



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2021 – 01:31 PM EDT

PDB ID : 2IER
Title : Crystal Structure of Aquifex aeolicus LpxC Complexed with Uridine 5'-Diphosphate
Authors : Gennadios, H.A.; Christianson, D.W.
Deposited on : 2006-09-19
Resolution : 2.70 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

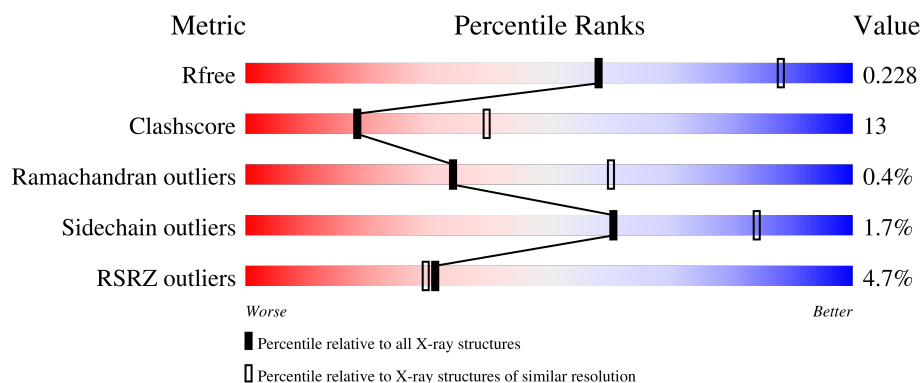
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	271	<div> <div>4%</div> <div>72%</div> <div>27%</div> </div>
1	B	271	<div> <div>5%</div> <div>71%</div> <div>27%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	GOL	A	501	-	X	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4492 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 UDP-3-O-[3-hydroxymyristoyl] N-acetylglucosamine deacetylase.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	0	0	0
			2157	1394	365	398			
1	B	267	Total	C	N	O	0	0	0
			2157	1394	365	398			

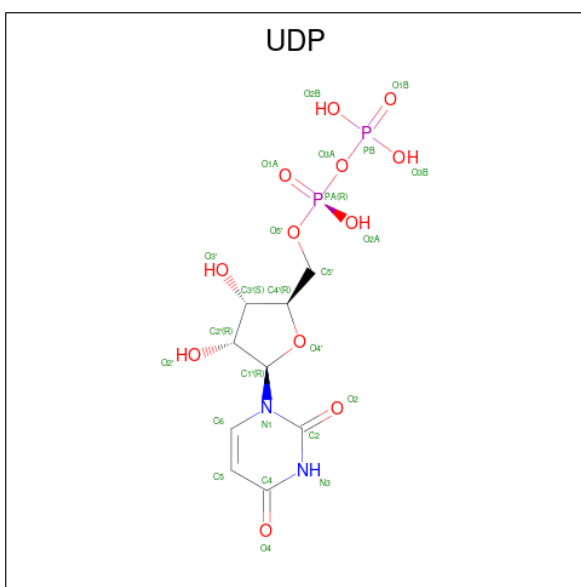
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	193	ALA	CYS	engineered mutation	UNP O67648
B	193	ALA	CYS	engineered mutation	UNP O67648

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

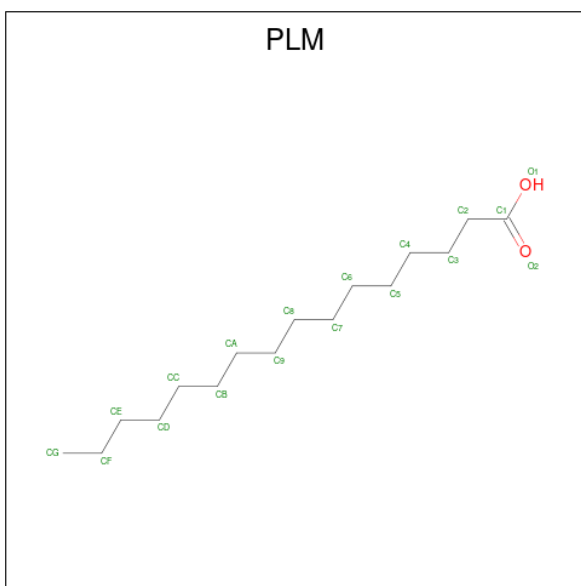
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	Zn	0	0
			5	5		
2	B	4	Total	Zn	0	0
			4	4		

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



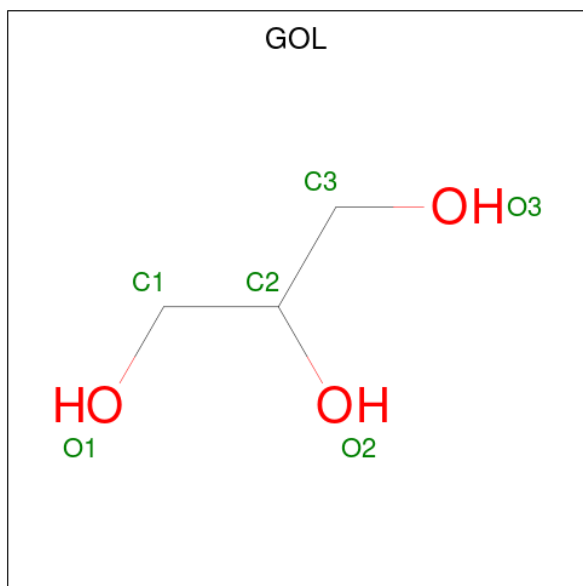
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 25	C 9	N 2	O 12	P 2	0	0
3	B	1	Total 25	C 9	N 2	O 12	P 2	0	0

- Molecule 4 is PALMITIC ACID (three-letter code: PLM) (formula: $\text{C}_{16}\text{H}_{32}\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			18	16	2		
4	B	1	Total	C	O	0	0
			18	16	2		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	42	Total	O	0	0
			42	42		
6	B	35	Total	O	0	0
			35	35		

4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	101.21Å 101.21Å 122.24Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.70 29.84 – 2.70	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.70) 99.9 (29.84-2.70)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.49 (at 2.68Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.213 , 0.241 0.195 , 0.228	Depositor DCC
R_{free} test set	1920 reflections (9.81%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtriage
Anisotropy	0.158	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 41.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.046 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4492	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% 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: ZN, UDP, PLM, GOL

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	0.38	0/2203	0.60	0/2966
1	B	0.39	0/2203	0.60	0/2966
All	All	0.38	0/4406	0.60	0/5932

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	2157	0	2176	60	0
1	B	2157	0	2176	59	0
2	A	5	0	0	0	0
2	B	4	0	0	0	0
3	A	25	0	11	2	0
3	B	25	0	11	1	0
4	A	18	0	30	2	0
4	B	18	0	31	2	0
5	A	6	0	4	2	0
6	A	42	0	0	2	0
6	B	35	0	0	0	0
All	All	4492	0	4439	118	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 13.

All (118) 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:133:GLU:H	1:B:280:LYS:HD3	1.22	1.01
1:A:133:GLU:H	1:A:280:LYS:HD3	1.22	1.00
1:A:133:GLU:N	1:A:280:LYS:HD3	1.97	0.79
1:B:133:GLU:N	1:B:280:LYS:HD3	1.97	0.78
1:B:158:GLU:HG3	1:B:172:LYS:HB3	1.66	0.78
1:B:216:LEU:HD21	1:B:229:LEU:HD22	1.68	0.76
1:A:216:LEU:HD21	1:A:229:LEU:HD22	1.66	0.76
1:A:158:GLU:HG3	1:A:172:LYS:HB3	1.66	0.75
1:A:255:LYS:HZ1	5:A:501:GOL:H11	1.53	0.73
1:A:216:LEU:HD21	1:A:229:LEU:CD2	2.20	0.71
3:A:801:UDP:H5'1	4:A:701:PLM:HG1	1.73	0.69
1:B:216:LEU:HD21	1:B:229:LEU:CD2	2.22	0.69
1:A:133:GLU:H	1:A:280:LYS:CD	2.04	0.66
1:B:145:ILE:HG21	1:B:271:LEU:HD22	1.78	0.65
1:B:133:GLU:H	1:B:280:LYS:CD	2.05	0.64
1:A:145:ILE:HG21	1:A:271:LEU:HD22	1.80	0.62
1:B:222:LYS:HE3	1:B:224:TYR:OH	2.01	0.61
1:A:222:LYS:HE3	1:A:224:TYR:OH	2.01	0.61
1:B:198:ILE:HD12	4:B:702:PLM:H82	1.83	0.61
1:B:26:LEU:HD11	1:B:101:ILE:HG23	1.83	0.60
1:A:26:LEU:HD11	1:A:101:ILE:HG23	1.83	0.59
1:B:167:PHE:HD2	1:B:221:ASP:HA	1.68	0.58
1:A:262:ARG:HG2	3:A:801:UDP:O4	2.05	0.57
1:A:144:LEU:C	1:A:144:LEU:HD23	2.25	0.57
1:A:167:PHE:HD2	1:A:221:ASP:HA	1.70	0.56
1:B:144:LEU:C	1:B:144:LEU:HD23	2.26	0.56
1:A:106:GLY:O	1:A:191:THR:HG23	2.06	0.55
1:A:197:GLU:HA	6:A:815:HOH:O	2.06	0.55
1:B:194:PHE:O	1:B:197:GLU:HG3	2.07	0.55
1:A:198:ILE:HD12	4:A:701:PLM:H71	1.88	0.55
1:A:37:ILE:CD1	1:A:92:VAL:HG22	2.37	0.54
1:A:194:PHE:O	1:A:197:GLU:HG3	2.07	0.54
1:A:255:LYS:NZ	5:A:501:GOL:H11	2.21	0.54
1:B:145:ILE:CG2	1:B:271:LEU:HD22	2.38	0.54
1:B:106:GLY:O	1:B:191:THR:HG23	2.08	0.54
1:A:134:GLU:HG3	1:A:278:LYS:NZ	2.23	0.53
1:A:145:ILE:CG2	1:A:271:LEU:HD22	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ILE:CD1	1:B:92:VAL:HG22	2.39	0.53
1:B:134:GLU:HG3	1:B:278:LYS:NZ	2.24	0.52
1:B:40:PHE:CZ	1:B:43:GLY:HA2	2.44	0.52
1:B:133:GLU:CD	1:B:280:LYS:HG3	2.31	0.51
1:A:132:VAL:HA	1:A:280:LYS:HD3	1.92	0.51
1:A:146:LYS:HE2	1:A:148:GLU:OE1	2.11	0.51
1:A:133:GLU:CD	1:A:280:LYS:HG3	2.31	0.51
1:A:40:PHE:CZ	1:A:43:GLY:HA2	2.46	0.50
1:B:5:LYS:HD3	1:B:120:ILE:HG21	1.93	0.50
1:A:37:ILE:HD12	1:A:92:VAL:HG22	1.92	0.50
1:B:146:LYS:HE2	1:B:148:GLU:OE1	2.11	0.50
1:B:132:VAL:HA	1:B:280:LYS:HD3	1.93	0.50
1:B:262:ARG:HG2	3:B:802:UDP:O4	2.11	0.50
1:B:277:LYS:C	1:B:279:GLN:H	2.16	0.49
1:A:277:LYS:C	1:A:279:GLN:H	2.16	0.49
1:A:278:LYS:O	1:A:280:LYS:N	2.46	0.49
1:A:133:GLU:HB2	1:A:280:LYS:HE2	1.95	0.48
1:B:154:GLU:OE2	1:B:257:LYS:HE3	2.14	0.48
1:A:5:LYS:HD3	1:A:120:ILE:HG21	1.94	0.48
1:B:37:ILE:HD12	1:B:92:VAL:HG22	1.95	0.48
1:B:194:PHE:HB2	1:B:197:GLU:HG2	1.95	0.48
1:B:130:PHE:CD2	1:B:250:LEU:HD13	2.49	0.48
1:B:278:LYS:O	1:B:280:LYS:N	2.47	0.47
1:A:7:VAL:HG23	1:A:30:PRO:HG3	1.95	0.47
1:B:155:VAL:HG21	1:B:244:ILE:HD13	1.96	0.47
1:A:112:TYR:O	1:A:116:ARG:HB2	2.15	0.47
1:A:194:PHE:HB2	1:A:197:GLU:HG2	1.95	0.47
1:A:130:PHE:CD2	1:A:250:LEU:HD13	2.49	0.47
1:A:154:GLU:OE2	1:A:257:LYS:HE3	2.14	0.47
1:A:155:VAL:HG21	1:A:244:ILE:HD13	1.96	0.47
1:B:133:GLU:HB2	1:B:280:LYS:HE2	1.96	0.47
1:B:112:TYR:O	1:B:116:ARG:HB2	2.15	0.47
1:A:26:LEU:C	1:A:27:ILE:HD12	2.35	0.46
1:B:129:TYR:CD2	1:B:253:PRO:HB2	2.51	0.46
1:B:210:GLY:HA3	4:B:702:PLM:H92	1.97	0.46
1:B:7:VAL:HG23	1:B:30:PRO:HG3	1.97	0.46
1:A:109:TRP:HB2	1:A:188:LEU:HD23	1.97	0.46
1:A:125:ARG:HG3	6:A:811:HOH:O	2.16	0.46
1:B:58:HIS:O	1:B:59:SER:HB3	2.16	0.46
1:A:20:THR:HB	1:A:22:GLU:HG3	1.99	0.45
1:A:124:ASN:HB3	1:B:97:ILE:CG2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:LEU:C	1:B:27:ILE:HD12	2.36	0.45
1:A:134:GLU:HG3	1:A:278:LYS:HZ3	1.81	0.45
1:B:20:THR:HB	1:B:22:GLU:HG3	1.98	0.45
1:B:109:TRP:HB2	1:B:188:LEU:HD23	1.98	0.44
1:B:191:THR:HG22	1:B:238:HIS:CE1	2.52	0.44
1:A:27:ILE:HD12	1:A:27:ILE:N	2.33	0.44
1:B:5:LYS:HD3	1:B:120:ILE:CG2	2.47	0.44
1:A:58:HIS:O	1:A:59:SER:HB3	2.18	0.44
1:B:129:TYR:CE2	1:B:255:LYS:HE2	2.53	0.44
1:A:129:TYR:CE2	1:A:255:LYS:HE2	2.52	0.44
1:A:129:TYR:CD2	1:A:253:PRO:HB2	2.53	0.43
1:A:191:THR:HG22	1:A:238:HIS:CE1	2.53	0.43
1:A:278:LYS:O	1:A:279:GLN:C	2.56	0.43
1:A:5:LYS:HD3	1:A:120:ILE:CG2	2.48	0.43
1:A:61:ASP:OD1	1:A:75:LYS:HA	2.19	0.43
1:B:278:LYS:O	1:B:279:GLN:C	2.56	0.43
1:A:216:LEU:HD11	1:A:229:LEU:HD23	2.01	0.42
1:B:134:GLU:H	1:B:280:LYS:HZ1	1.66	0.42
1:B:216:LEU:HD11	1:B:229:LEU:HD23	2.00	0.42
1:A:129:TYR:CD2	1:A:255:LYS:HE2	2.55	0.42
1:A:183:GLU:H	1:A:183:GLU:CD	2.23	0.42
1:B:183:GLU:H	1:B:183:GLU:CD	2.23	0.42
1:A:126:GLU:OE1	1:B:27:ILE:HG12	2.20	0.42
1:B:27:ILE:HD12	1:B:27:ILE:N	2.35	0.42
1:B:199:GLU:O	1:B:203:LYS:HB2	2.20	0.42
1:A:195:ASP:OD1	1:A:196:TRP:N	2.53	0.41
1:A:184:GLU:HA	1:A:187:VAL:HG22	2.02	0.41
1:A:265:HIS:O	1:A:269:VAL:HG23	2.21	0.41
1:B:184:GLU:HA	1:B:187:VAL:HG22	2.02	0.41
1:A:157:TYR:O	1:A:172:LYS:HA	2.20	0.41
1:A:199:GLU:O	1:A:203:LYS:HB2	2.20	0.41
1:B:134:GLU:HG3	1:B:278:LYS:HZ3	1.85	0.41
1:B:129:TYR:CD2	1:B:255:LYS:HE2	2.55	0.41
1:B:60:THR:O	1:B:77:VAL:HG22	2.21	0.41
1:B:157:TYR:O	1:B:172:LYS:HA	2.20	0.41
1:B:10:LYS:HD2	1:B:29:HIS:CE1	2.55	0.40
1:B:79:HIS:HE1	1:B:191:THR:CG2	2.35	0.40
1:B:195:ASP:OD1	1:B:196:TRP:N	2.54	0.40
1:B:61:ASP:OD1	1:B:75:LYS:HA	2.21	0.40
1:A:222:LYS:HE3	1:A:224:TYR:CZ	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	265/271 (98%)	250 (94%)	14 (5%)	1 (0%)	34	60
1	B	265/271 (98%)	249 (94%)	15 (6%)	1 (0%)	34	60
All	All	530/542 (98%)	499 (94%)	29 (6%)	2 (0%)	34	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	279	GLN
1	B	279	GLN

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	234/238 (98%)	230 (98%)	4 (2%)	60	84
1	B	234/238 (98%)	230 (98%)	4 (2%)	60	84
All	All	468/476 (98%)	460 (98%)	8 (2%)	60	84

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	232	GLU
1	A	248	TYR
1	A	274	GLU
1	A	279	GLN

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Mol	Chain	Res	Type
1	B	232	GLU
1	B	248	TYR
1	B	274	GLU
1	B	279	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 9 are monoatomic - leaving 5 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	UDP	B	802	2	20,26,26	1.70	3 (15%)	25,40,40	1.47	3 (12%)
5	GOL	A	501	-	5,5,5	4.53	5 (100%)	5,5,5	5.79	3 (60%)
4	PLM	A	701	-	14,17,17	0.50	0	13,17,17	1.62	2 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PLM	B	702	-	14,17,17	0.47	0	13,17,17	1.61	2 (15%)
3	UDP	A	801	2	20,26,26	1.72	4 (20%)	25,40,40	1.50	4 (16%)

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	UDP	B	802	2	-	8/14/32/32	0/2/2/2
5	GOL	A	501	-	-	2/4/4/4	-
4	PLM	A	701	-	-	8/13/15/15	-
4	PLM	B	702	-	-	9/13/15/15	-
3	UDP	A	801	2	-	6/14/32/32	0/2/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	501	GOL	C3-C2	-7.17	1.22	1.51
5	A	501	GOL	O1-C1	4.48	1.61	1.42
3	B	802	UDP	C6-N1	4.18	1.41	1.35
3	A	801	UDP	C6-N1	4.18	1.41	1.35
3	A	801	UDP	C4-N3	3.94	1.39	1.33
3	B	802	UDP	C4-N3	3.71	1.39	1.33
5	A	501	GOL	O3-C3	3.61	1.57	1.42
3	B	802	UDP	C3'-C4'	3.22	1.61	1.53
5	A	501	GOL	O2-C2	-3.10	1.34	1.43
3	A	801	UDP	C3'-C4'	3.00	1.60	1.53
5	A	501	GOL	C1-C2	-2.88	1.39	1.51
3	A	801	UDP	O4'-C1'	2.04	1.43	1.41

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	501	GOL	O3-C3-C2	10.56	160.84	110.20
5	A	501	GOL	O2-C2-C3	6.79	139.03	109.12
4	B	702	PLM	CF-CE-CD	4.88	160.63	115.30
4	A	701	PLM	CF-CE-CD	4.86	160.40	115.30
3	A	801	UDP	O3'-C3'-C4'	-4.44	98.22	111.05
3	B	802	UDP	O3'-C3'-C4'	-4.32	98.56	111.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	501	GOL	O1-C1-C2	3.16	125.36	110.20
3	B	802	UDP	C5'-C4'-C3'	-3.03	103.81	115.18
3	A	801	UDP	C5'-C4'-C3'	-2.99	103.96	115.18
4	B	702	PLM	CG-CF-CE	2.75	134.31	113.42
4	A	701	PLM	CG-CF-CE	2.74	134.25	113.42
3	B	802	UDP	C5-C4-N3	-2.10	118.70	123.31
3	A	801	UDP	O3B-PB-O3A	2.07	111.56	104.64
3	A	801	UDP	C5-C4-N3	-2.06	118.77	123.31

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	801	UDP	C5'-O5'-PA-O1A
3	A	801	UDP	C5'-O5'-PA-O2A
3	B	802	UDP	C5'-O5'-PA-O1A
3	B	802	UDP	PA-O3A-PB-O3B
5	A	501	GOL	C1-C2-C3-O3
3	A	801	UDP	C3'-C4'-C5'-O5'
3	B	802	UDP	C3'-C4'-C5'-O5'
3	A	801	UDP	O4'-C4'-C5'-O5'
3	B	802	UDP	O4'-C4'-C5'-O5'
4	B	702	PLM	CA-CB-CC-CD
4	B	702	PLM	C8-C9-CA-CB
4	A	701	PLM	C6-C7-C8-C9
4	A	701	PLM	CB-CC-CD-CE
4	B	702	PLM	CC-CD-CE-CF
4	A	701	PLM	C7-C8-C9-CA
4	B	702	PLM	C5-C6-C7-C8
4	A	701	PLM	C3-C4-C5-C6
4	A	701	PLM	C9-CA-CB-CC
5	A	501	GOL	O1-C1-C2-O2
4	B	702	PLM	C2-C3-C4-C5
4	B	702	PLM	C6-C7-C8-C9
4	A	701	PLM	CC-CD-CE-CF
4	B	702	PLM	C9-CA-CB-CC
3	A	801	UDP	C4'-C5'-O5'-PA
3	B	802	UDP	C5'-O5'-PA-O3A
3	B	802	UDP	C5'-O5'-PA-O2A
4	B	702	PLM	C3-C4-C5-C6
4	A	701	PLM	C5-C6-C7-C8
3	B	802	UDP	C4'-C5'-O5'-PA

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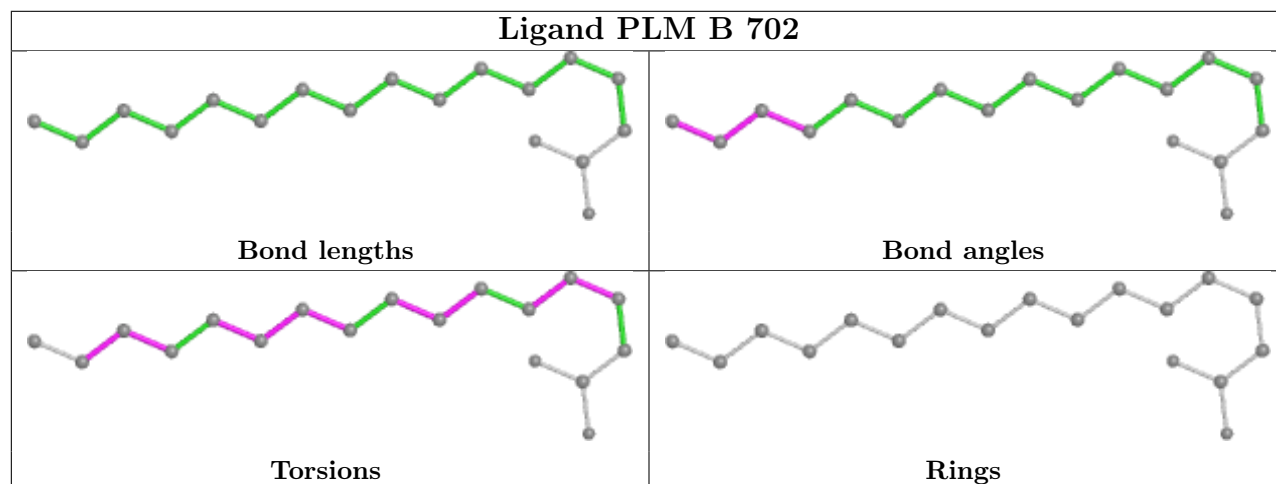
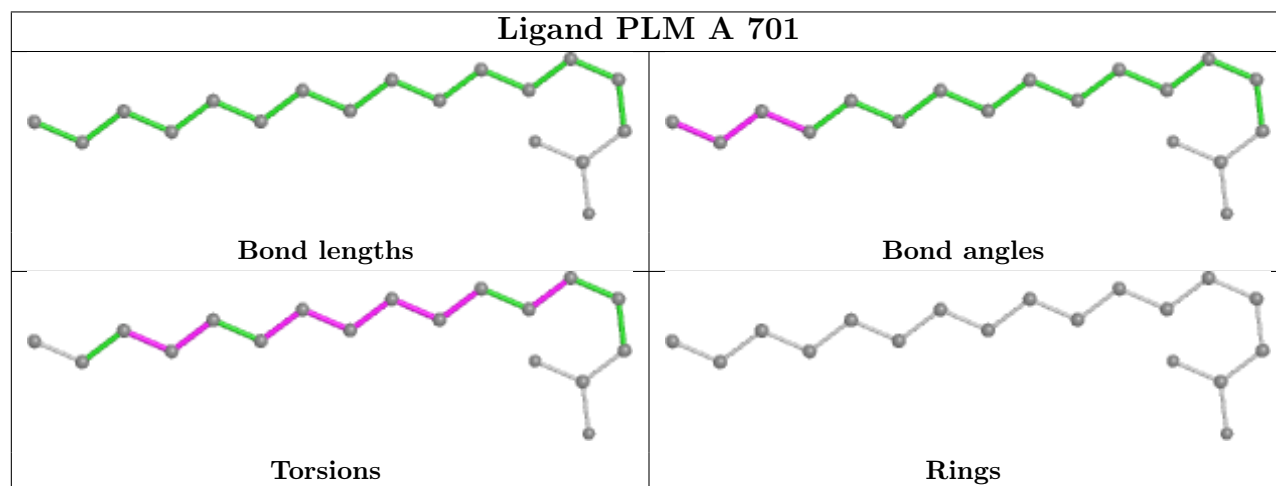
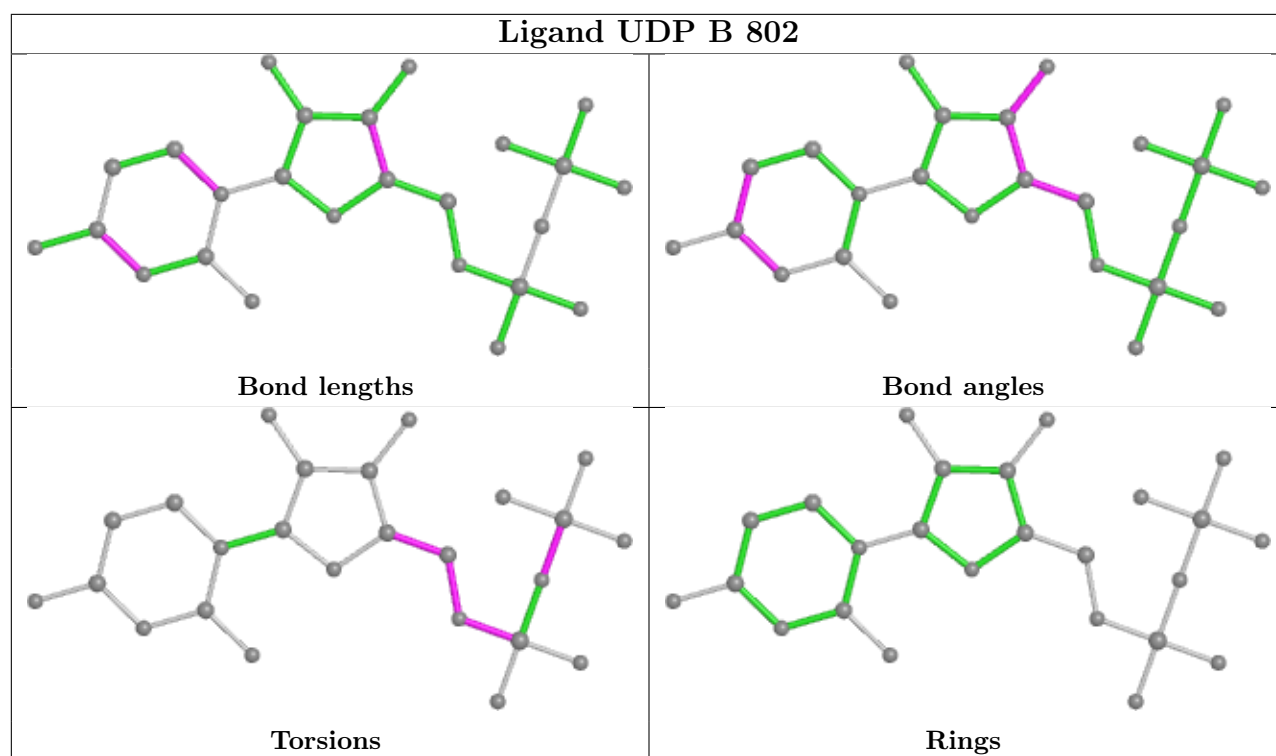
Mol	Chain	Res	Type	Atoms
4	A	701	PLM	C8-C9-CA-CB
3	A	801	UDP	C5'-O5'-PA-O3A
4	B	702	PLM	CD-CE-CF-CG
3	B	802	UDP	PA-O3A-PB-O1B

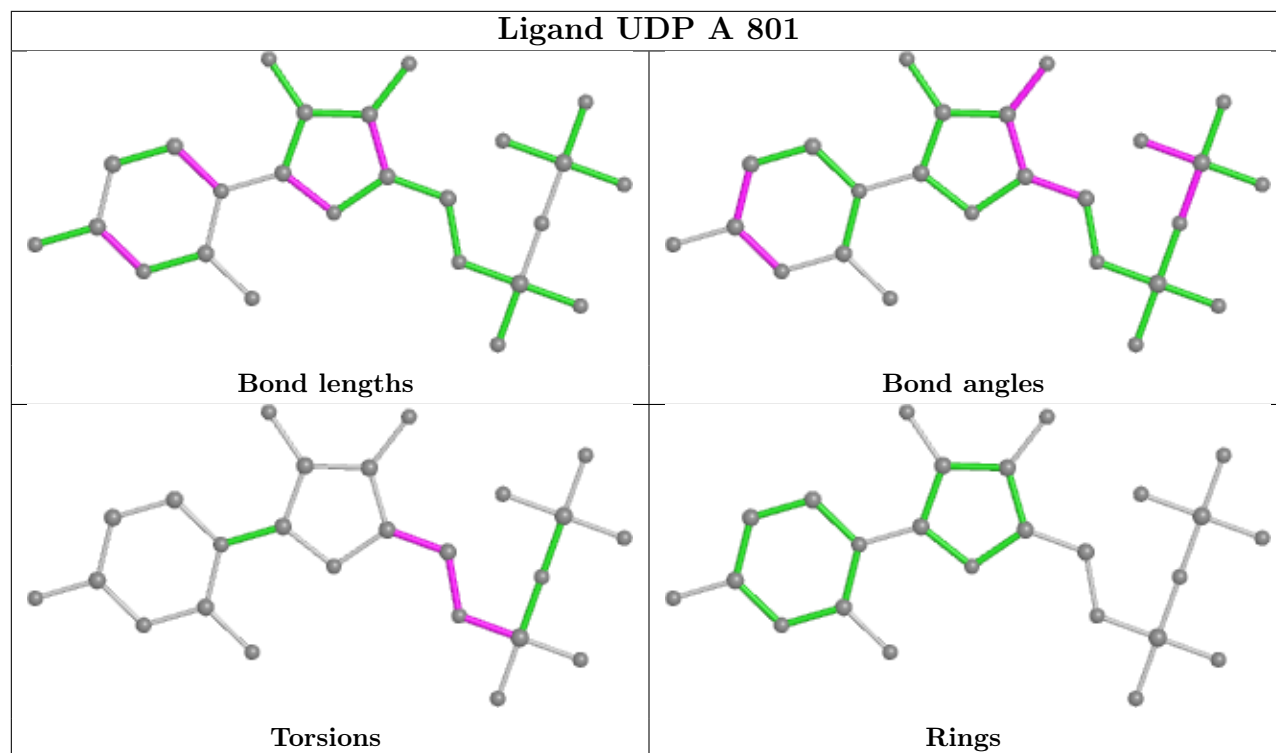
There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	802	UDP	1	0
5	A	501	GOL	2	0
4	A	701	PLM	2	0
4	B	702	PLM	2	0
3	A	801	UDP	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	267/271 (98%)	0.23	12 (4%) 33 31	7, 23, 40, 64	6 (2%)
1	B	267/271 (98%)	0.25	13 (4%) 29 28	7, 23, 41, 65	6 (2%)
All	All	534/542 (98%)	0.24	25 (4%) 31 30	7, 23, 41, 65	12 (2%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	141	GLU	4.8
1	B	151	ASP	4.3
1	B	124	ASN	4.1
1	A	23	TYR	4.0
1	B	190	ARG	4.0
1	B	141	GLU	3.7
1	B	279	GLN	3.5
1	A	279	GLN	3.1
1	A	220	LYS	3.0
1	A	190	ARG	3.0
1	A	73	ARG	2.8
1	A	32	LYS	2.7
1	B	280	LYS	2.6
1	B	23	TYR	2.5
1	B	113	GLU	2.4
1	A	135	PRO	2.4
1	B	12	SER	2.4
1	A	142	GLY	2.3
1	A	113	GLU	2.3
1	B	123	GLN	2.2
1	A	139	GLU	2.1
1	B	73	ARG	2.1
1	A	280	LYS	2.1
1	B	142	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	3	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

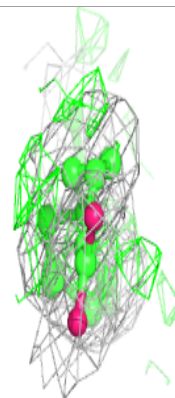
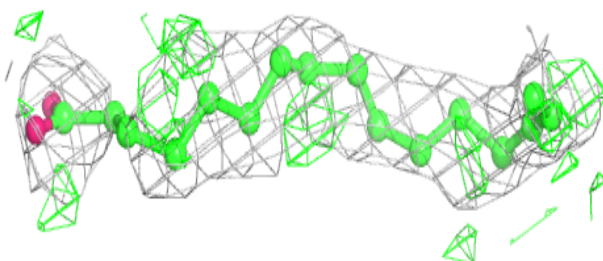
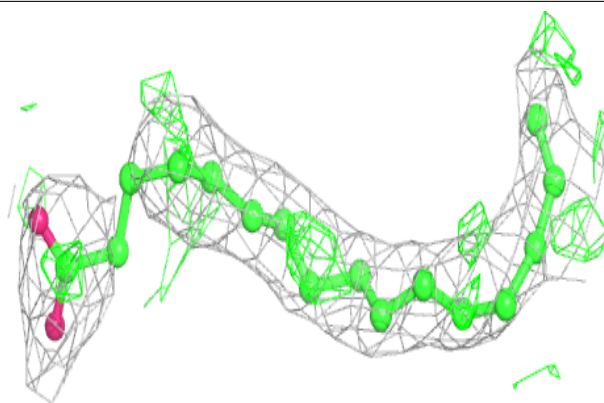
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	A	608	1/1	0.59	0.32	29,29,29,29	1
2	ZN	B	609	1/1	0.66	0.19	27,27,27,27	1
2	ZN	B	604	1/1	0.78	0.19	39,39,39,39	1
4	PLM	A	701	18/18	0.83	0.54	10,12,21,22	18
5	GOL	A	501	6/6	0.85	0.28	28,37,39,41	0
2	ZN	B	607	1/1	0.86	0.34	36,36,36,36	1
4	PLM	B	702	18/18	0.88	0.52	10,14,25,25	18
3	UDP	B	802	25/25	0.89	0.21	48,50,52,53	0
2	ZN	A	606	1/1	0.90	0.12	26,26,26,26	1
2	ZN	A	603	1/1	0.91	0.16	21,21,21,21	1
3	UDP	A	801	25/25	0.92	0.19	39,42,46,47	0
2	ZN	A	605	1/1	0.97	0.05	27,27,27,27	0
2	ZN	B	602	1/1	0.98	0.04	27,27,27,27	0
2	ZN	A	601	1/1	0.99	0.07	26,26,26,26	0

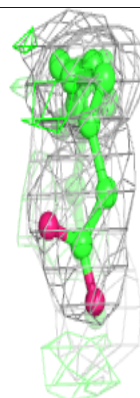
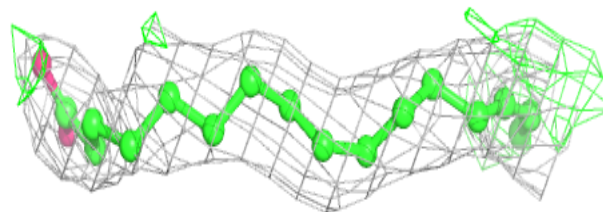
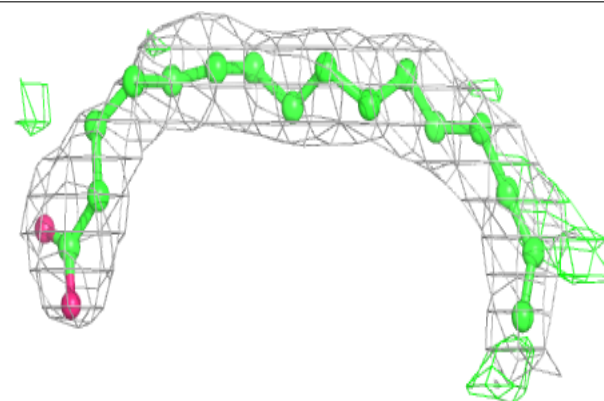
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PLM 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)

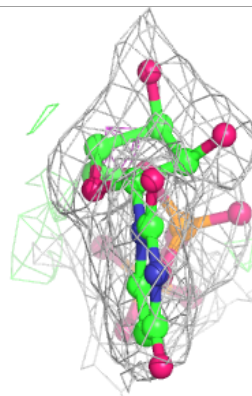
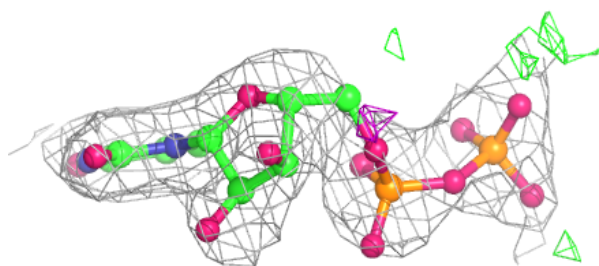
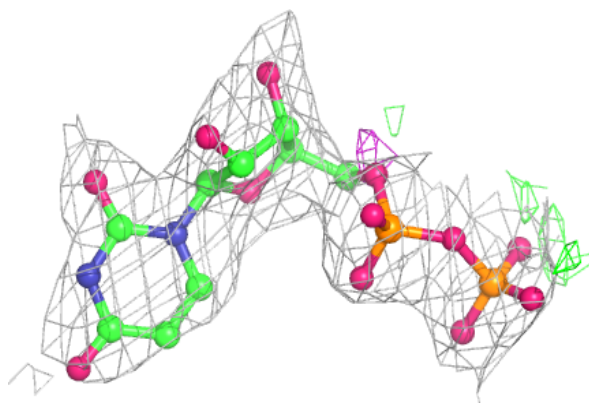
**Electron density around PLM B 702:**

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

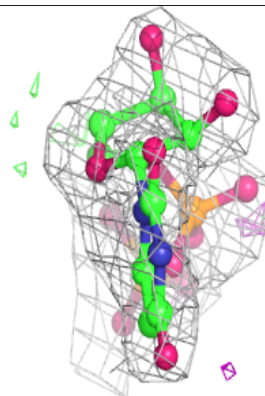
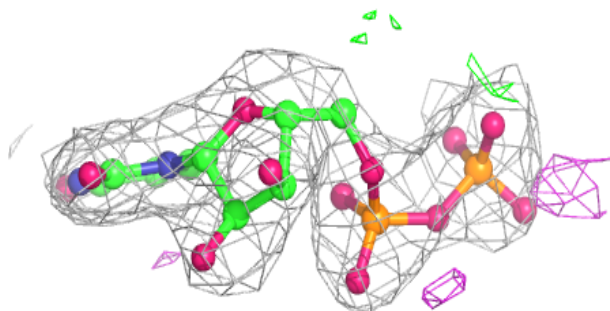
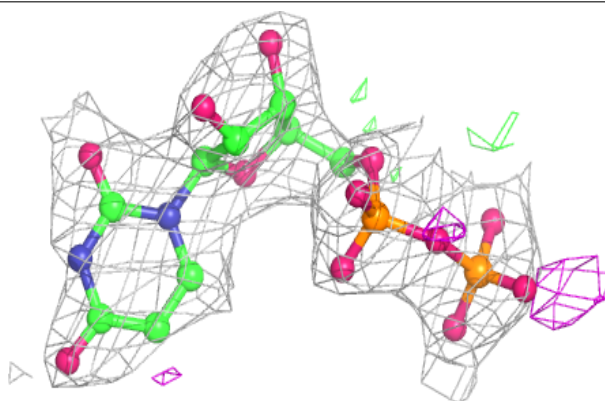


Electron density around UDP B 802:

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

**Electron density around UDP A 801:**

$2mF_o - DF_c$ (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.