Full wwPDB X-ray Structure Validation Report ⓘ

Jun 19, 2020 – 08:04 pm BST

PDB ID	:	5E9S
Title	:	Crystal structure of substrate-bound glutamate transporter homologue GltTk
Authors	:	Guskov, A.; Slotboom, D.J.
Deposited on	:	2015-10-15
Resolution	:	2.80 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:

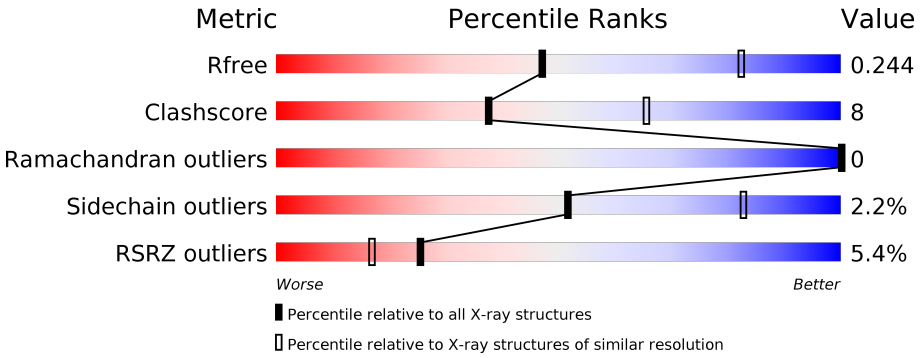
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 <=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	438	4% 76% 21%
1	B	438	6% 80% 17%
1	C	438	6% 78% 19%

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
4	DMU	A	505	-	-	-	X
5	1PE	B	506	-	-	-	X
5	1PE	B	507	-	-	-	X
6	PEG	B	511	-	-	-	X
6	PEG	B	516	-	-	-	X
6	PEG	B	519	-	-	-	X
6	PEG	B	522	-	-	-	X
6	PEG	C	509	-	-	-	X
6	PEG	C	513	-	-	-	X
6	PEG	C	514	-	-	-	X
8	P6G	C	520	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10213 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 Proton/glutamate symporter, SDF family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	0	1	0
			3196	2108	518	553	17			
1	B	427	Total	C	N	O	S	0	0	0
			3188	2103	517	552	16			
1	C	426	Total	C	N	O	S	0	1	0
			3186	2102	515	552	17			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	431	HIS	-	expression tag	UNP Q5JID0
A	432	HIS	-	expression tag	UNP Q5JID0
A	433	HIS	-	expression tag	UNP Q5JID0
A	434	HIS	-	expression tag	UNP Q5JID0
A	435	HIS	-	expression tag	UNP Q5JID0
A	436	HIS	-	expression tag	UNP Q5JID0
A	437	HIS	-	expression tag	UNP Q5JID0
A	438	HIS	-	expression tag	UNP Q5JID0
B	431	HIS	-	expression tag	UNP Q5JID0
B	432	HIS	-	expression tag	UNP Q5JID0
B	433	HIS	-	expression tag	UNP Q5JID0
B	434	HIS	-	expression tag	UNP Q5JID0
B	435	HIS	-	expression tag	UNP Q5JID0
B	436	HIS	-	expression tag	UNP Q5JID0
B	437	HIS	-	expression tag	UNP Q5JID0
B	438	HIS	-	expression tag	UNP Q5JID0
C	431	HIS	-	expression tag	UNP Q5JID0
C	432	HIS	-	expression tag	UNP Q5JID0
C	433	HIS	-	expression tag	UNP Q5JID0
C	434	HIS	-	expression tag	UNP Q5JID0
C	435	HIS	-	expression tag	UNP Q5JID0
C	436	HIS	-	expression tag	UNP Q5JID0
C	437	HIS	-	expression tag	UNP Q5JID0

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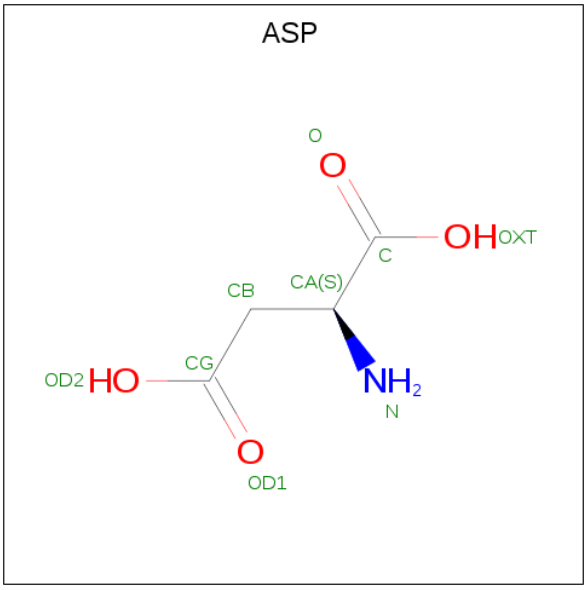
Chain	Residue	Modelled	Actual	Comment	Reference
C	438	HIS	-	expression tag	UNP Q5JID0

- Molecule 2 is ASPARTIC ACID (three-letter code: ASP) (formula: $C_4H_7NO_4$).

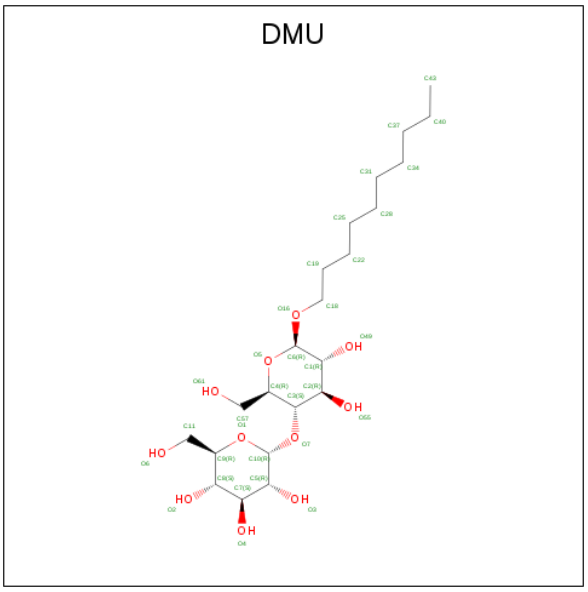


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N O 9 4 1 4	0	0
2	B	1	Total C N O 9 4 1 4	0	0
2	C	1	Total C N O 9 4 1 4	0	0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

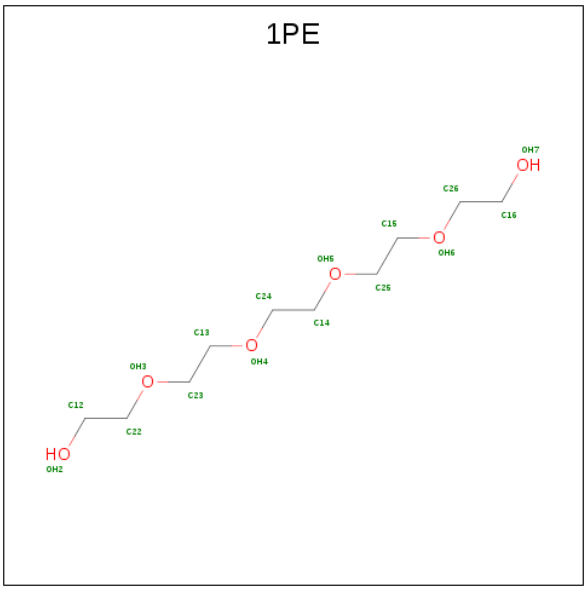
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	3	Total Na 3 3	0	0
3	A	3	Total Na 3 3	0	0
3	C	3	Total Na 3 3	0	0

- Molecule 4 is DECYL-BETA-D-MALTOPYRANOSIDE (three-letter code: DMU) (formula: $C_{22}H_{42}O_{11}$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 33 22 11	0	0
4	B	1	Total C O 33 22 11	0	0
4	C	1	Total C O 33 22 11	0	0

- Molecule 5 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 16 10 6	0	0

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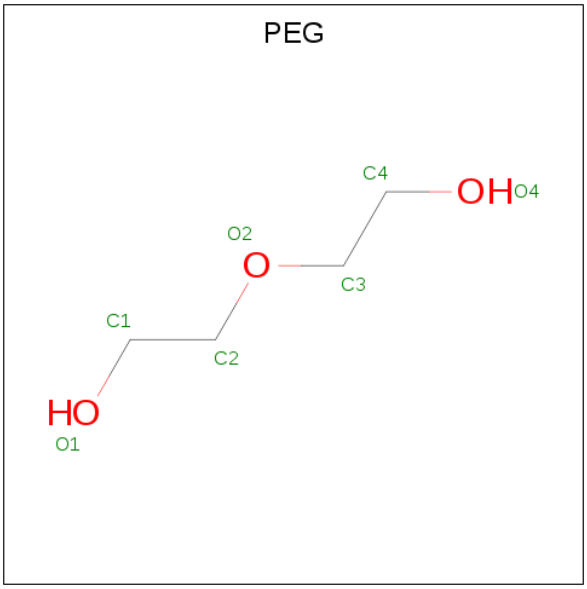
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 16 10 6	0	0
5	B	1	Total C O 16 10 6	0	0
5	B	1	Total C O 16 10 6	0	0
5	B	1	Total C O 16 10 6	0	0
5	C	1	Total C O 16 10 6	0	0
5	C	1	Total C O 16 10 6	0	0

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	A	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	B	1	Total C O 7 4 3	0	0
6	C	1	Total C O 7 4 3	0	0
6	C	1	Total C O 7 4 3	0	0

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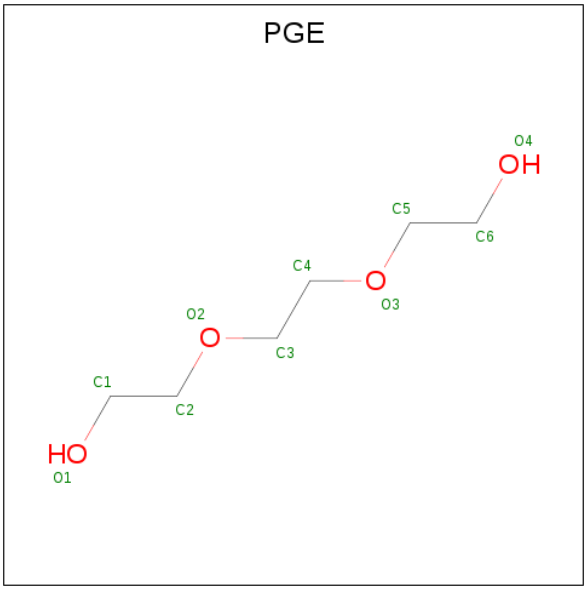
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	1	Total C O 7 4 3	0	0
6	C	1	Total C O 7 4 3	0	0
6	C	1	Total C O 7 4 3	0	0
6	C	1	Total C O 7 4 3	0	0
6	C	1	Total C O 7 4 3	0	0
6	C	1	Total C O 7 4 3	0	0
6	C	1	Total C O 7 4 3	0	0

- Molecule 7 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 10 6 4	0	0
7	A	1	Total C O 10 6 4	0	0
7	A	1	Total C O 10 6 4	0	0
7	A	1	Total C O 10 6 4	0	0

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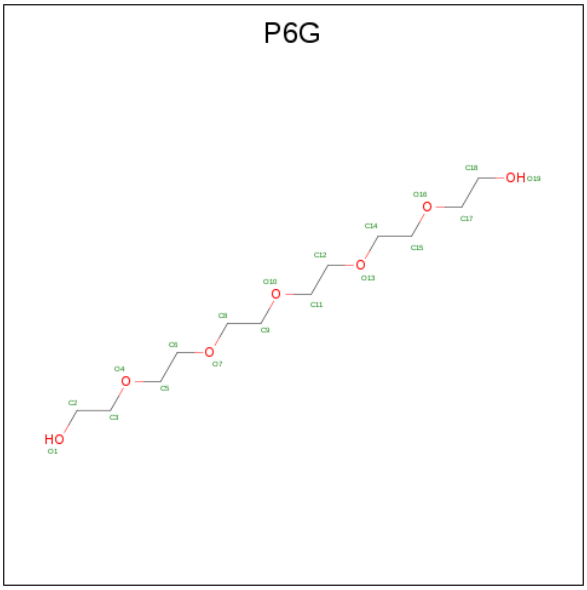
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total C O 10 6 4	0	0
7	C	1	Total C O 10 6 4	0	0
7	C	1	Total C O 10 6 4	0	0

- Molecule 8 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: $C_{12}H_{26}O_7$).



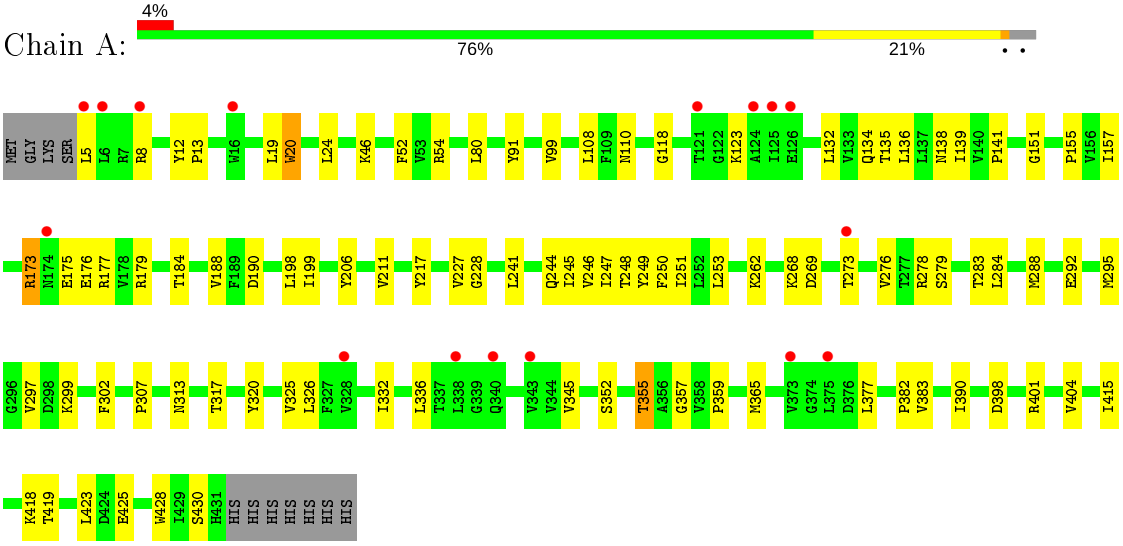
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 19 12 7	0	0
8	B	1	Total C O 19 12 7	0	0
8	B	1	Total C O 19 12 7	0	0
8	C	1	Total C O 19 12 7	0	0
8	C	1	Total C O 19 12 7	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

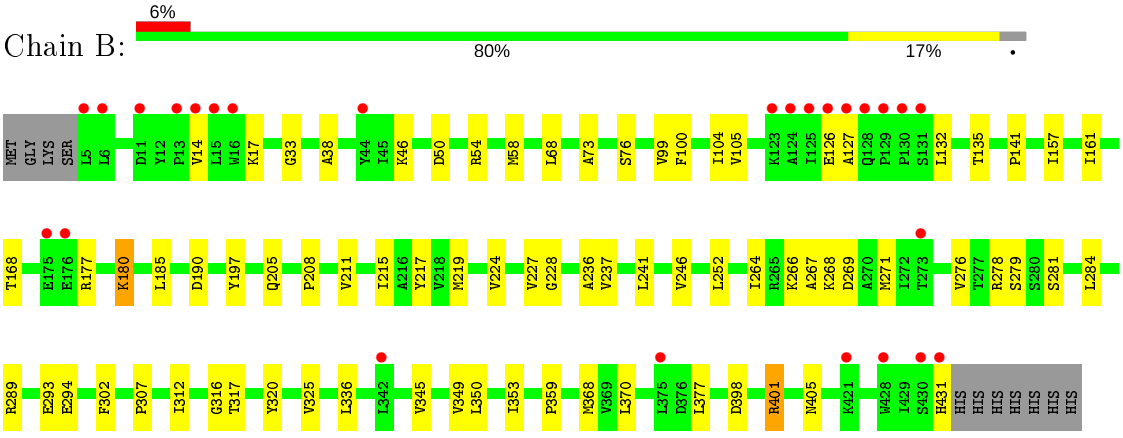
- Molecule 1: Proton/glutamate symporter, SDF family

Chain A:



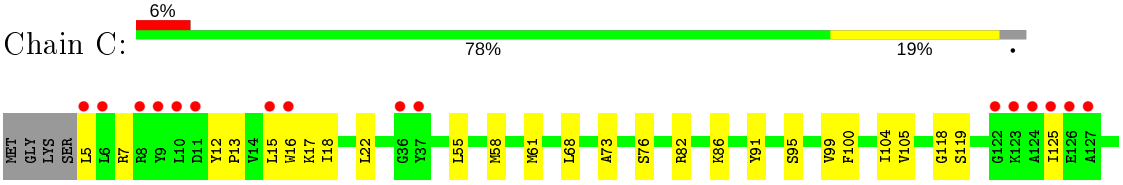
- Molecule 1: Proton/glutamate symporter, SDF family

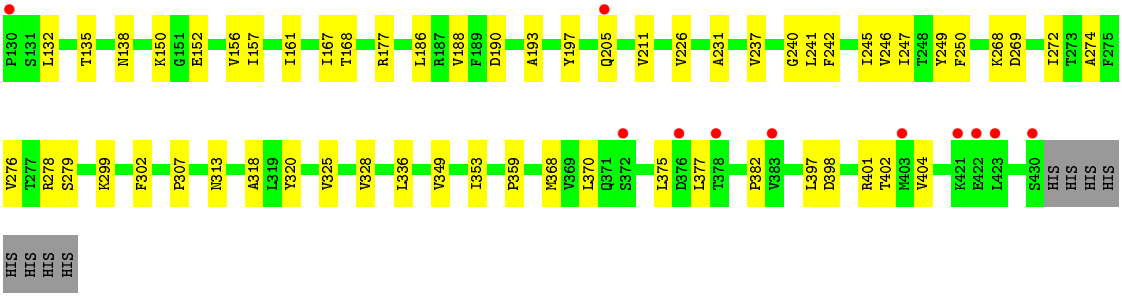
Chain B:



- Molecule 1: Proton/glutamate symporter, SDF family

Chain C:




P130 S131 L132 T135 N138 K150 G151 E152 V156 I157 I161 I167 T168 R177 L186 R187 V188 F189 D190 A193 Y197 Q205 V211 V226 A231 V237 G240 L241 F242 I245 V246 I247 T248 Y249 F250 K268 D269 I272 T273 A274 F275
V276 T277 R278 S279 K299 F302 P307 N313 A318 I319 Y320 V325 V328 L336 V349 I353 P359 M368 V369 L370 Q371 S372 L375 D376 L377 T378 P382 V383 L397 D398 R401 T402 N403 V404 K421 E422 L423 S430 HIS HIS HIS HIS
HIS HIS HIS HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	117.52Å 117.52Å 310.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.36 – 2.80 48.36 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.4 (48.36-2.80) 97.4 (48.36-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.04 (at 2.81Å)	Xtriage
Refinement program	PHENIX (1.10_2155)	Depositor
R, R_{free}	0.213 , 0.242 0.215 , 0.244	Depositor DCC
R_{free} test set	2009 reflections (3.32%)	wwPDB-VP
Wilson B-factor (Å^2)	83.7	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 65.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	10213	wwPDB-VP
Average B, all atoms (Å^2)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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: PGE, NA, 1PE, DMU, P6G, PEG

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	0.43	0/3256	0.64	0/4431
1	B	0.42	0/3248	0.61	0/4421
1	C	0.40	0/3245	0.63	0/4416
All	All	0.41	0/9749	0.63	0/13268

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	118	GLY	Peptide

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	3196	0	3402	64	0
1	B	3188	0	3394	45	0
1	C	3186	0	3395	53	0
2	A	9	0	3	2	0
2	B	9	0	3	3	0
2	C	9	0	3	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
4	A	33	0	42	0	0
4	B	33	0	42	1	0
4	C	33	0	42	2	0
5	A	32	0	44	2	0
5	B	48	0	66	4	0
5	C	32	0	44	0	0
6	A	70	0	100	1	0
6	B	98	0	140	3	0
6	C	63	0	90	5	0
7	A	40	0	56	2	0
7	B	10	0	14	0	0
7	C	20	0	28	0	0
8	A	19	0	26	0	0
8	B	38	0	52	2	0
8	C	38	0	52	1	0
All	All	10213	0	11038	163	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 8.

All (163) 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)
5:B:508:1PE:H241	6:B:522:PEG:H22	1.63	0.81
1:A:357:GLY:H	7:A:520:PGE:H12	1.47	0.79
1:A:110:ASN:HB2	5:A:507:1PE:H251	1.65	0.78
1:A:177:ARG:NH1	1:B:190:ASP:OD2	2.20	0.74
1:C:241:LEU:HB3	1:C:404:VAL:HG21	1.72	0.70
1:B:157:ILE:HD11	1:B:307:PRO:HB2	1.73	0.69
1:C:12:TYR:HD1	1:C:17:LYS:HD3	1.57	0.69
1:C:157:ILE:HD11	1:C:307:PRO:HB2	1.74	0.69
1:A:246:VAL:HG12	1:A:247:ILE:HD12	1.74	0.69
1:A:157:ILE:HD11	1:A:307:PRO:HB2	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:15:LEU:HB3	1:C:16:TRP:HD1	1.58	0.68
1:B:14:VAL:HG13	1:B:17:LYS:HE2	1.77	0.66
1:C:119:SER:HA	1:C:382:PRO:HG2	1.78	0.65
1:A:190:ASP:OD2	1:C:177:ARG:NH1	2.30	0.65
1:B:99:VAL:HG11	1:B:345:VAL:HA	1.79	0.64
1:A:355:THR:HG22	1:A:365:MET:HG3	1.79	0.64
1:A:262:LYS:HE2	1:A:428:TRP:O	2.00	0.62
1:A:5:LEU:HD23	1:A:8:ARG:HD3	1.80	0.61
1:C:246:VAL:HG23	1:C:247:ILE:HD12	1.82	0.61
5:B:507:1PE:H221	6:B:509:PEG:H32	1.82	0.60
1:C:274:ALA:HB1	1:C:402:THR:HG22	1.83	0.59
1:C:325:VAL:HG12	1:C:336:LEU:HD11	1.82	0.59
1:C:167:ILE:HG21	1:C:186:LEU:HD13	1.86	0.58
1:C:55:LEU:O	1:C:58:MET:HG2	2.03	0.58
1:A:279:SER:HB2	1:A:359:PRO:HD3	1.85	0.57
1:B:17:LYS:HE3	1:B:208:PRO:HG2	1.85	0.57
1:C:246:VAL:O	1:C:250:PHE:HB2	2.05	0.57
1:C:237:VAL:O	1:C:241:LEU:HG	2.05	0.56
1:B:177:ARG:NH1	1:C:190:ASP:OD2	2.38	0.56
1:A:418:LYS:NZ	1:A:425:GLU:OE2	2.40	0.55
1:B:325:VAL:HG22	1:B:370:LEU:HD12	1.89	0.55
1:A:284:LEU:HD22	1:A:288:MET:HE2	1.88	0.55
1:A:325:VAL:HG12	1:A:336:LEU:HD11	1.87	0.55
1:C:12:TYR:CD1	1:C:17:LYS:HD3	2.39	0.55
1:B:205:GLN:NE2	4:B:505:DMU:O49	2.36	0.55
1:B:278:ARG:HD2	1:B:398:ASP:HB3	1.89	0.55
1:A:20:TRP:HA	1:A:20:TRP:CE3	2.41	0.54
1:A:269:ASP:O	1:A:273:THR:HG23	2.08	0.54
1:A:262:LYS:NZ	1:A:430:SER:OG	2.38	0.54
1:C:197:TYR:HH	1:C:302:PHE:HE1	1.53	0.54
1:C:325:VAL:HG22	1:C:370:LEU:HD12	1.90	0.54
1:B:73:ALA:O	1:B:168:THR:HG21	2.09	0.53
1:A:20:TRP:HA	1:A:20:TRP:HE3	1.73	0.53
1:B:266:LYS:HB3	1:B:294:GLU:HB3	1.91	0.52
1:C:278:ARG:HD2	1:C:398:ASP:HB3	1.90	0.52
1:A:99:VAL:HG11	1:A:345:VAL:HA	1.92	0.51
1:A:401:ARG:NH1	2:A:501:ASP:OD1	2.43	0.51
1:C:99:VAL:HG22	1:C:318:ALA:HB1	1.93	0.51
1:A:12:TYR:CG	1:A:13:PRO:HD2	2.46	0.51
1:A:52:PHE:HB2	1:A:206:TYR:CE2	2.45	0.51
1:B:126:GLU:HB2	1:B:127:ALA:HA	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:ARG:HD2	1:A:398:ASP:HB3	1.93	0.50
1:C:226:VAL:HG13	1:C:231:ALA:HA	1.93	0.50
2:B:501:ASP:OD1	2:B:501:ASP:N	2.42	0.50
1:B:267:ALA:O	1:B:271:MET:HG3	2.12	0.50
1:A:184:THR:O	1:A:188:VAL:HG23	2.12	0.49
1:A:295:MET:HB2	1:A:297:VAL:HG23	1.94	0.49
1:A:54:ARG:NH2	1:C:138:ASN:HA	2.28	0.49
1:A:211:VAL:HG22	1:A:276:VAL:HG21	1.94	0.49
1:B:325:VAL:HG12	1:B:336:LEU:HD11	1.94	0.48
1:A:141:PRO:HD2	1:B:58:MET:HE3	1.94	0.48
1:A:132:LEU:O	1:A:135:THR:HB	2.14	0.48
1:A:245:ILE:HA	1:A:249:TYR:CD2	2.49	0.48
1:B:317:THR:OG1	2:B:501:ASP:OXT	2.30	0.48
1:A:108:LEU:HA	5:A:507:1PE:H242	1.95	0.48
1:B:279:SER:HB2	1:B:359:PRO:HD3	1.95	0.47
1:A:199:ILE:HG23	6:C:510:PEG:H11	1.96	0.47
1:A:423:LEU:HD11	1:A:428:TRP:HE1	1.79	0.47
1:A:246:VAL:O	1:A:250:PHE:HB2	2.15	0.47
6:C:508:PEG:H22	6:C:509:PEG:H21	1.95	0.47
1:C:73:ALA:O	1:C:168:THR:OG1	2.32	0.47
1:A:139:ILE:O	1:A:155:PRO:HA	2.15	0.47
1:A:317:THR:OG1	2:A:501:ASP:OD2	2.29	0.47
1:C:105:VAL:HG21	1:C:240:GLY:HA2	1.97	0.47
1:C:132:LEU:O	1:C:135:THR:N	2.47	0.47
1:C:242:PHE:O	1:C:246:VAL:HG22	2.14	0.47
1:B:197:TYR:HH	1:B:302:PHE:HE1	1.63	0.47
1:C:22:LEU:HD22	1:C:272:ILE:HG12	1.95	0.47
1:C:150:LYS:HD3	1:C:152:GLU:OE2	2.14	0.47
1:C:91:TYR:CG	1:C:313:ASN:HB2	2.50	0.46
1:A:151:GLY:HA3	7:A:521:PGE:H4	1.98	0.46
1:B:281:SER:O	1:B:284:LEU:HB2	2.15	0.46
1:A:118:GLY:O	1:A:382:PRO:HG2	2.15	0.46
1:A:227:VAL:HG13	1:A:228:GLY:N	2.31	0.45
1:B:132:LEU:O	1:B:135:THR:HB	2.16	0.45
1:B:353:ILE:HG21	5:B:506:1PE:H252	1.97	0.45
1:C:68:LEU:HD21	1:C:161:ILE:HG13	1.98	0.45
1:A:173:ARG:NH2	1:A:175:GLU:OE1	2.50	0.45
1:C:5:LEU:HG	1:C:7:ARG:H	1.81	0.45
1:A:176:GLU:HG3	1:A:179:ARG:NH2	2.32	0.45
1:A:268:LYS:HG3	1:A:269:ASP:N	2.32	0.44
1:B:405:ASN:ND2	2:B:501:ASP:HB2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:LYS:HD2	1:A:217:TYR:CG	2.53	0.44
1:C:349:VAL:O	1:C:353:ILE:HG13	2.17	0.44
8:C:520:P6G:H92	8:C:520:P6G:H141	2.00	0.44
1:A:377:LEU:HD23	1:A:377:LEU:HA	1.53	0.44
1:B:211:VAL:O	1:B:215:ILE:HG22	2.18	0.44
5:B:508:1PE:H131	6:B:522:PEG:H32	1.98	0.44
1:A:123:LYS:HE3	1:A:123:LYS:HB3	1.71	0.44
1:A:273:THR:OG1	1:A:283:THR:HG23	2.16	0.44
1:A:332:ILE:HD13	1:A:383:VAL:HG22	1.99	0.44
1:C:193:ALA:O	1:C:197:TYR:HD2	2.00	0.44
1:C:245:ILE:HA	1:C:249:TYR:CD2	2.53	0.44
1:A:198:LEU:HD22	6:C:510:PEG:H41	2.00	0.44
1:C:211:VAL:HG22	1:C:276:VAL:HG21	1.99	0.44
1:B:237:VAL:O	1:B:241:LEU:HG	2.18	0.43
1:C:186:LEU:HD23	6:C:512:PEG:H32	2.00	0.43
1:C:272:ILE:O	1:C:276:VAL:HG22	2.18	0.43
1:C:268:LYS:HG3	1:C:269:ASP:N	2.33	0.43
1:C:61:MET:CE	1:C:156:VAL:HG21	2.48	0.43
1:C:279:SER:HB2	1:C:359:PRO:HD3	2.00	0.43
1:C:99:VAL:CG2	1:C:318:ALA:HB1	2.49	0.43
1:B:276:VAL:O	1:B:278:ARG:NH1	2.52	0.43
1:B:33:GLY:HA2	1:B:38:ALA:HB2	1.99	0.43
1:B:312:ILE:HG21	8:B:524:P6G:H182	2.01	0.43
1:A:317:THR:HG21	1:A:355:THR:HG21	2.01	0.43
1:A:141:PRO:O	1:B:58:MET:HB3	2.18	0.43
1:A:80:LEU:HD12	1:A:80:LEU:HA	1.72	0.42
1:C:328:VAL:HG12	1:C:375:LEU:HD13	2.01	0.42
1:C:397:LEU:O	1:C:401:ARG:HG2	2.19	0.42
1:A:247:ILE:HG23	1:A:251:ILE:HD12	2.01	0.42
8:B:525:P6G:H21	8:B:525:P6G:H52	1.84	0.42
1:A:134:GLN:HG3	1:A:138:ASN:HD21	1.84	0.42
1:A:91:TYR:CG	1:A:313:ASN:HB2	2.55	0.42
1:B:289:ARG:NE	1:B:293:GLU:OE1	2.52	0.42
1:B:316:GLY:C	1:B:401:ARG:HG3	2.40	0.42
1:C:302:PHE:CD1	1:C:302:PHE:C	2.94	0.42
6:C:515:PEG:H32	6:C:515:PEG:H11	1.78	0.42
1:C:12:TYR:CG	1:C:13:PRO:HD2	2.55	0.41
4:C:505:DMU:H35	4:C:505:DMU:H30	2.01	0.41
1:A:352:SER:HA	1:A:355:THR:HG23	2.01	0.41
1:A:326:LEU:HD23	1:A:336:LEU:HD12	2.02	0.41
1:B:105:VAL:HG11	1:B:236:ALA:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:PHE:CE2	1:B:104:ILE:HD11	2.55	0.41
1:B:180:LYS:O	1:B:180:LYS:HD2	2.20	0.41
1:A:292:GLU:HB2	1:A:302:PHE:CZ	2.55	0.41
1:B:141:PRO:HG2	1:C:58:MET:HE3	2.01	0.41
1:A:244:GLN:O	1:A:248:THR:HB	2.20	0.41
1:B:350:LEU:HD23	1:B:350:LEU:HA	1.83	0.41
1:B:268:LYS:HG3	1:B:269:ASP:N	2.35	0.41
1:C:377:LEU:HA	1:C:377:LEU:HD23	1.67	0.41
6:A:511:PEG:H31	6:A:511:PEG:H12	1.88	0.41
1:B:349:VAL:O	1:B:353:ILE:HG13	2.20	0.41
1:C:95:SER:O	1:C:99:VAL:HG23	2.20	0.41
1:A:136:LEU:O	1:A:139:ILE:HB	2.21	0.41
1:B:252:LEU:HD23	1:B:252:LEU:HA	1.96	0.41
1:B:50:ASP:O	1:B:54:ARG:HG3	2.20	0.41
1:A:241:LEU:HB3	1:A:404:VAL:HG21	2.03	0.41
1:A:415:ILE:O	1:A:419:THR:HG23	2.21	0.41
1:B:227:VAL:HG13	1:B:228:GLY:N	2.36	0.41
1:B:46:LYS:HD2	1:B:217:TYR:CD2	2.55	0.41
1:A:20:TRP:O	1:A:24:LEU:HG	2.21	0.40
1:B:215:ILE:HD11	1:B:219:MET:HE2	2.03	0.40
1:C:205:GLN:HB3	4:C:505:DMU:H13	2.03	0.40
1:A:253:LEU:HA	1:A:253:LEU:HD23	1.92	0.40
1:B:68:LEU:HD21	1:B:161:ILE:HG13	2.02	0.40
1:B:141:PRO:O	1:C:58:MET:HB3	2.21	0.40
1:B:185:LEU:HD13	1:C:188:VAL:HG13	2.03	0.40
1:C:299:LYS:HG3	1:C:302:PHE:CZ	2.56	0.40
1:A:299:LYS:HB2	1:A:299:LYS:HE2	1.89	0.40
1:A:390:ILE:HD13	1:A:390:ILE:HA	1.83	0.40
1:C:100:PHE:CE2	1:C:104:ILE:HD11	2.56	0.40
1:C:82:ARG:HG2	1:C:86:LYS:HE3	2.03	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	426/438 (97%)	418 (98%)	8 (2%)	0	100	100
1	B	425/438 (97%)	415 (98%)	10 (2%)	0	100	100
1	C	425/438 (97%)	413 (97%)	12 (3%)	0	100	100
All	All	1276/1314 (97%)	1246 (98%)	30 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	336/345 (97%)	331 (98%)	5 (2%)	65	89
1	B	335/345 (97%)	325 (97%)	10 (3%)	41	75
1	C	335/345 (97%)	329 (98%)	6 (2%)	59	86
All	All	1006/1035 (97%)	985 (98%)	21 (2%)	52	84

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	20	TRP
1	A	173	ARG
1	A	320	TYR
1	A	355	THR
1	B	76	SER
1	B	180	LYS
1	B	224	VAL
1	B	246	VAL
1	B	264	ILE
1	B	320	TYR
1	B	368	MET
1	B	377	LEU

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Mol	Chain	Res	Type
1	B	401	ARG
1	B	431	HIS
1	C	18	ILE
1	C	76	SER
1	C	125	ILE
1	C	320	TYR
1	C	368[A]	MET
1	C	368[B]	MET

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

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 67 ligands modelled in this entry, 9 are monoatomic - leaving 58 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	PEG	B	519	-	6,6,6	0.49	0	5,5,5	0.30	0
6	PEG	C	510	-	6,6,6	0.48	0	5,5,5	0.39	0
6	PEG	B	520	-	6,6,6	0.53	0	5,5,5	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
7	PGE	A	520	-	9,9,9	0.45	0	8,8,8	0.35	0
6	PEG	B	515	-	6,6,6	0.49	0	5,5,5	0.42	0
4	DMU	A	505	-	34,34,34	1.68	9 (26%)	45,45,45	1.32	7 (15%)
6	PEG	C	515	-	6,6,6	0.50	0	5,5,5	0.27	0
7	PGE	A	518	-	9,9,9	0.35	0	8,8,8	0.34	0
6	PEG	B	510	-	6,6,6	0.52	0	5,5,5	0.28	0
6	PEG	C	516	-	6,6,6	0.45	0	5,5,5	0.30	0
7	PGE	B	523	-	9,9,9	0.43	0	8,8,8	0.27	0
8	P6G	B	524	-	18,18,18	0.54	0	17,17,17	0.47	0
6	PEG	A	513	-	6,6,6	0.50	0	5,5,5	0.36	0
5	1PE	A	506	-	15,15,15	0.54	0	14,14,14	0.36	0
6	PEG	C	509	-	6,6,6	0.53	0	5,5,5	0.35	0
5	1PE	C	506	-	15,15,15	0.52	0	14,14,14	0.58	0
6	PEG	A	510	-	6,6,6	0.49	0	5,5,5	0.32	0
4	DMU	B	505	-	34,34,34	1.61	8 (23%)	45,45,45	1.05	4 (8%)
8	P6G	C	519	-	18,18,18	0.60	0	17,17,17	0.55	0
6	PEG	A	509	-	6,6,6	0.49	0	5,5,5	0.29	0
8	P6G	B	525	-	18,18,18	0.56	0	17,17,17	0.49	0
6	PEG	A	512	-	6,6,6	0.46	0	5,5,5	0.38	0
6	PEG	A	516	-	6,6,6	0.53	0	5,5,5	0.30	0
6	PEG	C	508	-	6,6,6	0.50	0	5,5,5	0.43	0
6	PEG	A	517	-	6,6,6	0.56	0	5,5,5	0.80	0
8	P6G	C	520	-	18,18,18	0.54	0	17,17,17	0.50	0
6	PEG	A	515	-	6,6,6	0.51	0	5,5,5	0.26	0
6	PEG	A	514	-	6,6,6	0.51	0	5,5,5	0.25	0
6	PEG	A	508	-	6,6,6	0.47	0	5,5,5	0.31	0
6	PEG	C	514	-	6,6,6	0.51	0	5,5,5	0.38	0
6	PEG	B	521	-	6,6,6	0.49	0	5,5,5	0.30	0
6	PEG	A	511	-	6,6,6	0.51	0	5,5,5	0.48	0
7	PGE	C	518	-	9,9,9	0.31	0	8,8,8	0.58	0
6	PEG	B	509	-	6,6,6	0.52	0	5,5,5	0.31	0
6	PEG	B	522	-	6,6,6	0.51	0	5,5,5	0.28	0
6	PEG	B	514	-	6,6,6	0.49	0	5,5,5	0.31	0
6	PEG	B	513	-	6,6,6	0.51	0	5,5,5	0.38	0
7	PGE	C	517	-	9,9,9	0.34	0	8,8,8	0.26	0
6	PEG	B	517	-	6,6,6	0.50	0	5,5,5	0.24	0
5	1PE	B	506	-	15,15,15	0.56	0	14,14,14	0.49	0
5	1PE	B	508	-	15,15,15	0.54	0	14,14,14	0.24	0
5	1PE	B	507	-	15,15,15	0.55	0	14,14,14	0.33	0
6	PEG	B	512	-	6,6,6	0.54	0	5,5,5	0.32	0
6	PEG	B	518	-	6,6,6	0.49	0	5,5,5	0.38	0
6	PEG	B	511	-	6,6,6	0.48	0	5,5,5	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
6	PEG	C	511	-	6,6,6	0.52	0	5,5,5	0.38	0
4	DMU	C	505	-	34,34,34	1.60	9 (26%)	45,45,45	1.06	3 (6%)
5	1PE	A	507	-	15,15,15	0.54	0	14,14,14	0.42	0
8	P6G	A	522	-	18,18,18	0.59	0	17,17,17	0.48	0
7	PGE	A	519	-	9,9,9	0.38	0	8,8,8	0.32	0
7	PGE	A	521	-	9,9,9	0.36	0	8,8,8	0.40	0
5	1PE	C	507	-	15,15,15	0.54	0	14,14,14	0.22	0
6	PEG	C	512	-	6,6,6	0.51	0	5,5,5	0.35	0
6	PEG	B	516	-	6,6,6	0.53	0	5,5,5	0.28	0
6	PEG	C	513	-	6,6,6	0.50	0	5,5,5	0.36	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
6	PEG	B	519	-	-	2/4/4/4	-
6	PEG	C	510	-	-	3/4/4/4	-
6	PEG	B	520	-	-	2/4/4/4	-
7	PGE	A	520	-	-	3/7/7/7	-
6	PEG	B	515	-	-	2/4/4/4	-
4	DMU	A	505	-	-	7/19/59/59	0/2/2/2
6	PEG	C	515	-	-	4/4/4/4	-
7	PGE	A	518	-	-	6/7/7/7	-
6	PEG	B	510	-	-	2/4/4/4	-
6	PEG	C	516	-	-	1/4/4/4	-
7	PGE	B	523	-	-	4/7/7/7	-
8	P6G	B	524	-	-	7/16/16/16	-
6	PEG	A	513	-	-	1/4/4/4	-
5	1PE	A	506	-	-	7/13/13/13	-
6	PEG	C	509	-	-	1/4/4/4	-
5	1PE	C	506	-	-	5/13/13/13	-
6	PEG	A	510	-	-	2/4/4/4	-
4	DMU	B	505	-	-	10/19/59/59	0/2/2/2
8	P6G	C	519	-	-	9/16/16/16	-
6	PEG	A	509	-	-	2/4/4/4	-
8	P6G	B	525	-	-	10/16/16/16	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PEG	A	512	-	-	3/4/4/4	-
6	PEG	A	516	-	-	3/4/4/4	-
6	PEG	C	508	-	-	1/4/4/4	-
6	PEG	A	517	-	-	2/4/4/4	-
8	P6G	C	520	-	-	11/16/16/16	-
6	PEG	A	515	-	-	2/4/4/4	-
6	PEG	A	514	-	-	2/4/4/4	-
6	PEG	A	508	-	-	2/4/4/4	-
6	PEG	C	514	-	-	2/4/4/4	-
6	PEG	B	521	-	-	2/4/4/4	-
6	PEG	A	511	-	-	3/4/4/4	-
7	PGE	C	518	-	-	3/7/7/7	-
6	PEG	B	509	-	-	0/4/4/4	-
6	PEG	B	522	-	-	1/4/4/4	-
6	PEG	B	514	-	-	2/4/4/4	-
6	PEG	B	513	-	-	2/4/4/4	-
7	PGE	C	517	-	-	2/7/7/7	-
6	PEG	B	517	-	-	3/4/4/4	-
5	1PE	B	506	-	-	9/13/13/13	-
5	1PE	B	508	-	-	6/13/13/13	-
5	1PE	B	507	-	-	8/13/13/13	-
6	PEG	B	512	-	-	1/4/4/4	-
6	PEG	B	518	-	-	1/4/4/4	-
6	PEG	B	511	-	-	1/4/4/4	-
6	PEG	C	511	-	-	3/4/4/4	-
4	DMU	C	505	-	-	10/19/59/59	0/2/2/2
5	1PE	A	507	-	-	9/13/13/13	-
8	P6G	A	522	-	-	9/16/16/16	-
7	PGE	A	519	-	-	4/7/7/7	-
7	PGE	A	521	-	-	2/7/7/7	-
5	1PE	C	507	-	-	7/13/13/13	-
6	PEG	C	512	-	-	3/4/4/4	-
6	PEG	B	516	-	-	4/4/4/4	-
6	PEG	C	513	-	-	0/4/4/4	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	505	DMU	O1-C9	4.74	1.55	1.44
4	B	505	DMU	O1-C9	4.67	1.55	1.44
4	C	505	DMU	O1-C9	4.64	1.55	1.44
4	C	505	DMU	C11-C9	-3.33	1.40	1.51
4	B	505	DMU	C11-C9	-3.15	1.41	1.51
4	A	505	DMU	C11-C9	-3.14	1.41	1.51
4	A	505	DMU	O5-C6	3.02	1.49	1.41
4	A	505	DMU	O5-C4	2.82	1.51	1.44
4	B	505	DMU	O5-C6	2.65	1.48	1.41
4	B	505	DMU	O4-C7	2.65	1.49	1.43
4	B	505	DMU	C8-C9	2.59	1.58	1.53
4	A	505	DMU	O4-C7	2.58	1.49	1.43
4	C	505	DMU	O4-C7	2.48	1.48	1.43
4	B	505	DMU	O5-C4	2.38	1.50	1.44
4	C	505	DMU	C8-C9	2.37	1.58	1.53
4	A	505	DMU	O1-C10	2.36	1.47	1.41
4	C	505	DMU	O5-C6	2.33	1.47	1.41
4	A	505	DMU	C8-C9	2.32	1.57	1.53
4	C	505	DMU	C7-C5	-2.32	1.46	1.52
4	C	505	DMU	O5-C4	2.29	1.49	1.44
4	B	505	DMU	O3-C5	2.25	1.48	1.43
4	A	505	DMU	C7-C5	-2.12	1.46	1.52
4	A	505	DMU	O3-C5	2.12	1.48	1.43
4	B	505	DMU	C7-C5	-2.11	1.47	1.52
4	C	505	DMU	O1-C10	2.09	1.47	1.41
4	C	505	DMU	O3-C5	2.03	1.47	1.43

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	505	DMU	O5-C4-C3	3.79	117.73	109.75
4	A	505	DMU	C18-O16-C6	3.28	119.29	113.84
4	C	505	DMU	O5-C4-C3	3.14	116.37	109.75
4	B	505	DMU	O5-C4-C3	2.92	115.91	109.75
4	B	505	DMU	C7-C8-C9	2.84	115.30	110.24
4	A	505	DMU	O55-C2-C1	-2.82	103.82	110.35
4	A	505	DMU	C7-C8-C9	2.35	114.43	110.24
4	A	505	DMU	C6-O5-C4	2.30	118.21	113.69
4	A	505	DMU	C57-C4-C3	-2.28	106.69	113.33
4	A	505	DMU	C2-C3-C4	2.28	116.14	110.93
4	B	505	DMU	C18-O16-C6	2.13	117.38	113.84
4	B	505	DMU	C10-O7-C3	-2.13	112.69	117.96
4	C	505	DMU	C1-C2-C3	2.08	114.43	109.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	C	505	DMU	C10-O7-C3	-2.04	112.92	117.96

There are no chirality outliers.

All (213) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	505	DMU	O5-C6-O16-C18
4	B	505	DMU	O5-C6-O16-C18
4	C	505	DMU	O5-C6-O16-C18
4	A	505	DMU	C2-C3-O7-C10
4	B	505	DMU	O6-C11-C9-O1
8	A	522	P6G	C2-C3-O4-C5
6	B	512	PEG	C1-C2-O2-C3
6	C	515	PEG	O2-C3-C4-O4
7	A	518	PGE	O1-C1-C2-O2
6	B	513	PEG	O2-C3-C4-O4
7	A	519	PGE	O3-C5-C6-O4
6	B	513	PEG	C4-C3-O2-C2
8	B	525	P6G	C2-C3-O4-C5
4	B	505	DMU	O6-C11-C9-C8
4	B	505	DMU	O5-C4-C57-O61
7	A	518	PGE	O2-C3-C4-O3
8	B	524	P6G	O7-C8-C9-O10
8	B	525	P6G	O4-C5-C6-O7
5	A	507	1PE	OH4-C13-C23-OH3
5	B	507	1PE	OH4-C13-C23-OH3
4	C	505	DMU	O6-C11-C9-O1
8	C	519	P6G	O7-C8-C9-O10
5	B	506	1PE	OH4-C13-C23-OH3
4	C	505	DMU	O6-C11-C9-C8
8	C	520	P6G	O4-C5-C6-O7
4	B	505	DMU	C3-C4-C57-O61
7	A	519	PGE	O1-C1-C2-O2
7	A	521	PGE	O1-C1-C2-O2
4	A	505	DMU	C3-C4-C57-O61
8	A	522	P6G	O7-C8-C9-O10
5	B	508	1PE	OH4-C13-C23-OH3
5	B	508	1PE	OH6-C15-C25-OH5
8	C	520	P6G	O7-C8-C9-O10
8	B	524	P6G	O13-C14-C15-O16
4	C	505	DMU	O16-C18-C19-C22
6	C	510	PEG	O2-C3-C4-O4

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Mol	Chain	Res	Type	Atoms
5	A	506	1PE	OH7-C16-C26-OH6
5	C	506	1PE	OH2-C12-C22-OH3
6	C	511	PEG	O1-C1-C2-O2
6	A	512	PEG	O1-C1-C2-O2
6	A	508	PEG	O1-C1-C2-O2
6	B	510	PEG	O2-C3-C4-O4
5	B	508	1PE	OH7-C16-C26-OH6
5	B	507	1PE	OH7-C16-C26-OH6
7	B	523	PGE	O2-C3-C4-O3
5	A	506	1PE	OH4-C13-C23-OH3
8	B	525	P6G	O10-C11-C12-O13
6	A	517	PEG	C1-C2-O2-C3
6	A	509	PEG	O2-C3-C4-O4
8	C	520	P6G	O16-C17-C18-O19
6	B	521	PEG	O1-C1-C2-O2
6	C	512	PEG	O1-C1-C2-O2
8	C	519	P6G	O10-C11-C12-O13
4	B	505	DMU	C1-C6-O16-C18
4	C	505	DMU	C1-C6-O16-C18
6	C	515	PEG	C1-C2-O2-C3
7	A	520	PGE	O3-C5-C6-O4
6	A	516	PEG	O2-C3-C4-O4
8	B	525	P6G	O1-C2-C3-O4
6	A	514	PEG	O2-C3-C4-O4
7	C	518	PGE	O1-C1-C2-O2
6	B	514	PEG	O1-C1-C2-O2
5	C	507	1PE	OH7-C16-C26-OH6
4	B	505	DMU	C31-C34-C37-C40
4	C	505	DMU	C18-C19-C22-C25
4	A	505	DMU	C31-C34-C37-C40
8	B	525	P6G	O13-C14-C15-O16
5	A	506	1PE	OH2-C12-C22-OH3
6	A	510	PEG	O1-C1-C2-O2
8	B	524	P6G	O1-C2-C3-O4
7	C	518	PGE	O2-C3-C4-O3
5	A	507	1PE	OH5-C14-C24-OH4
4	A	505	DMU	O6-C11-C9-O1
6	A	517	PEG	C4-C3-O2-C2
5	A	506	1PE	OH6-C15-C25-OH5
6	A	516	PEG	O1-C1-C2-O2
8	B	524	P6G	O16-C17-C18-O19
6	B	516	PEG	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
8	A	522	P6G	O1-C2-C3-O4
8	C	519	P6G	O13-C14-C15-O16
6	B	515	PEG	O2-C3-C4-O4
7	B	523	PGE	O3-C5-C6-O4
6	B	522	PEG	O1-C1-C2-O2
6	B	516	PEG	O2-C3-C4-O4
7	C	517	PGE	O3-C5-C6-O4
6	B	510	PEG	O1-C1-C2-O2
6	C	512	PEG	O2-C3-C4-O4
5	C	506	1PE	OH5-C14-C24-OH4
4	B	505	DMU	O16-C18-C19-C22
7	A	519	PGE	C4-C3-O2-C2
5	B	506	1PE	OH5-C14-C24-OH4
4	C	505	DMU	C22-C25-C28-C31
8	C	520	P6G	C12-C11-O10-C9
8	A	522	P6G	O10-C11-C12-O13
8	B	525	P6G	O16-C17-C18-O19
6	B	517	PEG	O2-C3-C4-O4
5	C	507	1PE	OH2-C12-C22-OH3
5	A	506	1PE	OH5-C14-C24-OH4
4	B	505	DMU	C34-C37-C40-C43
8	C	519	P6G	C2-C3-O4-C5
6	B	514	PEG	C4-C3-O2-C2
5	B	507	1PE	OH6-C15-C25-OH5
7	A	521	PGE	O2-C3-C4-O3
5	A	507	1PE	OH2-C12-C22-OH3
4	B	505	DMU	C28-C31-C34-C37
6	A	511	PEG	C1-C2-O2-C3
6	B	521	PEG	C1-C2-O2-C3
8	C	519	P6G	O4-C5-C6-O7
5	B	508	1PE	C15-C25-OH5-C14
6	C	512	PEG	C1-C2-O2-C3
6	A	512	PEG	O2-C3-C4-O4
6	C	514	PEG	O2-C3-C4-O4
6	B	518	PEG	O2-C3-C4-O4
8	A	522	P6G	C15-C14-O13-C12
7	A	518	PGE	C1-C2-O2-C3
8	C	520	P6G	C11-C12-O13-C14
5	A	507	1PE	C24-C14-OH5-C25
7	B	523	PGE	C6-C5-O3-C4
5	A	507	1PE	C15-C25-OH5-C14
7	C	517	PGE	C3-C4-O3-C5

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Mol	Chain	Res	Type	Atoms
5	C	506	1PE	C12-C22-OH3-C23
6	A	514	PEG	C4-C3-O2-C2
6	C	515	PEG	C4-C3-O2-C2
5	A	506	1PE	C13-C23-OH3-C22
6	A	508	PEG	C1-C2-O2-C3
5	A	507	1PE	C12-C22-OH3-C23
5	B	506	1PE	C25-C15-OH6-C26
8	C	520	P6G	C5-C6-O7-C8
8	B	525	P6G	C8-C9-O10-C11
5	B	506	1PE	C24-C14-OH5-C25
4	A	505	DMU	O5-C4-C57-O61
5	B	507	1PE	C14-C24-OH4-C13
6	C	511	PEG	C4-C3-O2-C2
6	A	512	PEG	C4-C3-O2-C2
5	A	507	1PE	C25-C15-OH6-C26
6	C	511	PEG	C1-C2-O2-C3
5	B	507	1PE	C23-C13-OH4-C24
8	B	524	P6G	C9-C8-O7-C6
5	C	507	1PE	C14-C24-OH4-C13
6	C	516	PEG	O1-C1-C2-O2
7	B	523	PGE	C4-C3-O2-C2
8	B	525	P6G	C18-C17-O16-C15
8	A	522	P6G	C11-C12-O13-C14
6	A	510	PEG	C1-C2-O2-C3
6	A	516	PEG	C1-C2-O2-C3
6	C	515	PEG	O1-C1-C2-O2
8	C	520	P6G	O13-C14-C15-O16
8	C	520	P6G	C18-C17-O16-C15
8	C	520	P6G	C14-C15-O16-C17
5	A	506	1PE	C12-C22-OH3-C23
4	C	505	DMU	C25-C28-C31-C34
6	B	517	PEG	C1-C2-O2-C3
7	C	518	PGE	C1-C2-O2-C3
5	C	506	1PE	C14-C24-OH4-C13
5	B	506	1PE	C23-C13-OH4-C24
6	C	510	PEG	C4-C3-O2-C2
7	A	518	PGE	O3-C5-C6-O4
6	B	515	PEG	C4-C3-O2-C2
5	B	506	1PE	C12-C22-OH3-C23
8	C	520	P6G	C2-C3-O4-C5
8	C	520	P6G	C6-C5-O4-C3
7	A	518	PGE	C3-C4-O3-C5

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Mol	Chain	Res	Type	Atoms
4	C	505	DMU	O1-C10-O7-C3
5	B	506	1PE	C13-C23-OH3-C22
6	A	511	PEG	C4-C3-O2-C2
6	B	519	PEG	C4-C3-O2-C2
6	B	516	PEG	C1-C2-O2-C3
7	A	520	PGE	O2-C3-C4-O3
4	A	505	DMU	C18-C19-C22-C25
7	A	520	PGE	C1-C2-O2-C3
6	A	513	PEG	C1-C2-O2-C3
6	C	510	PEG	C1-C2-O2-C3
5	C	507	1PE	OH4-C13-C23-OH3
5	B	506	1PE	C14-C24-OH4-C13
8	C	519	P6G	C8-C9-O10-C11
6	B	519	PEG	O1-C1-C2-O2
6	C	508	PEG	C4-C3-O2-C2
6	A	509	PEG	C4-C3-O2-C2
5	C	507	1PE	C15-C25-OH5-C14
5	B	508	1PE	C25-C15-OH6-C26
5	B	506	1PE	OH7-C16-C26-OH6
8	A	522	P6G	O16-C17-C18-O19
5	B	507	1PE	C16-C26-OH6-C15
8	C	519	P6G	C9-C8-O7-C6
8	B	524	P6G	C11-C12-O13-C14
7	A	518	PGE	C6-C5-O3-C4
5	B	508	1PE	C16-C26-OH6-C15
6	B	511	PEG	C4-C3-O2-C2
5	A	507	1PE	C14-C24-OH4-C13
6	B	520	PEG	C1-C2-O2-C3
6	B	517	PEG	O1-C1-C2-O2
5	A	507	1PE	OH6-C15-C25-OH5
5	B	507	1PE	C24-C14-OH5-C25
6	C	509	PEG	O1-C1-C2-O2
8	B	525	P6G	C5-C6-O7-C8
4	C	505	DMU	C5-C10-O7-C3
8	A	522	P6G	O13-C14-C15-O16
5	C	507	1PE	C13-C23-OH3-C22
6	A	515	PEG	O1-C1-C2-O2
8	B	525	P6G	O7-C8-C9-O10
6	A	511	PEG	O1-C1-C2-O2
8	C	519	P6G	O1-C2-C3-O4
6	B	516	PEG	C4-C3-O2-C2
7	A	519	PGE	O2-C3-C4-O3

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Mol	Chain	Res	Type	Atoms
6	B	520	PEG	O1-C1-C2-O2
5	C	507	1PE	C25-C15-OH6-C26
8	B	524	P6G	O4-C5-C6-O7
5	B	507	1PE	OH5-C14-C24-OH4
5	C	506	1PE	OH6-C15-C25-OH5
6	C	514	PEG	C1-C2-O2-C3
8	A	522	P6G	C5-C6-O7-C8
8	C	519	P6G	O16-C17-C18-O19
6	A	515	PEG	O2-C3-C4-O4

There are no ring outliers.

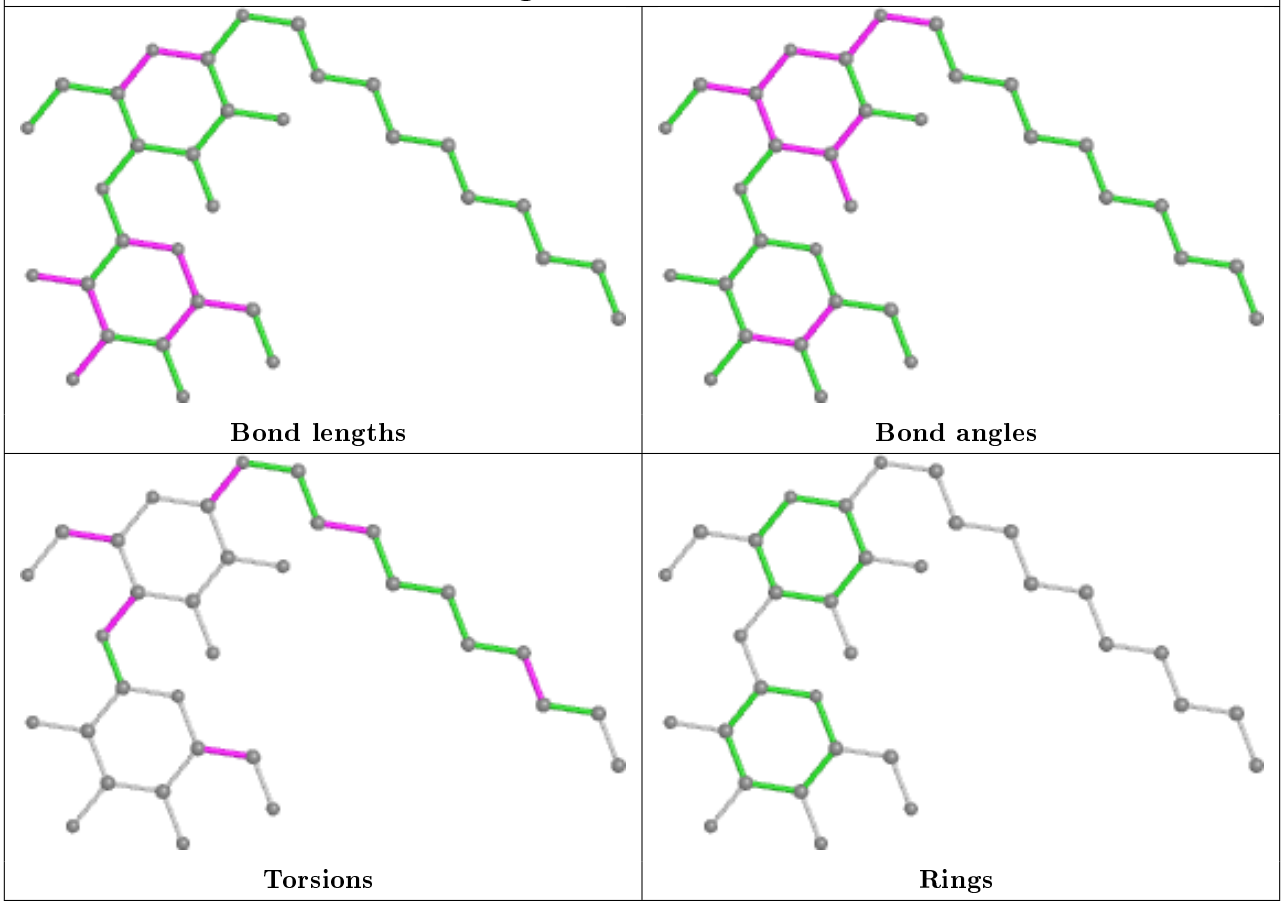
19 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	510	PEG	2	0
7	A	520	PGE	1	0
6	C	515	PEG	1	0
8	B	524	P6G	1	0
6	C	509	PEG	1	0
4	B	505	DMU	1	0
8	B	525	P6G	1	0
6	C	508	PEG	1	0
8	C	520	P6G	1	0
6	A	511	PEG	1	0
6	B	509	PEG	1	0
6	B	522	PEG	2	0
5	B	506	1PE	1	0
5	B	508	1PE	2	0
5	B	507	1PE	1	0
4	C	505	DMU	2	0
5	A	507	1PE	2	0
7	A	521	PGE	1	0
6	C	512	PEG	1	0

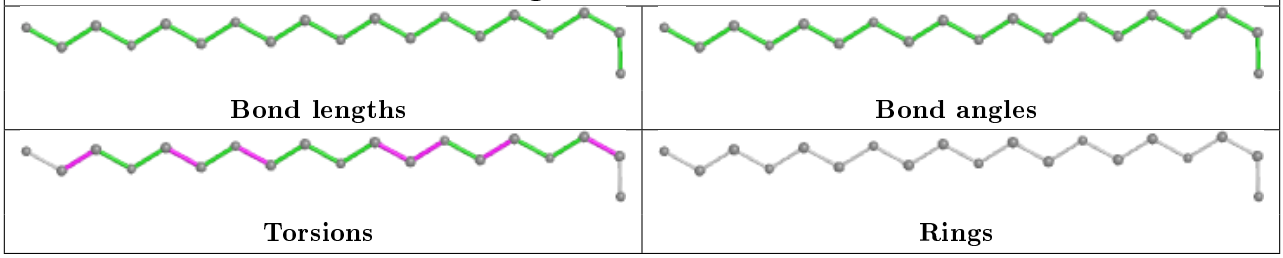
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 DMU A 505



Ligand P6G B 524



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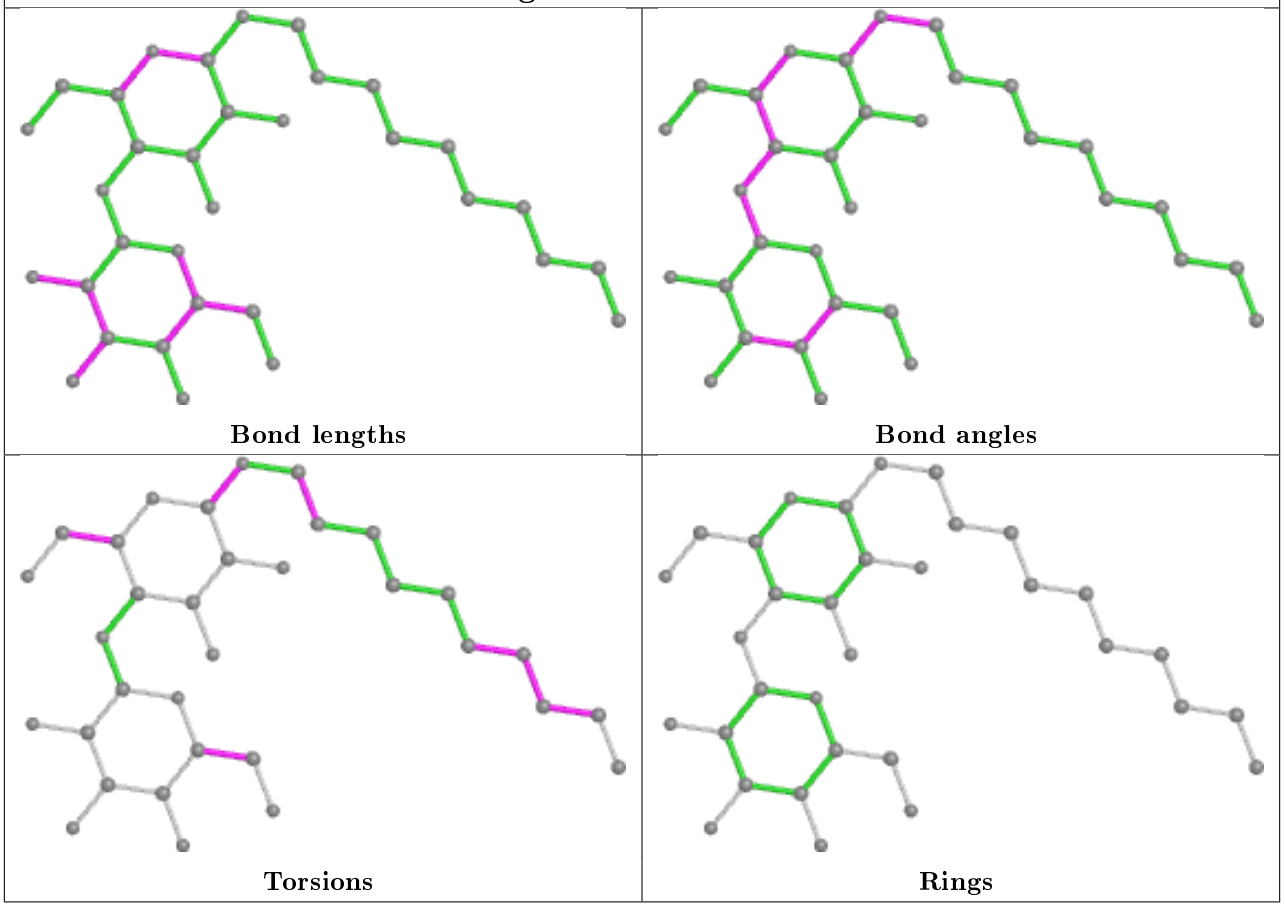

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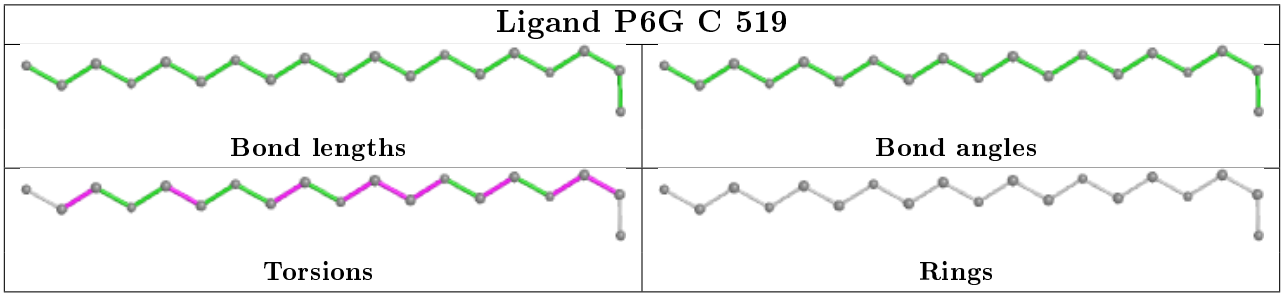

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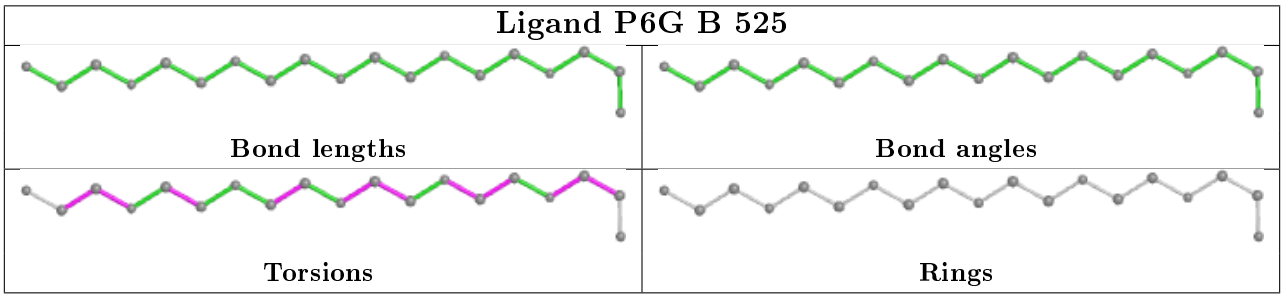
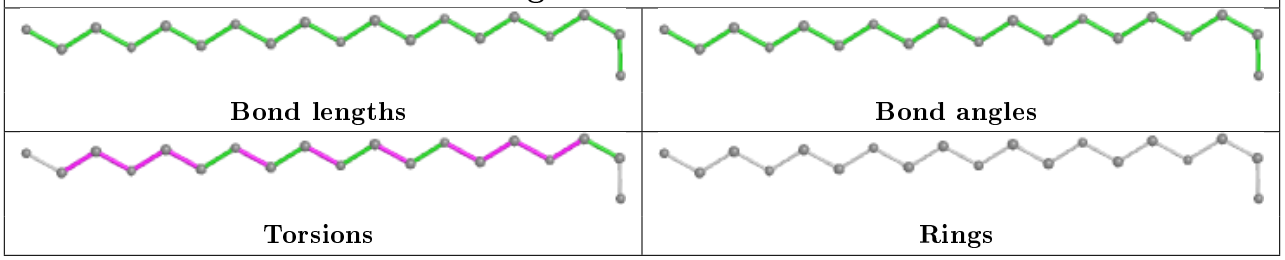
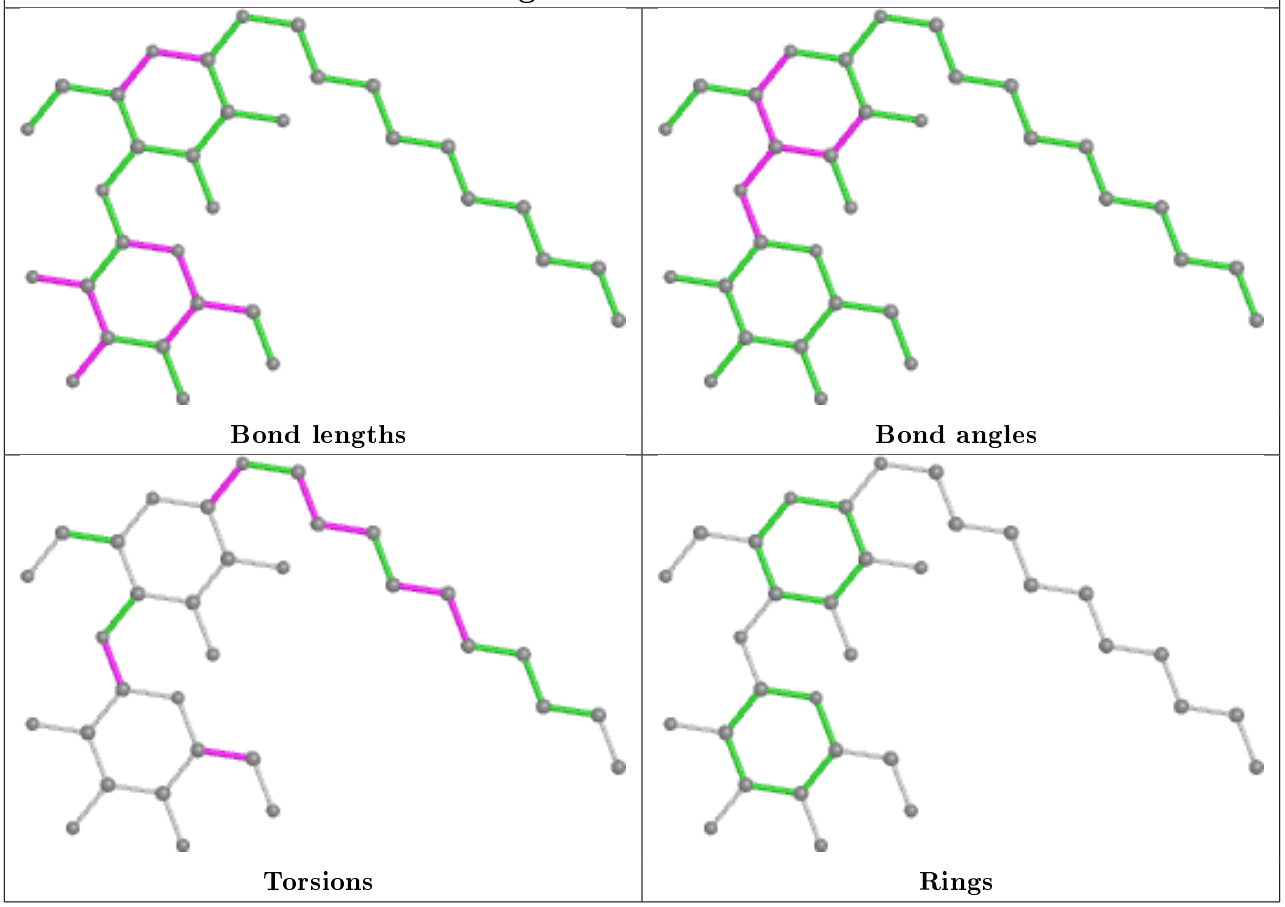


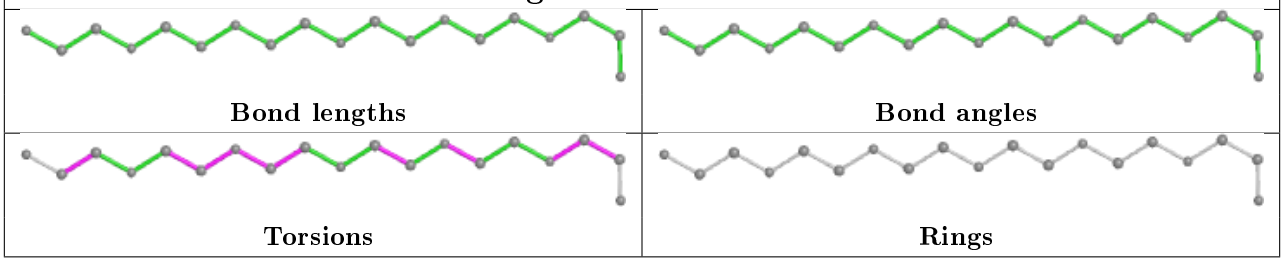
Ligand P6G C 520



Ligand DMU C 505



Ligand P6G A 522



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues

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	427/438 (97%)	-0.18	16 (3%)	41	31	55, 84, 153, 243	0
1	B	427/438 (97%)	0.06	26 (6%)	21	13	58, 90, 159, 234	0
1	C	426/438 (97%)	-0.08	27 (6%)	20	12	56, 94, 146, 212	0
All	All	1280/1314 (97%)	-0.07	69 (5%)	25	17	55, 89, 154, 243	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	430	SER	8.6
1	B	127	ALA	8.4
1	B	126	GLU	7.6
1	B	124	ALA	6.8
1	B	11	ASP	5.9
1	B	125	ILE	5.9
1	C	123	LYS	5.7
1	B	16	TRP	5.5
1	C	125	ILE	5.1
1	B	14	VAL	4.9
1	B	123	LYS	4.9
1	C	124	ALA	4.8
1	A	125	ILE	4.7
1	C	16	TRP	4.6
1	B	5	LEU	4.2
1	B	130	PRO	4.1
1	B	15	LEU	3.9
1	A	124	ALA	3.9
1	C	8	ARG	3.9
1	A	375	LEU	3.8
1	B	13	PRO	3.8
1	C	122	GLY	3.7
1	B	128	GLN	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	422	GLU	3.7
1	B	176	GLU	3.6
1	C	9	TYR	3.6
1	B	6	LEU	3.5
1	B	129	PRO	3.5
1	A	5	LEU	3.5
1	C	6	LEU	3.4
1	B	431	HIS	3.1
1	C	423	LEU	3.1
1	B	428	TRP	3.1
1	C	378	THR	3.0
1	C	376	ASP	3.0
1	A	373	VAL	3.0
1	A	121	THR	2.9
1	B	421	LYS	2.9
1	A	126	GLU	2.9
1	C	372	SER	2.8
1	C	130	PRO	2.8
1	C	126	GLU	2.8
1	A	8	ARG	2.8
1	B	44	TYR	2.8
1	C	37	TYR	2.8
1	A	6	LEU	2.8
1	C	127	ALA	2.7
1	A	16	TRP	2.7
1	C	15	LEU	2.7
1	A	340	GLN	2.6
1	C	421	LYS	2.6
1	B	430	SER	2.6
1	A	338	LEU	2.6
1	C	10	LEU	2.6
1	A	328	VAL	2.5
1	C	383	VAL	2.5
1	B	342	LEU	2.5
1	B	375	LEU	2.3
1	A	343	VAL	2.2
1	B	131	SER	2.2
1	A	174	ASN	2.2
1	C	5	LEU	2.1
1	B	273	THR	2.1
1	C	11	ASP	2.1
1	B	175	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	403	MET	2.0
1	C	36	GLY	2.0
1	A	273	THR	2.0
1	C	205	GLN	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 carbohydrates 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
6	PEG	B	520	7/7	0.54	0.30	120,125,133,138	0
6	PEG	C	514	7/7	0.62	0.84	122,124,129,130	0
5	1PE	B	507	16/16	0.63	0.60	80,121,137,138	0
6	PEG	A	514	7/7	0.64	0.34	119,128,132,132	0
5	1PE	B	508	16/16	0.66	0.36	121,139,151,151	0
8	P6G	C	519	19/19	0.67	0.30	83,109,120,122	0
6	PEG	C	513	7/7	0.67	0.46	113,118,124,125	0
6	PEG	A	511	7/7	0.68	0.21	111,116,121,121	0
6	PEG	B	515	7/7	0.68	0.14	109,110,118,120	0
6	PEG	B	521	7/7	0.68	0.24	124,124,128,129	0
6	PEG	B	519	7/7	0.69	0.58	132,133,142,142	0
5	1PE	C	507	16/16	0.69	0.15	127,143,149,151	0
8	P6G	B	525	19/19	0.69	0.25	95,125,130,131	0
7	PGE	B	523	10/10	0.70	0.23	108,130,138,139	0
5	1PE	B	506	16/16	0.72	0.47	94,113,121,125	0
6	PEG	B	517	7/7	0.73	0.22	110,113,118,119	0
4	DMU	A	505	33/33	0.73	0.58	108,156,169,172	0
6	PEG	B	522	7/7	0.73	0.56	120,125,133,134	0
7	PGE	C	518	10/10	0.74	0.36	124,133,135,137	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
8	P6G	A	522	19/19	0.74	0.26	99,124,135,137	0
8	P6G	C	520	19/19	0.76	0.41	111,120,138,139	0
6	PEG	B	511	7/7	0.76	0.55	109,122,129,132	0
5	1PE	A	507	16/16	0.76	0.30	111,124,133,133	0
6	PEG	A	513	7/7	0.76	0.26	125,130,134,137	0
6	PEG	C	515	7/7	0.77	0.16	121,121,123,123	0
6	PEG	B	512	7/7	0.78	0.32	100,106,116,117	0
6	PEG	C	509	7/7	0.78	0.57	92,105,108,108	0
6	PEG	B	516	7/7	0.79	0.42	105,118,124,124	0
7	PGE	A	520	10/10	0.79	0.25	97,102,114,117	0
6	PEG	C	511	7/7	0.80	0.29	115,118,120,120	0
5	1PE	A	506	16/16	0.81	0.30	107,117,124,125	0
6	PEG	C	512	7/7	0.81	0.14	106,115,124,124	0
8	P6G	B	524	19/19	0.81	0.22	78,96,112,114	0
6	PEG	A	515	7/7	0.81	0.18	128,133,136,138	0
6	PEG	A	510	7/7	0.82	0.36	105,106,110,113	0
6	PEG	A	517	7/7	0.83	0.18	112,123,133,133	0
6	PEG	C	516	7/7	0.83	0.39	130,134,140,144	0
6	PEG	B	513	7/7	0.84	0.30	98,105,121,123	0
7	PGE	A	521	10/10	0.84	0.23	125,136,140,141	0
6	PEG	A	512	7/7	0.84	0.31	96,112,119,119	0
3	NA	B	504	1/1	0.84	0.10	112,112,112,112	0
5	1PE	C	506	16/16	0.84	0.41	91,137,155,156	0
6	PEG	A	509	7/7	0.85	0.41	87,88,95,95	0
4	DMU	B	505	33/33	0.85	0.32	93,142,154,161	0
7	PGE	C	517	10/10	0.86	0.28	124,132,137,137	0
7	PGE	A	519	10/10	0.86	0.21	81,87,108,113	0
3	NA	C	504	1/1	0.87	0.14	85,85,85,85	0
6	PEG	C	510	7/7	0.88	0.61	102,109,111,115	0
6	PEG	C	508	7/7	0.88	0.52	93,96,97,97	0
6	PEG	B	518	7/7	0.88	0.44	103,114,121,125	0
4	DMU	C	505	33/33	0.88	0.64	81,140,149,151	0
6	PEG	B	514	7/7	0.88	0.13	104,110,117,118	0
6	PEG	B	509	7/7	0.89	0.31	98,101,106,108	0
3	NA	C	503	1/1	0.89	0.09	66,66,66,66	0
7	PGE	A	518	10/10	0.90	0.15	78,94,102,104	0
6	PEG	A	516	7/7	0.91	0.83	105,108,133,140	0
6	PEG	B	510	7/7	0.92	0.30	95,98,100,102	0
6	PEG	A	508	7/7	0.93	0.37	87,89,92,93	0
3	NA	A	504	1/1	0.93	0.11	86,86,86,86	0
3	NA	B	502	1/1	0.93	0.08	68,68,68,68	0
3	NA	A	503	1/1	0.94	0.12	65,65,65,65	0

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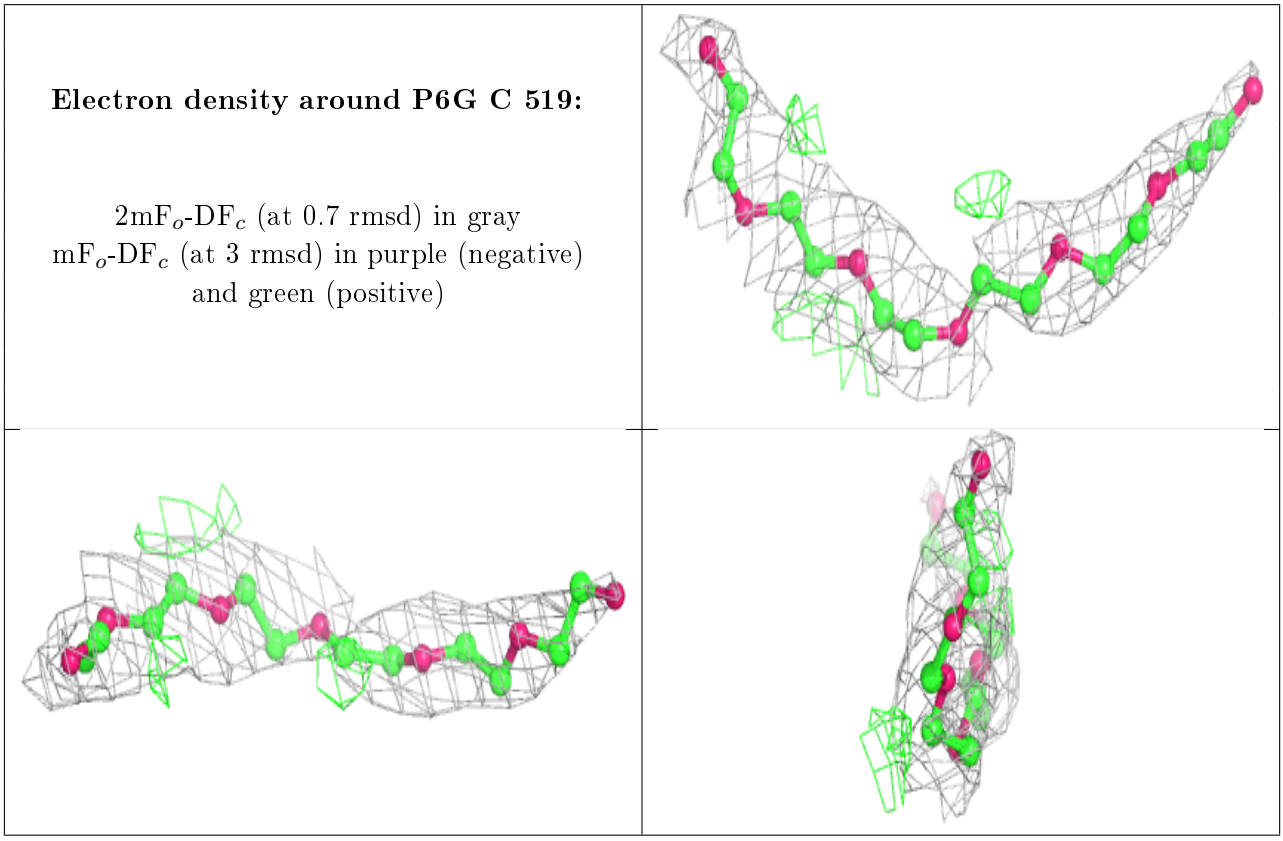
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ASP	A	501	9/9	0.95	0.21	63,70,79,87	0
3	NA	B	503	1/1	0.95	0.17	61,61,61,61	0
2	ASP	B	501	9/9	0.96	0.19	63,73,87,104	0
3	NA	A	502	1/1	0.97	0.24	63,63,63,63	0
2	ASP	C	501	9/9	0.98	0.18	71,76,81,86	0
3	NA	C	502	1/1	0.99	0.28	75,75,75,75	0

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.

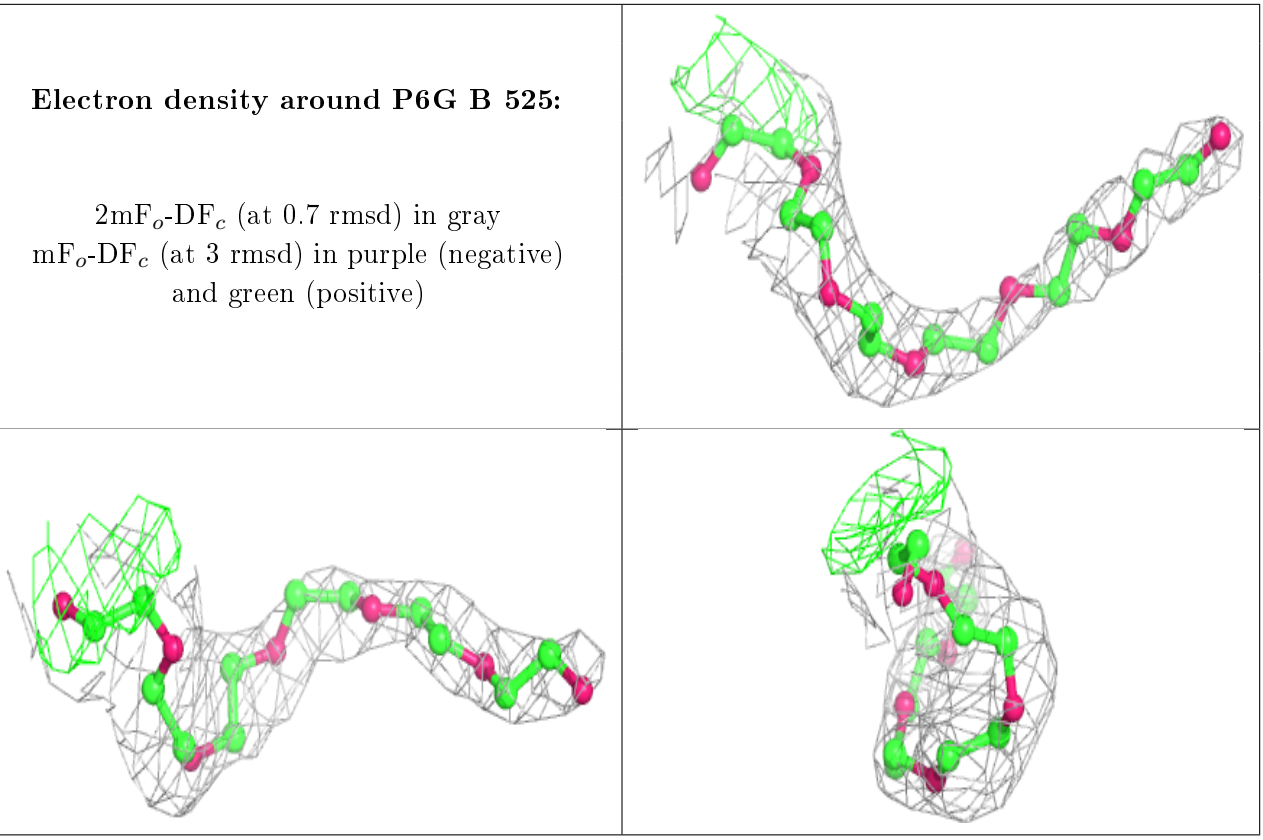
Electron density around P6G C 519:

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



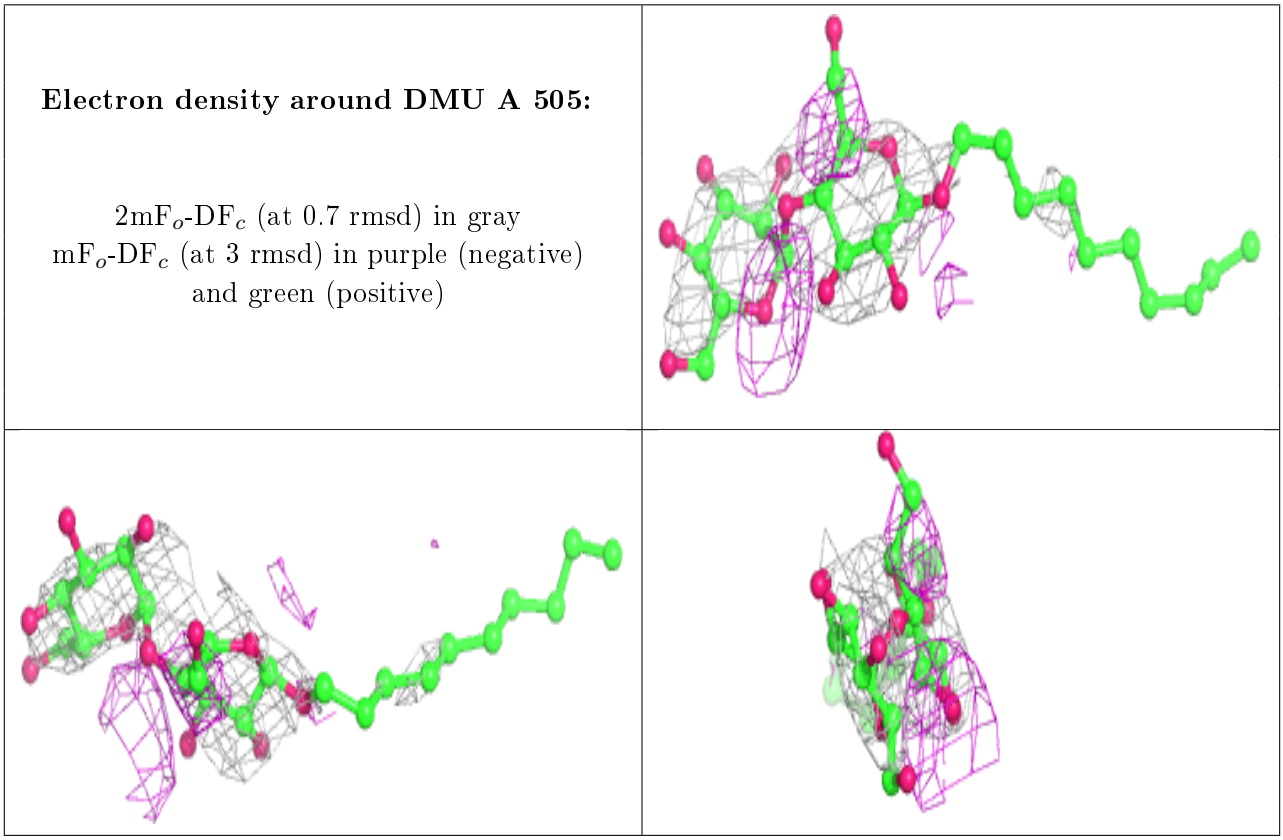
Electron density around P6G B 525:

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



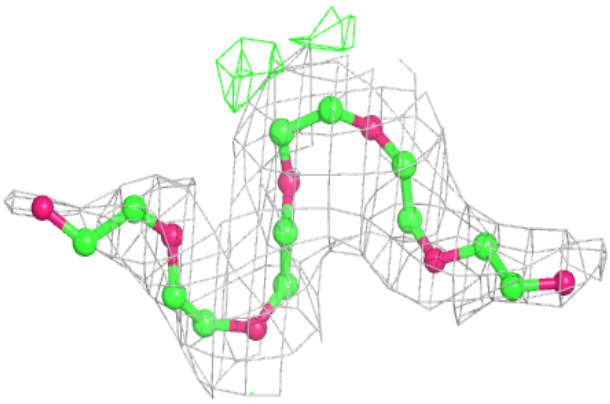
Electron density around DMU A 505:

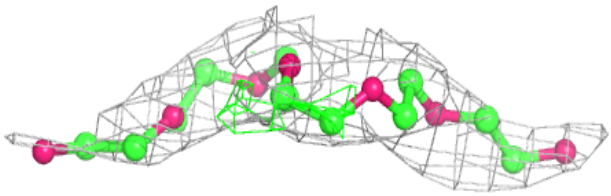
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

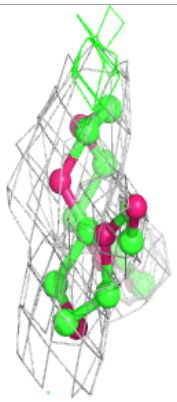


Electron density around P6G A 522:

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

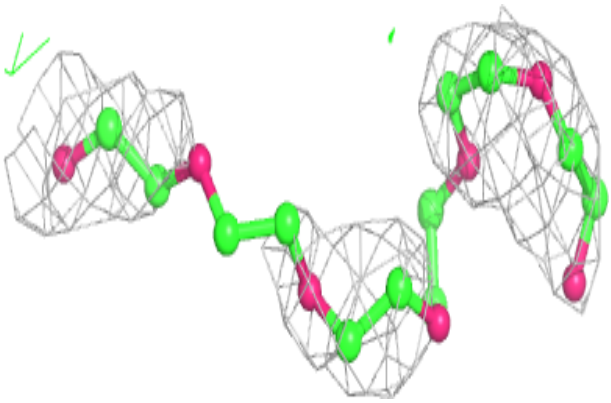


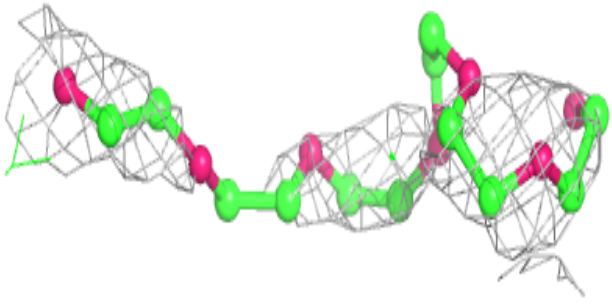


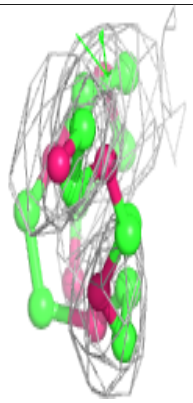


Electron density around P6G C 520:

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

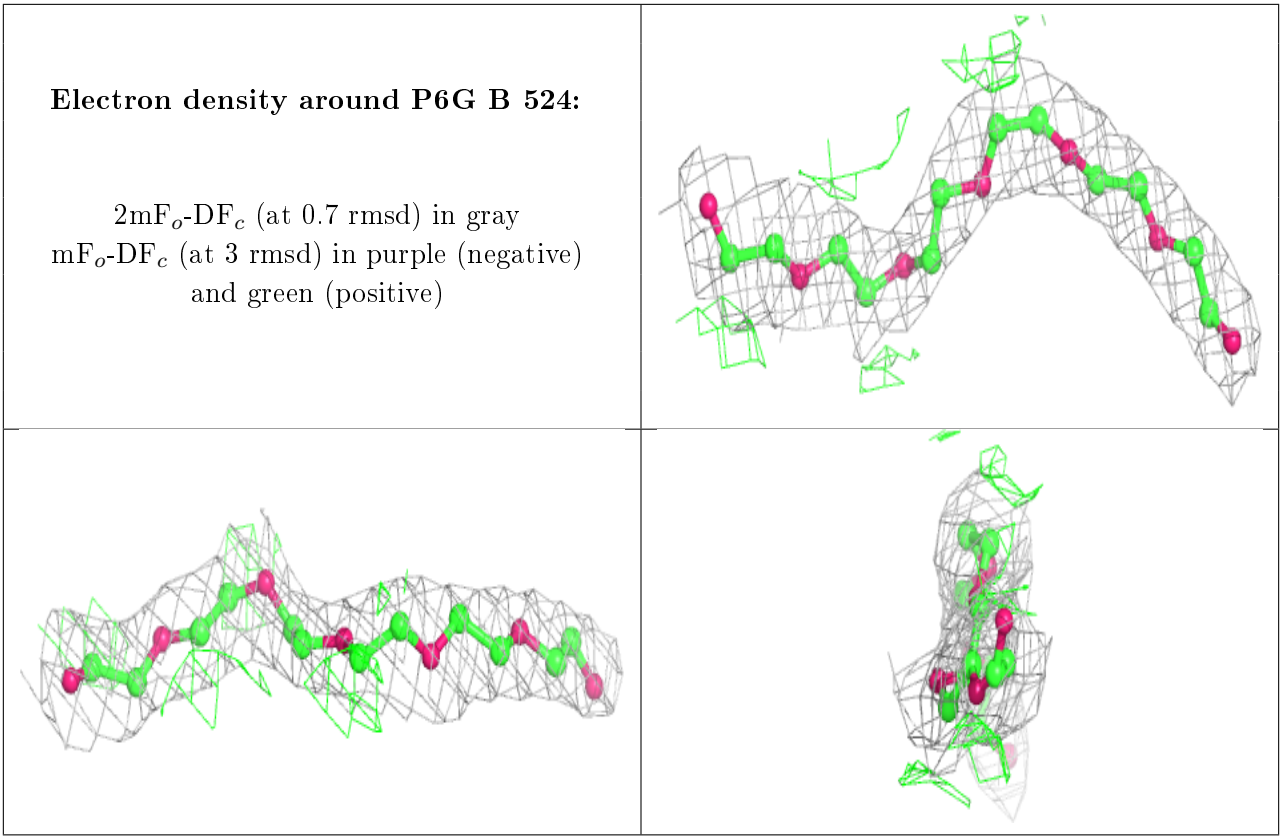






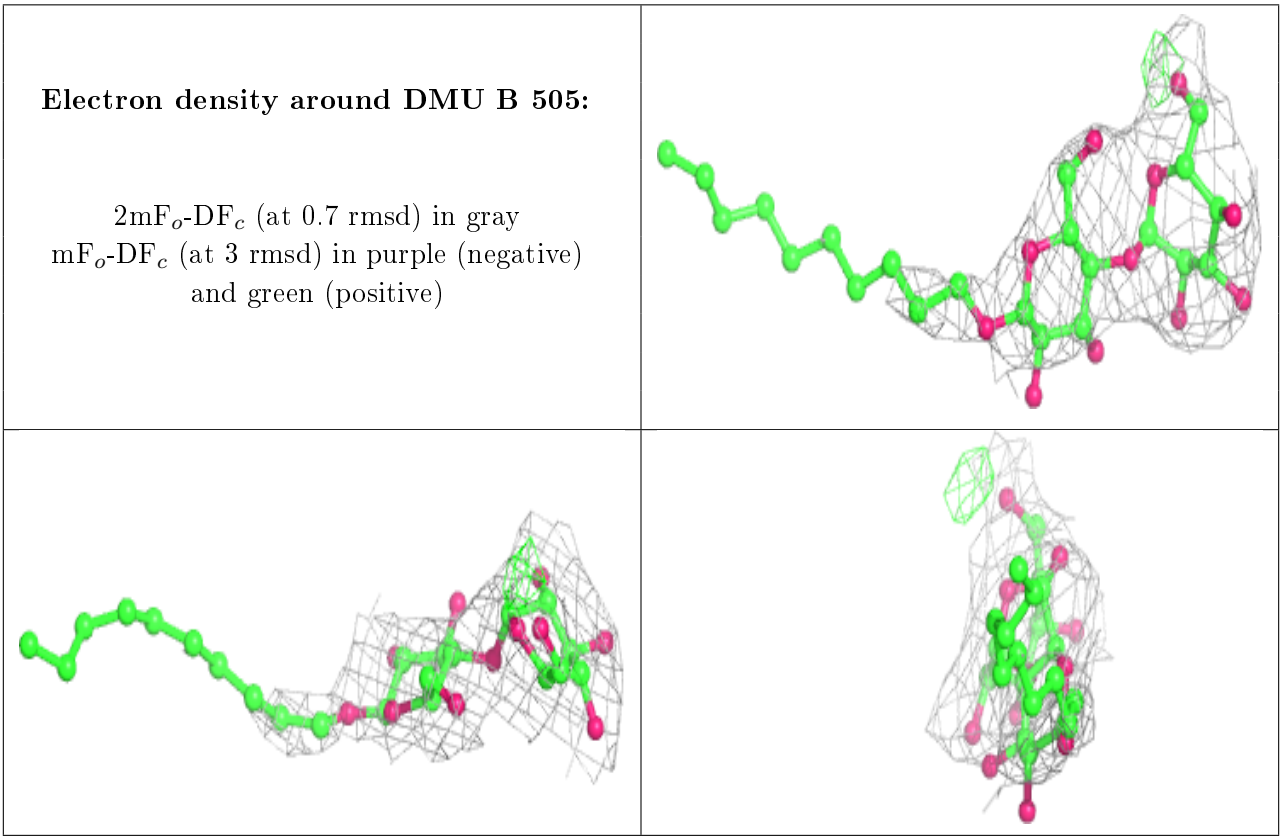
Electron density around P6G B 524:

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



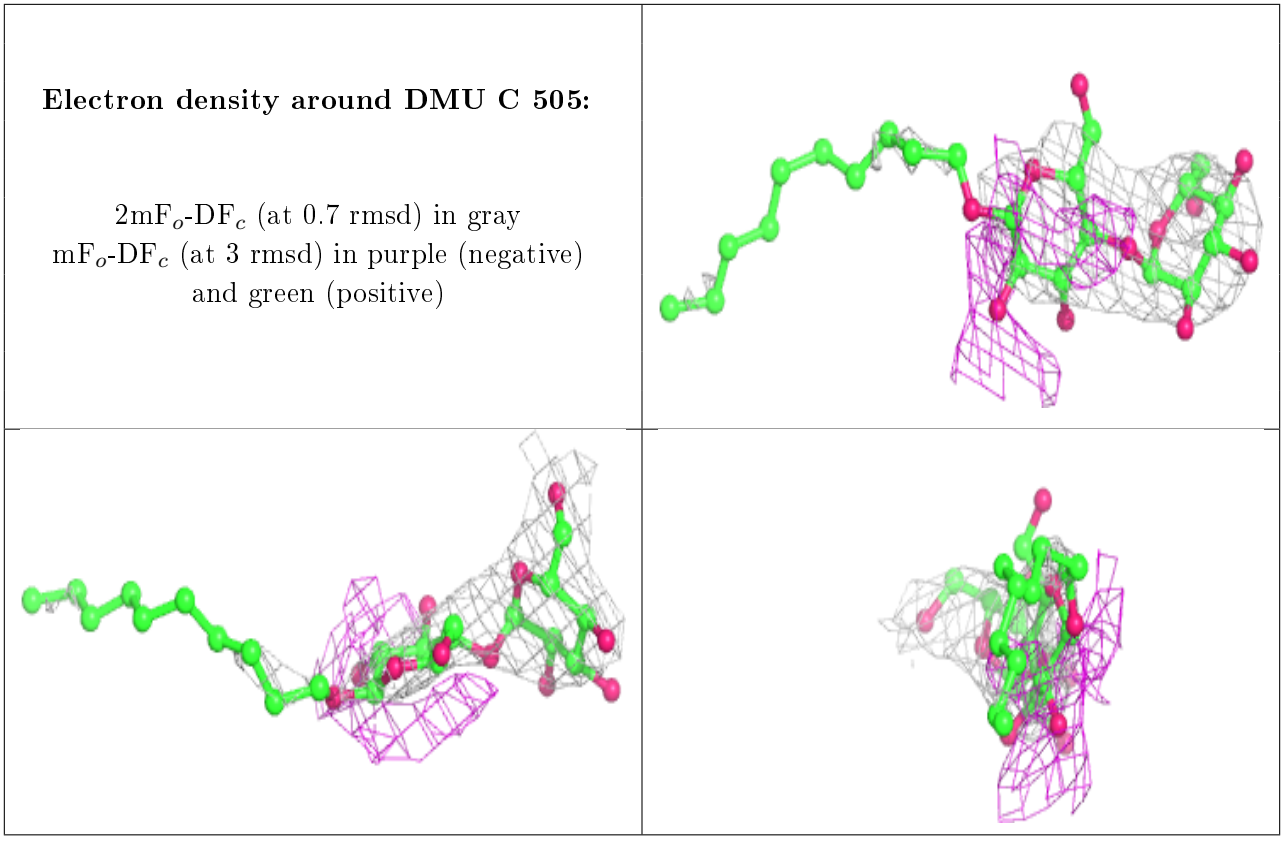
Electron density around DMU B 505:

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



Electron density around DMU C 505:

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



6.5 Other polymers

There are no such residues in this entry.