Full wwPDB X-ray Structure Validation Report (i)

Nov 14, 2023 – 09:39 PM JST

PDB ID	:	6AFZ
Title	:	Proton pyrophosphatase-E225H mutant
Authors	:	Tsai, J.-Y.; Li, K.-M.; Sun, Y.-J.
Deposited on	:	2018-08-08
Resolution	:	2.48 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 (i) 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 (i)) were used in the production of this report:

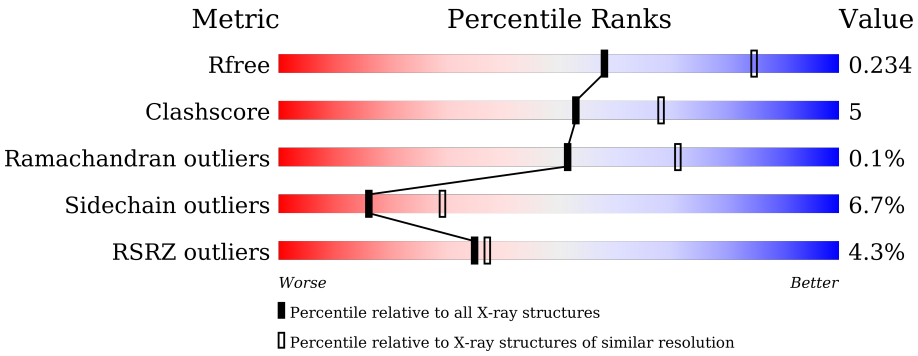
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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	5857 (2.50-2.46)
Clashscore	141614	6594 (2.50-2.46)
Ramachandran outliers	138981	6469 (2.50-2.46)
Sidechain outliers	138945	6471 (2.50-2.46)
RSRZ outliers	127900	5738 (2.50-2.46)

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 <=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	766	3% 82% 13%
1	B	766	5% 81% 14%

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
2	PO4	A	1001	-	-	X	-
2	PO4	A	1002	-	-	X	-
2	PO4	B	1002	-	-	X	-
5	1PG	A	1014	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11218 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.

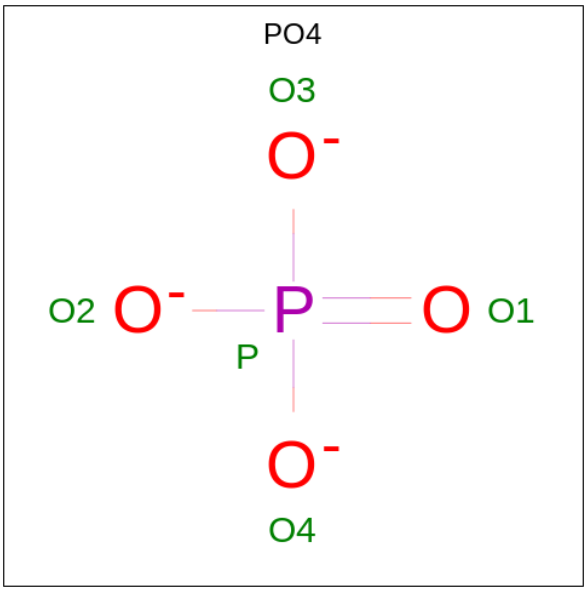
- Molecule 1 is a protein called Pyrophosphate-energized vacuolar membrane proton pump.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	740	Total	C	N	O	S	0	0	0
			5432	3541	864	998	29			
1	B	740	Total	C	N	O	S	0	0	0
			5432	3541	864	998	29			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	225	HIS	GLU	engineered mutation	UNP P21616
B	225	HIS	GLU	engineered mutation	UNP P21616

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0

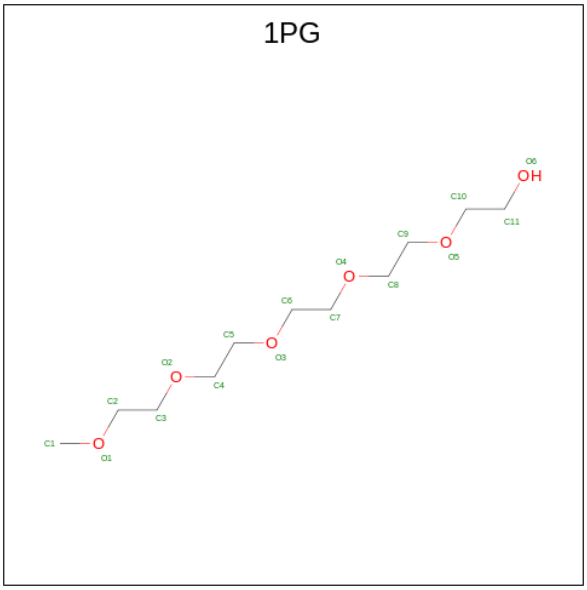
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	5	Total Mg 5 5	0	0
3	B	5	Total Mg 5 5	0	0

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total K 1 1	0	0
4	B	1	Total K 1 1	0	0

- Molecule 5 is 2-(2-{2-[2-(2-METHOXY-ETHOXY)-ETHOXY]-ETHOXY}-ETHOXY)-ETHANOL (three-letter code: 1PG) (formula: $C_{11}H_{24}O_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 17 11 6	0	0
5	A	1	Total C O 17 11 6	0	0
5	A	1	Total C O 17 11 6	0	0
5	A	1	Total C O 17 11 6	0	0
5	A	1	Total C O 17 11 6	0	0
5	A	1	Total C O 10 7 3	0	0
5	B	1	Total C O 17 11 6	0	0

- Molecule 6 is water.

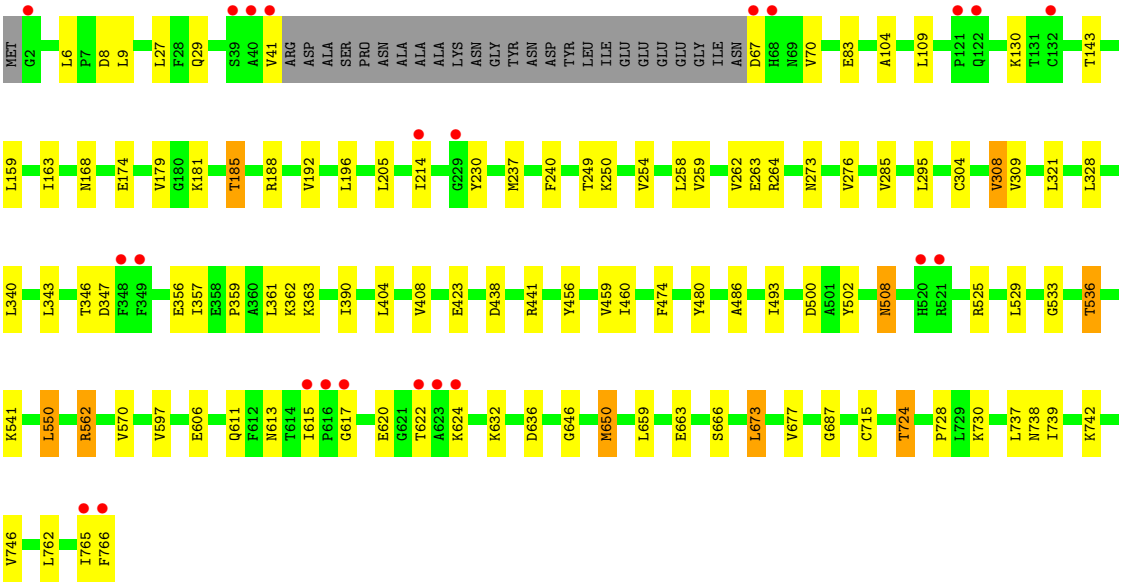
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	136	Total O 136 136	0	0
6	B	74	Total O 74 74	0	0

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.

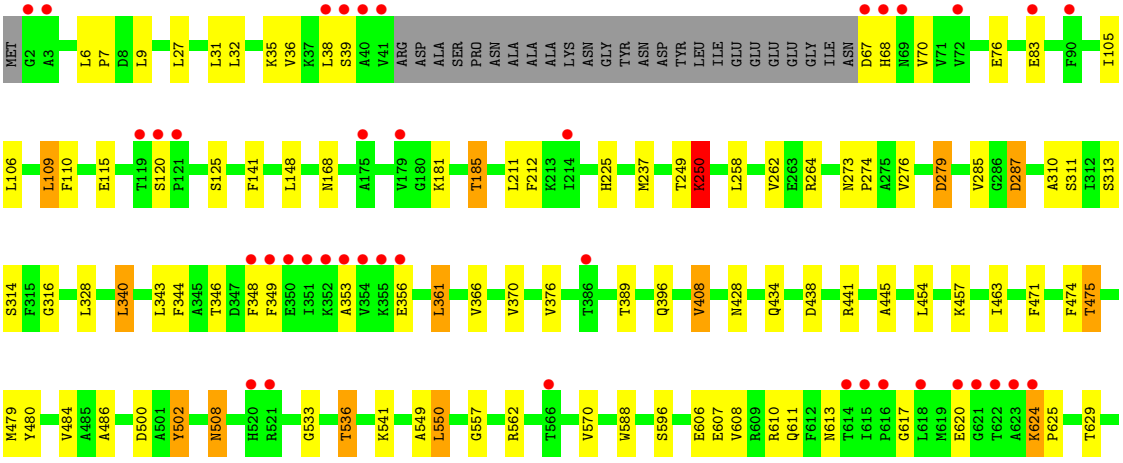
• Molecule 1: Pyrophosphate-energized vacuolar membrane proton pump

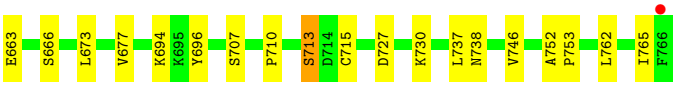
Chain A:



• Molecule 1: Pyrophosphate-energized vacuolar membrane proton pump

Chain B:




E663
S666
L673
V677
K694
K695
Y696
S707
P710
S713
D714
C715
D727
K730
L737
N738
V746
A752
P753
L762
I765
F766

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	215.60Å 88.02Å 158.19Å 90.00° 124.99° 90.00°	Depositor
Resolution (Å)	25.70 – 2.48 25.70 – 2.48	Depositor EDS
% Data completeness (in resolution range)	96.7 (25.70-2.48) 85.5 (25.70-2.48)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.73 (at 2.47Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.201 , 0.235 0.201 , 0.234	Depositor DCC
R_{free} test set	2000 reflections (2.41%)	wwPDB-VP
Wilson B-factor (Å^2)	48.6	Xtriage
Anisotropy	0.711	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 57.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	11218	wwPDB-VP
Average B, all atoms (Å^2)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: MG, 1PG, PO4, K

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	0.45	0/5539	0.57	1/7527 (0.0%)
1	B	0.41	0/5539	0.54	0/7527
All	All	0.43	0/11078	0.55	1/15054 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	673	LEU	CA-CB-CG	5.34	127.58	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	5432	0	5577	56	0
1	B	5432	0	5577	64	0
2	A	10	0	0	2	0
2	B	10	0	0	2	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	95	0	133	3	0
5	B	17	0	24	0	0
6	A	136	0	0	2	0
6	B	74	0	0	5	0
All	All	11218	0	11311	120	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 5.

All (120) 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:343:LEU:HA	1:B:346:THR:HG22	1.70	0.73
1:A:606:GLU:OE1	1:B:441:ARG:NH2	2.20	0.73
1:A:441:ARG:NH2	1:B:606:GLU:OE1	2.23	0.70
1:B:250:LYS:HG3	1:B:727:ASP:HB3	1.72	0.69
1:B:715:CYS:SG	6:B:1146:HOH:O	2.49	0.69
1:A:159:LEU:HD13	1:A:196:LEU:HD13	1.76	0.67
1:A:273:ASN:HB3	1:A:276:VAL:HG23	1.76	0.67
1:A:570:VAL:HG22	1:B:570:VAL:HG22	1.76	0.66
1:A:597:VAL:HG11	1:A:728:PRO:HB3	1.78	0.64
1:A:357:ILE:HD13	1:A:529:LEU:HD23	1.80	0.64
1:A:343:LEU:HA	1:A:346:THR:HG22	1.81	0.63
2:B:1002:PO4:O4	6:B:1101:HOH:O	2.16	0.62
1:B:249:THR:HG23	1:B:250:LYS:HD2	1.81	0.61
1:B:273:ASN:HB3	1:B:276:VAL:HG23	1.83	0.61
1:A:663:GLU:O	1:A:666:SER:HB2	2.02	0.58
1:B:249:THR:O	1:B:250:LYS:HB2	2.02	0.58
1:A:174:GLU:HG3	1:A:185:THR:HG21	1.86	0.57
1:A:254:VAL:HG13	1:A:724:THR:HG23	1.86	0.57
1:B:471:PHE:O	1:B:475:THR:OG1	2.23	0.56
2:B:1001:PO4:P	2:B:1002:PO4:O1	2.63	0.56
1:A:500:ASP:HA	1:A:536:THR:HG23	1.88	0.55
1:B:617:GLY:HA2	1:B:620:GLU:OE1	2.07	0.55
1:B:250:LYS:NZ	6:B:1103:HOH:O	2.40	0.54
1:A:258:LEU:HD21	1:A:724:THR:HG21	1.90	0.54
1:A:533:GLY:O	1:A:536:THR:HG22	2.08	0.54
1:B:32:LEU:O	1:B:35:LYS:HB2	2.08	0.53
1:A:240:PHE:CE2	1:A:650:MET:HE1	2.44	0.53
1:A:486:ALA:HB2	1:A:550:LEU:HB3	1.90	0.53
1:A:214:ILE:HG22	5:A:1014:1PG:H21	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:GLN:OE1	1:A:188:ARG:NH1	2.39	0.51
1:B:361:LEU:HD21	1:B:502:TYR:CE2	2.46	0.51
1:B:264:ARG:NH2	1:B:613:ASN:OD1	2.44	0.51
1:A:460:ILE:HD11	1:B:588:TRP:CD1	2.46	0.51
1:A:273:ASN:HB3	1:A:276:VAL:CG2	2.42	0.50
1:B:610:ARG:HH21	1:B:611:GLN:HG2	1.77	0.50
1:A:562:ARG:HD2	6:A:1110:HOH:O	2.11	0.50
1:A:687:GLY:HA3	1:A:730:LYS:HG2	1.94	0.49
1:A:632:LYS:HD3	1:A:636:ASP:OD2	2.12	0.49
1:A:570:VAL:HG22	1:B:570:VAL:CG2	2.42	0.49
1:B:438:ASP:OD2	1:B:441:ARG:NH1	2.45	0.49
1:A:163:ILE:HG12	1:A:192:VAL:HB	1.95	0.48
1:B:287:ASP:N	1:B:287:ASP:OD1	2.47	0.48
1:B:500:ASP:HA	1:B:536:THR:HG23	1.94	0.48
1:A:295:LEU:HD12	1:A:493:ILE:HD11	1.96	0.47
1:A:304:CYS:O	1:A:308:VAL:HG22	2.14	0.47
1:A:362:LYS:HD3	1:A:423:GLU:OE2	2.14	0.47
1:B:710:PRO:O	1:B:713:SER:HB3	2.14	0.47
1:A:715:CYS:SG	6:A:1163:HOH:O	2.48	0.47
1:B:533:GLY:O	1:B:536:THR:HG22	2.15	0.47
1:A:264:ARG:NH2	1:A:613:ASN:OD1	2.47	0.47
2:A:1001:PO4:O3	2:A:1002:PO4:O1	2.31	0.47
5:A:1013:1PG:H51	5:A:1013:1PG:H72	1.59	0.47
1:B:389:THR:HB	1:B:396:GLN:HB3	1.98	0.46
1:B:311:SER:O	1:B:316:GLY:HA3	2.16	0.46
1:A:739:ILE:HA	1:A:742:LYS:HE2	1.97	0.46
1:A:438:ASP:OD2	1:A:441:ARG:NH1	2.48	0.46
1:B:310:ALA:HB2	1:B:480:TYR:CE1	2.50	0.46
1:B:141:PHE:HB3	1:B:212:PHE:CE2	2.50	0.46
1:B:344:PHE:HA	1:B:348:PHE:HB2	1.98	0.45
1:B:663:GLU:O	1:B:666:SER:HB2	2.16	0.45
1:B:31:LEU:HD12	1:B:31:LEU:HA	1.77	0.45
1:A:724:THR:HG22	1:B:445:ALA:HB1	1.99	0.45
1:B:366:VAL:O	1:B:370:VAL:HG23	2.17	0.45
1:A:456:TYR:O	1:A:459:VAL:HG22	2.17	0.44
1:B:479:MET:SD	1:B:557:GLY:HA3	2.57	0.44
1:A:168:ASN:HB2	1:A:508:ASN:HB3	1.99	0.44
1:B:225:HIS:CD2	1:B:562:ARG:HE	2.36	0.44
1:A:404:LEU:HD23	1:A:404:LEU:HA	1.82	0.44
1:B:434:GLN:HG2	1:B:696:TYR:CZ	2.53	0.44
1:B:624:LYS:HD2	1:B:625:PRO:HD2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:THR:HG23	1:A:250:LYS:HE2	1.99	0.44
1:B:274:PRO:HD2	1:B:608:VAL:HG13	1.99	0.44
1:B:314:SER:HB3	1:B:389:THR:OG1	2.17	0.44
1:B:225:HIS:CD2	1:B:562:ARG:NE	2.86	0.43
1:A:258:LEU:O	1:A:262:VAL:HB	2.19	0.43
1:B:168:ASN:HB2	1:B:508:ASN:HB3	1.99	0.43
1:A:363:LYS:HD3	1:A:363:LYS:HA	1.72	0.43
1:B:454:LEU:O	1:B:457:LYS:HB3	2.18	0.43
1:A:346:THR:HG23	1:A:347:ASP:CG	2.39	0.43
1:A:179:VAL:HG13	1:A:525:ARG:HG2	2.01	0.43
1:A:659:LEU:HB3	1:A:762:LEU:HD22	2.00	0.43
2:A:1001:PO4:P	2:A:1002:PO4:O1	2.77	0.43
1:B:274:PRO:HG2	1:B:608:VAL:HG22	2.01	0.43
1:A:390:ILE:HD13	1:A:390:ILE:HA	1.89	0.42
1:B:279:ASP:OD1	6:B:1102:HOH:O	2.21	0.42
1:B:730:LYS:HE3	6:B:1133:HOH:O	2.19	0.42
1:B:550:LEU:HD12	1:B:550:LEU:HA	1.92	0.42
1:B:106:LEU:HD22	1:B:110:PHE:CZ	2.55	0.42
1:B:258:LEU:O	1:B:262:VAL:HB	2.20	0.42
1:B:752:ALA:HB3	1:B:753:PRO:HD3	2.02	0.42
1:A:611:GLN:HB3	1:A:615:ILE:HD12	2.02	0.42
1:A:104:ALA:HB1	1:A:143:THR:HG23	2.01	0.42
5:A:1010:1PG:H91	5:A:1010:1PG:H72	1.74	0.42
1:B:273:ASN:HB3	1:B:276:VAL:CG2	2.49	0.41
1:B:463:ILE:HD13	1:B:463:ILE:HA	1.84	0.41
1:B:549:ALA:HA	1:B:673:LEU:HD11	2.02	0.41
1:B:273:ASN:HA	1:B:274:PRO:HD2	1.96	0.41
1:B:313:SER:HB3	1:B:389:THR:O	2.20	0.41
1:B:376:VAL:HG11	1:B:408:VAL:HG22	2.02	0.41
1:B:428:ASN:HB2	1:B:696:TYR:CD1	2.55	0.41
1:A:196:LEU:HD23	1:A:196:LEU:HA	1.78	0.41
1:A:356:GLU:C	1:A:359:PRO:HD2	2.41	0.41
1:B:211:LEU:HD23	1:B:211:LEU:HA	1.91	0.41
1:A:8:ASP:CG	1:A:321:LEU:HB3	2.40	0.41
1:B:607:GLU:OE2	1:B:629:THR:HG21	2.20	0.41
1:A:67:ASP:O	1:A:70:VAL:HG12	2.20	0.41
1:A:205:LEU:HB2	1:A:230:TYR:CE2	2.56	0.41
1:B:6:LEU:HD12	1:B:7:PRO:HD2	2.02	0.41
1:B:353:ALA:HB3	1:B:356:GLU:HG3	2.03	0.41
1:A:259:VAL:O	1:A:263:GLU:HB2	2.21	0.41
1:B:181:LYS:O	1:B:185:THR:HG22	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:340:LEU:HD12	1:B:340:LEU:HA	1.89	0.41
1:A:240:PHE:CE1	1:A:646:GLY:HA3	2.56	0.41
1:A:617:GLY:HA2	1:A:620:GLU:HB2	2.03	0.41
1:A:650:MET:HE2	1:A:650:MET:HB2	1.77	0.40
1:A:762:LEU:O	1:A:766:PHE:HD2	2.04	0.40
1:A:309:VAL:HG11	1:A:480:TYR:HB2	2.03	0.40
1:B:486:ALA:HB2	1:B:550:LEU:HB3	2.03	0.40
1:B:105:ILE:HG22	1:B:109:LEU:HD22	2.03	0.40
1:B:694:LYS:HB3	1:B:694:LYS:HE2	1.65	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	736/766 (96%)	714 (97%)	22 (3%)	0	100 100
1	B	736/766 (96%)	703 (96%)	31 (4%)	2 (0%)	41 59
All	All	1472/1532 (96%)	1417 (96%)	53 (4%)	2 (0%)	51 71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	250	LYS
1	B	36	VAL

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	568/588 (97%)	535 (94%)	33 (6%)	20	36
1	B	568/588 (97%)	525 (92%)	43 (8%)	13	24
All	All	1136/1176 (97%)	1060 (93%)	76 (7%)	16	29

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	9	LEU
1	A	27	LEU
1	A	41	VAL
1	A	83	GLU
1	A	109	LEU
1	A	130	LYS
1	A	181	LYS
1	A	185	THR
1	A	237	MET
1	A	285	VAL
1	A	308	VAL
1	A	328	LEU
1	A	340	LEU
1	A	361	LEU
1	A	408	VAL
1	A	474	PHE
1	A	502	TYR
1	A	508	ASN
1	A	536	THR
1	A	541	LYS
1	A	550	LEU
1	A	562	ARG
1	A	622	THR
1	A	624	LYS
1	A	650	MET
1	A	673	LEU
1	A	677	VAL
1	A	724	THR
1	A	737	LEU
1	A	738	ASN
1	A	746	VAL
1	A	765	ILE

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Mol	Chain	Res	Type
1	B	9	LEU
1	B	27	LEU
1	B	38	LEU
1	B	39	SER
1	B	67	ASP
1	B	68	HIS
1	B	70	VAL
1	B	76	GLU
1	B	83	GLU
1	B	109	LEU
1	B	115	GLU
1	B	120	SER
1	B	125	SER
1	B	148	LEU
1	B	185	THR
1	B	237	MET
1	B	250	LYS
1	B	279	ASP
1	B	285	VAL
1	B	287	ASP
1	B	328	LEU
1	B	340	LEU
1	B	349	PHE
1	B	361	LEU
1	B	408	VAL
1	B	474	PHE
1	B	475	THR
1	B	484	VAL
1	B	502	TYR
1	B	508	ASN
1	B	536	THR
1	B	541	LYS
1	B	550	LEU
1	B	596	SER
1	B	624	LYS
1	B	677	VAL
1	B	707	SER
1	B	713	SER
1	B	737	LEU
1	B	738	ASN
1	B	746	VAL
1	B	762	LEU

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Mol	Chain	Res	Type
1	B	765	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 23 ligands modelled in this entry, 12 are monoatomic - leaving 11 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	$\#\|Z\| > 2$
5	1PG	A	1011	-	16,16,16	0.85	0	15,15,15	0.32	0
5	1PG	A	1010	-	16,16,16	0.87	0	15,15,15	0.45	0
2	PO4	B	1001	4,3	4,4,4	0.97	0	6,6,6	1.89	1 (16%)
2	PO4	A	1001	3	4,4,4	1.39	1 (25%)	6,6,6	0.40	0
2	PO4	B	1002	3	4,4,4	2.25	2 (50%)	6,6,6	1.06	0
5	1PG	A	1013	-	16,16,16	0.85	0	15,15,15	0.42	0
5	1PG	A	1009	-	16,16,16	0.82	0	15,15,15	0.44	0
5	1PG	B	1009	-	16,16,16	0.86	0	15,15,15	0.42	0
5	1PG	A	1014	-	9,9,16	0.72	0	8,8,15	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	1PG	A	1012	-	16,16,16	0.82	0	15,15,15	0.39	0
2	PO4	A	1002	3	4,4,4	0.51	0	6,6,6	0.77	0

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
5	1PG	A	1011	-	-	6/14/14/14	-
5	1PG	A	1010	-	-	10/14/14/14	-
5	1PG	A	1013	-	-	9/14/14/14	-
5	1PG	A	1009	-	-	9/14/14/14	-
5	1PG	B	1009	-	-	5/14/14/14	-
5	1PG	A	1014	-	-	3/7/7/14	-
5	1PG	A	1012	-	-	10/14/14/14	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1002	PO4	P-O4	-2.86	1.46	1.54
2	A	1001	PO4	P-O1	2.62	1.57	1.50
2	B	1002	PO4	P-O3	-2.59	1.46	1.54

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1001	PO4	O3-P-O1	-3.64	97.58	110.89

There are no chirality outliers.

All (52) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1013	1PG	C7-C6-O3-C5
5	A	1010	1PG	C2-C3-O2-C4
5	A	1012	1PG	C11-C10-O5-C9
5	A	1010	1PG	C9-C8-O4-C7
5	A	1011	1PG	O4-C8-C9-O5
5	A	1013	1PG	O4-C8-C9-O5
5	A	1014	1PG	O1-C2-C3-O2

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Mol	Chain	Res	Type	Atoms
5	B	1009	1PG	O3-C6-C7-O4
5	A	1010	1PG	O3-C6-C7-O4
5	A	1013	1PG	O3-C6-C7-O4
5	A	1011	1PG	O1-C2-C3-O2
5	A	1009	1PG	O1-C2-C3-O2
5	A	1012	1PG	O4-C8-C9-O5
5	A	1010	1PG	C8-C9-O5-C10
5	A	1013	1PG	O2-C4-C5-O3
5	A	1013	1PG	O5-C10-C11-O6
5	A	1012	1PG	O1-C2-C3-O2
5	A	1013	1PG	C5-C4-O2-C3
5	A	1009	1PG	O3-C6-C7-O4
5	A	1010	1PG	O1-C2-C3-O2
5	A	1013	1PG	O1-C2-C3-O2
5	A	1009	1PG	O2-C4-C5-O3
5	A	1010	1PG	O4-C8-C9-O5
5	A	1014	1PG	C7-C6-O3-C5
5	B	1009	1PG	C5-C4-O2-C3
5	B	1009	1PG	C4-C5-O3-C6
5	A	1010	1PG	C4-C5-O3-C6
5	A	1009	1PG	C8-C9-O5-C10
5	A	1014	1PG	C4-C5-O3-C6
5	A	1010	1PG	C6-C7-O4-C8
5	A	1013	1PG	C8-C9-O5-C10
5	A	1009	1PG	C6-C7-O4-C8
5	A	1012	1PG	C5-C4-O2-C3
5	A	1013	1PG	C6-C7-O4-C8
5	A	1012	1PG	C9-C8-O4-C7
5	A	1010	1PG	C11-C10-O5-C9
5	A	1009	1PG	C7-C6-O3-C5
5	A	1011	1PG	C6-C7-O4-C8
5	A	1009	1PG	C5-C4-O2-C3
5	A	1012	1PG	C2-C3-O2-C4
5	A	1011	1PG	C11-C10-O5-C9
5	B	1009	1PG	O2-C4-C5-O3
5	A	1012	1PG	C6-C7-O4-C8
5	A	1012	1PG	O5-C10-C11-O6
5	A	1011	1PG	O3-C6-C7-O4
5	A	1009	1PG	C2-C3-O2-C4
5	A	1012	1PG	O2-C4-C5-O3
5	A	1009	1PG	C3-C2-O1-C1
5	A	1012	1PG	C7-C6-O3-C5

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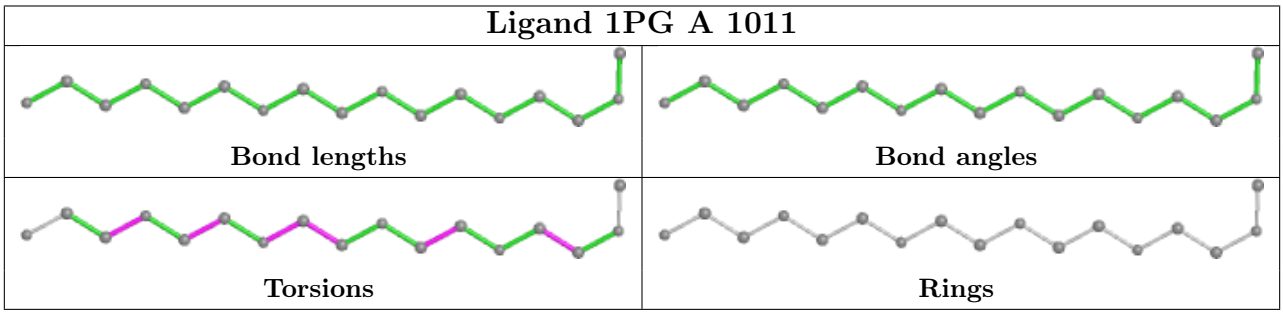
Mol	Chain	Res	Type	Atoms
5	B	1009	1PG	O4-C8-C9-O5
5	A	1011	1PG	O2-C4-C5-O3
5	A	1010	1PG	O2-C4-C5-O3

There are no ring outliers.

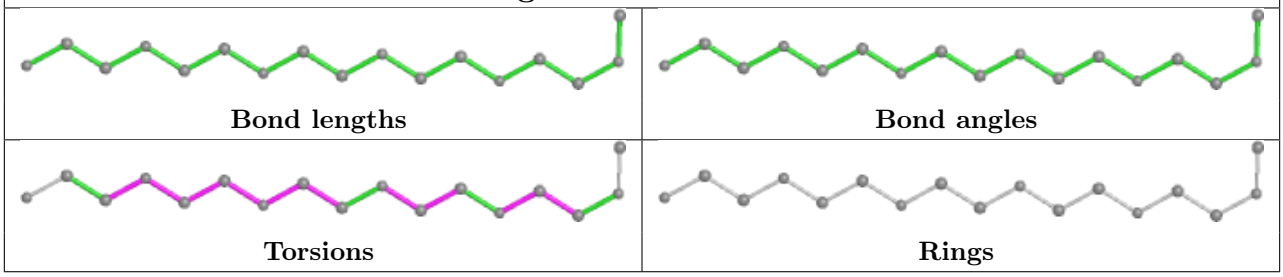
7 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1010	1PG	1	0
2	B	1001	PO4	1	0
2	A	1001	PO4	2	0
2	B	1002	PO4	2	0
5	A	1013	1PG	1	0
5	A	1014	1PG	1	0
2	A	1002	PO4	2	0

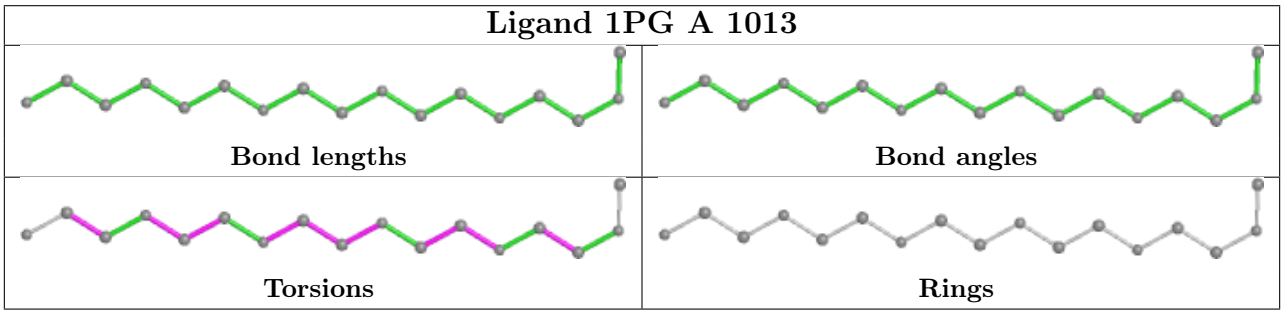
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



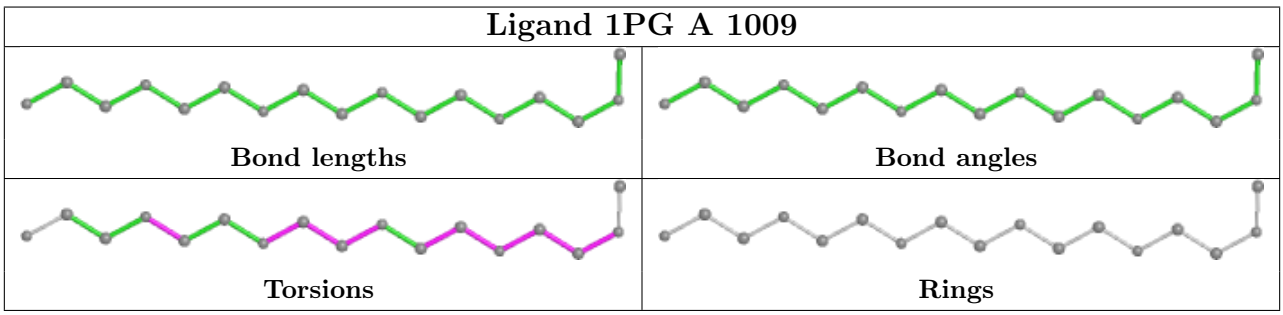
Ligand 1PG A 1010



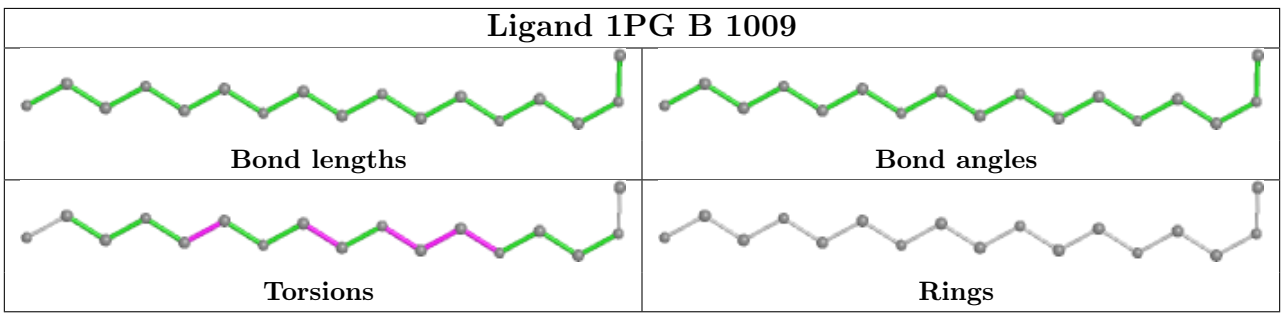
Ligand 1PG A 1013



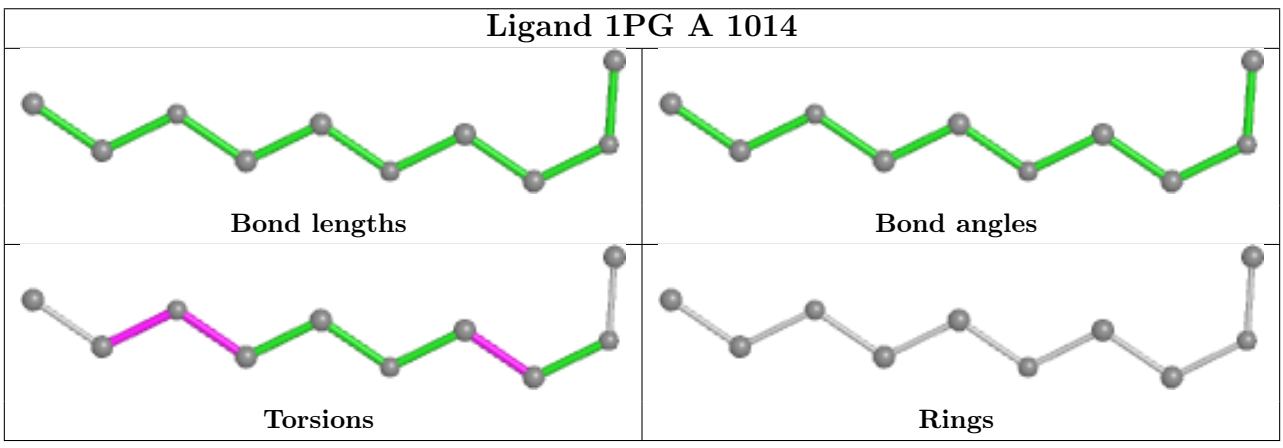
Ligand 1PG A 1009

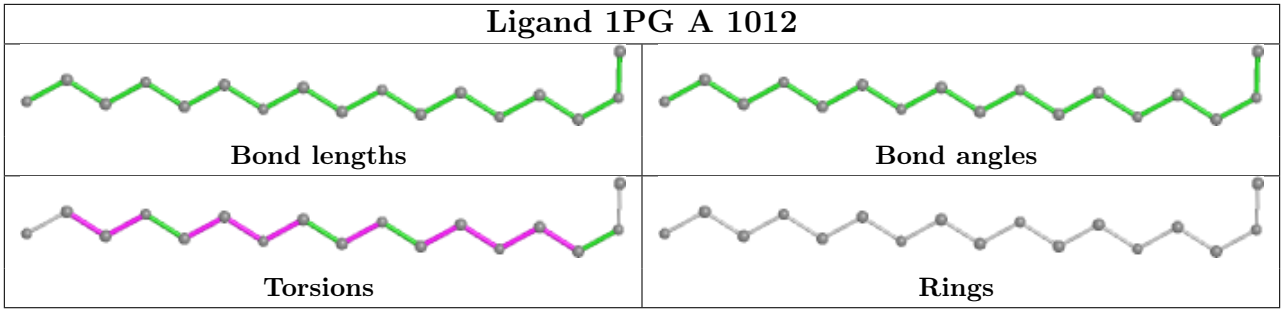


Ligand 1PG B 1009



Ligand 1PG A 1014





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, 95^{th} 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	740/766 (96%)	-0.25	23 (3%)	49	51	31, 44, 61, 95	0
1	B	740/766 (96%)	-0.05	41 (5%)	25	26	31, 53, 81, 102	0
All	All	1480/1532 (96%)	-0.15	64 (4%)	35	37	31, 47, 75, 102	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	41	VAL	11.9
1	A	41	VAL	10.6
1	A	40	ALA	10.1
1	B	40	ALA	8.3
1	B	2	GLY	8.0
1	B	39	SER	7.0
1	B	623	ALA	6.9
1	A	39	SER	5.6
1	B	615	ILE	5.0
1	A	624	LYS	4.8
1	B	624	LYS	4.8
1	B	38	LEU	4.5
1	B	616	PRO	4.5
1	A	623	ALA	4.5
1	A	348	PHE	4.4
1	A	520	HIS	4.3
1	A	2	GLY	4.1
1	A	616	PRO	3.9
1	B	622	THR	3.8
1	A	68	HIS	3.8
1	B	520	HIS	3.8
1	A	122	GLN	3.7
1	B	353	ALA	3.7
1	B	349	PHE	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	614	THR	3.7
1	B	214	ILE	3.6
1	B	3	ALA	3.6
1	B	67	ASP	3.5
1	B	69	ASN	3.5
1	B	348	PHE	3.4
1	B	352	LYS	3.3
1	A	67	ASP	3.3
1	A	766	PHE	3.2
1	B	72	VAL	3.1
1	A	617	GLY	3.1
1	A	132	CYS	3.0
1	B	68	HIS	3.0
1	B	351	ILE	2.9
1	B	350	GLU	2.9
1	B	766	PHE	2.8
1	B	179	VAL	2.7
1	A	214	ILE	2.6
1	A	349	PHE	2.6
1	A	521	ARG	2.6
1	A	121	PRO	2.5
1	B	175	ALA	2.5
1	B	356	GLU	2.4
1	A	615	ILE	2.4
1	B	621	GLY	2.3
1	A	622	THR	2.3
1	B	120	SER	2.3
1	B	119	THR	2.3
1	B	355	LYS	2.2
1	B	354	VAL	2.2
1	B	566	THR	2.2
1	B	386	THR	2.1
1	B	90	PHE	2.1
1	B	521	ARG	2.1
1	A	229	GLY	2.1
1	B	83	GLU	2.1
1	B	121	PRO	2.1
1	A	765	ILE	2.1
1	B	618	LEU	2.1
1	B	620	GLU	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)

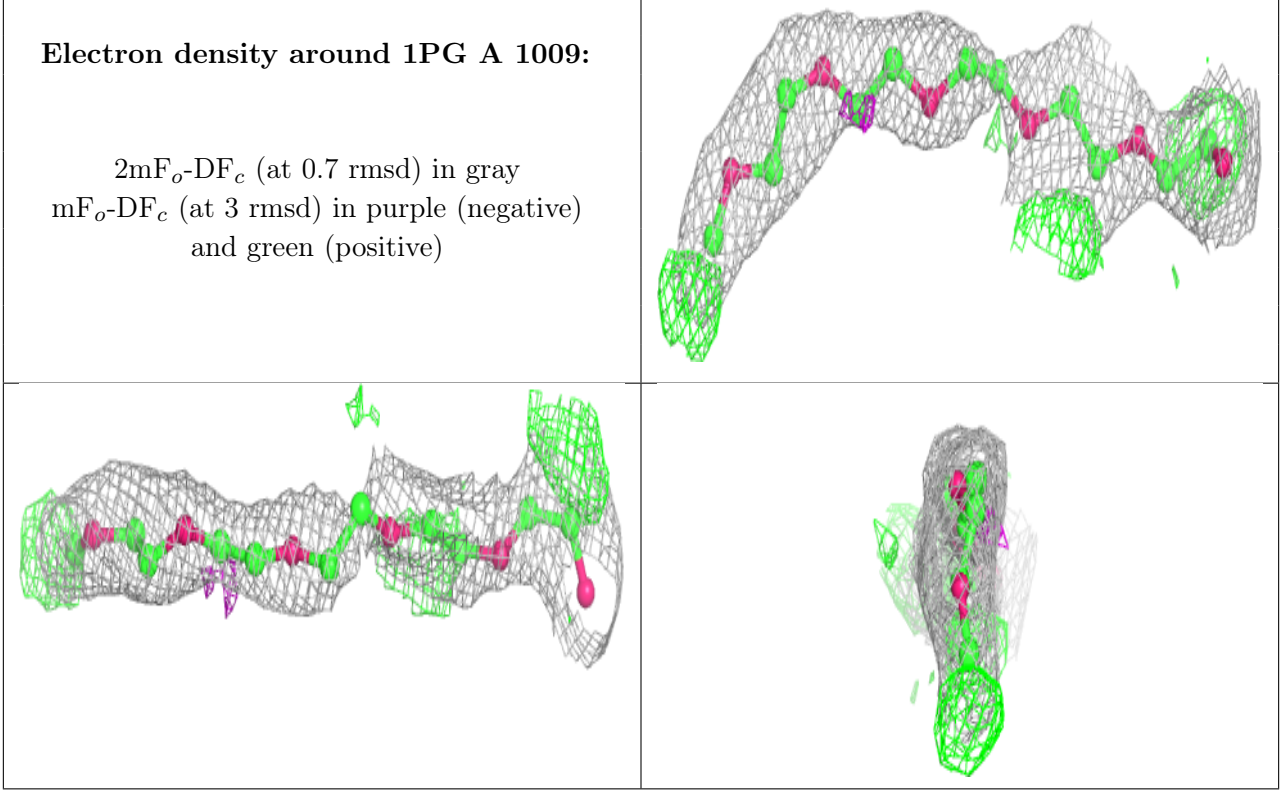
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
5	1PG	A	1009	17/17	0.72	0.27	48,59,78,81	0
5	1PG	A	1010	17/17	0.73	0.23	54,70,74,82	0
5	1PG	B	1009	17/17	0.75	0.32	52,64,78,82	0
5	1PG	A	1012	17/17	0.77	0.35	56,74,81,85	0
5	1PG	A	1014	10/17	0.78	0.46	61,65,75,79	0
3	MG	B	1006	1/1	0.79	0.11	54,54,54,54	0
5	1PG	A	1011	17/17	0.83	0.22	57,66,84,85	0
5	1PG	A	1013	17/17	0.85	0.36	48,58,69,73	0
3	MG	B	1007	1/1	0.90	0.08	56,56,56,56	0
3	MG	B	1003	1/1	0.91	0.11	52,52,52,52	0
4	K	B	1008	1/1	0.92	0.12	63,63,63,63	0
3	MG	A	1004	1/1	0.93	0.17	46,46,46,46	0
2	PO4	B	1001	5/5	0.93	0.10	49,52,61,65	0
3	MG	A	1006	1/1	0.94	0.14	42,42,42,42	0
3	MG	A	1007	1/1	0.95	0.12	42,42,42,42	0
3	MG	A	1005	1/1	0.95	0.21	37,37,37,37	0
2	PO4	A	1001	5/5	0.95	0.15	33,39,44,60	0
2	PO4	B	1002	5/5	0.97	0.09	45,49,54,55	0
3	MG	B	1005	1/1	0.97	0.09	50,50,50,50	0
3	MG	A	1003	1/1	0.97	0.14	36,36,36,36	0
4	K	A	1008	1/1	0.98	0.08	48,48,48,48	0
3	MG	B	1004	1/1	0.98	0.12	58,58,58,58	0
2	PO4	A	1002	5/5	0.98	0.17	31,37,40,41	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.

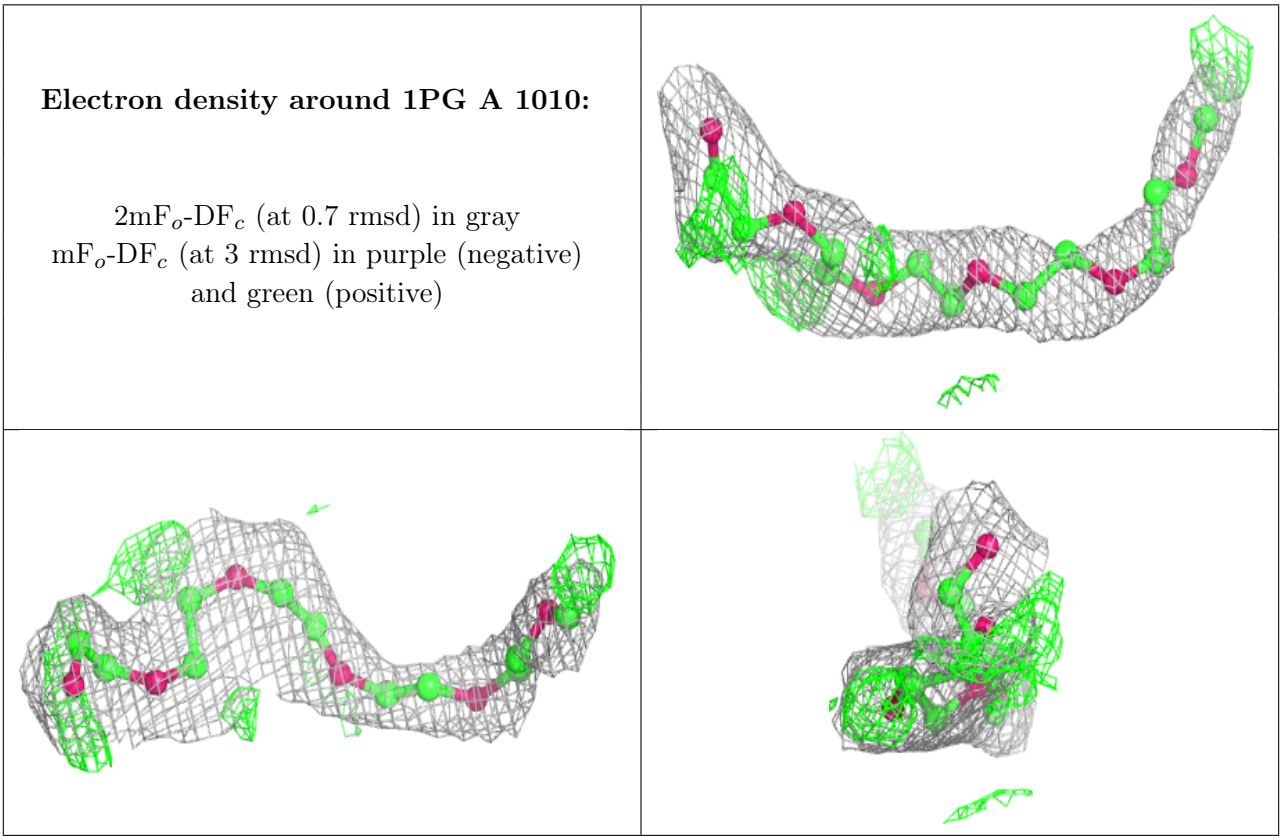
Electron density around 1PG A 1009:

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



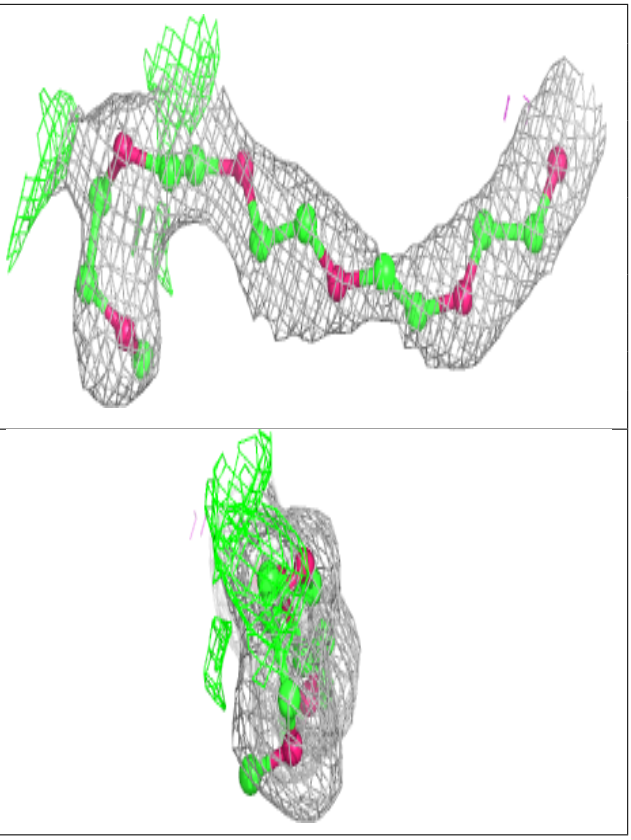
Electron density around 1PG A 1010:

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



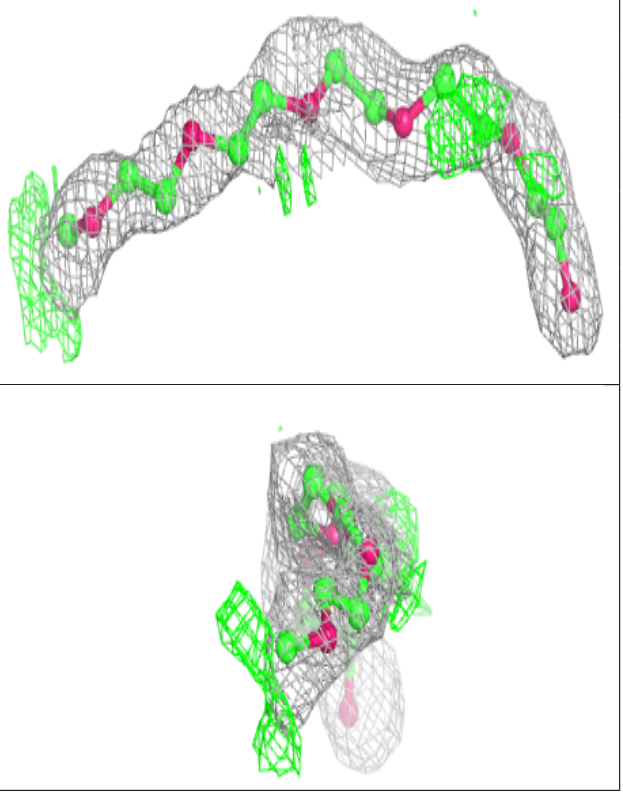
Electron density around 1PG B 1009:

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



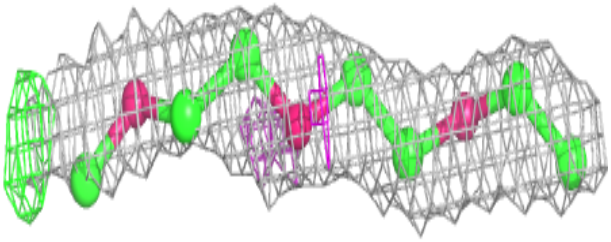
Electron density around 1PG A 1012:

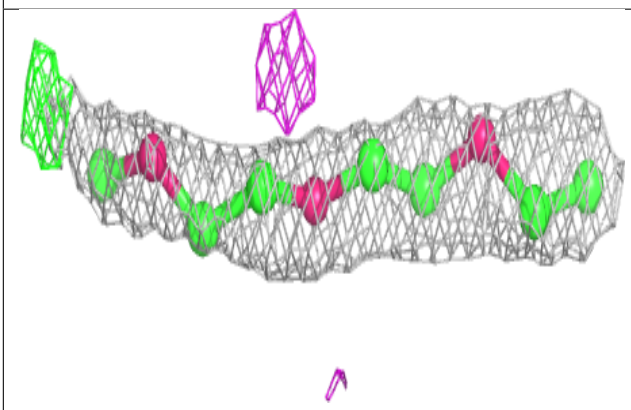
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

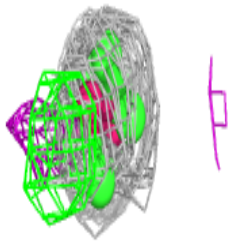


Electron density around 1PG A 1014:

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

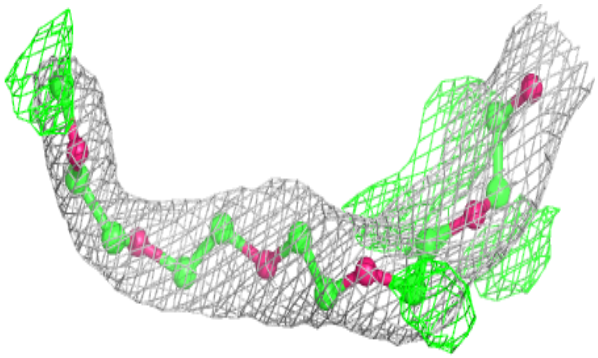


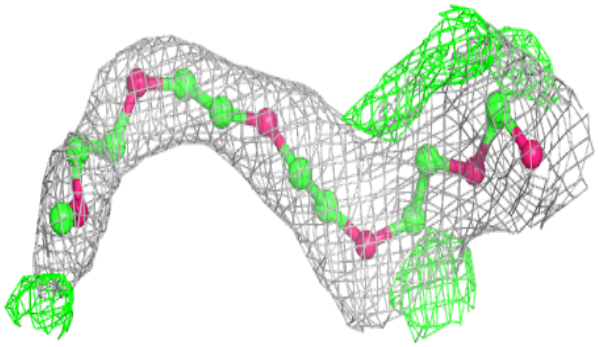


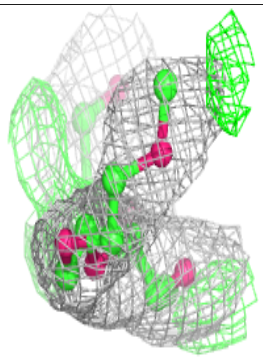


Electron density around 1PG A 1011:

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

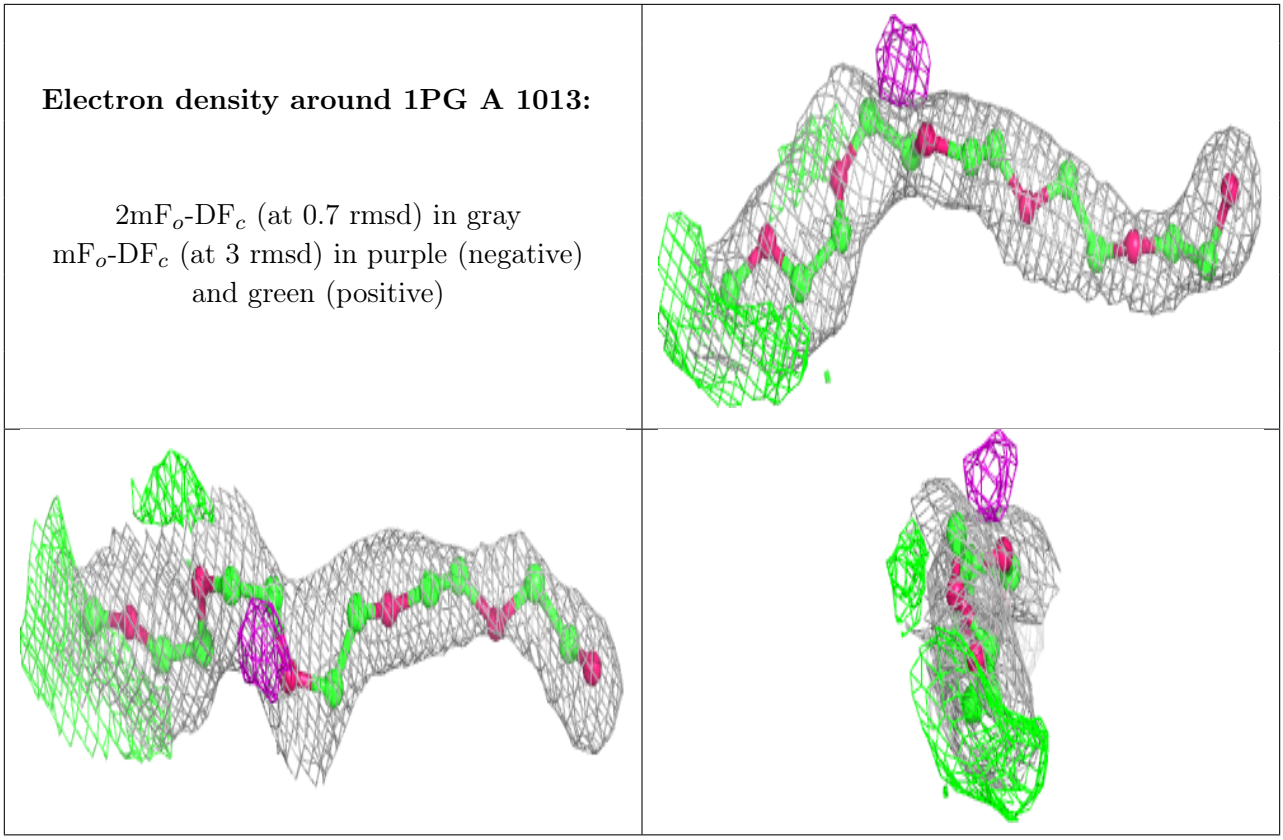






Electron density around 1PG A 1013:

$2mF_o$-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.