



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 05:56 am BST

PDB ID : 6BTZ
Title : Crystal structure of the PI3KC2alpha C2 domain in space group C121
Authors : Chen, K.-E.; Collins, B.M.
Deposited on : 2017-12-08
Resolution : 1.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

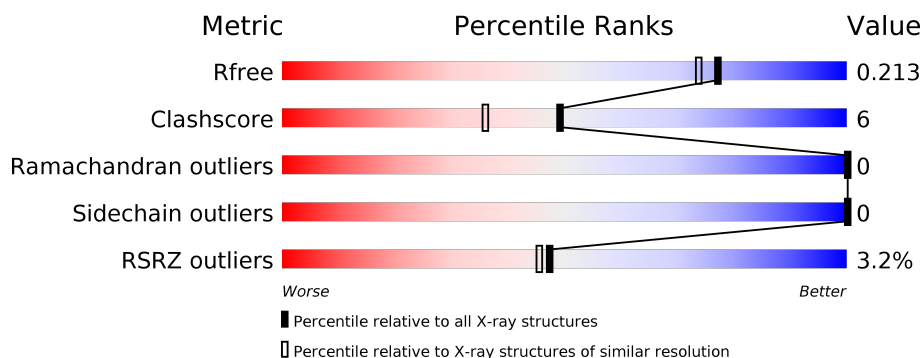
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	131	<div> <div>3%</div> <div>88%</div> <div>8%</div> <div>5%</div> </div>
1	B	131	<div> <div>3%</div> <div>83%</div> <div>12%</div> <div>5%</div> </div>
1	C	131	<div> <div>2%</div> <div>85%</div> <div>11%</div> <div>•</div> </div>
1	D	131	<div> <div>4%</div> <div>92%</div> <div>5%</div> <div>•</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	O4B	D	1704	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4717 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 Phosphatidylinositol 4-phosphate 3-kinase C2 domain-containing subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	125	Total	C	N	O	S	11	3	0
			1023	657	176	187	3			
1	B	125	Total	C	N	O	S	15	2	0
			1019	656	174	186	3			
1	C	126	Total	C	N	O	S	16	1	0
			1017	651	174	188	4			
1	D	126	Total	C	N	O	S	23	2	0
			1023	657	175	188	3			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP O00443
A	-1	ASN	-	expression tag	UNP O00443
A	0	ALA	-	expression tag	UNP O00443
B	-2	SER	-	expression tag	UNP O00443
B	-1	ASN	-	expression tag	UNP O00443
B	0	ALA	-	expression tag	UNP O00443
C	-2	SER	-	expression tag	UNP O00443
C	-1	ASN	-	expression tag	UNP O00443
C	0	ALA	-	expression tag	UNP O00443
D	-2	SER	-	expression tag	UNP O00443
D	-1	ASN	-	expression tag	UNP O00443
D	0	ALA	-	expression tag	UNP O00443

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



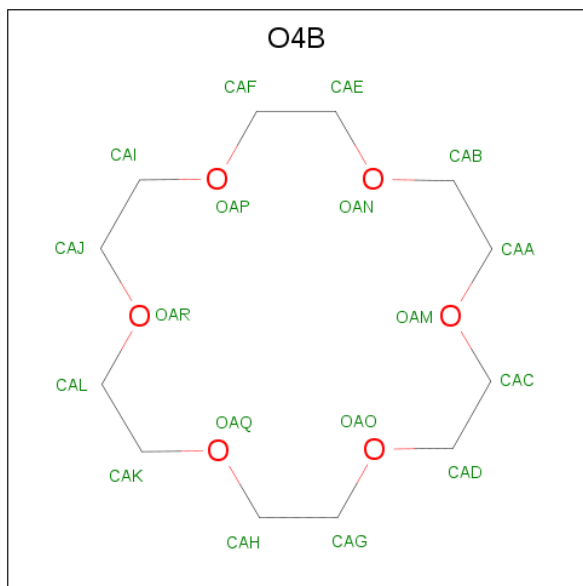
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 1,4,7,10,13,16-HEXAOXACYCLOOCTADECANE (three-letter code: O4B) (formula: $C_{12}H_{24}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			18	12	6		
3	A	1	Total	C	O	3	0
			18	12	6		
3	B	1	Total	C	O	0	0
			18	12	6		
3	C	1	Total	C	O	0	0
			18	12	6		
3	D	1	Total	C	O	0	0
			18	12	6		
3	D	1	Total	C	O	3	0
			18	12	6		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		

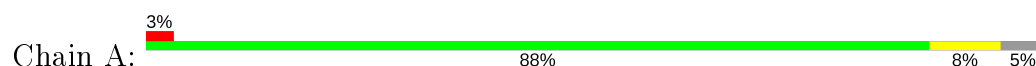
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	97	Total	O	0	0
			97	97		
5	B	103	Total	O	0	0
			103	103		
5	C	107	Total	O	0	0
			107	107		
5	D	98	Total	O	0	0
			98	98		

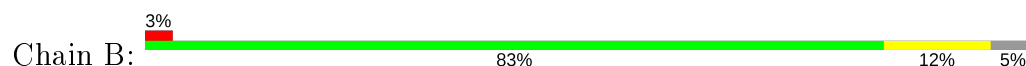
3 Residue-property plots [i](#)

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

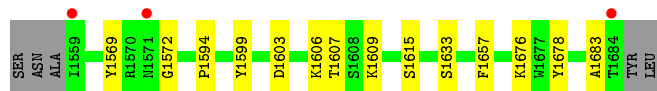
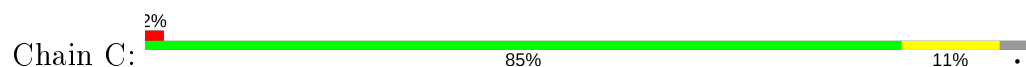
- Molecule 1: Phosphatidylinositol 4-phosphate 3-kinase C2 domain-containing subunit alpha



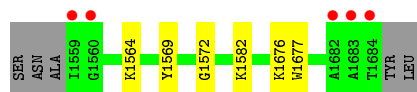
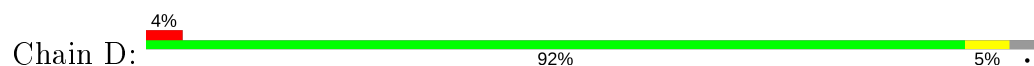
- Molecule 1: Phosphatidylinositol 4-phosphate 3-kinase C2 domain-containing subunit alpha



- Molecule 1: Phosphatidylinositol 4-phosphate 3-kinase C2 domain-containing subunit alpha



- Molecule 1: Phosphatidylinositol 4-phosphate 3-kinase C2 domain-containing subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	105.35Å 61.55Å 90.50Å 90.00° 92.58° 90.00°	Depositor
Resolution (Å)	46.40 – 1.85 46.39 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.40-1.85) 99.7 (46.39-1.85)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.15 (at 1.84Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.187 , 0.215 0.187 , 0.213	Depositor DCC
R_{free} test set	2525 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	17.4	Xtriage
Anisotropy	0.619	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.010 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.011 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.021 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.013 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.146 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4717	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.70% 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: GOL, O4B, SO4

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	0.40	0/1051	0.55	0/1415
1	B	0.36	0/1045	0.55	0/1408
1	C	0.36	0/1039	0.58	0/1401
1	D	0.35	0/1048	0.55	0/1413
All	All	0.37	0/4183	0.56	0/5637

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	1023	0	1069	12	0
1	B	1019	0	1057	17	0
1	C	1017	0	1051	9	0
1	D	1023	0	1066	12	0
2	A	20	0	0	0	0
2	B	25	0	0	0	0
2	C	25	0	0	0	0
2	D	10	0	0	0	0
3	A	36	0	48	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	18	0	24	3	0
3	C	18	0	24	1	0
3	D	36	0	48	11	0
4	A	18	0	24	2	0
4	B	18	0	24	6	0
4	C	6	0	8	1	0
5	A	97	0	0	2	0
5	B	103	0	0	1	0
5	C	107	0	0	0	0
5	D	98	0	0	0	0
All	All	4717	0	4443	51	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 6.

All (51) 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:1603:ASP:O	1:C:1606:LYS:HE2	1.60	0.99
1:B:1651:SER:HB3	4:B:1707:GOL:H12	1.66	0.77
1:B:1630:SER:H	4:B:1709:GOL:H32	1.50	0.75
1:B:1676:LYS:HE2	3:B:1706:O4B:HAK2	1.70	0.73
1:B:1676:LYS:HZ1	3:B:1706:O4B:HAC2	1.54	0.72
1:B:1634:LYS:HB2	4:B:1708:GOL:H2	1.72	0.72
1:A:1677:TRP:HD1	3:D:1704:O4B:HAH2	1.55	0.71
4:A:1707:GOL:H11	1:D:1564:LYS:HZ1	1.57	0.70
1:C:1676:LYS:HE2	3:C:1706:O4B:HAB2	1.74	0.69
1:D:1582:LYS:NZ	3:D:1704:O4B:HAK2	2.09	0.68
1:B:1587:GLU:HB2	1:B:1656[B]:PHE:CE2	2.30	0.67
1:B:1564:LYS:HG2	1:B:1579:MET:HE1	1.76	0.66
1:B:1676:LYS:NZ	3:B:1706:O4B:HAC2	2.11	0.66
1:A:1582:LYS:NZ	3:A:1706:O4B:HAH1	2.11	0.66
1:C:1633:SER:HB2	4:C:1707:GOL:H31	1.80	0.63
1:A:1582:LYS:HZ2	3:A:1706:O4B:HAC1	1.64	0.62
3:A:1706:O4B:HAA1	5:A:1802:HOH:O	2.01	0.60
1:A:1582:LYS:NZ	3:A:1706:O4B:HAC1	2.17	0.59
1:B:1630:SER:H	4:B:1709:GOL:C3	2.14	0.58
1:D:1676:LYS:HZ1	3:D:1703:O4B:HAH2	1.69	0.57
1:D:1676:LYS:HD2	3:D:1703:O4B:HAL1	1.86	0.57
1:D:1582:LYS:HD2	3:D:1704:O4B:HAJ2	1.86	0.56
3:A:1706:O4B:HAK1	1:D:1677:TRP:HD1	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1603:ASP:O	1:A:1606[B]:LYS:HE2	2.07	0.55
4:A:1707:GOL:H11	1:D:1564:LYS:NZ	2.22	0.55
1:D:1582:LYS:NZ	3:D:1704:O4B:HAA1	2.25	0.52
1:A:1582:LYS:HZ2	3:A:1706:O4B:HAH1	1.74	0.52
1:D:1582:LYS:HZ2	3:D:1704:O4B:HAK2	1.77	0.50
1:A:1566:SER:HB3	1:A:1577:MET:HB3	1.95	0.49
5:A:1854:HOH:O	3:D:1704:O4B:HAH1	2.12	0.49
1:B:1566:SER:HB3	1:B:1577:MET:HB3	1.94	0.48
1:A:1678:TYR:HE2	3:A:1705:O4B:HAH1	1.79	0.48
1:D:1582:LYS:HZ1	3:D:1704:O4B:HAA1	1.78	0.48
1:B:1652:LEU:HB3	1:C:1607:THR:HA	1.96	0.48
1:C:1657:PHE:HD2	1:C:1683:ALA:HB2	1.79	0.47
1:C:1676:LYS:HE3	1:C:1678:TYR:CE1	2.50	0.47
1:A:1676:LYS:NZ	3:A:1705:O4B:HAH2	2.30	0.46
1:B:1569:TYR:OH	1:B:1572:GLY:HA2	2.16	0.45
1:A:1569:TYR:OH	1:A:1572:GLY:HA2	2.16	0.45
1:C:1594:PRO:HD2	1:C:1615:SER:O	2.18	0.44
1:D:1569:TYR:OH	1:D:1572:GLY:HA2	2.18	0.43
1:D:1582:LYS:CE	3:D:1704:O4B:HAA1	2.48	0.43
1:B:1564:LYS:HG2	1:B:1579:MET:CE	2.45	0.43
1:A:1677:TRP:CD1	3:D:1704:O4B:HAH2	2.43	0.42
1:A:1678:TYR:CE2	3:A:1705:O4B:HAH1	2.54	0.42
1:B:1616:ARG:HH22	4:B:1707:GOL:H2	1.85	0.42
1:B:1648:SER:HB3	1:B:1656[B]:PHE:CE1	2.55	0.42
1:C:1569:TYR:OH	1:C:1572:GLY:HA2	2.21	0.41
1:B:1559:ILE:O	5:B:1801:HOH:O	2.22	0.41
1:C:1599:TYR:OH	1:C:1609:LYS:HD3	2.21	0.41
1:B:1630:SER:N	4:B:1709:GOL:H32	2.26	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	126/131 (96%)	123 (98%)	3 (2%)	0	100	100
1	B	125/131 (95%)	123 (98%)	2 (2%)	0	100	100
1	C	125/131 (95%)	122 (98%)	3 (2%)	0	100	100
1	D	126/131 (96%)	123 (98%)	3 (2%)	0	100	100
All	All	502/524 (96%)	491 (98%)	11 (2%)	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	116/118 (98%)	116 (100%)	0	100	100
1	B	115/118 (98%)	115 (100%)	0	100	100
1	C	115/118 (98%)	115 (100%)	0	100	100
1	D	116/118 (98%)	116 (100%)	0	100	100
All	All	462/472 (98%)	462 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

29 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	$\# Z > 2$
2	SO4	B	1705	-	4,4,4	0.12	0	6,6,6	0.10	0
2	SO4	C	1703	-	4,4,4	0.15	0	6,6,6	0.08	0
4	GOL	A	1709	-	5,5,5	0.39	0	5,5,5	0.36	0
4	GOL	B	1709	-	5,5,5	0.39	0	5,5,5	0.36	0
2	SO4	D	1701	-	4,4,4	0.16	0	6,6,6	0.20	0
2	SO4	A	1702	-	4,4,4	0.15	0	6,6,6	0.18	0
2	SO4	B	1704	-	4,4,4	0.16	0	6,6,6	0.14	0
3	O4B	D	1704	-	18,18,18	0.66	0	18,18,18	0.57	0
2	SO4	A	1704	-	4,4,4	0.14	0	6,6,6	0.07	0
2	SO4	B	1701	-	4,4,4	0.20	0	6,6,6	0.21	0
2	SO4	C	1702	-	4,4,4	0.19	0	6,6,6	0.12	0
3	O4B	A	1706	-	18,18,18	0.73	0	18,18,18	0.61	0
3	O4B	B	1706	-	18,18,18	0.71	0	18,18,18	0.55	0
2	SO4	B	1702	-	4,4,4	0.16	0	6,6,6	0.18	0
2	SO4	C	1705	-	4,4,4	0.15	0	6,6,6	0.08	0
4	GOL	A	1707	-	5,5,5	0.40	0	5,5,5	0.14	0
2	SO4	C	1704	-	4,4,4	0.13	0	6,6,6	0.17	0
3	O4B	D	1703	-	18,18,18	0.62	0	18,18,18	0.57	0
2	SO4	D	1702	-	4,4,4	0.14	0	6,6,6	0.19	0
2	SO4	B	1703	-	4,4,4	0.16	0	6,6,6	0.06	0
2	SO4	C	1701	-	4,4,4	0.14	0	6,6,6	0.16	0
3	O4B	A	1705	-	18,18,18	0.62	0	18,18,18	0.56	0
2	SO4	A	1701	-	4,4,4	0.12	0	6,6,6	0.16	0
4	GOL	B	1708	-	5,5,5	0.40	0	5,5,5	0.38	0
3	O4B	C	1706	-	18,18,18	0.65	0	18,18,18	0.59	0
2	SO4	A	1703	-	4,4,4	0.18	0	6,6,6	0.23	0
4	GOL	B	1707	-	5,5,5	0.27	0	5,5,5	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	A	1708	-	5,5,5	0.37	0	5,5,5	0.44	0
4	GOL	C	1707	-	5,5,5	0.33	0	5,5,5	0.55	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
3	O4B	C	1706	-	-	9/18/18/18	0/1/1/1
3	O4B	D	1703	-	-	8/18/18/18	0/1/1/1
3	O4B	A	1705	-	-	8/18/18/18	0/1/1/1
4	GOL	B	1707	-	-	2/4/4/4	-
4	GOL	A	1709	-	-	1/4/4/4	-
4	GOL	B	1709	-	-	4/4/4/4	-
3	O4B	A	1706	-	-	9/18/18/18	1/1/1/1
3	O4B	B	1706	-	-	8/18/18/18	0/1/1/1
4	GOL	A	1707	-	-	2/4/4/4	-
4	GOL	A	1708	-	-	2/4/4/4	-
4	GOL	C	1707	-	-	0/4/4/4	-
3	O4B	D	1704	-	-	9/18/18/18	1/1/1/1
4	GOL	B	1708	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (66) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1709	GOL	O1-C1-C2-C3
4	A	1707	GOL	O1-C1-C2-C3
4	B	1708	GOL	O1-C1-C2-C3
4	B	1708	GOL	C1-C2-C3-O3
4	B	1708	GOL	O2-C2-C3-O3
4	B	1707	GOL	O1-C1-C2-C3
3	D	1704	O4B	OAQ-CAK-CAL-OAR
3	B	1706	O4B	OAO-CAG-CAH-OAQ
3	B	1706	O4B	OAP-CAI-CAJ-OAR

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Mol	Chain	Res	Type	Atoms
3	A	1705	O4B	OAQ-CAG-CAH-OAQ
3	C	1706	O4B	OAQ-CAK-CAL-OAR
3	B	1706	O4B	OAM-CAA-CAB-OAN
3	C	1706	O4B	OAP-CAI-CAJ-OAR
3	D	1703	O4B	OAQ-CAK-CAL-OAR
3	D	1704	O4B	OAM-CAA-CAB-OAN
3	A	1705	O4B	OAM-CAA-CAB-OAN
3	A	1706	O4B	OAQ-CAK-CAL-OAR
3	A	1706	O4B	OAM-CAA-CAB-OAN
3	A	1705	O4B	OAM-CAC-CAD-OAO
3	B	1706	O4B	OAQ-CAK-CAL-OAR
3	A	1705	O4B	OAQ-CAK-CAL-OAR
4	B	1709	GOL	C1-C2-C3-O3
4	B	1709	GOL	O1-C1-C2-O2
4	B	1709	GOL	O2-C2-C3-O3
4	A	1707	GOL	O1-C1-C2-O2
4	B	1708	GOL	O1-C1-C2-O2
4	B	1707	GOL	O1-C1-C2-O2
3	C	1706	O4B	OAM-CAA-CAB-OAN
3	D	1703	O4B	OAM-CAA-CAB-OAN
3	C	1706	O4B	OAM-CAC-CAD-OAO
3	B	1706	O4B	CAC-CAD-OAO-CAG
3	D	1703	O4B	CAH-CAG-OAO-CAD
3	A	1705	O4B	CAH-CAG-OAO-CAD
3	C	1706	O4B	CAH-CAG-OAO-CAD
3	A	1706	O4B	CAK-CAL-OAR-CAJ
3	A	1706	O4B	CAH-CAG-OAO-CAD
3	B	1706	O4B	CAK-CAL-OAR-CAJ
3	D	1703	O4B	CAC-CAD-OAO-CAG
3	D	1704	O4B	CAA-CAB-OAN-CAE
3	D	1704	O4B	CAH-CAG-OAO-CAD
3	A	1706	O4B	CAC-CAD-OAO-CAG
3	A	1705	O4B	CAL-CAK-OAQ-CAH
3	C	1706	O4B	CAI-CAJ-OAR-CAL
3	D	1704	O4B	CAC-CAD-OAO-CAG
3	D	1703	O4B	CAA-CAB-OAN-CAE
4	A	1709	GOL	O1-C1-C2-C3
3	A	1705	O4B	CAD-CAC-OAM-CAA
3	B	1706	O4B	CAB-CAA-OAM-CAC
3	D	1704	O4B	CAL-CAK-OAQ-CAH
3	A	1706	O4B	CAA-CAB-OAN-CAE
3	D	1703	O4B	OAM-CAC-CAD-OAO

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Mol	Chain	Res	Type	Atoms
3	C	1706	O4B	CAK-CAL-OAR-CAJ
3	B	1706	O4B	OAM-CAC-CAD-OAO
3	A	1705	O4B	OAN-CAE-CAF-OAP
4	A	1708	GOL	C1-C2-C3-O3
4	A	1708	GOL	O2-C2-C3-O3
3	A	1706	O4B	OAM-CAC-CAD-OAO
3	A	1706	O4B	OAO-CAG-CAH-OAQ
3	D	1703	O4B	OAO-CAG-CAH-OAQ
3	C	1706	O4B	CAF-CAE-OAN-CAB
3	D	1704	O4B	CAK-CAL-OAR-CAJ
3	C	1706	O4B	CAL-CAK-OAQ-CAH
3	D	1704	O4B	OAO-CAG-CAH-OAQ
3	D	1703	O4B	CAB-CAA-OAM-CAC
3	D	1704	O4B	OAM-CAC-CAD-OAO
3	A	1706	O4B	CAD-CAC-OAM-CAA

All (2) ring outliers are listed below:

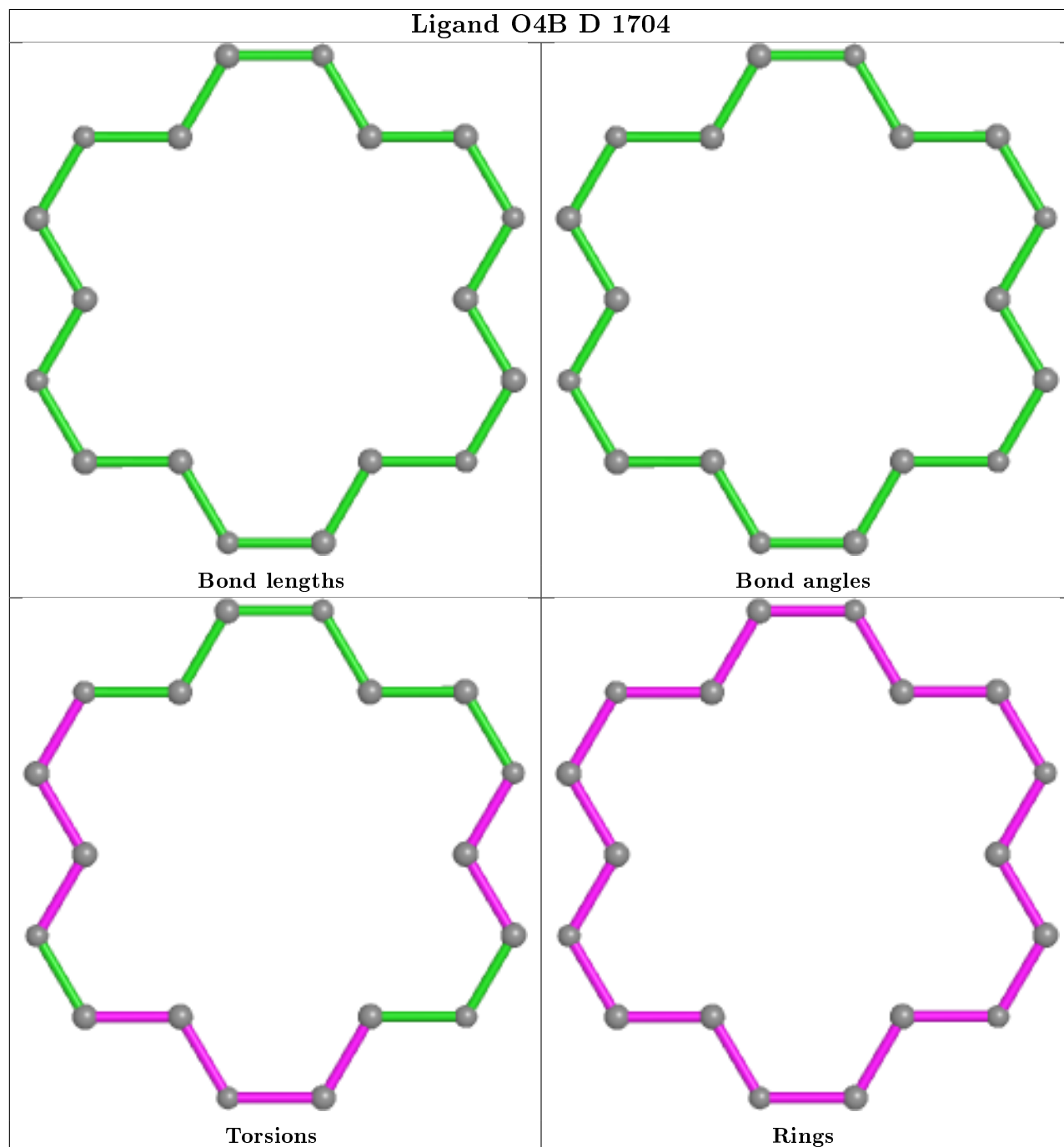
Mol	Chain	Res	Type	Atoms
3	D	1704	O4B	CAA-CAB-CAC-CAD-CAE-CAF-CAG-CAH-CAI-CAJ-CAK-CAL-OAM-O
3	A	1706	O4B	CAA-CAB-CAC-CAD-CAE-CAF-CAG-CAH-CAI-CAJ-CAK-CAL-OAM-O

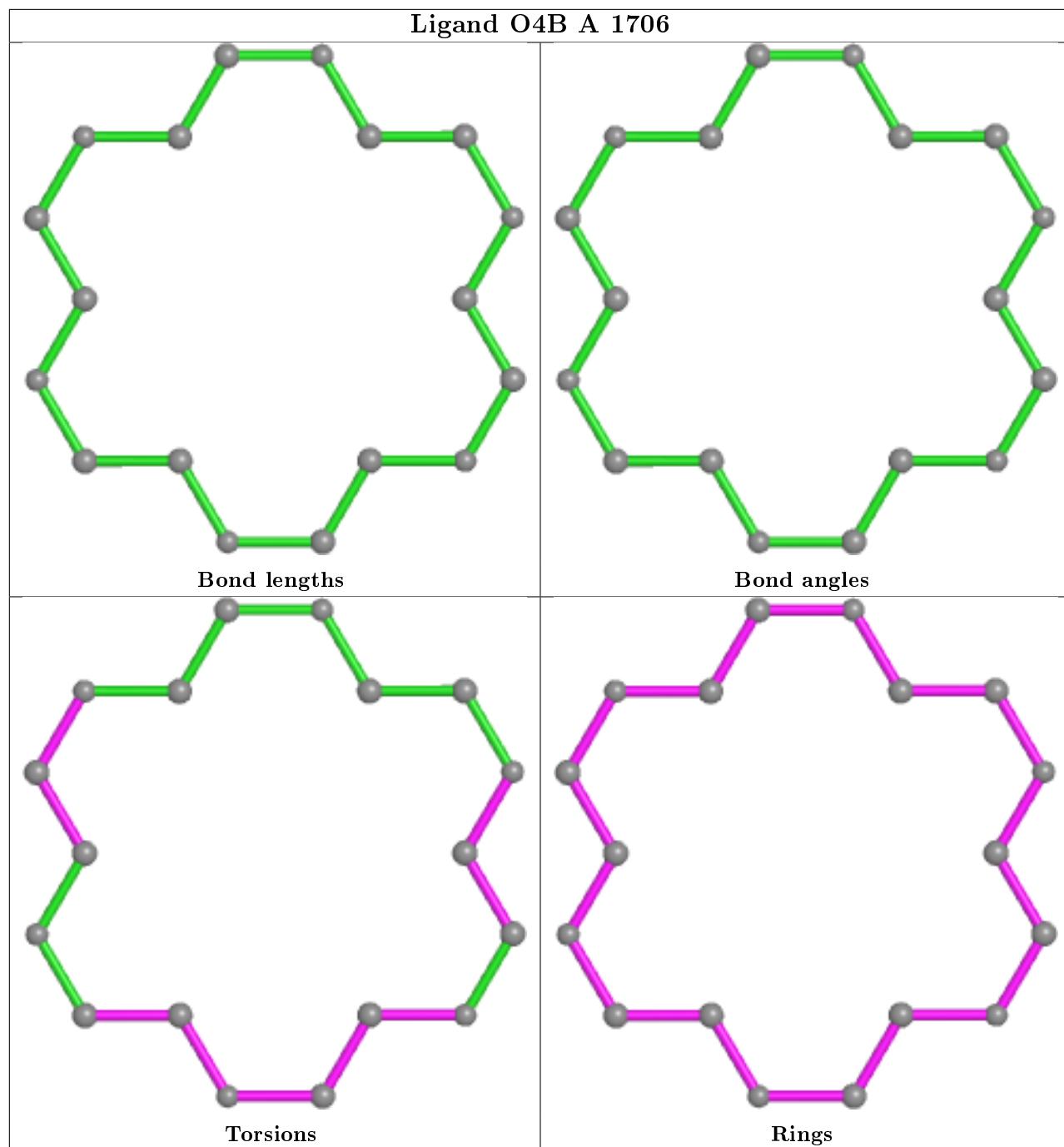
11 monomers are involved in 33 short contacts:

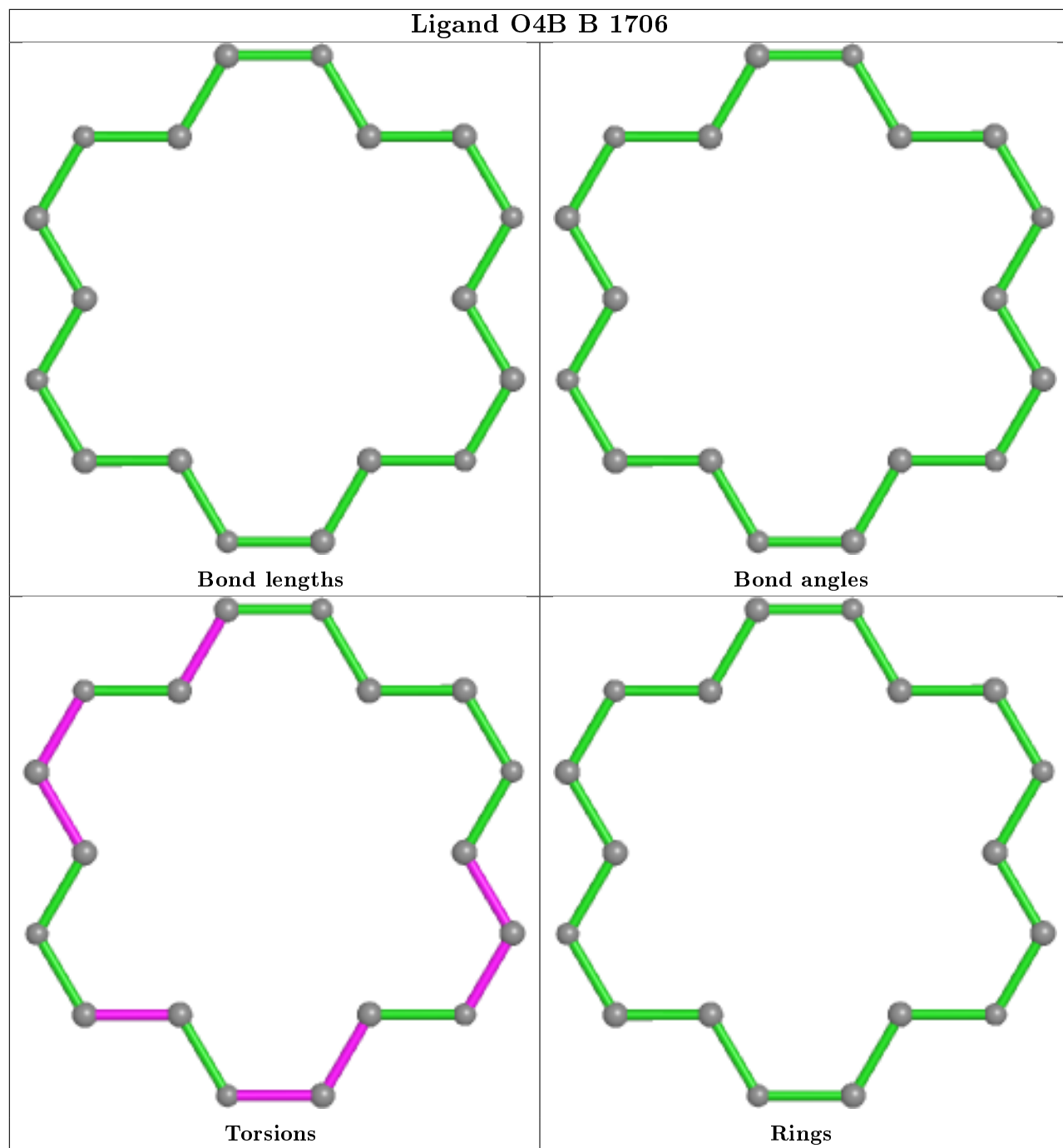
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1709	GOL	3	0
3	D	1704	O4B	9	0
3	A	1706	O4B	6	0
3	B	1706	O4B	3	0
4	A	1707	GOL	2	0
3	D	1703	O4B	2	0
3	A	1705	O4B	3	0
4	B	1708	GOL	1	0
3	C	1706	O4B	1	0
4	B	1707	GOL	2	0
4	C	1707	GOL	1	0

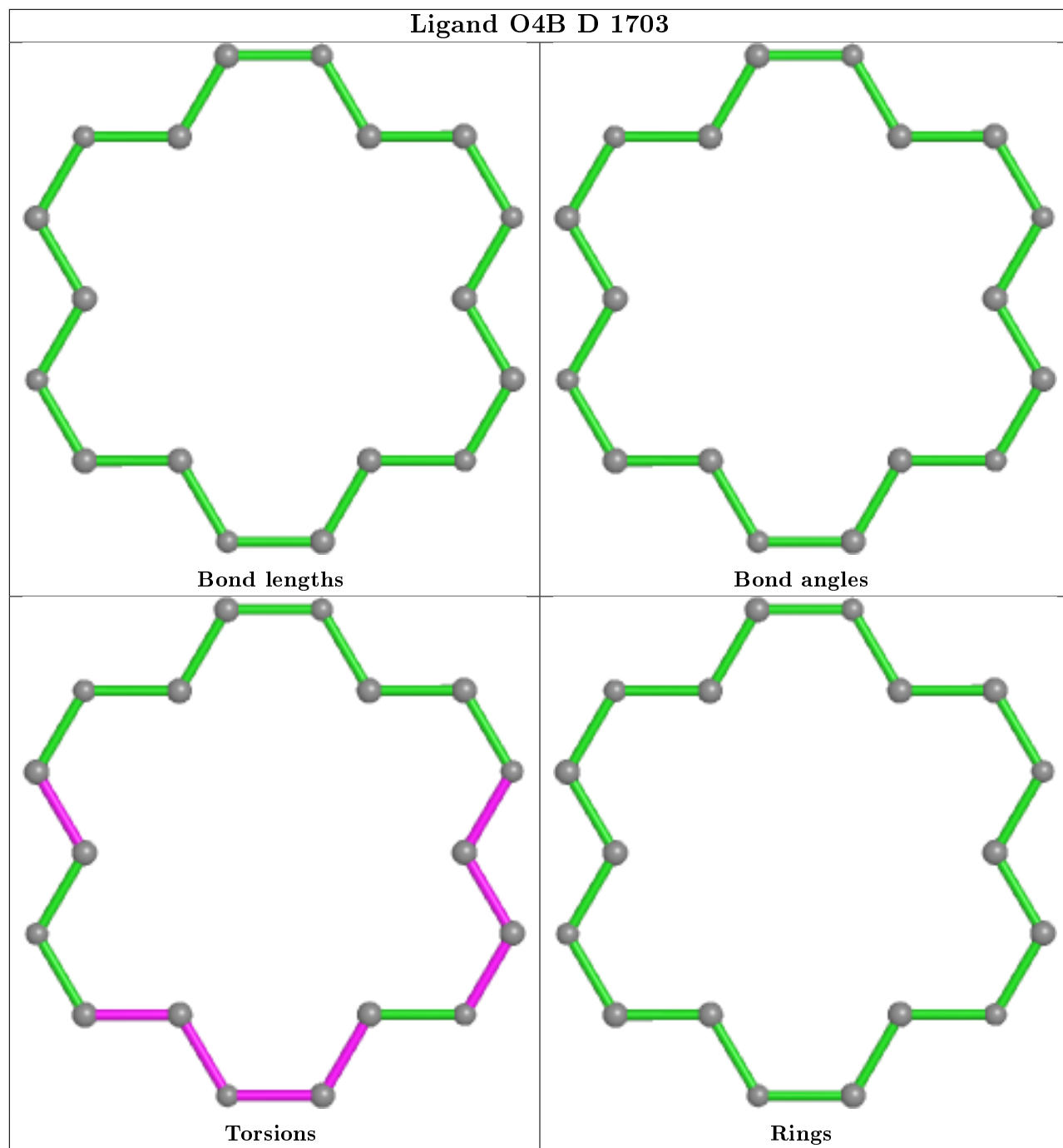
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

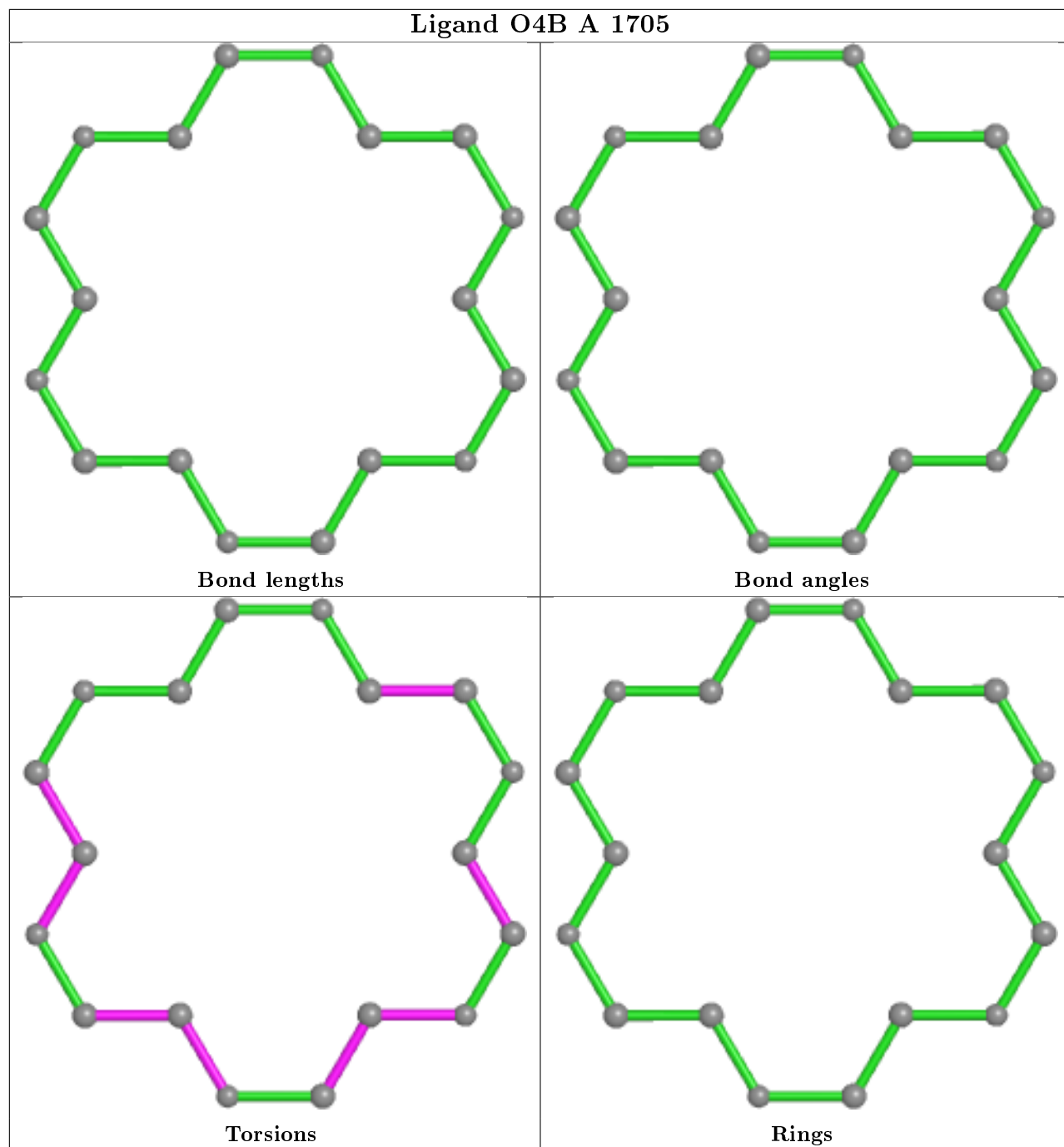
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD 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.

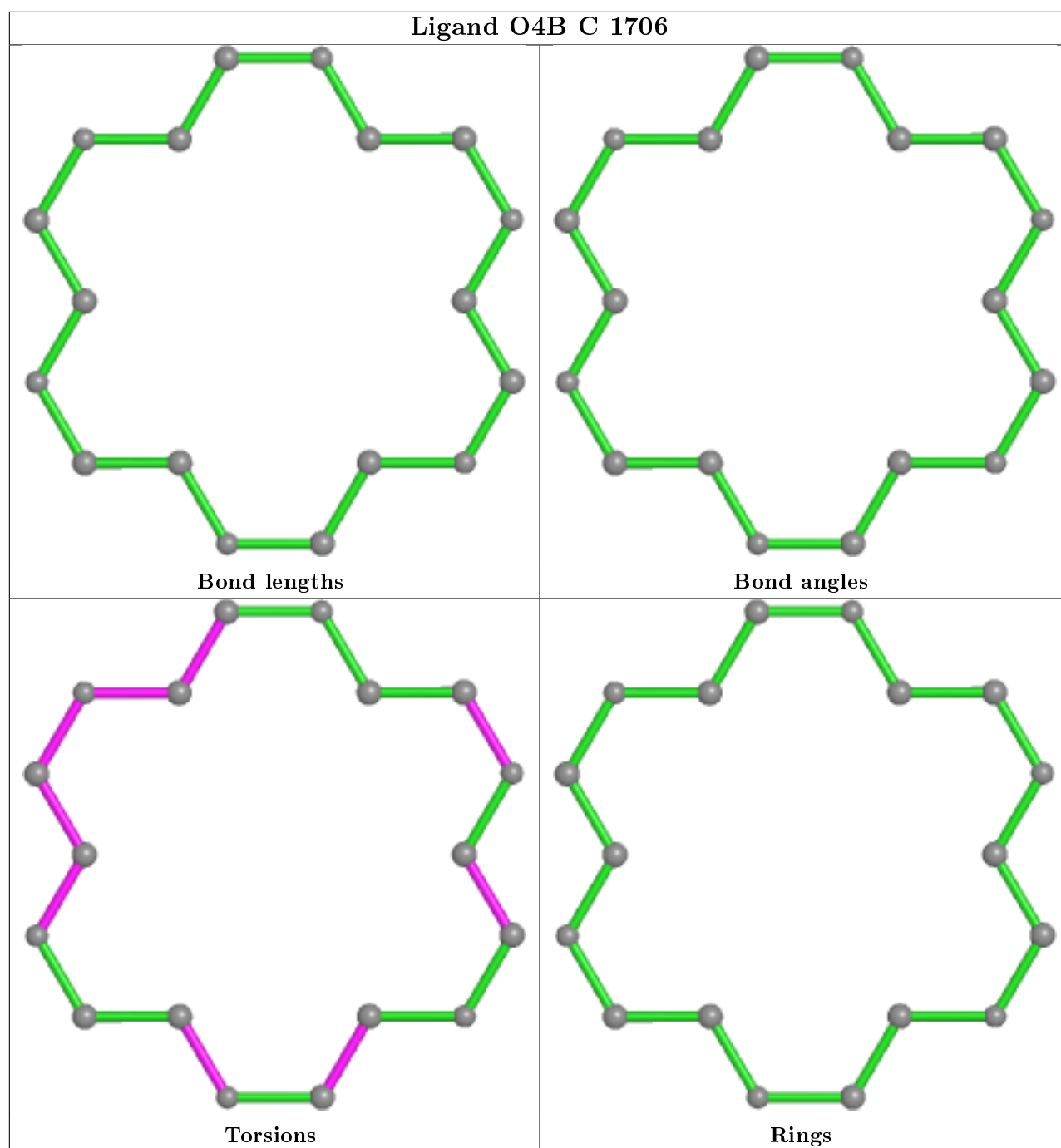












5.7 Other polymers [i](#)

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 ⓘ

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	125/131 (95%)	0.07	4 (3%)	47	45	12, 17, 31, 53	3 (2%)
1	B	125/131 (95%)	-0.01	4 (3%)	47	45	10, 18, 33, 52	6 (4%)
1	C	126/131 (96%)	-0.05	3 (2%)	59	57	10, 18, 33, 47	6 (4%)
1	D	126/131 (96%)	0.05	5 (3%)	38	36	12, 18, 32, 56	6 (4%)
All	All	502/524 (95%)	0.02	16 (3%)	47	45	10, 18, 34, 56	21 (4%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1684	THR	7.8
1	A	1683	ALA	6.8
1	D	1683	ALA	6.3
1	D	1684	THR	6.3
1	B	1559	ILE	5.0
1	B	1683	ALA	4.8
1	A	1559	ILE	4.5
1	C	1559	ILE	4.2
1	D	1559	ILE	3.9
1	A	1682	ALA	3.5
1	D	1560	GLY	3.3
1	B	1682	ALA	2.9
1	C	1571	ASN	2.8
1	D	1682	ALA	2.7
1	A	1571	ASN	2.6
1	B	1560	GLY	2.1

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 carbohydrates 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.

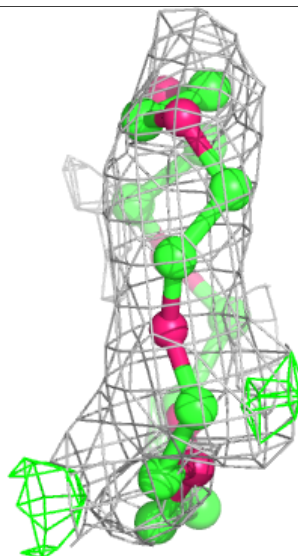
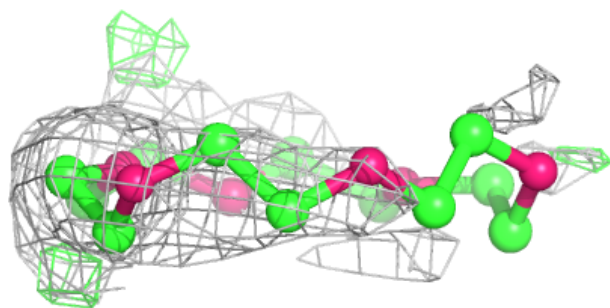
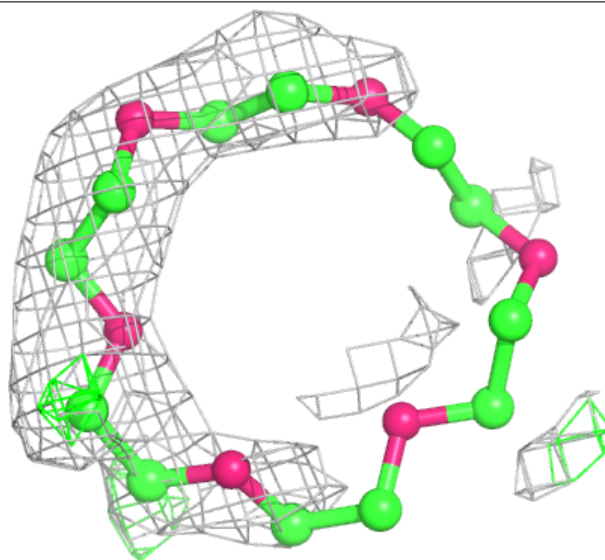
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	GOL	B	1707	6/6	0.67	0.24	35,37,41,41	0
4	GOL	A	1707	6/6	0.72	0.20	26,36,37,38	0
4	GOL	C	1707	6/6	0.75	0.25	34,43,44,45	0
3	O4B	A	1706	18/18	0.78	0.31	35,53,60,61	3
4	GOL	B	1709	6/6	0.80	0.25	29,30,39,41	0
4	GOL	A	1708	6/6	0.81	0.17	33,44,56,56	0
3	O4B	D	1704	18/18	0.82	0.28	33,52,60,61	3
3	O4B	B	1706	18/18	0.83	0.22	36,43,60,61	0
3	O4B	C	1706	18/18	0.84	0.23	28,44,59,64	0
4	GOL	A	1709	6/6	0.85	0.27	27,34,41,42	0
2	SO4	C	1705	5/5	0.86	0.15	38,43,55,58	0
4	GOL	B	1708	6/6	0.87	0.15	35,36,40,41	0
3	O4B	A	1705	18/18	0.88	0.17	31,41,51,55	0
3	O4B	D	1703	18/18	0.88	0.21	30,45,56,61	0
2	SO4	C	1703	5/5	0.91	0.26	49,51,55,57	0
2	SO4	B	1704	5/5	0.94	0.18	32,34,49,54	0
2	SO4	B	1703	5/5	0.95	0.17	42,45,51,57	0
2	SO4	C	1704	5/5	0.95	0.18	34,35,48,60	0
2	SO4	A	1701	5/5	0.95	0.12	31,38,45,48	0
2	SO4	B	1705	5/5	0.95	0.23	28,30,41,47	0
2	SO4	C	1702	5/5	0.96	0.20	30,31,39,45	0
2	SO4	A	1704	5/5	0.97	0.18	42,44,46,56	0
2	SO4	D	1702	5/5	0.97	0.15	33,36,45,50	0
2	SO4	A	1702	5/5	0.98	0.12	31,36,38,46	0
2	SO4	B	1702	5/5	0.98	0.12	28,30,37,38	0
2	SO4	D	1701	5/5	0.98	0.12	29,34,40,42	0
2	SO4	B	1701	5/5	0.99	0.11	16,17,19,20	0
2	SO4	C	1701	5/5	0.99	0.13	27,28,36,36	0
2	SO4	A	1703	5/5	0.99	0.10	16,17,19,22	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.

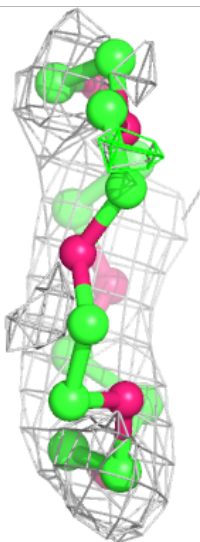
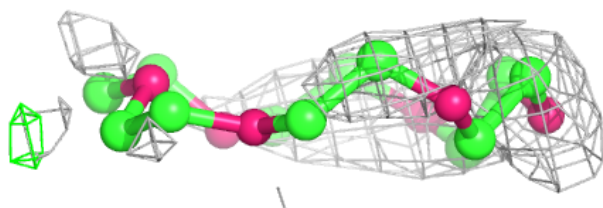
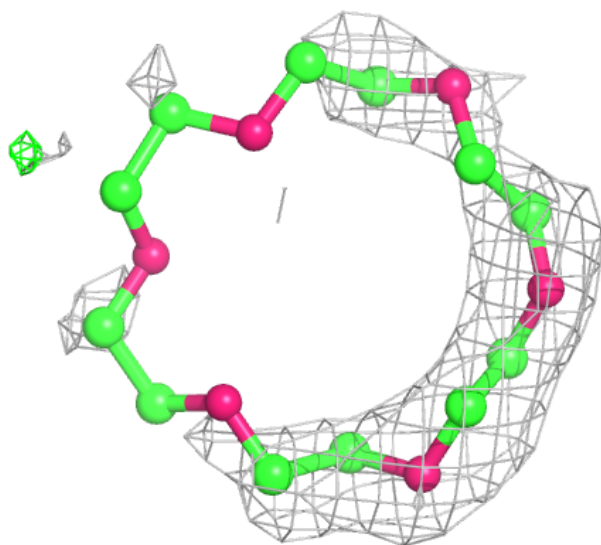
Electron density around O4B A 1706:

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



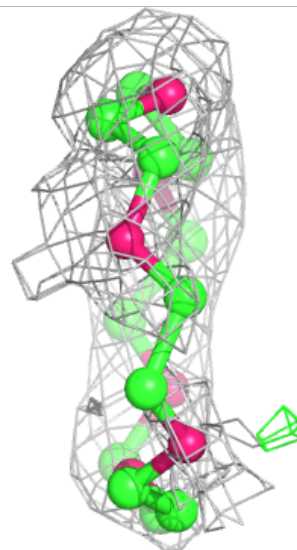
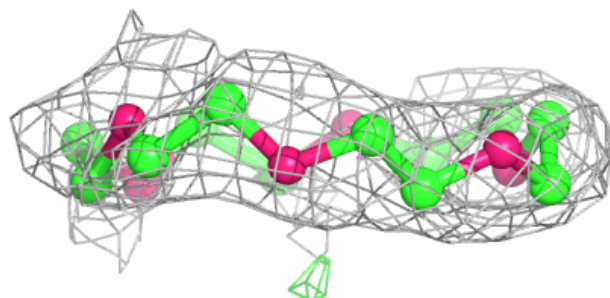
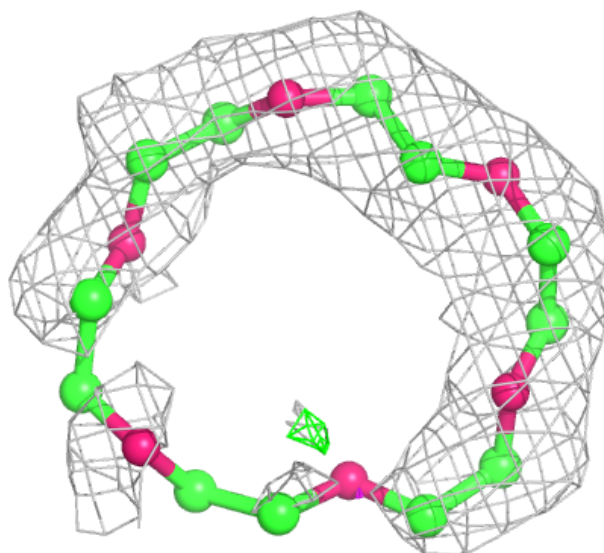
Electron density around O4B D 1704:

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



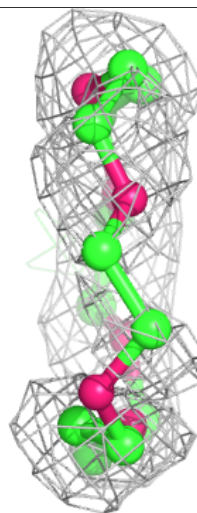
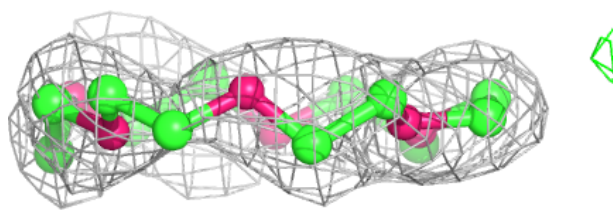
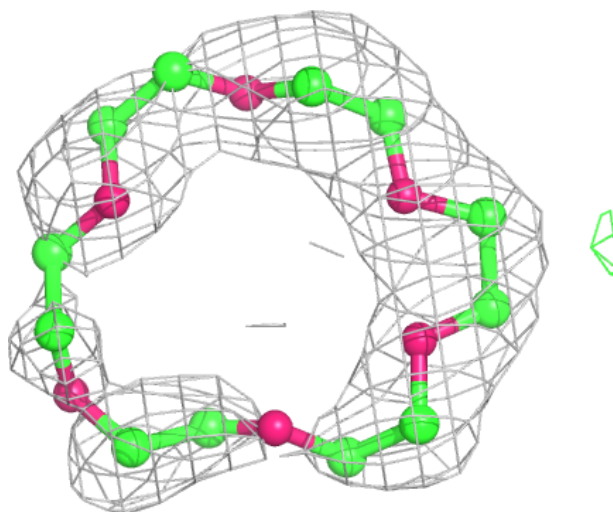
Electron density around O4B B 1706:

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



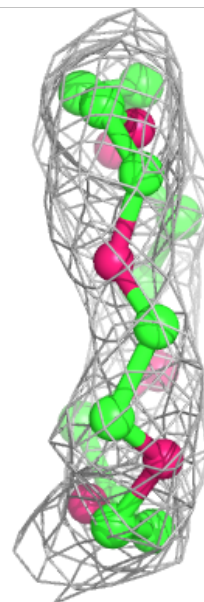
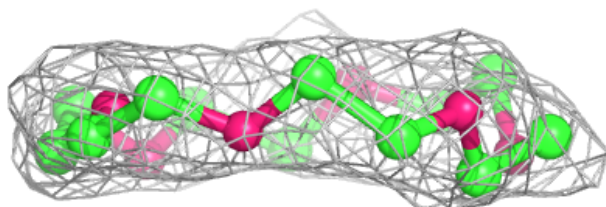
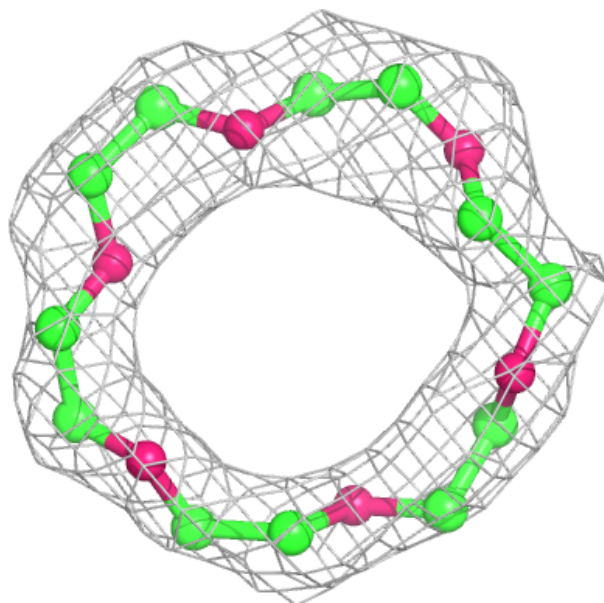
Electron density around O4B C 1706:

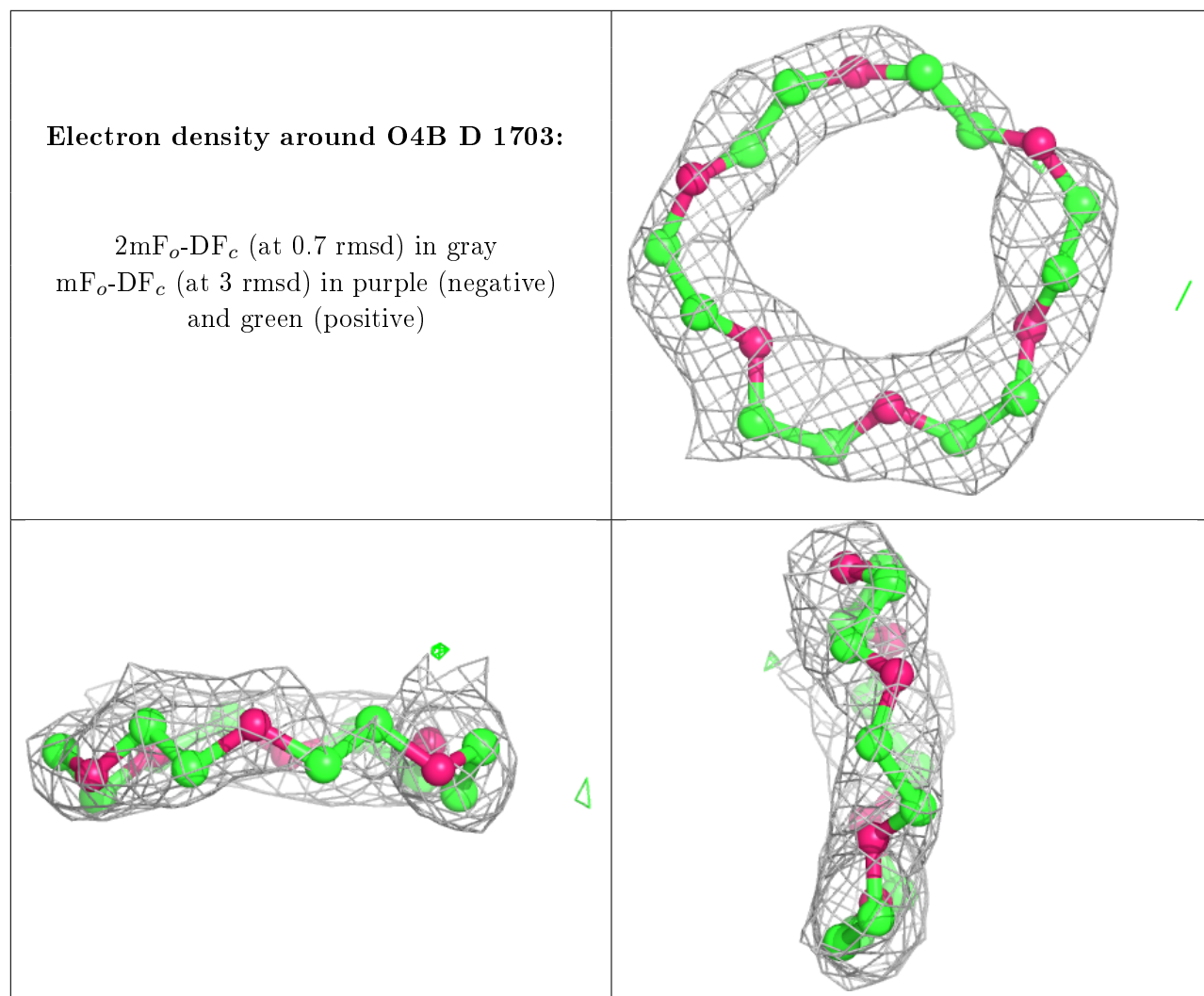
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around O4B A 1705:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.