Full wwPDB X-ray Structure Validation Report (i)

Aug 5, 2026 – 10:37 PM EDT

PDB ID	:	3PUV / pdb_00003puv
Title	:	Crystal Structure of an outward-facing MBP-Maltose transporter complex bound to ADP-VO4
Authors	:	Oldham, M.L.; Chen, J.
Deposited on	:	2010-12-06
Resolution	:	2.40 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:

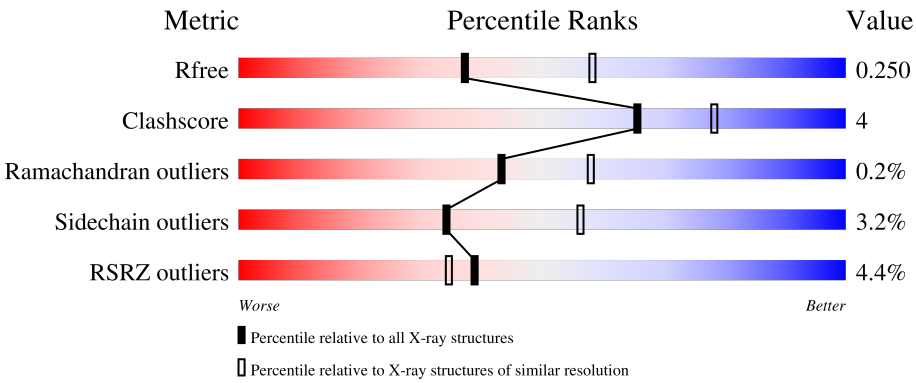
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

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 2.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	E	378	2%, 92%, 6% • •
2	F	514	6%, 85%, 9% • 5%
3	G	296	2%, 90%, 7% • •
4	A	381	2%, 85%, 11% • •
4	B	381	8%, 80%, 12% • 6%

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Mol	Chain	Length	Quality of chain
5	C	2	50% 50%

2 Entry composition (i)

There are 11 unique types of molecules in this entry. The entry contains 15093 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 Maltose-binding periplasmic protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	E	374	2901	1869	473	553	6	0	1	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	371	ALA	-	expression tag	UNP P0AEX9
E	372	SER	-	expression tag	UNP P0AEX9
E	373	ALA	-	expression tag	UNP P0AEX9
E	374	SER	-	expression tag	UNP P0AEX9
E	375	HIS	-	expression tag	UNP P0AEX9
E	376	HIS	-	expression tag	UNP P0AEX9
E	377	HIS	-	expression tag	UNP P0AEX9
E	378	HIS	-	expression tag	UNP P0AEX9

- Molecule 2 is a protein called Maltose transport system permease protein malF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	F	490	3821	2511	608	685	17	0	0	0

- Molecule 3 is a protein called Maltose transport system permease protein malG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	G	288	2221	1487	355	371	8	0	0	0

- Molecule 4 is a protein called Maltose/maltodextrin import ATP-binding protein MalK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	A	371	2910	1839	523	534	14	0	4	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	B	358	2798	1771	501	513	13	0	1	0

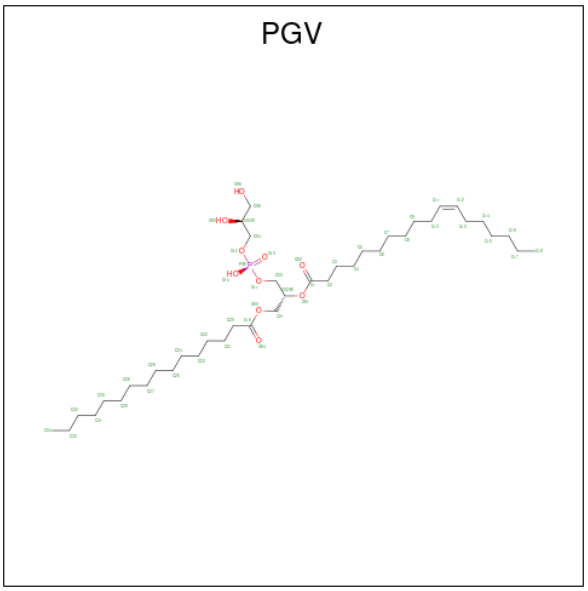
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	372	ALA	-	expression tag	UNP P68187
A	373	SER	-	expression tag	UNP P68187
A	374	ALA	-	expression tag	UNP P68187
A	375	SER	-	expression tag	UNP P68187
A	376	HIS	-	expression tag	UNP P68187
A	377	HIS	-	expression tag	UNP P68187
A	378	HIS	-	expression tag	UNP P68187
A	379	HIS	-	expression tag	UNP P68187
A	380	HIS	-	expression tag	UNP P68187
A	381	HIS	-	expression tag	UNP P68187
B	372	ALA	-	expression tag	UNP P68187
B	373	SER	-	expression tag	UNP P68187
B	374	ALA	-	expression tag	UNP P68187
B	375	SER	-	expression tag	UNP P68187
B	376	HIS	-	expression tag	UNP P68187
B	377	HIS	-	expression tag	UNP P68187
B	378	HIS	-	expression tag	UNP P68187
B	379	HIS	-	expression tag	UNP P68187
B	380	HIS	-	expression tag	UNP P68187
B	381	HIS	-	expression tag	UNP P68187

- Molecule 5 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.

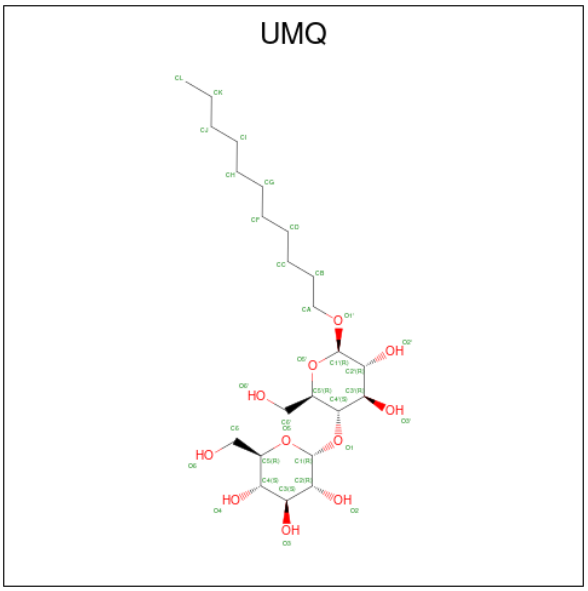
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
5	C	2	23	12	11	0	0	0

- Molecule 6 is (1R)-2-{[{[(2S)-2,3-DIHYDROXYPROPYL]OXY}(HYDROXY)PHOSPHORYL]OXY}-1-[(PALMITOYLOXY)METHYL]ETHYL (11E)-OCTADEC-11-ENOATE (CCD ID: PGV) (formula: $C_{40}H_{77}O_{10}P$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	F	1	Total C O P 51 40 10 1	0	0
6	F	1	Total C 9 9	0	0
6	F	1	Total C 8 8	0	0
6	F	1	Total C 13 13	0	0
6	G	1	Total C 8 8	0	0
6	G	1	Total C 12 12	0	0
6	G	1	Total C 10 10	0	0
6	G	1	Total C 7 7	0	0
6	G	1	Total C 9 9	0	0
6	G	1	Total C 12 12	0	0

- Molecule 7 is UNDECYL-MALTOSIDE (CCD ID: UMQ) (formula: $C_{23}H_{44}O_{11}$).

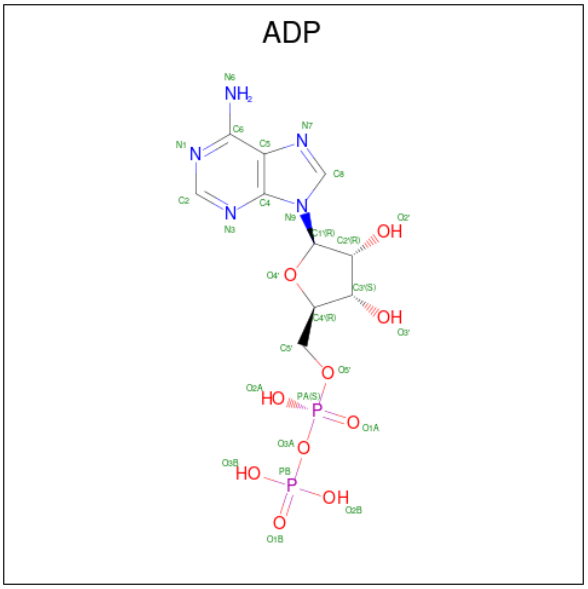


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
7	F	1	34	23	11	0	0

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

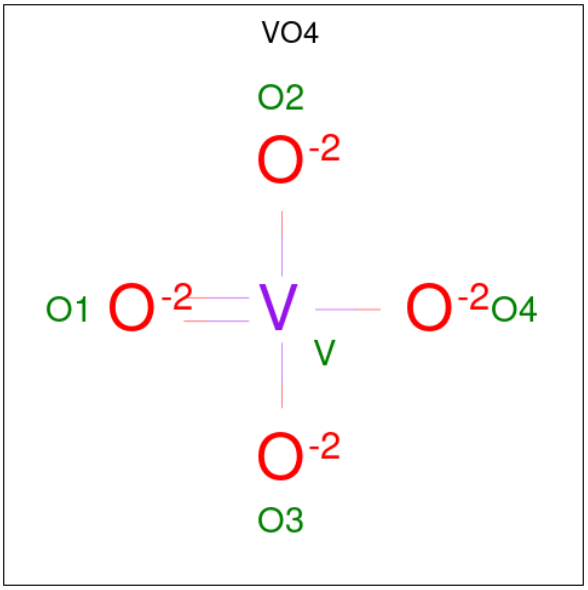
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
8	A	1	1	1	0	0
8	B	1	Total	Mg	0	0
			1	1		

- Molecule 9 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
9	A	1	27	10	5	10	2	0	0
			Total	C	N	O	P		
9	B	1	27	10	5	10	2	0	0

- Molecule 10 is VANADATE ION (CCD ID: VO4) (formula: O_4V).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	V		
10	A	1	5	4	1	0	0
			Total	O	V		
10	B	1	5	4	1	0	0

- Molecule 11 is water.

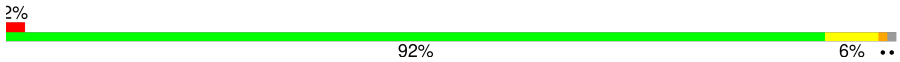
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
11	E	49	49	49	0	0
			Total	O		
11	F	34	34	34	0	0
			Total	O		
11	G	31	31	31	0	0
			Total	O		
11	A	38	38	38	0	0
			Total	O		
11	B	28	28	28	0	0

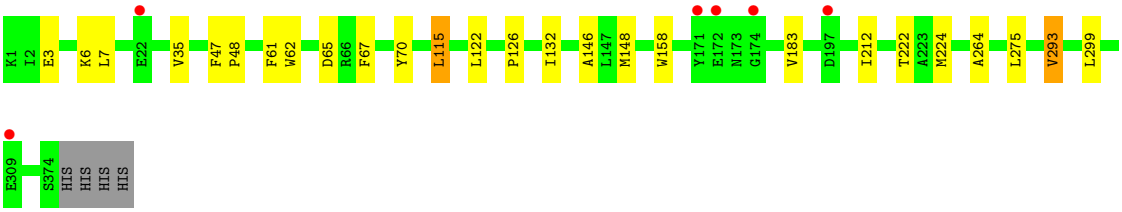
3 Residue-property plots (i)

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: Maltose-binding periplasmic protein

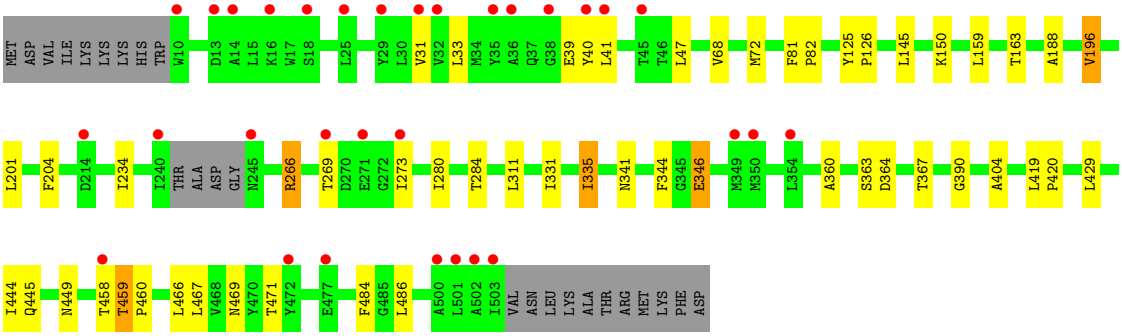
Chain E:





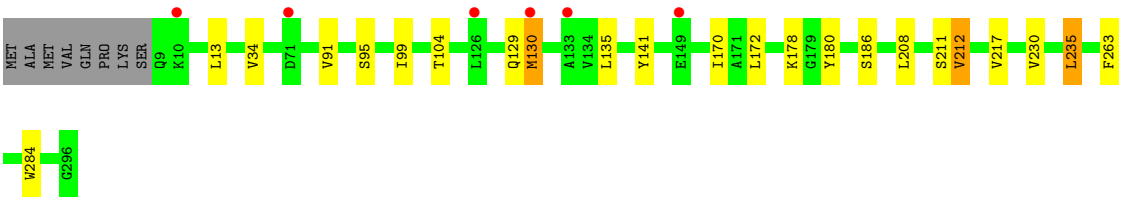
• Molecule 2: Maltose transport system permease protein malF

Chain F:



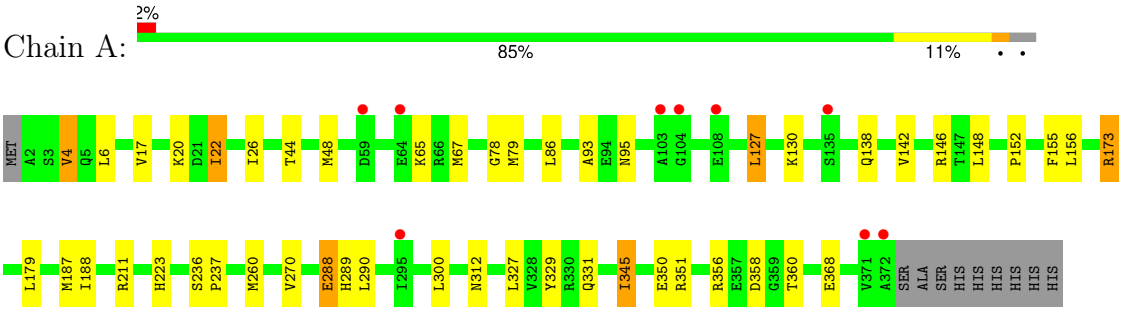
• Molecule 3: Maltose transport system permease protein malG

Chain G:



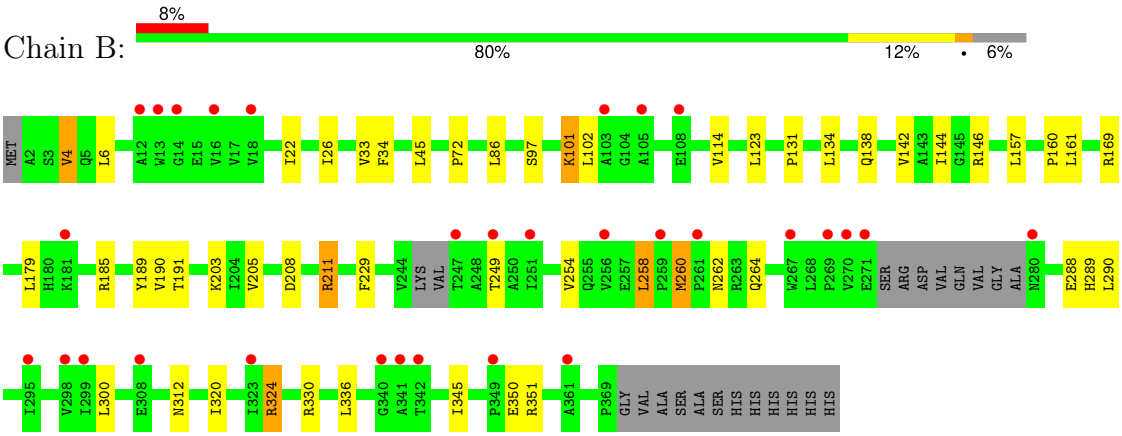
• Molecule 4: Maltose/maltodextrin import ATP-binding protein MalK

Chain A:



- Molecule 4: Maltose/maltodextrin import ATP-binding protein MalK

Chain B:



- Molecule 5: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain C:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	81.79Å 97.12Å 112.79Å 85.58° 78.71° 72.69°	Depositor
Resolution (Å)	20.00 – 2.40 20.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	63.2 (20.00-2.40) 63.1 (20.00-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.83 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.223 , 0.253 0.221 , 0.250	Depositor DCC
R_{free} test set	4002 reflections (3.16%)	wwPDB-VP
Wilson B-factor (Å^2)	41.9	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 38.1	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	0.92	EDS
Total number of atoms	15093	wwPDB-VP
Average B, all atoms (Å^2)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: UMQ, MG, PGV, VO4, GLC, ADP

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	E	0.43	1/2973 (0.0%)	0.79	0/4036
2	F	0.43	0/3916	0.79	2/5330 (0.0%)
3	G	0.45	0/2282	0.80	0/3119
4	A	0.43	0/2960	0.77	0/4012
4	B	0.43	0/2846	0.77	2/3856 (0.1%)
All	All	0.43	1/14977 (0.0%)	0.78	4/20353 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	132	ILE	CA-CB	6.43	1.57	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	459	THR	CA-C-N	5.90	127.22	119.84
2	F	459	THR	C-N-CA	5.90	127.22	119.84
4	B	260	MET	CA-C-N	5.18	125.48	119.47
4	B	260	MET	C-N-CA	5.18	125.48	119.47

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2901	0	2882	15	0
2	F	3821	0	3849	29	0
3	G	2221	0	2308	14	0
4	A	2910	0	2978	34	0
4	B	2798	0	2856	32	0
5	C	23	0	21	0	0
6	F	81	0	122	0	0
6	G	58	0	82	0	0
7	F	34	0	44	0	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
9	A	27	0	12	0	0
9	B	27	0	12	0	0
10	A	5	0	0	0	0
10	B	5	0	0	0	0
11	A	38	0	0	0	0
11	B	28	0	0	0	0
11	E	49	0	0	0	0
11	F	34	0	0	0	0
11	G	31	0	0	0	0
All	All	15093	0	15166	114	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 4.

All (114) 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)
4:A:173[A]:ARG:HB2	4:A:173[A]:ARG:NH1	1.54	1.22
4:B:211[A]:ARG:CG	4:B:211[A]:ARG:HH11	1.67	1.07
4:A:173[A]:ARG:CG	4:A:173[A]:ARG:HH11	1.67	1.05
4:A:173[A]:ARG:HH11	4:A:173[A]:ARG:CB	1.72	1.02
4:A:173[A]:ARG:NH1	4:A:173[A]:ARG:CB	2.24	1.00
2:F:266:ARG:HH11	2:F:266:ARG:HG2	1.30	0.96
4:A:173[A]:ARG:HH11	4:A:173[A]:ARG:HG3	1.33	0.91
4:B:211[A]:ARG:HH11	4:B:211[A]:ARG:HG3	1.37	0.89
4:B:211[A]:ARG:HH11	4:B:211[A]:ARG:HG2	1.40	0.86
4:A:173[A]:ARG:HB2	4:A:173[A]:ARG:HH11	1.28	0.84
3:G:91:VAL:O	3:G:95:SER:HB2	1.87	0.74
4:A:260[A]:MET:HE2	4:A:300:LEU:HD22	1.73	0.71
2:F:159:LEU:HD11	2:F:188:ALA:HB1	1.77	0.67
4:B:211[A]:ARG:HG2	4:B:211[A]:ARG:NH1	2.07	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:288:GLU:HG2	4:B:312:ASN:HB2	1.80	0.64
2:F:266:ARG:HG2	2:F:266:ARG:NH1	2.06	0.63
4:A:356:ARG:HH11	4:A:360:THR:HG23	1.65	0.61
4:A:223:HIS:CE1	4:A:368:GLU:HG2	2.36	0.60
4:A:329:TYR:CE2	4:A:331:GLN:HG2	2.37	0.60
4:B:33:VAL:HG12	4:B:189:TYR:HB3	1.84	0.60
2:F:444:ILE:HG13	2:F:466:LEU:HG	1.84	0.59
4:B:131:PRO:HA	4:B:134:LEU:HD12	1.86	0.57
2:F:68:VAL:HG12	2:F:72:MET:HG3	1.88	0.56
1:E:115:LEU:HD21	1:E:224:MET:HE3	1.88	0.56
4:A:79:MET:HG3	4:A:156:LEU:HB2	1.88	0.56
4:A:173[A]:ARG:NH1	4:A:173[A]:ARG:HG3	2.10	0.55
4:A:6:LEU:HD22	4:A:22:ILE:HD11	1.88	0.55
4:B:211[A]:ARG:HG3	4:B:211[A]:ARG:NH1	2.11	0.55
4:A:155:PHE:HB2	4:A:187:MET:HG2	1.87	0.55
4:A:20:LYS:HE2	4:A:211:ARG:HD3	1.90	0.54
4:A:173[A]:ARG:CB	4:A:173[A]:ARG:CZ	2.83	0.53
1:E:3:GLU:HB3	1:E:6:LYS:HE2	1.91	0.53
4:B:288:GLU:HG3	4:B:330:ARG:HD3	1.90	0.53
3:G:99:ILE:HG23	3:G:170:ILE:HG22	1.91	0.53
4:B:97:SER:HB3	4:B:114:VAL:HG21	1.90	0.52
2:F:429:LEU:HD23	3:G:172:LEU:HD22	1.92	0.52
1:E:293[A]:VAL:HG23	1:E:299:LEU:HD11	1.92	0.51
4:B:34:PHE:HB2	4:B:190:VAL:HG22	1.90	0.51
4:A:93:ALA:HB2	4:A:127:LEU:HG	1.92	0.51
4:A:260[A]:MET:SD	4:A:260[A]:MET:N	2.83	0.51
4:B:4:VAL:HG13	4:B:26:ILE:HB	1.92	0.51
4:A:4:VAL:HG13	4:A:26:ILE:HB	1.92	0.51
4:A:236:SER:HA	4:A:237:PRO:C	2.35	0.51
2:F:419:LEU:HB3	2:F:420:PRO:HD3	1.93	0.50
4:A:4:VAL:HG11	4:A:188:ILE:HD11	1.94	0.50
1:E:115:LEU:HD11	1:E:224:MET:HE3	1.93	0.49
1:E:146:ALA:O	1:E:224:MET:HG2	2.12	0.49
2:F:404:ALA:HB1	4:B:72:PRO:HB2	1.94	0.49
2:F:284:THR:HG22	2:F:466:LEU:HA	1.95	0.49
2:F:273:ILE:HG22	2:F:459:THR:HG21	1.95	0.48
4:B:144:ILE:HD11	4:B:160:PRO:O	2.13	0.48
1:E:122:LEU:HD21	1:E:126:PRO:HD3	1.95	0.48
3:G:180:TYR:CE2	3:G:211:SER:HA	2.48	0.48
4:B:6:LEU:HD22	4:B:22:ILE:HD11	1.96	0.47
4:B:324:ARG:HD3	4:B:324:ARG:H	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:312:ASN:HB2	4:B:288:GLU:HG2	1.96	0.47
4:A:358:ASP:OD1	4:A:360:THR:HG22	2.14	0.47
2:F:196:VAL:CG1	2:F:204:PHE:HB3	2.45	0.47
4:B:138:GLN:O	4:B:142:VAL:HG23	2.15	0.47
4:A:86:LEU:HA	4:A:146:ARG:HH21	1.80	0.47
3:G:129:GLN:HB3	3:G:130:MET:HE2	1.97	0.46
2:F:484:PHE:HB3	3:G:135:LEU:HD11	1.96	0.46
4:B:211[A]:ARG:HD2	4:B:211[A]:ARG:C	2.40	0.46
2:F:364:ASP:HB3	2:F:367:THR:OG1	2.14	0.46
2:F:360:ALA:HB1	2:F:363:SER:HB2	1.96	0.46
4:A:138:GLN:O	4:A:142:VAL:HG23	2.16	0.46
4:A:290:LEU:HD22	4:A:345:ILE:HD13	1.98	0.46
3:G:104:THR:HG21	3:G:208:LEU:HD21	1.99	0.45
2:F:39:GLU:O	2:F:41:LEU:N	2.50	0.45
4:B:249:THR:HG22	4:B:254:VAL:HG13	1.98	0.45
4:B:86:LEU:HA	4:B:146:ARG:NH2	2.31	0.45
2:F:335:ILE:HD13	3:G:34:VAL:HG22	1.99	0.45
4:B:290:LEU:HD22	4:B:345:ILE:HD13	1.99	0.44
1:E:62:TRP:HB3	1:E:67:PHE:HE1	1.82	0.44
2:F:331:ILE:HD12	3:G:263:PHE:HD2	1.83	0.44
1:E:61:PHE:CE2	1:E:264:ALA:HB2	2.52	0.44
4:B:34:PHE:HD2	4:B:205:VAL:HB	1.83	0.44
4:B:189:TYR:CE2	4:B:191:THR:HB	2.52	0.44
4:A:44:THR:HG22	4:A:48:MET:HE2	2.00	0.43
2:F:390:GLY:HA3	3:G:217:VAL:HG21	2.00	0.43
4:B:161:LEU:HB3	4:B:169:ARG:HG2	2.01	0.43
1:E:48:PRO:HG3	1:E:70:TYR:HE1	1.83	0.43
2:F:346:GLU:CD	2:F:346:GLU:H	2.25	0.43
2:F:341:ASN:O	2:F:344:PHE:O	2.37	0.43
4:B:260:MET:HG3	4:B:262:ASN:OD1	2.19	0.43
3:G:212:VAL:HG22	3:G:284:TRP:CE3	2.54	0.43
4:B:208:ASP:HB2	4:B:229:PHE:CE2	2.54	0.42
1:E:48:PRO:HG3	1:E:70:TYR:CE1	2.55	0.42
4:A:152:PRO:HD2	4:A:155:PHE:CE1	2.54	0.42
1:E:47:PHE:HB3	1:E:48:PRO:HD3	2.01	0.42
2:F:280:ILE:O	2:F:284:THR:HG23	2.20	0.42
2:F:81:PHE:HB3	2:F:82:PRO:HD3	2.02	0.42
1:E:7:LEU:HB2	1:E:35:VAL:HG22	2.01	0.42
4:B:157:LEU:HB3	4:B:160:PRO:HG3	2.01	0.42
3:G:180:TYR:HE2	3:G:211:SER:HA	1.85	0.42
4:A:148:LEU:HD22	4:A:179:LEU:HD22	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:284:THR:HG21	2:F:467:LEU:H	1.84	0.41
4:B:289:HIS:CG	4:B:351:ARG:HD2	2.55	0.41
2:F:266:ARG:HD3	2:F:486:LEU:HD13	2.02	0.41
2:F:445:GLN:HG2	2:F:469:ASN:ND2	2.35	0.41
1:E:275:LEU:HD23	2:F:201:LEU:HD11	2.02	0.41
3:G:141:TYR:HB2	3:G:235:LEU:HD11	2.01	0.41
4:A:78:GLY:HA3	4:A:152:PRO:HG3	2.02	0.41
2:F:125:TYR:HA	2:F:126:PRO:HD2	1.92	0.41
2:F:335:ILE:CD1	3:G:34:VAL:HG22	2.50	0.41
4:A:65:LYS:O	4:A:67:MET:HG2	2.21	0.41
4:B:211[A]:ARG:HD2	4:B:211[A]:ARG:O	2.20	0.41
2:F:31:VAL:HG11	2:F:47:LEU:HD13	2.02	0.41
1:E:148:MET:HB2	1:E:222:THR:HG21	2.03	0.41
1:E:158:TRP:HH2	1:E:183:VAL:HG12	1.84	0.41
4:B:258:LEU:HD23	4:B:264:GLN:HB3	2.02	0.41
4:A:95:ASN:O	4:A:146:ARG:HG3	2.22	0.40
4:A:289:HIS:CG	4:A:351:ARG:HD2	2.56	0.40
4:B:97:SER:O	4:B:101:LYS:HB2	2.22	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	E	373/378 (99%)	358 (96%)	15 (4%)	0	100	100
2	F	486/514 (95%)	468 (96%)	16 (3%)	2 (0%)	30	43
3	G	286/296 (97%)	282 (99%)	3 (1%)	1 (0%)	36	50
4	A	373/381 (98%)	363 (97%)	10 (3%)	0	100	100
4	B	353/381 (93%)	337 (96%)	16 (4%)	0	100	100
All	All	1871/1950 (96%)	1808 (97%)	60 (3%)	3 (0%)	43	58

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	40	TYR
3	G	230	VAL
2	F	460	PRO

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	E	298/303 (98%)	293 (98%)	5 (2%)	53	74
2	F	402/424 (95%)	388 (96%)	14 (4%)	32	53
3	G	230/237 (97%)	224 (97%)	6 (3%)	40	63
4	A	318/323 (98%)	306 (96%)	12 (4%)	29	49
4	B	306/323 (95%)	290 (95%)	16 (5%)	21	36
All	All	1554/1610 (96%)	1501 (97%)	53 (3%)	34	54

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	65	ASP
1	E	115	LEU
1	E	212	ILE
1	E	293[A]	VAL
1	E	293[B]	VAL
2	F	33	LEU
2	F	145	LEU
2	F	150	LYS
2	F	163	THR
2	F	196	VAL
2	F	234	ILE
2	F	266	ARG
2	F	269	THR
2	F	311	LEU
2	F	335	ILE

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Mol	Chain	Res	Type
2	F	346	GLU
2	F	449	ASN
2	F	458	THR
2	F	471	THR
3	G	13	LEU
3	G	130	MET
3	G	178	LYS
3	G	186	SER
3	G	212	VAL
3	G	235	LEU
4	A	4	VAL
4	A	17	VAL
4	A	22	ILE
4	A	127	LEU
4	A	130	LYS
4	A	173[A]	ARG
4	A	173[B]	ARG
4	A	270	VAL
4	A	288	GLU
4	A	327	LEU
4	A	345	ILE
4	A	350	GLU
4	B	4	VAL
4	B	45	LEU
4	B	101	LYS
4	B	102	LEU
4	B	123	LEU
4	B	179	LEU
4	B	185	ARG
4	B	203	LYS
4	B	211[A]	ARG
4	B	211[B]	ARG
4	B	258	LEU
4	B	300	LEU
4	B	320	ILE
4	B	324	ARG
4	B	336	LEU
4	B	350	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	49	GLN
1	E	218	ASN
1	E	349	ASN
1	E	355	GLN
2	F	115	GLN
2	F	232	ASN
2	F	437	ASN
2	F	440	ASN
3	G	282	GLN
4	A	82	GLN
4	A	262	ASN
4	A	305	GLN
4	A	309	GLN
4	A	319	GLN
4	A	326	ASN
4	B	111	ASN
4	B	331	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

2 monosaccharides 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	$\#\|Z\| > 2$
5	GLC	C	1	5	12,12,12	0.50	0	17,17,17	0.73	0
5	GLC	C	2	5	11,11,12	0.40	0	15,15,17	0.87	1 (6%)

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	GLC	C	1	5	-	0/2/22/22	0/1/1/1
5	GLC	C	2	5	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	C	2	GLC	C1-O5-C5	2.25	115.20	112.19

There are no chirality outliers.

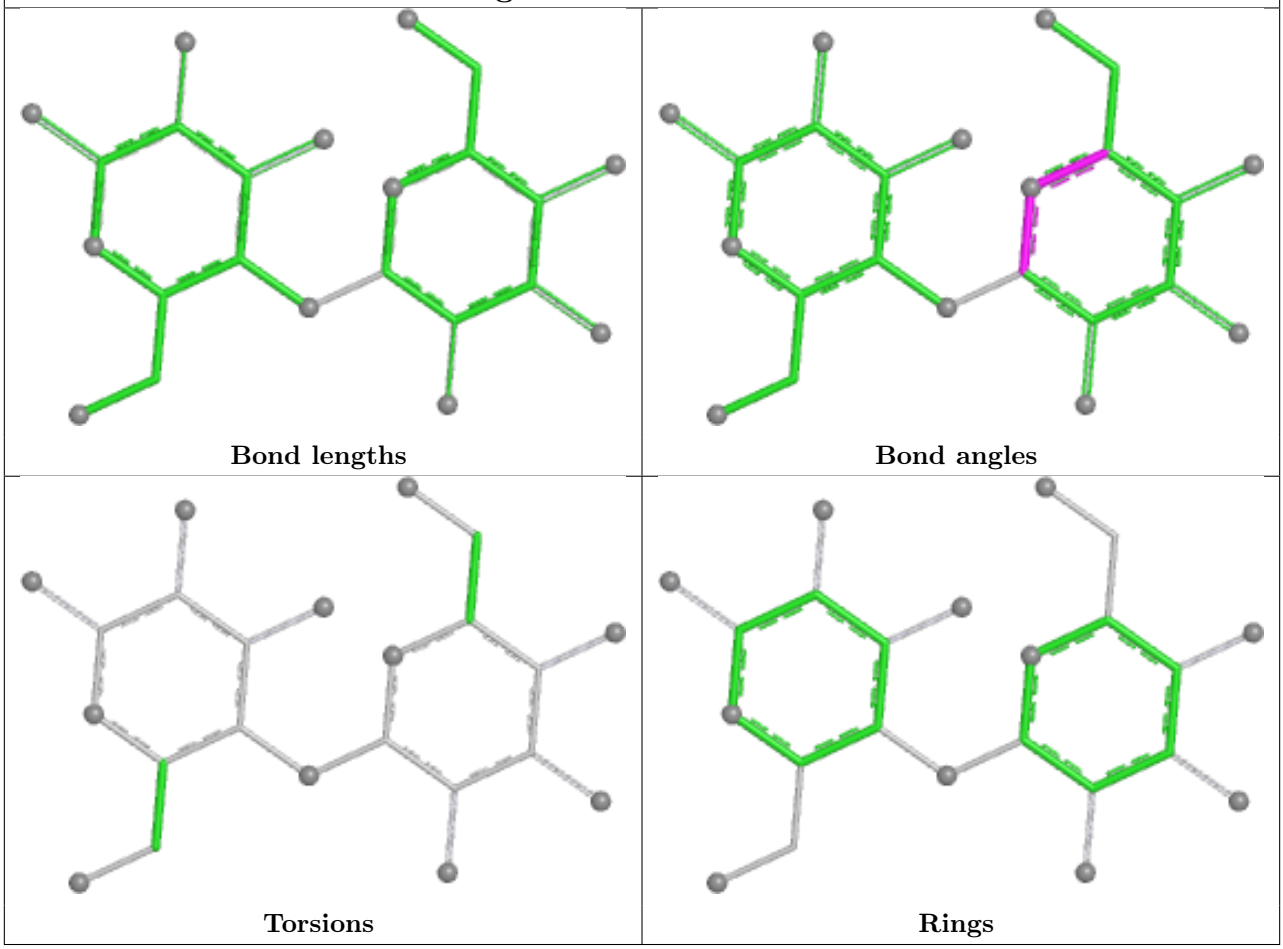
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain C



5.6 Ligand geometry (i)

Of 17 ligands modelled in this entry, 2 are monoatomic - leaving 15 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	$\#\|Z\| > 2$
6	PGV	F	4001	-	50,50,50	1.09	3 (6%)	53,56,56	1.03	3 (5%)
9	ADP	B	2502	10,8	28,29,29	0.98	2 (7%)	43,45,45	1.78	8 (18%)
6	PGV	F	4008	-	7,7,50	0.30	0	6,6,56	0.43	0
6	PGV	G	4004	-	11,11,50	1.04	1 (9%)	10,10,56	1.11	1 (10%)
6	PGV	G	4003	-	7,7,50	0.28	0	6,6,56	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
6	PGV	G	4007	-	8,8,50	1.20	1 (12%)	7,7,56	1.29	2 (28%)
10	VO4	A	3001	9,8	0,4,4	-	-	-		
6	PGV	G	4005	-	9,9,50	1.14	1 (11%)	8,8,56	1.19	1 (12%)
9	ADP	A	2501	10,8	28,29,29	0.98	2 (7%)	43,45,45	1.74	8 (18%)
6	PGV	F	4002	-	8,8,50	0.30	0	7,7,56	0.46	0
6	PGV	G	4009	-	11,11,50	1.20	1 (9%)	10,10,56	0.83	0
10	VO4	B	3002	9,8	0,4,4	-	-	-		
7	UMQ	F	5004	-	35,35,35	0.44	0	46,46,46	0.68	0
6	PGV	G	4006	-	6,6,50	1.39	1 (16%)	5,5,56	1.47	1 (20%)
6	PGV	F	4010	-	12,12,50	1.15	1 (8%)	11,11,56	1.03	1 (9%)

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
6	PGV	F	4001	-	-	28/55/55/55	-
9	ADP	B	2502	10,8	-	3/16/32/32	0/3/3/3
6	PGV	F	4008	-	-	2/5/5/55	-
6	PGV	G	4004	-	-	2/9/9/55	-
6	PGV	G	4003	-	-	1/5/5/55	-
6	PGV	G	4007	-	-	3/6/6/55	-
6	PGV	G	4005	-	-	4/7/7/55	-
9	ADP	A	2501	10,8	-	2/16/32/32	0/3/3/3
6	PGV	F	4002	-	-	4/6/6/55	-
6	PGV	G	4009	-	-	7/9/9/55	-
7	UMQ	F	5004	-	-	4/20/60/60	0/2/2/2
6	PGV	G	4006	-	-	2/4/4/55	-
6	PGV	F	4010	-	-	7/10/10/55	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	F	4001	PGV	O01-C1	4.26	1.46	1.34
6	F	4001	PGV	O03-C19	4.21	1.45	1.33
6	G	4009	PGV	C12-C11	3.81	1.53	1.31
6	F	4010	PGV	C12-C11	3.79	1.53	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	F	4001	PGV	C12-C11	3.77	1.53	1.31
6	G	4006	PGV	C12-C11	3.24	1.53	1.29
6	G	4005	PGV	C11-C12	3.22	1.53	1.29
6	G	4004	PGV	C11-C12	3.22	1.52	1.29
6	G	4007	PGV	C11-C12	3.21	1.52	1.29
9	A	2501	ADP	C8-N7	2.18	1.35	1.31
9	B	2502	ADP	C5-N7	-2.18	1.35	1.39
9	A	2501	ADP	C5-N7	-2.11	1.35	1.39
9	B	2502	ADP	C8-N7	2.07	1.35	1.31

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	B	2502	ADP	C5-C4-N3	-5.25	119.48	126.72
9	B	2502	ADP	N3-C2-N1	-5.10	120.86	128.58
9	A	2501	ADP	C5-C4-N3	-4.98	119.87	126.72
9	A	2501	ADP	N3-C2-N1	-4.97	121.06	128.58
6	F	4001	PGV	O01-C1-C2	4.04	120.23	111.48
9	B	2502	ADP	C2-N3-C4	3.60	120.62	111.83
9	B	2502	ADP	N3-C4-N9	3.46	133.06	127.17
9	A	2501	ADP	C2-N3-C4	3.39	120.12	111.83
9	A	2501	ADP	N9-C8-N7	-3.39	109.12	113.94
9	A	2501	ADP	N3-C4-N9	3.35	132.87	127.17
9	B	2502	ADP	N9-C8-N7	-3.29	109.26	113.94
6	F	4001	PGV	O03-C19-C20	2.90	120.67	111.83
6	G	4006	PGV	C13-C12-C11	-2.62	110.65	130.48
6	G	4007	PGV	C10-C11-C12	-2.58	110.96	130.48
6	G	4004	PGV	C10-C11-C12	-2.56	111.12	130.48
9	A	2501	ADP	C4-N9-C8	2.54	108.40	105.74
6	G	4005	PGV	C10-C11-C12	-2.51	111.49	130.48
9	B	2502	ADP	C4-C5-N7	-2.49	107.74	110.58
9	B	2502	ADP	C5-N7-C8	2.48	107.35	103.45
9	A	2501	ADP	C5-N7-C8	2.44	107.28	103.45
6	F	4010	PGV	C13-C12-C11	-2.43	111.78	126.42
9	B	2502	ADP	C4-N9-C8	2.34	108.19	105.74
9	A	2501	ADP	C4-C5-N7	-2.32	107.93	110.58
6	F	4001	PGV	O03-C19-O04	-2.29	117.89	123.63
6	G	4007	PGV	C13-C12-C11	-2.01	110.74	126.43

There are no chirality outliers.

All (69) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	F	4001	PGV	C03-O11-P-O12
6	F	4001	PGV	C03-O11-P-O13
6	F	4001	PGV	C04-O12-P-O11
6	F	4001	PGV	C04-O12-P-O14
6	F	4001	PGV	C2-C1-O01-C02
6	G	4009	PGV	C12-C13-C14-C15
9	A	2501	ADP	PA-O3A-PB-O2B
9	B	2502	ADP	PA-O3A-PB-O2B
9	B	2502	ADP	PA-O3A-PB-O3B
6	F	4001	PGV	O02-C1-O01-C02
6	F	4001	PGV	C10-C11-C12-C13
6	F	4010	PGV	C10-C11-C12-C13
6	F	4001	PGV	O04-C19-O03-C01
6	F	4001	PGV	C20-C19-O03-C01
6	G	4009	PGV	C10-C11-C12-C13
6	F	4001	PGV	C04-C05-C06-O06
6	F	4001	PGV	C30-C31-C32-C33
7	F	5004	UMQ	CC-CD-CF-CG
6	F	4010	PGV	C6-C7-C8-C9
6	F	4001	PGV	C2-C3-C4-C5
7	F	5004	UMQ	CF-CG-CH-CI
6	F	4001	PGV	C29-C30-C31-C32
6	F	4002	PGV	C4-C5-C6-C7
6	G	4005	PGV	C6-C7-C8-C9
6	F	4001	PGV	C6-C7-C8-C9
6	F	4001	PGV	C20-C21-C22-C23
6	F	4001	PGV	C11-C10-C9-C8
6	F	4010	PGV	C11-C10-C9-C8
6	F	4001	PGV	C24-C25-C26-C27
7	F	5004	UMQ	CD-CF-CG-CH
6	F	4008	PGV	C4-C5-C6-C7
6	G	4009	PGV	C11-C10-C9-C8
6	F	4001	PGV	C22-C23-C24-C25
6	F	4010	PGV	C5-C6-C7-C8
6	F	4001	PGV	C19-C20-C21-C22
6	F	4002	PGV	C6-C7-C8-C9
6	F	4001	PGV	C13-C14-C15-C16
6	F	4008	PGV	C5-C6-C7-C8
6	G	4005	PGV	C4-C5-C6-C7
6	G	4009	PGV	C7-C8-C9-C10
6	F	4002	PGV	C2-C3-C4-C5
6	G	4004	PGV	C10-C11-C12-C13
6	G	4005	PGV	C10-C11-C12-C13

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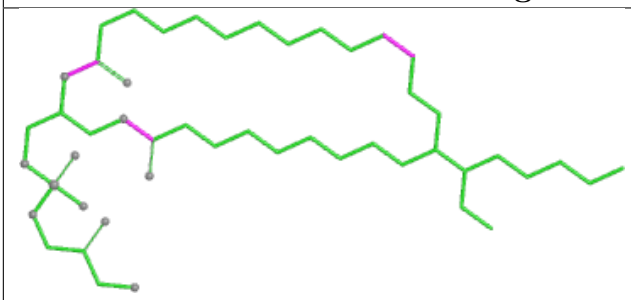
Mol	Chain	Res	Type	Atoms
6	G	4006	PGV	C10-C11-C12-C13
6	G	4007	PGV	C10-C11-C12-C13
6	G	4003	PGV	C6-C7-C8-C9
6	G	4007	PGV	C6-C7-C8-C9
6	G	4007	PGV	C5-C6-C7-C8
6	G	4005	PGV	C7-C8-C9-C10
6	F	4001	PGV	O05-C05-C06-O06
6	G	4009	PGV	C4-C5-C6-C7
6	G	4004	PGV	C11-C10-C9-C8
6	F	4001	PGV	C27-C28-C29-C30
7	F	5004	UMQ	CB-CC-CD-CF
6	F	4001	PGV	C03-O11-P-O14
6	F	4010	PGV	C9-C10-C11-C12
6	F	4001	PGV	O03-C01-C02-O01
6	G	4009	PGV	C11-C12-C13-C14
6	G	4009	PGV	C9-C10-C11-C12
6	F	4002	PGV	C3-C4-C5-C6
9	A	2501	ADP	PA-O3A-PB-O3B
6	F	4001	PGV	O03-C01-C02-C03
6	F	4001	PGV	C3-C4-C5-C6
6	F	4001	PGV	C26-C27-C28-C29
6	G	4006	PGV	C11-C12-C13-C14
9	B	2502	ADP	PA-O3A-PB-O1B
6	F	4010	PGV	C11-C12-C13-C14
6	F	4010	PGV	C4-C5-C6-C7
6	F	4001	PGV	C7-C8-C9-C10

There are no ring outliers.

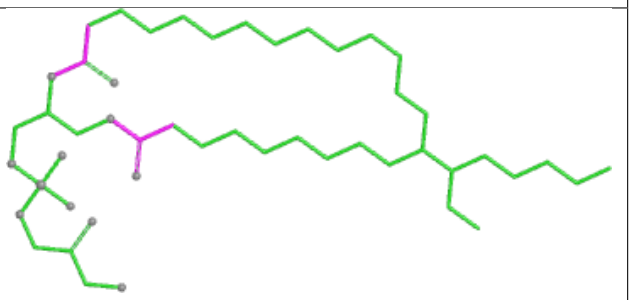
No monomer is involved in short contacts.

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.

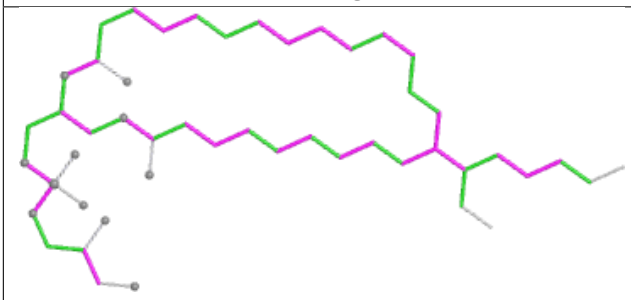
Ligand PGV F 4001



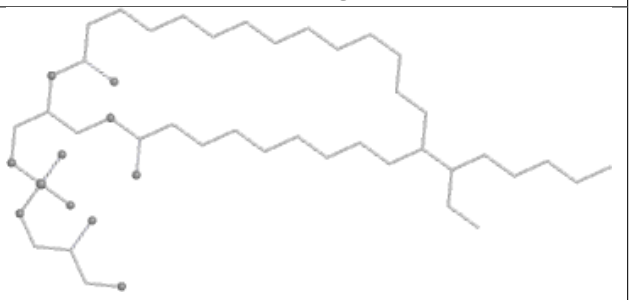
Bond lengths



Bond angles

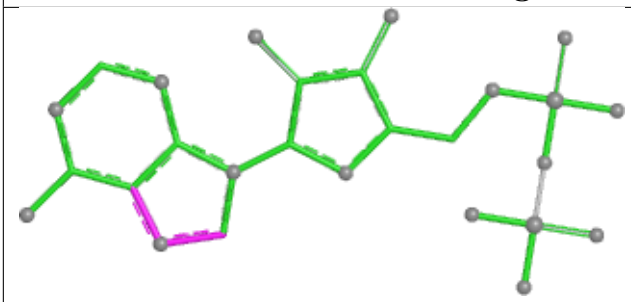


Torsions

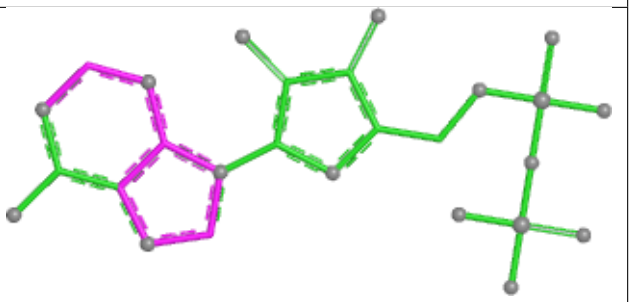


Rings

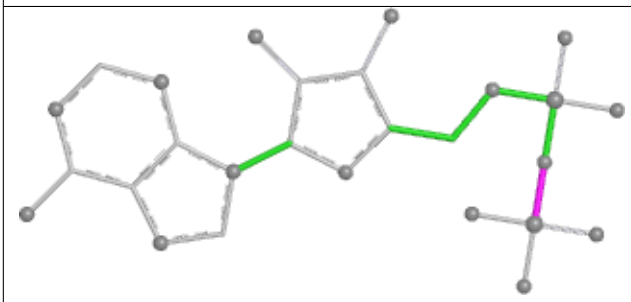
Ligand ADP B 2502



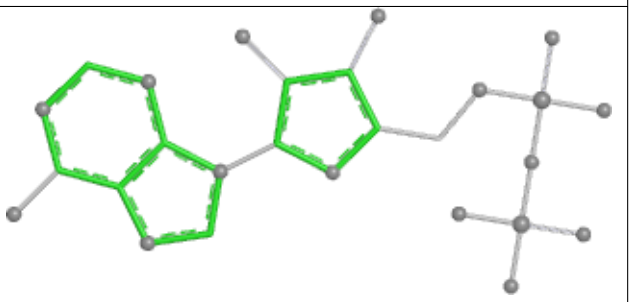
Bond lengths



Bond angles



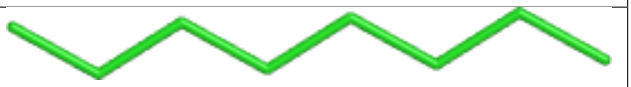
Torsions



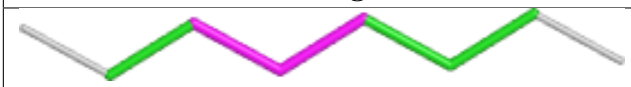
Rings

Ligand PGV F 4008

Bond lengths



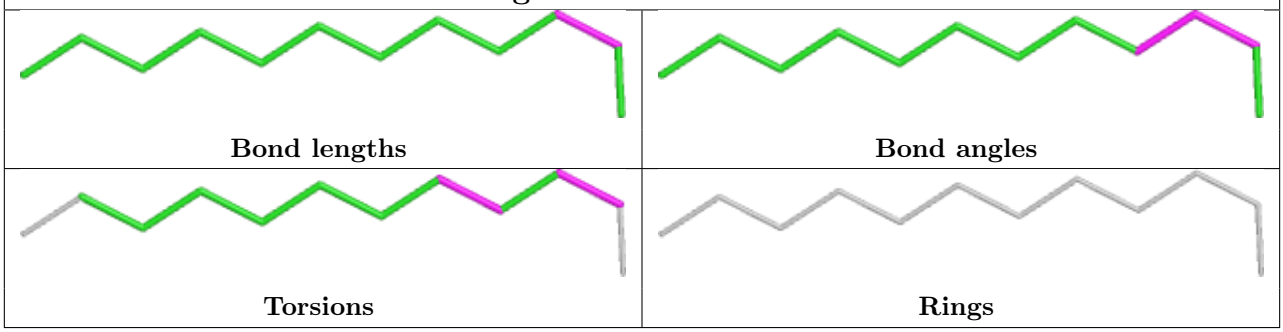
Bond angles



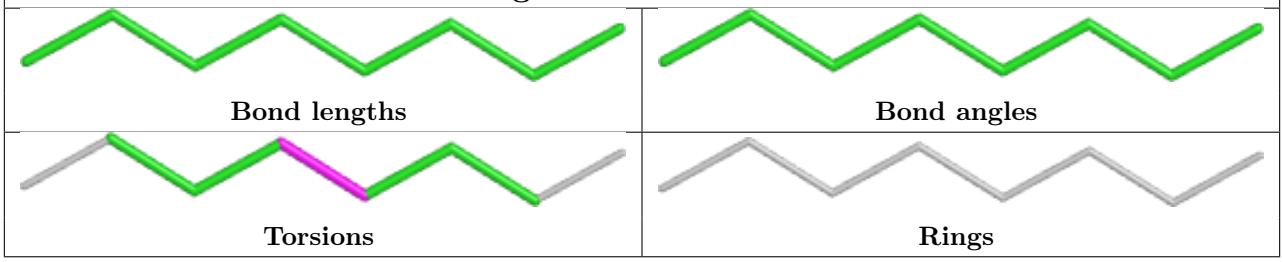
Torsions

Rings

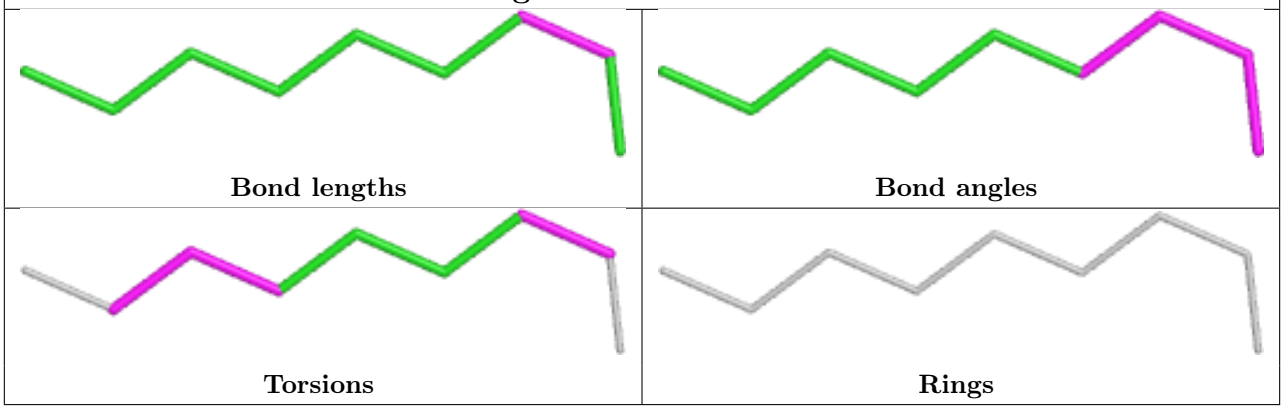
Ligand PGV G 4004



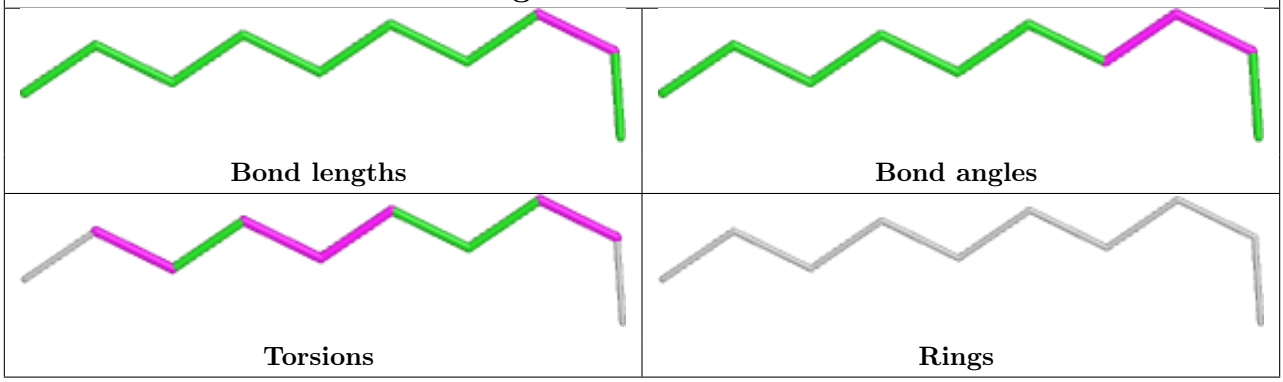
Ligand PGV G 4003



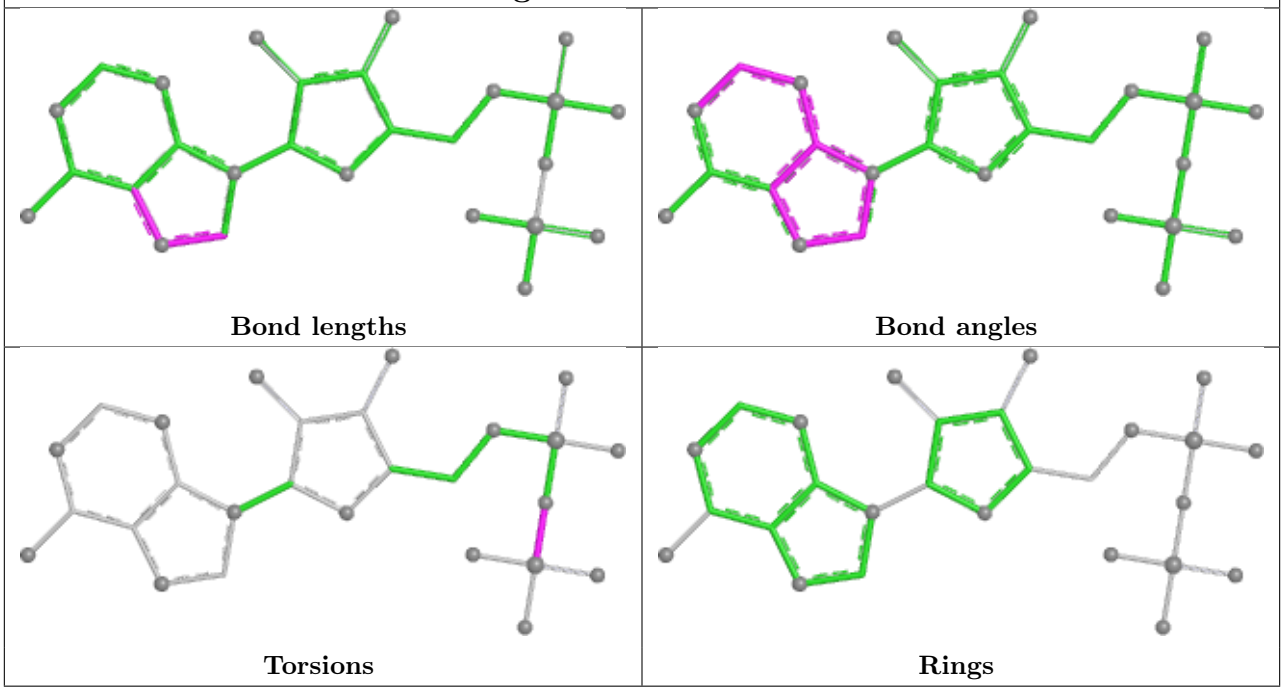
Ligand PGV G 4007



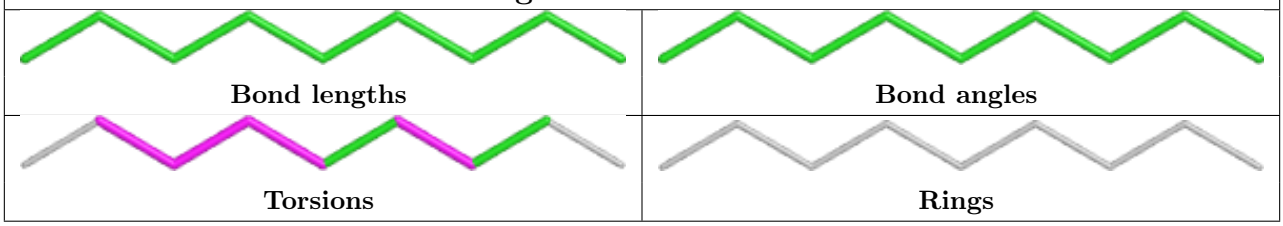
Ligand PGV G 4005



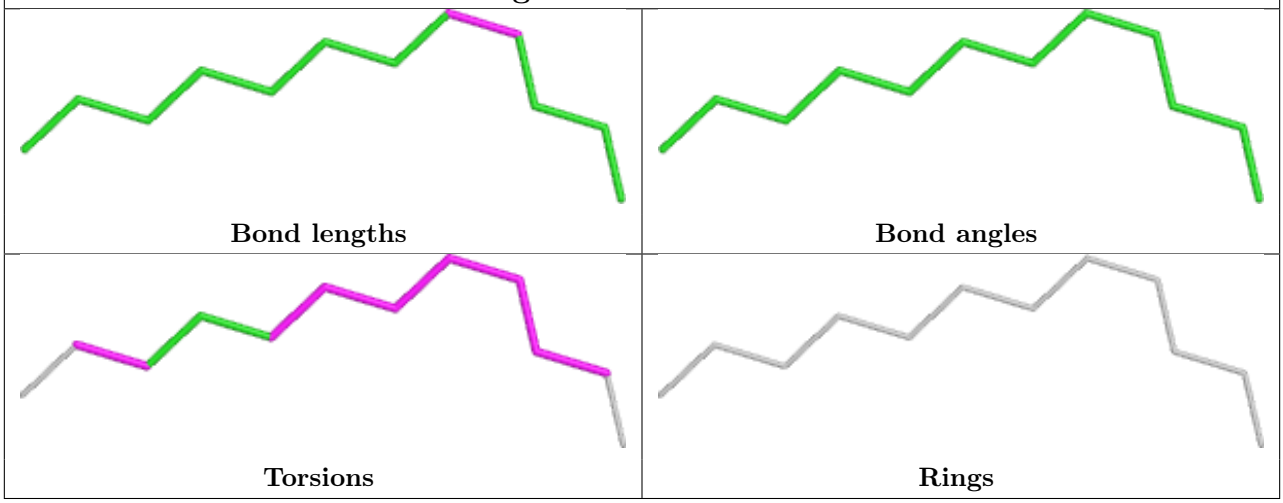
Ligand ADP A 2501



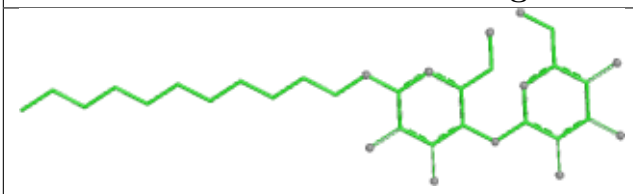
Ligand PGV F 4002



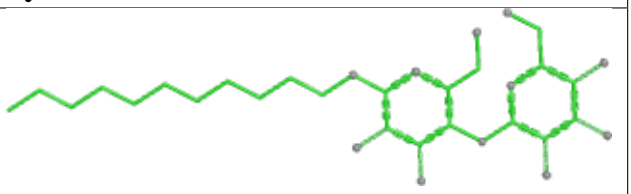
Ligand PGV G 4009



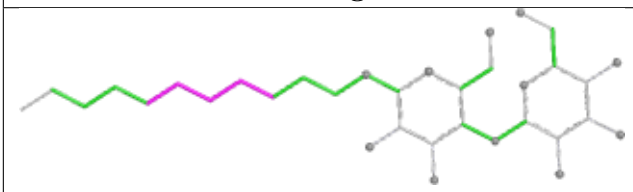
Ligand UMQ F 5004



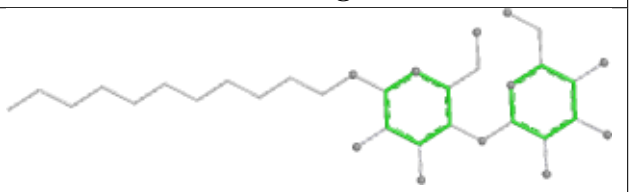
Bond lengths



Bond angles

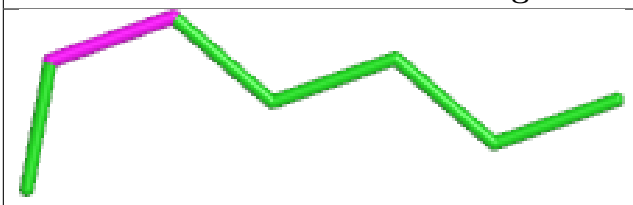


Torsions

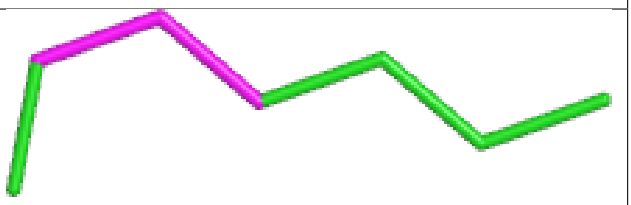


Rings

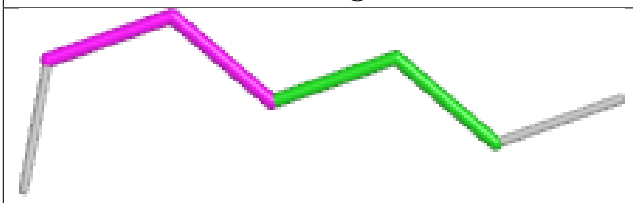
Ligand PGV G 4006



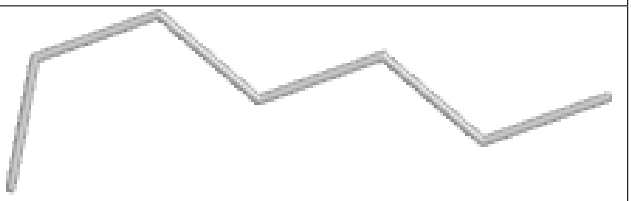
Bond lengths



Bond angles



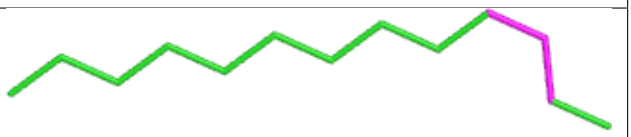
Torsions



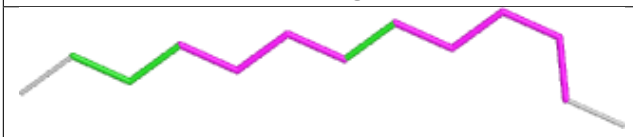
Rings

Ligand PGV F 4010

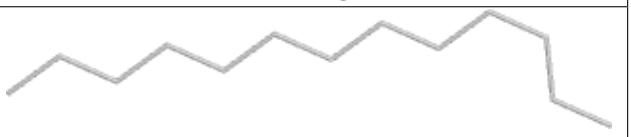
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	E	374/378 (98%)	0.18	6 (1%)	70	66	26, 60, 97, 110	2 (0%)
2	F	490/514 (95%)	0.46	31 (6%)	26	22	24, 62, 115, 163	0
3	G	288/296 (97%)	0.02	6 (2%)	63	59	17, 39, 77, 102	0
4	A	371/381 (97%)	0.19	9 (2%)	59	55	11, 54, 89, 117	4 (1%)
4	B	358/381 (93%)	0.55	30 (8%)	17	14	23, 69, 146, 171	1 (0%)
All	All	1881/1950 (96%)	0.30	82 (4%)	39	34	11, 58, 109, 171	7 (0%)

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	10	TRP	7.7
4	A	104	GLY	5.4
4	A	295	ILE	5.3
4	B	249	THR	4.4
4	B	361	ALA	4.2
3	G	133	ALA	4.1
4	B	251	ILE	4.0
4	A	103	ALA	3.9
2	F	16	LYS	3.7
4	B	323	ILE	3.6
4	B	259	PRO	3.5
4	A	372	ALA	3.5
2	F	45	THR	3.4
2	F	14	ALA	3.4
2	F	271	GLU	3.2
2	F	503	ILE	3.2
2	F	40	TYR	3.2
4	B	261	PRO	3.1
2	F	214	ASP	3.1
4	B	271	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
4	B	247	THR	2.9
2	F	472	TYR	2.9
4	B	299	ILE	2.9
4	B	270	VAL	2.8
4	B	13	TRP	2.8
1	E	174	GLY	2.8
2	F	354	LEU	2.8
3	G	149	GLU	2.8
3	G	71	ASP	2.7
2	F	273	ILE	2.7
4	A	371	VAL	2.7
2	F	501	LEU	2.7
4	B	256	VAL	2.7
2	F	29	TYR	2.6
4	B	103	ALA	2.6
4	B	295	ILE	2.6
1	E	197	ASP	2.6
4	A	135	SER	2.6
4	B	341	ALA	2.5
2	F	32	VAL	2.5
4	B	269	PRO	2.5
2	F	25	LEU	2.5
3	G	10	LYS	2.5
2	F	36	ALA	2.5
3	G	130	MET	2.5
4	A	108	GLU	2.4
2	F	269	THR	2.4
2	F	458	THR	2.4
4	B	267	TRP	2.4
4	B	298	VAL	2.4
2	F	500	ALA	2.4
4	B	342	THR	2.4
2	F	38	GLY	2.4
4	B	340	GLY	2.4
1	E	309	GLU	2.3
4	B	308	GLU	2.3
4	B	105	ALA	2.3
3	G	126	LEU	2.3
2	F	35	TYR	2.3
4	B	181	LYS	2.2
4	B	16	VAL	2.2
2	F	18	SER	2.2

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Mol	Chain	Res	Type	RSRZ
4	B	14	GLY	2.2
4	A	59	ASP	2.2
2	F	41	LEU	2.2
2	F	349	MET	2.1
4	B	12	ALA	2.1
4	A	64	GLU	2.1
2	F	477	GLU	2.1
2	F	13	ASP	2.1
4	B	18	VAL	2.1
1	E	22	GLU	2.1
2	F	240	ILE	2.1
2	F	31	VAL	2.0
1	E	172	GLU	2.0
4	B	349	PRO	2.0
2	F	350	MET	2.0
2	F	502	ALA	2.0
4	B	108	GLU	2.0
2	F	245	ASN	2.0
4	B	280	ASN	2.0
1	E	171	TYR	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)

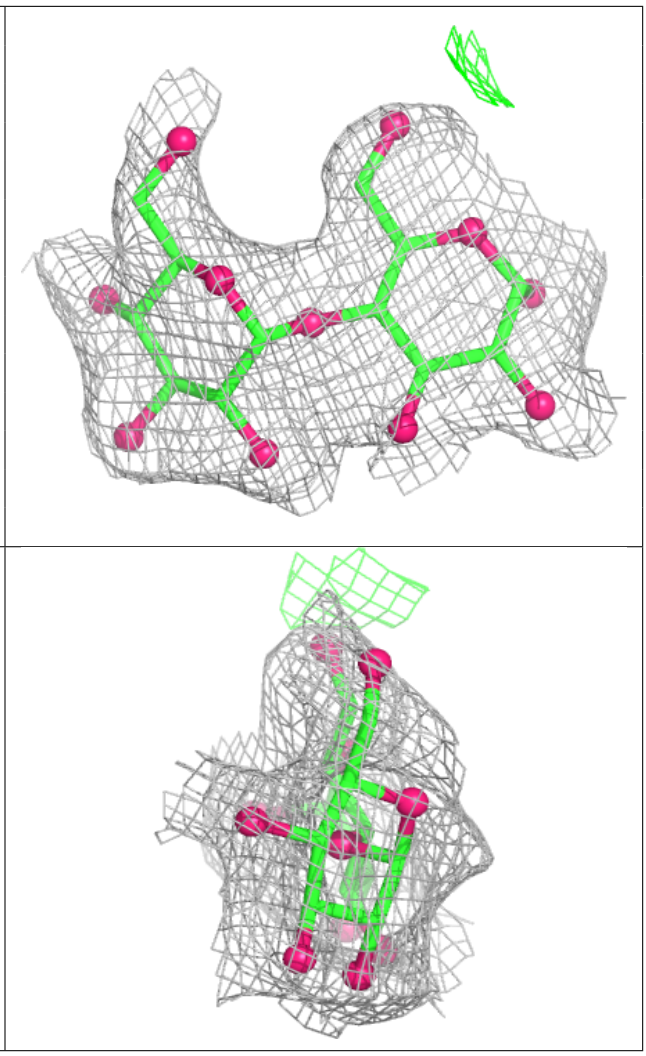
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	GLC	C	1	12/12	-	-	39,41,42,42	0
5	GLC	C	2	11/12	-	-	38,39,40,41	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

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



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
6	PGV	G	4005	10/51	0.56	0.17	58,65,70,72	0
6	PGV	G	4004	12/51	0.64	0.22	56,61,67,70	0
6	PGV	F	4008	8/51	0.64	0.19	53,58,61,64	0
6	PGV	G	4009	12/51	0.66	0.19	60,61,63,63	0
6	PGV	G	4007	9/51	0.71	0.17	55,62,72,72	0
6	PGV	G	4003	8/51	0.71	0.22	74,78,78,78	0
6	PGV	G	4006	7/51	0.75	0.17	61,63,68,70	0
6	PGV	F	4002	9/51	0.78	0.15	37,43,49,52	0
6	PGV	F	4001	51/51	0.86	0.14	64,70,77,77	0

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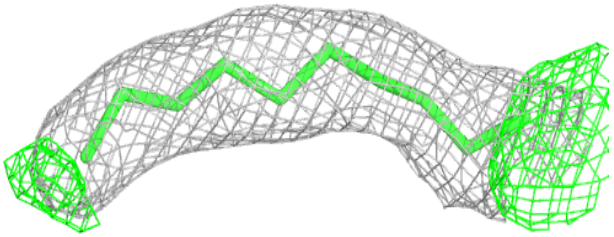
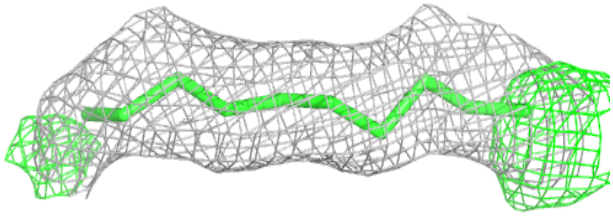
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
6	PGV	F	4010	13/51	0.91	0.14	74,79,85,85	0
7	UMQ	F	5004	34/34	0.92	0.10	63,76,82,84	0
9	ADP	B	2502	27/27	0.97	0.07	33,48,57,60	0
9	ADP	A	2501	27/27	0.98	0.06	21,32,40,43	0
8	MG	B	1502	1/1	0.99	0.03	28,28,28,28	0
10	VO4	A	3001	5/5	0.99	0.06	18,19,20,21	0
10	VO4	B	3002	5/5	0.99	0.04	29,29,31,31	0
8	MG	A	1501	1/1	1.00	0.02	17,17,17,17	0

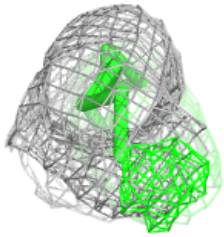
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PGV G 4005:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

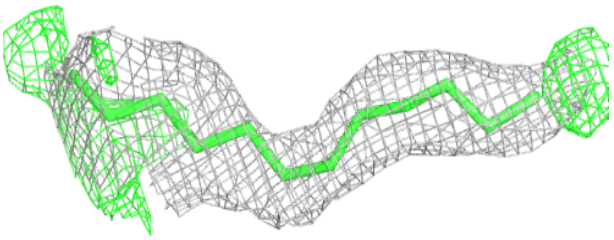


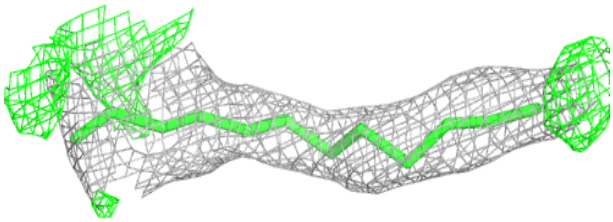


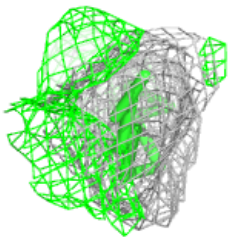


Electron density around PGV G 4004:

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

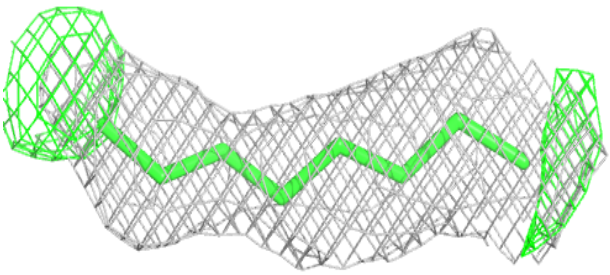


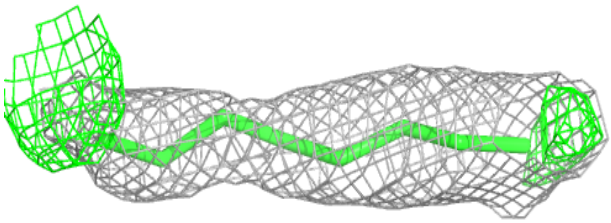


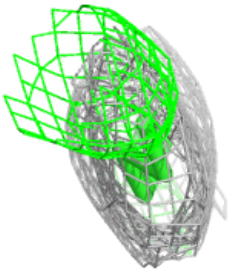


Electron density around PGV F 4008:

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

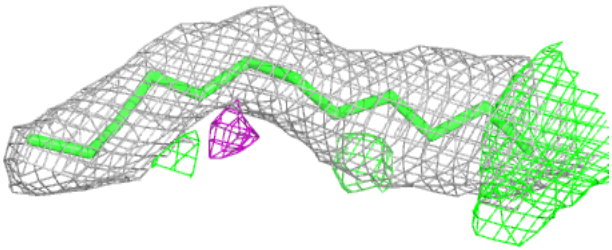


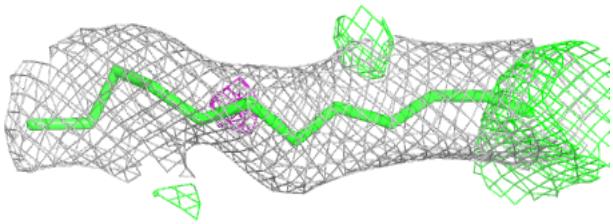


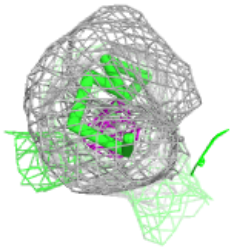


Electron density around PGV G 4009:

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

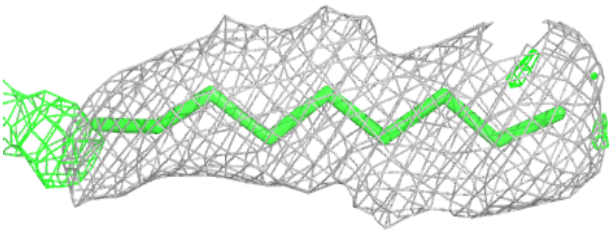


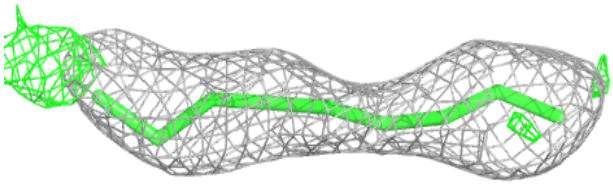


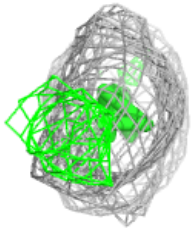


Electron density around PGV G 4007:

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

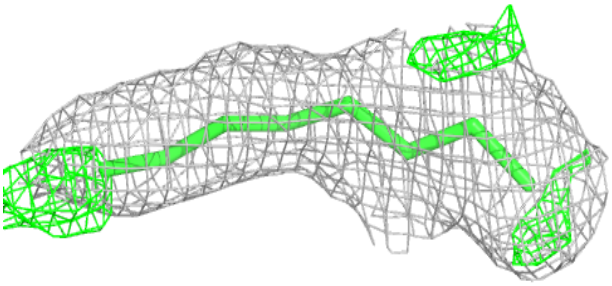


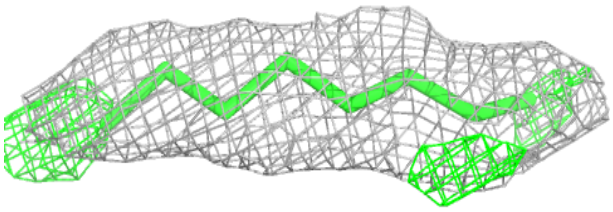


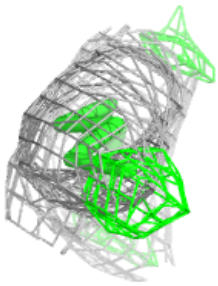


Electron density around PGV G 4003:

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

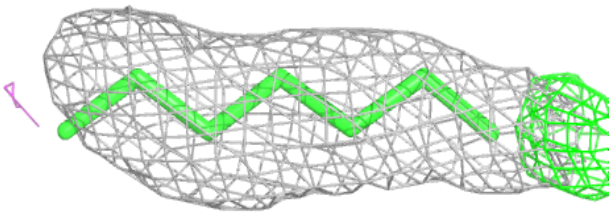


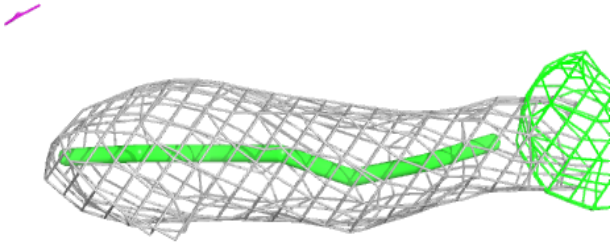


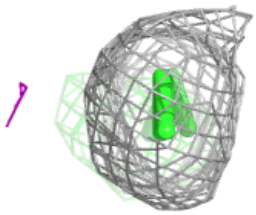


Electron density around PGV G 4006:

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

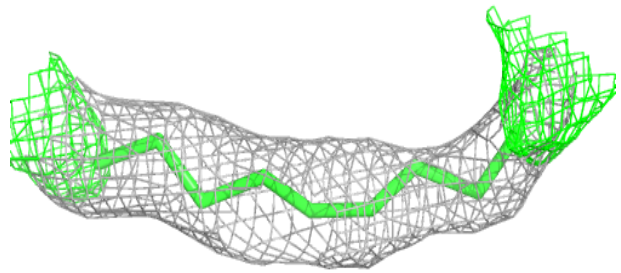


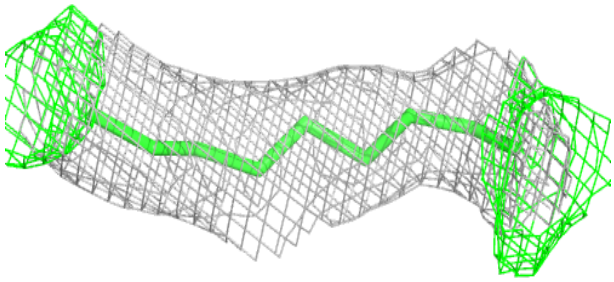


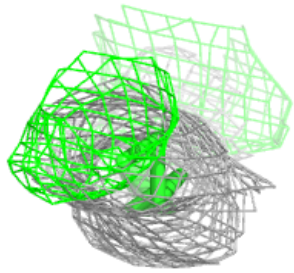


Electron density around PGV F 4002:

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

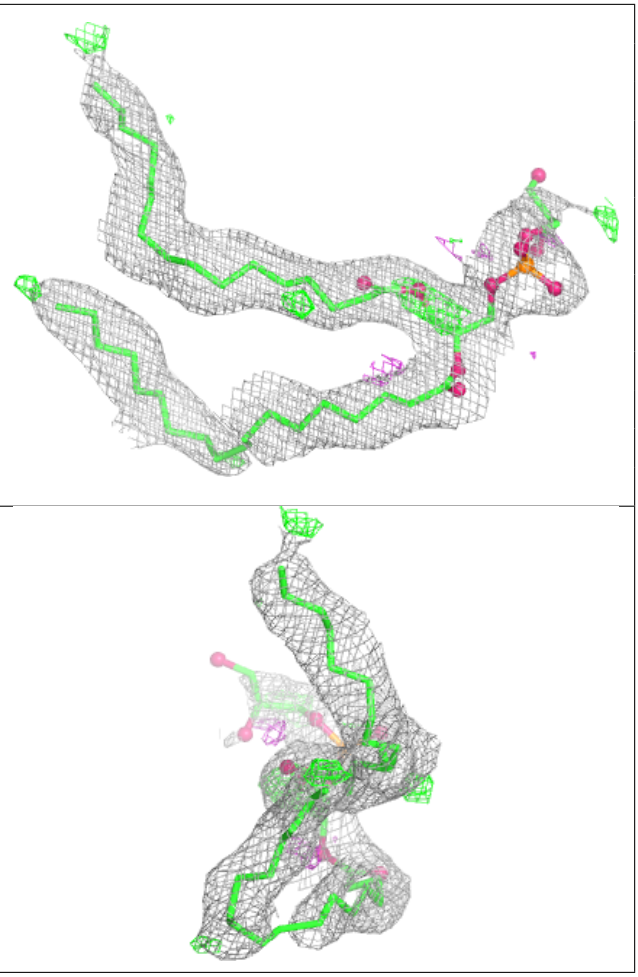






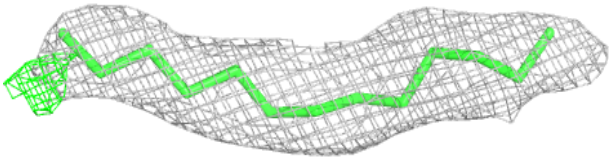
Electron density around PGV F 4001:

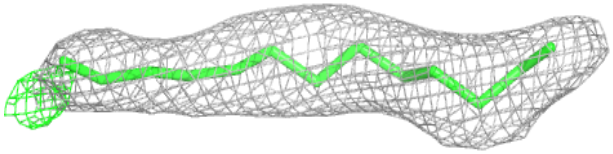
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

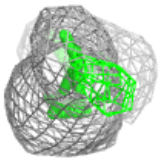


Electron density around PGV F 4010:

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

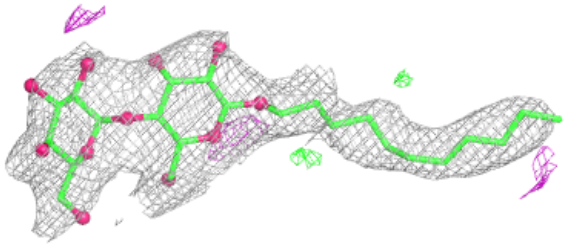


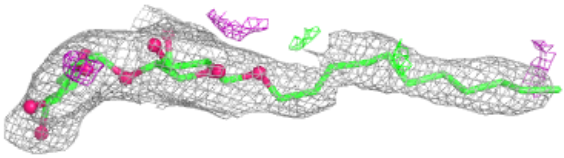


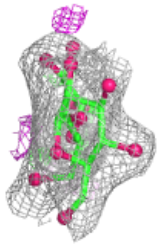


Electron density around UMQ F 5004:

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

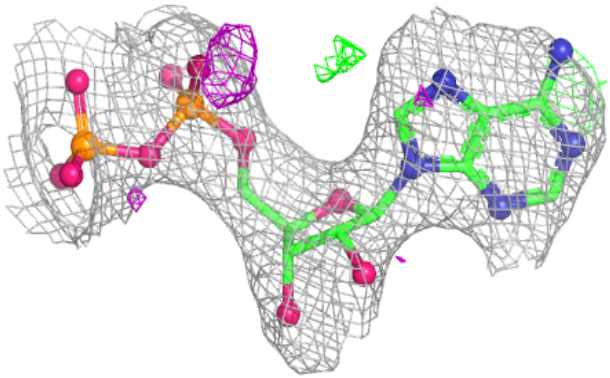


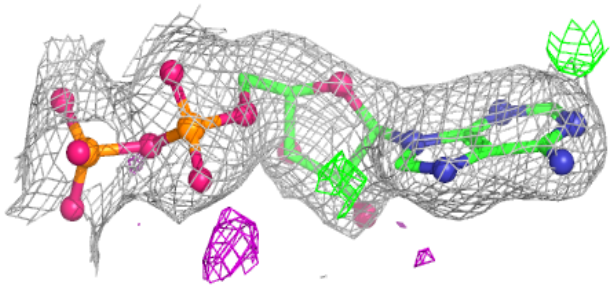


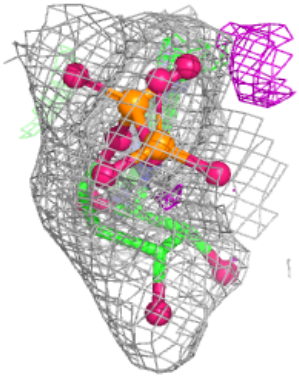


Electron density around ADP B 2502:

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

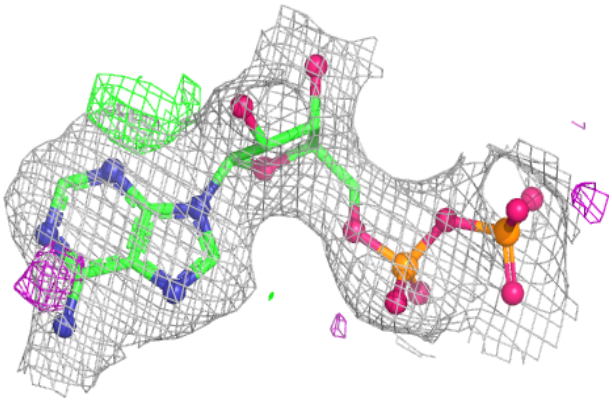


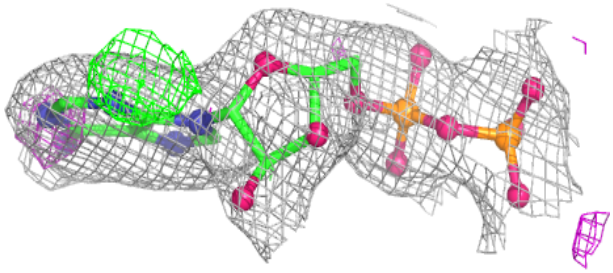


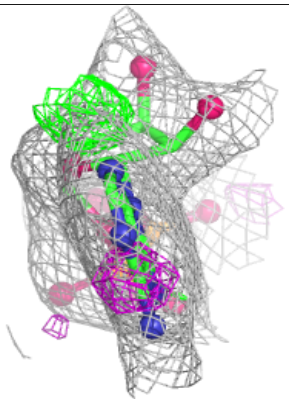


Electron density around ADP A 2501:

$2mF_o$-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.