



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:47 PM UTC

PDB ID : 8GKT / pdb_00008gkt
Title : Human mitochondrial serine hydroxymethyltransferase (SHMT2) in complex with PLP, glycine and AGF320 inhibitor
Authors : Katinas, J.M.; Dann III, C.E.
Deposited on : 2023-03-20
Resolution : 2.64 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

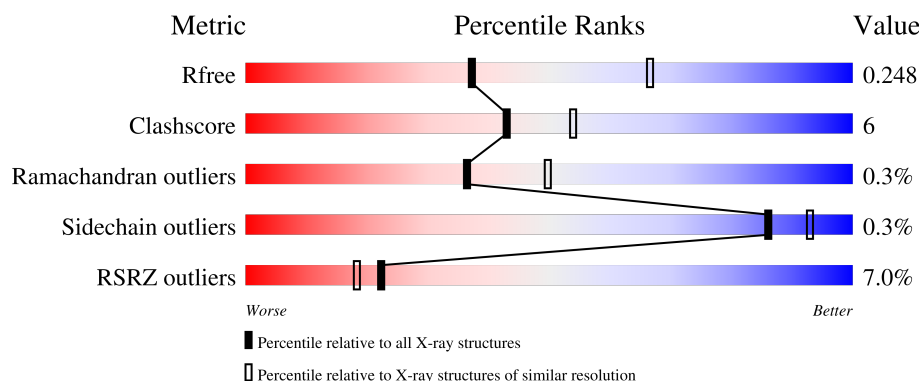
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	493	
1	B	493	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7327 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, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	457	Total	C	N	O	S	0	0	0
			3548	2235	636	661	16			
1	B	456	Total	C	N	O	S	0	0	0
			3439	2169	605	649	16			

There are 34 discrepancies between the modelled and reference sequences:

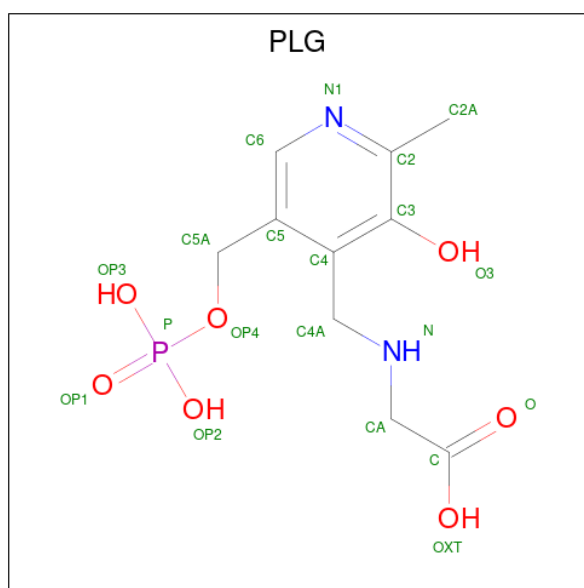
Chain	Residue	Modelled	Actual	Comment	Reference
A	12	MET	-	initiating methionine	UNP P34897
A	13	GLY	-	expression tag	UNP P34897
A	14	SER	-	expression tag	UNP P34897
A	15	SER	-	expression tag	UNP P34897
A	16	HIS	-	expression tag	UNP P34897
A	17	HIS	-	expression tag	UNP P34897
A	18	HIS	-	expression tag	UNP P34897
A	19	HIS	-	expression tag	UNP P34897
A	20	HIS	-	expression tag	UNP P34897
A	21	HIS	-	expression tag	UNP P34897
A	22	SER	-	expression tag	UNP P34897
A	23	SER	-	expression tag	UNP P34897
A	24	GLY	-	expression tag	UNP P34897
A	25	LEU	-	expression tag	UNP P34897
A	26	VAL	-	expression tag	UNP P34897
A	27	PRO	-	expression tag	UNP P34897
A	28	ARG	-	expression tag	UNP P34897
B	12	MET	-	initiating methionine	UNP P34897
B	13	GLY	-	expression tag	UNP P34897
B	14	SER	-	expression tag	UNP P34897
B	15	SER	-	expression tag	UNP P34897
B	16	HIS	-	expression tag	UNP P34897
B	17	HIS	-	expression tag	UNP P34897
B	18	HIS	-	expression tag	UNP P34897
B	19	HIS	-	expression tag	UNP P34897

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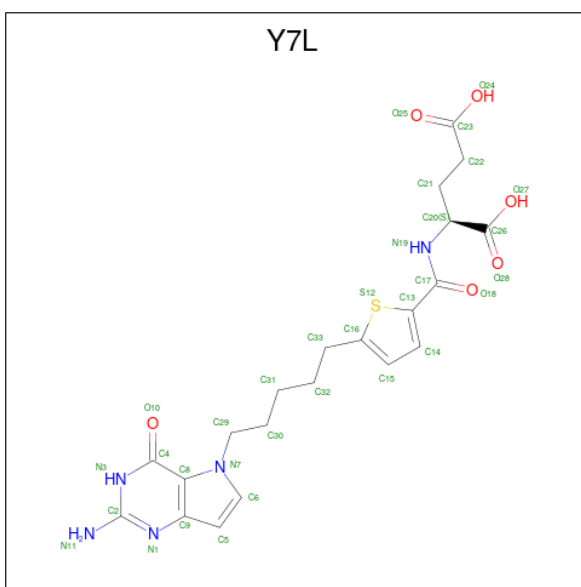
Chain	Residue	Modelled	Actual	Comment	Reference
B	20	HIS	-	expression tag	UNP P34897
B	21	HIS	-	expression tag	UNP P34897
B	22	SER	-	expression tag	UNP P34897
B	23	SER	-	expression tag	UNP P34897
B	24	GLY	-	expression tag	UNP P34897
B	25	LEU	-	expression tag	UNP P34897
B	26	VAL	-	expression tag	UNP P34897
B	27	PRO	-	expression tag	UNP P34897
B	28	ARG	-	expression tag	UNP P34897

- Molecule 2 is N-GLYCINE-[3-HYDROXY-2-METHYL-5-PHOSPHONOOXYMETHYL-PYRIDIN-4-YL-METHANE] (CCD ID: PLG) (formula: $C_{10}H_{15}N_2O_7P$).



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		

- Molecule 3 is N-{5-[5-(2-amino-4-oxo-3,4-dihydro-5H-pyrrolo[3,2-d]pyrimidin-5-yl)pentyl]thiophene-2-carbonyl}-L-glutamic acid (CCD ID: Y7L) (formula: $C_{21}H_{25}N_5O_6S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 66	C 42	N 10	O 12	S 2	0	1
3	B	1	Total 66	C 42	N 10	O 12	S 2	0	1

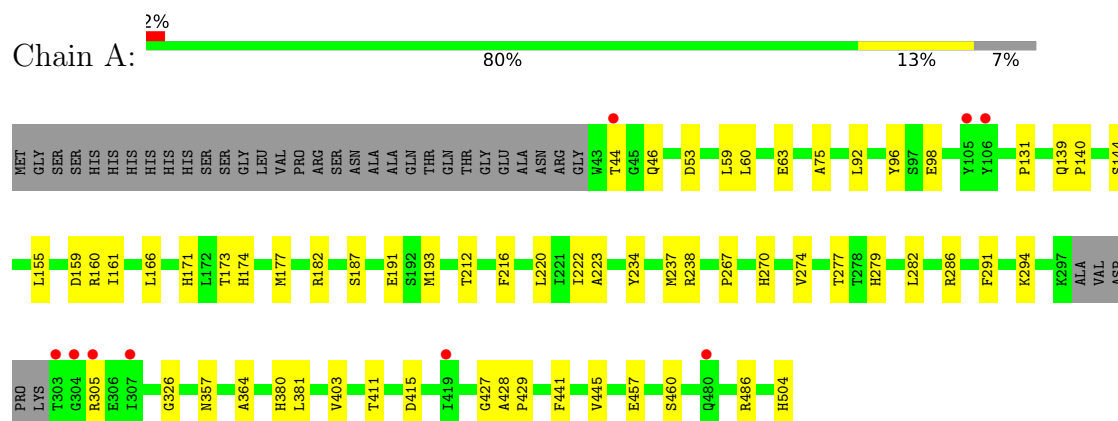
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	119	Total O 119 119	0	0
4	B	49	Total O 49 49	0	0

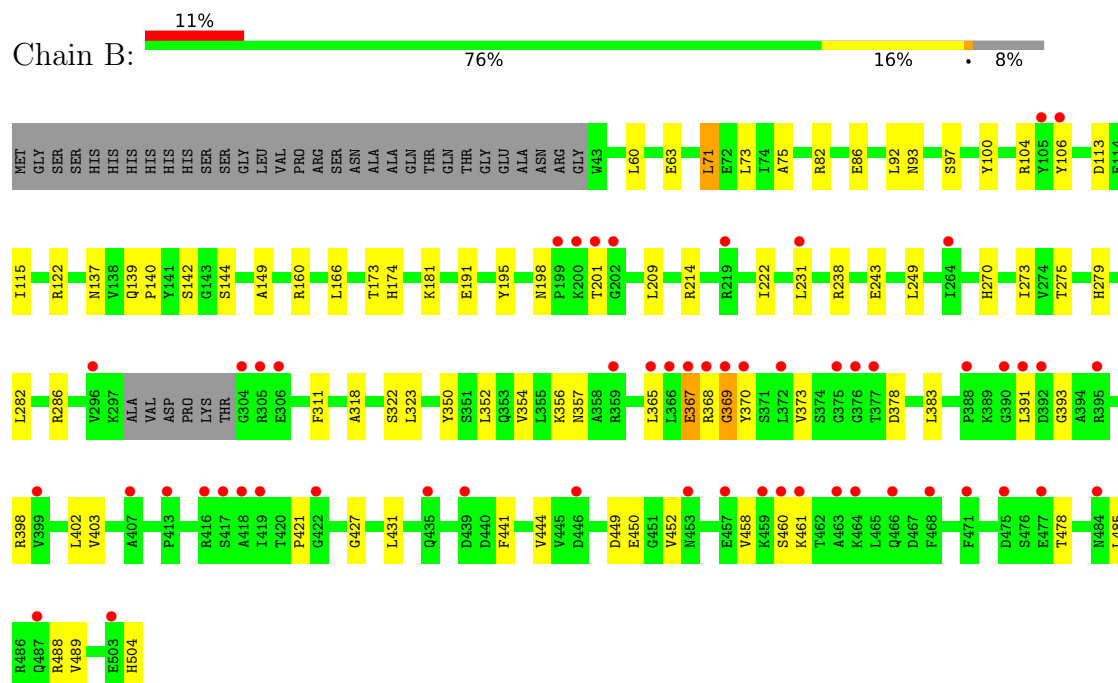
3 Residue-property plots [i](#)

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, mitochondrial



- Molecule 1: Serine hydroxymethyltransferase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	158.20Å 158.20Å 207.28Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.65 – 2.64 48.65 – 2.64	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.65-2.64) 99.9 (48.65-2.64)	Depositor EDS
R_{merge}	0.34	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.13-2998	Depositor
R, R_{free}	0.201 , 0.247 0.201 , 0.248	Depositor DCC
R_{free} test set	2243 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.389	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 64.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7327	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% 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: Y7L, 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	0.17	0/3619	0.41	0/4897
1	B	0.18	0/3509	0.40	0/4764
All	All	0.18	0/7128	0.40	0/9661

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

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	3548	0	3513	40	0
1	B	3439	0	3317	51	0
2	A	20	0	11	0	0
2	B	20	0	11	3	0
3	A	66	0	0	1	0
3	B	66	0	0	2	0
4	A	119	0	0	0	0
4	B	49	0	0	0	0
All	All	7327	0	6852	87	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 6.

All (87) 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:367:GLU:O	1:B:369:GLY:N	2.19	0.76
1:B:142:SER:OG	2:B:601:PLG:OP1	2.06	0.73
1:A:305:ARG:HD3	1:A:305:ARG:H	1.59	0.67
1:B:82:ARG:O	1:B:86:GLU:HG3	1.99	0.62
1:B:449:ASP:HA	1:B:452:VAL:HG12	1.81	0.62
1:A:277:THR:HG23	1:A:279:HIS:CE1	2.33	0.62
1:B:142:SER:HG	2:B:601:PLG:P	2.26	0.59
1:B:198:ASN:HB3	1:B:201:THR:HB	1.86	0.58
1:B:144:SER:HA	1:B:173:THR:HG21	1.85	0.57
1:A:59:LEU:HB3	1:B:115:ILE:HD11	1.87	0.57
1:A:174:HIS:CD2	1:A:222:ILE:HG21	2.41	0.56
1:B:485:LEU:O	1:B:489:VAL:HG23	2.06	0.56
1:A:279:HIS:ND1	1:A:286:ARG:HA	2.21	0.55
1:B:249:LEU:HD11	1:B:275:THR:HG23	1.88	0.55
1:A:457:GLU:O	1:A:460:SER:OG	2.26	0.53
1:B:181:LYS:NZ	3:B:602[B]:Y7L:O24	2.32	0.53
1:A:44:THR:HG23	1:A:46:GLN:H	1.74	0.52
1:B:231:LEU:HD13	1:B:378:ASP:HB2	1.92	0.52
1:B:365:LEU:O	1:B:370:TYR:HD2	1.92	0.52
1:B:403:VAL:O	1:B:489:VAL:HG11	2.09	0.51
1:B:97:SER:HA	1:B:106:TYR:HE1	1.75	0.51
1:A:161:ILE:HD13	1:A:220:LEU:HB3	1.92	0.51
1:B:71:LEU:HB3	1:B:431:LEU:HD11	1.92	0.51
1:B:75:ALA:HA	1:B:427:GLY:HA3	1.92	0.51
1:A:63:GLU:HG2	1:B:92:LEU:HD23	1.93	0.51
1:A:160:ARG:HA	1:A:191:GLU:O	2.11	0.50
1:A:60:LEU:HD23	1:A:504:HIS:HB2	1.92	0.50
1:A:144:SER:HA	1:A:173:THR:HG21	1.94	0.49
1:A:177:MET:HG3	1:A:187:SER:HB2	1.92	0.49
1:B:357:ASN:HB3	1:B:441:PHE:CD2	2.47	0.49
1:B:166:LEU:O	3:B:602[B]:Y7L:N11	2.46	0.49
1:B:238:ARG:HD3	1:B:270:HIS:O	2.13	0.48
1:B:60:LEU:HD23	1:B:504:HIS:HB2	1.96	0.48
1:B:279:HIS:ND1	1:B:286:ARG:HA	2.30	0.47
1:B:352:LEU:O	1:B:356:LYS:HG3	2.13	0.47
1:B:450:GLU:CD	1:B:488:ARG:HH12	2.20	0.47
1:A:403:VAL:HA	1:A:486:ARG:HB2	1.96	0.47
1:A:171:HIS:ND1	1:A:173:THR:OG1	2.43	0.47
1:A:223:ALA:HB2	1:A:237:MET:HE2	1.97	0.47
1:B:214:ARG:NH2	1:B:243:GLU:OE1	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:MET:HG2	1:A:182:ARG:HA	1.97	0.47
1:B:73:LEU:O	1:B:427:GLY:N	2.47	0.47
1:B:398:ARG:O	1:B:402:LEU:HD23	2.15	0.47
1:A:160:ARG:HB3	1:A:216:PHE:CE2	2.50	0.46
1:A:193:MET:HE1	1:A:212:THR:HB	1.97	0.46
1:A:411:THR:HG23	1:A:415:ASP:HB2	1.98	0.46
1:B:174:HIS:CD2	1:B:222:ILE:HG21	2.50	0.46
1:B:441:PHE:HA	1:B:444:VAL:HG22	1.97	0.46
1:B:350:TYR:O	1:B:354:VAL:HG23	2.16	0.46
1:B:391:LEU:HD23	1:B:452:VAL:HG23	1.98	0.46
1:A:53:ASP:OD2	1:B:122:ARG:NH1	2.45	0.45
1:B:458:VAL:CG2	1:B:478:THR:HG21	2.46	0.45
1:A:277:THR:HG23	1:A:279:HIS:ND1	2.31	0.45
1:A:357:ASN:HB3	1:A:441:PHE:CD2	2.52	0.45
1:A:282:LEU:HD23	1:A:282:LEU:HA	1.80	0.45
1:B:398:ARG:CZ	1:B:402:LEU:HD21	2.46	0.44
1:B:450:GLU:OE1	1:B:488:ARG:NH1	2.51	0.44
1:A:131:PRO:O	1:A:294:LYS:NZ	2.50	0.44
1:B:100:TYR:O	1:B:104:ARG:HG2	2.17	0.44
1:A:364:ALA:HB3	1:A:445:VAL:HG21	2.00	0.43
1:A:92:LEU:HD23	1:B:63:GLU:HG2	2.00	0.43
1:B:149:ALA:HA	1:B:323:LEU:HD12	2.00	0.42
1:A:155:LEU:HB3	1:A:159:ASP:HB2	2.02	0.42
1:B:460:SER:OG	1:B:461:LYS:N	2.52	0.42
1:B:160:ARG:HA	1:B:191:GLU:O	2.19	0.42
2:B:601:PLG:O3	2:B:601:PLG:N	2.49	0.42
1:A:44:THR:OG1	1:A:46:GLN:HG3	2.19	0.42
1:B:195:TYR:CD2	1:B:209:LEU:HD13	2.54	0.42
1:A:380:HIS:CE1	1:A:381:LEU:HD23	2.55	0.42
1:A:234:TYR:CZ	1:A:267:PRO:HB3	2.55	0.42
1:B:137:ASN:OD1	1:B:140:PRO:HD3	2.20	0.42
1:B:104:ARG:NE	1:B:113:ASP:OD1	2.44	0.42
1:A:428:ALA:N	1:A:429:PRO:CD	2.84	0.41
1:A:75:ALA:HA	1:A:427:GLY:HA3	2.02	0.41
1:A:139:GLN:N	1:A:140:PRO:CD	2.84	0.41
1:A:238:ARG:HD2	1:A:270:HIS:O	2.19	0.41
1:A:166:LEU:O	3:A:602[A]:Y7L:N11	2.53	0.41
1:B:282:LEU:HD23	1:B:282:LEU:HA	1.83	0.41
1:A:98:GLU:HG3	1:A:326:GLY:HA2	2.01	0.41
1:A:286:ARG:HG3	1:B:93:ASN:OD1	2.21	0.41
1:B:273:ILE:HD11	1:B:311:PHE:CG	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:274:VAL:HB	1:A:291:PHE:HB2	2.02	0.41
1:B:318:ALA:HA	1:B:322:SER:HB2	2.02	0.41
1:B:139:GLN:N	1:B:140:PRO:CD	2.84	0.41
1:B:393:GLY:N	1:B:421:PRO:O	2.43	0.40
1:B:373:VAL:N	1:B:383:LEU:O	2.52	0.40
1:A:294:LYS:HB3	1:A:294:LYS:HE2	1.89	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	453/493 (92%)	437 (96%)	16 (4%)	0	100	100
1	B	452/493 (92%)	428 (95%)	21 (5%)	3 (1%)	18	27
All	All	905/986 (92%)	865 (96%)	37 (4%)	3 (0%)	36	50

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	367	GLU
1	B	368	ARG
1	B	369	GLY

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	371/405 (92%)	370 (100%)	1 (0%)	86	93
1	B	348/405 (86%)	347 (100%)	1 (0%)	86	93
All	All	719/810 (89%)	717 (100%)	2 (0%)	86	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	96	TYR
1	B	71	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	207	ASN
1	A	353	GLN
1	A	357	ASN
1	A	466	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	PLG	B	601	-	20,20,20	2.55	3 (15%)	26,28,28	1.67	2 (7%)
3	Y7L	A	602[A]	-	34,35,35	2.55	9 (26%)	44,48,48	2.28	13 (29%)
3	Y7L	A	602[B]	-	34,35,35	2.61	9 (26%)	44,48,48	2.49	16 (36%)
3	Y7L	B	602[A]	-	34,35,35	2.57	9 (26%)	44,48,48	2.23	12 (27%)
3	Y7L	B	602[B]	-	34,35,35	2.58	9 (26%)	44,48,48	2.26	11 (25%)
2	PLG	A	601	-	20,20,20	2.66	3 (15%)	26,28,28	1.79	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
2	PLG	B	601	-	-	2/12/12/12	0/1/1/1
3	Y7L	A	602[A]	-	-	4/25/25/25	0/3/3/3
3	Y7L	A	602[B]	-	-	7/25/25/25	0/3/3/3
3	Y7L	B	602[A]	-	-	14/25/25/25	0/3/3/3
3	Y7L	B	602[B]	-	-	18/25/25/25	0/3/3/3
2	PLG	A	601	-	-	2/12/12/12	0/1/1/1

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602[B]	Y7L	O10-C4	8.85	1.40	1.23
3	B	602[B]	Y7L	O10-C4	8.82	1.40	1.23
3	B	602[A]	Y7L	O10-C4	8.82	1.40	1.23
3	A	602[A]	Y7L	O10-C4	8.75	1.40	1.23
2	B	601	PLG	C3-C2	7.50	1.48	1.41
2	A	601	PLG	C3-C2	7.33	1.48	1.41
2	A	601	PLG	C5-C4	6.27	1.49	1.40
2	A	601	PLG	C3-C4	6.26	1.49	1.40
3	A	602[B]	Y7L	C17-N19	6.02	1.46	1.34
3	B	602[B]	Y7L	C17-N19	5.83	1.46	1.34
3	B	602[A]	Y7L	C17-N19	5.75	1.46	1.34
2	B	601	PLG	C3-C4	5.72	1.48	1.40
2	B	601	PLG	C5-C4	5.71	1.48	1.40
3	A	602[A]	Y7L	C17-N19	5.68	1.46	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602[A]	Y7L	O25-C23	5.60	1.40	1.22
3	B	602[B]	Y7L	O25-C23	5.60	1.40	1.22
3	A	602[B]	Y7L	O25-C23	5.60	1.40	1.22
3	B	602[A]	Y7L	O25-C23	5.59	1.40	1.22
3	B	602[B]	Y7L	C2-N11	4.88	1.45	1.34
3	B	602[A]	Y7L	C2-N11	4.84	1.45	1.34
3	A	602[B]	Y7L	C2-N11	4.83	1.45	1.34
3	A	602[A]	Y7L	C2-N11	4.81	1.45	1.34
3	A	602[B]	Y7L	C4-N3	-3.17	1.32	1.38
3	A	602[A]	Y7L	C4-N3	-3.14	1.33	1.38
3	A	602[B]	Y7L	C17-C13	3.07	1.54	1.48
3	B	602[A]	Y7L	C4-N3	-3.05	1.33	1.38
3	B	602[B]	Y7L	C4-N3	-3.01	1.33	1.38
3	A	602[A]	Y7L	O24-C23	3.00	1.40	1.30
3	B	602[A]	Y7L	O24-C23	2.97	1.40	1.30
3	B	602[B]	Y7L	O24-C23	2.94	1.40	1.30
3	A	602[B]	Y7L	O24-C23	2.93	1.40	1.30
3	B	602[B]	Y7L	C17-C13	2.76	1.54	1.48
3	B	602[A]	Y7L	C17-C13	2.75	1.54	1.48
3	A	602[A]	Y7L	C17-C13	2.65	1.53	1.48
3	B	602[A]	Y7L	C6-N7	-2.44	1.32	1.38
3	B	602[B]	Y7L	C6-N7	-2.38	1.32	1.38
3	B	602[B]	Y7L	O18-C17	-2.37	1.19	1.23
3	A	602[B]	Y7L	O18-C17	-2.37	1.19	1.23
3	A	602[B]	Y7L	C6-N7	-2.36	1.32	1.38
3	B	602[A]	Y7L	O18-C17	-2.35	1.19	1.23
3	A	602[A]	Y7L	C6-N7	-2.33	1.32	1.38
3	A	602[A]	Y7L	O18-C17	-2.29	1.19	1.23

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602[B]	Y7L	C8-C9-N1	-6.77	118.32	124.81
3	A	602[A]	Y7L	C8-C9-N1	-6.75	118.34	124.81
3	B	602[B]	Y7L	C8-C9-N1	-6.66	118.42	124.81
3	B	602[A]	Y7L	C8-C9-N1	-6.59	118.50	124.81
2	A	601	PLG	C4A-N-CA	6.47	120.25	112.72
3	A	602[B]	Y7L	C16-S12-C13	6.37	100.25	92.63
2	B	601	PLG	C4A-N-CA	5.83	119.51	112.72
3	B	602[B]	Y7L	C16-S12-C13	5.74	99.50	92.63
3	B	602[A]	Y7L	C16-S12-C13	5.66	99.41	92.63
3	A	602[A]	Y7L	C16-S12-C13	5.26	98.93	92.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602[B]	Y7L	C13-C17-N19	5.15	125.82	116.86
3	A	602[B]	Y7L	C8-C4-N3	5.10	119.92	110.94
3	B	602[B]	Y7L	C8-C4-N3	5.10	119.91	110.94
3	A	602[A]	Y7L	C8-C4-N3	5.08	119.88	110.94
3	B	602[A]	Y7L	C8-C4-N3	5.02	119.78	110.94
3	A	602[A]	Y7L	C17-C13-S12	4.62	129.26	117.18
3	B	602[B]	Y7L	C9-C8-C4	-4.18	118.02	121.36
3	B	602[A]	Y7L	C9-C8-C4	-4.18	118.03	121.36
3	A	602[A]	Y7L	C9-C8-C4	-4.17	118.04	121.36
3	A	602[B]	Y7L	C9-C8-C4	-4.14	118.06	121.36
3	B	602[B]	Y7L	C17-C13-S12	4.00	127.63	117.18
3	A	602[B]	Y7L	C5-C9-N1	3.72	135.59	123.36
3	A	602[B]	Y7L	C14-C13-S12	-3.70	105.11	110.49
3	B	602[B]	Y7L	C5-C9-N1	3.67	135.43	123.36
3	A	602[A]	Y7L	C5-C9-N1	3.66	135.41	123.36
3	B	602[A]	Y7L	C5-C9-N1	3.65	135.39	123.36
3	B	602[A]	Y7L	C17-C13-S12	3.48	126.27	117.18
3	A	602[B]	Y7L	O18-C17-N19	-3.41	117.10	123.09
3	A	602[B]	Y7L	C9-N1-C2	3.25	121.72	116.63
3	A	602[A]	Y7L	C9-N1-C2	3.24	121.72	116.63
3	B	602[B]	Y7L	C9-N1-C2	3.20	121.65	116.63
3	B	602[B]	Y7L	C14-C13-S12	-3.19	105.86	110.49
3	B	602[A]	Y7L	C9-N1-C2	3.15	121.57	116.63
3	B	602[A]	Y7L	C14-C13-S12	-3.08	106.02	110.49
3	A	602[A]	Y7L	C14-C13-S12	-3.07	106.02	110.49
3	B	602[B]	Y7L	C2-N3-C4	-2.82	119.99	125.11
3	A	602[A]	Y7L	C33-C16-S12	2.82	128.24	120.24
3	A	602[B]	Y7L	C17-C13-S12	2.79	124.47	117.18
3	A	602[B]	Y7L	C2-N3-C4	-2.77	120.09	125.11
3	A	602[A]	Y7L	C2-N3-C4	-2.75	120.13	125.11
3	B	602[A]	Y7L	C2-N3-C4	-2.73	120.17	125.11
2	A	601	PLG	C6-N1-C2	2.72	124.13	119.20
3	A	602[B]	Y7L	C20-N19-C17	2.72	126.57	122.00
3	B	602[B]	Y7L	O10-C4-C8	-2.71	120.97	127.62
3	B	602[A]	Y7L	O10-C4-C8	-2.69	121.03	127.62
3	A	602[B]	Y7L	O10-C4-C8	-2.54	121.38	127.62
3	B	602[B]	Y7L	C33-C16-S12	2.52	127.39	120.24
3	A	602[A]	Y7L	O10-C4-C8	-2.52	121.43	127.62
3	B	602[A]	Y7L	C33-C16-S12	2.40	127.05	120.24
2	B	601	PLG	C6-N1-C2	2.33	123.43	119.20
3	A	602[B]	Y7L	C33-C16-S12	2.22	126.54	120.24
3	A	602[A]	Y7L	O27-C26-C20	2.15	120.79	113.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602[A]	Y7L	O27-C26-O28	-2.12	119.27	124.08
3	A	602[B]	Y7L	C15-C16-S12	-2.10	105.88	110.10
3	A	602[B]	Y7L	O27-C26-C20	2.07	120.50	113.51
3	B	602[A]	Y7L	O27-C26-O28	-2.05	119.42	124.08
2	A	601	PLG	C6-C5-C4	2.02	119.59	118.06

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	PLG	C3-C4-C4A-N
2	A	601	PLG	C5-C4-C4A-N
2	B	601	PLG	C3-C4-C4A-N
2	B	601	PLG	C5-C4-C4A-N
3	A	602[B]	Y7L	C13-C17-N19-C20
3	A	602[B]	Y7L	O18-C17-N19-C20
3	A	602[B]	Y7L	C26-C20-C21-C22
3	B	602[B]	Y7L	S12-C13-C17-O18
3	B	602[B]	Y7L	C26-C20-C21-C22
3	B	602[A]	Y7L	N19-C20-C21-C22
3	B	602[B]	Y7L	N19-C20-C21-C22
3	B	602[B]	Y7L	C31-C32-C33-C16
3	B	602[A]	Y7L	C20-C21-C22-C23
3	B	602[B]	Y7L	C14-C13-C17-O18
3	B	602[B]	Y7L	C14-C13-C17-N19
3	A	602[B]	Y7L	C30-C31-C32-C33
3	B	602[B]	Y7L	S12-C13-C17-N19
3	B	602[A]	Y7L	C26-C20-C21-C22
3	A	602[B]	Y7L	N19-C20-C21-C22
3	B	602[B]	Y7L	C30-C31-C32-C33
3	B	602[B]	Y7L	C29-C30-C31-C32
3	A	602[A]	Y7L	C30-C31-C32-C33
3	B	602[B]	Y7L	N19-C20-C26-O27
3	B	602[B]	Y7L	N19-C20-C26-O28
3	B	602[A]	Y7L	N7-C29-C30-C31
3	B	602[A]	Y7L	N19-C20-C26-O27
3	B	602[A]	Y7L	S12-C13-C17-N19
3	B	602[B]	Y7L	C21-C20-C26-O27
3	B	602[B]	Y7L	C21-C20-C26-O28
3	B	602[B]	Y7L	C15-C16-C33-C32
3	B	602[B]	Y7L	S12-C16-C33-C32
3	B	602[A]	Y7L	C14-C13-C17-N19

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Mol	Chain	Res	Type	Atoms
3	B	602[A]	Y7L	N19-C20-C26-O28
3	A	602[A]	Y7L	N19-C20-C26-O27
3	A	602[B]	Y7L	C21-C22-C23-O24
3	B	602[A]	Y7L	S12-C13-C17-O18
3	A	602[A]	Y7L	N19-C20-C26-O28
3	B	602[A]	Y7L	C14-C13-C17-O18
3	B	602[A]	Y7L	C21-C22-C23-O24
3	A	602[B]	Y7L	C21-C22-C23-O25
3	B	602[A]	Y7L	C21-C22-C23-O25
3	B	602[A]	Y7L	C26-C20-N19-C17
3	B	602[B]	Y7L	C21-C22-C23-O24
3	B	602[B]	Y7L	C21-C22-C23-O25
3	B	602[B]	Y7L	C26-C20-N19-C17
3	B	602[A]	Y7L	C29-C30-C31-C32
3	A	602[A]	Y7L	C20-C21-C22-C23

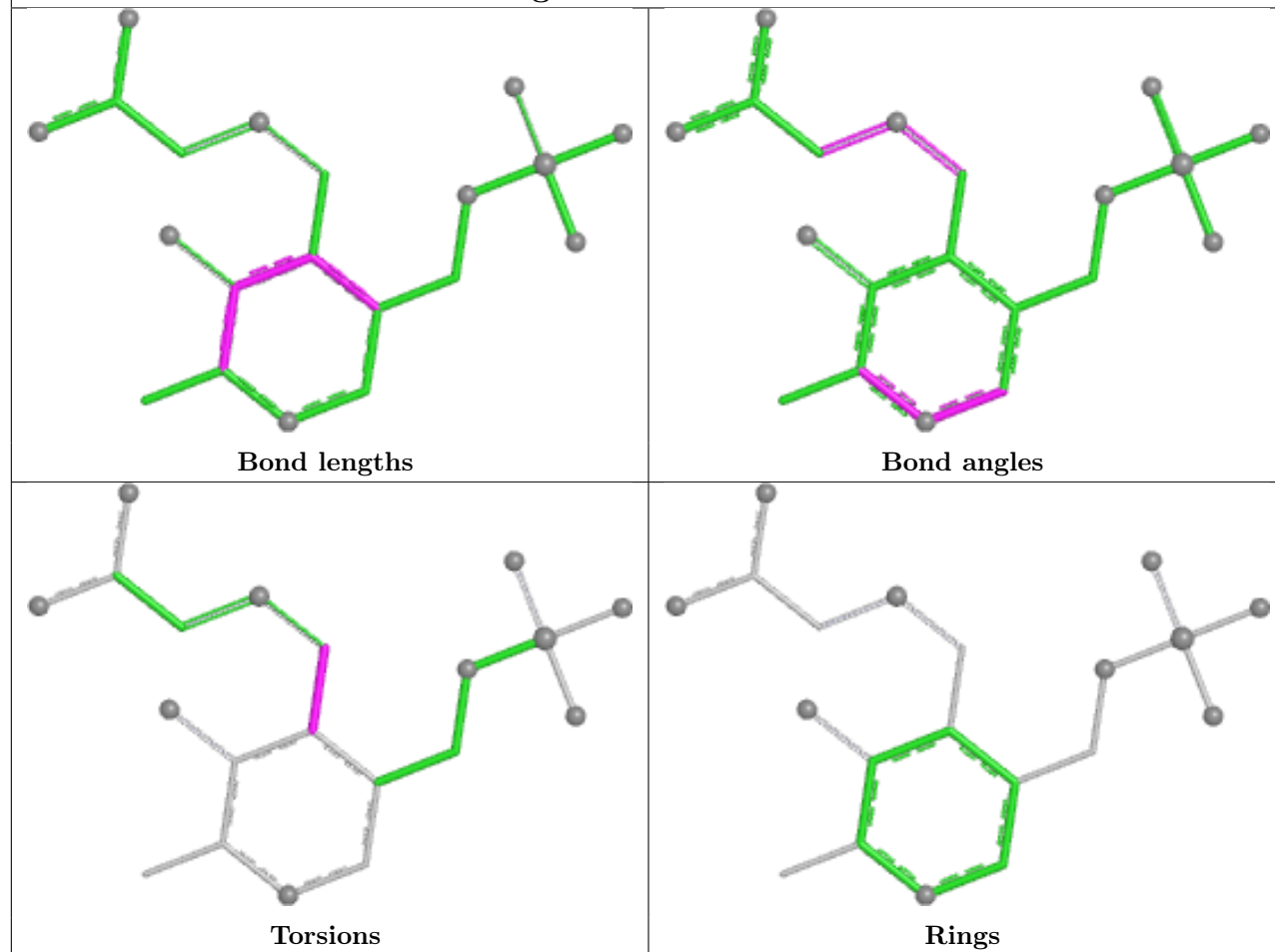
There are no ring outliers.

3 monomers are involved in 6 short contacts:

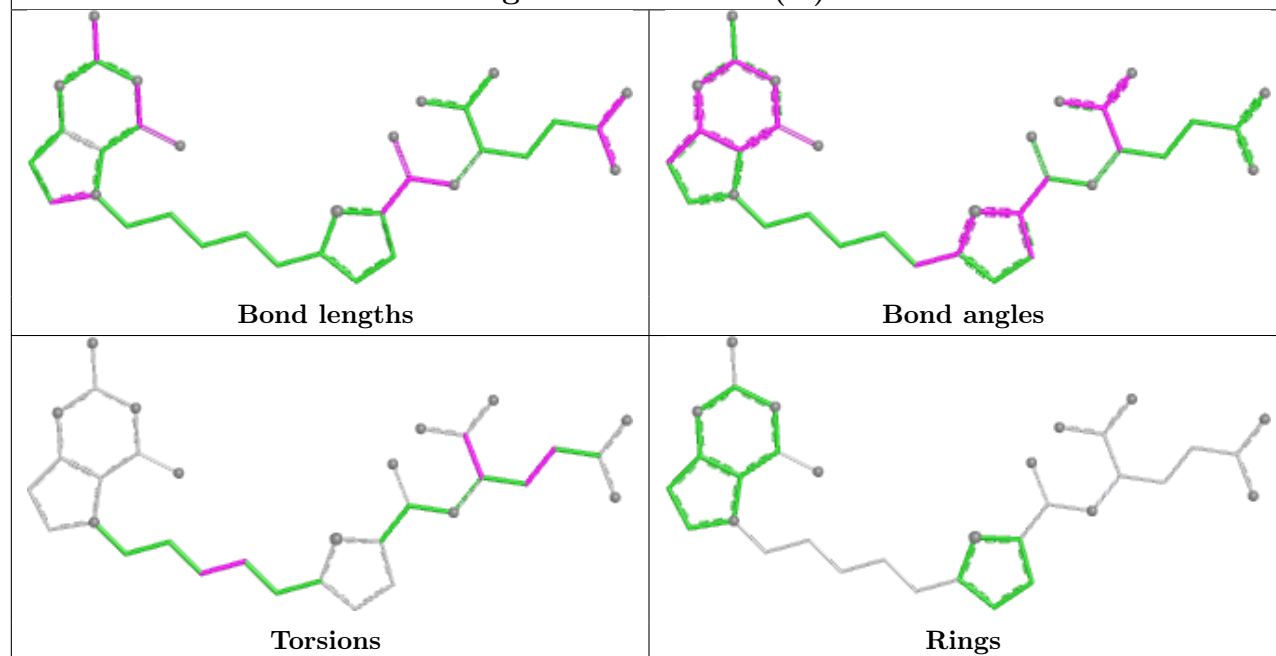
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	PLG	3	0
3	A	602[A]	Y7L	1	0
3	B	602[B]	Y7L	2	0

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.

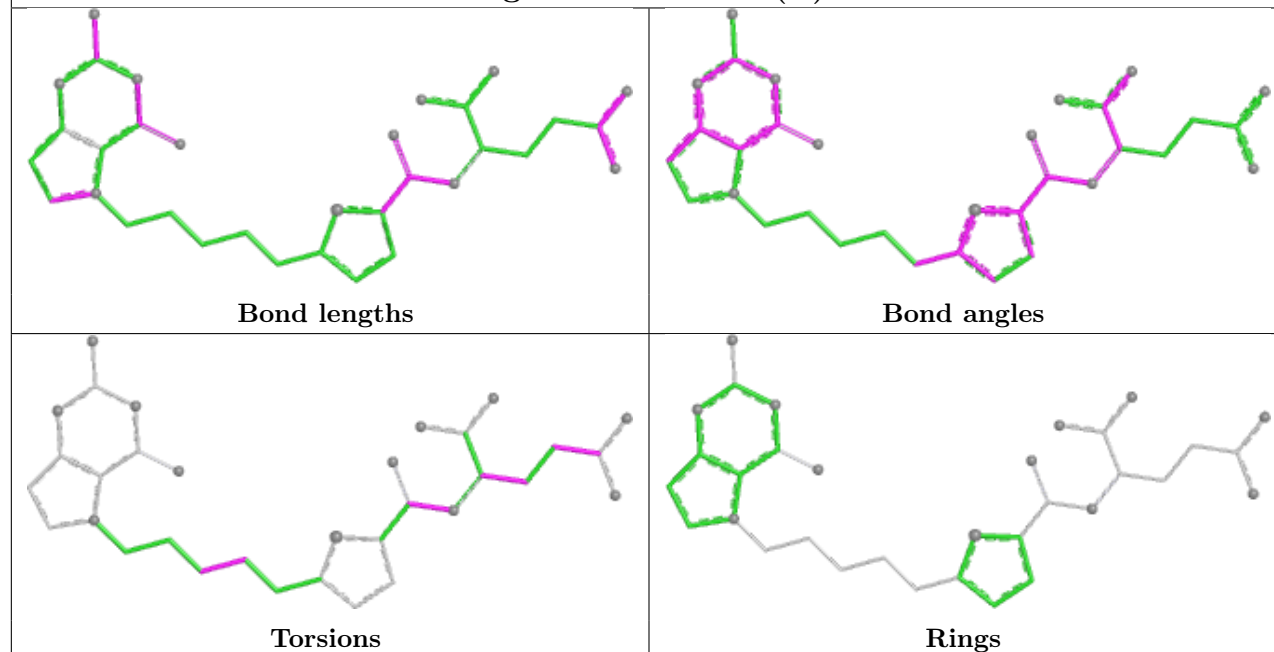
Ligand PLG B 601



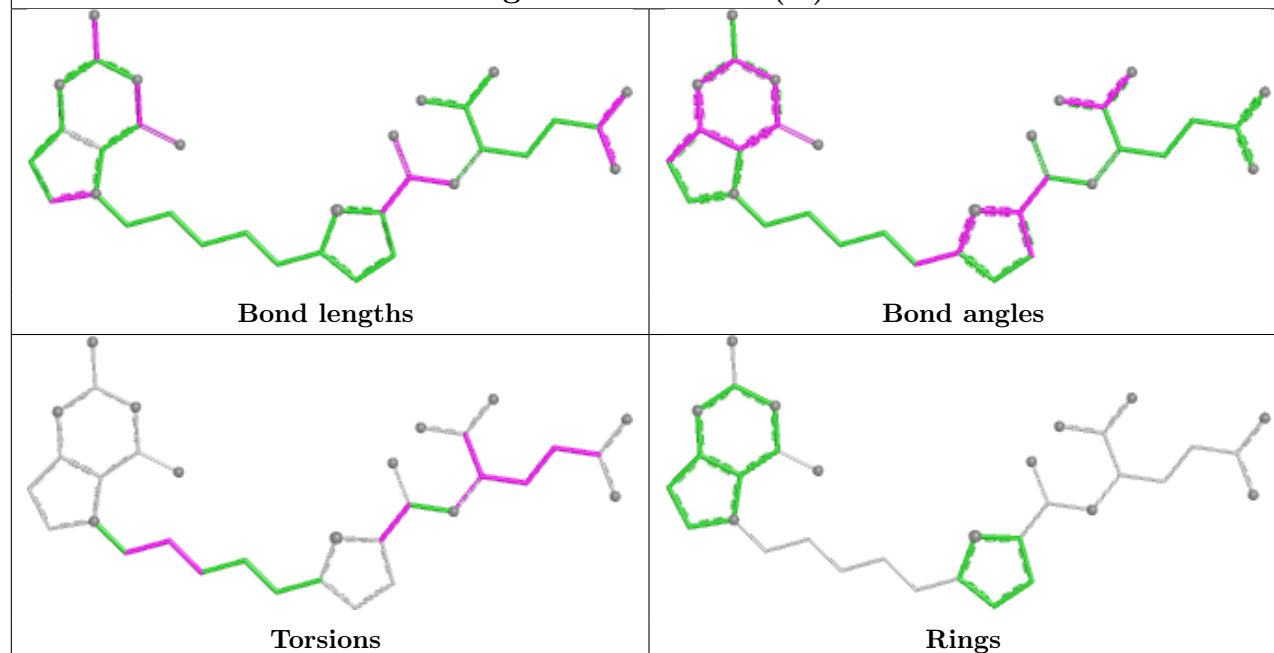
Ligand Y7L A 602 (A)



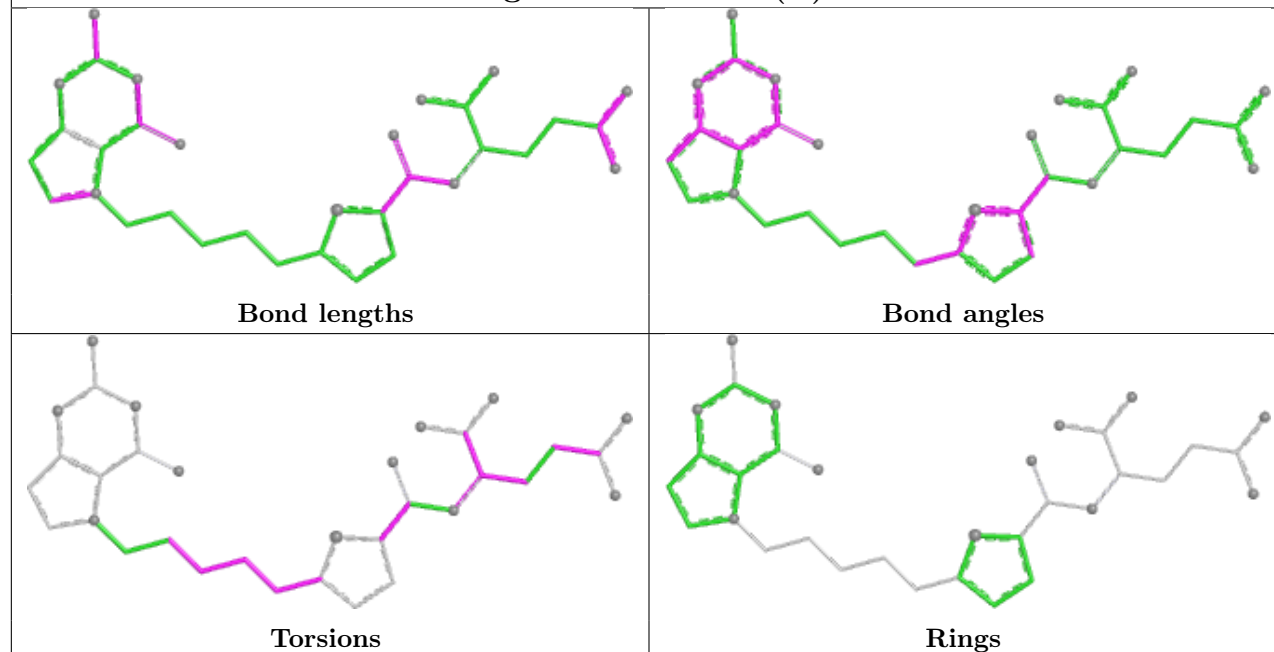
Ligand Y7L A 602 (B)



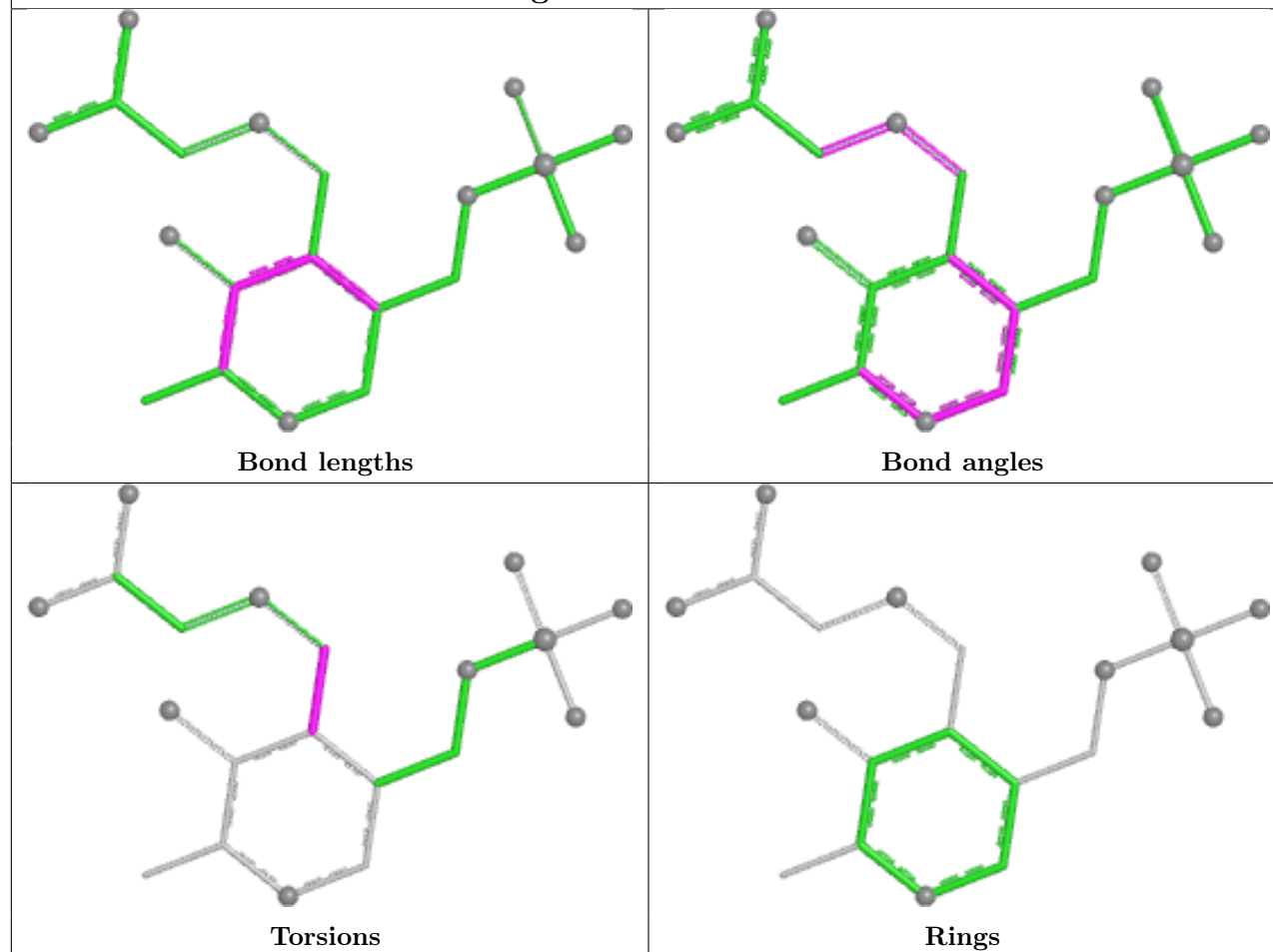
Ligand Y7L B 602 (A)



Ligand Y7L B 602 (B)



Ligand PLG A 601



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	457/493 (92%)	-0.29	9 (1%) 65 61	24, 42, 75, 126	0
1	B	456/493 (92%)	0.61	55 (12%) 8 7	29, 71, 113, 136	0
All	All	913/986 (92%)	0.16	64 (7%) 22 18	24, 54, 109, 136	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	GLY	6.8
1	A	303	THR	5.5
1	A	419	ILE	5.2
1	B	484	ASN	4.4
1	B	368	ARG	4.4
1	B	105	TYR	4.3
1	B	106	TYR	4.3
1	B	306	GLU	4.2
1	B	200	LYS	4.2
1	B	305	ARG	4.1
1	B	471	PHE	4.1
1	B	219	ARG	4.0
1	B	446	ASP	3.9
1	B	477	GLU	3.8
1	B	369	GLY	3.6
1	A	307	ILE	3.6
1	B	418	ALA	3.5
1	A	305	ARG	3.4
1	B	367	GLU	3.3
1	B	376	GLY	3.3
1	A	106	TYR	3.3
1	B	391	LEU	3.1
1	B	463	ALA	3.1
1	B	435	GLN	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	419	ILE	3.0
1	B	377	THR	3.0
1	B	457	GLU	3.0
1	B	459	LYS	2.9
1	A	105	TYR	2.9
1	B	370	TYR	2.9
1	B	399	VAL	2.8
1	B	460	SER	2.8
1	B	392	ASP	2.7
1	B	372	LEU	2.7
1	B	466	GLN	2.7
1	B	366	LEU	2.6
1	B	201	THR	2.5
1	B	464	LYS	2.5
1	A	304	GLY	2.5
1	B	375	GLY	2.5
1	B	468	PHE	2.5
1	B	202	GLY	2.5
1	B	365	LEU	2.5
1	B	416	ARG	2.4
1	B	199	PRO	2.4
1	B	461	LYS	2.4
1	B	487	GLN	2.4
1	B	503	GLU	2.4
1	B	413	PRO	2.3
1	B	388	PRO	2.3
1	B	407	ALA	2.3
1	B	422	GLY	2.3
1	B	453	ASN	2.2
1	B	417	SER	2.2
1	B	231	LEU	2.2
1	B	359	ARG	2.2
1	B	390	GLY	2.2
1	B	264	ILE	2.1
1	A	44	THR	2.1
1	B	296	VAL	2.1
1	B	475	ASP	2.0
1	A	480	GLN	2.0
1	B	439	ASP	2.0
1	B	395	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

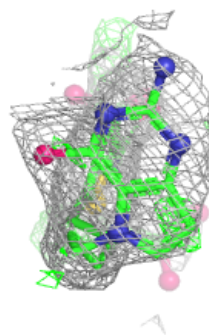
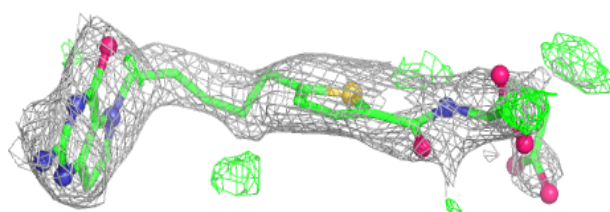
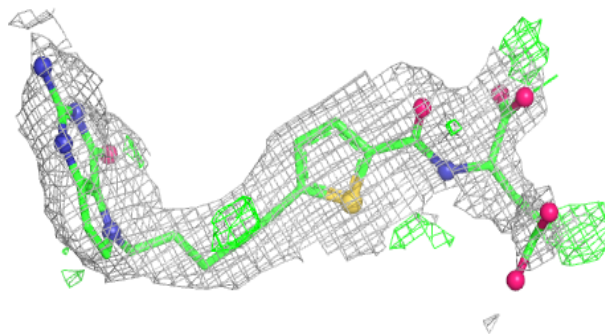
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	Y7L	B	602[A]	33/33	0.81	0.23	88,101,108,110	33
3	Y7L	B	602[B]	33/33	0.81	0.23	92,101,108,109	33
3	Y7L	A	602[A]	33/33	0.85	0.16	30,51,63,63	33
3	Y7L	A	602[B]	33/33	0.85	0.16	29,49,62,63	33
2	PLG	A	601	20/20	0.97	0.08	29,39,49,51	0
2	PLG	B	601	20/20	0.97	0.09	41,53,65,66	0

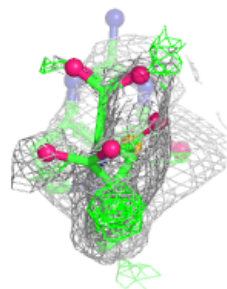
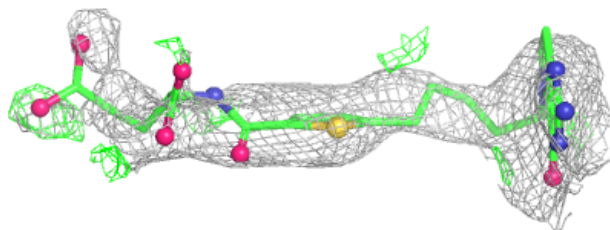
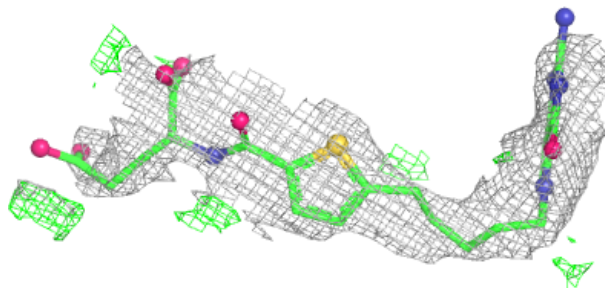
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Y7L B 602 (A):

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

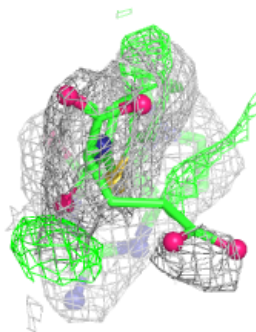
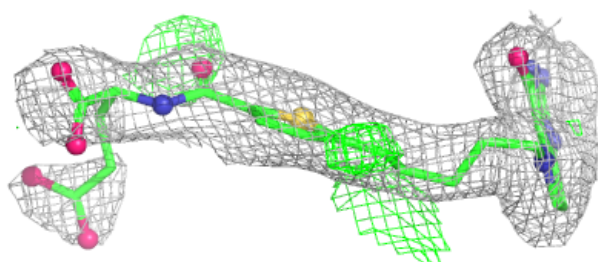
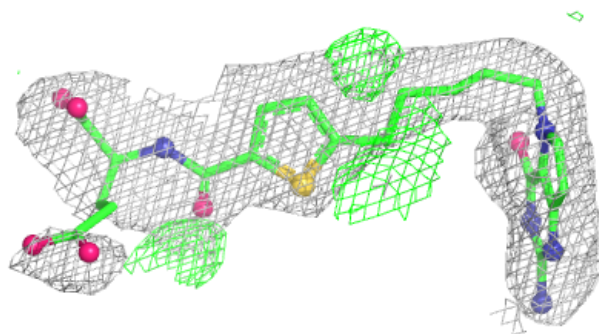
**Electron density around Y7L B 602 (B):**

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

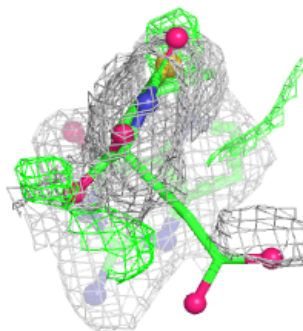
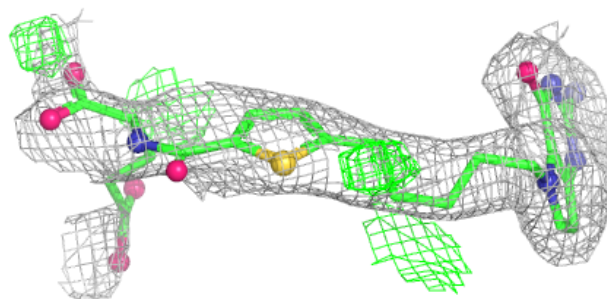
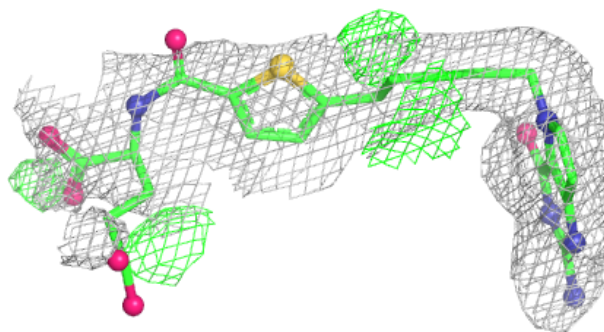


Electron density around Y7L A 602 (A):

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

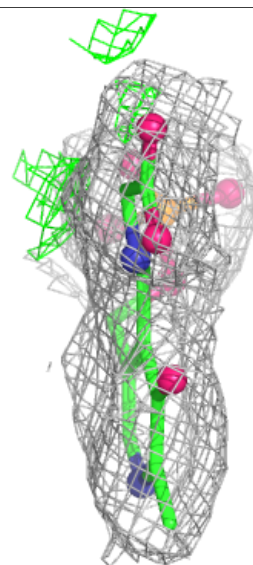
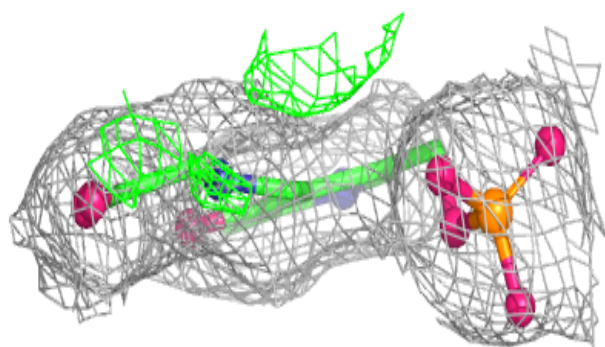
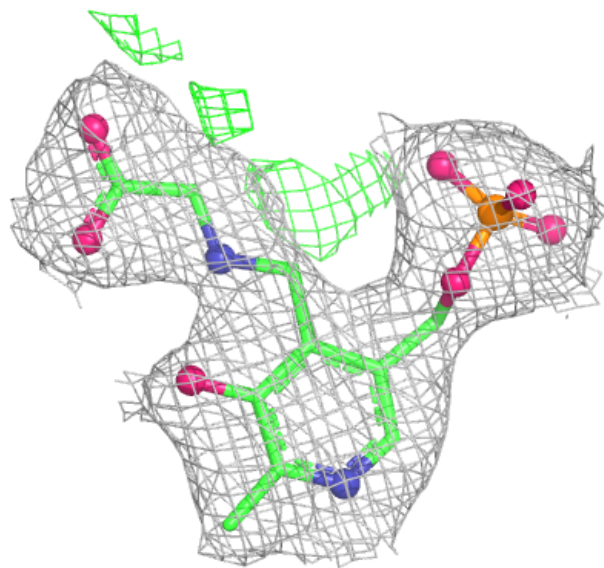
**Electron density around Y7L A 602 (B):**

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



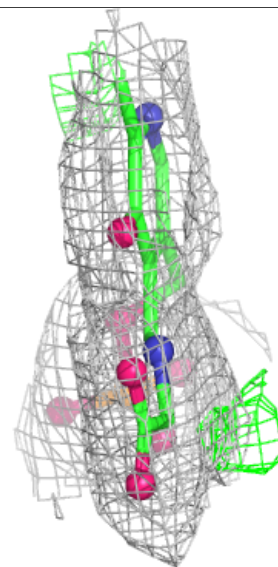
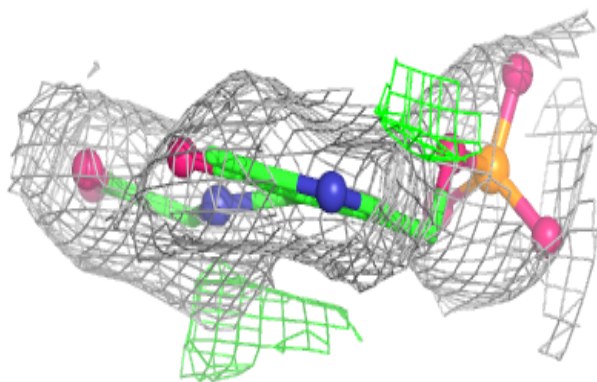
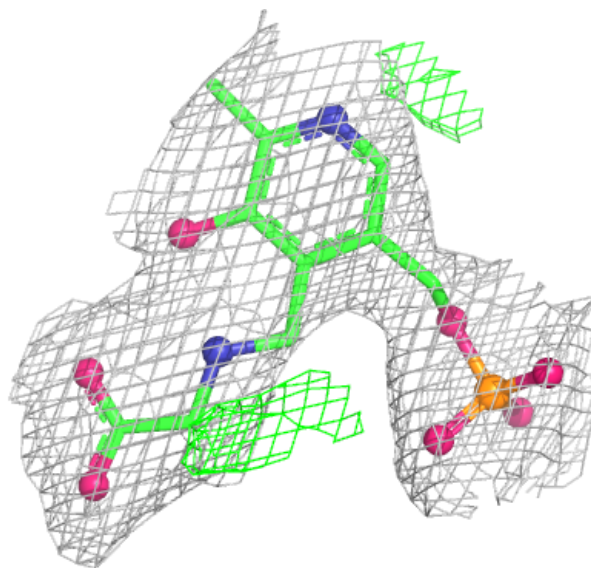
Electron density around PLG A 601:

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



Electron density around PLG B 601:

$2mF_o-DF_c$ (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.