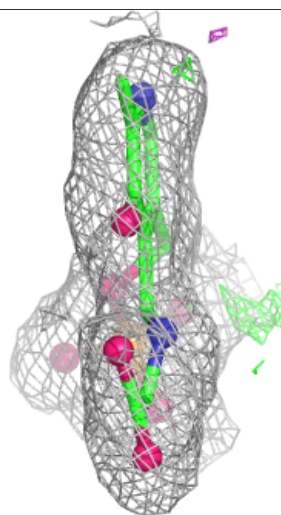
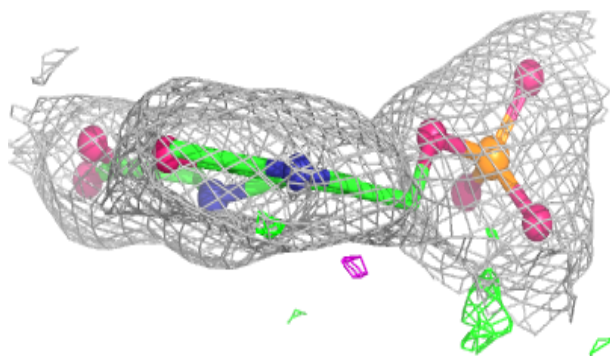
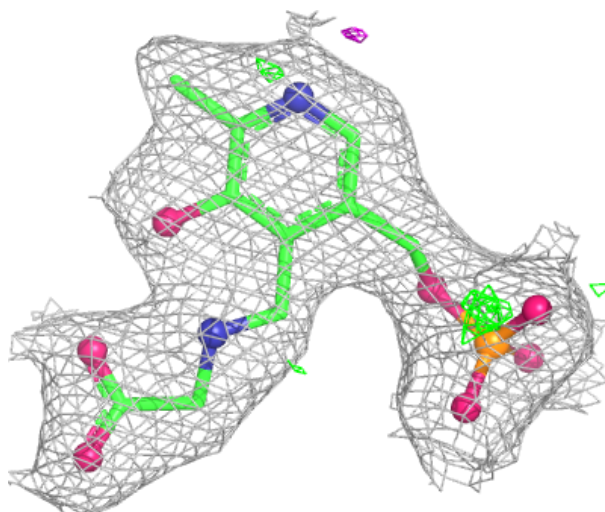
Electron density around PLG D 501:

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



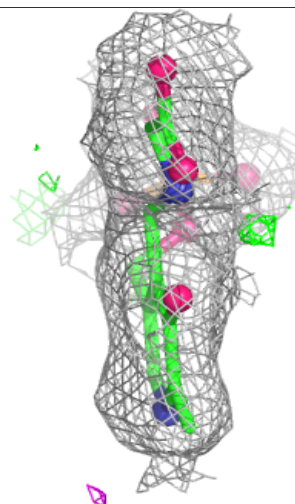
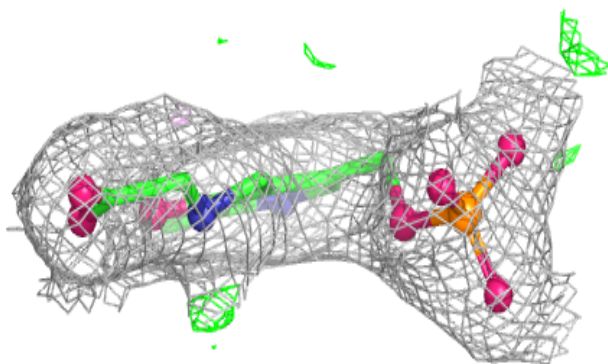
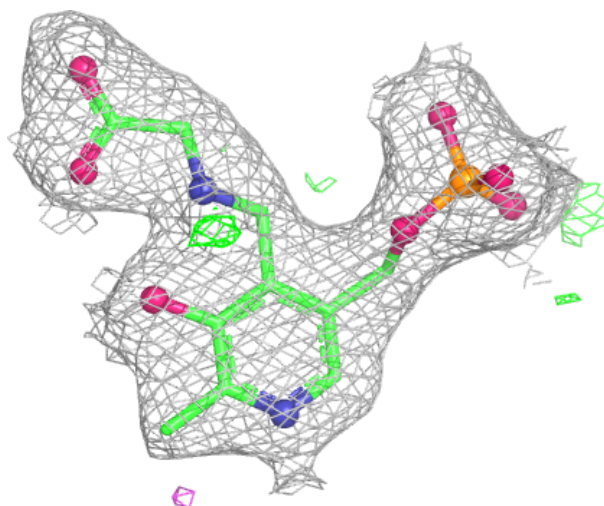
Electron density around PLG B 501:

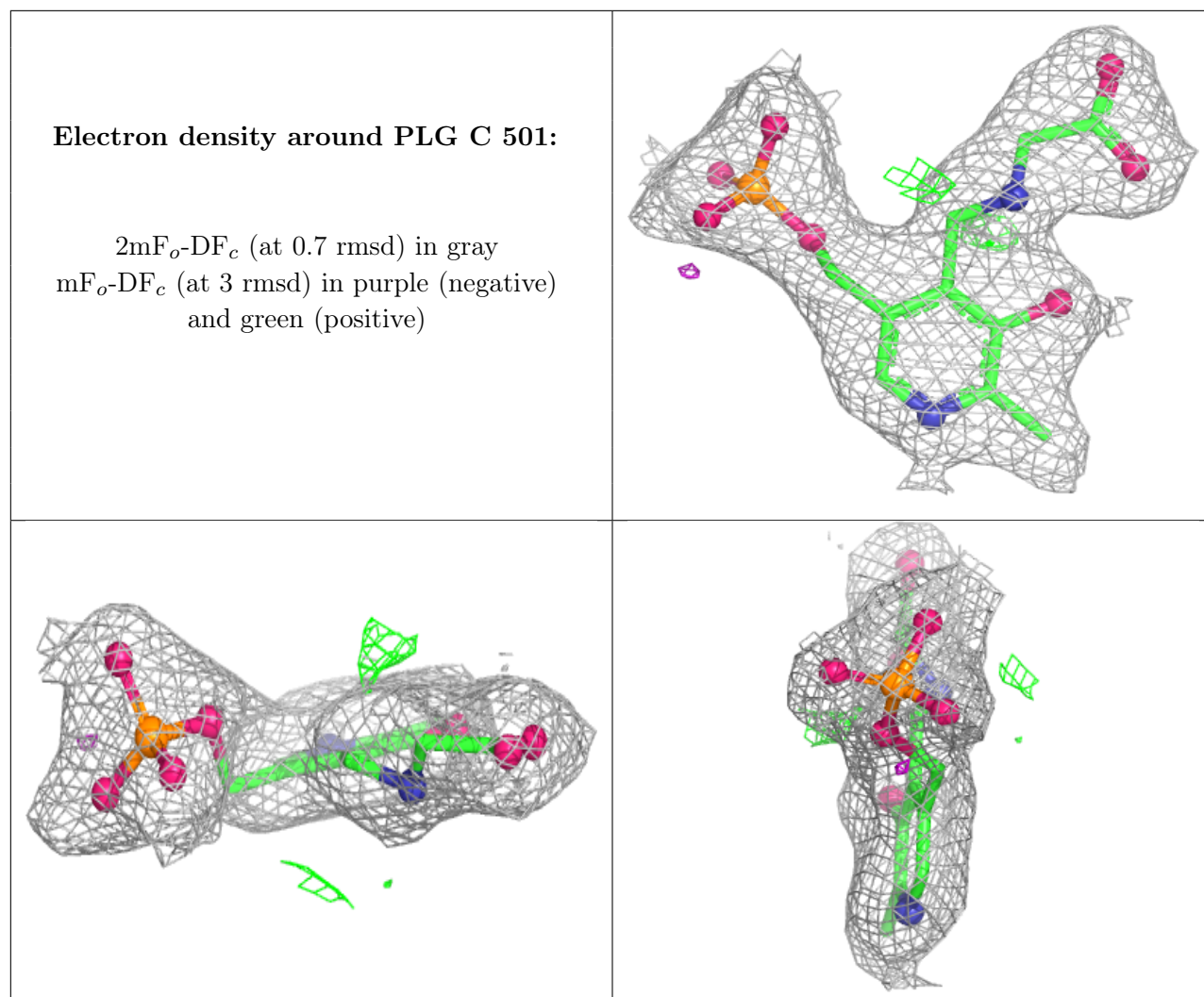
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





6.5 Other polymers ⓘ

There are no such residues in this entry.