



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:08 PM UTC

PDB ID : 6JGG / pdb_00006jgg
Title : Crystal structure of barley exohydrolaseI W434F mutant in complex with methyl 2-thio-beta-sophoroside.
Authors : Luang, S.; Streltsov, V.A.; Hrmova, M.
Deposited on : 2019-02-14
Resolution : 1.90 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

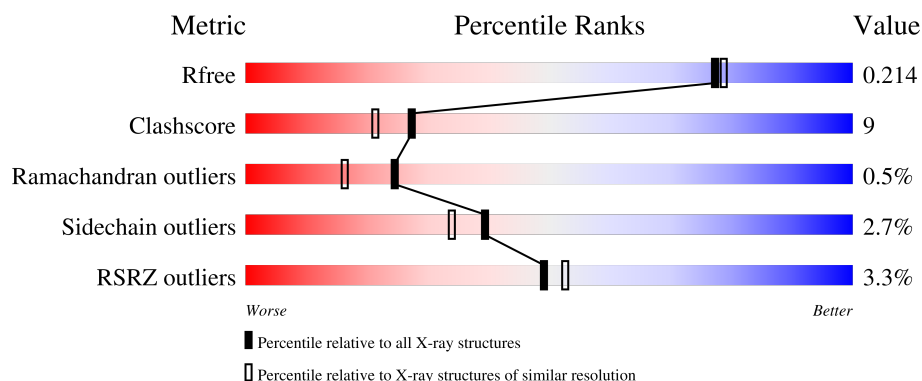
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	609	3% 86% 13%
2	B	2	50% 50%

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
5	ACT	A	710	-	-	X	-
5	ACT	A	716	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5359 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 BETA-D-GLUCAN GLUCOHYDROLASE ISOENZYME EXO1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	606	Total	C	N	O	S	20	10	0
			4615	2917	801	870	27			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	HIS	-	expression tag	UNP A0A287SCR5
A	-2	HIS	-	expression tag	UNP A0A287SCR5
A	-1	ALA	-	expression tag	UNP A0A287SCR5
A	0	ALA	-	expression tag	UNP A0A287SCR5
A	320	LYS	ASN	See sequence details	UNP A0A287SCR5
A	434	PHE	TRP	engineered mutation	UNP A0A287SCR5

- Molecule 2 is an oligosaccharide called beta-D-glucopyranose-(1-2)-methyl 2-thio-beta-D-glucopyranoside.

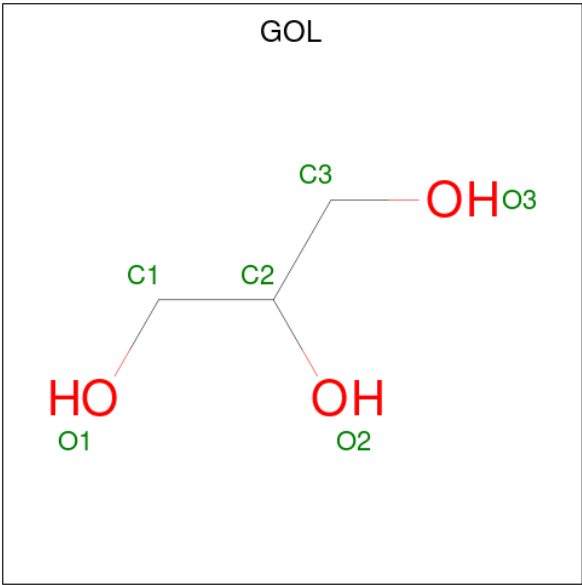
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	2	Total	C	O	S	0	1	0
			25	14	10	1			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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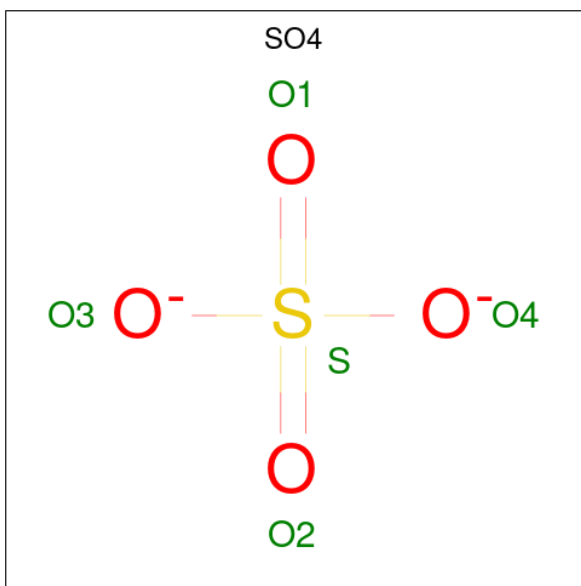
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



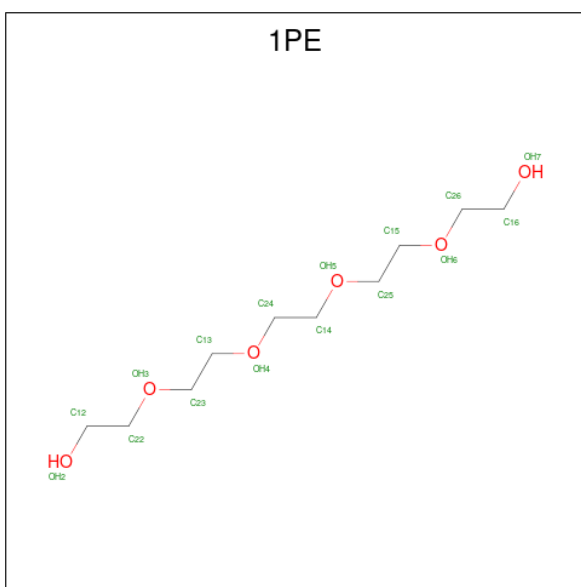
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 8 5 3	0	0
7	A	1	Total C O 6 4 2	0	0
7	A	1	Total C O 6 4 2	0	0
7	A	1	Total C O 9 6 3	0	0
7	A	1	Total C O 9 6 3	0	0
7	A	1	Total C O 4 2 2	0	0
7	A	1	Total C O 4 2 2	0	0
7	A	1	Total C O 4 2 2	0	0
7	A	1	Total C O 5 3 2	0	0
7	A	1	Total C O 6 4 2	0	0
7	A	1	Total C O 4 2 2	0	0
7	A	1	Total C O 5 3 2	0	0
7	A	1	Total C O 6 4 2	0	0
7	A	1	Total C O 6 4 2	0	0
7	A	1	Total C O 9 5 4	0	1
7	A	1	Total C O 16 10 6	0	0
7	A	1	Total C O 12 6 6	0	1

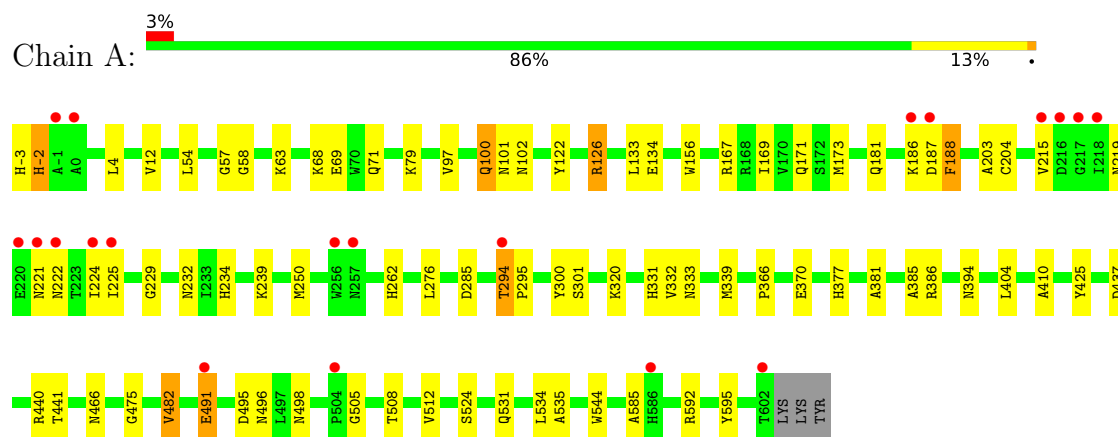
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	505	Total O 505 505	0	0

3 Residue-property plots [i](#)

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

- Molecule 1: BETA-D-GLUCAN GLUCOHYDROLASE ISOENZYME EXO1



- Molecule 2: beta-D-glucopyranose-(1-2)-methyl 2-thio-beta-D-glucopyranoside



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	100.72Å 100.72Å 180.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	87.99 – 1.90 87.99 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.5 (87.99-1.90) 98.6 (87.99-1.90)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.81 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.161 , 0.202 0.176 , 0.214	Depositor DCC
R_{free} test set	3669 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	27.6	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5359	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 1PE, SO4, NAG, U2A, BGC, GOL, ACT

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	1.23	2/4760 (0.0%)	1.15	11/6462 (0.2%)

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	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	-2	HIS	C-N	10.26	1.48	1.33
1	A	69	GLU	N-CA	5.95	1.53	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	-2	HIS	O-C-N	-8.89	110.77	122.59
1	A	126	ARG	NE-CZ-NH2	-6.15	113.67	119.20
1	A	475	GLY	N-CA-C	6.14	121.59	114.16
1	A	126	ARG	NE-CZ-NH1	6.11	127.61	121.50
1	A	188	PHE	N-CA-C	5.94	123.46	110.80
1	A	495	ASP	N-CA-C	-5.88	101.34	110.10
1	A	385	ALA	N-CA-C	-5.86	104.80	111.07
1	A	410	ALA	CA-C-N	-5.58	112.76	118.85
1	A	410	ALA	C-N-CA	-5.58	112.76	118.85
1	A	381	ALA	N-CA-C	-5.45	105.42	111.36
1	A	300	TYR	N-CA-C	-5.24	105.74	111.82

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	-3	HIS	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	4615	0	4577	68	0
2	B	25	0	10	4	0
3	A	14	0	13	3	0
4	A	30	0	40	4	0
5	A	36	0	27	9	0
6	A	15	0	0	0	0
7	A	119	0	141	22	0
8	A	505	0	0	23	0
All	All	5359	0	4808	88	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 9.

All (88) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:ASN:HD21	3:A:701:NAG:C1	1.11	1.62
1:A:221:ASN:ND2	3:A:701:NAG:C1	1.85	1.34
1:A:466[A]:ASN:ND2	1:A:508:THR:OG1	1.82	1.12
1:A:394:ASN:HD21	1:A:404:LEU:H	1.20	0.86
7:A:720:1PE:H222	8:A:811:HOH:O	1.78	0.84
1:A:156:TRP:HE1	1:A:219:ASN:HD22	1.27	0.81
7:A:723:1PE:C14	8:A:1253:HOH:O	2.29	0.80
1:A:100:GLN:HE21	1:A:100:GLN:HA	1.49	0.78
1:A:167:ARG:HH21	1:A:171:GLN:HE22	1.30	0.78
7:A:736[A]:1PE:H222	8:A:839:HOH:O	1.84	0.77
1:A:592:ARG:NH2	5:A:716:ACT:H2	2.01	0.76
1:A:221:ASN:ND2	3:A:701:NAG:O5	2.20	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:736[B]:1PE:H222	8:A:839:HOH:O	1.93	0.69
1:A:122:TYR:CE2	1:A:126:ARG:HD2	2.28	0.68
4:A:707:GOL:O2	8:A:803:HOH:O	2.11	0.68
1:A:181:GLN:HE21	1:A:203:ALA:H	1.38	0.68
5:A:710:ACT:H3	8:A:1217:HOH:O	1.91	0.68
1:A:100:GLN:HE21	1:A:100:GLN:CA	2.07	0.68
1:A:377:HIS:HE1	5:A:710:ACT:CH3	2.07	0.67
1:A:592:ARG:H	5:A:716:ACT:H3	1.61	0.66
1:A:97:VAL:H	1:A:101:ASN:HD21	1.42	0.66
1:A:491:GLU:HB3	8:A:999:HOH:O	1.96	0.65
1:A:167:ARG:HH21	1:A:171:GLN:NE2	1.93	0.64
1:A:134:GLU:OE2	1:A:377:HIS:HD2	1.80	0.64
1:A:496:ASN:HD22	7:A:734[B]:1PE:H131	1.63	0.62
1:A:97:VAL:H	1:A:101:ASN:ND2	1.95	0.62
1:A:425:TYR:OH	7:A:725:1PE:H121	1.99	0.61
1:A:333:ASN:HD21	4:A:705:GOL:C3	2.14	0.61
1:A:57:GLY:H	2:B:1[A]:U2A:C7	2.14	0.60
1:A:262:HIS:HE1	1:A:285:ASP:H	1.49	0.60
1:A:68:LYS:HE3	8:A:827:HOH:O	2.01	0.59
1:A:239:LYS:HD3	1:A:276:LEU:HD22	1.84	0.59
7:A:735:1PE:H222	8:A:1144:HOH:O	2.02	0.58
7:A:732:1PE:H232	8:A:834:HOH:O	2.05	0.57
1:A:57:GLY:H	2:B:1[B]:U2A:C7	2.17	0.57
1:A:58:GLY:H	1:A:102:ASN:ND2	2.02	0.56
1:A:12:VAL:HB	7:A:733:1PE:H121	1.87	0.56
7:A:734[A]:1PE:H132	8:A:1260:HOH:O	2.06	0.55
1:A:134:GLU:OE2	1:A:377:HIS:CD2	2.61	0.54
1:A:156:TRP:HE1	1:A:219:ASN:ND2	2.00	0.54
1:A:386:ARG:HH22	4:A:704:GOL:H11	1.73	0.54
1:A:370[A]:GLU:HB3	8:A:802:HOH:O	2.07	0.53
1:A:167:ARG:NH2	1:A:171:GLN:HE22	2.04	0.53
1:A:301:SER:O	1:A:331:HIS:HE1	1.92	0.52
1:A:524:SER:O	1:A:544:TRP:HA	2.10	0.52
1:A:437:ASP:HB3	1:A:441:THR:HG21	1.91	0.52
1:A:592:ARG:HH21	5:A:716:ACT:H2	1.70	0.51
1:A:181:GLN:NE2	1:A:203:ALA:H	2.07	0.51
1:A:262:HIS:CE1	1:A:285:ASP:H	2.28	0.50
1:A:496:ASN:ND2	7:A:734[B]:1PE:H131	2.25	0.49
7:A:736[B]:1PE:H221	8:A:1089:HOH:O	2.12	0.49
1:A:535:ALA:HB2	7:A:735:1PE:H242	1.95	0.48
1:A:534:LEU:HD13	7:A:723:1PE:H131	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:736[A]:1PE:H221	8:A:1089:HOH:O	2.14	0.48
1:A:234:HIS:HE1	8:A:1231:HOH:O	1.97	0.47
1:A:234:HIS:HD2	8:A:1118:HOH:O	1.96	0.47
1:A:100:GLN:HA	1:A:100:GLN:NE2	2.25	0.47
1:A:531:GLN:HG2	7:A:735:1PE:H222	1.98	0.46
1:A:332:VAL:HG11	1:A:339:MET:HE1	1.97	0.46
1:A:585:ALA:HB3	7:A:727:1PE:H221	1.98	0.45
1:A:219:ASN:HD21	4:A:703:GOL:C1	2.28	0.45
1:A:221:ASN:CG	1:A:222:ASN:H	2.24	0.45
1:A:595:TYR:CD2	5:A:716:ACT:H1	2.51	0.45
1:A:122:TYR:CZ	1:A:126:ARG:HD2	2.51	0.45
1:A:482[B]:VAL:HG22	1:A:512:VAL:HG11	1.99	0.45
1:A:482[B]:VAL:CG2	1:A:512:VAL:HG11	2.47	0.44
1:A:133:LEU:HD13	5:A:710:ACT:H1	1.99	0.44
1:A:225:ILE:HD11	1:A:229:GLY:HA3	1.99	0.44
1:A:215:VAL:HG21	1:A:224:ILE:HD12	2.00	0.43
7:A:720:1PE:H231	8:A:1198:HOH:O	2.18	0.43
1:A:215:VAL:CG2	1:A:224:ILE:HD12	2.48	0.43
1:A:531:GLN:CB	7:A:723:1PE:H242	2.49	0.43
8:A:823:HOH:O	2:B:1[A]:U2A:C7	2.68	0.42
1:A:169:ILE:O	1:A:173[A]:MET:HG2	2.20	0.41
1:A:71:GLN:HE22	1:A:366:PRO:HA	1.85	0.41
1:A:377:HIS:HE1	5:A:710:ACT:H1	1.83	0.41
7:A:735:1PE:C16	8:A:1193:HOH:O	2.66	0.41
8:A:823:HOH:O	2:B:1[B]:U2A:C7	2.69	0.41
1:A:294:THR:HA	1:A:295:PRO:C	2.45	0.41
1:A:100:GLN:CA	1:A:100:GLN:NE2	2.80	0.41
1:A:204:CYS:SG	1:A:250:MET:HB2	2.61	0.41
1:A:4:LEU:C	1:A:4:LEU:HD23	2.46	0.41
1:A:133:LEU:CD1	5:A:710:ACT:H1	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	612/609 (100%)	592 (97%)	17 (3%)	3 (0%)	24	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	-2	HIS
1	A	188	PHE
1	A	505	GLY

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	495/490 (101%)	480 (97%)	15 (3%)	36	30

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	LEU
1	A	63	LYS
1	A	79	LYS
1	A	100	GLN
1	A	186	LYS
1	A	187	ASP
1	A	232[A]	ASN
1	A	232[B]	ASN
1	A	294	THR
1	A	320	LYS
1	A	440[A]	ARG
1	A	482[A]	VAL
1	A	482[B]	VAL
1	A	491	GLU
1	A	498	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	100	GLN
1	A	101	ASN
1	A	102	ASN
1	A	171	GLN
1	A	181	GLN
1	A	219	ASN
1	A	221	ASN
1	A	234	HIS
1	A	262	HIS
1	A	331	HIS
1	A	333	ASN
1	A	377	HIS
1	A	394	ASN
1	A	487	HIS
1	A	581	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	A	701	-	14,14,15	0.51	0	17,19,21	1.39	3 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	701	-	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	NAG	C1-C2-N2	2.89	114.98	110.43
3	A	701	NAG	C2-N2-C7	2.78	126.63	122.90
3	A	701	NAG	C3-C4-C5	2.05	113.95	110.23

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	701	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	701	NAG	3	0

5.5 Carbohydrates [i](#)

3 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	U2A	B	1[A]	-	12,13,13	1.99	2 (16%)	14,18,18	3.24	5 (35%)
2	U2A	B	1[B]	-	12,13,13	1.99	2 (16%)	14,18,18	3.13	5 (35%)
2	BGC	B	2	2	11,11,12	0.78	0	15,15,17	1.32	3 (20%)

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
2	U2A	B	1[A]	-	-	4/4/24/24	0/1/1/1
2	U2A	B	1[B]	-	-	4/4/24/24	0/1/1/1
2	BGC	B	2	2	-	0/2/19/22	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1[A]	U2A	O1-C1	4.66	1.48	1.40
2	B	1[B]	U2A	O1-C1	4.66	1.48	1.40
2	B	1[A]	U2A	C1-C2	3.83	1.59	1.52
2	B	1[B]	U2A	C1-C2	3.83	1.59	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1[A]	U2A	O1-C1-C2	7.73	119.83	108.10
2	B	1[B]	U2A	O1-C1-C2	7.73	119.83	108.10
2	B	1[A]	U2A	O5-C1-O1	-5.79	97.46	110.94
2	B	1[B]	U2A	O5-C1-O1	-5.79	97.46	110.94
2	B	1[A]	U2A	C7-O1-C1	4.05	119.42	113.26
2	B	1[A]	U2A	O4-C4-C3	-3.56	101.99	110.38
2	B	1[B]	U2A	O4-C4-C3	-3.56	101.99	110.38
2	B	1[A]	U2A	O3-C3-C2	3.31	115.93	109.17
2	B	1[B]	U2A	O3-C3-C2	3.31	115.93	109.17
2	B	2	BGC	O5-C5-C4	-2.81	103.99	110.83
2	B	2	BGC	C1-C2-C3	2.71	113.59	109.64
2	B	2	BGC	C1-O5-C5	2.66	115.75	112.19
2	B	1[B]	U2A	C7-O1-C1	2.50	117.06	113.26

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1[A]	U2A	C2-C1-O1-C7
2	B	1[B]	U2A	C2-C1-O1-C7
2	B	1[A]	U2A	O5-C1-O1-C7
2	B	1[B]	U2A	O5-C1-O1-C7

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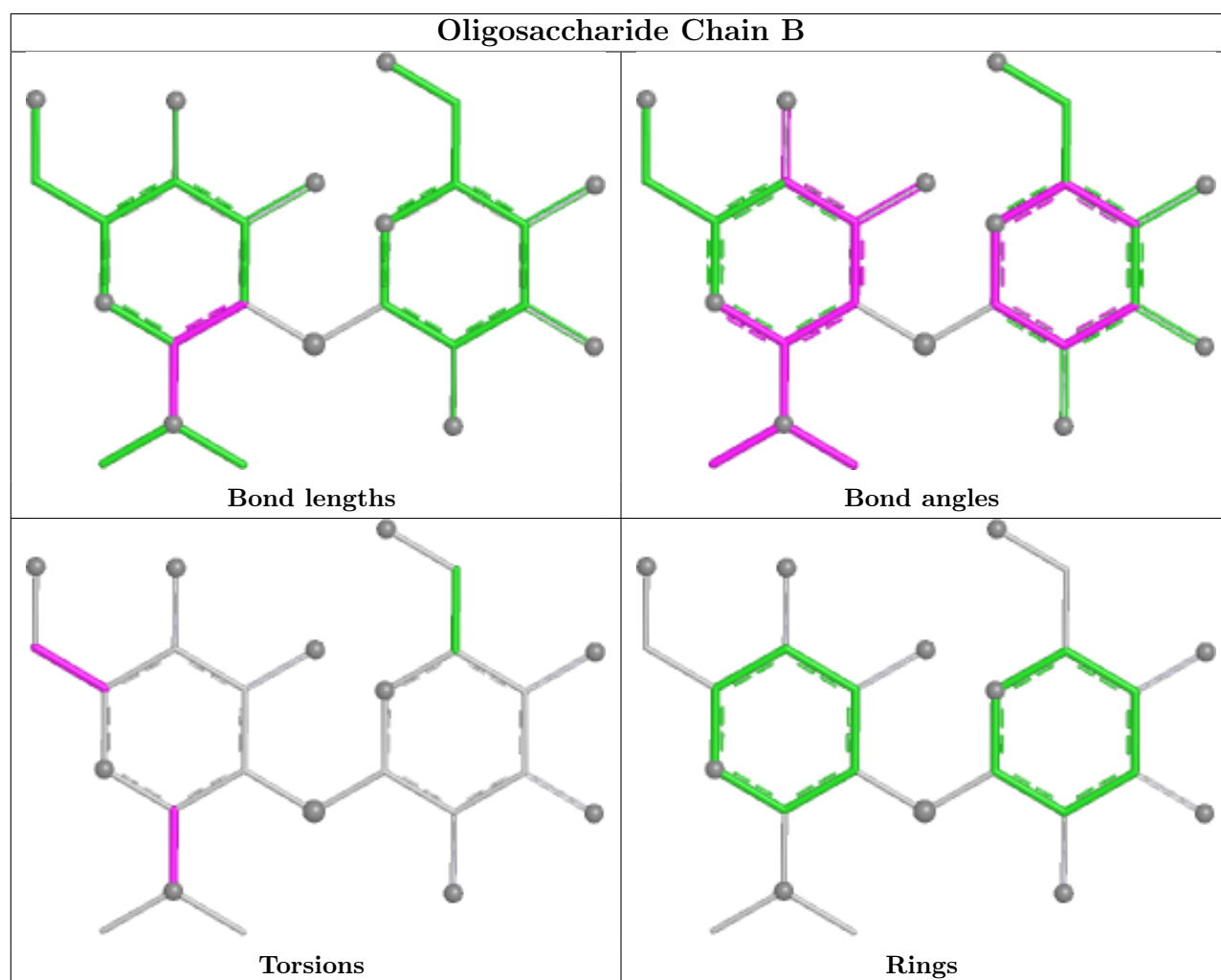
Mol	Chain	Res	Type	Atoms
2	B	1[A]	U2A	O5-C5-C6-O6
2	B	1[B]	U2A	O5-C5-C6-O6
2	B	1[A]	U2A	C4-C5-C6-O6
2	B	1[B]	U2A	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1[B]	U2A	2	0
2	B	1[A]	U2A	2	0

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



5.6 Ligand geometry

37 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	GOL	A	705	-	5,5,5	0.52	0	5,5,5	1.00	0
4	GOL	A	704	-	5,5,5	0.70	0	5,5,5	0.47	0
7	1PE	A	723	-	8,8,15	0.84	0	7,7,14	1.18	0
4	GOL	A	706	-	5,5,5	0.61	0	5,5,5	0.98	0
7	1PE	A	733	-	5,5,15	0.59	0	4,4,14	1.07	0
7	1PE	A	728	-	4,4,15	0.58	0	3,3,14	0.17	0
5	ACT	A	710	-	3,3,3	0.90	0	3,3,3	0.80	0
5	ACT	A	712	-	3,3,3	1.16	0	3,3,3	0.56	0
6	SO4	A	718	-	4,4,4	0.53	0	6,6,6	1.25	1 (16%)
7	1PE	A	721	-	5,5,15	0.76	0	4,4,14	0.87	0
7	1PE	A	736[A]	-	6,6,15	0.92	0	5,5,14	1.00	0
7	1PE	A	732	-	5,5,15	1.44	0	4,4,14	2.17	1 (25%)
5	ACT	A	711	-	3,3,3	1.14	0	3,3,3	0.99	0
7	1PE	A	722	-	5,5,15	0.78	0	4,4,14	0.83	0
5	ACT	A	714	-	3,3,3	1.00	0	3,3,3	0.62	0
7	1PE	A	729	-	5,5,15	0.67	0	4,4,14	0.72	0
7	1PE	A	720	-	7,7,15	0.45	0	6,6,14	1.03	0
7	1PE	A	730	-	3,3,15	0.32	0	2,2,14	0.97	0
7	1PE	A	731	-	4,4,15	0.76	0	3,3,14	0.85	0
7	1PE	A	726	-	3,3,15	0.66	0	2,2,14	0.70	0
7	1PE	A	736[B]	-	6,6,15	0.93	0	5,5,14	0.74	0
6	SO4	A	719	-	4,4,4	0.58	0	6,6,6	0.83	0
7	1PE	A	735	-	15,15,15	0.69	0	14,14,14	0.58	0
7	1PE	A	725	-	3,3,15	0.26	0	2,2,14	0.54	0
5	ACT	A	709	-	3,3,3	1.19	0	3,3,3	0.66	0
5	ACT	A	708	-	3,3,3	0.97	0	3,3,3	0.14	0
4	GOL	A	707	-	5,5,5	0.65	0	5,5,5	0.71	0
3	NAG	A	701	-	14,14,15	0.51	0	17,19,21	1.39	3 (17%)
5	ACT	A	716	-	3,3,3	0.64	0	3,3,3	1.24	0
4	GOL	A	703	-	5,5,5	0.74	0	5,5,5	2.43	3 (60%)
7	1PE	A	734[A]	-	6,6,15	0.67	0	5,5,14	2.16	2 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	ACT	A	715	-	3,3,3	0.91	0	3,3,3	0.32	0
5	ACT	A	713	-	3,3,3	0.98	0	3,3,3	0.40	0
7	1PE	A	734[B]	-	6,6,15	0.59	0	5,5,14	1.31	1 (20%)
7	1PE	A	727	-	3,3,15	0.37	0	2,2,14	0.72	0
7	1PE	A	724	-	8,8,15	0.91	0	7,7,14	0.95	0
6	SO4	A	717	-	4,4,4	1.01	0	6,6,6	1.73	1 (16%)

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
4	GOL	A	705	-	-	2/4/4/4	-
4	GOL	A	704	-	-	2/4/4/4	-
7	1PE	A	723	-	-	6/6/6/13	-
4	GOL	A	706	-	-	3/4/4/4	-
7	1PE	A	733	-	-	2/3/3/13	-
7	1PE	A	728	-	-	0/2/2/13	-
7	1PE	A	721	-	-	2/3/3/13	-
7	1PE	A	736[A]	-	-	1/4/4/13	-
7	1PE	A	732	-	-	1/3/3/13	-
7	1PE	A	722	-	-	1/3/3/13	-
7	1PE	A	729	-	-	3/3/3/13	-
7	1PE	A	720	-	-	3/5/5/13	-
7	1PE	A	730	-	-	1/1/1/13	-
7	1PE	A	731	-	-	1/2/2/13	-
7	1PE	A	726	-	-	1/1/1/13	-
7	1PE	A	736[B]	-	-	4/4/4/13	-
7	1PE	A	735	-	-	7/13/13/13	-
7	1PE	A	725	-	-	0/1/1/13	-
4	GOL	A	707	-	-	0/4/4/4	-
3	NAG	A	701	-	-	1/6/23/26	0/1/1/1
4	GOL	A	703	-	-	0/4/4/4	-
7	1PE	A	734[A]	-	-	2/4/4/13	-
7	1PE	A	734[B]	-	-	3/4/4/13	-
7	1PE	A	727	-	-	1/1/1/13	-
7	1PE	A	724	-	-	4/6/6/13	-

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	734[A]	1PE	OH3-C23-C13	-3.61	94.21	110.11
7	A	732	1PE	C22-OH3-C23	3.60	125.70	113.06
6	A	717	SO4	O4-S-O1	-3.30	92.28	109.56
4	A	703	GOL	O2-C2-C3	3.20	122.44	109.18
4	A	703	GOL	C3-C2-C1	-3.04	100.63	111.80
7	A	734[A]	1PE	OH4-C13-C23	-3.01	94.10	111.82
3	A	701	NAG	C1-C2-N2	2.89	114.98	110.43
3	A	701	NAG	C2-N2-C7	2.78	126.63	122.90
6	A	718	SO4	O4-S-O2	2.60	123.14	109.56
7	A	734[B]	1PE	OH3-C23-C13	-2.47	99.23	110.11
4	A	703	GOL	O3-C3-C2	2.24	120.46	110.38
3	A	701	NAG	C3-C4-C5	2.05	113.95	110.23

There are no chirality outliers.

All (51) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	704	GOL	O1-C1-C2-C3
4	A	705	GOL	C1-C2-C3-O3
4	A	705	GOL	O2-C2-C3-O3
4	A	706	GOL	O1-C1-C2-C3
7	A	723	1PE	C13-C23-OH3-C22
7	A	724	1PE	C16-C26-OH6-C15
7	A	723	1PE	OH4-C13-C23-OH3
7	A	720	1PE	OH4-C13-C23-OH3
7	A	729	1PE	OH2-C12-C22-OH3
7	A	736[A]	1PE	OH2-C12-C22-OH3
7	A	735	1PE	OH4-C13-C23-OH3
7	A	720	1PE	OH2-C12-C22-OH3
7	A	733	1PE	OH2-C12-C22-OH3
7	A	734[A]	1PE	OH2-C12-C22-OH3
7	A	734[B]	1PE	OH2-C12-C22-OH3
7	A	735	1PE	C24-C14-OH5-C25
7	A	723	1PE	OH2-C12-C22-OH3
7	A	736[B]	1PE	OH2-C12-C22-OH3
4	A	704	GOL	O1-C1-C2-O2
7	A	733	1PE	C13-C23-OH3-C22
7	A	732	1PE	OH4-C13-C23-OH3
7	A	721	1PE	OH4-C13-C23-OH3
7	A	734[B]	1PE	OH4-C13-C23-OH3

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Mol	Chain	Res	Type	Atoms
7	A	726	1PE	OH7-C16-C26-OH6
7	A	721	1PE	C12-C22-OH3-C23
7	A	720	1PE	C13-C23-OH3-C22
7	A	724	1PE	OH6-C15-C25-OH5
7	A	729	1PE	C13-C23-OH3-C22
7	A	722	1PE	C12-C22-OH3-C23
3	A	701	NAG	C4-C5-C6-O6
7	A	735	1PE	C12-C22-OH3-C23
7	A	724	1PE	C15-C25-OH5-C14
7	A	736[B]	1PE	C13-C23-OH3-C22
7	A	736[B]	1PE	OH4-C13-C23-OH3
7	A	736[B]	1PE	C12-C22-OH3-C23
7	A	727	1PE	OH2-C12-C22-OH3
7	A	730	1PE	OH2-C12-C22-OH3
7	A	734[B]	1PE	C13-C23-OH3-C22
7	A	723	1PE	C23-C13-OH4-C24
7	A	735	1PE	OH5-C14-C24-OH4
7	A	723	1PE	C12-C22-OH3-C23
7	A	734[A]	1PE	OH4-C13-C23-OH3
7	A	735	1PE	OH2-C12-C22-OH3
4	A	706	GOL	C1-C2-C3-O3
7	A	735	1PE	C25-C15-OH6-C26
4	A	706	GOL	O1-C1-C2-O2
7	A	724	1PE	C24-C14-OH5-C25
7	A	723	1PE	C14-C24-OH4-C13
7	A	731	1PE	C12-C22-OH3-C23
7	A	729	1PE	C12-C22-OH3-C23
7	A	735	1PE	C23-C13-OH4-C24

There are no ring outliers.

18 monomers are involved in 38 short contacts:

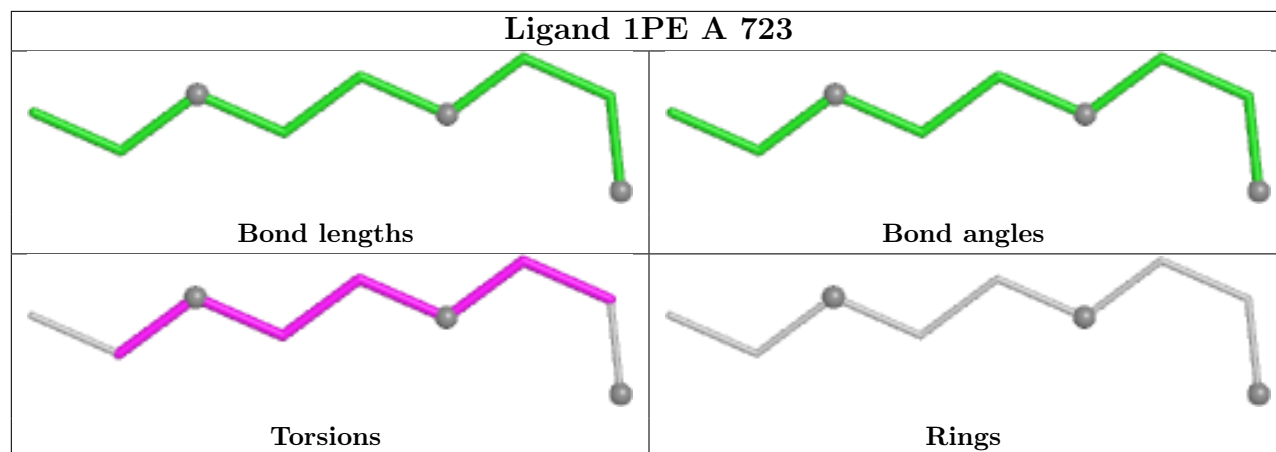
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	705	GOL	1	0
4	A	704	GOL	1	0
7	A	723	1PE	3	0
7	A	733	1PE	1	0
5	A	710	ACT	5	0
7	A	736[A]	1PE	2	0
7	A	732	1PE	1	0
7	A	720	1PE	2	0
7	A	736[B]	1PE	2	0

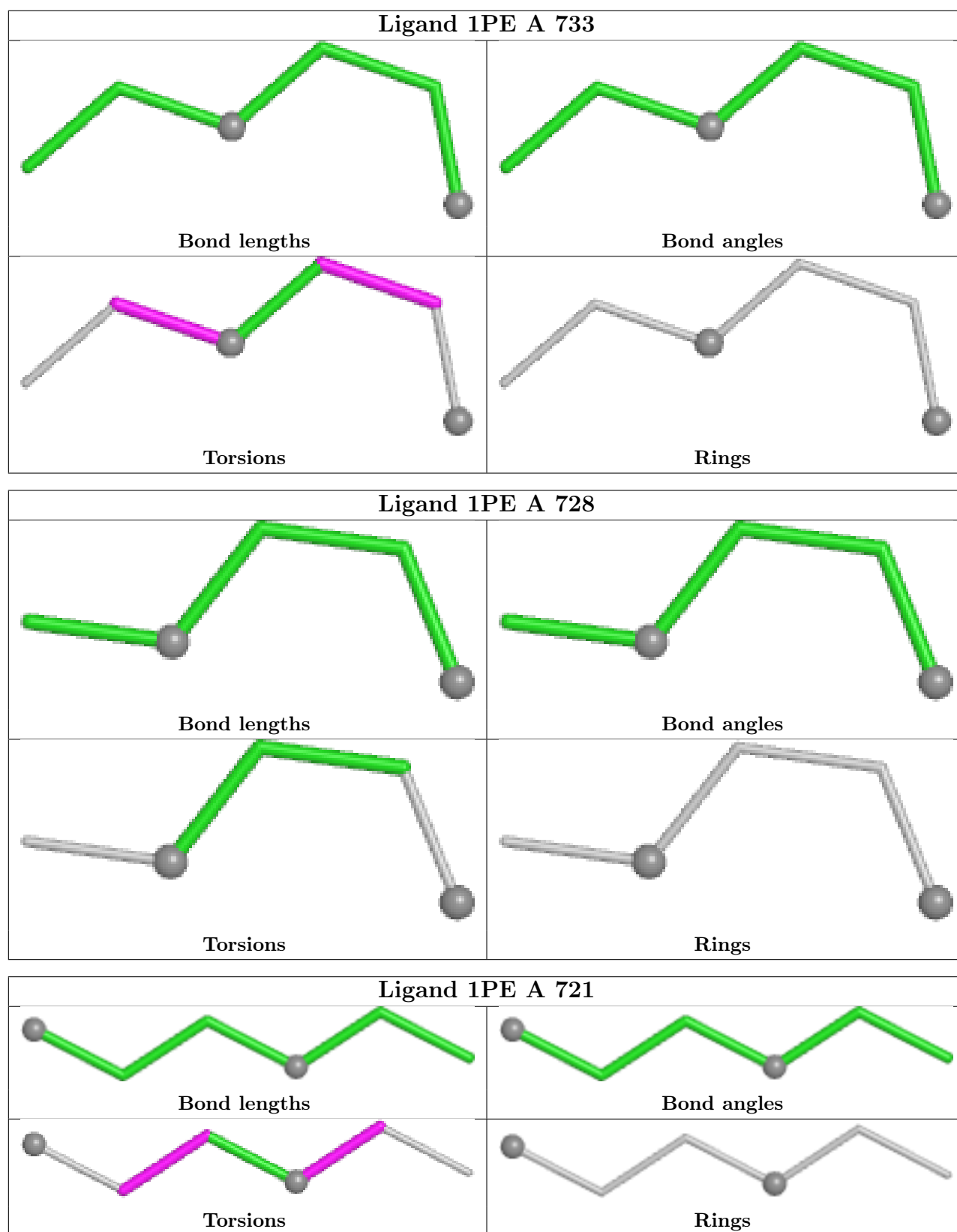
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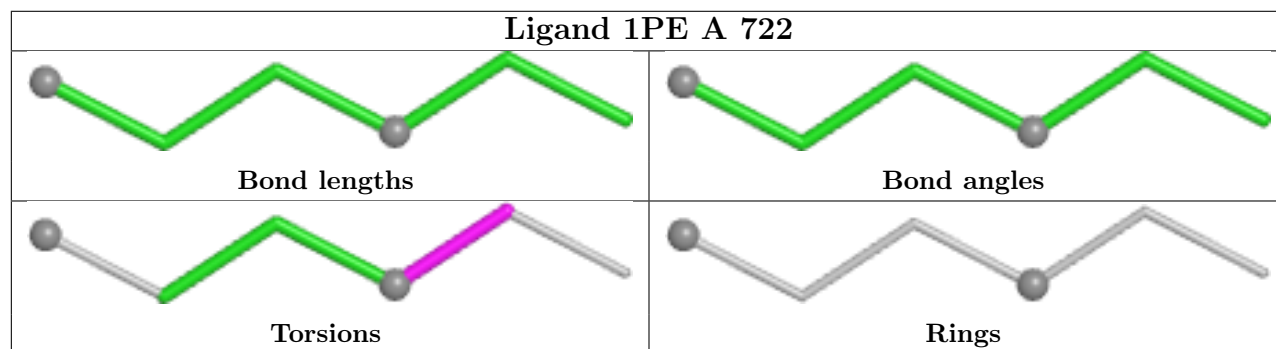
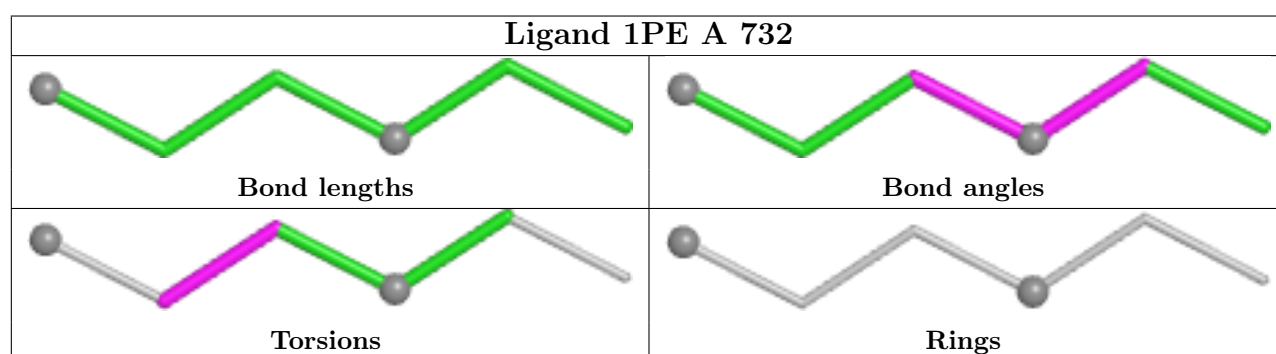
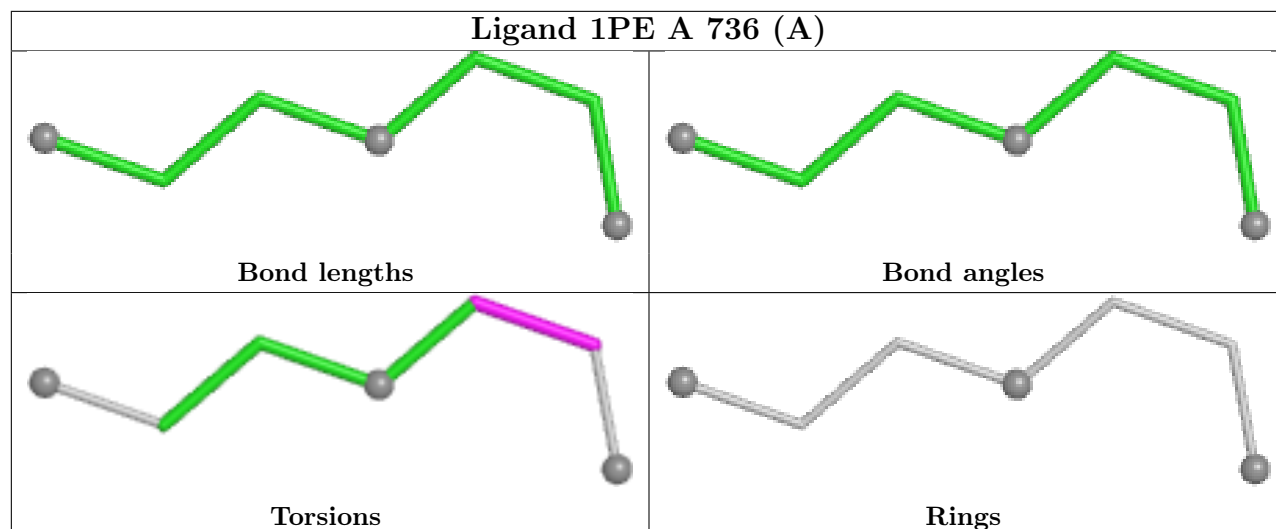
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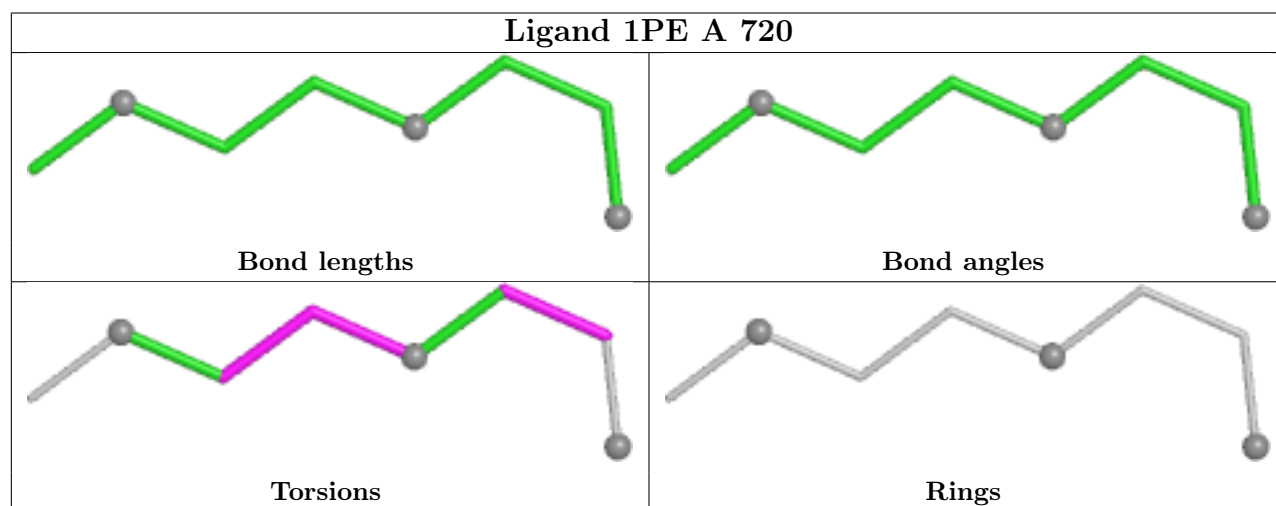
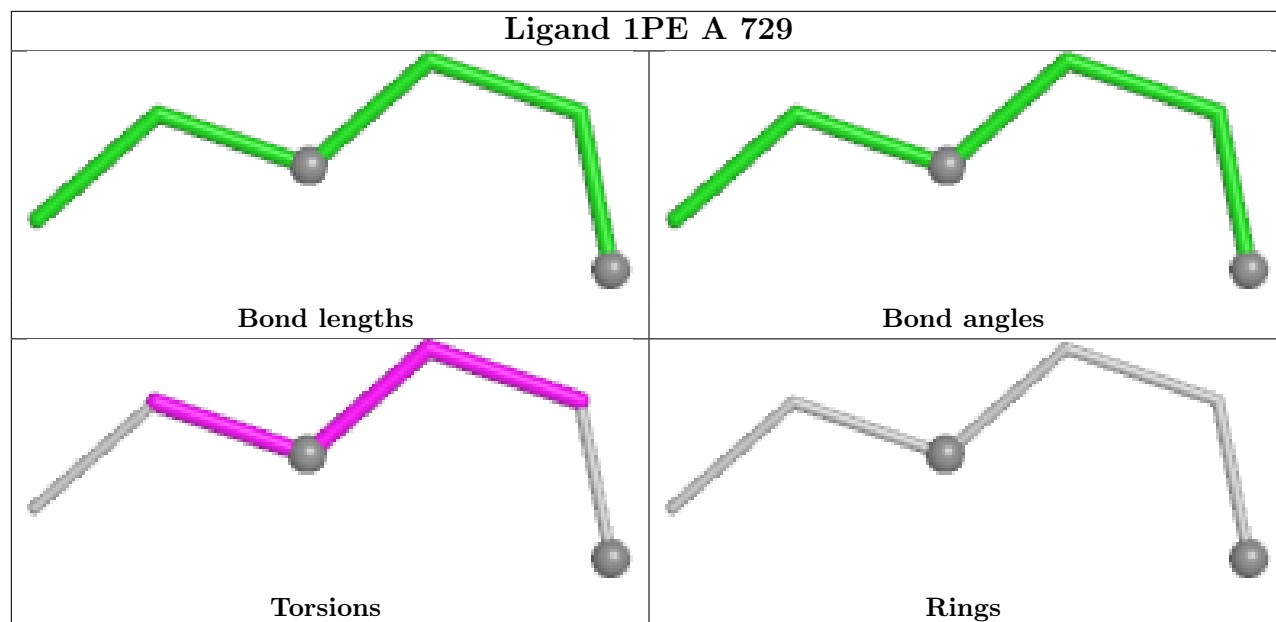
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	735	1PE	4	0
7	A	725	1PE	1	0
4	A	707	GOL	1	0
3	A	701	NAG	3	0
5	A	716	ACT	4	0
4	A	703	GOL	1	0
7	A	734[A]	1PE	1	0
7	A	734[B]	1PE	4	0
7	A	727	1PE	1	0

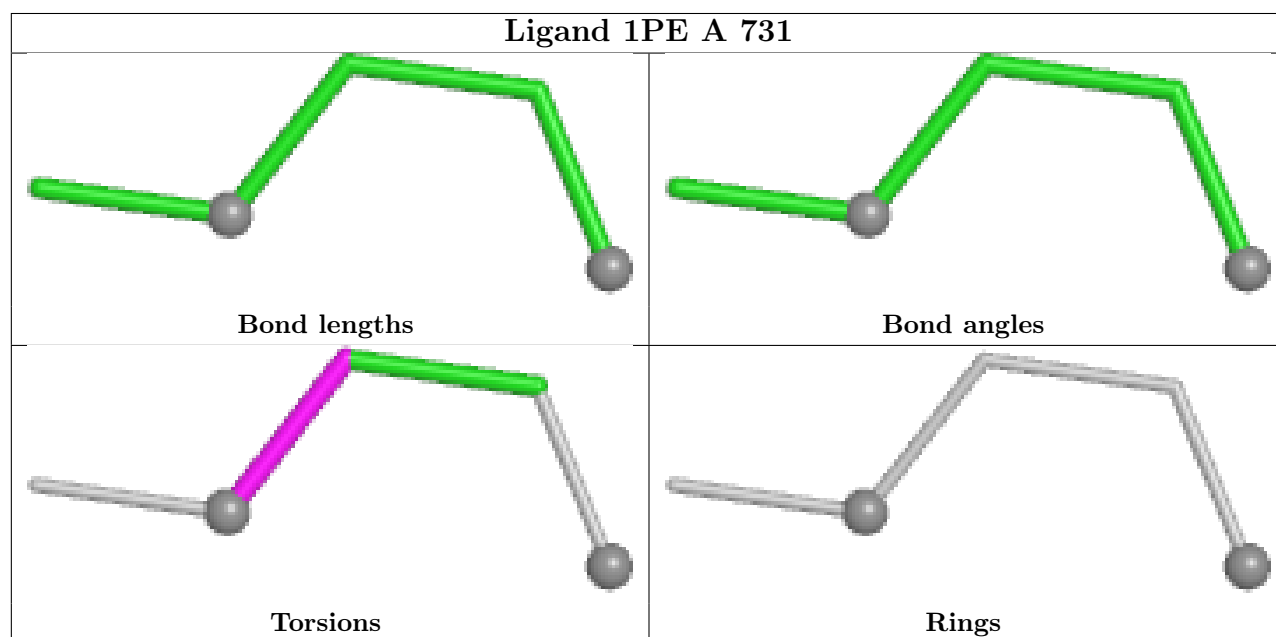
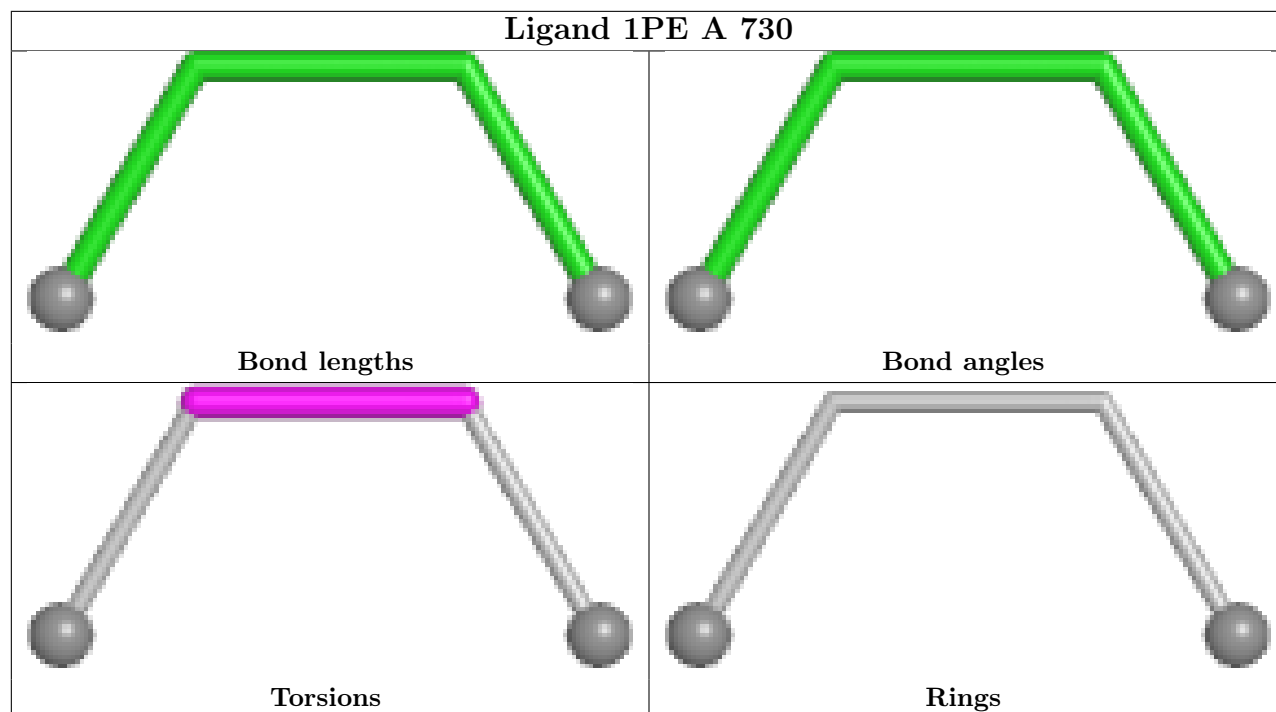
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

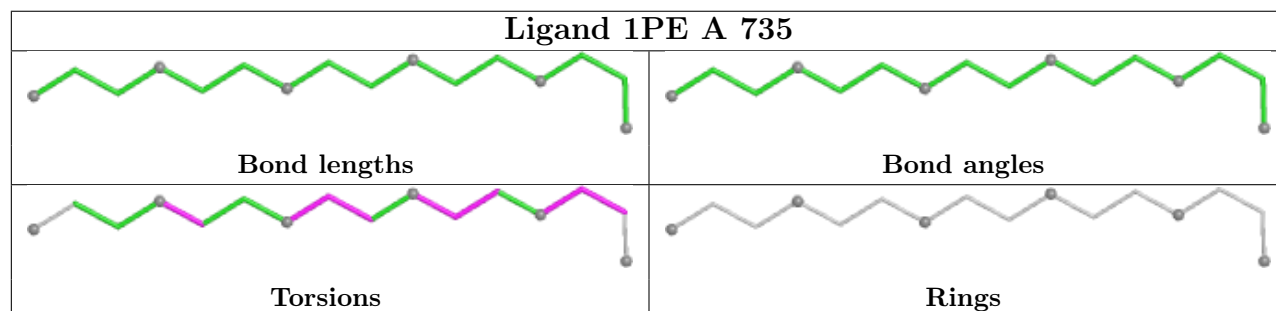
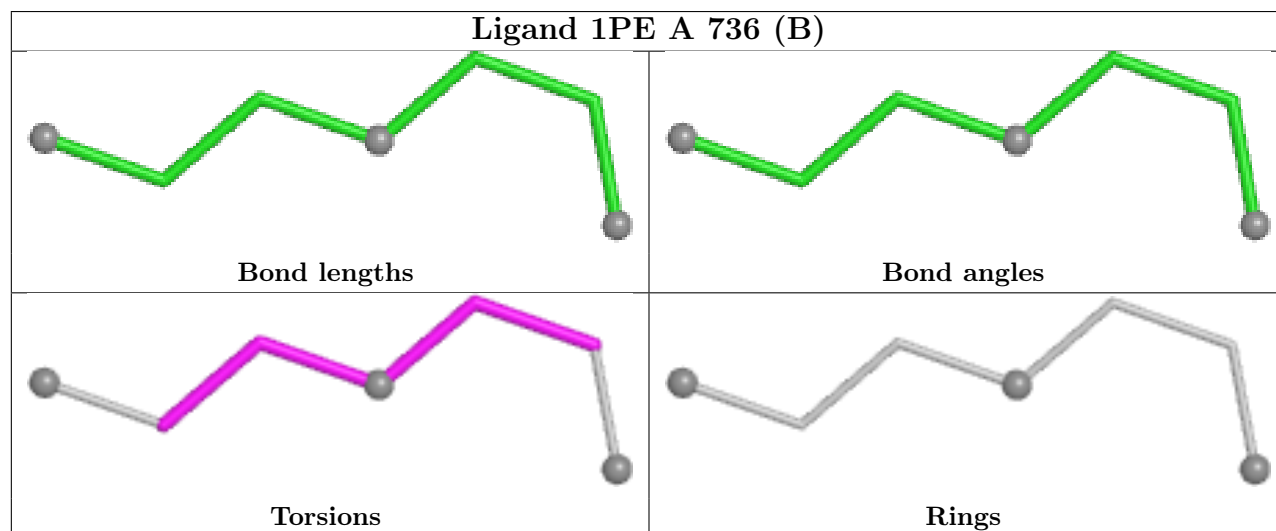
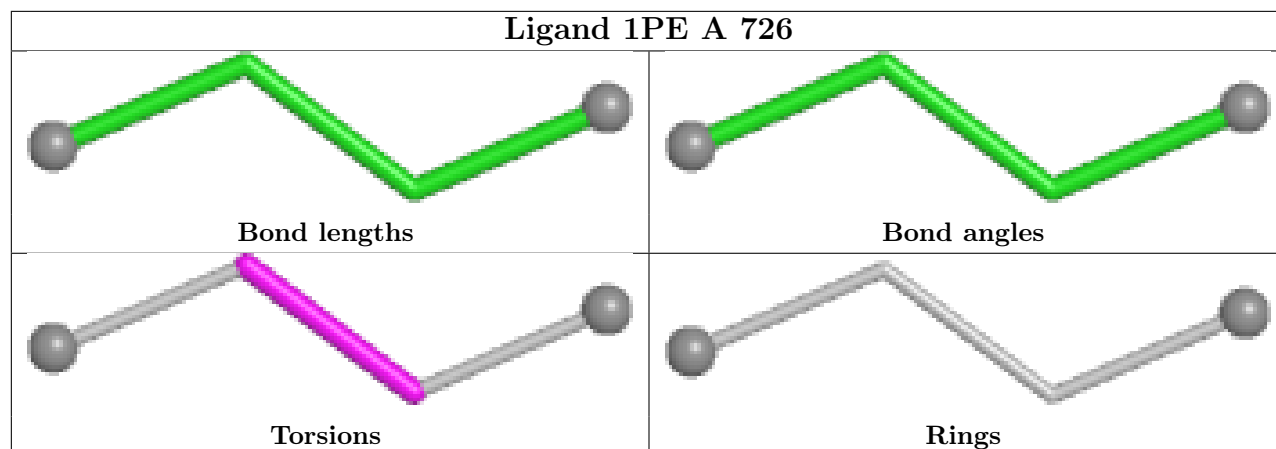


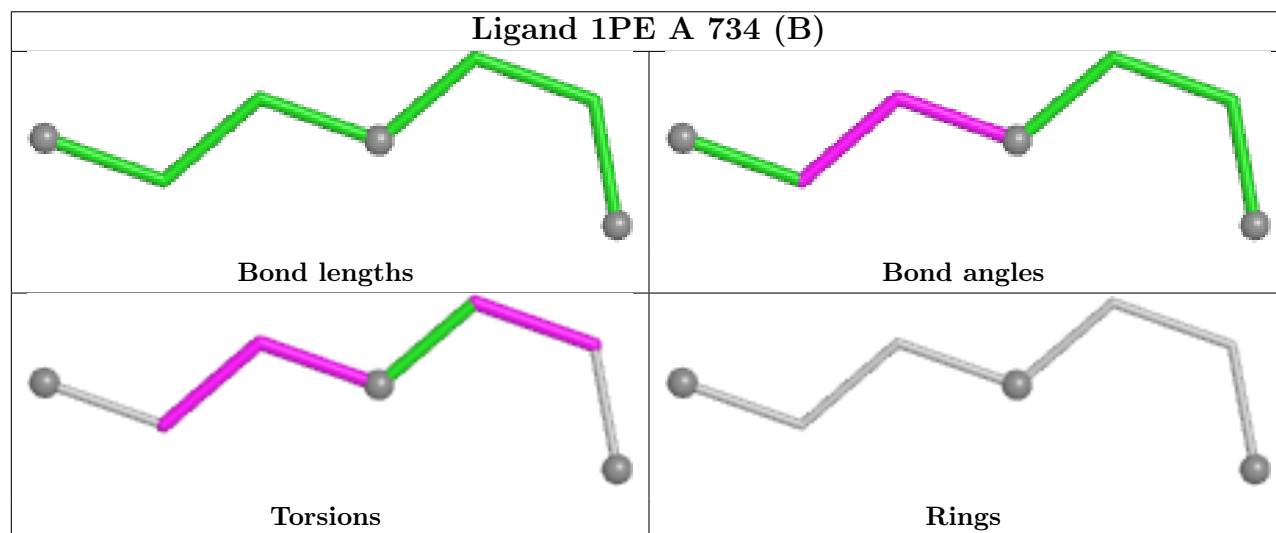
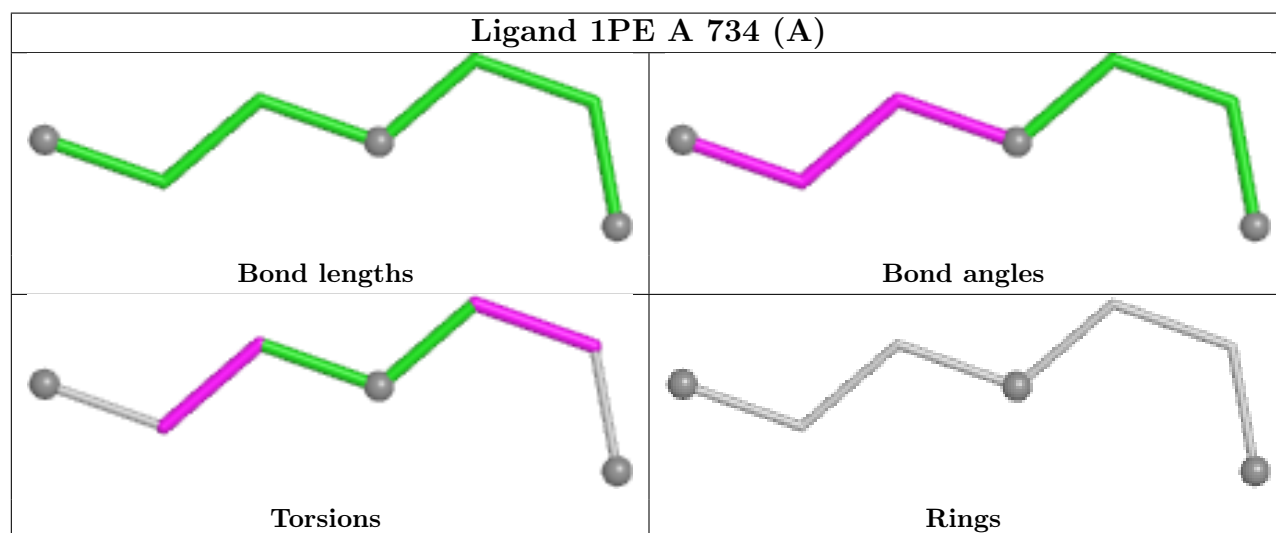
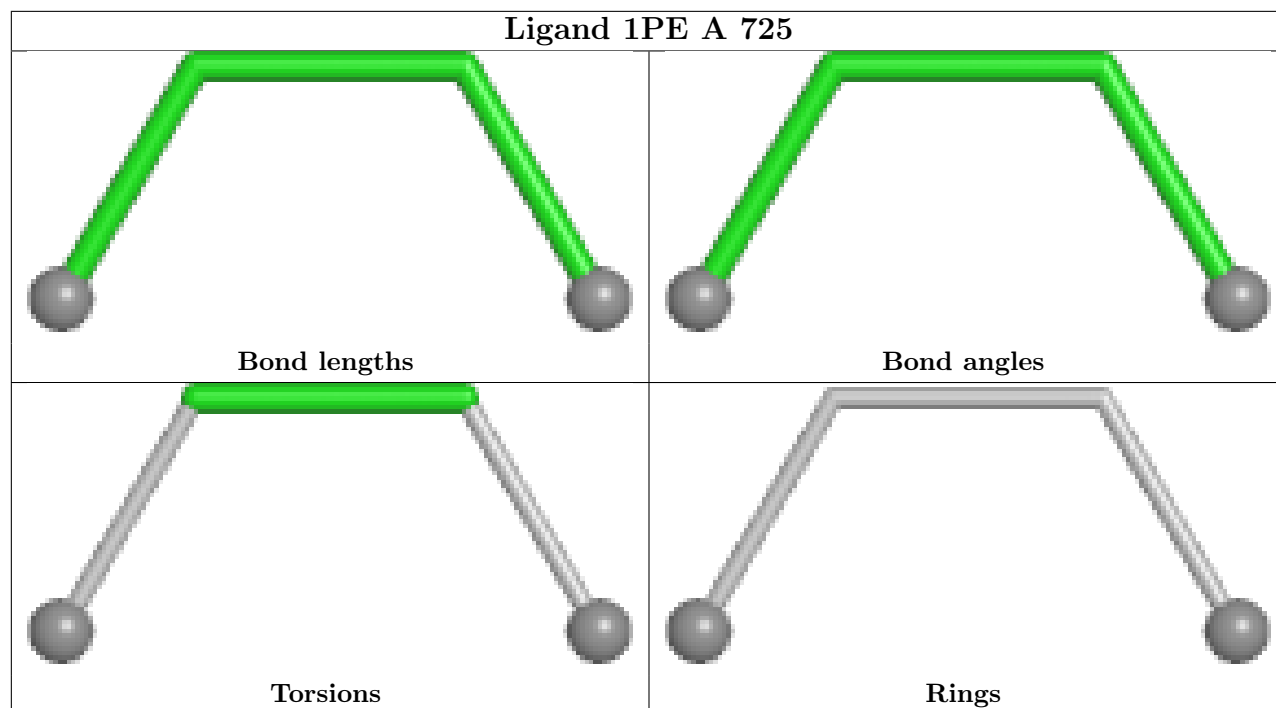


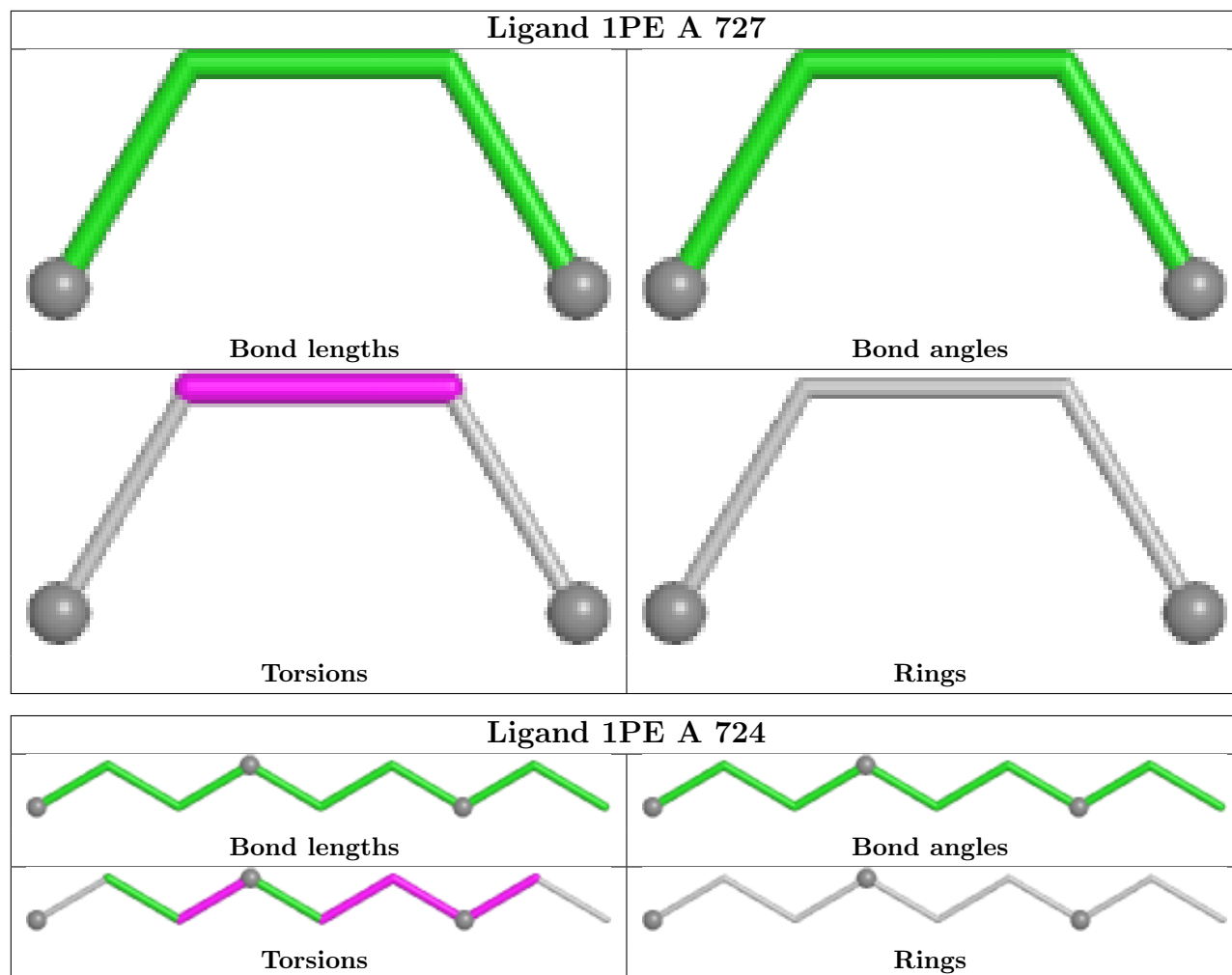












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	604/609 (99%)	-0.25	20 (3%)	49 52	18, 26, 48, 90	7 (1%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	218	ILE	5.1
1	A	586	HIS	5.1
1	A	215	VAL	4.4
1	A	-1	ALA	4.1
1	A	0	ALA	3.7
1	A	224	ILE	3.3
1	A	217	GLY	3.1
1	A	602	THR	2.8
1	A	186	LYS	2.6
1	A	221	ASN	2.6
1	A	222	ASN	2.6
1	A	256	TRP	2.5
1	A	216	ASP	2.4
1	A	491	GLU	2.2
1	A	220	GLU	2.1
1	A	187	ASP	2.1
1	A	504	PRO	2.1
1	A	225	ILE	2.1
1	A	257	ASN	2.0
1	A	294	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column

labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

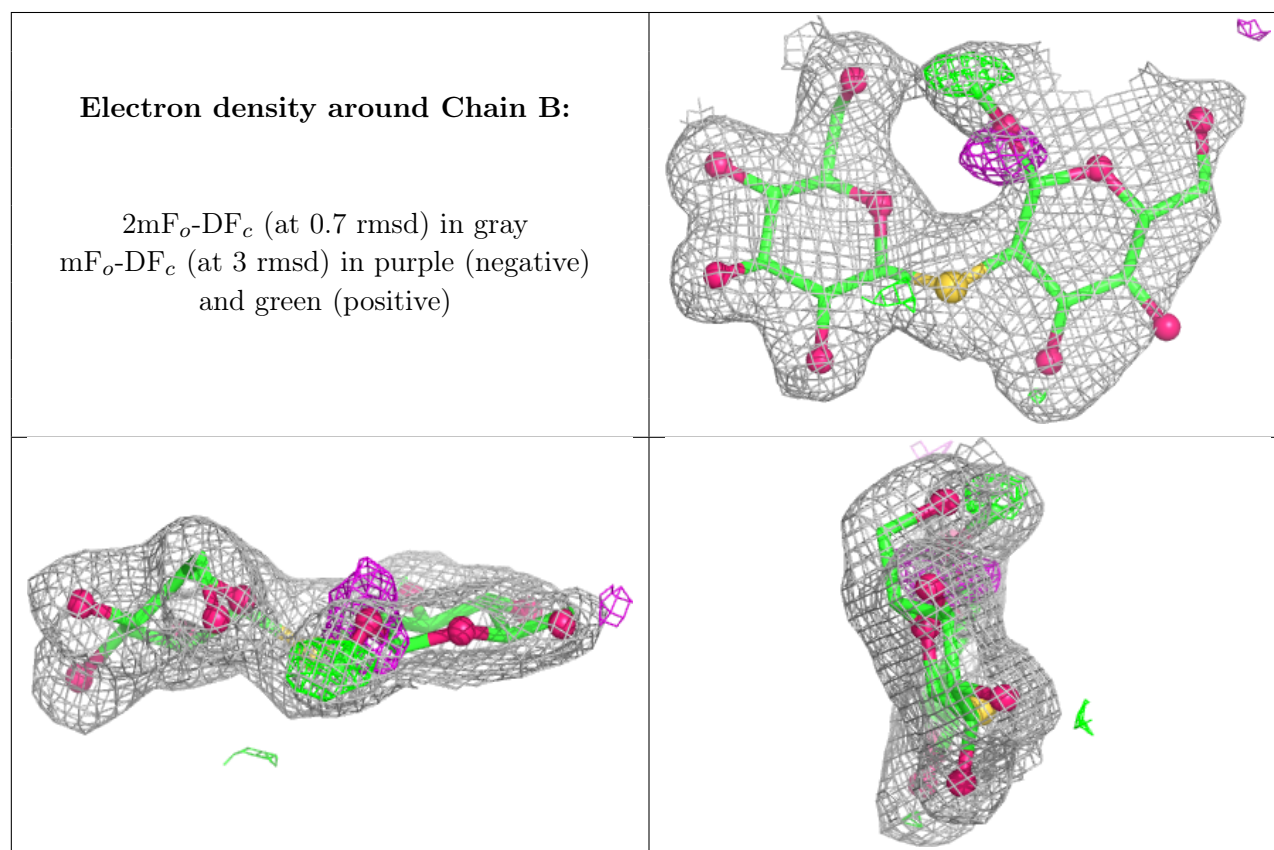
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	A	701	14/15	0.74	0.17	65,83,98,102	0

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	U2A	B	1[A]	13/13	0.92	0.14	34,56,68,76	1
2	U2A	B	1[B]	13/13	0.92	0.14	40,56,68,76	1
2	BGC	B	2	11/12	0.97	0.06	26,29,35,39	0

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



6.4 Ligands ⓘ

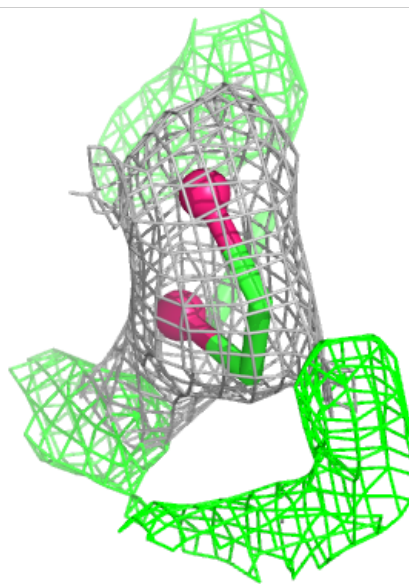
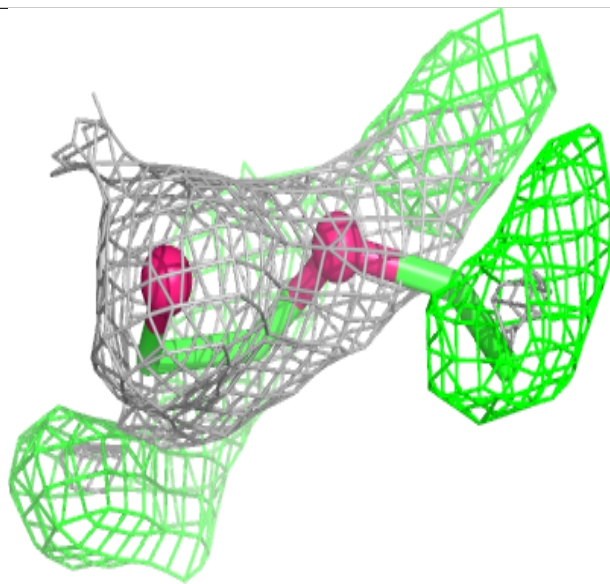
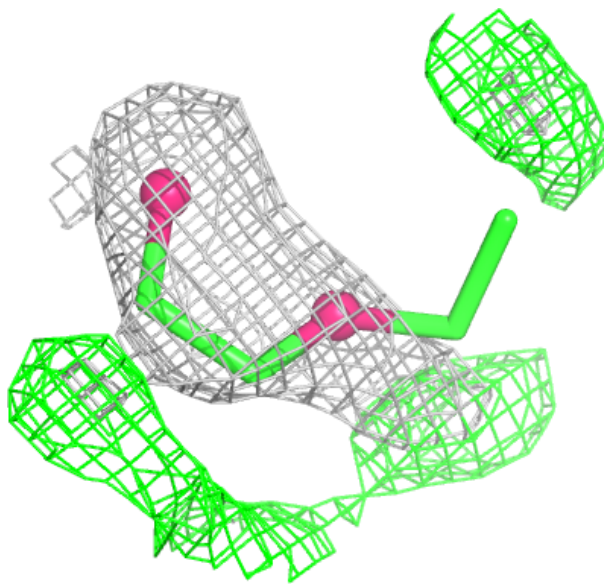
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	1PE	A	729	6/16	0.69	0.30	62,79,85,90	0
5	ACT	A	711	4/4	0.70	0.21	51,53,62,65	0
3	NAG	A	701	14/15	0.74	0.17	65,83,98,102	0
7	1PE	A	732	6/16	0.75	0.23	37,51,55,59	0
5	ACT	A	714	4/4	0.78	0.19	60,66,67,68	0
7	1PE	A	721	6/16	0.78	0.25	50,58,72,72	0
7	1PE	A	736[A]	7/16	0.78	0.21	47,49,54,56	5
7	1PE	A	736[B]	7/16	0.78	0.21	39,43,49,49	5
5	ACT	A	712	4/4	0.79	0.21	58,59,61,64	0
7	1PE	A	735	16/16	0.79	0.21	61,71,84,85	0
7	1PE	A	726	4/16	0.80	0.20	59,65,66,66	0
7	1PE	A	730	4/16	0.80	0.19	62,64,64,67	0
7	1PE	A	727	4/16	0.83	0.16	50,57,59,68	0
7	1PE	A	723	9/16	0.84	0.22	41,70,80,83	0
5	ACT	A	710	4/4	0.85	0.14	36,49,51,53	0
5	ACT	A	713	4/4	0.85	0.17	61,62,68,68	0
4	GOL	A	707	6/6	0.85	0.18	66,74,75,78	0
7	1PE	A	731	5/16	0.85	0.21	48,54,57,61	0
7	1PE	A	722	6/16	0.86	0.19	53,64,67,68	0
5	ACT	A	716	4/4	0.87	0.17	51,58,67,68	0
4	GOL	A	706	6/6	0.88	0.16	42,56,59,62	0
4	GOL	A	704	6/6	0.88	0.14	47,53,62,62	0
7	1PE	A	720	8/16	0.88	0.17	43,52,59,59	0
7	1PE	A	728	5/16	0.89	0.17	52,55,61,66	0
5	ACT	A	708	4/4	0.90	0.14	51,53,54,56	0
5	ACT	A	715	4/4	0.90	0.14	60,60,64,65	0
7	1PE	A	733	6/16	0.90	0.16	42,51,55,60	0
7	1PE	A	724	9/16	0.91	0.17	36,52,61,67	0
7	1PE	A	734[B]	7/16	0.92	0.15	38,45,55,60	2
6	SO4	A	719	5/5	0.92	0.12	30,43,46,47	5
5	ACT	A	709	4/4	0.92	0.17	32,45,46,57	0
7	1PE	A	734[A]	7/16	0.92	0.15	26,42,55,60	2
7	1PE	A	725	4/16	0.93	0.12	53,56,58,59	0
6	SO4	A	718	5/5	0.93	0.09	39,40,54,55	5
4	GOL	A	705	6/6	0.95	0.10	25,41,48,49	0
4	GOL	A	703	6/6	0.95	0.09	36,41,42,45	0
6	SO4	A	717	5/5	0.95	0.09	22,29,35,38	5

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.

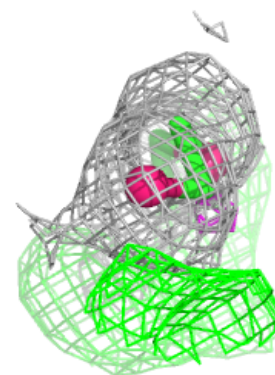
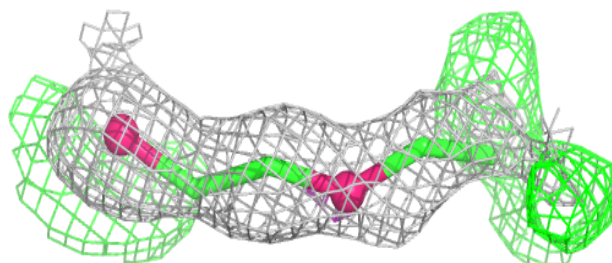
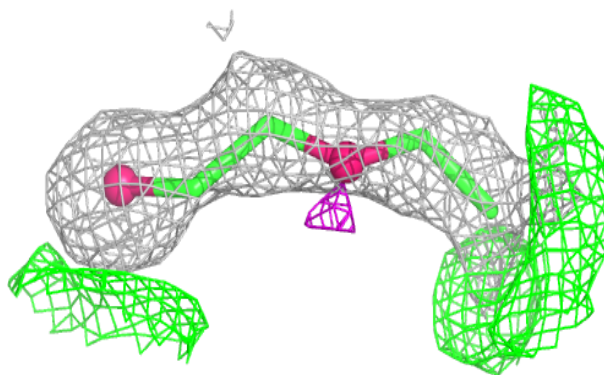
Electron density around 1PE A 729:

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and green (positive)

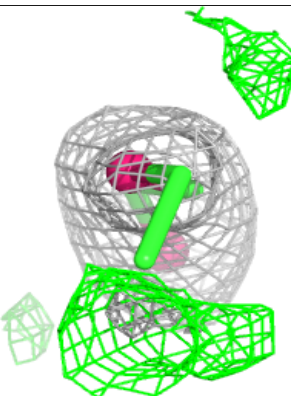
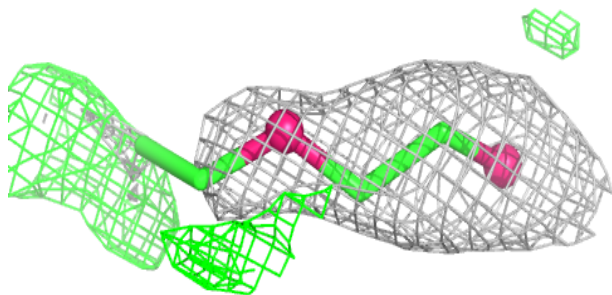
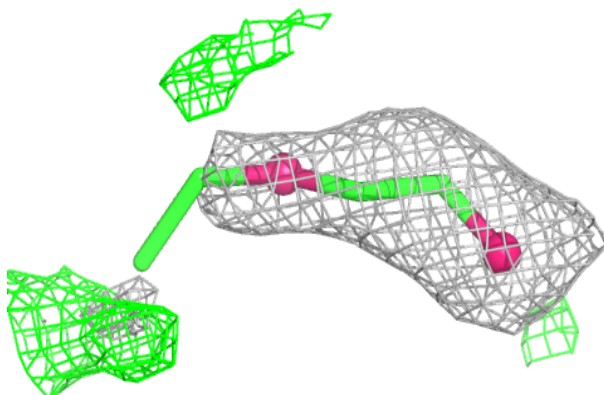


Electron density around 1PE A 732:

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and green (positive)

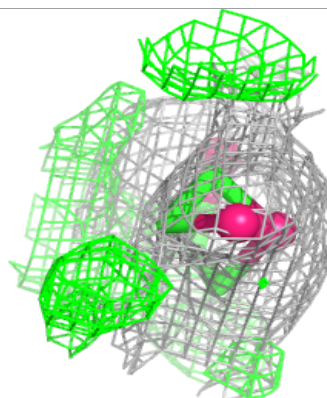
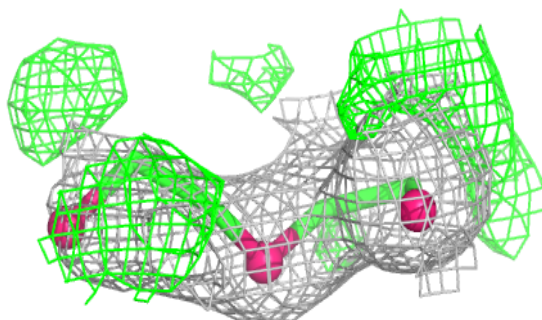
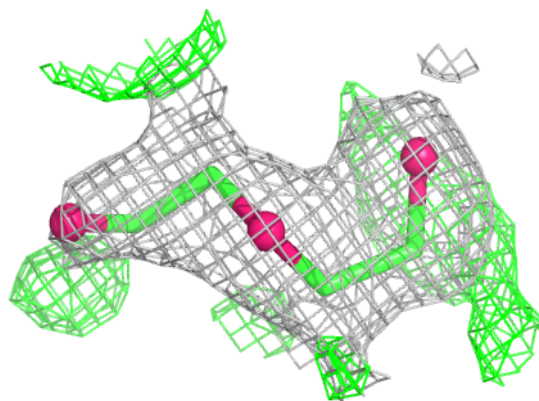
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and green (positive)

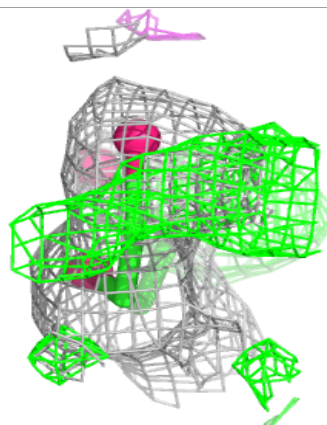
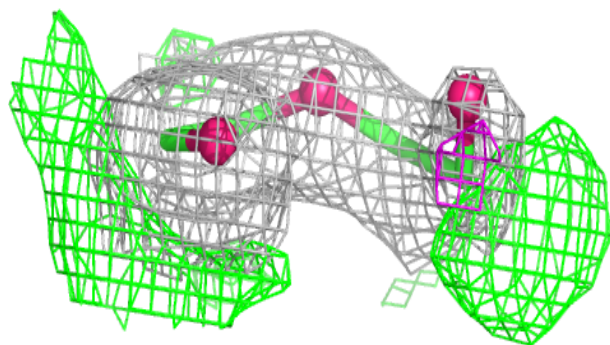
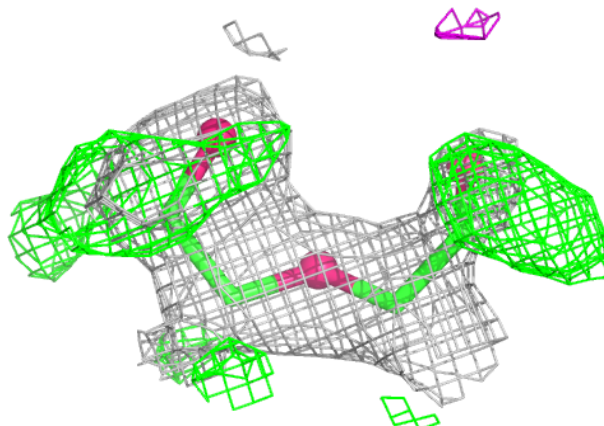


Electron density around 1PE A 736 (A):

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and green (positive)

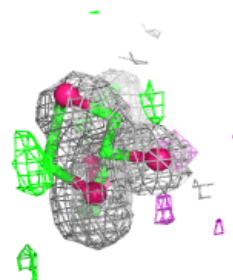
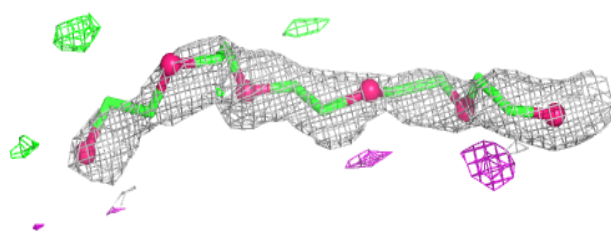
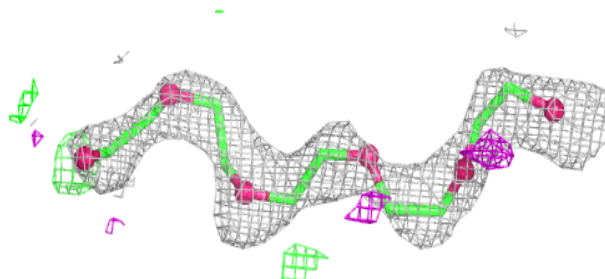
**Electron density around 1PE A 736 (B):**

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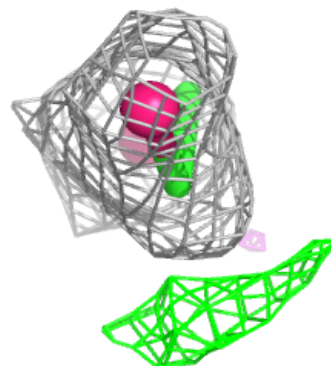
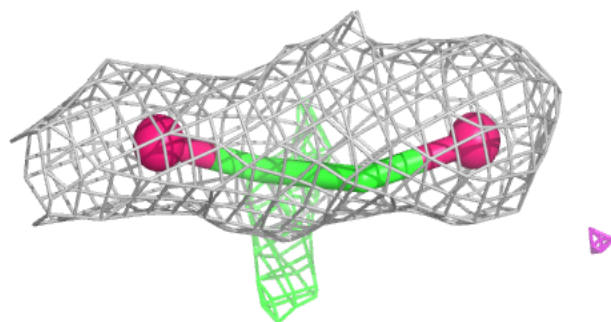
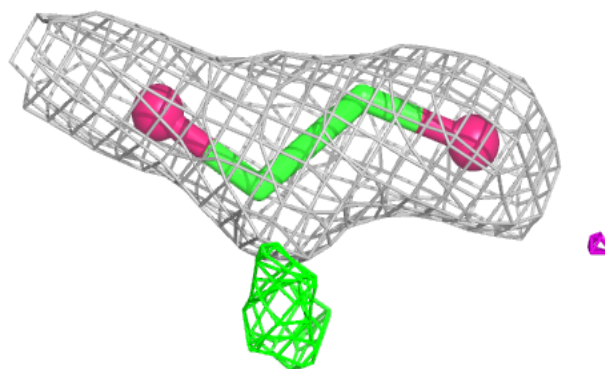


Electron density around 1PE A 735:

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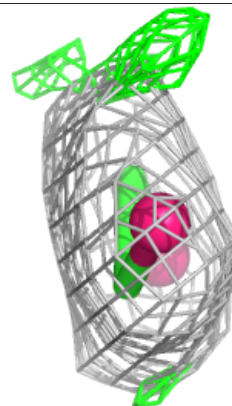
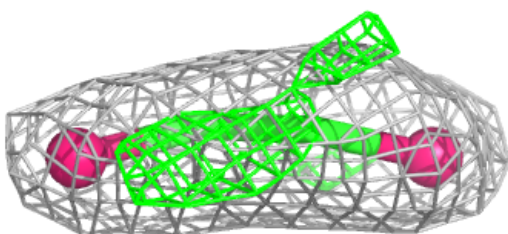
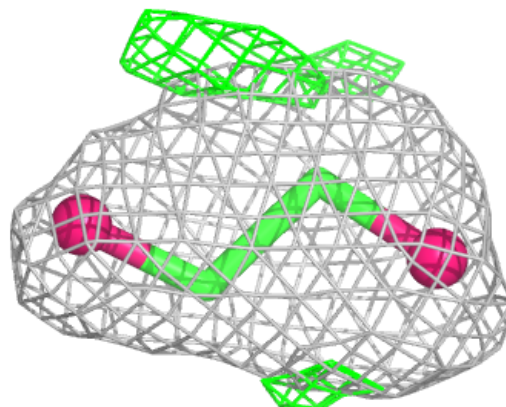
**Electron density around 1PE A 726:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



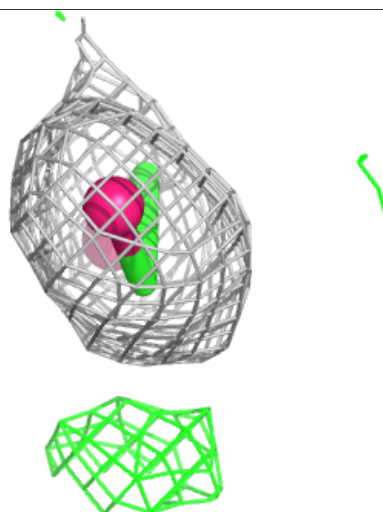
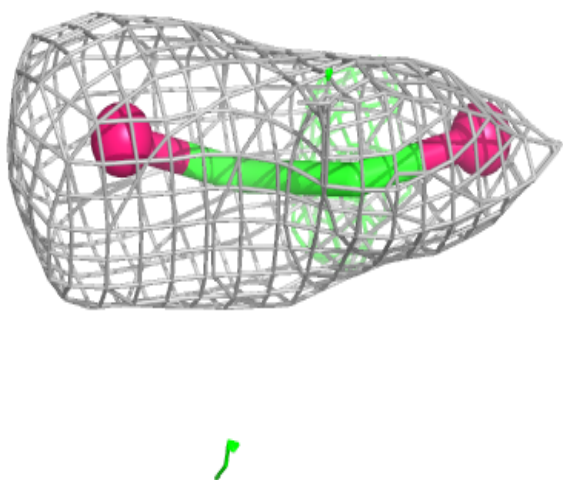
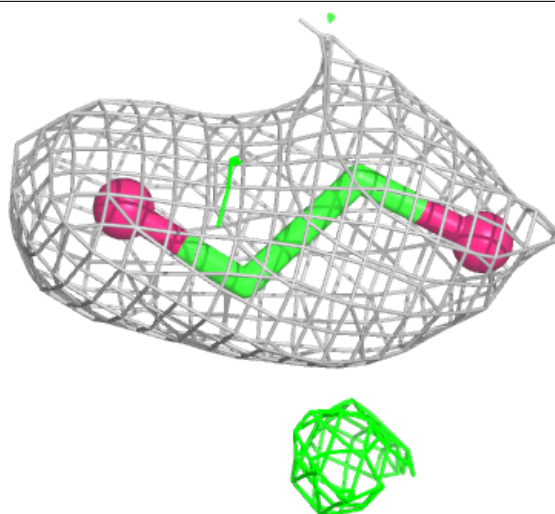
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and green (positive)



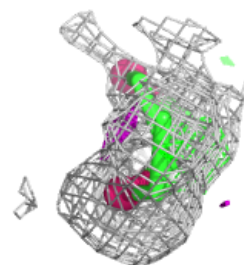
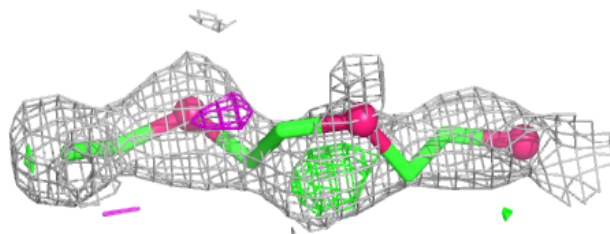
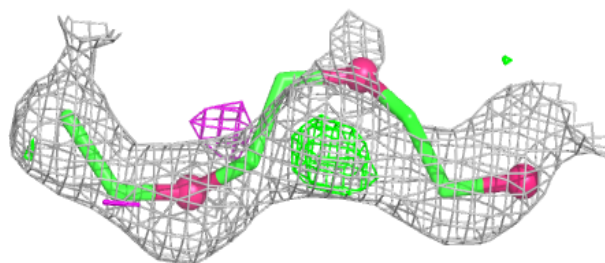
Electron density around 1PE A 727:

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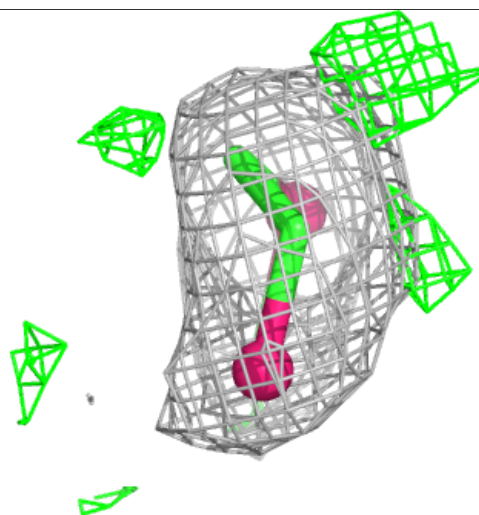
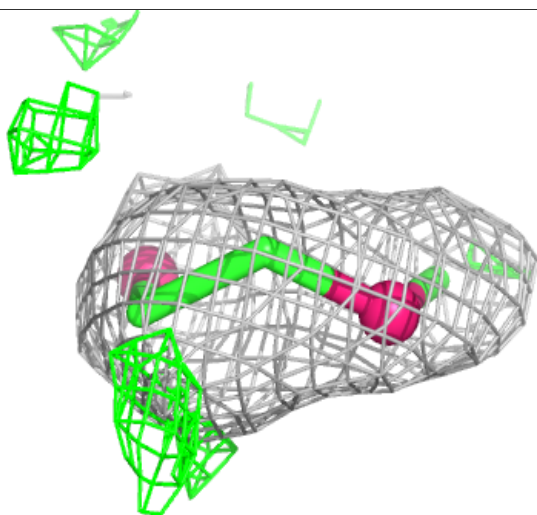
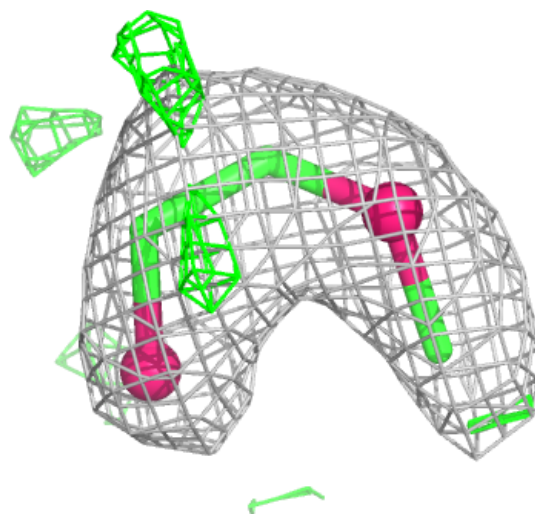
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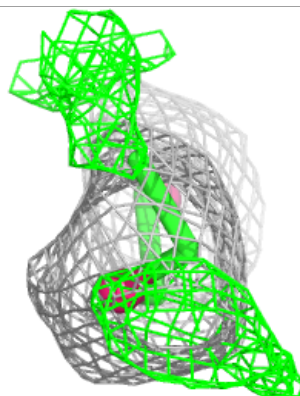
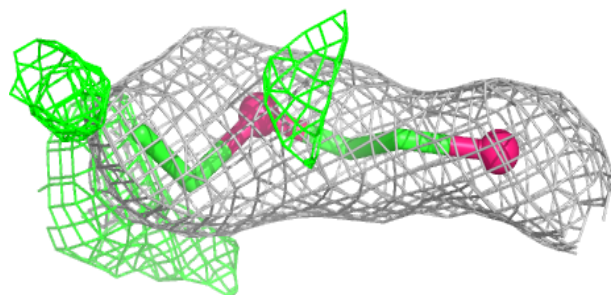
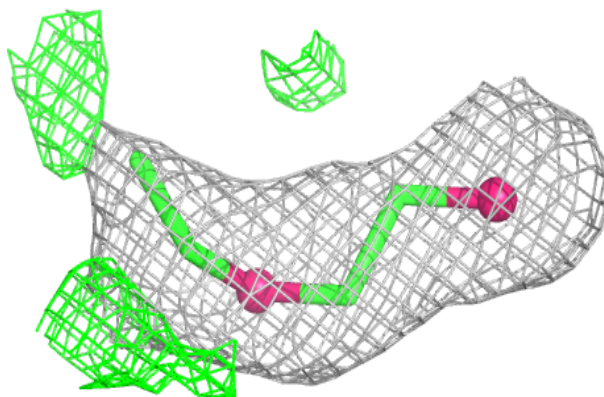
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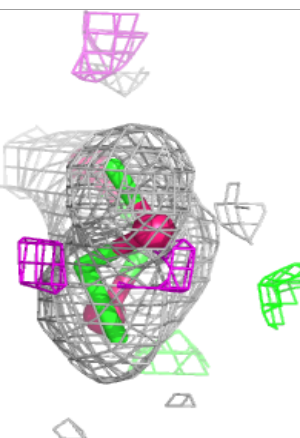
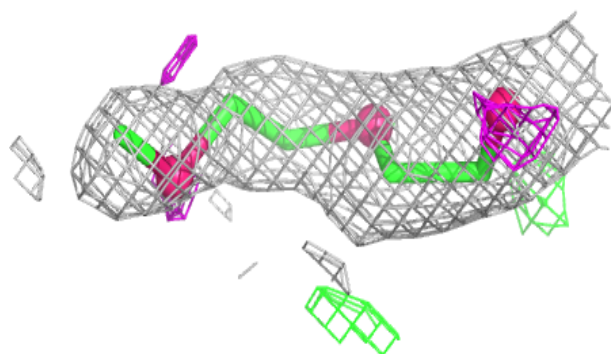
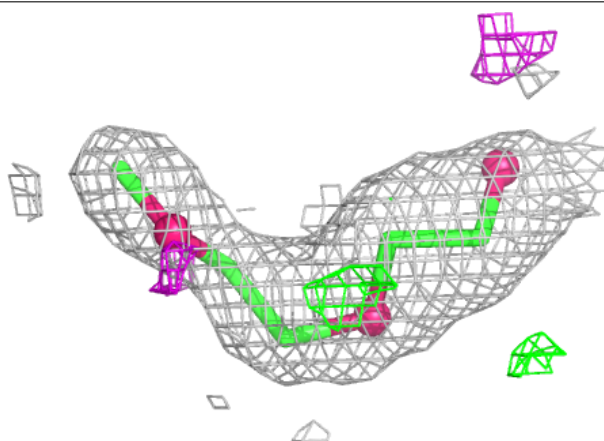


Electron density around 1PE A 722:

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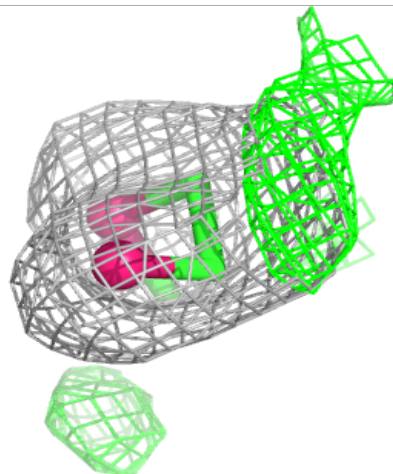
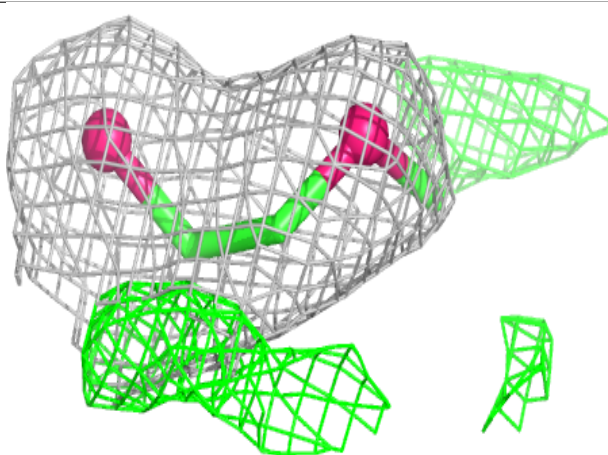
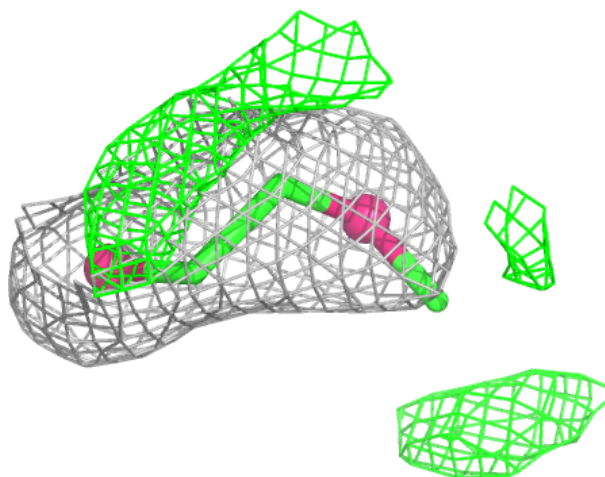
**Electron density around 1PE A 720:**

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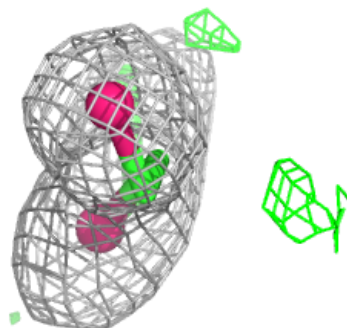
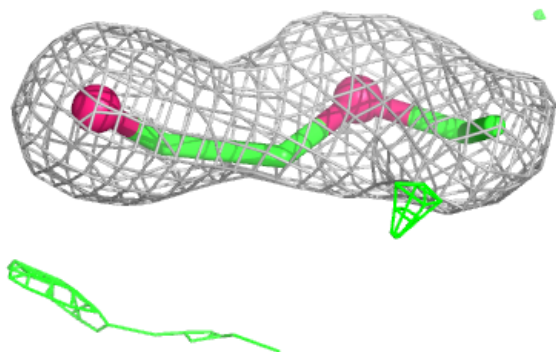
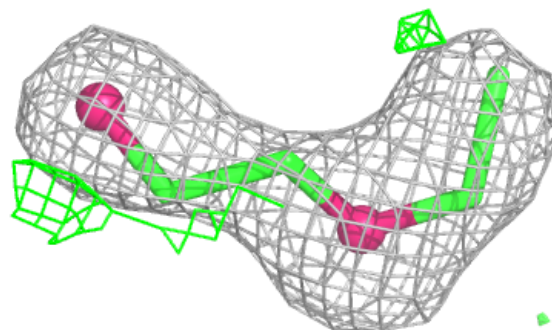
Electron density around 1PE A 728:

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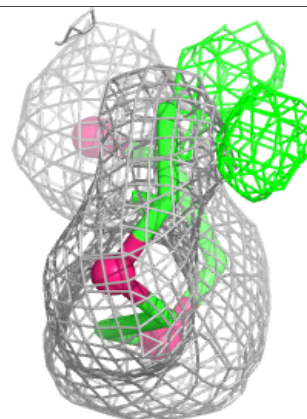
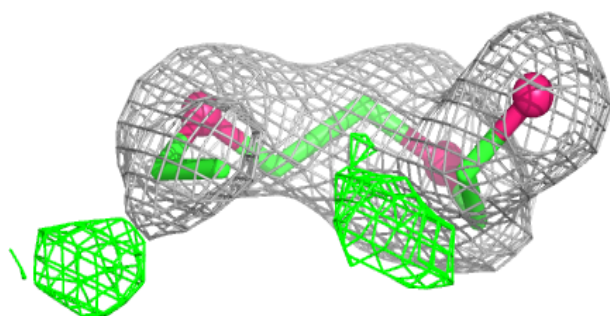
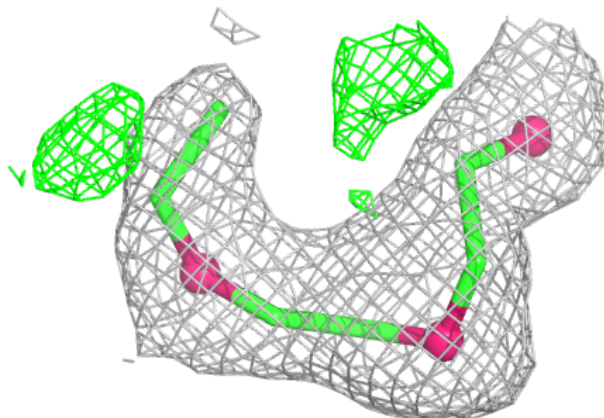


Electron density around 1PE A 733:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

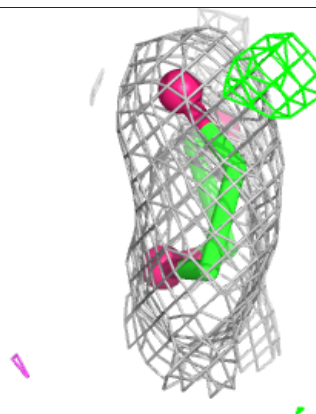
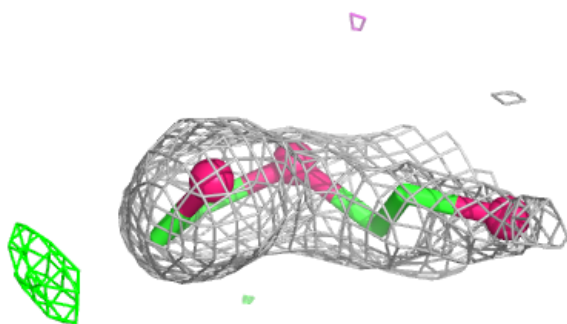
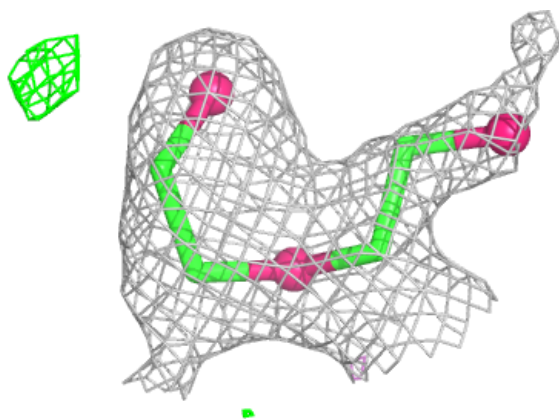
**Electron density around 1PE A 724:**

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

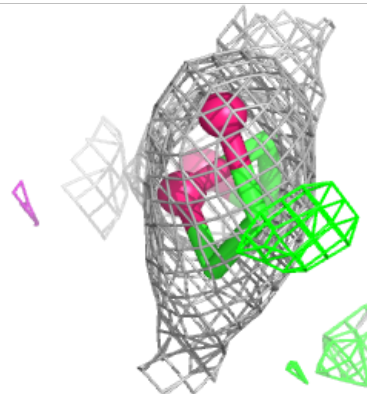
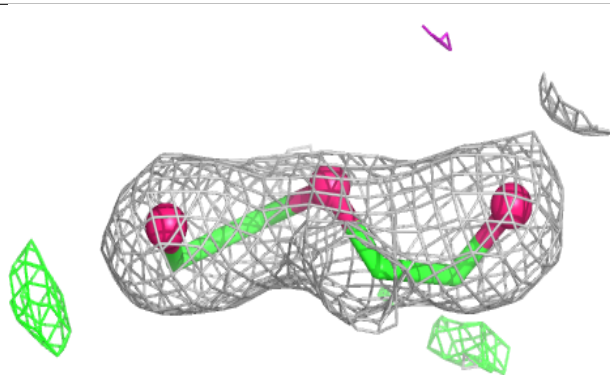
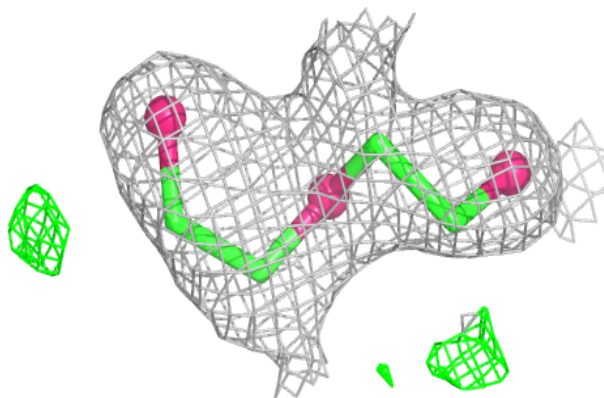


Electron density around 1PE A 734 (B):

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

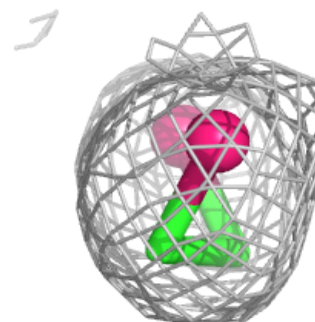
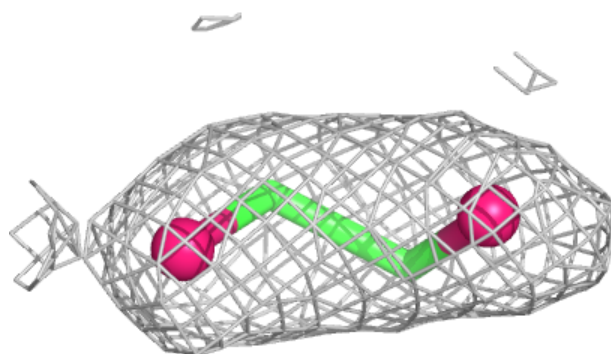
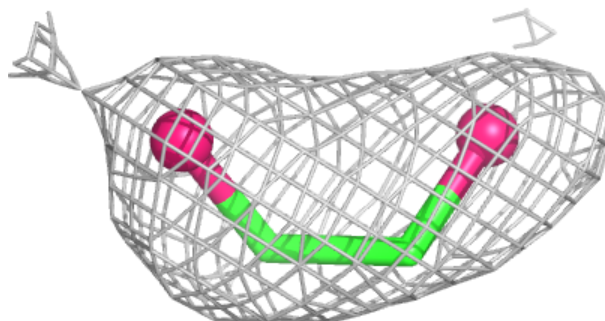
**Electron density around 1PE A 734 (A):**

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



Electron density around 1PE A 725:

$2mF_o - DF_c$ (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.