



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 8, 2025 – 10:11 AM EDT

PDB ID : 9DLT / pdb_00009dlt
Title : STRUCTURE OF SERINE HYDROXYMETHYLTRANSFERASE 5 FROM
GLYCINE MAX CULTIVAR ESSEX COMPLEXED WITH PLP-GLYCINE
Authors : Beamer, L.J.; Owuocha, L.F.
Deposited on : 2024-09-11
Resolution : 1.90 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0rc1
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

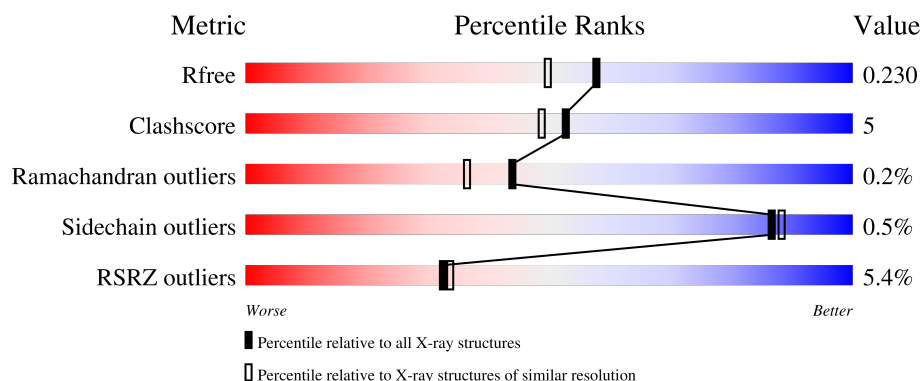
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	164625	7293 (1.90-1.90)
Clashscore	180529	8090 (1.90-1.90)
Ramachandran outliers	177936	8022 (1.90-1.90)
Sidechain outliers	177891	8022 (1.90-1.90)
RSRZ outliers	164620	7292 (1.90-1.90)

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>6%</div> <div> <div></div> <div>82%</div> <div>11%</div> <div>6%</div> </div> </div>
1	B	496	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>6%</div> </div> </div>
1	C	496	<div> <div>6%</div> <div> <div></div> <div>85%</div> <div>8%</div> <div>7%</div> </div> </div>
1	D	496	<div> <div>9%</div> <div> <div></div> <div>82%</div> <div>11%</div> <div>7%</div> </div> </div>
1	E	496	<div> <div>3%</div> <div> <div></div> <div>85%</div> <div>9%</div> <div>6%</div> </div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	496	<div><div></div><div>5%</div><div>86%</div><div>7%</div><div>6%</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 22924 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	466	Total	C	N	O	S	5	2	0
			3544	2256	605	666	17			
1	B	465	Total	C	N	O	S	0	4	0
			3566	2264	611	674	17			
1	C	462	Total	C	N	O	S	0	5	0
			3544	2250	609	668	17			
1	D	462	Total	C	N	O	S	0	3	0
			3489	2218	595	659	17			
1	E	466	Total	C	N	O	S	0	8	0
			3620	2304	617	681	18			
1	F	465	Total	C	N	O	S	0	12	0
			3618	2308	619	673	18			

There are 150 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	initiating methionine	UNP A0A0R4J3C9
A	-23	GLY	-	expression tag	UNP A0A0R4J3C9
A	-22	MET	-	expression tag	UNP A0A0R4J3C9
A	-21	HIS	-	expression tag	UNP A0A0R4J3C9
A	-20	HIS	-	expression tag	UNP A0A0R4J3C9
A	-19	HIS	-	expression tag	UNP A0A0R4J3C9
A	-18	HIS	-	expression tag	UNP A0A0R4J3C9
A	-17	HIS	-	expression tag	UNP A0A0R4J3C9
A	-16	HIS	-	expression tag	UNP A0A0R4J3C9
A	-15	SER	-	expression tag	UNP A0A0R4J3C9
A	-14	SER	-	expression tag	UNP A0A0R4J3C9
A	-13	GLY	-	expression tag	UNP A0A0R4J3C9
A	-12	VAL	-	expression tag	UNP A0A0R4J3C9
A	-11	ASP	-	expression tag	UNP A0A0R4J3C9
A	-10	LEU	-	expression tag	UNP A0A0R4J3C9
A	-9	GLY	-	expression tag	UNP A0A0R4J3C9
A	-8	THR	-	expression tag	UNP A0A0R4J3C9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLU	-	expression tag	UNP A0A0R4J3C9
A	-6	ASN	-	expression tag	UNP A0A0R4J3C9
A	-5	LEU	-	expression tag	UNP A0A0R4J3C9
A	-4	TYR	-	expression tag	UNP A0A0R4J3C9
A	-3	PHE	-	expression tag	UNP A0A0R4J3C9
A	-2	GLN	-	expression tag	UNP A0A0R4J3C9
A	-1	SER	-	expression tag	UNP A0A0R4J3C9
A	0	ASN	-	expression tag	UNP A0A0R4J3C9
B	-24	MET	-	initiating methionine	UNP A0A0R4J3C9
B	-23	GLY	-	expression tag	UNP A0A0R4J3C9
B	-22	MET	-	expression tag	UNP A0A0R4J3C9
B	-21	HIS	-	expression tag	UNP A0A0R4J3C9
B	-20	HIS	-	expression tag	UNP A0A0R4J3C9
B	-19	HIS	-	expression tag	UNP A0A0R4J3C9
B	-18	HIS	-	expression tag	UNP A0A0R4J3C9
B	-17	HIS	-	expression tag	UNP A0A0R4J3C9
B	-16	HIS	-	expression tag	UNP A0A0R4J3C9
B	-15	SER	-	expression tag	UNP A0A0R4J3C9
B	-14	SER	-	expression tag	UNP A0A0R4J3C9
B	-13	GLY	-	expression tag	UNP A0A0R4J3C9
B	-12	VAL	-	expression tag	UNP A0A0R4J3C9
B	-11	ASP	-	expression tag	UNP A0A0R4J3C9
B	-10	LEU	-	expression tag	UNP A0A0R4J3C9
B	-9	GLY	-	expression tag	UNP A0A0R4J3C9
B	-8	THR	-	expression tag	UNP A0A0R4J3C9
B	-7	GLU	-	expression tag	UNP A0A0R4J3C9
B	-6	ASN	-	expression tag	UNP A0A0R4J3C9
B	-5	LEU	-	expression tag	UNP A0A0R4J3C9
B	-4	TYR	-	expression tag	UNP A0A0R4J3C9
B	-3	PHE	-	expression tag	UNP A0A0R4J3C9
B	-2	GLN	-	expression tag	UNP A0A0R4J3C9
B	-1	SER	-	expression tag	UNP A0A0R4J3C9
B	0	ASN	-	expression tag	UNP A0A0R4J3C9
C	-24	MET	-	initiating methionine	UNP A0A0R4J3C9
C	-23	GLY	-	expression tag	UNP A0A0R4J3C9
C	-22	MET	-	expression tag	UNP A0A0R4J3C9
C	-21	HIS	-	expression tag	UNP A0A0R4J3C9
C	-20	HIS	-	expression tag	UNP A0A0R4J3C9
C	-19	HIS	-	expression tag	UNP A0A0R4J3C9
C	-18	HIS	-	expression tag	UNP A0A0R4J3C9
C	-17	HIS	-	expression tag	UNP A0A0R4J3C9
C	-16	HIS	-	expression tag	UNP A0A0R4J3C9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-15	SER	-	expression tag	UNP A0A0R4J3C9
C	-14	SER	-	expression tag	UNP A0A0R4J3C9
C	-13	GLY	-	expression tag	UNP A0A0R4J3C9
C	-12	VAL	-	expression tag	UNP A0A0R4J3C9
C	-11	ASP	-	expression tag	UNP A0A0R4J3C9
C	-10	LEU	-	expression tag	UNP A0A0R4J3C9
C	-9	GLY	-	expression tag	UNP A0A0R4J3C9
C	-8	THR	-	expression tag	UNP A0A0R4J3C9
C	-7	GLU	-	expression tag	UNP A0A0R4J3C9
C	-6	ASN	-	expression tag	UNP A0A0R4J3C9
C	-5	LEU	-	expression tag	UNP A0A0R4J3C9
C	-4	TYR	-	expression tag	UNP A0A0R4J3C9
C	-3	PHE	-	expression tag	UNP A0A0R4J3C9
C	-2	GLN	-	expression tag	UNP A0A0R4J3C9
C	-1	SER	-	expression tag	UNP A0A0R4J3C9
C	0	ASN	-	expression tag	UNP A0A0R4J3C9
D	-24	MET	-	initiating methionine	UNP A0A0R4J3C9
D	-23	GLY	-	expression tag	UNP A0A0R4J3C9
D	-22	MET	-	expression tag	UNP A0A0R4J3C9
D	-21	HIS	-	expression tag	UNP A0A0R4J3C9
D	-20	HIS	-	expression tag	UNP A0A0R4J3C9
D	-19	HIS	-	expression tag	UNP A0A0R4J3C9
D	-18	HIS	-	expression tag	UNP A0A0R4J3C9
D	-17	HIS	-	expression tag	UNP A0A0R4J3C9
D	-16	HIS	-	expression tag	UNP A0A0R4J3C9
D	-15	SER	-	expression tag	UNP A0A0R4J3C9
D	-14	SER	-	expression tag	UNP A0A0R4J3C9
D	-13	GLY	-	expression tag	UNP A0A0R4J3C9
D	-12	VAL	-	expression tag	UNP A0A0R4J3C9
D	-11	ASP	-	expression tag	UNP A0A0R4J3C9
D	-10	LEU	-	expression tag	UNP A0A0R4J3C9
D	-9	GLY	-	expression tag	UNP A0A0R4J3C9
D	-8	THR	-	expression tag	UNP A0A0R4J3C9
D	-7	GLU	-	expression tag	UNP A0A0R4J3C9
D	-6	ASN	-	expression tag	UNP A0A0R4J3C9
D	-5	LEU	-	expression tag	UNP A0A0R4J3C9
D	-4	TYR	-	expression tag	UNP A0A0R4J3C9
D	-3	PHE	-	expression tag	UNP A0A0R4J3C9
D	-2	GLN	-	expression tag	UNP A0A0R4J3C9
D	-1	SER	-	expression tag	UNP A0A0R4J3C9
D	0	ASN	-	expression tag	UNP A0A0R4J3C9
E	-24	MET	-	initiating methionine	UNP A0A0R4J3C9

Continued on next page...

Continued from previous page...

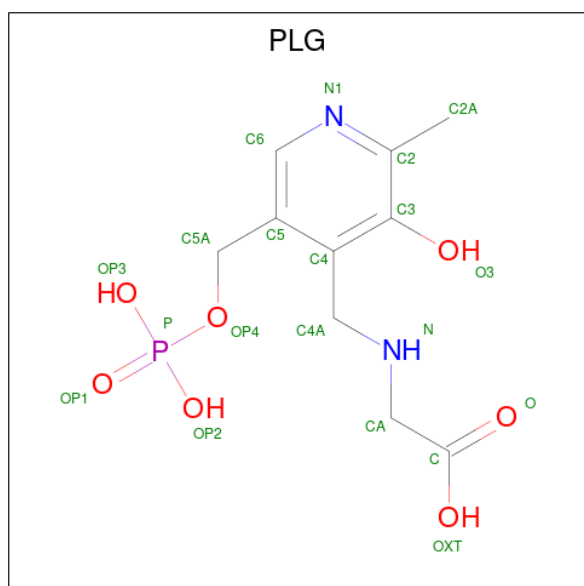
Chain	Residue	Modelled	Actual	Comment	Reference
E	-23	GLY	-	expression tag	UNP A0A0R4J3C9
E	-22	MET	-	expression tag	UNP A0A0R4J3C9
E	-21	HIS	-	expression tag	UNP A0A0R4J3C9
E	-20	HIS	-	expression tag	UNP A0A0R4J3C9
E	-19	HIS	-	expression tag	UNP A0A0R4J3C9
E	-18	HIS	-	expression tag	UNP A0A0R4J3C9
E	-17	HIS	-	expression tag	UNP A0A0R4J3C9
E	-16	HIS	-	expression tag	UNP A0A0R4J3C9
E	-15	SER	-	expression tag	UNP A0A0R4J3C9
E	-14	SER	-	expression tag	UNP A0A0R4J3C9
E	-13	GLY	-	expression tag	UNP A0A0R4J3C9
E	-12	VAL	-	expression tag	UNP A0A0R4J3C9
E	-11	ASP	-	expression tag	UNP A0A0R4J3C9
E	-10	LEU	-	expression tag	UNP A0A0R4J3C9
E	-9	GLY	-	expression tag	UNP A0A0R4J3C9
E	-8	THR	-	expression tag	UNP A0A0R4J3C9
E	-7	GLU	-	expression tag	UNP A0A0R4J3C9
E	-6	ASN	-	expression tag	UNP A0A0R4J3C9
E	-5	LEU	-	expression tag	UNP A0A0R4J3C9
E	-4	TYR	-	expression tag	UNP A0A0R4J3C9
E	-3	PHE	-	expression tag	UNP A0A0R4J3C9
E	-2	GLN	-	expression tag	UNP A0A0R4J3C9
E	-1	SER	-	expression tag	UNP A0A0R4J3C9
E	0	ASN	-	expression tag	UNP A0A0R4J3C9
F	-24	MET	-	initiating methionine	UNP A0A0R4J3C9
F	-23	GLY	-	expression tag	UNP A0A0R4J3C9
F	-22	MET	-	expression tag	UNP A0A0R4J3C9
F	-21	HIS	-	expression tag	UNP A0A0R4J3C9
F	-20	HIS	-	expression tag	UNP A0A0R4J3C9
F	-19	HIS	-	expression tag	UNP A0A0R4J3C9
F	-18	HIS	-	expression tag	UNP A0A0R4J3C9
F	-17	HIS	-	expression tag	UNP A0A0R4J3C9
F	-16	HIS	-	expression tag	UNP A0A0R4J3C9
F	-15	SER	-	expression tag	UNP A0A0R4J3C9
F	-14	SER	-	expression tag	UNP A0A0R4J3C9
F	-13	GLY	-	expression tag	UNP A0A0R4J3C9
F	-12	VAL	-	expression tag	UNP A0A0R4J3C9
F	-11	ASP	-	expression tag	UNP A0A0R4J3C9
F	-10	LEU	-	expression tag	UNP A0A0R4J3C9
F	-9	GLY	-	expression tag	UNP A0A0R4J3C9
F	-8	THR	-	expression tag	UNP A0A0R4J3C9
F	-7	GLU	-	expression tag	UNP A0A0R4J3C9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
F	-6	ASN	-	expression tag	UNP A0A0R4J3C9
F	-5	LEU	-	expression tag	UNP A0A0R4J3C9
F	-4	TYR	-	expression tag	UNP A0A0R4J3C9
F	-3	PHE	-	expression tag	UNP A0A0R4J3C9
F	-2	GLN	-	expression tag	UNP A0A0R4J3C9
F	-1	SER	-	expression tag	UNP A0A0R4J3C9
F	0	ASN	-	expression tag	UNP A0A0R4J3C9

- Molecule 2 is N-GLYCINE-[3-HYDROXY-2-METHYL-5-PHOSPHONOXYMETHYL-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		
2	F	1	Total	C	N	O	P	0	0
			20	10	2	7	1		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

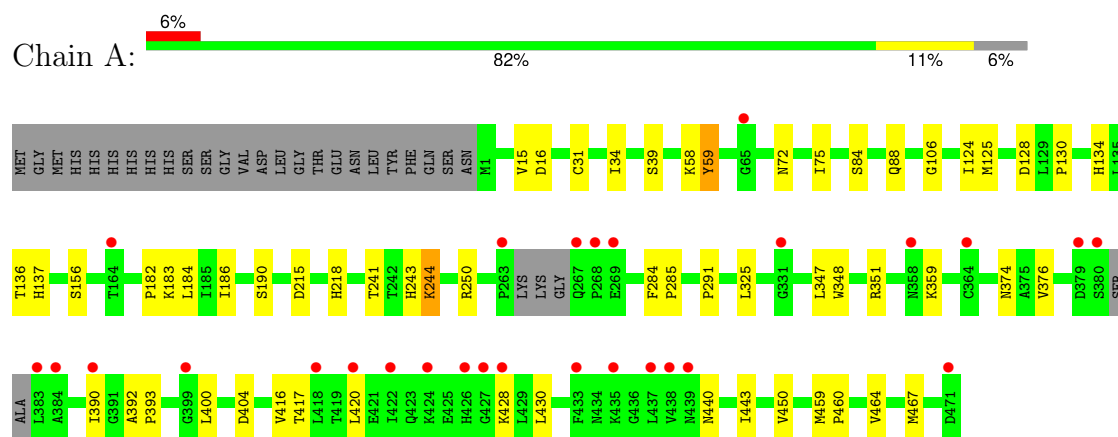
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	201	Total	O	0	0
			201	201		
4	B	253	Total	O	0	0
			253	253		
4	C	222	Total	O	0	0
			222	222		
4	D	165	Total	O	0	0
			165	165		
4	E	295	Total	O	0	0
			295	295		
4	F	275	Total	O	0	0
			275	275		

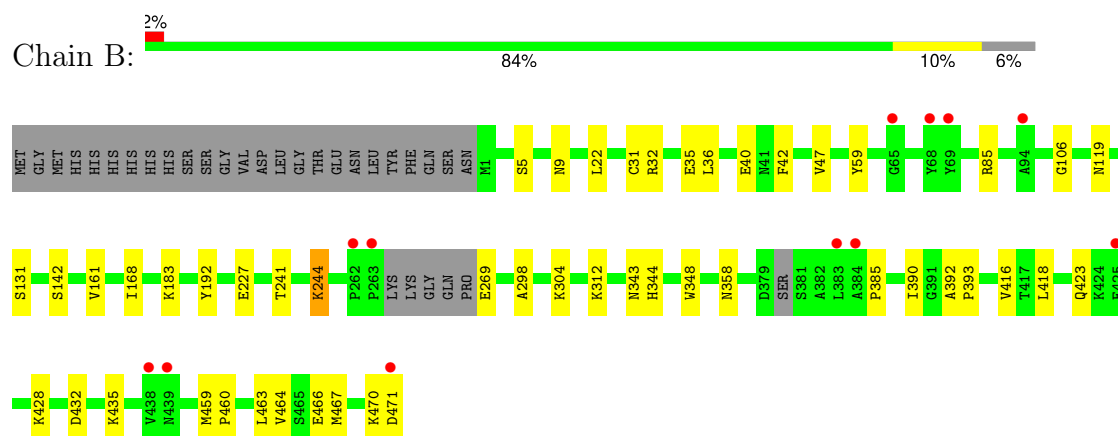
3 Residue-property plots [i](#)

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.

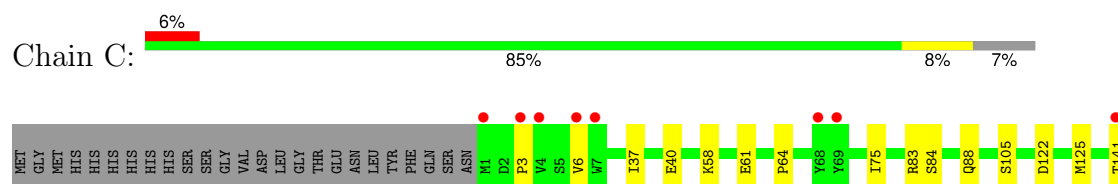
• Molecule 1: Serine hydroxymethyltransferase

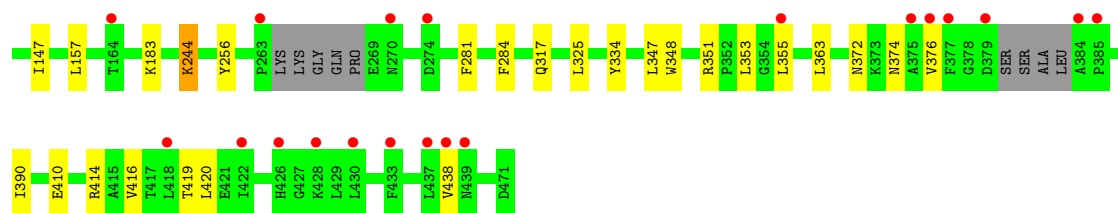


• Molecule 1: Serine hydroxymethyltransferase

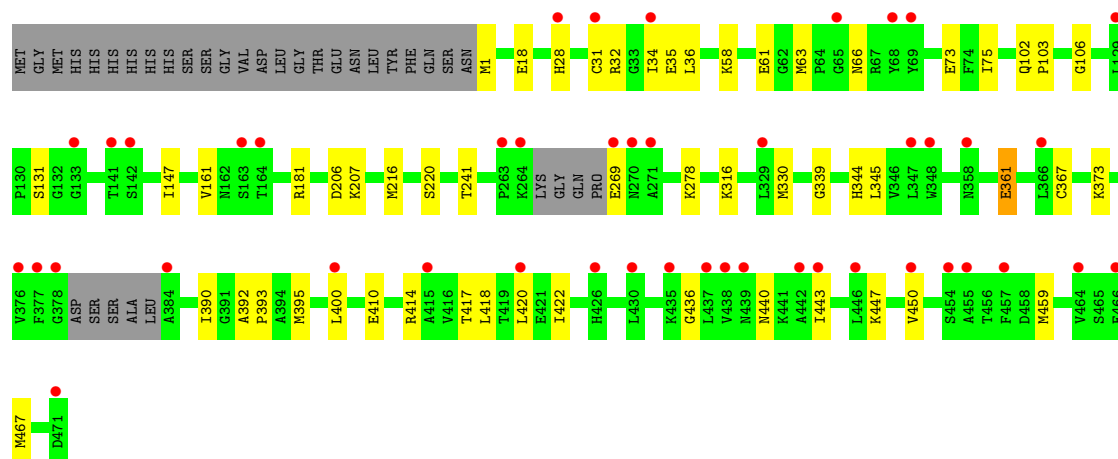
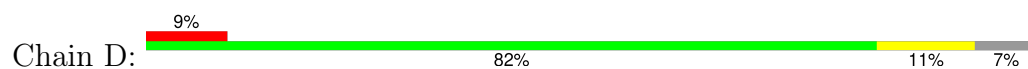


• Molecule 1: Serine hydroxymethyltransferase

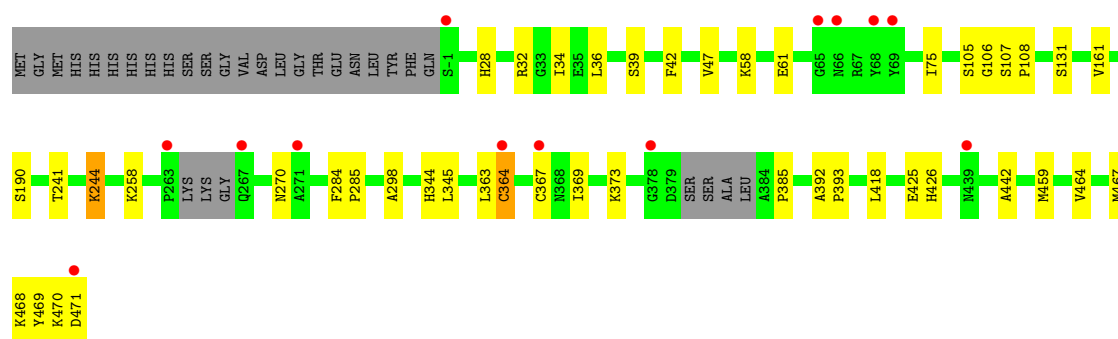
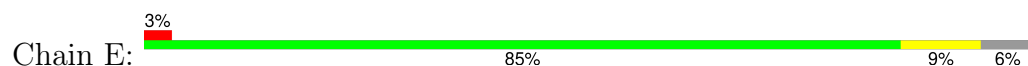




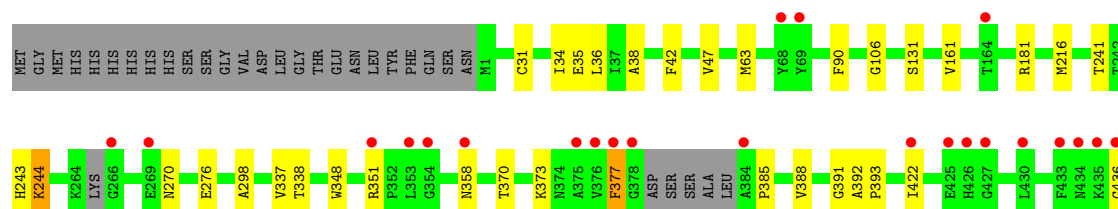
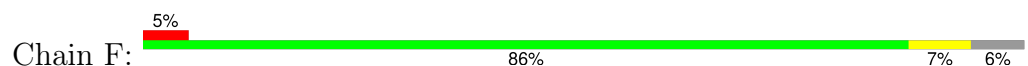
• Molecule 1: Serine hydroxymethyltransferase

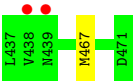


• Molecule 1: Serine hydroxymethyltransferase



• 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	174.83Å 174.83Å 184.11Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.67 – 1.90 48.67 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.67-1.90) 100.0 (48.67-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.199 , 0.231 0.199 , 0.230	Depositor DCC
R_{free} test set	12847 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	27.6	Xtriage
Anisotropy	0.687	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	22924	wwPDB-VP
Average B, all atoms (Å ²)	36.0	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 37.42 % 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 4.3030e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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, EDO

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	0.32	0/3636	0.54	1/4935 (0.0%)
1	B	0.33	0/3659	0.55	0/4959
1	C	0.33	0/3637	0.54	0/4930
1	D	0.32	0/3579	0.51	0/4858
1	E	0.37	0/3727	0.60	2/5046 (0.0%)
1	F	0.37	0/3740	0.56	0/5061
All	All	0.34	0/21978	0.55	3/29789 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	363	LEU	CA-C-N	-8.52	106.96	122.38
1	E	363	LEU	C-N-CA	-8.52	106.96	122.38
1	A	428	LYS	CB-CG-CD	5.36	123.62	111.30

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	3544	0	3408	44	0
1	B	3566	0	3455	37	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3544	0	3423	25	0
1	D	3489	0	3327	45	0
1	E	3620	0	3542	30	0
1	F	3618	0	3561	27	0
2	A	20	0	12	4	0
2	B	20	0	12	1	0
2	C	20	0	12	3	0
2	D	20	0	12	1	0
2	E	20	0	12	3	0
2	F	20	0	12	3	0
3	B	4	0	6	0	0
3	C	4	0	6	0	0
3	E	4	0	6	1	0
4	A	201	0	0	5	0
4	B	253	0	0	3	0
4	C	222	0	0	3	0
4	D	165	0	0	7	0
4	E	295	0	0	3	0
4	F	275	0	0	4	1
All	All	22924	0	20806	209	1

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 5.

The worst 5 of 209 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:D:35:GLU:H	1:D:395:MET:HE2	1.24	0.99
1:B:183:LYS:NZ	4:B:601:HOH:O	2.14	0.80
1:A:183:LYS:NZ	4:A:801:HOH:O	2.05	0.79
1:C:84:SER:O	1:C:88:GLN:HG3	1.84	0.77
1:B:358:ASN:HD21	1:B:385:PRO:HD2	1.50	0.77

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:869:HOH:O	4:F:870:HOH:O[6_345]	2.17	0.03

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%)	450 (97%)	12 (3%)	1 (0%)	44	36
1	B	463/496 (93%)	453 (98%)	9 (2%)	1 (0%)	44	36
1	C	461/496 (93%)	448 (97%)	12 (3%)	1 (0%)	44	36
1	D	459/496 (92%)	444 (97%)	15 (3%)	0	100	100
1	E	468/496 (94%)	458 (98%)	9 (2%)	1 (0%)	44	36
1	F	472/496 (95%)	455 (96%)	14 (3%)	3 (1%)	22	13
All	All	2786/2976 (94%)	2708 (97%)	71 (2%)	7 (0%)	44	29

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	377	PHE
1	A	244	LYS
1	E	244	LYS
1	B	244	LYS
1	C	244	LYS

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	362/407 (89%)	361 (100%)	1 (0%)	91	92
1	B	369/407 (91%)	368 (100%)	1 (0%)	91	92
1	C	365/407 (90%)	363 (100%)	2 (0%)	86	88

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	352/407 (86%)	348 (99%)	4 (1%)	70	71
1	E	380/407 (93%)	377 (99%)	3 (1%)	79	80
1	F	378/407 (93%)	378 (100%)	0	100	100
All	All	2206/2442 (90%)	2195 (100%)	11 (0%)	86	88

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	361	GLU
1	E	105	SER
1	E	367	CYS
1	E	364	CYS
1	D	73[A]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	292	HIS
1	F	305	GLN
1	F	295	GLN
1	C	358	ASN
1	E	20	HIS

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

9 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$
3	EDO	E	502	-	3,3,3	0.43	0	2,2,2	0.53	0
2	PLG	C	701	-	20,20,20	1.34	4 (20%)	26,28,28	1.27	2 (7%)
3	EDO	B	501	-	3,3,3	0.46	0	2,2,2	0.34	0
2	PLG	A	701	-	20,20,20	1.31	1 (5%)	26,28,28	1.29	3 (11%)
2	PLG	D	601	-	20,20,20	1.31	2 (10%)	26,28,28	1.33	2 (7%)
3	EDO	C	702	-	3,3,3	0.46	0	2,2,2	0.42	0
2	PLG	F	501	-	20,20,20	1.37	3 (15%)	26,28,28	1.47	4 (15%)
2	PLG	E	501	-	20,20,20	1.30	1 (5%)	26,28,28	1.41	3 (11%)
2	PLG	B	502	-	20,20,20	1.26	2 (10%)	26,28,28	1.30	3 (11%)

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
3	EDO	E	502	-	-	1/1/1/1	-
2	PLG	C	701	-	-	5/12/12/12	0/1/1/1
3	EDO	B	501	-	-	0/1/1/1	-
2	PLG	A	701	-	-	5/12/12/12	0/1/1/1
2	PLG	D	601	-	-	5/12/12/12	0/1/1/1
3	EDO	C	702	-	-	0/1/1/1	-
2	PLG	F	501	-	-	4/12/12/12	0/1/1/1
2	PLG	E	501	-	-	5/12/12/12	0/1/1/1
2	PLG	B	502	-	-	7/12/12/12	0/1/1/1

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	502	PLG	C5-C4	-2.81	1.36	1.40
2	E	501	PLG	C5-C4	-2.65	1.36	1.40
2	D	601	PLG	C2A-C2	2.40	1.54	1.50
2	D	601	PLG	C5-C4	-2.38	1.37	1.40
2	A	701	PLG	C2A-C2	2.28	1.54	1.50

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	501	PLG	C6-C5-C4	3.99	121.08	118.06
2	E	501	PLG	C5-C6-N1	-3.37	118.34	123.83
2	C	701	PLG	C6-C5-C4	3.18	120.47	118.06
2	D	601	PLG	C6-C5-C4	3.09	120.40	118.06
2	D	601	PLG	C5-C6-N1	-2.87	119.17	123.83

There are no chirality outliers.

5 of 32 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	PLG	C3-C4-C4A-N
2	A	701	PLG	C5-C4-C4A-N
2	A	701	PLG	C5A-OP4-P-OP1
2	A	701	PLG	C5A-OP4-P-OP3
2	B	502	PLG	C5-C4-C4A-N

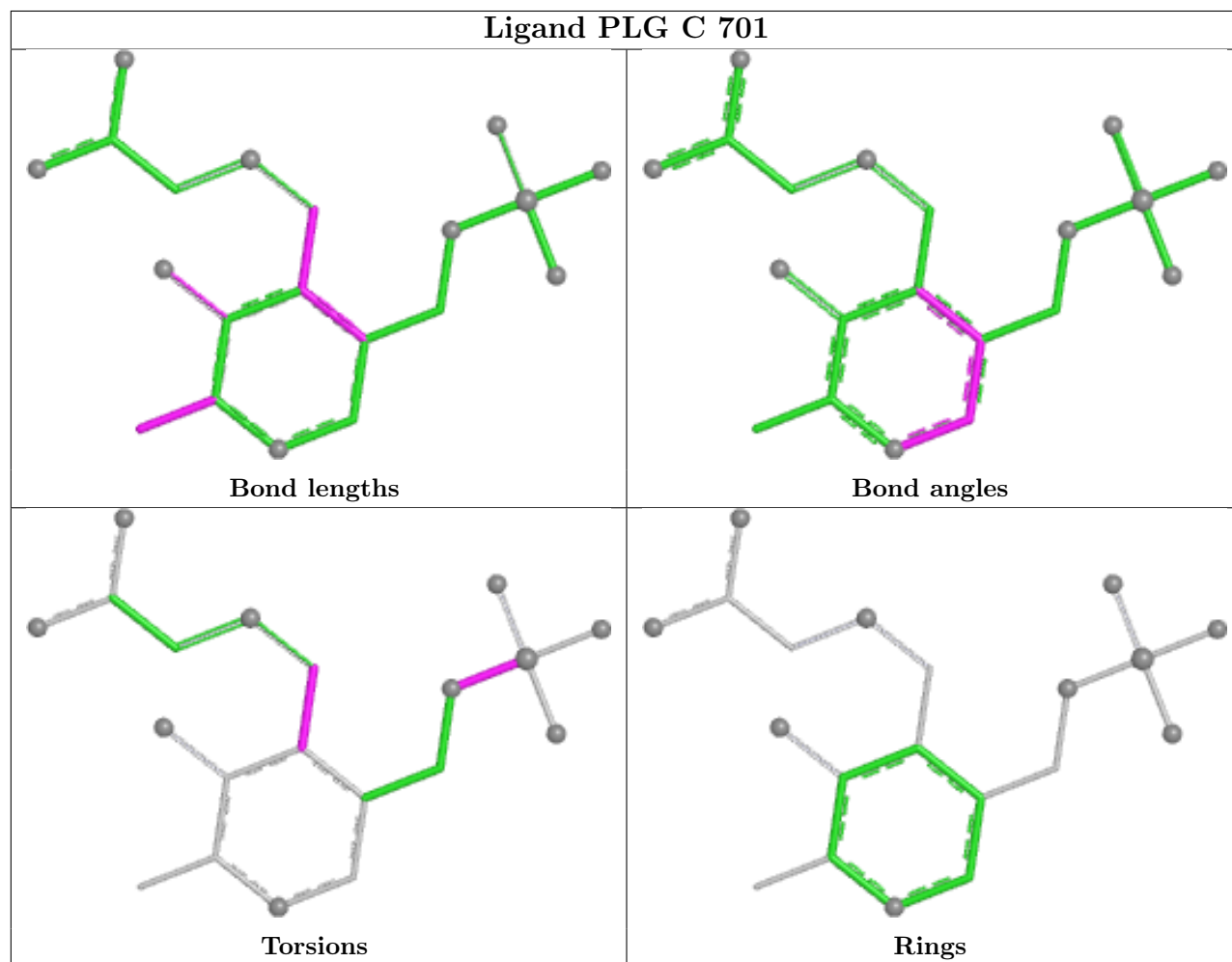
There are no ring outliers.

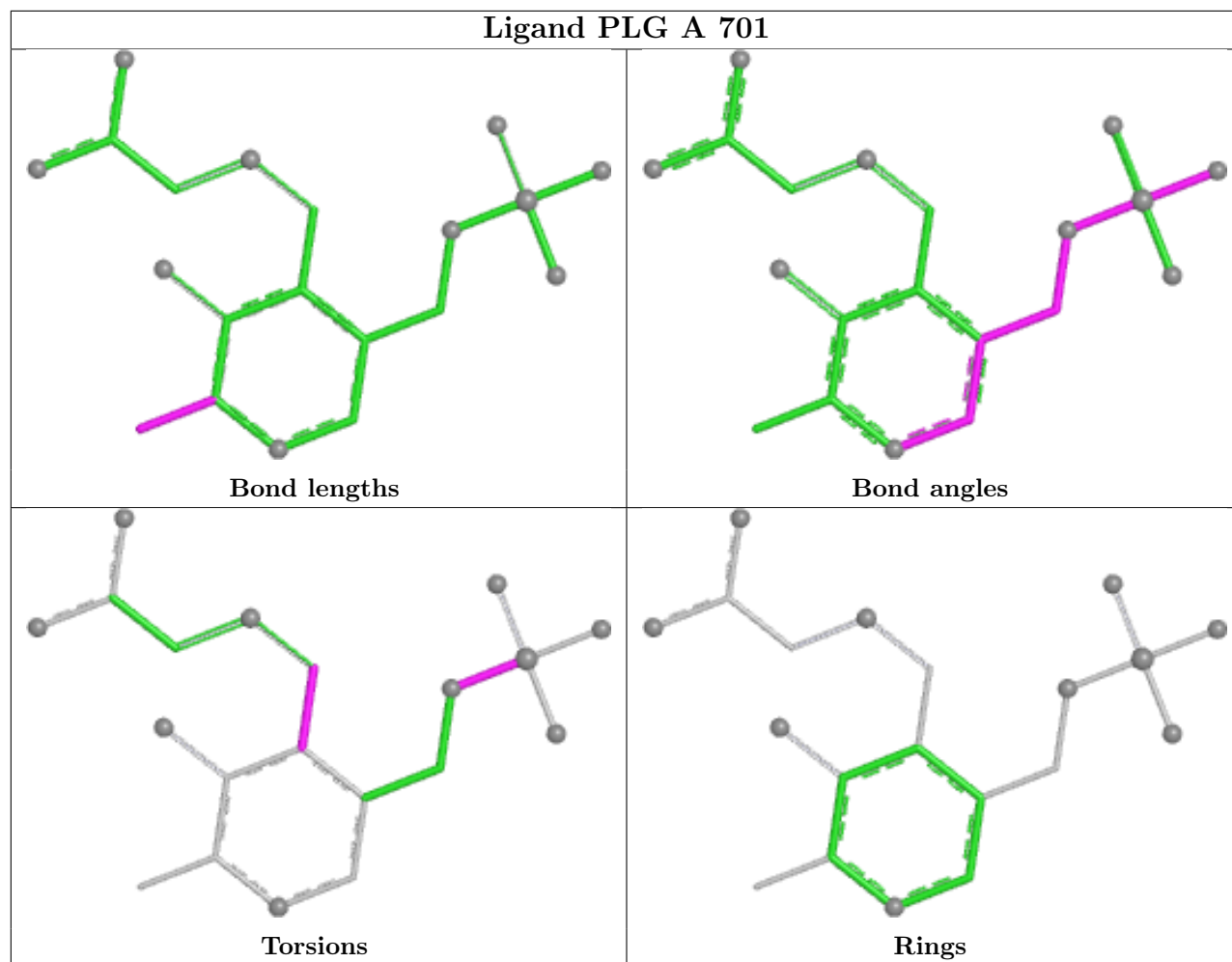
7 monomers are involved in 16 short contacts:

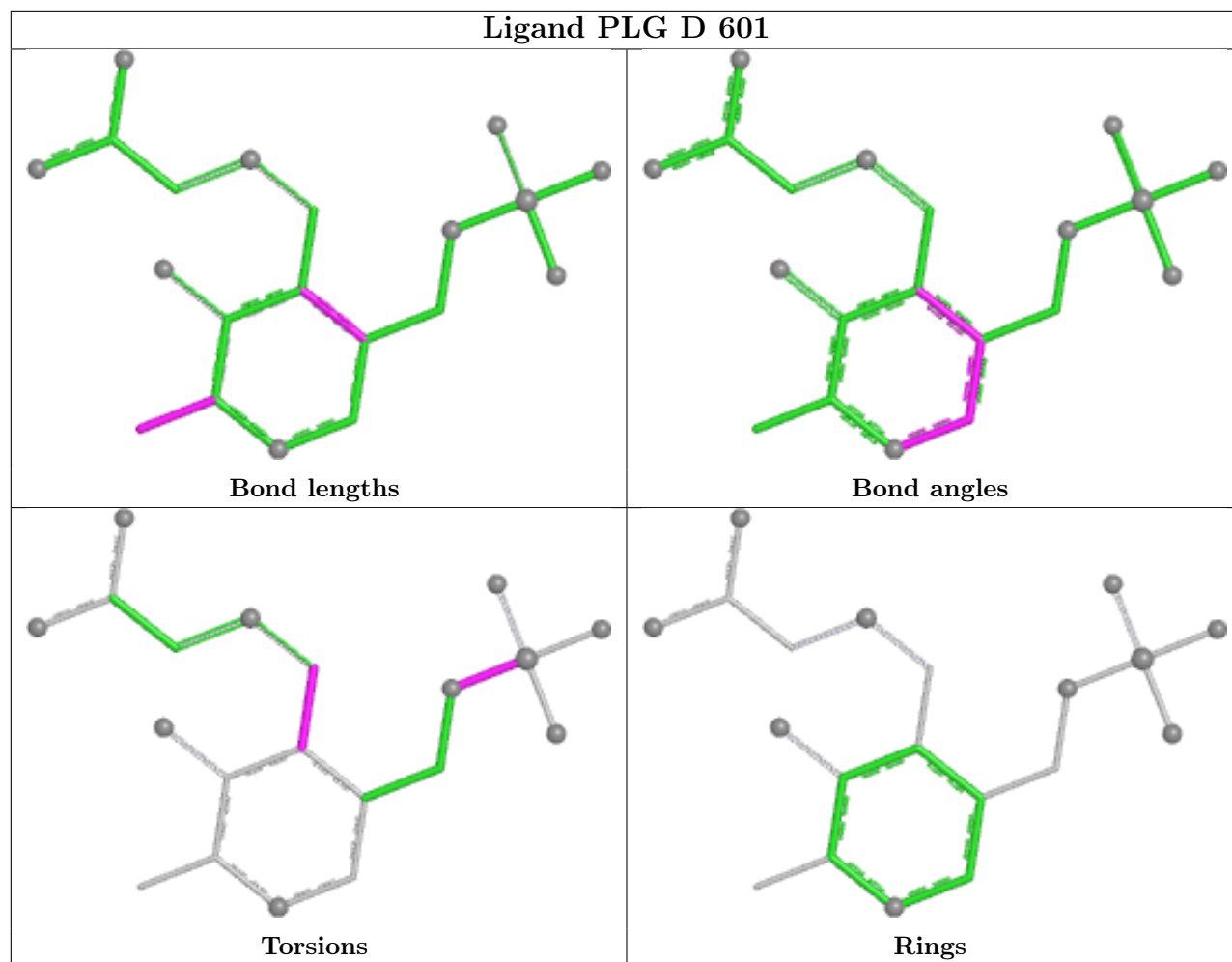
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	502	EDO	1	0
2	C	701	PLG	3	0
2	A	701	PLG	4	0
2	D	601	PLG	1	0
2	F	501	PLG	3	0
2	E	501	PLG	3	0
2	B	502	PLG	1	0

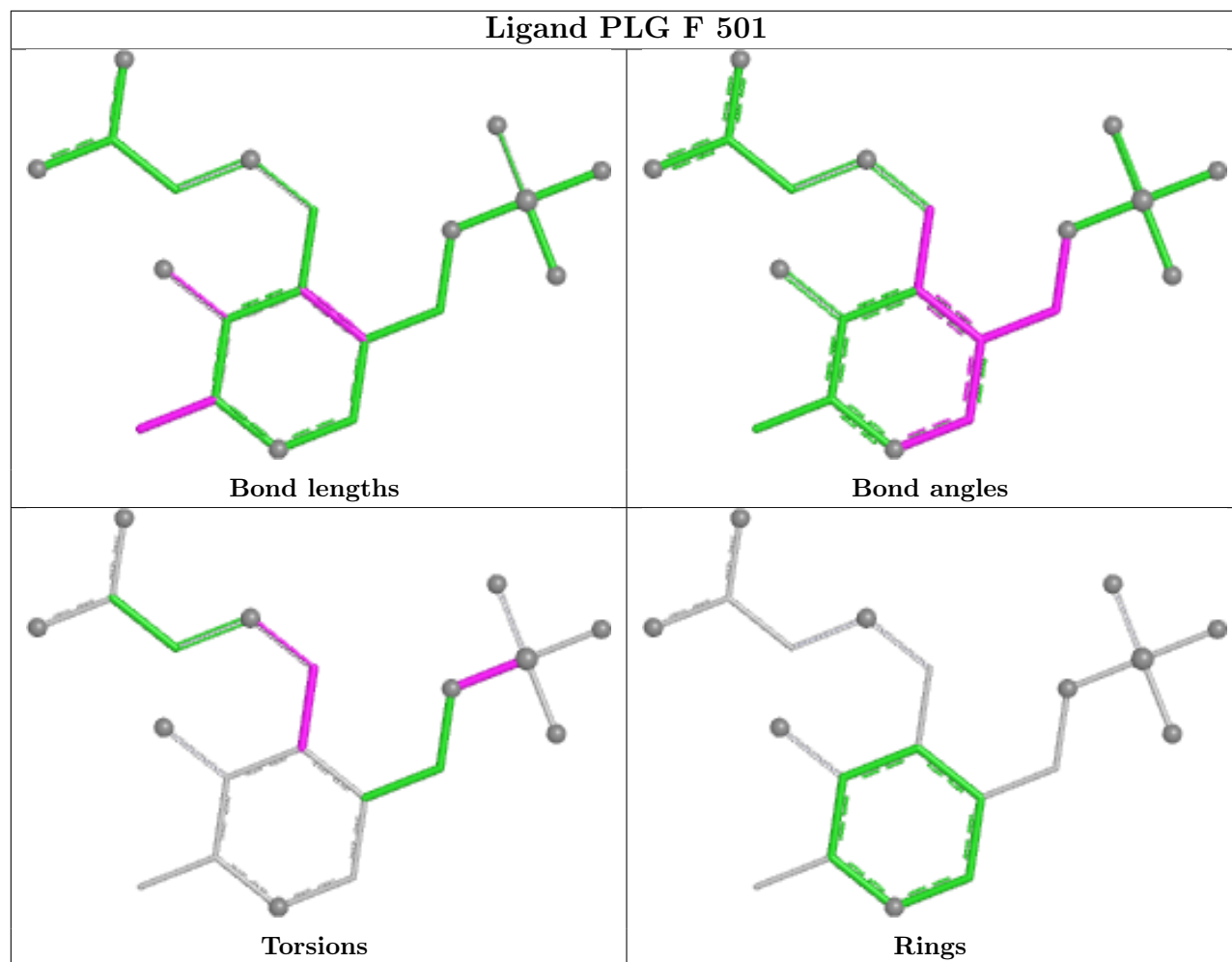
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

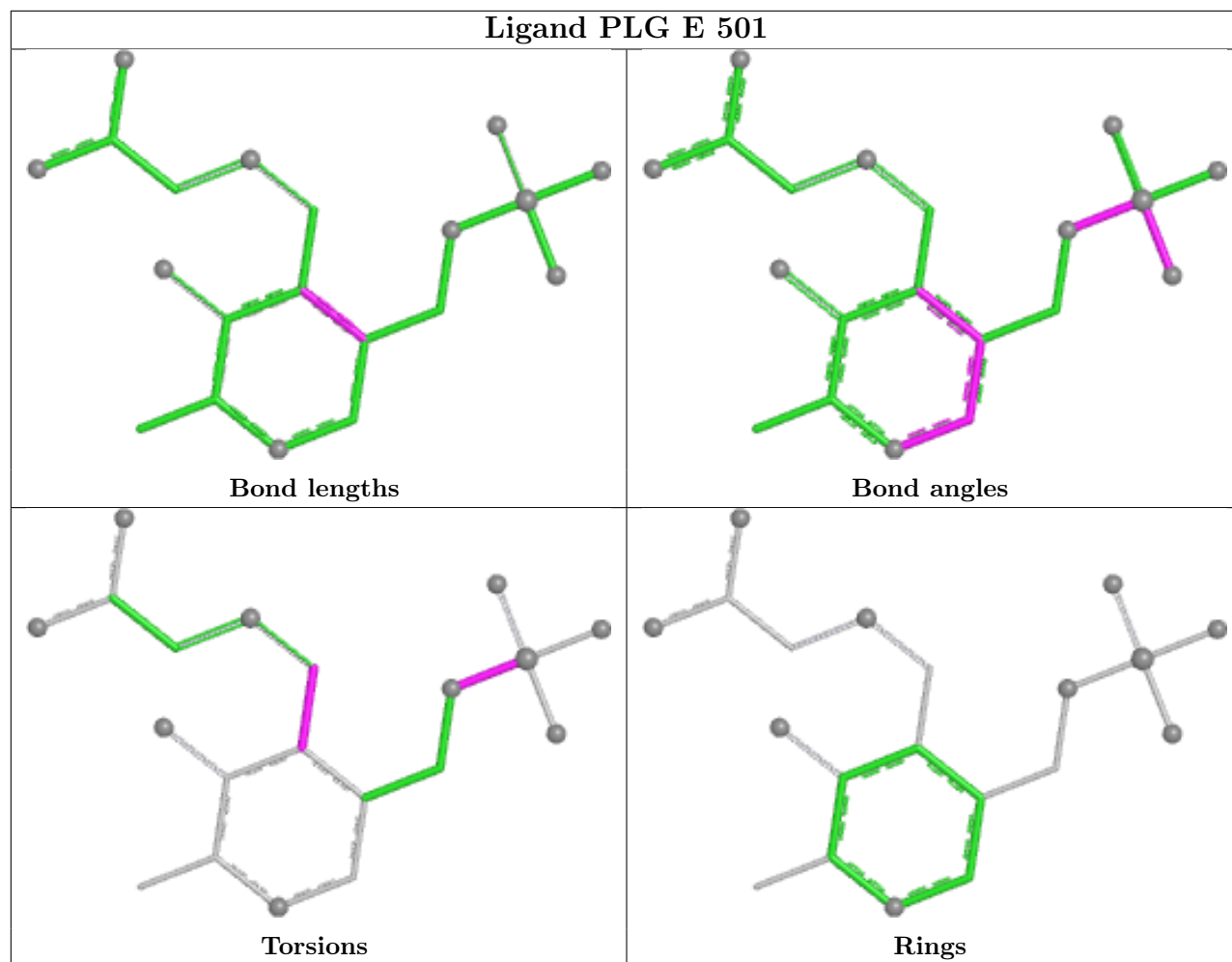
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

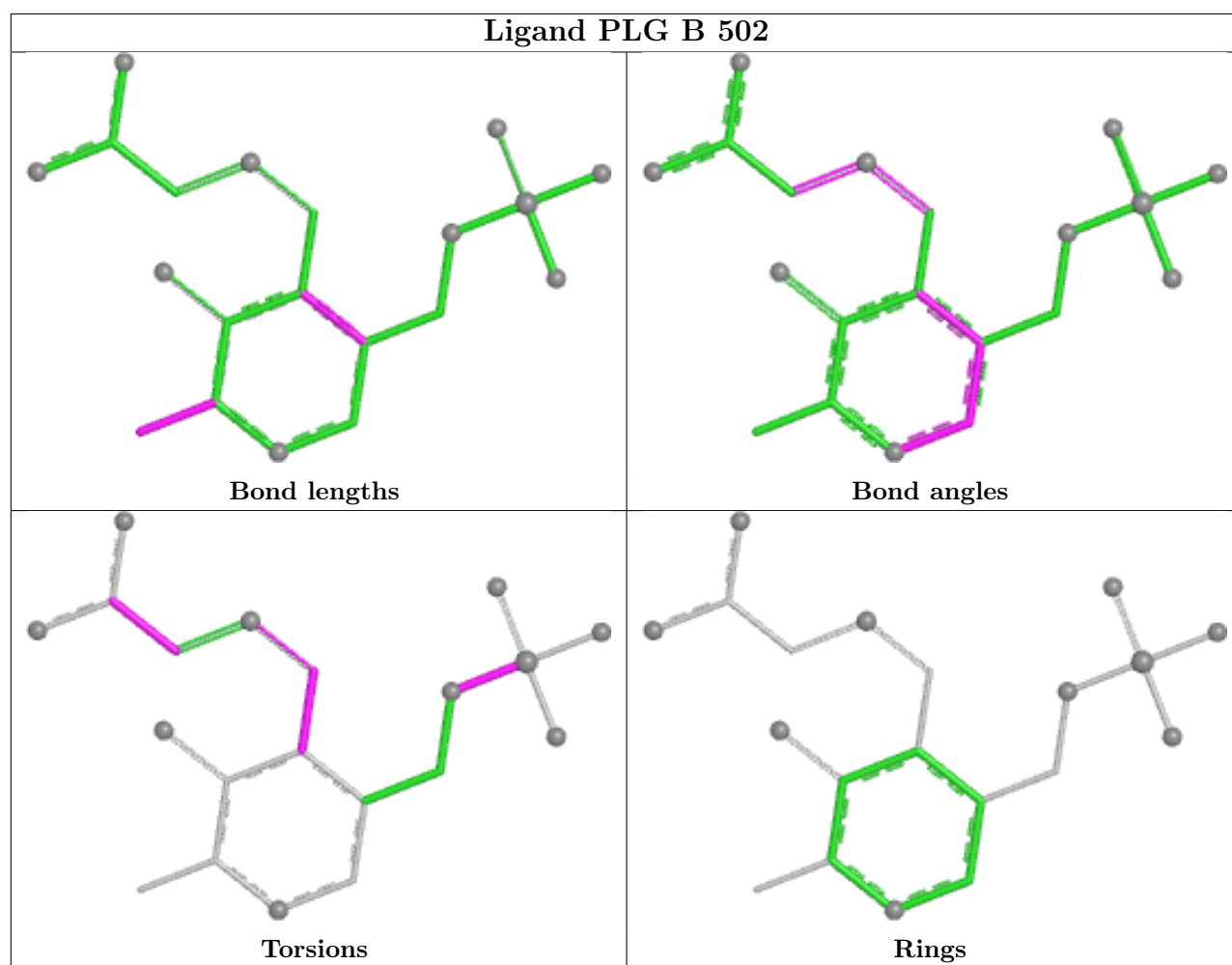












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	466/496 (93%)	0.55	28 (6%) 29 30	11, 36, 63, 73	2 (0%)
1	B	465/496 (93%)	0.35	12 (2%) 57 59	20, 33, 52, 70	4 (0%)
1	C	462/496 (93%)	0.41	28 (6%) 28 29	15, 34, 64, 82	5 (1%)
1	D	462/496 (93%)	0.80	45 (9%) 15 15	21, 39, 73, 85	3 (0%)
1	E	466/496 (93%)	0.26	13 (2%) 55 57	16, 29, 48, 69	8 (1%)
1	F	465/496 (93%)	0.34	25 (5%) 32 34	16, 30, 60, 76	12 (2%)
All	All	2786/2976 (93%)	0.45	151 (5%) 32 34	11, 33, 61, 85	34 (1%)

The worst 5 of 151 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	364	CYS	10.2
1	D	377	PHE	5.8
1	A	383	LEU	5.1
1	D	438	VAL	5.0
1	F	377	PHE	4.4

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

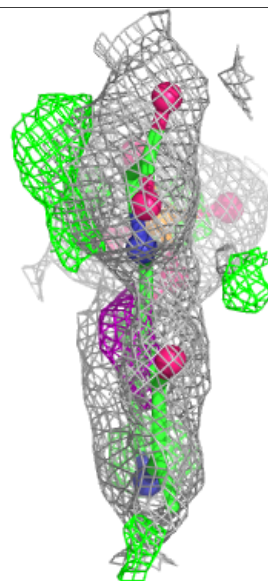
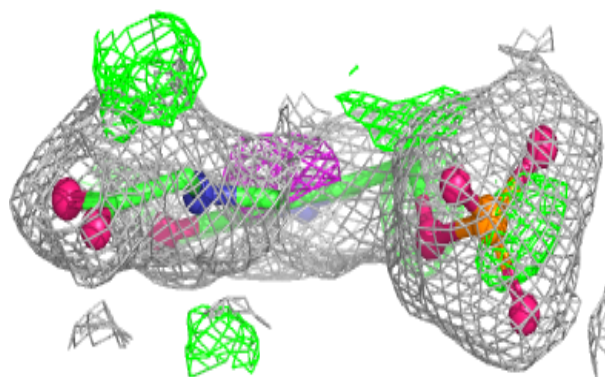
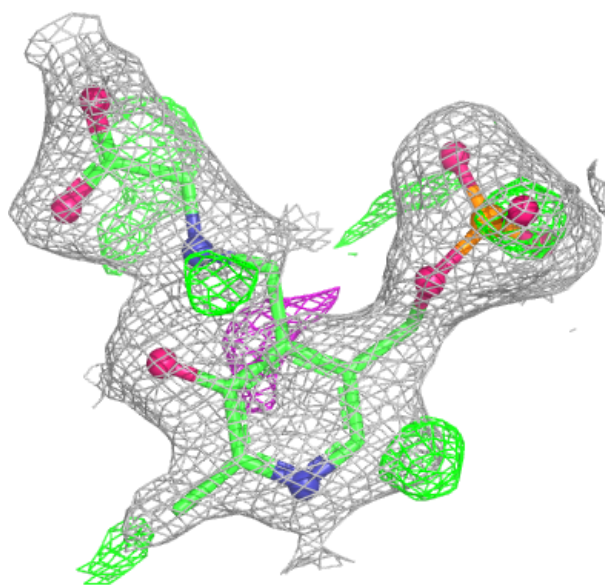
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PLG	A	701	20/20	0.90	0.15	28,38,45,51	20
2	PLG	D	601	20/20	0.90	0.15	34,42,48,50	20
3	EDO	E	502	4/4	0.90	0.12	30,31,33,39	4
2	PLG	F	501	20/20	0.93	0.13	26,34,41,42	20
3	EDO	B	501	4/4	0.93	0.14	32,36,36,37	0
3	EDO	C	702	4/4	0.93	0.10	31,33,34,37	0
2	PLG	C	701	20/20	0.93	0.12	32,41,48,49	0
2	PLG	B	502	20/20	0.94	0.11	28,35,42,46	20
2	PLG	E	501	20/20	0.94	0.10	26,31,37,38	20

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.

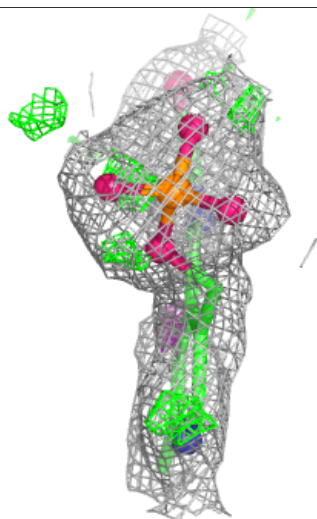
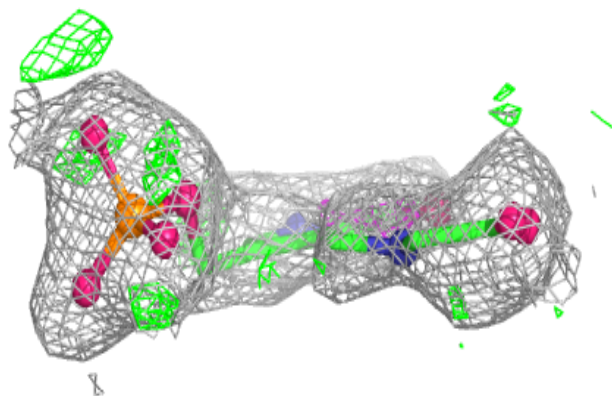
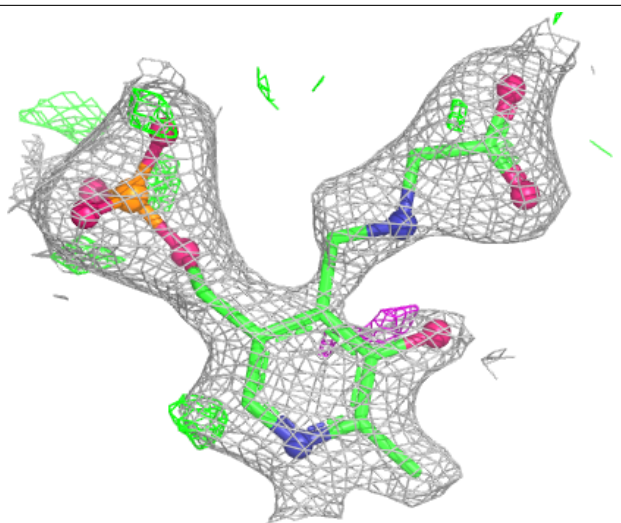
Electron density around PLG A 701:

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



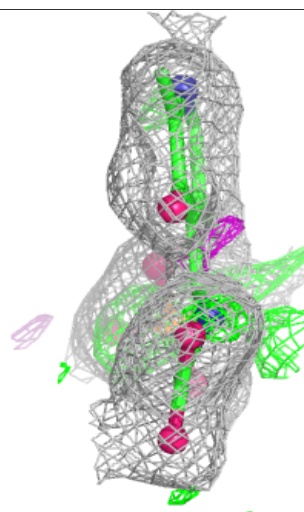
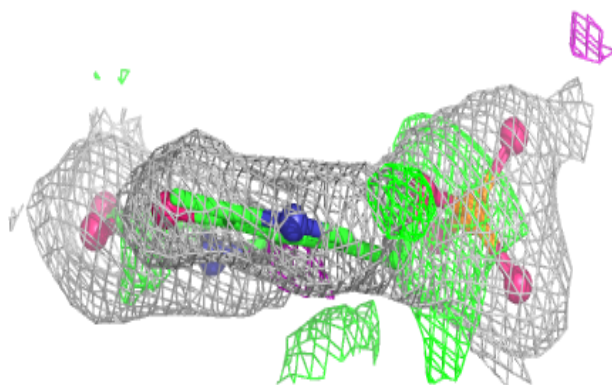
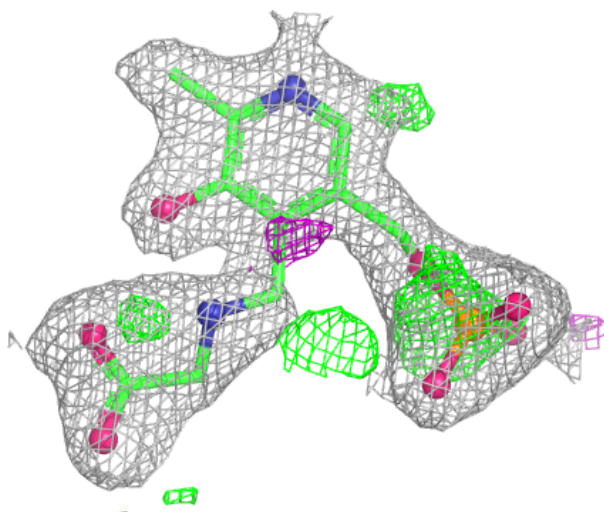
Electron density around PLG D 601:

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



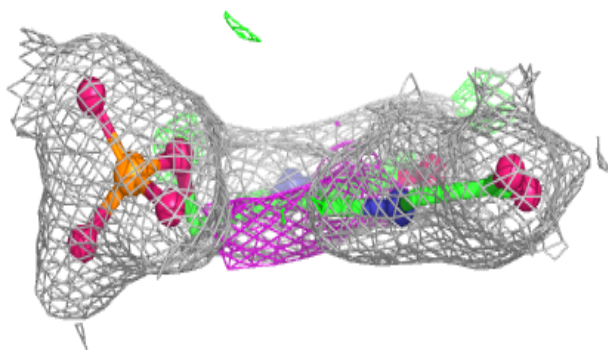
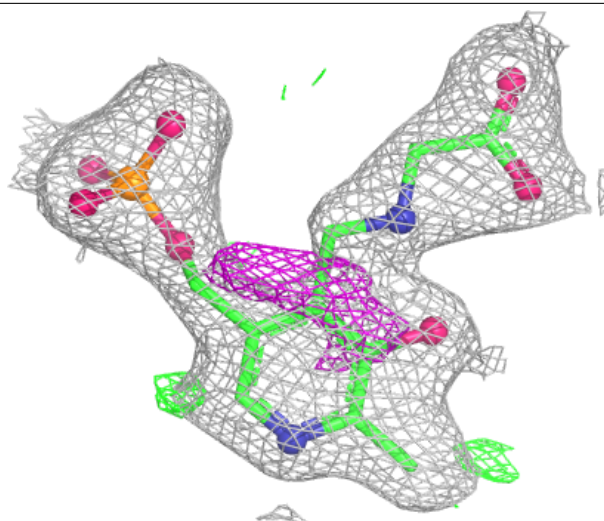
Electron density around PLG F 501:

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



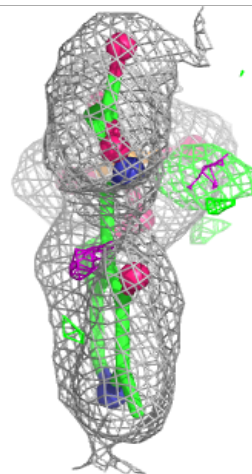
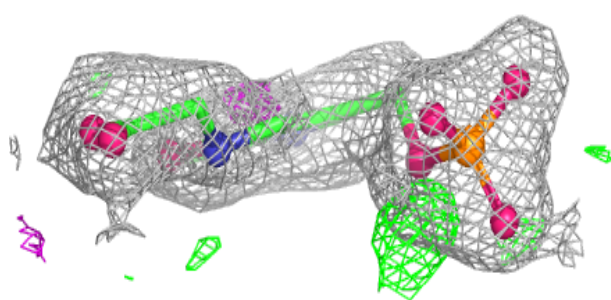
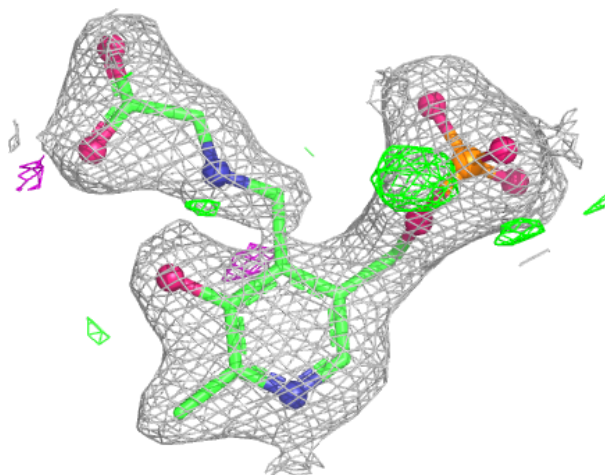
Electron density around PLG C 701:

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



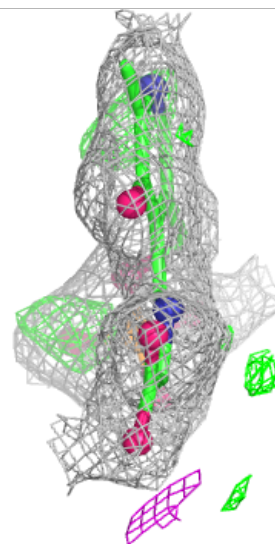
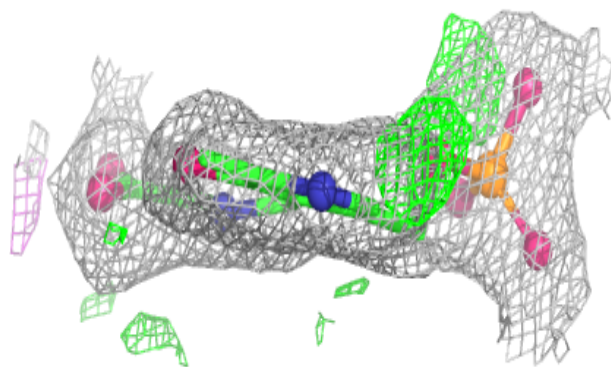
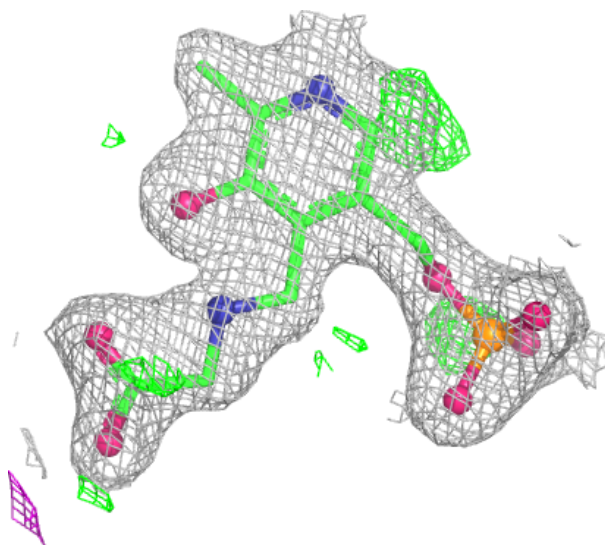
Electron density around PLG B 502:

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



Electron density around PLG 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.