



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:39 AM UTC

PDB ID : 6ZL4 / pdb_00006zl4
Title : the structure of glutamate transporter homologue GltTk in complex with the photo switchable compound (cis)
Authors : Arkhipova, V.; Slotboom, D.J.; Guskov, A.
Deposited on : 2020-06-30
Resolution : 3.00 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) 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	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

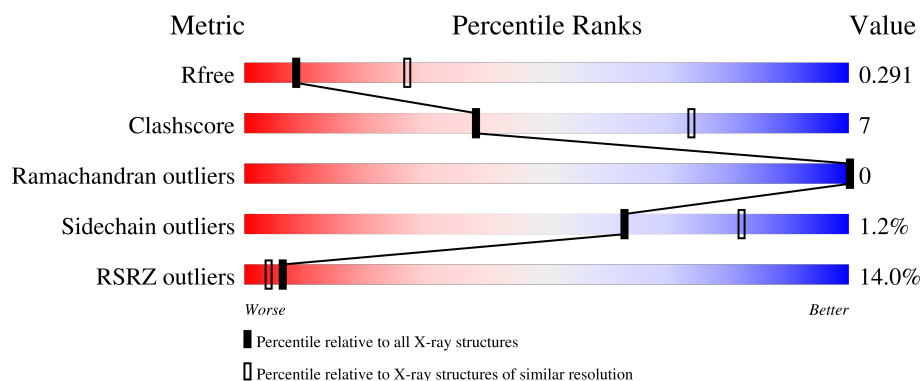
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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

Mol	Chain	Length	Quality of chain
1	A	438	
1	B	438	
1	C	438	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 9881 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	424	Total	C	N	O	S	0	0	0
			3161	2085	511	549	16			
1	B	423	Total	C	N	O	S	0	1	0
			3159	2084	511	548	16			
1	C	423	Total	C	N	O	S	0	1	0
			3159	2084	511	548	16			

There are 24 discrepancies between the modelled and reference sequences:

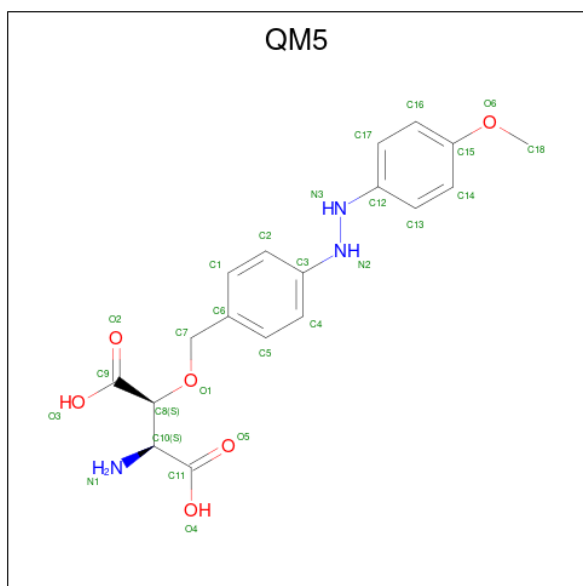
Chain	Residue	Modelled	Actual	Comment	Reference
A	431	HIS	-	expression tag	UNP Q5JID0
A	432	HIS	-	expression tag	UNP Q5JID0
A	433	HIS	-	expression tag	UNP Q5JID0
A	434	HIS	-	expression tag	UNP Q5JID0
A	435	HIS	-	expression tag	UNP Q5JID0
A	436	HIS	-	expression tag	UNP Q5JID0
A	437	HIS	-	expression tag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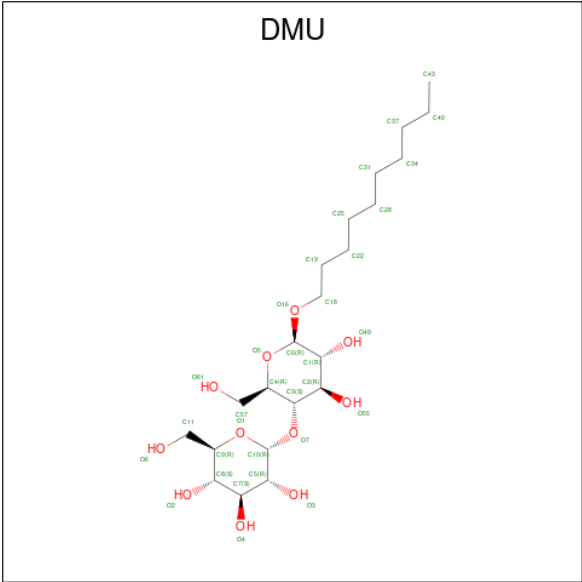
Chain	Residue	Modelled	Actual	Comment	Reference
C	438	HIS	-	expression tag	UNP Q5JID0

- Molecule 2 is (2 {S},3 {S})-2-azanyl-3-[[4-[2-(4-methoxyphenyl)hydrazinyl]phenyl]methoxy]butanedioic acid (CCD ID: QM5) (formula: C₁₈H₂₁N₃O₆).



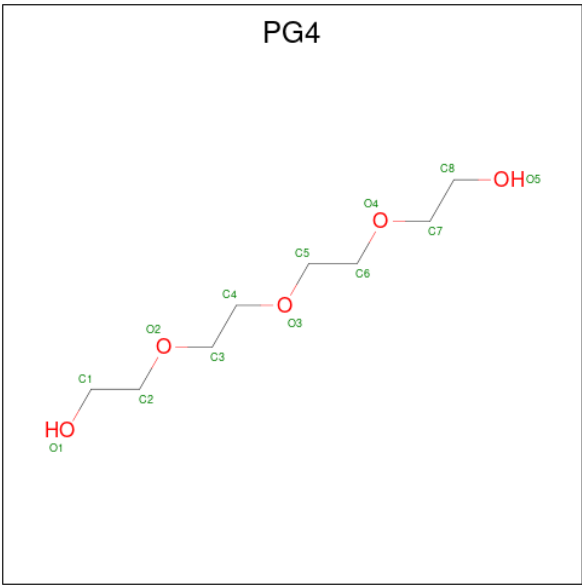
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			27	18	3	6		
2	B	1	Total	C	N	O	0	0
			27	18	3	6		
2	C	1	Total	C	N	O	0	0
			27	18	3	6		

- Molecule 3 is DECYL-BETA-D-MALTOPYRANOSIDE (CCD ID: DMU) (formula: C₂₂H₄₂O₁₁).



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

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



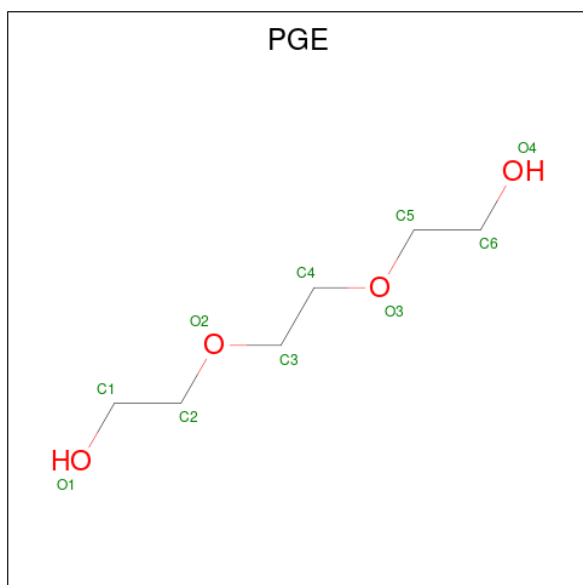
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	8	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	8	5		
4	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



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

- Molecule 6 is SODIUM ION (CCD ID: NA) (formula: Na).

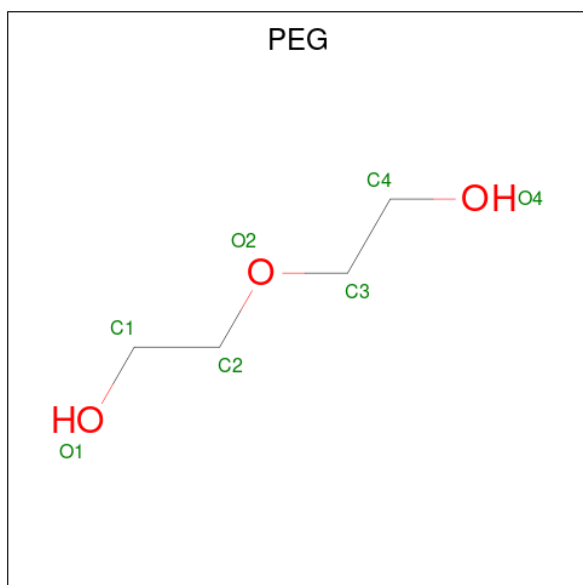
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	3	Total	Na	0	0
			3	3		

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

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



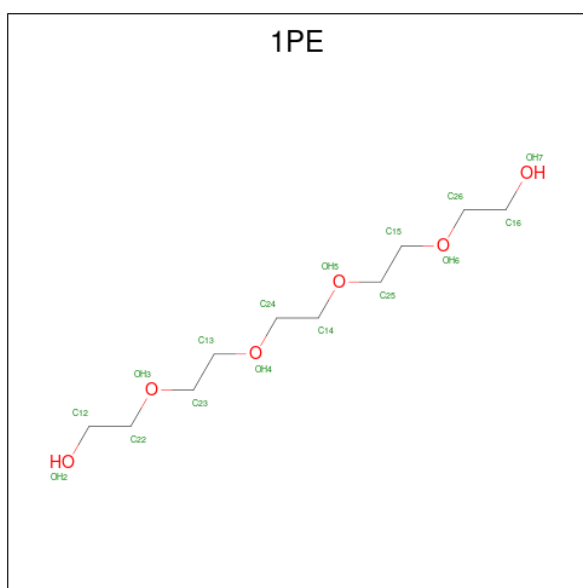
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total 7	C 4	O 3	0	0
7	A	1	Total 7	C 4	O 3	0	0
7	B	1	Total 7	C 4	O 3	0	0
7	B	1	Total 7	C 4	O 3	0	0
7	B	1	Total 7	C 4	O 3	0	0
7	B	1	Total 7	C 4	O 3	0	0
7	B	1	Total 7	C 4	O 3	0	0
7	C	1	Total 7	C 4	O 3	0	0
7	C	1	Total 7	C 4	O 3	0	0

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

- Molecule 8 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).

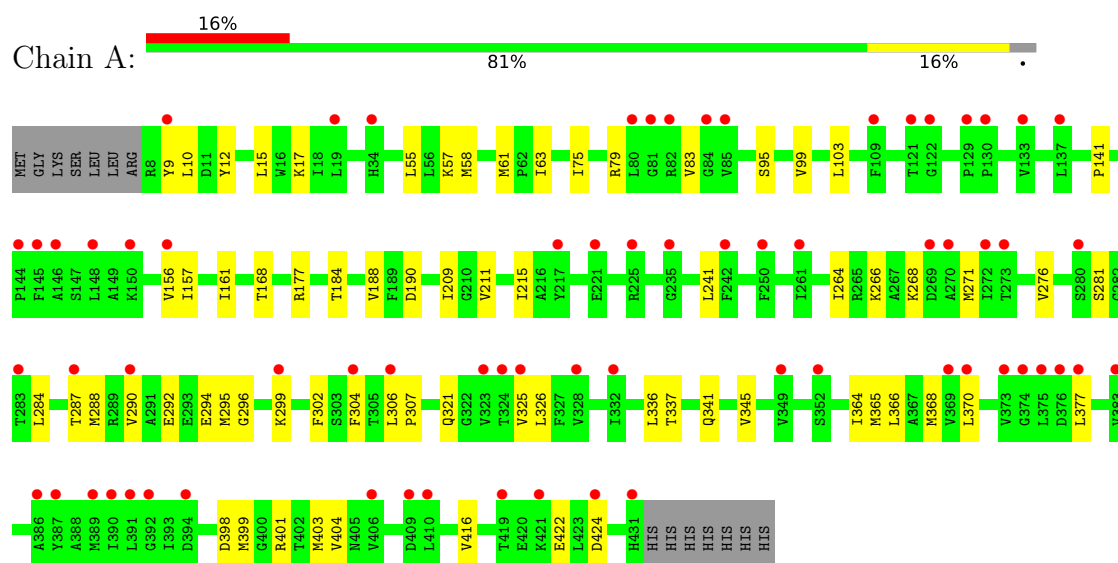


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	C	1	Total	C	O	0	0
			16	10	6		

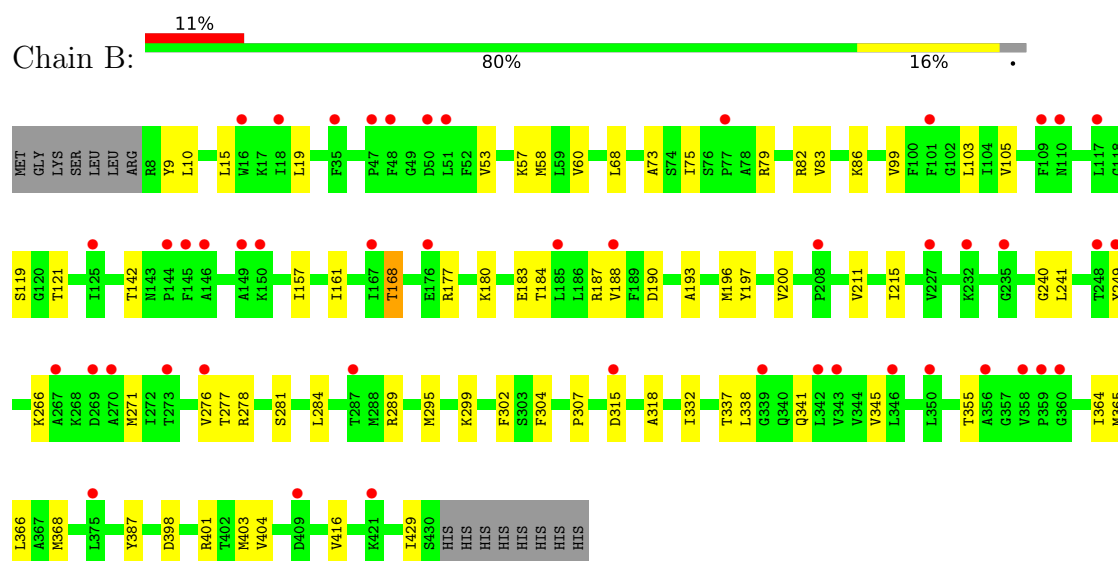
3 Residue-property plots [i](#)

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.

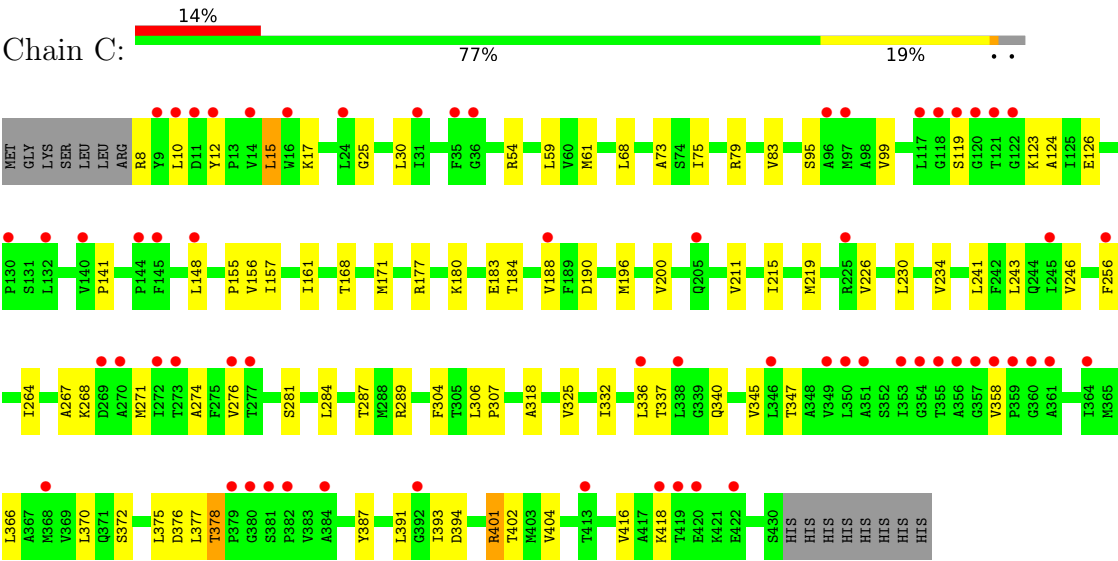
- Molecule 1: Proton/glutamate symporter, SDF family



- Molecule 1: Proton/glutamate symporter, SDF family



- Molecule 1: Proton/glutamate symporter, SDF family



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.21Å 117.21Å 309.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.23 – 3.00 48.23 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.23-3.00) 99.3 (48.23-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.96 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.255 , 0.288 0.261 , 0.291	Depositor DCC
R_{free} test set	2498 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	144.3	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 69.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9881	wwPDB-VP
Average B, all atoms (Å ²)	172.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, PGE, PG4, DMU, PEG, QM5, 1PE

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	0.18	0/3221	0.40	0/4385
1	B	0.19	0/3221	0.39	0/4384
1	C	0.18	0/3221	0.37	0/4384
All	All	0.19	0/9663	0.39	0/13153

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	A	3161	0	3359	44	0
1	B	3159	0	3365	46	0
1	C	3159	0	3365	53	0
2	A	27	0	0	0	0
2	B	27	0	0	0	0
2	C	27	0	0	1	0
3	A	33	0	42	1	0
3	B	33	0	42	1	0
3	C	33	0	42	1	0
4	A	39	0	54	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	20	0	28	0	0
5	B	30	0	42	0	0
5	C	10	0	14	0	0
6	A	3	0	0	0	0
6	B	3	0	0	0	0
6	C	3	0	0	0	0
7	A	14	0	20	0	0
7	B	35	0	50	0	0
7	C	49	0	70	1	0
8	C	16	0	22	1	0
All	All	9881	0	10515	140	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 7.

All (140) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:157:ILE:HD11	1:C:307:PRO:HB2	1.60	0.83
1:B:157:ILE:HD11	1:B:307:PRO:HB2	1.61	0.82
1:B:278:ARG:HD2	1:B:398:ASP:HB3	1.70	0.74
1:B:364:ILE:HG22	1:B:368:MET:HE2	1.69	0.73
1:A:211:VAL:HG22	1:A:276:VAL:HG21	1.72	0.70
1:A:364:ILE:HG22	1:A:368:MET:HE2	1.73	0.69
1:C:289:ARG:HH12	3:C:502:DMU:H36	1.59	0.67
1:C:401:ARG:NH2	2:C:501:QM5:O2	2.28	0.65
1:C:61:MET:HE1	1:C:156:VAL:HG21	1.78	0.64
1:B:177:ARG:NH1	1:C:190:ASP:OD2	2.33	0.61
1:C:30:LEU:HD21	1:C:219:MET:HE3	1.83	0.61
1:A:103:LEU:HD13	1:A:341:GLN:HB3	1.83	0.60
1:A:9:TYR:HE1	1:A:17:LYS:HD2	1.65	0.60
1:C:68:LEU:HD21	1:C:161:ILE:HG13	1.82	0.60
8:C:504:1PE:H242	7:C:511:PEG:H41	1.84	0.60
1:A:15:LEU:HD13	1:A:268:LYS:HZ1	1.67	0.59
1:A:370:LEU:HD12	1:A:377:LEU:HD11	1.85	0.58
1:C:211:VAL:HG22	1:C:276:VAL:HG21	1.84	0.58
1:C:211:VAL:HG13	1:C:276:VAL:HG11	1.86	0.57
1:A:211:VAL:HG13	1:A:276:VAL:HG11	1.85	0.57
1:B:365:MET:HA	1:B:368:MET:HE3	1.87	0.57
1:B:241:LEU:HB3	1:B:404:VAL:HG21	1.86	0.56
1:B:103:LEU:HD13	1:B:341:GLN:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:MET:HE1	1:A:156:VAL:HG21	1.87	0.56
1:C:12:TYR:HB3	1:C:17:LYS:HD3	1.87	0.56
1:A:266:LYS:HB2	1:A:295:MET:HE2	1.89	0.55
1:B:161:ILE:HD11	1:B:304:PHE:CD1	2.42	0.55
1:A:325:VAL:HG12	1:A:336:LEU:HD11	1.89	0.54
1:C:99:VAL:HG22	1:C:318:ALA:HB1	1.87	0.54
1:A:365:MET:HA	1:A:368:MET:HE3	1.88	0.54
1:B:75:ILE:HD11	1:B:79:ARG:HG2	1.89	0.54
1:B:99:VAL:HG11	1:B:345:VAL:HA	1.90	0.54
1:C:10:LEU:HD12	1:C:10:LEU:H	1.73	0.53
1:C:161:ILE:HD11	1:C:304:PHE:CD1	2.43	0.53
1:A:177:ARG:NH1	1:B:190:ASP:OD2	2.42	0.53
1:C:196:MET:O	1:C:200:VAL:HG23	2.09	0.53
1:C:264:ILE:HG13	1:C:271:MET:HE1	1.90	0.53
1:C:274:ALA:HB1	1:C:402:THR:HG22	1.90	0.52
1:B:398:ASP:HA	1:B:401:ARG:HG2	1.91	0.52
1:C:370:LEU:HB3	1:C:375:LEU:O	2.08	0.52
1:B:196:MET:O	1:B:200:VAL:HG23	2.10	0.52
1:C:211:VAL:O	1:C:215:ILE:HG22	2.10	0.52
1:B:83:VAL:HG13	1:B:416:VAL:HG11	1.92	0.51
1:A:157:ILE:HD11	1:A:307:PRO:HB2	1.93	0.51
1:C:267:ALA:HB3	1:C:271:MET:HE3	1.93	0.51
1:A:296:GLY:HA3	1:A:424:ASP:HB3	1.92	0.51
1:C:73:ALA:O	1:C:168:THR:OG1	2.28	0.51
1:C:366:LEU:HD23	1:C:387:TYR:HE1	1.74	0.51
1:B:289:ARG:HH22	3:B:505:DMU:H40	1.76	0.51
1:A:83:VAL:HG13	1:A:416:VAL:HG11	1.92	0.51
1:A:215:ILE:HD12	1:A:399:MET:HE1	1.93	0.51
1:A:215:ILE:HA	1:A:399:MET:HE1	1.93	0.50
1:C:241:LEU:HB3	1:C:404:VAL:HG21	1.93	0.50
1:B:99:VAL:HG22	1:B:318:ALA:HB1	1.94	0.50
1:C:184:THR:O	1:C:188:VAL:HG23	2.11	0.50
1:A:184:THR:O	1:A:188:VAL:HG23	2.11	0.50
1:B:10:LEU:H	1:B:10:LEU:HD12	1.77	0.49
1:C:99:VAL:HG11	1:C:345:VAL:HA	1.93	0.49
1:A:10:LEU:H	1:A:10:LEU:HD12	1.78	0.49
1:C:123:LYS:HG3	1:C:376:ASP:CG	2.37	0.49
1:A:290:VAL:O	1:A:294:GLU:HB2	2.13	0.48
1:C:370:LEU:HD12	1:C:377:LEU:HD11	1.94	0.48
1:C:325:VAL:HG12	1:C:336:LEU:HD11	1.95	0.48
1:A:99:VAL:HG11	1:A:345:VAL:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:ASP:HA	1:A:401:ARG:HG2	1.95	0.48
1:B:142:THR:HG23	1:C:54:ARG:HE	1.77	0.48
1:C:281:SER:O	1:C:284:LEU:HB2	2.13	0.48
1:B:284:LEU:HD11	1:B:307:PRO:HA	1.95	0.47
1:B:119:SER:HB2	1:B:332:ILE:HB	1.96	0.47
1:B:299:LYS:HA	1:B:302:PHE:CE2	2.49	0.47
1:A:63:ILE:HG23	1:A:288:MET:HE1	1.95	0.47
1:A:287:THR:HG22	1:A:306:LEU:HD13	1.97	0.47
1:C:171:MET:HE3	1:C:183:GLU:HA	1.96	0.47
1:B:19:LEU:HA	1:B:19:LEU:HD23	1.82	0.47
1:B:105:VAL:HG21	1:B:240:GLY:HA2	1.97	0.47
1:B:266:LYS:HB2	1:B:295:MET:HE2	1.97	0.47
1:A:161:ILE:HD11	1:A:304:PHE:CD1	2.50	0.46
1:C:337:THR:HG23	1:C:340:GLN:H	1.80	0.46
1:A:337:THR:O	1:A:341:GLN:HG3	2.16	0.46
1:C:83:VAL:HG13	1:C:416:VAL:HG11	1.98	0.46
1:C:226:VAL:HG13	1:C:234:VAL:HG21	1.97	0.46
1:B:366:LEU:HD23	1:B:387:TYR:HE1	1.81	0.46
1:A:281:SER:O	1:A:284:LEU:HB2	2.17	0.45
1:C:99:VAL:CG2	1:C:318:ALA:HB1	2.47	0.45
1:C:123:LYS:HG3	1:C:376:ASP:OD2	2.17	0.45
1:C:95:SER:O	1:C:99:VAL:HG23	2.15	0.45
1:B:53:VAL:HG22	1:B:277:THR:HG22	1.98	0.45
1:B:99:VAL:CG2	1:B:318:ALA:HB1	2.46	0.45
1:A:211:VAL:O	1:A:215:ILE:HG22	2.17	0.45
1:A:292:GLU:HB2	1:A:302:PHE:CZ	2.51	0.45
1:B:249:TYR:OH	1:B:315:ASP:OD2	2.28	0.45
1:B:73:ALA:O	1:B:168:THR:HG21	2.17	0.44
1:A:141:PRO:O	1:B:58:MET:HB2	2.17	0.44
1:B:68:LEU:HD23	1:B:157:ILE:HA	2.00	0.44
1:A:75:ILE:HD11	1:A:79:ARG:HG2	2.00	0.44
1:B:355:THR:HG21	1:B:365:MET:SD	2.58	0.44
1:B:57:LYS:HA	1:B:60:VAL:HG23	1.99	0.44
1:C:59:LEU:HD13	1:C:200:VAL:HG22	1.99	0.44
1:C:119:SER:HB2	1:C:332:ILE:HB	2.00	0.43
1:A:55:LEU:HA	1:A:58:MET:HE2	2.00	0.43
1:C:180:LYS:HA	1:C:180:LYS:HD2	1.86	0.43
1:A:190:ASP:OD2	1:C:177:ARG:NH1	2.52	0.43
1:C:75:ILE:HD11	1:C:79:ARG:HG2	2.01	0.43
1:A:57:LYS:HA	1:A:57:LYS:HD2	1.94	0.43
1:A:241:LEU:HB3	1:A:404:VAL:HG21	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:15:LEU:HB2	1:C:268:LYS:HE2	2.01	0.42
1:C:141:PRO:HB3	1:C:155:PRO:HB3	2.01	0.42
1:C:347:THR:HG21	1:C:372:SER:HB2	2.00	0.42
1:C:391:LEU:HA	1:C:394:ASP:HB2	2.00	0.42
1:A:299:LYS:HA	1:A:302:PHE:CE2	2.54	0.42
1:B:271:MET:O	1:B:403:MET:HG3	2.18	0.42
1:C:230:LEU:HB3	1:C:393:ILE:HD13	2.01	0.42
1:C:126:GLU:HG3	1:C:377:LEU:HB2	2.01	0.42
1:B:184:THR:O	1:B:188:VAL:HG23	2.18	0.42
1:B:211:VAL:O	1:B:215:ILE:HG22	2.19	0.42
1:B:211:VAL:HG22	1:B:276:VAL:HG21	2.01	0.42
1:A:271:MET:O	1:A:403:MET:HG3	2.20	0.42
1:A:284:LEU:HD23	1:A:284:LEU:HA	1.91	0.42
1:A:79:ARG:NH2	1:A:422:GLU:OE2	2.53	0.42
1:C:25:GLY:HA3	1:C:215:ILE:HG23	2.02	0.42
1:B:180:LYS:HD2	1:B:180:LYS:HA	1.89	0.41
1:B:281:SER:O	1:B:284:LEU:HB2	2.20	0.41
1:A:321:GLN:HG2	1:A:366:LEU:HD13	2.00	0.41
1:C:243:LEU:HA	1:C:246:VAL:HG22	2.02	0.41
1:C:287:THR:HG22	1:C:306:LEU:HD13	2.02	0.41
1:B:337:THR:O	1:B:341:GLN:HG3	2.20	0.41
1:B:403:MET:HE2	1:B:403:MET:HB3	1.94	0.41
1:C:124:ALA:HB3	1:C:378:THR:HG23	2.01	0.41
1:A:10:LEU:HD21	3:A:502:DMU:H9	2.03	0.41
1:A:326:LEU:HD23	1:A:336:LEU:HD12	2.02	0.41
4:A:503:PG4:H22	4:A:503:PG4:H82	2.03	0.41
1:B:15:LEU:H	1:B:15:LEU:HG	1.68	0.41
1:B:82:ARG:HG2	1:B:86:LYS:NZ	2.35	0.41
1:B:183:GLU:O	1:B:187[A]:ARG:HG3	2.21	0.41
1:A:12:TYR:HB3	1:A:17:LYS:HD3	2.03	0.40
1:B:193:ALA:O	1:B:197:TYR:HD1	2.03	0.40
1:C:61:MET:HE3	1:C:148:LEU:HD23	2.02	0.40
1:C:256:PHE:O	1:C:418:LYS:NZ	2.55	0.40
1:A:95:SER:O	1:A:99:VAL:HG23	2.22	0.40
1:B:338:LEU:HD12	1:B:338:LEU:HA	1.97	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	422/438 (96%)	409 (97%)	13 (3%)	0	100	100
1	B	422/438 (96%)	408 (97%)	14 (3%)	0	100	100
1	C	422/438 (96%)	408 (97%)	14 (3%)	0	100	100
All	All	1266/1314 (96%)	1225 (97%)	41 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	332/345 (96%)	329 (99%)	3 (1%)	70	85
1	B	332/345 (96%)	328 (99%)	4 (1%)	63	82
1	C	332/345 (96%)	327 (98%)	5 (2%)	57	80
All	All	996/1035 (96%)	984 (99%)	12 (1%)	63	82

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	168	THR
1	A	209	ILE
1	A	264	ILE
1	B	9	TYR
1	B	121	THR

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Mol	Chain	Res	Type
1	B	168	THR
1	B	429	ILE
1	C	8	ARG
1	C	15	LEU
1	C	358	VAL
1	C	378	THR
1	C	401	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	134	GLN
1	A	321	GLN
1	B	34	HIS
1	B	172	ASN
1	B	341	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](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 39 ligands modelled in this entry, 9 are monoatomic - leaving 30 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	PGE	A	505	-	9,9,9	0.32	0	8,8,8	0.26	0
3	DMU	A	502	-	34,34,34	1.52	8 (23%)	45,45,45	1.15	5 (11%)
7	PEG	B	512	-	6,6,6	0.51	0	5,5,5	0.26	0
5	PGE	B	502	-	9,9,9	0.33	0	8,8,8	0.25	0
5	PGE	B	507	-	9,9,9	0.32	0	8,8,8	0.32	0
5	PGE	B	506	-	9,9,9	0.32	0	8,8,8	0.31	0
8	1PE	C	504	-	15,15,15	0.54	0	14,14,14	0.21	0
7	PEG	A	512	-	6,6,6	0.51	0	5,5,5	0.26	0
4	PG4	A	506	-	12,12,12	0.54	0	11,11,11	0.24	0
7	PEG	C	509	-	6,6,6	0.52	0	5,5,5	0.41	0
7	PEG	B	504	-	6,6,6	0.50	0	5,5,5	0.26	0
2	QM5	A	501	6	27,28,28	1.41	4 (14%)	34,37,37	1.16	4 (11%)
4	PG4	A	503	-	12,12,12	0.54	0	11,11,11	0.24	0
7	PEG	A	511	-	6,6,6	0.50	0	5,5,5	0.26	0
5	PGE	C	503	-	9,9,9	0.32	0	8,8,8	0.31	0
3	DMU	C	502	-	34,34,34	1.50	7 (20%)	45,45,45	1.34	6 (13%)
7	PEG	C	510	-	6,6,6	0.51	0	5,5,5	0.25	0
7	PEG	C	513	-	6,6,6	0.50	0	5,5,5	0.27	0
7	PEG	C	512	-	6,6,6	0.51	0	5,5,5	0.26	0
7	PEG	B	513	-	6,6,6	0.50	0	5,5,5	0.29	0
7	PEG	B	511	-	6,6,6	0.50	0	5,5,5	0.22	0
2	QM5	C	501	-	27,28,28	1.43	4 (14%)	34,37,37	1.22	3 (8%)
3	DMU	B	505	-	34,34,34	1.50	7 (20%)	45,45,45	1.37	6 (13%)
4	PG4	A	504	-	12,12,12	0.54	0	11,11,11	0.18	0
2	QM5	B	501	-	27,28,28	1.46	4 (14%)	34,37,37	1.15	3 (8%)
7	PEG	C	514	-	6,6,6	0.50	0	5,5,5	0.24	0
7	PEG	C	511	-	6,6,6	0.50	0	5,5,5	0.34	0
5	PGE	A	507	-	9,9,9	0.32	0	8,8,8	0.31	0
7	PEG	B	503	-	6,6,6	0.51	0	5,5,5	0.24	0
7	PEG	C	508	-	6,6,6	0.51	0	5,5,5	0.30	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PGE	A	505	-	-	2/7/7/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DMU	A	502	-	-	8/19/59/59	0/2/2/2
7	PEG	B	512	-	-	2/4/4/4	-
5	PGE	B	502	-	-	3/7/7/7	-
5	PGE	B	507	-	-	4/7/7/7	-
5	PGE	B	506	-	-	5/7/7/7	-
8	1PE	C	504	-	-	9/13/13/13	-
7	PEG	A	512	-	-	0/4/4/4	-
4	PG4	A	506	-	-	4/10/10/10	-
7	PEG	C	509	-	-	2/4/4/4	-
7	PEG	B	504	-	-	1/4/4/4	-
2	QM5	A	501	6	-	10/24/24/24	0/2/2/2
4	PG4	A	503	-	-	5/10/10/10	-
7	PEG	A	511	-	-	0/4/4/4	-
5	PGE	C	503	-	-	4/7/7/7	-
3	DMU	C	502	-	-	7/19/59/59	0/2/2/2
7	PEG	C	510	-	-	2/4/4/4	-
7	PEG	C	513	-	-	1/4/4/4	-
7	PEG	C	512	-	-	1/4/4/4	-
7	PEG	B	513	-	-	1/4/4/4	-
7	PEG	B	511	-	-	3/4/4/4	-
2	QM5	C	501	-	-	11/24/24/24	0/2/2/2
3	DMU	B	505	-	-	9/19/59/59	0/2/2/2
4	PG4	A	504	-	-	3/10/10/10	-
2	QM5	B	501	-	-	7/24/24/24	0/2/2/2
7	PEG	C	514	-	-	2/4/4/4	-
7	PEG	C	511	-	-	2/4/4/4	-
5	PGE	A	507	-	-	3/7/7/7	-
7	PEG	B	503	-	-	3/4/4/4	-
7	PEG	C	508	-	-	0/4/4/4	-

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	505	DMU	O1-C9	4.01	1.54	1.44
3	A	502	DMU	O1-C9	4.00	1.54	1.44
3	C	502	DMU	O1-C9	3.85	1.53	1.44
2	B	501	QM5	C3-N2	3.60	1.50	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	QM5	C3-N2	3.52	1.50	1.39
2	C	501	QM5	C3-N2	3.42	1.50	1.39
2	B	501	QM5	N3-N2	3.01	1.48	1.39
2	A	501	QM5	N3-N2	2.91	1.48	1.39
3	C	502	DMU	C11-C9	-2.85	1.42	1.51
3	A	502	DMU	C11-C9	-2.84	1.42	1.51
2	C	501	QM5	N3-N2	2.84	1.47	1.39
3	B	505	DMU	C11-C9	-2.83	1.42	1.51
3	A	502	DMU	O1-C10	2.76	1.48	1.41
3	C	502	DMU	O1-C10	2.75	1.48	1.41
2	B	501	QM5	C12-N3	2.62	1.47	1.39
2	A	501	QM5	C12-N3	2.57	1.47	1.39
2	C	501	QM5	C12-N3	2.56	1.47	1.39
3	A	502	DMU	C7-C5	-2.51	1.45	1.52
3	B	505	DMU	O5-C6	2.47	1.48	1.41
3	C	502	DMU	C7-C5	-2.46	1.46	1.52
3	B	505	DMU	O3-C5	2.42	1.49	1.43
3	A	502	DMU	O5-C6	2.42	1.48	1.41
3	B	505	DMU	O1-C10	2.41	1.48	1.41
3	B	505	DMU	O4-C7	2.38	1.48	1.43
3	A	502	DMU	O4-C7	2.33	1.48	1.43
3	A	502	DMU	O3-C5	2.33	1.48	1.43
3	C	502	DMU	O3-C5	2.32	1.48	1.43
3	C	502	DMU	O4-C7	2.31	1.48	1.43
3	C	502	DMU	O5-C6	2.28	1.47	1.41
3	B	505	DMU	C7-C5	-2.27	1.46	1.52
2	A	501	QM5	C10-C11	2.19	1.57	1.52
2	B	501	QM5	C10-C11	2.15	1.56	1.52
2	C	501	QM5	C10-C11	2.12	1.56	1.52
3	A	502	DMU	O5-C4	2.01	1.49	1.44

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	502	DMU	O1-C10-C5	4.11	118.82	110.37
3	B	505	DMU	C8-C7-C5	3.53	117.03	110.83
2	C	501	QM5	O1-C8-C9	-3.50	107.04	112.26
3	B	505	DMU	C7-C8-C9	3.37	116.34	110.23
3	C	502	DMU	C10-C5-C7	3.24	116.83	110.01
3	C	502	DMU	C10-O7-C3	-3.11	110.60	117.98
3	B	505	DMU	C1-C2-C3	2.94	116.35	109.68
2	A	501	QM5	O1-C7-C6	-2.92	103.20	109.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	DMU	O1-C10-C5	2.89	116.31	110.37
3	B	505	DMU	C10-O7-C3	-2.87	111.18	117.98
2	C	501	QM5	O1-C7-C6	-2.77	103.56	109.91
3	B	505	DMU	C10-C5-C7	2.74	115.77	110.01
3	C	502	DMU	C10-O1-C9	2.68	118.96	113.72
3	A	502	DMU	C10-C5-C7	2.63	115.55	110.01
2	B	501	QM5	C7-O1-C8	2.60	117.21	113.55
2	C	501	QM5	C7-O1-C8	2.49	117.05	113.55
3	B	505	DMU	C6-C1-C2	2.46	115.19	110.01
3	C	502	DMU	C1-C2-C3	2.25	114.79	109.68
3	A	502	DMU	C10-O7-C3	-2.25	112.64	117.98
3	A	502	DMU	C10-O1-C9	2.23	118.08	113.72
2	B	501	QM5	O3-C9-C8	2.21	121.60	113.64
2	A	501	QM5	C7-O1-C8	2.19	116.64	113.55
2	A	501	QM5	O3-C9-C8	2.19	121.56	113.64
3	A	502	DMU	O5-C4-C3	2.15	114.17	109.72
3	C	502	DMU	O5-C4-C3	2.14	114.14	109.72
2	B	501	QM5	O1-C7-C6	-2.12	105.04	109.91
2	A	501	QM5	O3-C9-O2	-2.06	119.40	124.08

There are no chirality outliers.

All (114) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	QM5	O1-C8-C9-O2
2	A	501	QM5	O1-C8-C9-O3
2	A	501	QM5	C9-C8-O1-C7
2	A	501	QM5	C13-C12-N3-N2
2	A	501	QM5	C17-C12-N3-N2
2	A	501	QM5	C3-N2-N3-C12
2	B	501	QM5	O1-C8-C9-O2
2	B	501	QM5	O1-C8-C9-O3
2	B	501	QM5	C9-C8-O1-C7
2	B	501	QM5	C13-C12-N3-N2
2	B	501	QM5	C17-C12-N3-N2
2	B	501	QM5	C3-N2-N3-C12
2	C	501	QM5	O1-C8-C9-O2
2	C	501	QM5	O1-C8-C9-O3
2	C	501	QM5	C11-C10-C8-C9
2	C	501	QM5	C17-C12-N3-N2
2	C	501	QM5	C4-C3-N2-N3
2	C	501	QM5	C3-N2-N3-C12

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Mol	Chain	Res	Type	Atoms
7	C	509	PEG	C1-C2-O2-C3
3	A	502	DMU	O6-C11-C9-O1
2	C	501	QM5	C13-C12-N3-N2
2	C	501	QM5	C2-C3-N2-N3
3	C	502	DMU	O6-C11-C9-O1
3	B	505	DMU	O6-C11-C9-O1
3	C	502	DMU	O6-C11-C9-C8
3	B	505	DMU	O6-C11-C9-C8
5	A	507	PGE	O2-C3-C4-O3
5	B	506	PGE	O2-C3-C4-O3
3	A	502	DMU	O6-C11-C9-C8
5	C	503	PGE	O2-C3-C4-O3
4	A	506	PG4	O2-C3-C4-O3
4	A	506	PG4	O4-C7-C8-O5
8	C	504	1PE	OH6-C15-C25-OH5
5	B	502	PGE	O1-C1-C2-O2
5	B	506	PGE	O1-C1-C2-O2
3	C	502	DMU	C25-C28-C31-C34
3	B	505	DMU	C18-C19-C22-C25
3	A	502	DMU	C25-C28-C31-C34
4	A	503	PG4	O4-C7-C8-O5
8	C	504	1PE	OH2-C12-C22-OH3
3	B	505	DMU	O5-C4-C57-O61
7	B	513	PEG	O1-C1-C2-O2
3	B	505	DMU	C19-C22-C25-C28
5	A	507	PGE	O1-C1-C2-O2
7	C	509	PEG	O2-C3-C4-O4
4	A	503	PG4	O2-C3-C4-O3
3	B	505	DMU	C25-C28-C31-C34
3	C	502	DMU	C28-C31-C34-C37
3	A	502	DMU	O5-C4-C57-O61
7	C	510	PEG	O2-C3-C4-O4
5	B	507	PGE	O1-C1-C2-O2
2	A	501	QM5	C6-C7-O1-C8
3	B	505	DMU	C22-C25-C28-C31
7	B	511	PEG	O1-C1-C2-O2
3	C	502	DMU	C19-C18-O16-C6
8	C	504	1PE	OH4-C13-C23-OH3
7	B	511	PEG	C4-C3-O2-C2
3	A	502	DMU	C18-C19-C22-C25
5	B	507	PGE	O2-C3-C4-O3
7	C	511	PEG	C1-C2-O2-C3

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Mol	Chain	Res	Type	Atoms
7	C	511	PEG	O1-C1-C2-O2
2	B	501	QM5	C6-C7-O1-C8
7	C	510	PEG	C1-C2-O2-C3
5	B	506	PGE	C3-C4-O3-C5
8	C	504	1PE	C12-C22-OH3-C23
5	A	505	PGE	C1-C2-O2-C3
5	B	507	PGE	C4-C3-O2-C2
4	A	504	PG4	O2-C3-C4-O3
5	B	506	PGE	C1-C2-O2-C3
7	B	512	PEG	C1-C2-O2-C3
8	C	504	1PE	C24-C14-OH5-C25
4	A	503	PG4	C3-C4-O3-C5
2	C	501	QM5	N1-C10-C8-O1
7	B	503	PEG	C1-C2-O2-C3
7	B	503	PEG	C4-C3-O2-C2
5	C	503	PGE	O3-C5-C6-O4
3	C	502	DMU	O16-C18-C19-C22
7	C	513	PEG	C1-C2-O2-C3
8	C	504	1PE	C25-C15-OH6-C26
8	C	504	1PE	C16-C26-OH6-C15
7	B	504	PEG	C1-C2-O2-C3
2	A	501	QM5	C8-C10-C11-O5
8	C	504	1PE	OH7-C16-C26-OH6
5	C	503	PGE	C3-C4-O3-C5
5	A	505	PGE	C4-C3-O2-C2
3	A	502	DMU	C28-C31-C34-C37
5	B	502	PGE	C1-C2-O2-C3
5	B	507	PGE	C1-C2-O2-C3
7	C	514	PEG	C1-C2-O2-C3
3	A	502	DMU	O5-C6-O16-C18
3	B	505	DMU	O5-C6-O16-C18
3	B	505	DMU	C31-C34-C37-C40
3	C	502	DMU	C22-C25-C28-C31
5	C	503	PGE	C1-C2-O2-C3
7	B	511	PEG	C1-C2-O2-C3
4	A	503	PG4	C1-C2-O2-C3
4	A	506	PG4	C8-C7-O4-C6
2	A	501	QM5	C10-C8-C9-O3
4	A	506	PG4	C5-C6-O4-C7
3	A	502	DMU	C19-C22-C25-C28
5	B	506	PGE	C4-C3-O2-C2
7	C	512	PEG	O2-C3-C4-O4

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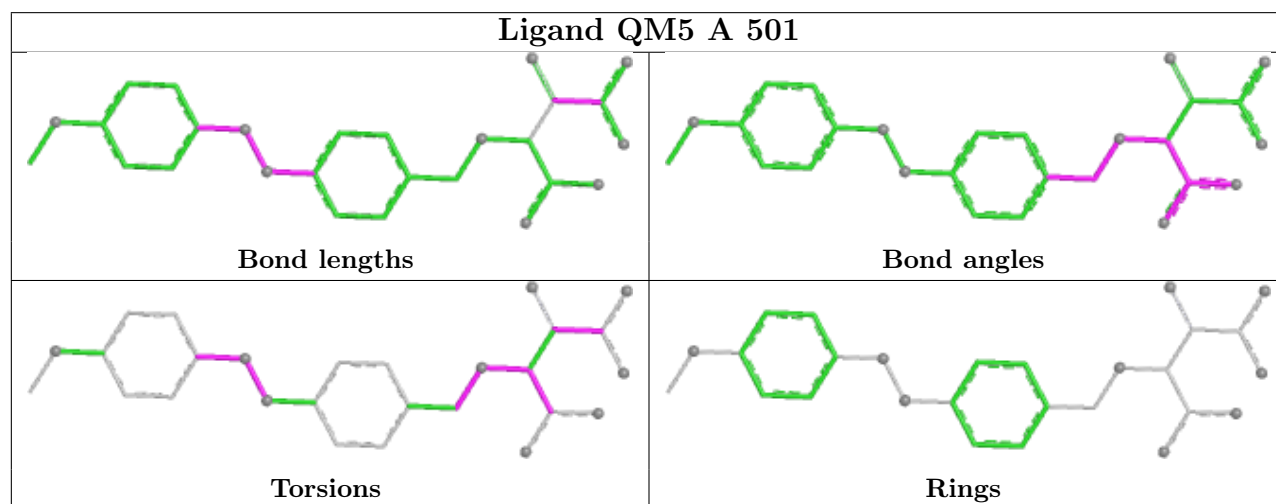
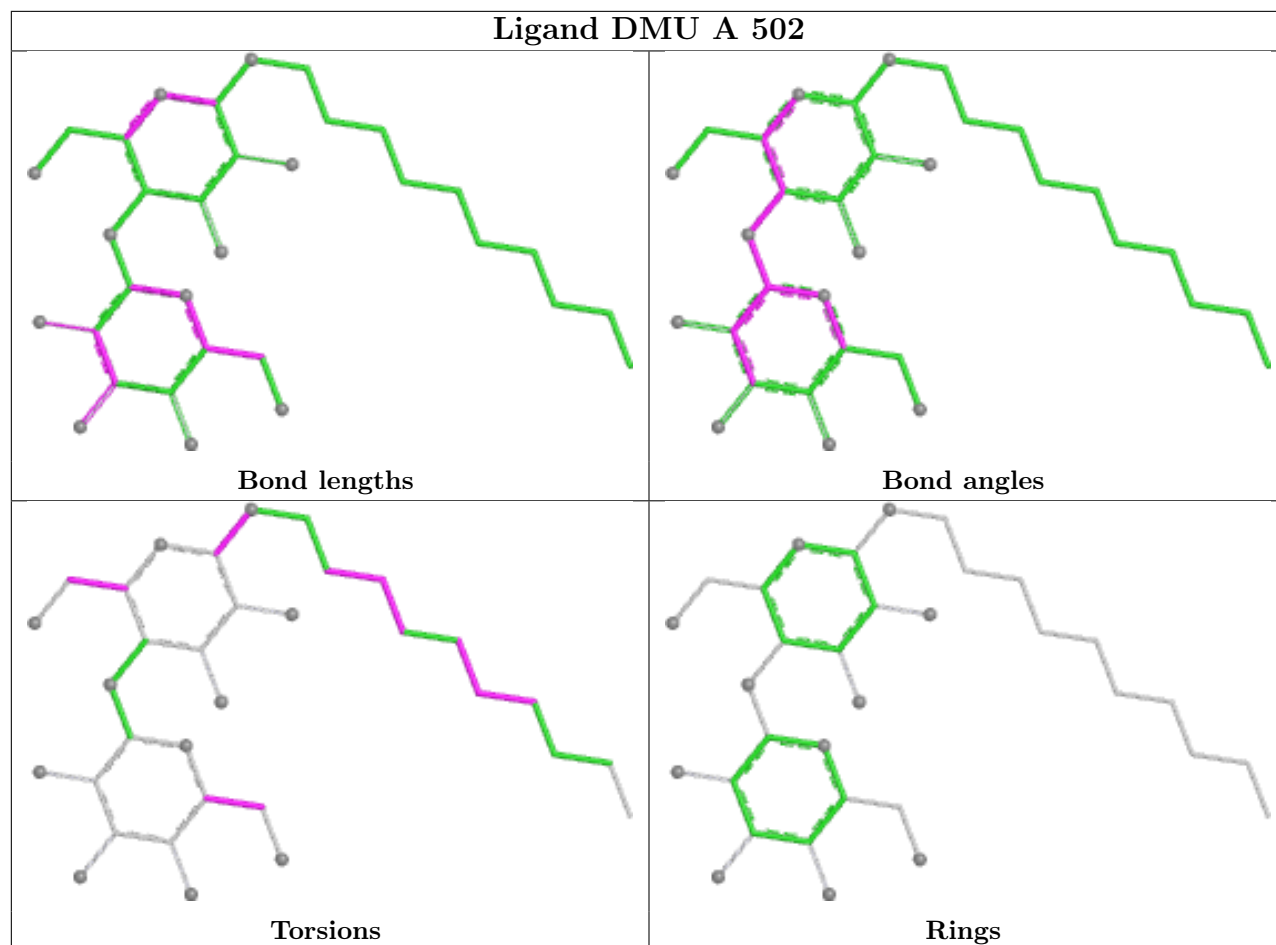
Mol	Chain	Res	Type	Atoms
4	A	504	PG4	C5-C6-O4-C7
5	A	507	PGE	C3-C4-O3-C5
4	A	504	PG4	O3-C5-C6-O4
5	B	502	PGE	C4-C3-O2-C2
7	B	503	PEG	O1-C1-C2-O2
2	C	501	QM5	C10-C8-O1-C7
4	A	503	PG4	C6-C5-O3-C4
2	C	501	QM5	N1-C10-C11-O4
7	B	512	PEG	C4-C3-O2-C2
8	C	504	1PE	OH5-C14-C24-OH4
7	C	514	PEG	C4-C3-O2-C2
2	A	501	QM5	C10-C8-C9-O2

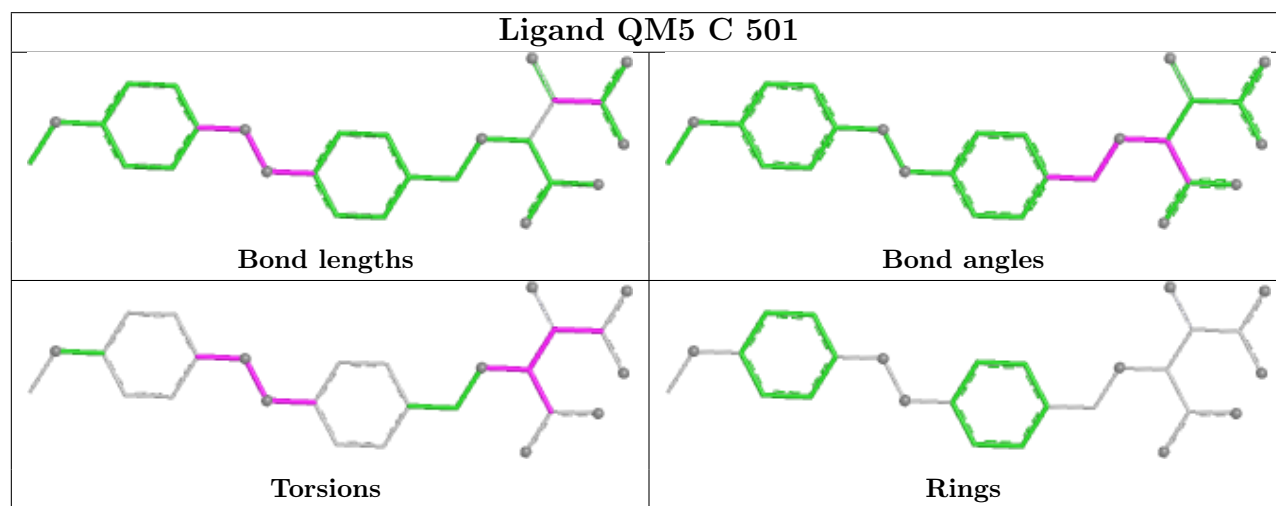
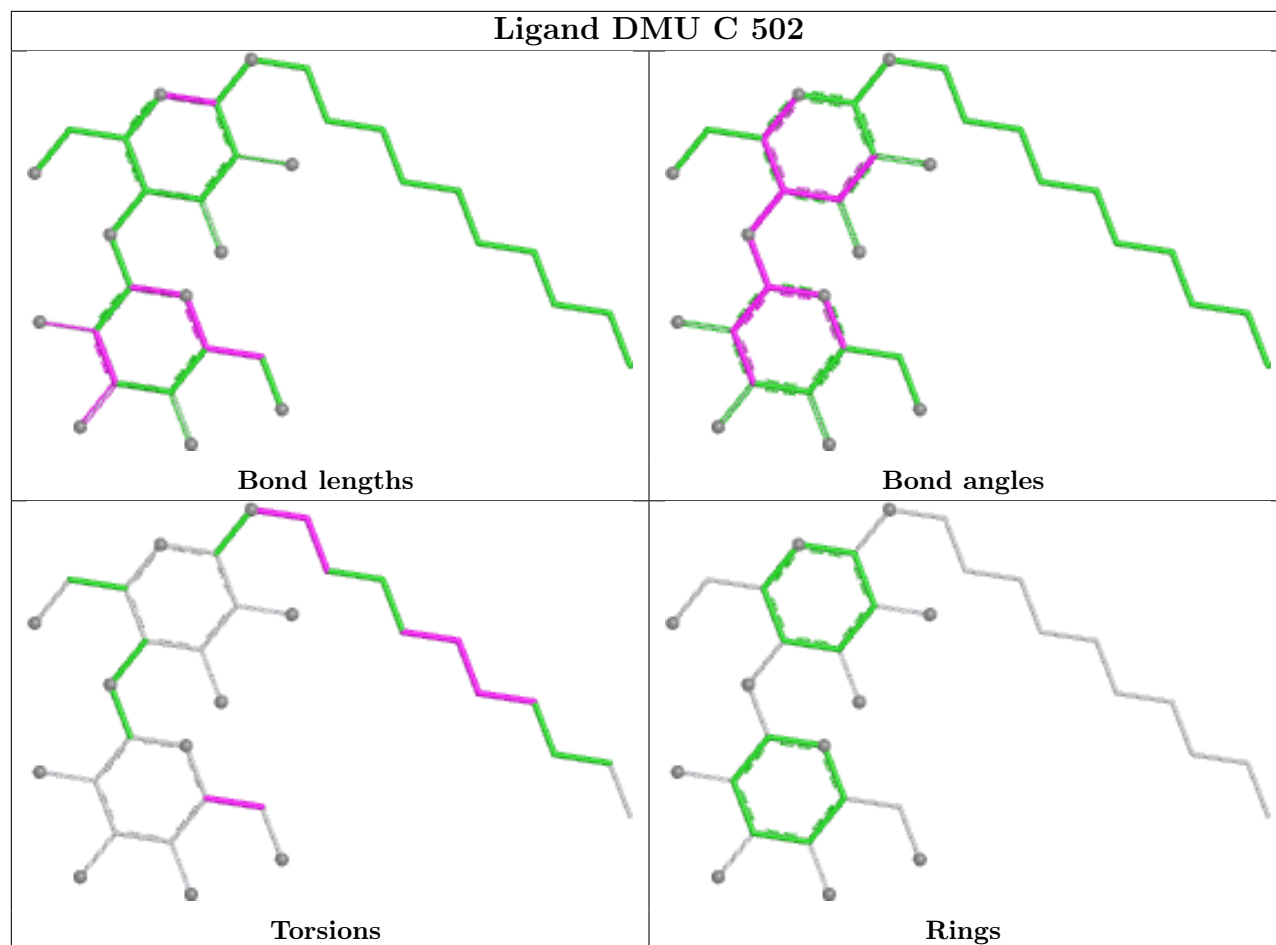
There are no ring outliers.

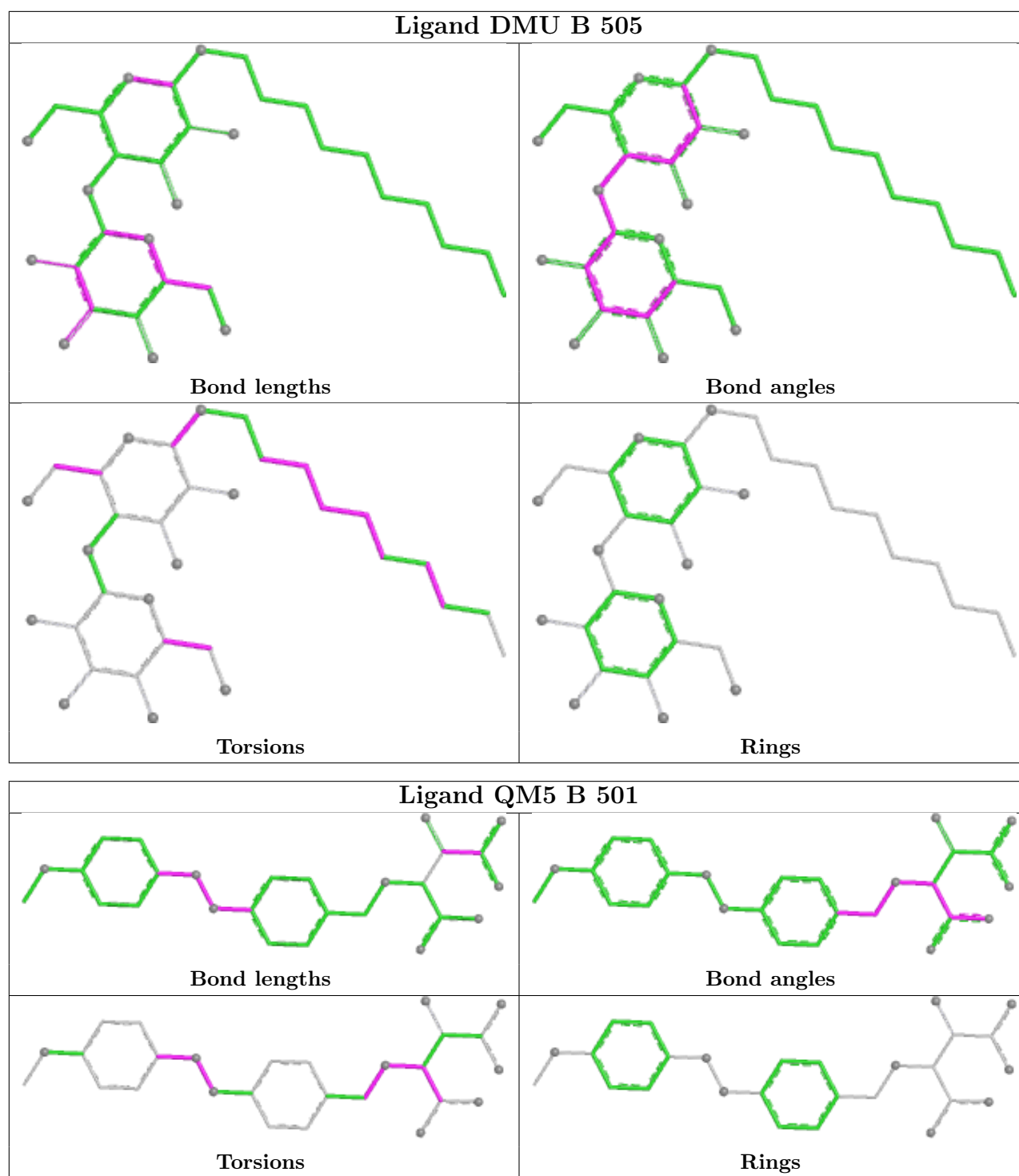
7 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	DMU	1	0
8	C	504	1PE	1	0
4	A	503	PG4	1	0
3	C	502	DMU	1	0
2	C	501	QM5	1	0
3	B	505	DMU	1	0
7	C	511	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 than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers ⓘ

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	424/438 (96%)	0.76	68 (16%) 5 3	125, 161, 225, 386	0
1	B	423/438 (96%)	0.58	47 (11%) 10 6	98, 160, 225, 372	1 (0%)
1	C	423/438 (96%)	0.73	63 (14%) 5 3	91, 175, 226, 328	1 (0%)
All	All	1270/1314 (96%)	0.69	178 (14%) 6 4	91, 165, 227, 386	2 (0%)

All (178) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	350	LEU	14.1
1	C	353	ILE	10.9
1	C	359	PRO	9.4
1	C	354	GLY	9.1
1	B	359	PRO	9.0
1	A	375	LEU	8.9
1	C	205	GLN	8.2
1	B	342	LEU	7.4
1	C	10	LEU	7.3
1	A	377	LEU	6.7
1	A	324	THR	6.6
1	B	346	LEU	6.5
1	A	387	TYR	6.3
1	A	325	VAL	6.2
1	C	121	THR	6.1
1	C	36	GLY	6.1
1	C	35	PHE	6.1
1	A	328	VAL	6.0
1	B	109	PHE	5.9
1	A	370	LEU	5.9
1	C	132	LEU	5.8
1	B	145	PHE	5.5
1	A	145	PHE	5.5

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Mol	Chain	Res	Type	RSRZ
1	C	145	PHE	5.4
1	C	357	GLY	5.1
1	B	146	ALA	5.1
1	C	346	LEU	5.0
1	C	419	THR	4.9
1	B	343	VAL	4.8
1	A	84	GLY	4.7
1	C	9	TYR	4.7
1	C	11	ASP	4.6
1	C	14	VAL	4.6
1	C	382	PRO	4.6
1	C	356	ALA	4.5
1	C	225	ARG	4.5
1	B	167	ILE	4.3
1	B	150	LYS	4.2
1	C	413	THR	4.2
1	C	380	GLY	4.2
1	A	410	LEU	4.1
1	A	323	VAL	4.1
1	C	122	GLY	4.0
1	A	121	THR	3.9
1	B	48	PHE	3.9
1	A	148	LEU	3.9
1	A	376	ASP	3.8
1	A	419	THR	3.8
1	A	391	LEU	3.8
1	B	421	LYS	3.7
1	C	144	PRO	3.7
1	C	384	ALA	3.7
1	C	130	PRO	3.6
1	B	149	ALA	3.6
1	C	148	LEU	3.6
1	A	273	THR	3.6
1	A	406	VAL	3.6
1	B	110	ASN	3.5
1	B	188	VAL	3.5
1	A	409	ASP	3.5
1	C	245	ILE	3.5
1	C	381	SER	3.5
1	A	349	VAL	3.4
1	A	392	GLY	3.4
1	B	35	PHE	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	85	VAL	3.4
1	A	390	ILE	3.4
1	B	269	ASP	3.4
1	A	287	THR	3.3
1	C	349	VAL	3.3
1	C	97	MET	3.3
1	A	374	GLY	3.3
1	B	360	GLY	3.3
1	A	144	PRO	3.2
1	A	389	MET	3.2
1	B	18	ILE	3.2
1	C	272	ILE	3.2
1	C	418	LYS	3.1
1	A	81	GLY	3.1
1	B	339	GLY	3.1
1	C	96	ALA	3.1
1	B	51	LEU	3.1
1	B	235	GLY	3.1
1	A	394	ASP	3.1
1	B	47	PRO	3.1
1	B	232	LYS	3.1
1	A	383	VAL	3.1
1	B	350	LEU	3.0
1	C	119	SER	3.0
1	B	315	ASP	3.0
1	B	125	ILE	3.0
1	C	31	ILE	3.0
1	A	221	GLU	3.0
1	B	144	PRO	3.0
1	B	50	ASP	3.0
1	A	133	VAL	3.0
1	C	120	GLY	3.0
1	A	82	ARG	2.9
1	A	217	TYR	2.9
1	C	117	LEU	2.9
1	A	270	ALA	2.9
1	A	373	VAL	2.9
1	C	355	THR	2.9
1	C	379	PRO	2.9
1	C	118	GLY	2.8
1	A	261	ILE	2.8
1	B	77	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	176	GLU	2.8
1	C	422	GLU	2.8
1	B	358	VAL	2.8
1	A	242	PHE	2.7
1	B	273	THR	2.7
1	B	287	THR	2.7
1	A	431	HIS	2.7
1	A	129	PRO	2.7
1	A	269	ASP	2.7
1	B	409	ASP	2.7
1	C	270	ALA	2.7
1	C	16	TRP	2.7
1	C	420	GLU	2.6
1	C	273	THR	2.6
1	B	16	TRP	2.6
1	A	352	SER	2.6
1	A	235	GLY	2.6
1	B	249	TYR	2.6
1	A	19	LEU	2.6
1	A	306	LEU	2.6
1	A	386	ALA	2.6
1	C	276	VAL	2.5
1	C	351	ALA	2.5
1	C	360	GLY	2.5
1	A	421	LYS	2.5
1	B	276	VAL	2.5
1	A	150	LYS	2.5
1	A	122	GLY	2.5
1	A	424	ASP	2.5
1	A	130	PRO	2.5
1	A	80	LEU	2.4
1	A	304	PHE	2.4
1	B	208	PRO	2.4
1	B	117	LEU	2.4
1	A	225	ARG	2.3
1	C	368	MET	2.3
1	A	146	ALA	2.3
1	C	256	PHE	2.3
1	C	338	LEU	2.3
1	B	248	THR	2.3
1	C	361	ALA	2.3
1	A	34	HIS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	250	PHE	2.3
1	C	277	THR	2.3
1	C	358	VAL	2.3
1	B	185	LEU	2.3
1	C	336	LEU	2.3
1	C	140	VAL	2.3
1	C	269	ASP	2.3
1	A	280	SER	2.2
1	B	267	ALA	2.2
1	B	375	LEU	2.2
1	A	290	VAL	2.2
1	A	283	THR	2.2
1	A	156	VAL	2.2
1	A	9	TYR	2.2
1	A	332	ILE	2.2
1	B	270	ALA	2.1
1	A	369	VAL	2.1
1	C	364	ILE	2.1
1	A	299	LYS	2.1
1	A	272	ILE	2.1
1	C	24	LEU	2.1
1	C	12	TYR	2.1
1	C	392	GLY	2.1
1	C	188	VAL	2.1
1	A	137	LEU	2.1
1	B	356	ALA	2.1
1	A	109	PHE	2.1
1	B	101	PHE	2.0
1	B	227	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

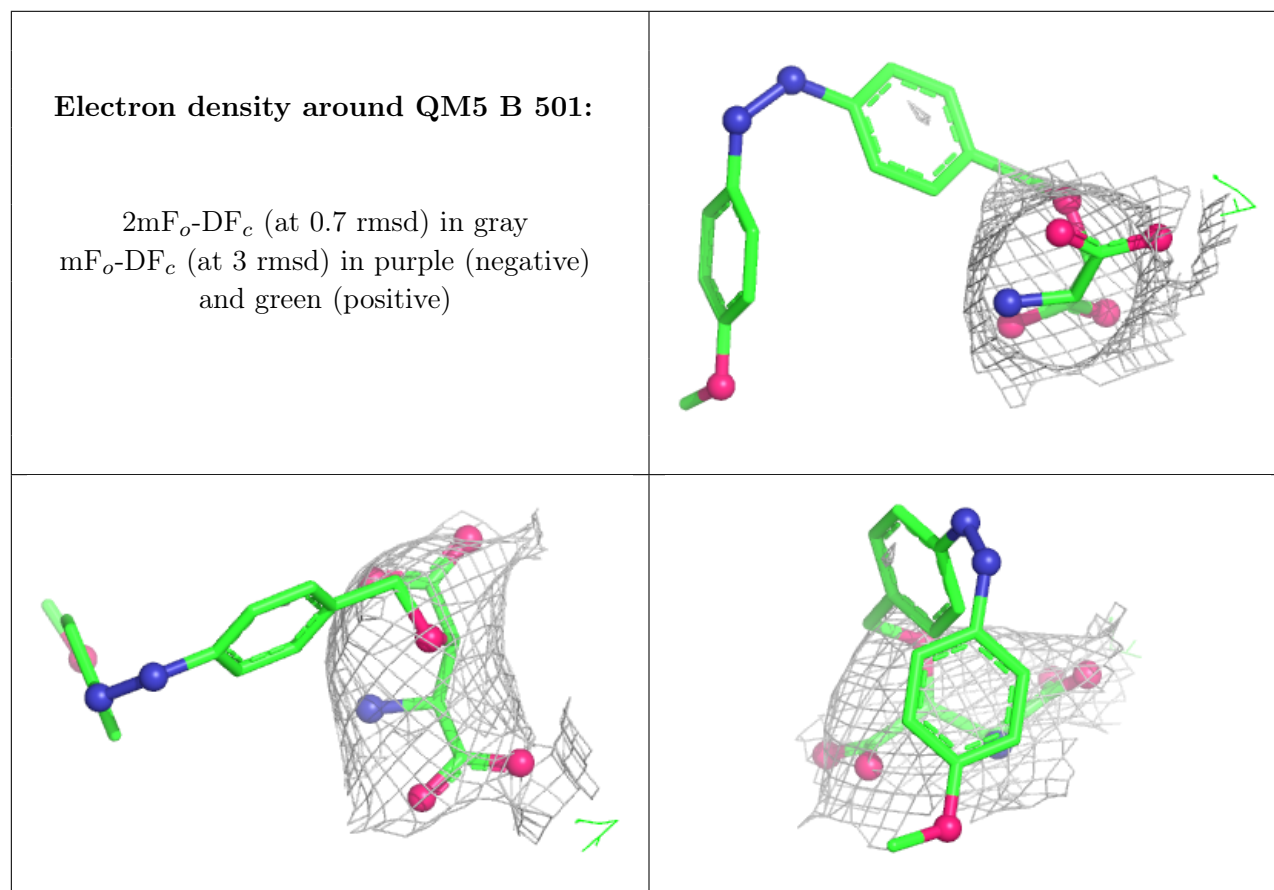
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	NA	C	507	1/1	0.21	0.36	224,224,224,224	0
4	PG4	A	503	13/13	0.45	0.12	193,199,209,210	0
6	NA	B	510	1/1	0.52	0.29	150,150,150,150	0
5	PGE	B	507	10/10	0.53	0.33	165,185,206,209	0
7	PEG	C	513	7/7	0.53	0.13	188,190,194,194	0
7	PEG	C	512	7/7	0.54	0.13	200,209,212,213	0
7	PEG	B	503	7/7	0.54	0.13	154,159,161,161	0
4	PG4	A	504	13/13	0.55	0.29	211,219,224,224	0
5	PGE	B	506	10/10	0.55	0.11	205,222,223,223	0
5	PGE	A	505	10/10	0.56	0.21	189,197,201,202	0
5	PGE	B	502	10/10	0.57	0.12	170,179,184,186	0
4	PG4	A	506	13/13	0.57	0.11	195,198,206,208	0
7	PEG	B	513	7/7	0.58	0.18	186,202,214,217	0
5	PGE	C	503	10/10	0.59	0.22	187,209,223,225	0
6	NA	B	508	1/1	0.61	0.30	147,147,147,147	0
7	PEG	A	512	7/7	0.61	0.24	144,163,184,191	0
7	PEG	C	509	7/7	0.62	0.11	141,143,157,161	0
7	PEG	C	514	7/7	0.63	0.24	192,194,198,199	0
6	NA	A	508	1/1	0.64	0.25	132,132,132,132	0
5	PGE	A	507	10/10	0.67	0.11	192,200,212,215	0
2	QM5	B	501	27/27	0.70	0.19	179,185,230,231	0
7	PEG	C	511	7/7	0.73	0.15	143,153,162,162	0
2	QM5	C	501	27/27	0.76	0.18	183,192,225,229	0
3	DMU	C	502	33/33	0.77	0.17	145,215,232,233	0
7	PEG	B	504	7/7	0.84	0.17	165,173,177,177	0
7	PEG	C	508	7/7	0.87	0.22	182,188,195,196	0
2	QM5	A	501	27/27	0.88	0.11	157,176,230,231	0
7	PEG	B	511	7/7	0.88	0.14	167,174,180,182	0
3	DMU	A	502	33/33	0.88	0.17	182,195,234,236	0
7	PEG	A	511	7/7	0.89	0.36	160,170,175,179	0
6	NA	B	509	1/1	0.89	0.18	169,169,169,169	0
6	NA	A	509	1/1	0.89	0.13	164,164,164,164	0
3	DMU	B	505	33/33	0.89	0.14	164,196,221,225	0
6	NA	C	505	1/1	0.90	0.14	138,138,138,138	0
6	NA	A	510	1/1	0.90	0.22	140,140,140,140	0
8	1PE	C	504	16/16	0.93	0.20	171,181,193,193	0
6	NA	C	506	1/1	0.94	0.09	191,191,191,191	0

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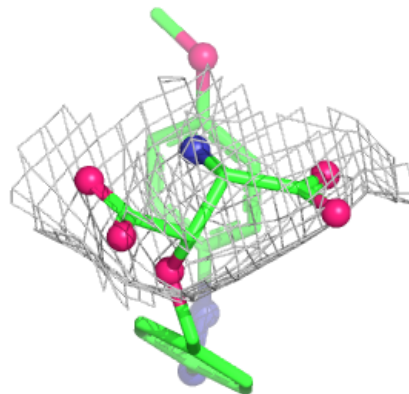
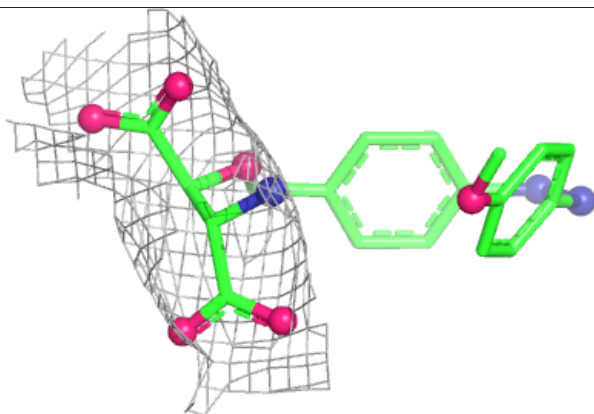
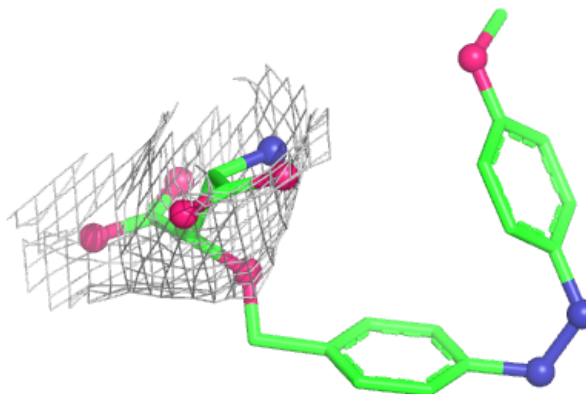
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	PEG	B	512	7/7	0.94	0.27	172,181,185,186	0
7	PEG	C	510	7/7	0.96	0.24	180,194,201,203	0

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.

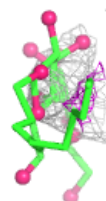
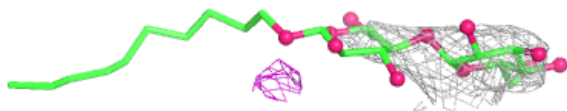
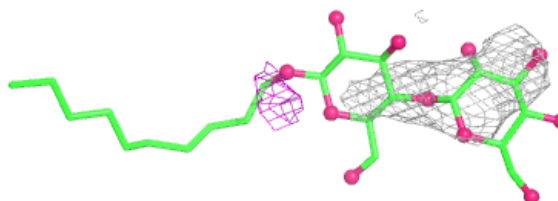


Electron density around QM5 C 501:

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

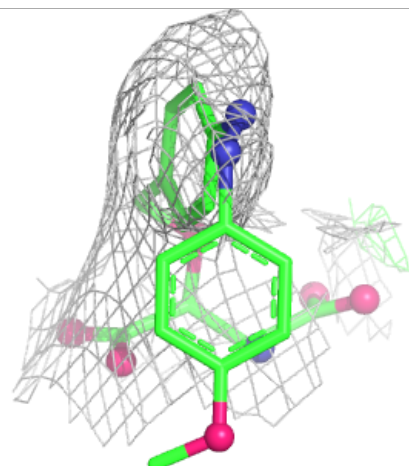
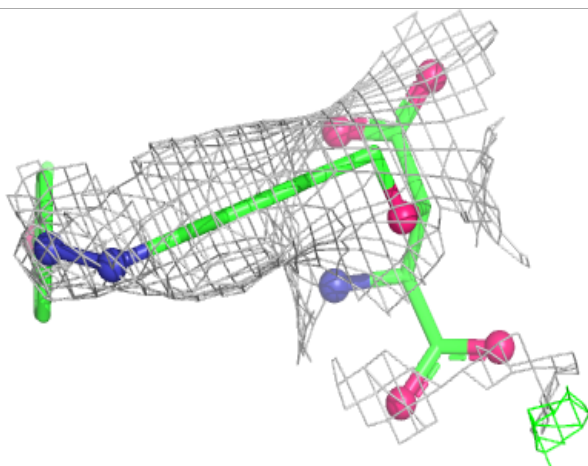
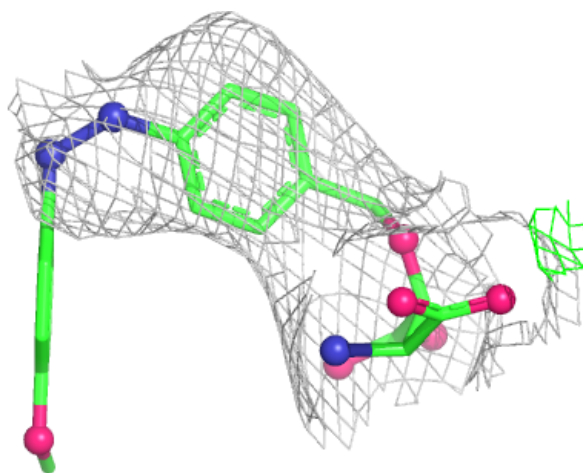
**Electron density around DMU C 502:**

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



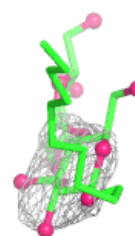
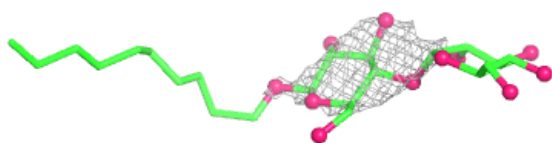
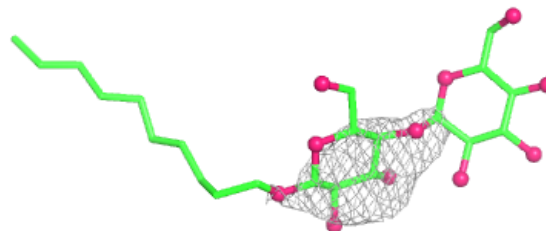
Electron density around QM5 A 501:

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

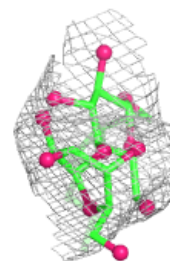
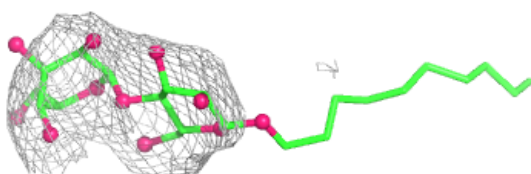
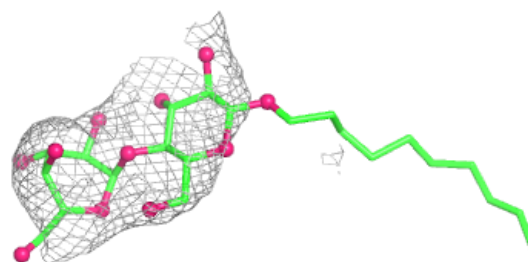


Electron density around DMU A 502:

$2mF_o-DF_c$ (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)



6.5 Other polymers [i](#)

There are no such residues in this entry.