



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 04:47 PM BST

PDB ID : 6BPM
Title : The crystal structure of the Ferric-Catecholate import receptor Fiu from K12
E. coli: Closed form (C21)
Authors : Grinter, R.
Deposited on : 2017-11-23
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

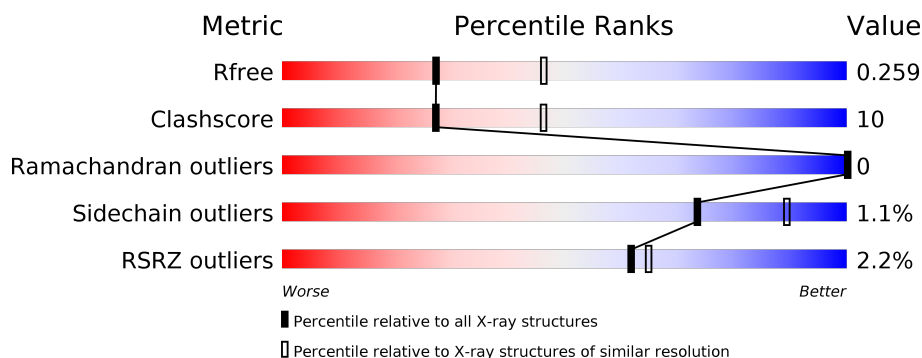
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	727	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>19%</div> <div>.</div> </div> </div>
1	C	727	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>19%</div> <div>..</div> </div> </div>

2 Entry composition ⓘ

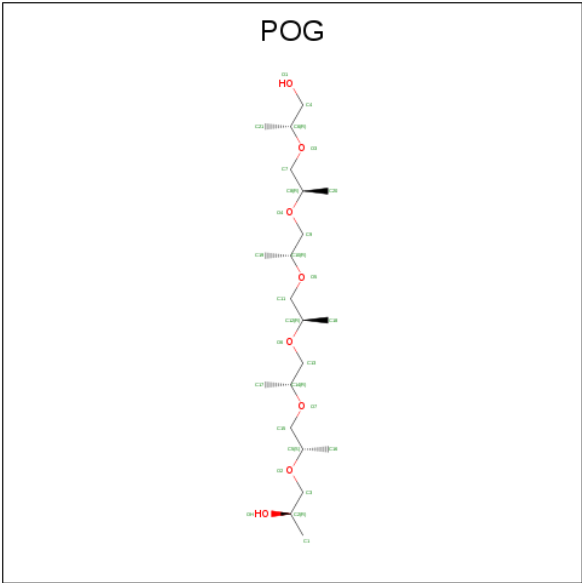
There are 4 unique types of molecules in this entry. The entry contains 11579 atoms, of which 320 are hydrogens and 0 are deuteriums.

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 Catecholate siderophore receptor Fiu.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	711	Total	C	N	O	S	0	1	0
			5417	3348	955	1102	12			
1	A	711	Total	C	N	O	S	0	2	0
			5427	3354	958	1103	12			

- Molecule 2 is (20S)-2,5,8,11,14,17-HEXAMETHYL-3,6,9,12,15,18-HEXAOXAHENICOSAN E-1,20-DIOL (three-letter code: POG) (formula: C₂₁H₄₄O₈).



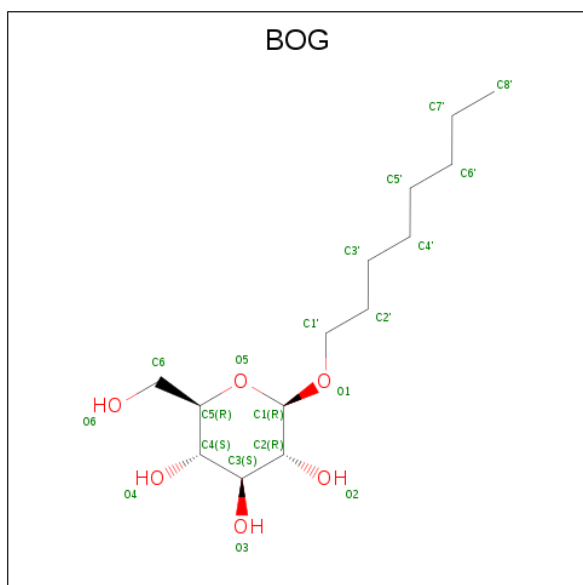
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	H	O	0	0
			73	21	44	8		
2	C	1	Total	C	H	O	0	0
			73	21	44	8		
2	C	1	Total	C	H	O	0	0
			73	21	44	8		
2	A	1	Total	C	H	O	0	0
			73	21	44	8		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			73	21	44	8		
2	A	1	Total	C	H	O	0	0
			73	21	44	8		

- Molecule 3 is octyl beta-D-glucopyranoside (three-letter code: BOG) (formula: $C_{14}H_{28}O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	H	O	0	0
			48	14	28	6		
3	A	1	Total	C	H	O	0	0
			48	14	28	6		

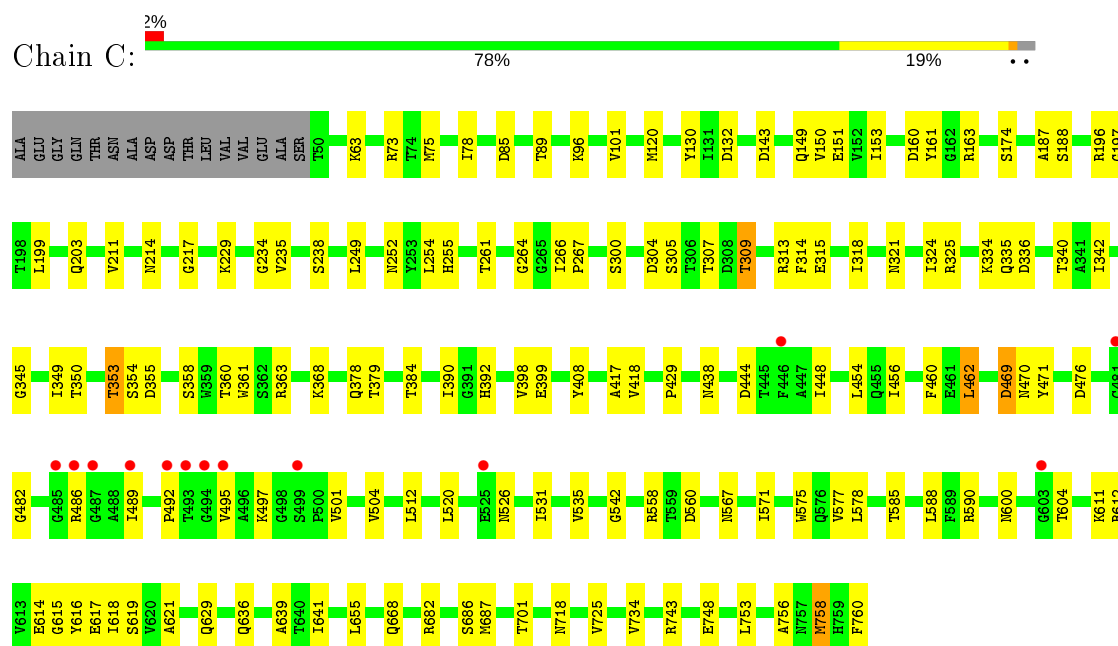
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	129	Total	O	0	0
			129	129		
4	A	72	Total	O	0	0
			72	72		

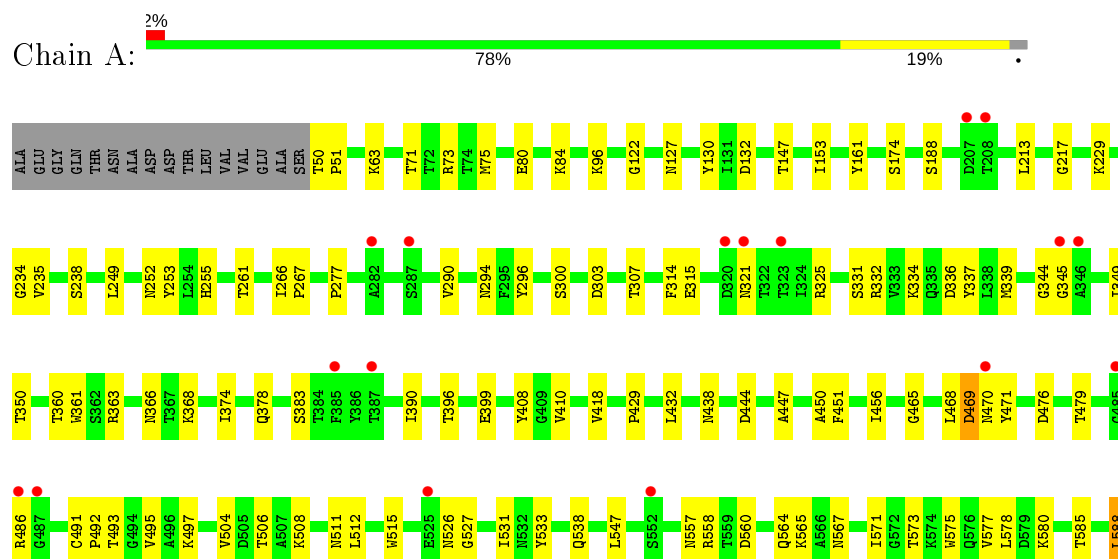
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Catecholate siderophore receptor Fiu



• Molecule 1: Catecholate siderophore receptor Fiu





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	191.43Å 75.67Å 136.87Å 90.00° 91.89° 90.00°	Depositor
Resolution (Å)	49.34 – 2.50 49.35 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.34-2.50) 99.9 (49.35-2.50)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.199 , 0.259 0.199 , 0.259	Depositor DCC
R_{free} test set	3282 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.921	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.3	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	0.93	EDS
Total number of atoms	11579	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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: POG, BOG

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	0.37	0/5544	0.61	1/7551 (0.0%)
1	C	0.43	0/5533	0.65	0/7536
All	All	0.40	0/11077	0.63	1/15087 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	588	LEU	CA-CB-CG	5.73	128.47	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5427	0	5153	94	0
1	C	5417	0	5147	107	0
2	A	87	132	132	27	0
2	C	87	132	132	12	0
3	A	20	28	28	0	0
3	C	20	28	28	0	0
4	A	72	0	0	3	0
4	C	129	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	11259	320	10620	218	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 10.

All (218) 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:A:531:ILE:HG22	1:A:571:ILE:HG12	1.30	1.13
1:C:187:ALA:HB1	2:C:801:POG:H161	1.27	1.13
1:C:254:LEU:HB3	1:C:309:THR:HG23	1.35	1.06
1:C:318:ILE:HD11	1:C:324:ILE:HD11	1.29	1.05
1:C:531:ILE:HG22	1:C:571:ILE:HG12	1.35	1.01
2:A:801:POG:H211	2:A:801:POG:H193	1.45	0.99
1:C:577:VAL:HG12	1:C:578:LEU:HG	1.43	0.97
1:C:353:THR:HG22	1:C:355:ASP:H	1.32	0.94
1:A:577:VAL:HG12	1:A:578:LEU:HG	1.48	0.94
1:C:75:MET:HG2	1:C:153:ILE:HG12	1.51	0.93
1:A:75:MET:HG3	1:A:153:ILE:HG12	1.54	0.89
1:C:254:LEU:HB3	1:C:309:THR:CG2	2.05	0.87
1:C:585:THR:OG1	1:C:619:SER:OG	1.93	0.87
1:C:492:PRO:O	1:C:495:VAL:HG12	1.80	0.82
1:C:318:ILE:HD11	1:C:324:ILE:CD1	2.12	0.80
1:A:470:ASN:OD1	1:A:511:ASN:ND2	2.15	0.79
1:C:585:THR:HG1	1:C:619:SER:HG	1.30	0.77
2:A:803:POG:H173	2:A:803:POG:H181	1.64	0.77
1:A:147:THR:HG23	4:A:964:HOH:O	1.85	0.77
1:A:396:THR:HG22	1:A:450:ALA:HA	1.66	0.76
1:C:143:ASP:OD2	1:C:196:ARG:NH2	2.21	0.73
1:A:249:LEU:HD13	1:A:314:PHE:CE1	2.23	0.72
1:C:492:PRO:HD2	1:C:495:VAL:HG11	1.71	0.72
1:C:249:LEU:HD13	1:C:314:PHE:CE1	2.23	0.72
1:A:217:GLY:HA3	2:A:801:POG:H132	1.74	0.70
1:C:75:MET:CG	1:C:153:ILE:HG12	2.23	0.68
1:A:588:LEU:HB3	2:A:802:POG:C19	2.23	0.68
1:C:686:SER:HB2	1:C:701:THR:O	1.93	0.68
1:C:482:GLY:O	1:C:497:LYS:NZ	2.23	0.68
1:A:580:LYS:O	1:A:580:LYS:HG2	1.94	0.67
1:C:78:ILE:HB	1:C:150:VAL:HG13	1.78	0.66
2:A:801:POG:C21	2:A:801:POG:H193	2.24	0.65
1:A:469:ASP:HB2	1:A:512:LEU:HB2	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:801:POG:H181	2:C:801:POG:H172	1.80	0.64
2:A:801:POG:H191	2:A:801:POG:H31	1.79	0.63
1:C:469:ASP:HB2	1:C:512:LEU:HB2	1.80	0.63
1:C:577:VAL:CG1	1:C:578:LEU:HG	2.24	0.62
1:C:336:ASP:HA	1:C:368:LYS:O	2.00	0.62
1:A:596:GLU:OE1	1:A:611:LYS:NZ	2.31	0.61
1:C:350:THR:HB	1:C:360:THR:OG1	2.01	0.61
1:A:266:ILE:O	1:A:743:ARG:HA	2.00	0.61
1:C:378:GLN:HG2	1:C:399:GLU:HG2	1.82	0.60
2:C:803:POG:H193	2:C:803:POG:H8	1.82	0.60
1:C:318:ILE:CD1	1:C:324:ILE:HD11	2.20	0.60
1:C:264:GLY:O	1:C:743:ARG:NH1	2.31	0.60
2:C:801:POG:H72	1:A:235:VAL:HG21	1.83	0.60
1:C:238:SER:HB3	1:C:252:ASN:HD22	1.67	0.59
1:A:558:ARG:HB3	1:A:560:ASP:OD1	2.02	0.59
1:C:615:GLY:HA3	1:C:636:GLN:O	2.02	0.59
1:C:489:ILE:HD11	1:C:501:VAL:HB	1.83	0.58
1:A:600:ASN:OD1	1:A:604:THR:HG22	2.03	0.58
1:A:315:GLU:HG2	1:A:325:ARG:HG2	1.85	0.58
1:A:588:LEU:HB3	2:A:802:POG:H191	1.85	0.58
2:A:801:POG:H201	2:A:801:POG:H10	1.85	0.58
2:A:803:POG:H181	2:A:803:POG:C17	2.32	0.57
1:C:590:ARG:NH2	1:C:614:GLU:OE1	2.23	0.57
1:A:336:ASP:HA	1:A:368:LYS:O	2.05	0.57
1:C:617:GLU:O	1:C:618:ILE:HD13	2.03	0.56
2:A:803:POG:C19	2:A:803:POG:H201	2.35	0.56
1:C:558:ARG:HB3	1:C:560:ASP:OD1	2.05	0.56
1:A:122:GLY:HA2	1:A:658:THR:OG1	2.05	0.56
1:C:78:ILE:HB	1:C:150:VAL:CG1	2.36	0.55
1:C:639:ALA:HB1	1:C:655:LEU:HD12	1.88	0.55
1:A:588:LEU:HB3	2:A:802:POG:H192	1.87	0.55
1:A:188:SER:HB3	1:A:753:LEU:HD11	1.88	0.55
1:A:585:THR:HB	1:A:619:SER:OG	2.07	0.55
2:A:803:POG:C18	2:A:803:POG:H173	2.35	0.53
1:A:96:LYS:HD3	1:A:682:ARG:CZ	2.38	0.53
1:A:217:GLY:HA3	2:A:801:POG:C13	2.37	0.53
1:C:345:GLY:O	1:C:349:ILE:HG13	2.07	0.53
1:A:728:LEU:HD13	1:A:728:LEU:O	2.09	0.52
1:A:491:CYS:SG	1:A:497:LYS:HA	2.49	0.52
1:A:71:THR:HG21	1:A:75:MET:CE	2.39	0.52
1:A:229:LYS:H	1:A:261:THR:HB	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:479:THR:HG22	4:A:968:HOH:O	2.08	0.52
1:C:748:GLU:OE1	1:C:748:GLU:N	2.30	0.52
1:A:564:GLN:HG3	1:A:596:GLU:OE2	2.10	0.52
1:C:460:PHE:CZ	1:C:462:LEU:HD11	2.45	0.52
1:A:337:TYR:CE2	1:A:368:LYS:HE2	2.45	0.52
1:C:315:GLU:HG2	1:C:325:ARG:HG2	1.91	0.52
1:C:378:GLN:HG2	1:C:399:GLU:CG	2.40	0.52
1:A:332:ARG:CZ	1:A:334:LYS:HE3	2.40	0.51
1:A:577:VAL:CG1	1:A:578:LEU:HG	2.31	0.51
1:A:294:ASN:HB3	1:A:296:TYR:CE1	2.45	0.51
1:C:701:THR:HG21	1:C:734:VAL:HG11	1.92	0.51
1:C:588:LEU:HD23	1:C:616:TYR:HB3	1.91	0.51
2:C:801:POG:C18	2:C:801:POG:H172	2.40	0.51
2:A:801:POG:H202	2:A:801:POG:H6	1.93	0.50
1:C:235:VAL:HG11	2:A:801:POG:H192	1.93	0.50
1:C:321:ASN:HB3	1:C:384:THR:HG22	1.93	0.50
1:A:748:GLU:N	1:A:748:GLU:OE1	2.29	0.50
1:A:331:SER:HB3	1:A:374:ILE:HG22	1.93	0.50
1:C:300:SER:HB2	1:C:429:PRO:HD2	1.94	0.50
1:A:447:ALA:HA	1:A:468:LEU:O	2.12	0.50
1:C:217:GLY:HA3	2:C:801:POG:H10	1.94	0.50
1:A:476:ASP:HA	1:A:504:VAL:O	2.12	0.50
1:A:615:GLY:HA2	2:A:802:POG:C17	2.42	0.50
1:A:344:GLY:HA3	1:A:361:TRP:CE2	2.47	0.50
1:C:235:VAL:HG21	2:A:801:POG:H32	1.94	0.50
1:C:85:ASP:HB3	1:C:668:GLN:OE1	2.10	0.49
1:A:253:TYR:HB2	2:A:803:POG:H111	1.94	0.49
1:C:353:THR:CG2	1:C:353:THR:O	2.61	0.49
1:C:470:ASN:HB3	2:C:803:POG:H14	1.95	0.49
1:C:611:LYS:HE3	1:C:641:ILE:HD13	1.95	0.49
1:A:615:GLY:HA2	2:A:802:POG:H173	1.95	0.49
2:C:801:POG:H182	1:A:235:VAL:HG11	1.95	0.49
1:C:229:LYS:H	1:C:261:THR:HB	1.77	0.49
1:C:353:THR:HG22	1:C:355:ASP:N	2.14	0.49
1:C:718:ASN:O	1:C:758:MET:HA	2.13	0.49
1:C:353:THR:CG2	1:C:355:ASP:HB3	2.43	0.48
1:A:390:ILE:HD12	1:A:456:ILE:HD13	1.95	0.48
1:C:254:LEU:HD23	1:C:309:THR:CG2	2.43	0.48
1:A:267:PRO:HD2	1:A:290:VAL:HG21	1.96	0.48
1:C:600:ASN:OD1	1:C:604:THR:HG22	2.13	0.48
1:C:132:ASP:OD2	1:C:174:SER:OG	2.23	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:533:TYR:OH	1:A:567:ASN:ND2	2.36	0.47
2:A:801:POG:C19	2:A:801:POG:H211	2.31	0.47
1:C:577:VAL:CG1	1:C:578:LEU:N	2.77	0.47
1:A:249:LEU:HD13	1:A:314:PHE:CD1	2.49	0.47
1:C:398:VAL:HG12	1:C:448:ILE:HG13	1.95	0.47
1:A:277:PRO:HA	1:A:695:VAL:HG12	1.96	0.47
1:A:378:GLN:HG2	1:A:399:GLU:HG2	1.96	0.47
2:A:803:POG:H193	2:A:803:POG:H8	1.95	0.47
1:A:300:SER:OG	1:A:429:PRO:HD2	2.15	0.47
1:C:96:LYS:HB3	1:C:682:ARG:HD3	1.97	0.47
1:A:444:ASP:O	1:A:471:TYR:HA	2.15	0.47
1:A:531:ILE:CG2	1:A:571:ILE:HG12	2.21	0.47
1:A:238:SER:HB3	1:A:252:ASN:HD22	1.80	0.46
1:A:80:GLU:O	1:A:84:LYS:HG2	2.15	0.46
1:C:75:MET:HG2	1:C:153:ILE:CG1	2.34	0.46
1:C:255:HIS:HA	1:C:307:THR:O	2.16	0.46
1:A:130:TYR:CE2	1:A:161:TYR:HB3	2.51	0.46
1:A:677:VAL:HG22	1:A:711:LEU:CD2	2.46	0.45
1:C:163:ARG:NH1	1:C:542:GLY:O	2.49	0.45
1:A:647:VAL:HG23	1:A:656:PRO:HD3	1.99	0.45
1:A:127:ASN:HD21	1:A:538:GLN:NE2	2.14	0.45
1:C:535:VAL:HG22	1:C:567:ASN:OD1	2.16	0.45
1:A:350:THR:HB	1:A:360:THR:OG1	2.17	0.45
1:A:573:THR:HG1	1:A:575:TRP:HE1	1.65	0.45
1:C:203:GLN:HG2	1:A:760:PHE:HB3	1.99	0.45
1:A:345:GLY:O	1:A:349:ILE:HG13	2.17	0.45
1:C:188:SER:HB3	1:C:753:LEU:HD11	1.99	0.45
1:A:396:THR:HG22	1:A:450:ALA:CA	2.43	0.44
1:C:63:LYS:NZ	1:C:160:ASP:O	2.45	0.44
1:C:217:GLY:HA3	2:C:801:POG:H112	1.99	0.44
1:A:718:ASN:O	1:A:758:MET:HA	2.16	0.44
1:A:71:THR:HG21	1:A:75:MET:HE3	1.98	0.44
1:C:267:PRO:HG3	1:C:342:ILE:HD12	1.98	0.44
1:C:130:TYR:CE2	1:C:161:TYR:HB3	2.52	0.44
1:A:565:LYS:NZ	1:A:565:LYS:HB3	2.32	0.44
1:C:353:THR:HG22	1:C:355:ASP:HB3	1.99	0.44
2:A:803:POG:H193	2:A:803:POG:H201	1.99	0.44
1:C:629:GLN:NE2	4:C:907:HOH:O	2.43	0.44
1:A:408:TYR:HA	1:A:438:ASN:OD1	2.17	0.44
1:A:132:ASP:OD2	1:A:174:SER:OG	2.25	0.43
1:A:321:ASN:O	1:A:383:SER:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:197:GLY:HA3	2:C:801:POG:H71	2.01	0.43
1:A:63:LYS:HB3	1:A:451:PHE:HE2	1.82	0.43
2:C:802:POG:H191	2:C:802:POG:H12	2.00	0.43
1:A:468:LEU:C	1:A:468:LEU:HD13	2.39	0.43
1:A:588:LEU:HD13	1:A:616:TYR:HB3	2.00	0.43
2:A:801:POG:O4	2:A:801:POG:H211	2.19	0.43
1:C:756:ALA:O	1:C:758:MET:CE	2.67	0.43
1:C:149:GLN:NE2	1:C:151:GLU:OE2	2.50	0.43
1:C:199:LEU:O	1:C:214:ASN:HA	2.18	0.43
1:C:255:HIS:CD2	2:A:801:POG:H181	2.54	0.43
1:C:305:SER:HA	1:C:334:LYS:O	2.18	0.42
1:C:321:ASN:HB3	1:C:384:THR:CG2	2.48	0.42
1:A:486:ARG:NH2	4:A:914:HOH:O	2.52	0.42
1:C:476:ASP:HA	1:C:504:VAL:O	2.19	0.42
2:A:803:POG:C8	2:A:803:POG:H193	2.49	0.42
1:C:360:THR:HB	1:C:417:ALA:HB1	2.01	0.42
1:C:266:ILE:O	1:C:743:ARG:HA	2.19	0.42
1:C:101:VAL:HG22	1:C:120:MET:HG3	2.00	0.42
1:C:267:PRO:HB3	1:C:361:TRP:CH2	2.55	0.42
1:C:211:VAL:HB	2:C:802:POG:H211	2.01	0.42
1:C:612:ARG:HD3	1:C:614:GLU:HG3	2.02	0.42
1:C:390:ILE:HD12	1:C:456:ILE:HD13	2.02	0.42
1:A:363:ARG:CZ	1:A:418:VAL:HB	2.50	0.42
1:A:492:PRO:HD2	1:A:495:VAL:HB	2.01	0.42
1:C:304:ASP:O	1:C:335:GLN:HA	2.19	0.42
1:C:760:PHE:HE2	2:A:803:POG:HOH	1.68	0.42
1:A:213:LEU:HB2	2:A:803:POG:H31	2.02	0.41
1:A:303:ASP:HA	1:A:337:TYR:HB3	2.02	0.41
1:A:526:ASN:O	1:A:575:TRP:HA	2.20	0.41
1:C:526:ASN:O	1:C:575:TRP:HA	2.20	0.41
1:A:234:GLY:HA2	1:A:255:HIS:O	2.21	0.41
1:A:527:GLY:HA3	1:A:575:TRP:CD2	2.54	0.41
1:C:520:LEU:HD23	1:C:520:LEU:C	2.40	0.41
1:A:615:GLY:HA3	1:A:636:GLN:O	2.20	0.41
1:A:410:VAL:HG21	1:A:432:LEU:HD22	2.02	0.41
1:A:465:GLY:O	1:A:515:TRP:HA	2.20	0.41
1:A:593:ILE:O	1:A:610:LYS:HA	2.20	0.41
1:C:234:GLY:HA2	1:C:255:HIS:O	2.21	0.41
1:C:254:LEU:HD23	1:C:309:THR:HG21	2.02	0.41
1:A:73:ARG:NH1	1:A:153:ILE:HG21	2.35	0.41
1:A:50:THR:HA	1:A:51:PRO:HD3	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:577:VAL:CG1	1:A:578:LEU:N	2.83	0.41
1:C:408:TYR:HA	1:C:438:ASN:OD1	2.21	0.41
1:A:255:HIS:HA	1:A:307:THR:O	2.21	0.41
1:A:636:GLN:HA	1:A:660:GLU:O	2.21	0.41
1:A:639:ALA:HB1	1:A:655:LEU:HD12	2.03	0.41
1:C:353:THR:HB	1:C:358:SER:OG	2.21	0.41
1:C:353:THR:HG23	1:C:353:THR:O	2.21	0.41
1:C:444:ASP:O	1:C:471:TYR:HA	2.19	0.41
1:A:547:LEU:HA	1:A:557:ASN:HB3	2.02	0.41
1:C:392:HIS:CE1	1:C:454:LEU:HD21	2.56	0.41
1:A:339:MET:HG3	1:A:366:ASN:HB2	2.02	0.40
1:C:73:ARG:NH1	1:C:153:ILE:HG21	2.36	0.40
1:C:89:THR:HB	1:C:196:ARG:CZ	2.51	0.40
1:C:621:ALA:HB2	4:C:957:HOH:O	2.21	0.40
1:A:471:TYR:OH	1:A:508:LYS:HE3	2.20	0.40
1:C:264:GLY:HA3	1:C:340:THR:O	2.21	0.40
1:C:353:THR:O	1:C:354:SER:OG	2.35	0.40
1:C:363:ARG:CZ	1:C:418:VAL:HB	2.51	0.40
1:C:325:ARG:O	1:C:379:THR:HA	2.20	0.40
1:C:313:ARG:NH1	1:C:315:GLU:OE2	2.51	0.40
1:C:489:ILE:CD1	1:C:501:VAL:HB	2.51	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	711/727 (98%)	695 (98%)	16 (2%)	0	100	100
1	C	710/727 (98%)	694 (98%)	16 (2%)	0	100	100
All	All	1421/1454 (98%)	1389 (98%)	32 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	581/591 (98%)	576 (99%)	5 (1%)	78	92
1	C	580/591 (98%)	572 (99%)	8 (1%)	67	86
All	All	1161/1182 (98%)	1148 (99%)	13 (1%)	73	89

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	309	THR
1	C	353	THR
1	C	462	LEU
1	C	469	ASP
1	C	486	ARG
1	C	687	MET
1	C	725	VAL
1	C	758	MET
1	A	469	ASP
1	A	493	THR
1	A	506	THR
1	A	617	GLU
1	A	687	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	511	ASN
1	A	538	GLN

5.3.3 RNA ⓘ

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](#)

8 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	BOG	A	804	-	20,20,20	0.88	0	25,25,25	1.32	2 (8%)
3	BOG	C	804	-	20,20,20	1.21	3 (15%)	25,25,25	1.35	3 (12%)
2	POG	A	802	-	21,28,28	1.21	0	26,34,34	1.53	3 (11%)
2	POG	C	801	-	21,28,28	1.10	0	26,34,34	1.01	2 (7%)
2	POG	A	801	-	21,28,28	1.10	1 (4%)	26,34,34	2.05	10 (38%)
2	POG	C	802	-	21,28,28	1.13	0	26,34,34	1.51	3 (11%)
2	POG	A	803	-	21,28,28	1.23	3 (14%)	26,34,34	1.73	5 (19%)
2	POG	C	803	-	21,28,28	1.21	1 (4%)	26,34,34	1.70	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	BOG	A	804	-	-	9/11/31/31	0/1/1/1
3	BOG	C	804	-	-	3/11/31/31	0/1/1/1
2	POG	A	802	-	-	13/32/32/32	-
2	POG	C	801	-	-	17/32/32/32	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	POG	A	801	-	-	17/32/32/32	-
2	POG	C	802	-	-	16/32/32/32	-
2	POG	A	803	-	-	9/32/32/32	-
2	POG	C	803	-	-	12/32/32/32	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	804	BOG	C4-C5	2.46	1.58	1.53
3	C	804	BOG	C4-C3	2.43	1.58	1.52
2	C	803	POG	O6-C13	-2.20	1.39	1.43
2	A	803	POG	O7-C15	-2.19	1.39	1.43
2	A	801	POG	O5-C11	-2.18	1.39	1.43
3	C	804	BOG	C1-C2	2.14	1.58	1.52
2	A	803	POG	O5-C11	-2.13	1.39	1.43
2	A	803	POG	O4-C9	-2.03	1.40	1.43

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	803	POG	C11-O5-C10	6.65	124.94	115.02
2	A	802	POG	C3-O2-C5	5.36	123.02	115.02
2	A	801	POG	C13-O6-C12	5.35	123.01	115.02
2	C	802	POG	C13-O6-C12	5.29	122.92	115.02
2	A	803	POG	C9-O4-C8	4.78	122.15	115.02
2	A	803	POG	C15-O7-C14	-4.32	108.58	115.02
2	A	802	POG	C11-O5-C10	4.13	121.18	115.02
2	A	801	POG	C11-O5-C10	3.74	120.61	115.02
2	A	803	POG	C3-O2-C5	3.56	120.34	115.02
3	A	804	BOG	C1-C2-C3	-3.25	103.23	110.00
3	C	804	BOG	O5-C1-O1	-3.11	102.61	109.97
2	A	801	POG	O5-C11-C12	-2.87	103.98	110.90
2	C	803	POG	C9-O4-C8	2.87	119.30	115.02
2	A	801	POG	C3-O2-C5	2.76	119.13	115.02
3	C	804	BOG	C1-C2-C3	-2.74	104.29	110.00
2	A	801	POG	C18-C12-C11	-2.68	104.92	112.63
2	A	801	POG	C16-C5-C15	-2.59	105.17	112.63
2	C	802	POG	C15-O7-C14	2.35	118.53	115.02
3	A	804	BOG	O2-C2-C3	-2.33	104.97	110.35
2	A	801	POG	O5-C10-C9	2.31	116.24	108.89
2	A	803	POG	C16-C5-C15	-2.29	106.06	112.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	POG	O2-C5-C15	2.28	116.15	108.89
2	A	802	POG	C19-C10-C9	-2.27	106.09	112.63
2	A	801	POG	O6-C12-C11	2.12	115.66	108.89
2	A	803	POG	O5-C11-C12	-2.11	105.80	110.90
2	C	803	POG	C20-C8-C7	-2.11	106.56	112.63
3	C	804	BOG	O2-C2-C3	-2.10	105.49	110.35
2	C	802	POG	O2-C3-C2	2.10	112.93	108.86
2	C	801	POG	C3-O2-C5	2.08	118.13	115.02
2	C	801	POG	C15-O7-C14	2.04	118.07	115.02
2	A	801	POG	C7-O3-C6	2.01	118.02	115.02

There are no chirality outliers.

All (96) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	804	BOG	C2-C1-O1-C1'
3	A	804	BOG	O5-C1-O1-C1'
2	A	802	POG	O1-C4-C6-O3
2	A	802	POG	O1-C4-C6-C21
2	A	802	POG	O3-C7-C8-O4
2	A	802	POG	O5-C10-C9-O4
2	A	802	POG	C19-C10-C9-O4
2	A	802	POG	O6-C13-C14-O7
2	A	802	POG	O6-C13-C14-C17
2	C	801	POG	OH-C2-C3-O2
2	C	801	POG	C21-C6-O3-C7
2	C	801	POG	O3-C7-C8-O4
2	C	801	POG	O3-C7-C8-C20
2	C	801	POG	C20-C8-O4-C9
2	C	801	POG	O5-C11-C12-O6
2	C	801	POG	O5-C11-C12-C18
2	C	801	POG	O6-C13-C14-O7
2	C	801	POG	O6-C13-C14-C17
2	C	801	POG	C5-C15-O7-C14
2	A	801	POG	C15-C5-O2-C3
2	A	801	POG	O1-C4-C6-O3
2	A	801	POG	O1-C4-C6-C21
2	A	801	POG	O3-C7-C8-O4
2	A	801	POG	O3-C7-C8-C20
2	A	801	POG	O5-C11-C12-O6
2	A	801	POG	C11-C12-O6-C13
2	A	801	POG	O6-C13-C14-O7

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Mol	Chain	Res	Type	Atoms
2	A	801	POG	O6-C13-C14-C17
2	C	802	POG	C2-C3-O2-C5
2	C	802	POG	O7-C15-C5-O2
2	C	802	POG	O7-C15-C5-C16
2	C	802	POG	O1-C4-C6-O3
2	C	802	POG	O1-C4-C6-C21
2	C	802	POG	O3-C7-C8-C20
2	C	802	POG	O5-C10-C9-O4
2	C	802	POG	C19-C10-C9-O4
2	C	802	POG	O5-C11-C12-O6
2	C	802	POG	O5-C11-C12-C18
2	C	802	POG	C18-C12-O6-C13
2	C	802	POG	C17-C14-O7-C15
2	A	803	POG	C16-C5-O2-C3
2	A	803	POG	O1-C4-C6-O3
2	A	803	POG	O3-C7-C8-O4
2	A	803	POG	O3-C7-C8-C20
2	A	803	POG	C20-C8-O4-C9
2	A	803	POG	C14-C13-O6-C12
2	C	803	POG	O7-C15-C5-O2
2	C	803	POG	O7-C15-C5-C16
2	C	803	POG	O3-C7-C8-O4
2	C	803	POG	C19-C10-O5-C11
2	A	803	POG	OH-C2-C3-O2
3	A	804	BOG	O5-C5-C6-O6
3	C	804	BOG	O5-C5-C6-O6
2	A	801	POG	OH-C2-C3-O2
2	C	803	POG	OH-C2-C3-O2
3	C	804	BOG	C4-C5-C6-O6
3	A	804	BOG	C1'-C2'-C3'-C4'
3	C	804	BOG	C5'-C6'-C7'-C8'
2	C	801	POG	C1-C2-C3-O2
2	A	803	POG	C1-C2-C3-O2
2	C	803	POG	C1-C2-C3-O2
2	C	801	POG	C2-C3-O2-C5
2	A	801	POG	C8-C7-O3-C6
2	C	803	POG	C8-C7-O3-C6
2	C	803	POG	C5-C15-O7-C14
2	A	802	POG	C15-C5-O2-C3
2	A	801	POG	C9-C10-O5-C11
2	A	802	POG	O7-C15-C5-C16
2	A	802	POG	O3-C7-C8-C20

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Mol	Chain	Res	Type	Atoms
2	C	801	POG	O7-C15-C5-C16
2	A	801	POG	O5-C11-C12-C18
2	A	802	POG	O7-C15-C5-O2
2	C	801	POG	O7-C15-C5-O2
2	C	802	POG	O3-C7-C8-O4
2	A	803	POG	O1-C4-C6-C21
2	A	801	POG	C21-C6-O3-C7
2	C	802	POG	C16-C5-O2-C3
2	C	803	POG	C20-C8-O4-C9
2	A	802	POG	C12-C11-O5-C10
2	C	802	POG	C10-C9-O4-C8
2	C	803	POG	C12-C11-O5-C10
3	A	804	BOG	C4-C5-C6-O6
3	A	804	BOG	C3'-C4'-C5'-C6'
2	C	801	POG	O5-C10-C9-O4
2	C	802	POG	O6-C13-C14-O7
2	C	803	POG	O5-C10-C9-O4
2	A	802	POG	C18-C12-O6-C13
2	C	801	POG	C19-C10-O5-C11
2	A	801	POG	C19-C10-O5-C11
2	A	801	POG	C18-C12-O6-C13
2	C	803	POG	C21-C6-O3-C7
3	A	804	BOG	O1-C1'-C2'-C3'
2	C	801	POG	C14-C13-O6-C12
2	A	801	POG	C14-C13-O6-C12
3	A	804	BOG	C2'-C1'-O1-C1
3	A	804	BOG	C2'-C3'-C4'-C5'

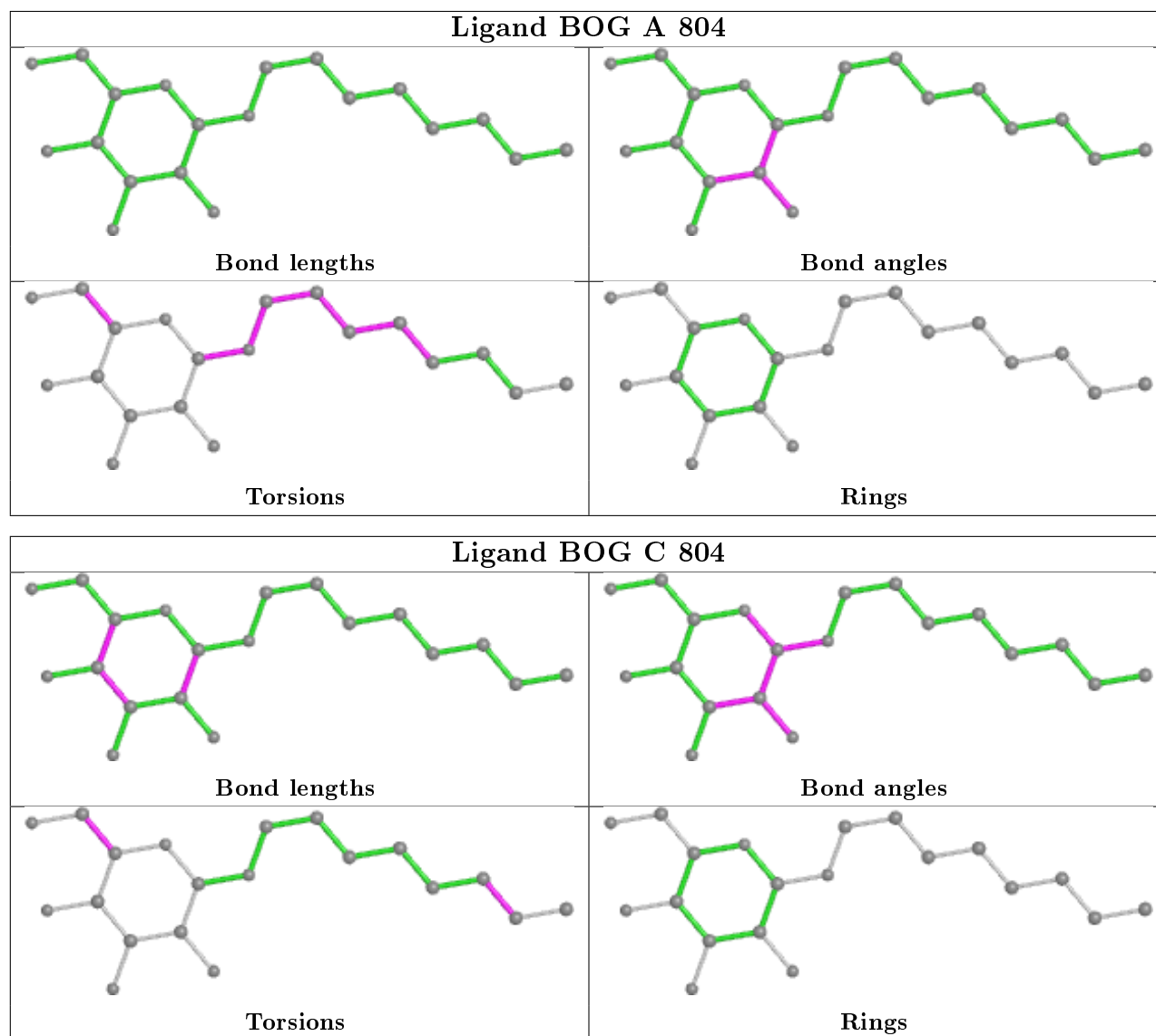
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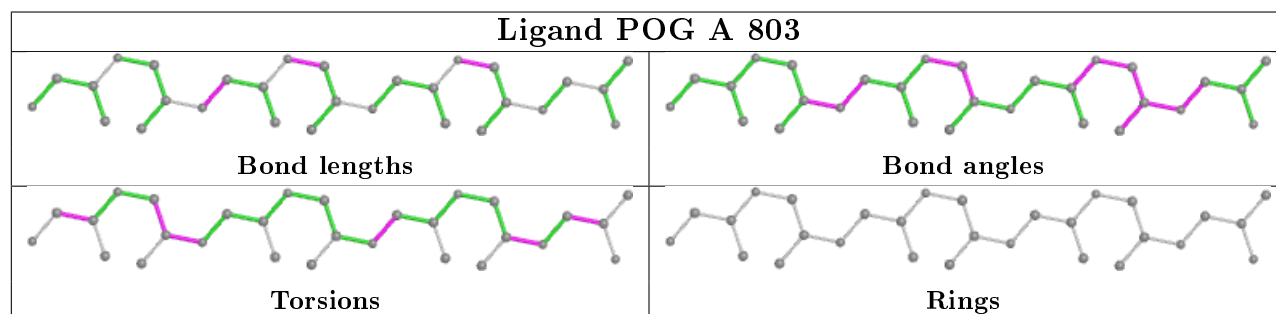
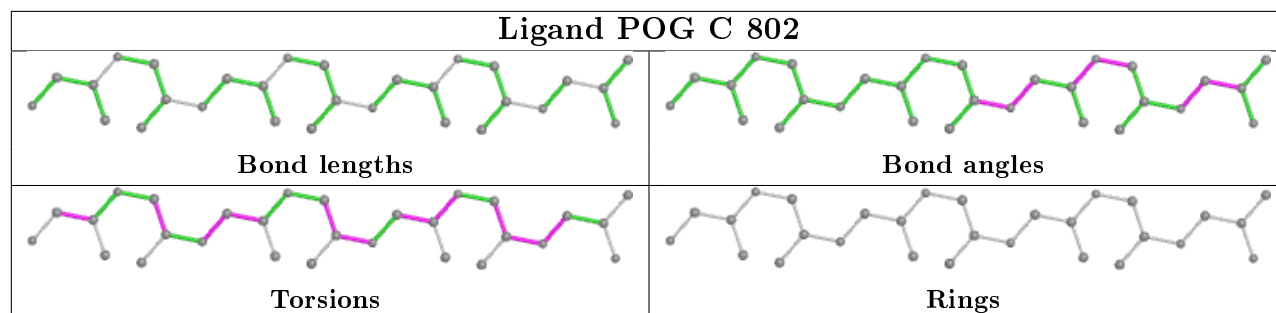
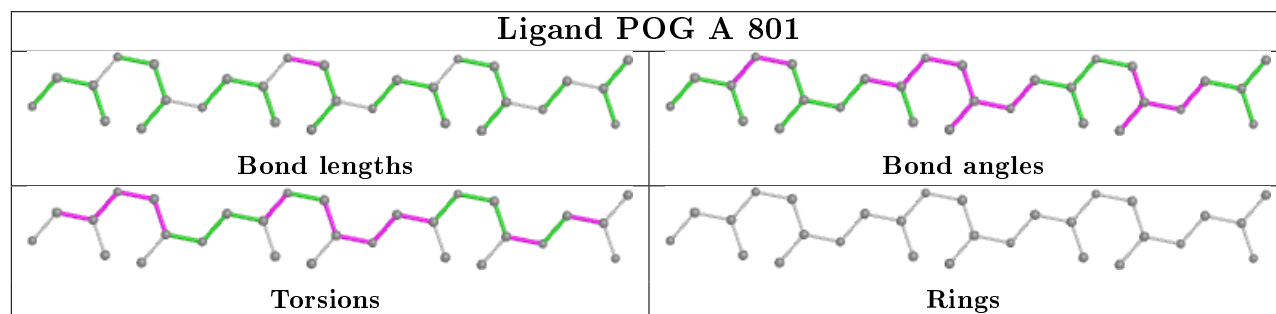
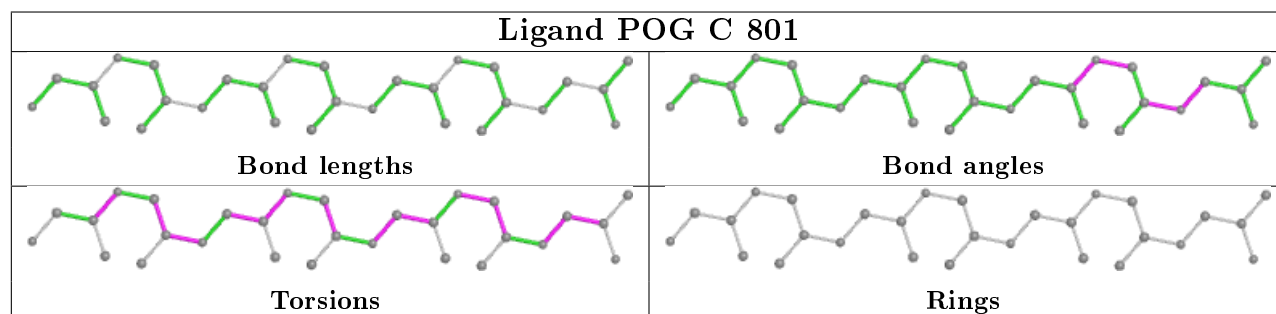
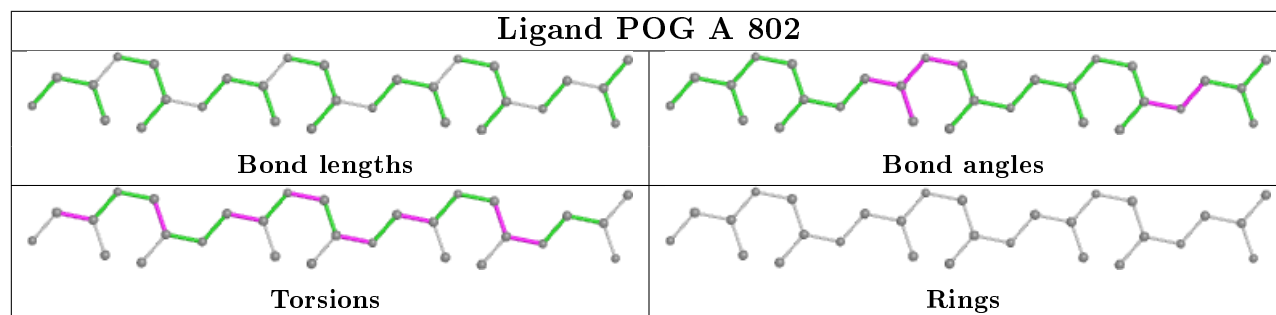
6 monomers are involved in 39 short contacts:

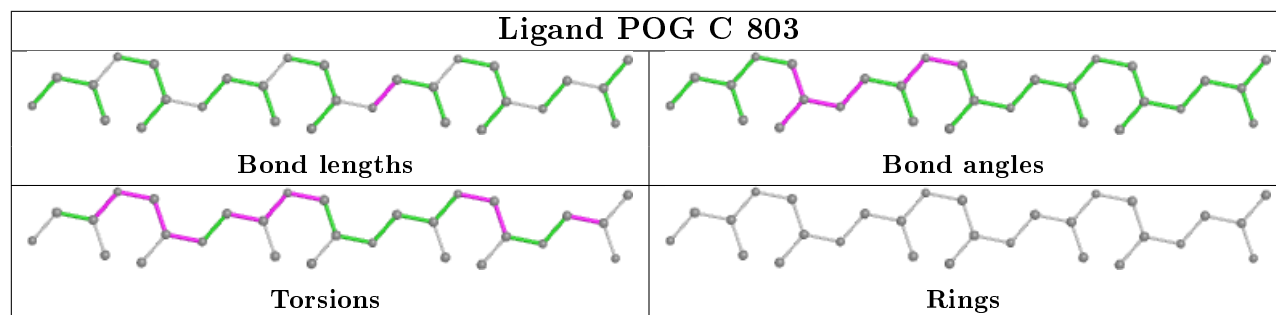
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	802	POG	5	0
2	C	801	POG	8	0
2	A	801	POG	12	0
2	C	802	POG	2	0
2	A	803	POG	10	0
2	C	803	POG	2	0

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	711/727 (97%)	0.10	18 (2%) 57 61	25, 51, 76, 119	0
1	C	711/727 (97%)	-0.11	13 (1%) 68 71	23, 38, 61, 110	0
All	All	1422/1454 (97%)	-0.01	31 (2%) 62 65	23, 44, 73, 119	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	486	ARG	5.0
1	C	494	GLY	4.5
1	C	495	VAL	3.8
1	A	486	ARG	3.6
1	A	207	ASP	3.6
1	A	712	GLY	3.3
1	A	323	THR	3.2
1	C	492	PRO	3.1
1	A	487	GLY	3.1
1	A	485	GLY	3.1
1	C	525	GLU	3.1
1	A	321	ASN	2.8
1	C	489	ILE	2.8
1	A	345	GLY	2.8
1	C	487	GLY	2.7
1	A	552	SER	2.7
1	A	282	ALA	2.6
1	C	493	THR	2.6
1	C	481	CYS	2.5
1	A	385	PHE	2.3
1	A	320	ASP	2.3
1	C	499	SER	2.3
1	A	346	ALA	2.3
1	C	603	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	525	GLU	2.2
1	A	208	THR	2.2
1	C	485	GLY	2.2
1	A	287	SER	2.2
1	A	470	ASN	2.1
1	A	387	THR	2.1
1	C	446	PHE	2.1

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](#)

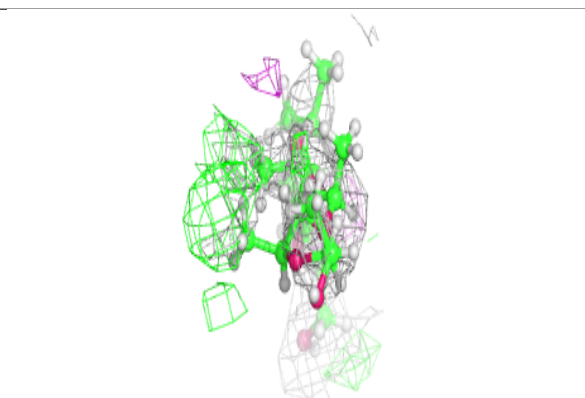
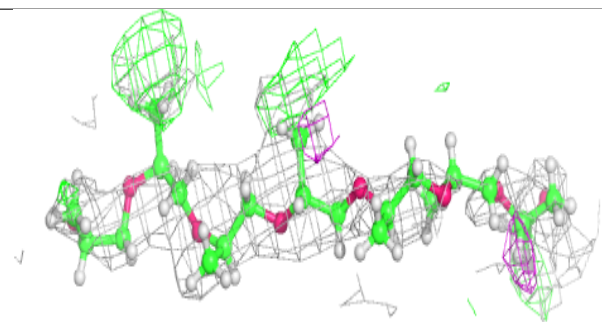
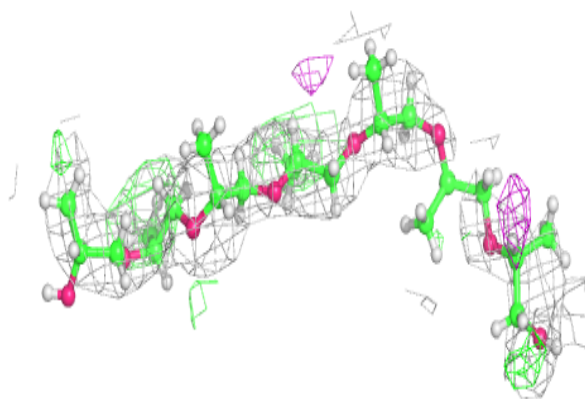
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	POG	A	802	29/29	0.64	0.34	66,116,158,185	0
2	POG	C	803	29/29	0.81	0.46	66,115,163,167	0
2	POG	C	802	29/29	0.82	0.46	71,110,137,139	0
2	POG	A	803	29/29	0.83	0.33	79,110,156,162	0
3	BOG	A	804	20/20	0.84	0.47	85,103,119,121	0
3	BOG	C	804	20/20	0.87	0.37	54,89,116,123	0
2	POG	A	801	29/29	0.89	0.23	25,71,139,153	0
2	POG	C	801	29/29	0.92	0.23	47,71,176,186	0

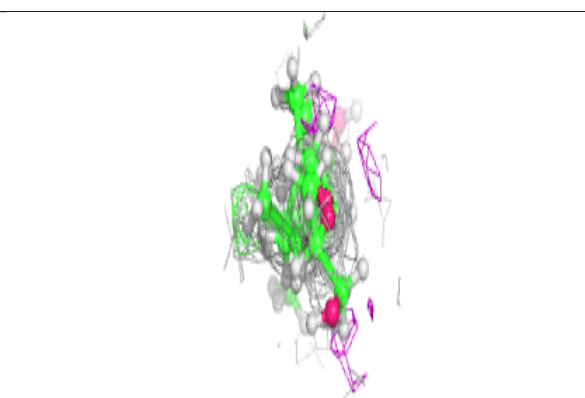
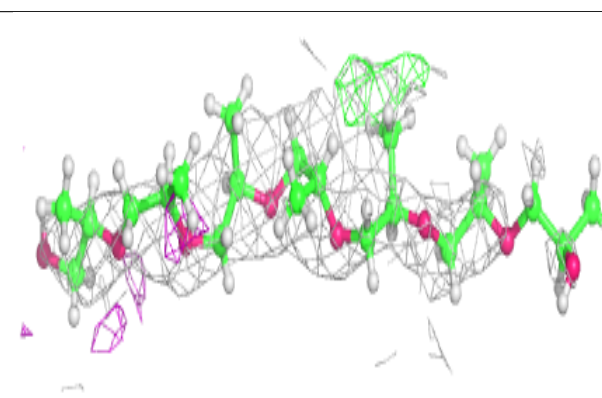
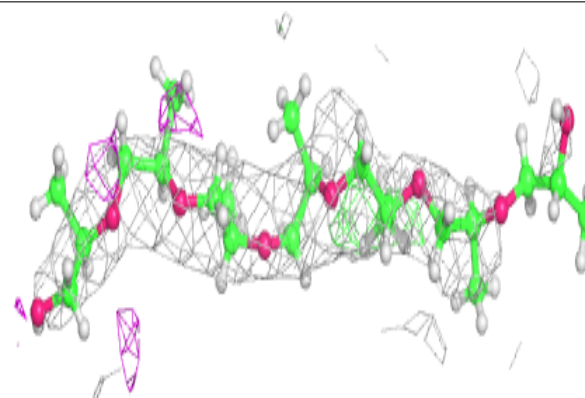
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around POG A 802:

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

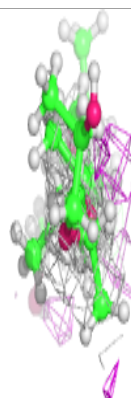
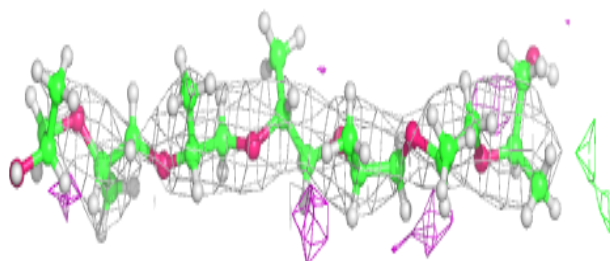
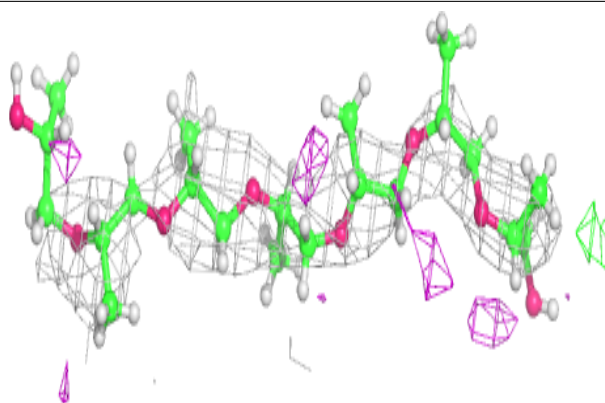
**Electron density around POG C 803:**

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

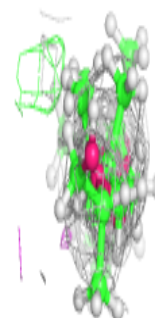
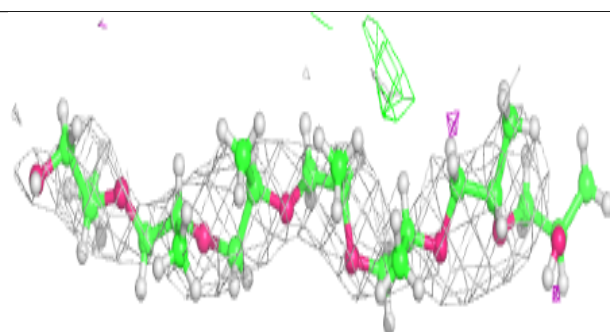
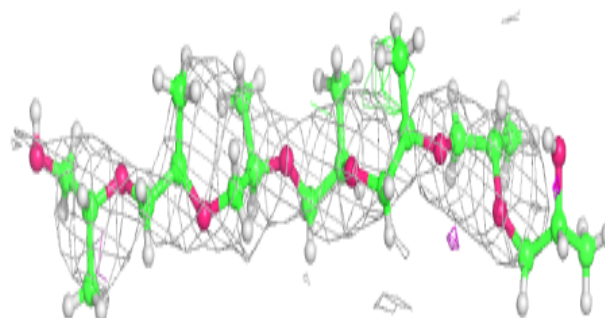


Electron density around POG C 802:

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

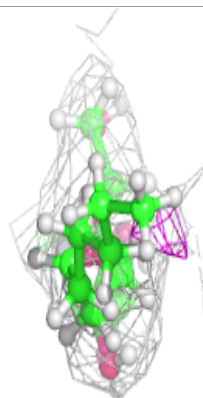
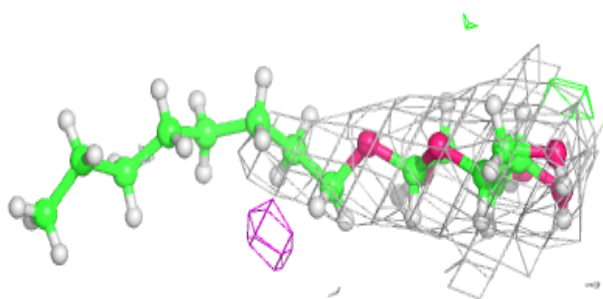
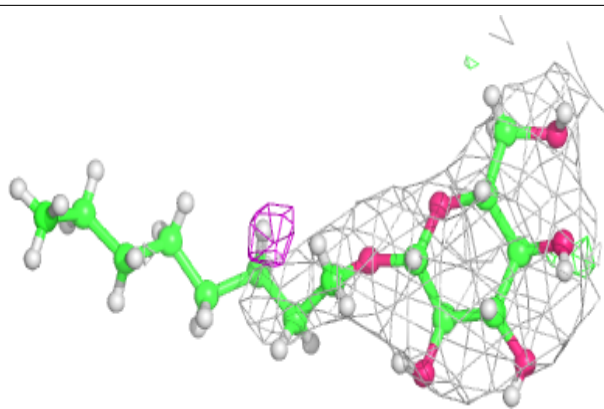
**Electron density around POG A 803:**

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

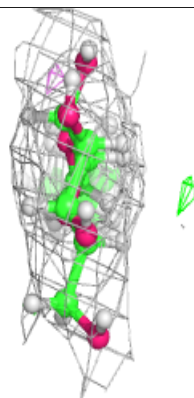
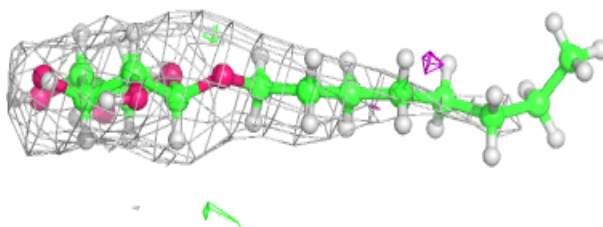
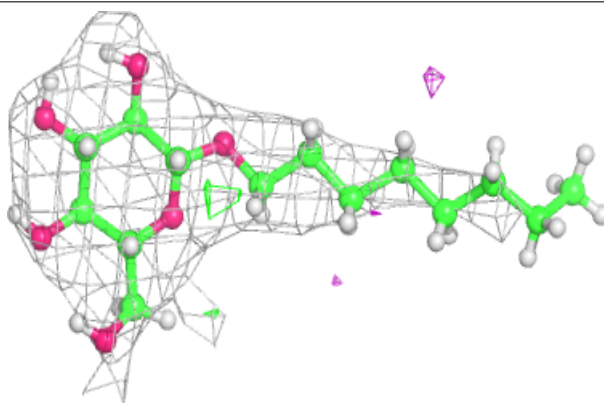


Electron density around BOG A 804:

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

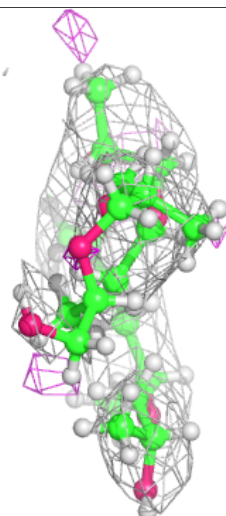
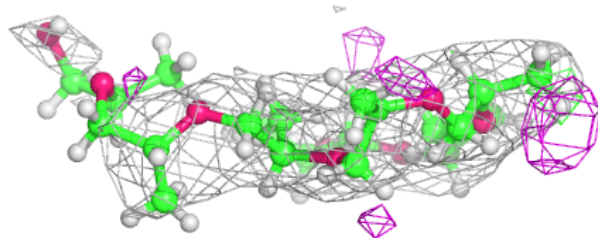
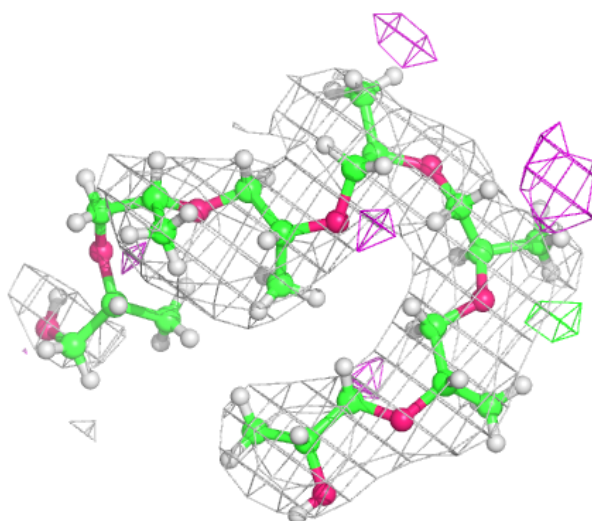
**Electron density around BOG C 804:**

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



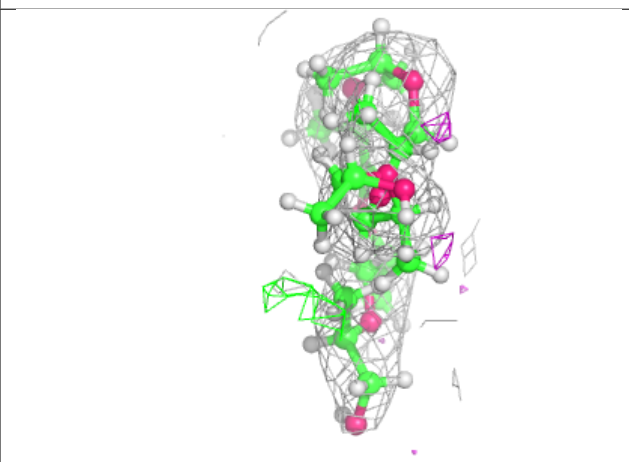
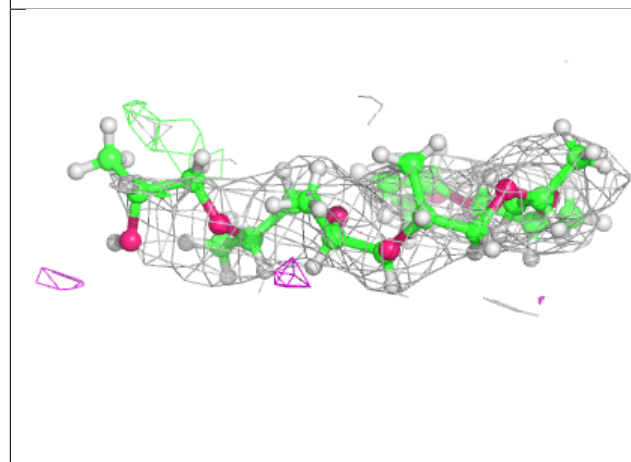
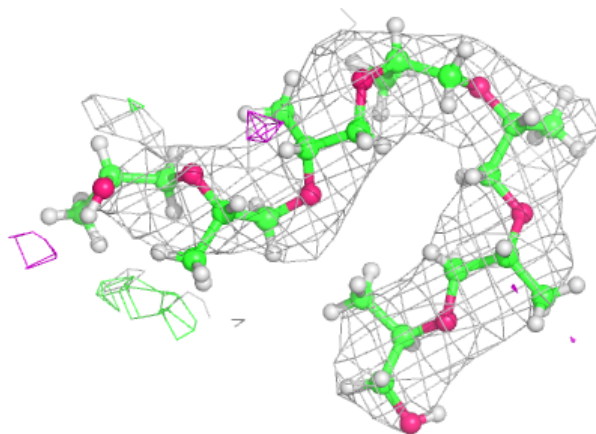
Electron density around POG A 801:

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



Electron density around POG C 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 [i](#)

There are no such residues in this entry.