



# Full wwPDB X-ray Structure Validation Report ⓘ

May 5, 2025 – 01:39 pm BST

PDB ID : 9F3I / pdb\_00009f3i  
Title : Crystal structure of Kluyveromyces lactis glucokinase in complex with glucose  
Authors : Weisse, R.H.; Strater, N.  
Deposited on : 2024-04-25  
Resolution : 2.65 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](mailto: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>.

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

MolProbity : 4-5-2 with Phenix2.0rc1  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.003 (Gargrove)  
Density-Fitness : 1.0.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.43.1

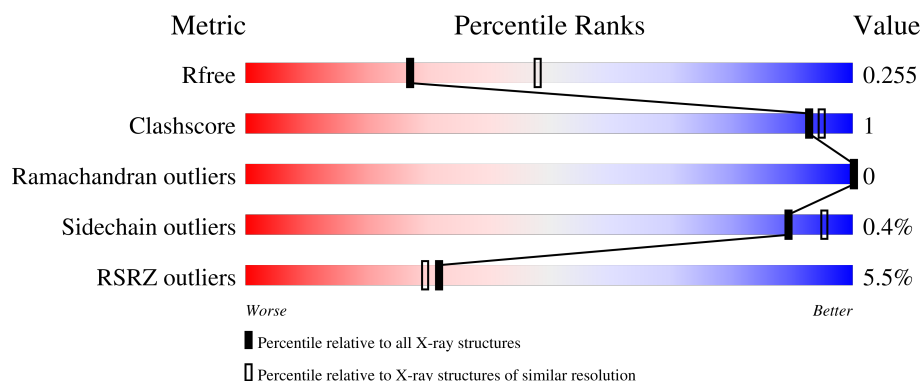
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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))
$R_{free}$	164625	1003 (2.66-2.66)
Clashscore	180529	1063 (2.66-2.66)
Ramachandran outliers	177936	1052 (2.66-2.66)
Sidechain outliers	177891	1052 (2.66-2.66)
RSRZ outliers	164620	1003 (2.66-2.66)

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  $\geq 3$ , 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	481	<div> <div>3%</div> <div>97%</div> </div>
1	B	481	<div> <div>8%</div> <div>96%</div> </div>
1	C	481	<div> <div>8%</div> <div>95%</div> </div>
1	D	481	<div> <div>3%</div> <div>96%</div> </div>

## 2 Entry composition [i](#)

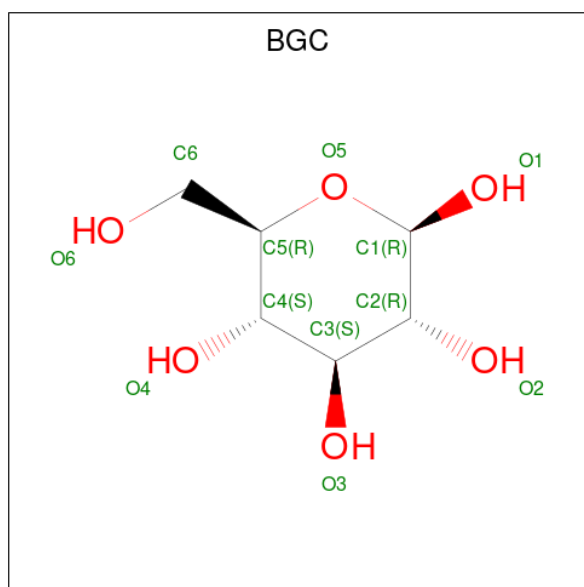
There are 4 unique types of molecules in this entry. The entry contains 30318 atoms, of which 15007 are hydrogens and 0 are deuteriums.

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 Glucokinase-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	480	Total	C	H	N	O	S	0	0	0
			7497	2381	3728	641	727	20			
1	B	480	Total	C	H	N	O	S	0	0	0
			7497	2381	3728	641	727	20			
1	C	480	Total	C	H	N	O	S	0	2	0
			7527	2389	3745	644	729	20			
1	D	480	Total	C	H	N	O	S	0	0	0
			7497	2381	3728	641	727	20			

- Molecule 2 is beta-D-glucopyranose (CCD ID: BGC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>) (labeled as "Ligand of Interest" by depositor).



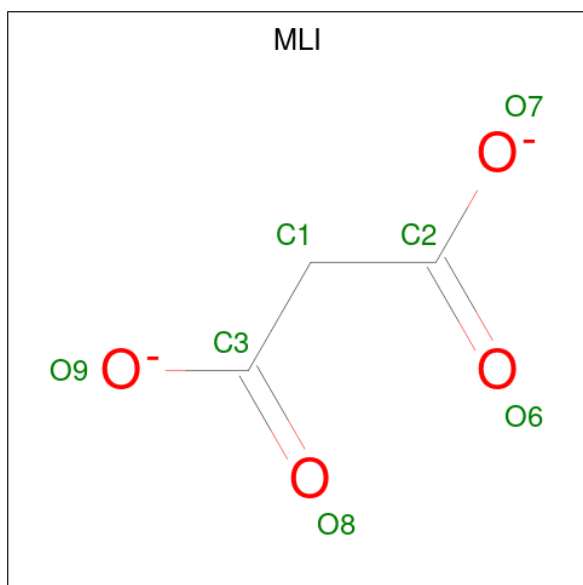
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			24	6	12	6		
2	B	1	Total	C	H	O	0	0
			24	6	12	6		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	H	O	0	0
			24	6	12	6		
2	D	1	Total	C	H	O	0	0
			24	6	12	6		

- Molecule 3 is MALONATE ION (CCD ID: MLI) (formula:  $C_3H_2O_4$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			9	3	2	4		
3	A	1	Total	C	H	O	0	0
			9	3	2	4		
3	A	1	Total	C	H	O	0	0
			9	3	2	4		
3	B	1	Total	C	H	O	0	0
			9	3	2	4		
3	B	1	Total	C	H	O	0	0
			9	3	2	4		
3	B	1	Total	C	H	O	0	0
			9	3	2	4		
3	C	1	Total	C	H	O	0	0
			9	3	2	4		
3	C	1	Total	C	H	O	0	0
			9	3	2	4		
3	C	1	Total	C	H	O	0	0
			9	3	2	4		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	H	O	0	0
			9	3	2	4		
3	D	1	Total	C	H	O	0	0
			9	3	2	4		
3	D	1	Total	C	H	O	0	0
			9	3	2	4		
3	D	1	Total	C	H	O	0	0
			9	3	2	4		
3	D	1	Total	C	H	O	0	0
			9	3	2	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	20	Total	O	0	0
			20	20		
4	B	18	Total	O	0	0
			18	18		
4	C	9	Total	O	0	0
			9	9		
4	D	22	Total	O	0	0
			22	22		

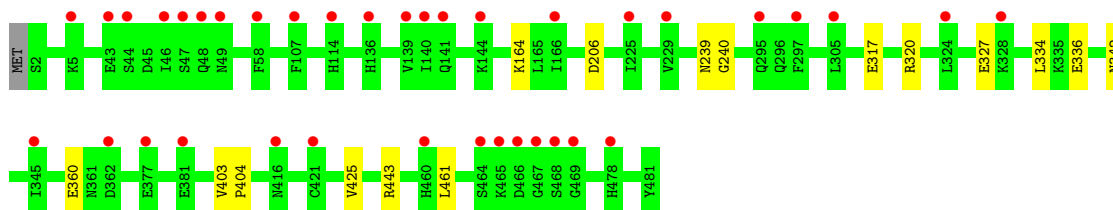
### 3 Residue-property plots [i](#)

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.

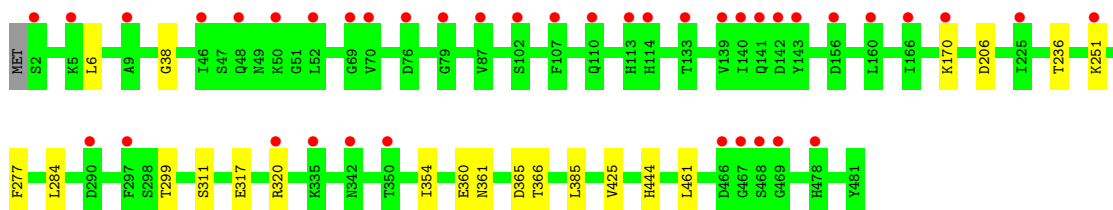
- Molecule 1: Glucokinase-1



- Molecule 1: Glucokinase-1



- Molecule 1: Glucokinase-1



- Molecule 1: Glucokinase-1



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	203.78Å 88.11Å 212.03Å 90.00° 90.10° 90.00°	Depositor
Resolution (Å)	64.72 – 2.65 64.72 – 2.65	Depositor EDS
% Data completeness (in resolution range)	97.8 (64.72-2.65) 98.2 (64.72-2.65)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.36 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.235 , 0.258 0.234 , 0.255	Depositor DCC
$R_{free}$ test set	5398 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	61.6	Xtriage
Anisotropy	0.412	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 60.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.027 for l,k,-h 0.011 for l,-k,h 0.014 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	30318	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MLI, BGC

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	0.12	0/3843	0.27	0/5199
1	B	0.11	0/3843	0.26	0/5199
1	C	0.10	0/3856	0.25	0/5215
1	D	0.16	0/3843	0.30	0/5199
All	All	0.12	0/15385	0.27	0/20812

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	3769	3728	3726	7	0
1	B	3769	3728	3726	9	0
1	C	3782	3745	3741	12	0
1	D	3769	3728	3726	13	0
2	A	12	12	12	0	0
2	B	12	12	12	0	0
2	C	12	12	12	0	0
2	D	12	12	12	0	0
3	A	21	6	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	21	6	6	2	0
3	C	28	8	8	2	0
3	D	35	10	10	3	0
4	A	20	0	0	0	0
4	B	18	0	0	2	0
4	C	9	0	0	1	0
4	D	22	0	0	0	0
All	All	15311	15007	14997	42	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 1.

All (42) 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:236:THR:O	1:C:311:SER:OG	2.04	0.74
1:D:53:PRO:O	1:D:251:LYS:NZ	2.32	0.62
1:C:170[B]:LYS:O	1:C:299:THR:OG1	2.16	0.59
3:B:503:MLI:O9	4:B:601:HOH:O	2.17	0.58
1:B:317:GLU:OE2	1:B:320:ARG:NH1	2.37	0.57
3:C:505:MLI:O7	1:D:457:ARG:NH2	2.39	0.55
1:D:23:LYS:NZ	1:D:283:HIS:O	2.39	0.55
1:D:248:ASN:ND2	3:D:506:MLI:O9	2.40	0.54
1:D:236:THR:OG1	3:D:503:MLI:O6	2.23	0.54
1:C:206:ASP:OD1	1:C:206:ASP:N	2.44	0.51
1:D:206:ASP:OD1	1:D:206:ASP:N	2.44	0.50
1:A:206:ASP:OD1	1:A:206:ASP:N	2.41	0.49
1:A:38:GLY:O	1:A:251:LYS:NZ	2.34	0.49
1:C:354:ILE:HG12	1:C:366:THR:HG23	1.96	0.48
1:B:425:VAL:HB	1:B:461:LEU:HD23	1.96	0.48
1:D:312:GLY:O	1:D:348:THR:OG1	2.26	0.47
1:D:425:VAL:HB	1:D:461:LEU:HD23	1.97	0.47
1:C:425:VAL:HB	1:C:461:LEU:HD23	1.97	0.47
1:A:420:HIS:ND1	3:B:502:MLI:O9	2.48	0.47
1:D:444:HIS:ND1	3:D:502:MLI:O9	2.39	0.46
1:B:327:GLU:OE1	1:B:334:LEU:HD12	2.16	0.46
1:C:38:GLY:O	1:C:251:LYS:NZ	2.42	0.45
1:B:239:ASN:OD1	1:B:240:GLY:N	2.50	0.45
1:C:317:GLU:OE2	1:C:320:ARG:NH1	2.50	0.44
1:B:336:GLU:N	1:B:336:GLU:OE1	2.51	0.44
1:B:360:GLU:OE1	1:B:360:GLU:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:477:LEU:C	1:D:477:LEU:HD23	2.43	0.44
1:B:206:ASP:OD1	1:B:206:ASP:N	2.46	0.43
1:B:443:ARG:NH2	4:B:602:HOH:O	2.37	0.43
1:A:409:LEU:HD13	1:A:415:LEU:HD21	2.00	0.42
1:D:317:GLU:O	1:D:321:ASN:ND2	2.48	0.42
1:D:439:ARG:HD2	1:D:463:ILE:HD11	2.01	0.42
1:C:6:LEU:HG	1:C:385:LEU:HD13	2.01	0.42
1:C:277:PHE:O	1:C:284:LEU:HD11	2.19	0.42
1:C:444:HIS:ND1	3:C:502:MLI:O7	2.51	0.42
1:A:403:VAL:HB	1:A:404:PRO:HD3	2.02	0.42
1:A:274:TRP:C	1:A:274:TRP:CD1	2.98	0.42
1:C:236:THR:HB	4:C:602:HOH:O	2.20	0.41
1:D:277:PHE:O	1:D:284:LEU:HD11	2.20	0.41
1:A:239:ASN:OD1	1:A:240:GLY:N	2.49	0.40
1:B:403:VAL:HB	1:B:404:PRO:HD3	2.03	0.40
1:C:360:GLU:OE1	1:C:360:GLU:N	2.52	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	478/481 (99%)	457 (96%)	21 (4%)	0	100	100
1	B	478/481 (99%)	459 (96%)	19 (4%)	0	100	100
1	C	480/481 (100%)	458 (95%)	22 (5%)	0	100	100
1	D	478/481 (99%)	457 (96%)	21 (4%)	0	100	100
All	All	1914/1924 (100%)	1831 (96%)	83 (4%)	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	421/422 (100%)	419 (100%)	2 (0%)	86	94
1	B	421/422 (100%)	419 (100%)	2 (0%)	86	94
1	C	422/422 (100%)	420 (100%)	2 (0%)	86	94
1	D	421/422 (100%)	421 (100%)	0	100	100
All	All	1685/1688 (100%)	1679 (100%)	6 (0%)	89	95

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	220	ASP
1	A	342	ASN
1	B	164	LYS
1	B	342	ASN
1	C	361	ASN
1	C	365	ASP

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	99	GLN
1	A	283	HIS
1	A	329	GLN
1	A	422	GLN
1	A	434	HIS
1	A	460	HIS
1	B	26	GLN
1	B	188	GLN
1	B	413	ASN
1	B	416	ASN
1	B	434	HIS
1	C	361	ASN
1	C	364	GLN

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Mol	Chain	Res	Type
1	C	434	HIS
1	D	92	ASN
1	D	180	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](#)

19 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	BGC	A	501	-	12,12,12	0.23	0	17,17,17	0.38	0
3	MLI	C	504	-	6,6,6	1.34	0	7,7,7	0.95	0
2	BGC	D	501	-	12,12,12	0.17	0	17,17,17	0.45	0
3	MLI	C	502	-	6,6,6	1.36	0	7,7,7	0.91	0
3	MLI	B	504	-	6,6,6	1.35	0	7,7,7	0.94	0
3	MLI	D	505	-	6,6,6	1.33	0	7,7,7	0.98	0
3	MLI	C	503	-	6,6,6	1.33	0	7,7,7	0.98	0
3	MLI	A	502	-	6,6,6	1.35	0	7,7,7	0.95	0
3	MLI	D	506	-	6,6,6	1.29	0	7,7,7	0.96	0
3	MLI	D	503	-	6,6,6	1.32	0	7,7,7	0.97	0
3	MLI	C	505	-	6,6,6	1.36	0	7,7,7	0.96	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	MLI	D	504	-	6,6,6	1.34	0	7,7,7	0.93	0
2	BGC	B	501	-	12,12,12	0.18	0	17,17,17	0.45	0
3	MLI	A	504	-	6,6,6	1.32	0	7,7,7	0.98	0
3	MLI	B	502	-	6,6,6	1.34	0	7,7,7	0.95	0
2	BGC	C	501	-	12,12,12	0.21	0	17,17,17	0.31	0
3	MLI	B	503	-	6,6,6	1.34	0	7,7,7	0.92	0
3	MLI	A	503	-	6,6,6	1.39	0	7,7,7	0.92	0
3	MLI	D	502	-	6,6,6	1.37	0	7,7,7	0.96	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
2	BGC	A	501	-	-	0/2/22/22	0/1/1/1
3	MLI	C	504	-	-	2/4/4/4	-
2	BGC	D	501	-	-	0/2/22/22	0/1/1/1
3	MLI	C	502	-	-	0/4/4/4	-
3	MLI	B	504	-	-	0/4/4/4	-
3	MLI	D	505	-	-	0/4/4/4	-
3	MLI	C	503	-	-	0/4/4/4	-
3	MLI	A	502	-	-	2/4/4/4	-
3	MLI	D	506	-	-	0/4/4/4	-
3	MLI	D	503	-	-	0/4/4/4	-
3	MLI	C	505	-	-	2/4/4/4	-
3	MLI	D	504	-	-	4/4/4/4	-
2	BGC	B	501	-	-	0/2/22/22	0/1/1/1
3	MLI	A	504	-	-	2/4/4/4	-
3	MLI	B	502	-	-	2/4/4/4	-
2	BGC	C	501	-	-	0/2/22/22	0/1/1/1
3	MLI	B	503	-	-	1/4/4/4	-
3	MLI	A	503	-	-	0/4/4/4	-
3	MLI	D	502	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (18) torsion outliers are listed below:

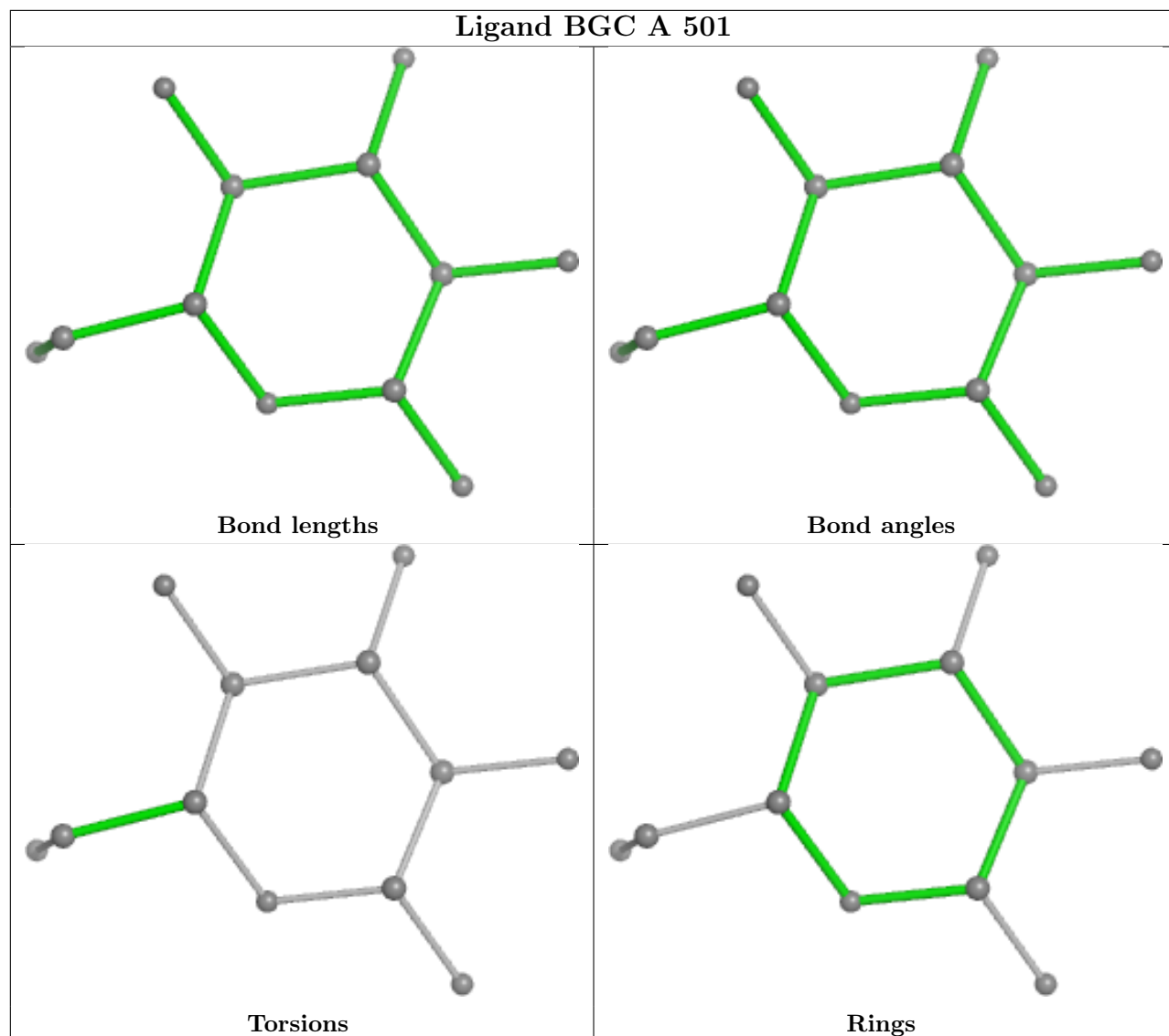
Mol	Chain	Res	Type	Atoms
3	D	504	MLI	C3-C1-C2-O6
3	D	504	MLI	C3-C1-C2-O7
3	A	502	MLI	C2-C1-C3-O9
3	B	502	MLI	C3-C1-C2-O7
3	D	502	MLI	C3-C1-C2-O6
3	D	502	MLI	C3-C1-C2-O7
3	D	504	MLI	C2-C1-C3-O8
3	D	504	MLI	C2-C1-C3-O9
3	B	502	MLI	C3-C1-C2-O6
3	C	505	MLI	C3-C1-C2-O6
3	A	502	MLI	C2-C1-C3-O8
3	C	505	MLI	C3-C1-C2-O7
3	C	504	MLI	C3-C1-C2-O7
3	A	504	MLI	C2-C1-C3-O8
3	A	504	MLI	C2-C1-C3-O9
3	C	504	MLI	C3-C1-C2-O6
3	B	503	MLI	C3-C1-C2-O7
3	D	502	MLI	C2-C1-C3-O9

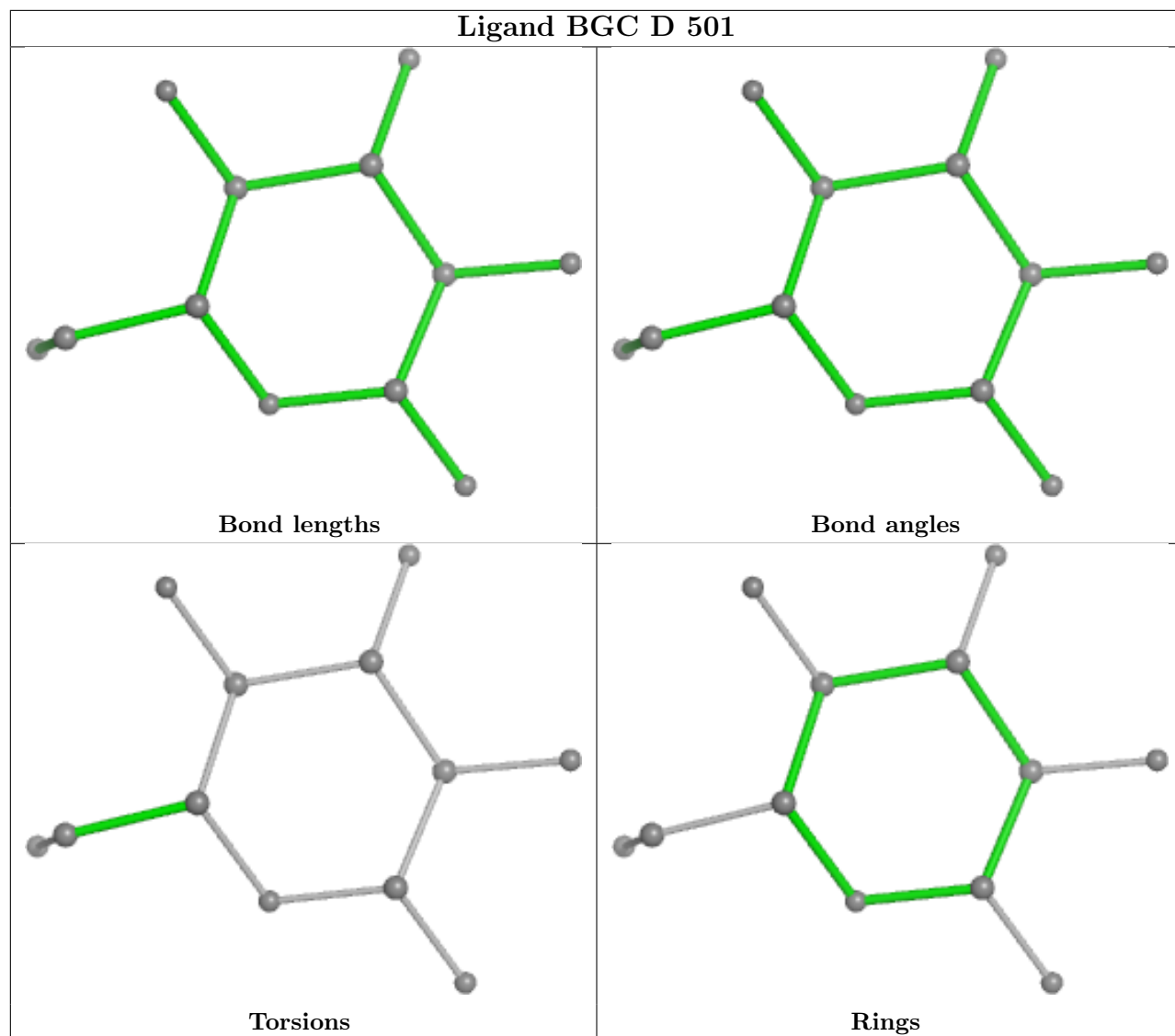
There are no ring outliers.

7 monomers are involved in 7 short contacts:

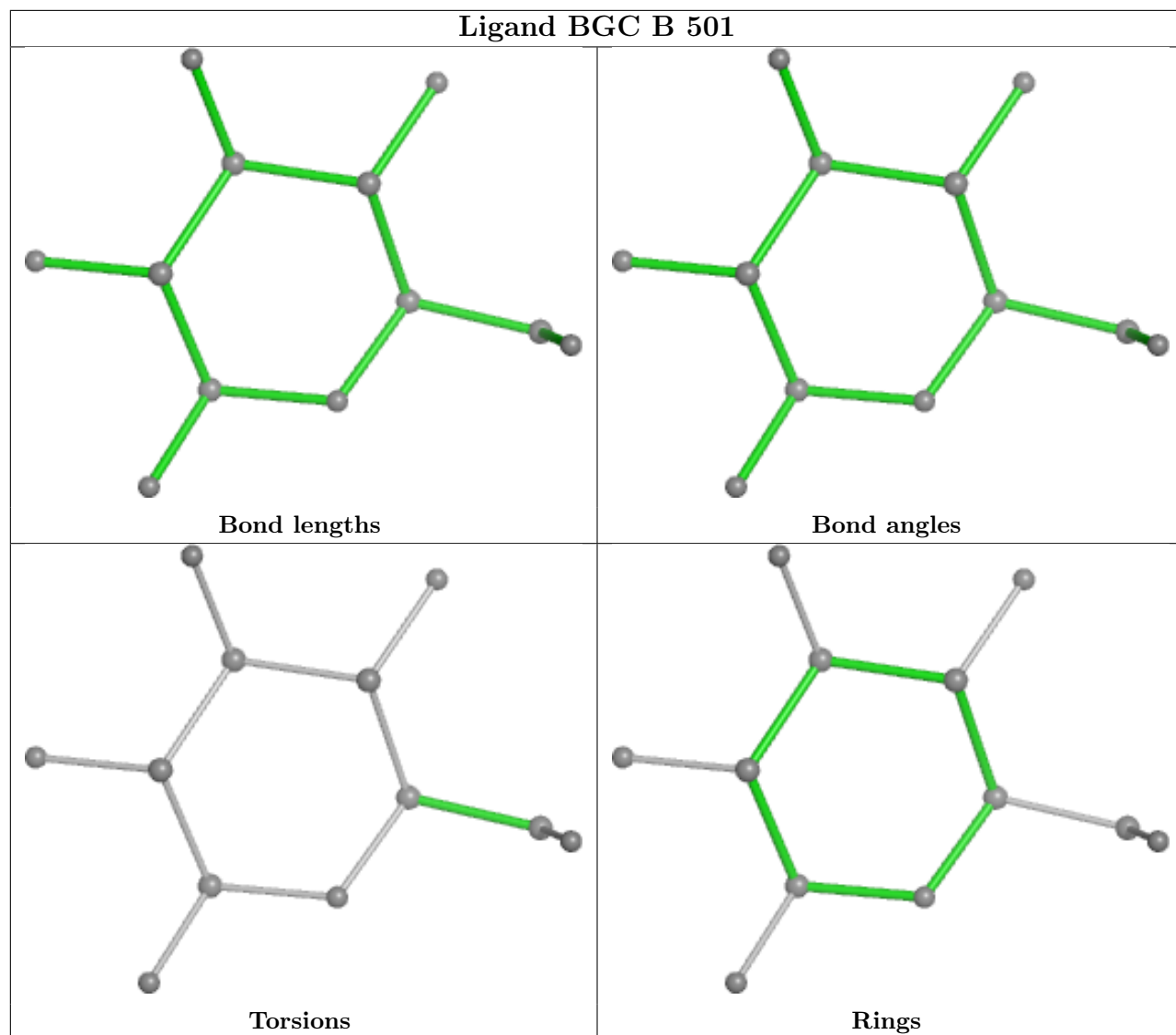
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	502	MLI	1	0
3	D	506	MLI	1	0
3	D	503	MLI	1	0
3	C	505	MLI	1	0
3	B	502	MLI	1	0
3	B	503	MLI	1	0
3	D	502	MLI	1	0

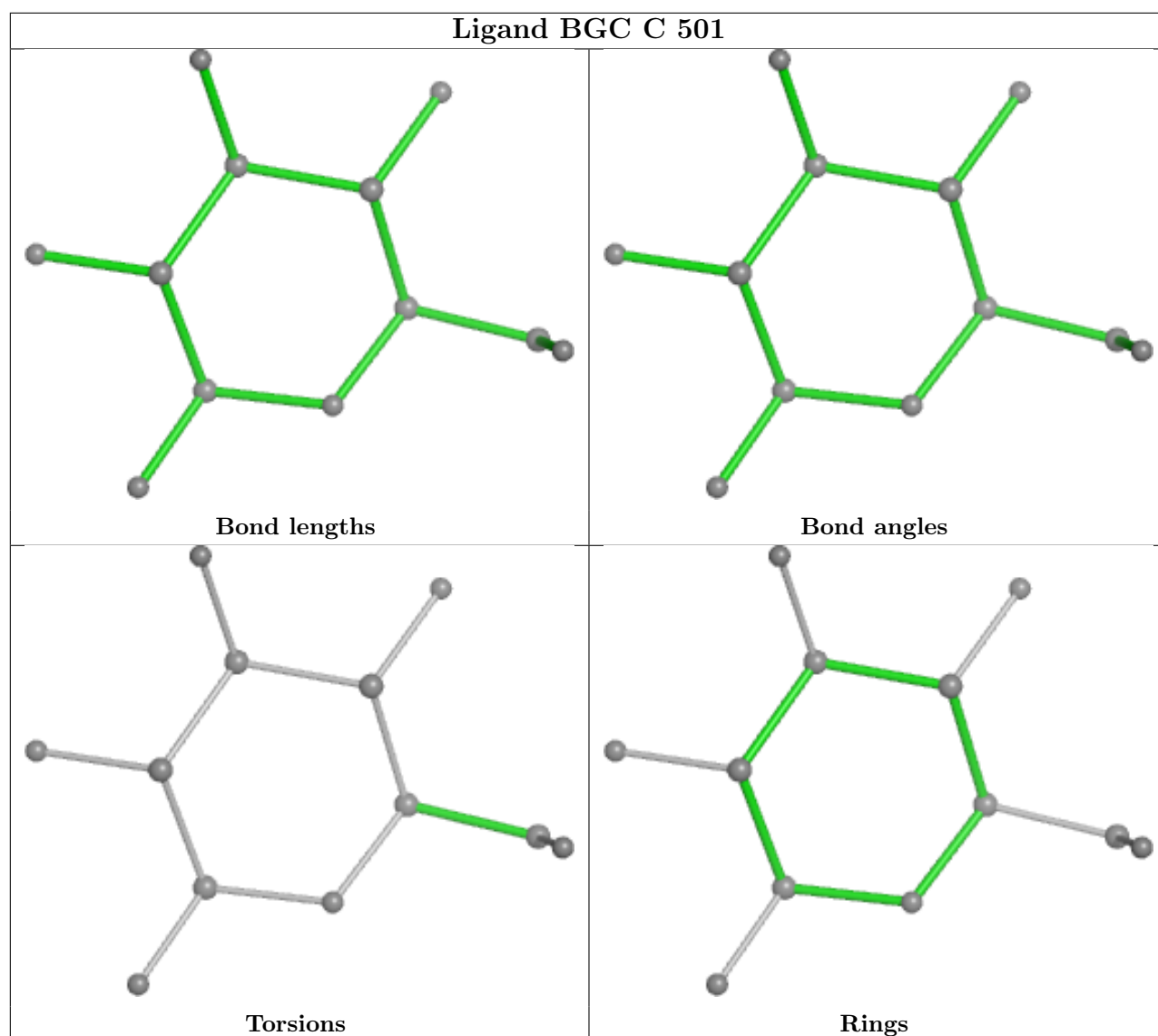
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	480/481 (99%)	0.46	15 (3%)	51	49	59, 80, 122, 155	0
1	B	480/481 (99%)	0.76	37 (7%)	21	19	68, 101, 139, 178	0
1	C	480/481 (99%)	0.81	40 (8%)	19	17	53, 107, 144, 175	2 (0%)
1	D	480/481 (99%)	0.20	14 (2%)	54	52	56, 75, 113, 159	0
All	All	1920/1924 (99%)	0.56	106 (5%)	32	29	53, 91, 132, 178	2 (0%)

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	139	VAL	5.0
1	B	140	ILE	4.7
1	A	415	LEU	4.6
1	C	107	PHE	4.4
1	D	467	GLY	4.3
1	B	466	ASP	4.3
1	C	2	SER	4.2
1	C	467	GLY	4.1
1	B	467	GLY	3.9
1	C	114	HIS	3.8
1	C	113	HIS	3.8
1	B	464	SER	3.5
1	C	166	ILE	3.4
1	C	320	ARG	3.4
1	C	46	ILE	3.4
1	B	468	SER	3.3
1	D	466	ASP	3.3
1	D	465	LYS	3.2
1	A	466	ASP	3.2
1	B	377	GLU	3.2
1	A	377	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	46	ILE	3.1
1	C	142	ASP	3.1
1	B	362	ASP	3.1
1	C	50	LYS	3.1
1	A	225	ILE	3.0
1	B	465	LYS	3.0
1	B	469	GLY	2.9
1	B	43	GLU	2.9
1	D	356	ILE	2.8
1	C	79	GLY	2.8
1	C	350	THR	2.8
1	B	141	GLN	2.8
1	C	48	GLN	2.8
1	C	76	ASP	2.8
1	A	139	VAL	2.7
1	C	140	ILE	2.7
1	D	336	GLU	2.7
1	C	170[A]	LYS	2.7
1	B	47	SER	2.7
1	B	166	ILE	2.7
1	B	139	VAL	2.6
1	D	139	VAL	2.6
1	A	2	SER	2.6
1	B	5	LYS	2.5
1	A	222	ALA	2.5
1	B	381	GLU	2.5
1	B	295	GLN	2.5
1	A	82	PHE	2.5
1	D	142	ASP	2.5
1	B	114	HIS	2.5
1	C	52	LEU	2.5
1	C	468	SER	2.4
1	C	141	GLN	2.4
1	B	345	ILE	2.4
1	C	469	GLY	2.4
1	A	413	ASN	2.4
1	C	69	GLY	2.4
1	D	377	GLU	2.4
1	B	305	LEU	2.4
1	C	70	VAL	2.4
1	D	478	HIS	2.4
1	C	225	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	133	THR	2.3
1	B	225	ILE	2.3
1	D	292	LYS	2.3
1	C	290	ASP	2.3
1	B	49	ASN	2.2
1	C	478	HIS	2.2
1	B	297	PHE	2.2
1	B	136	HIS	2.2
1	B	421	CYS	2.2
1	C	9	ALA	2.2
1	C	335	LYS	2.2
1	A	467	GLY	2.2
1	D	47	SER	2.2
1	B	48	GLN	2.2
1	B	107	PHE	2.2
1	C	251	LYS	2.2
1	A	160	LEU	2.2
1	B	460	HIS	2.2
1	C	156	ASP	2.2
1	C	110	GLN	2.2
1	B	324	LEU	2.2
1	C	160	LEU	2.2
1	B	44	SER	2.1
1	D	2	SER	2.1
1	C	5	LYS	2.1
1	C	297	PHE	2.1
1	C	102	SER	2.1
1	B	229	VAL	2.1
1	A	468	SER	2.1
1	B	144	LYS	2.1
1	D	218	SER	2.1
1	C	342	ASN	2.1
1	A	460	HIS	2.1
1	D	70	VAL	2.1
1	C	143	TYR	2.1
1	A	421	CYS	2.0
1	B	328	LYS	2.0
1	B	478	HIS	2.0
1	B	58	PHE	2.0
1	C	87	VAL	2.0
1	B	416	ASN	2.0
1	C	466	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	46	ILE	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

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

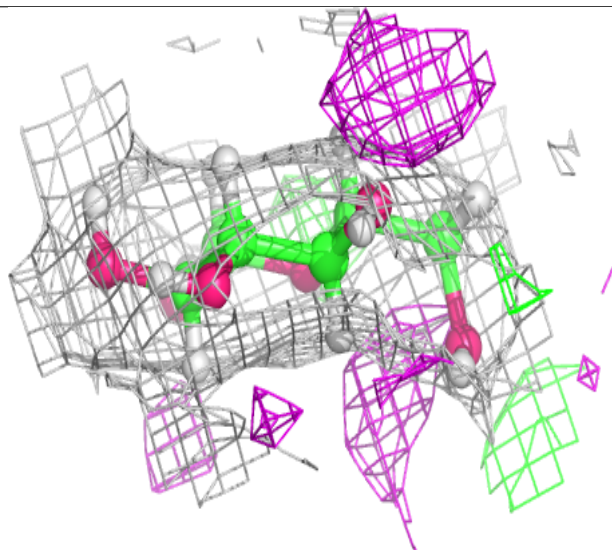
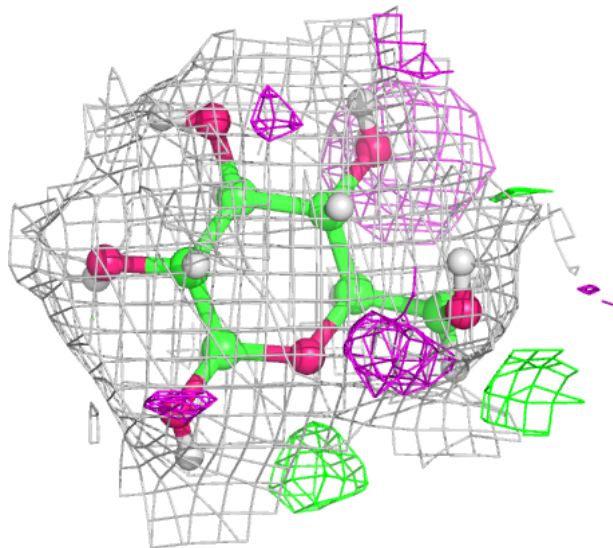
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	MLI	B	503	7/7	0.70	0.20	76,77,93,93	0
3	MLI	C	504	7/7	0.72	0.15	101,104,123,123	0
3	MLI	B	504	7/7	0.75	0.14	91,94,113,113	0
3	MLI	A	503	7/7	0.77	0.21	75,78,94,94	0
3	MLI	C	505	7/7	0.77	0.23	80,82,101,101	0
3	MLI	A	502	7/7	0.82	0.14	63,66,77,77	0
3	MLI	D	506	7/7	0.82	0.14	70,72,84,84	0
3	MLI	D	505	7/7	0.83	0.25	70,72,89,89	0
3	MLI	A	504	7/7	0.84	0.16	69,74,93,93	0
3	MLI	D	503	7/7	0.85	0.13	60,63,76,76	0
3	MLI	D	504	7/7	0.86	0.13	73,75,88,88	0
3	MLI	B	502	7/7	0.89	0.10	72,73,87,87	0
3	MLI	C	502	7/7	0.89	0.11	68,76,92,92	0
3	MLI	C	503	7/7	0.89	0.15	83,87,106,106	0
2	BGC	D	501	12/12	0.90	0.10	46,53,61,64	0
2	BGC	B	501	12/12	0.91	0.10	67,78,90,95	0
3	MLI	D	502	7/7	0.91	0.12	62,69,84,84	0
2	BGC	C	501	12/12	0.91	0.13	76,87,102,105	0
2	BGC	A	501	12/12	0.93	0.10	52,56,66,68	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.

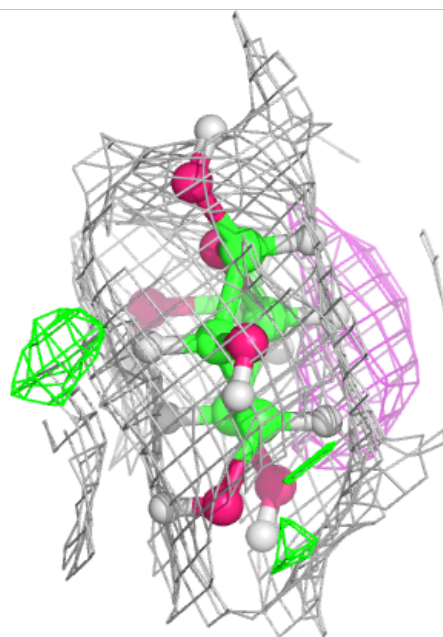
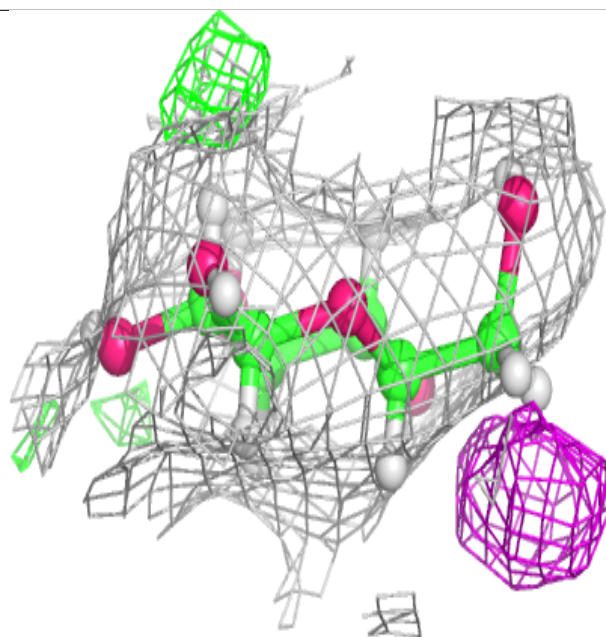
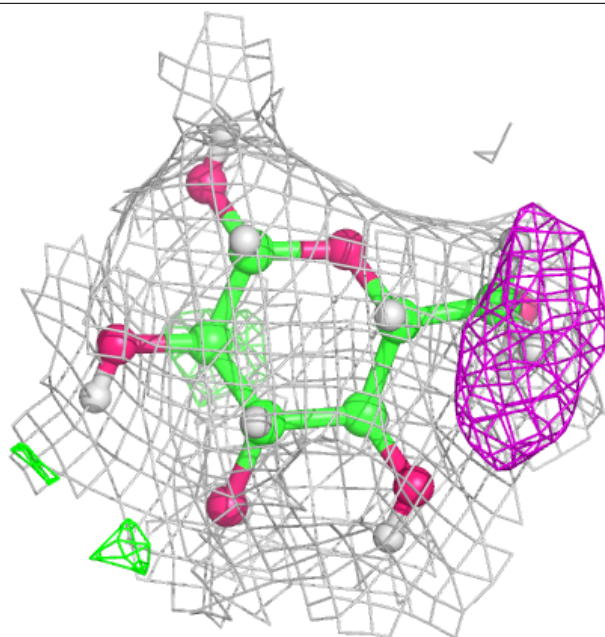
**Electron density around BGC D 501:**

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



**Electron density around BGC B 501:**

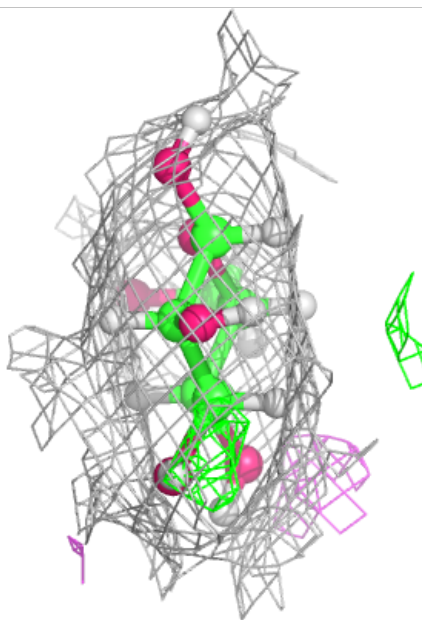
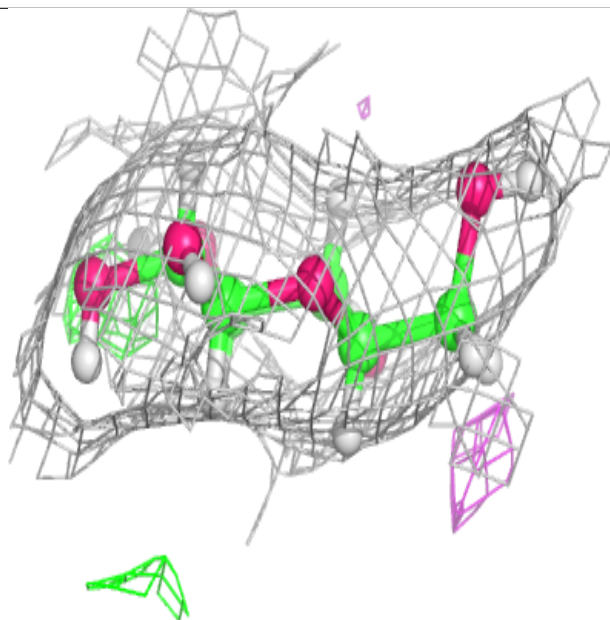
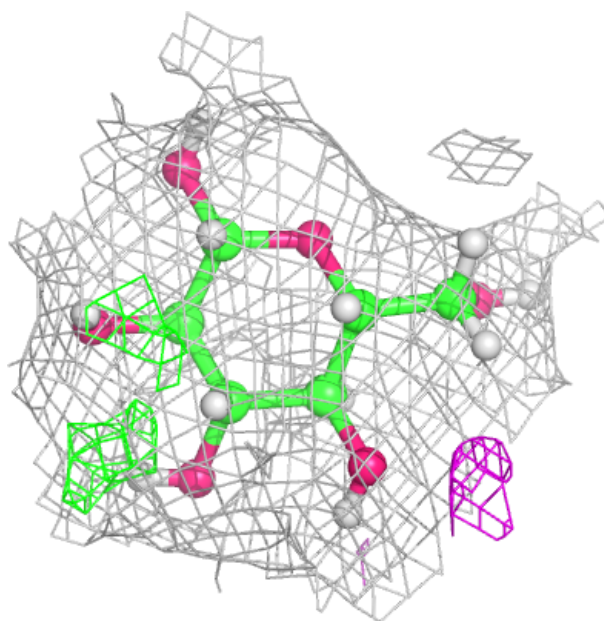
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





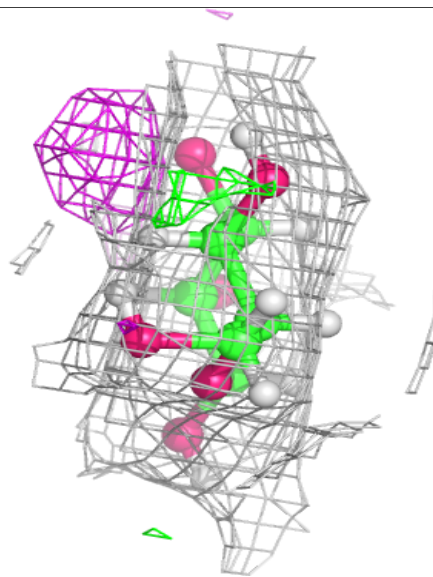
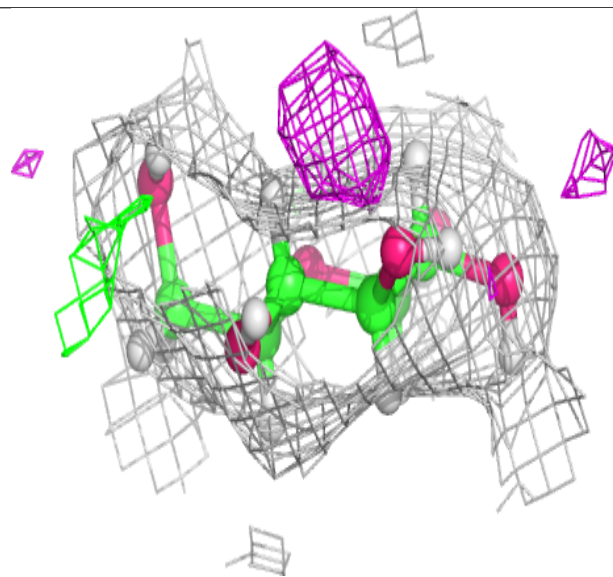
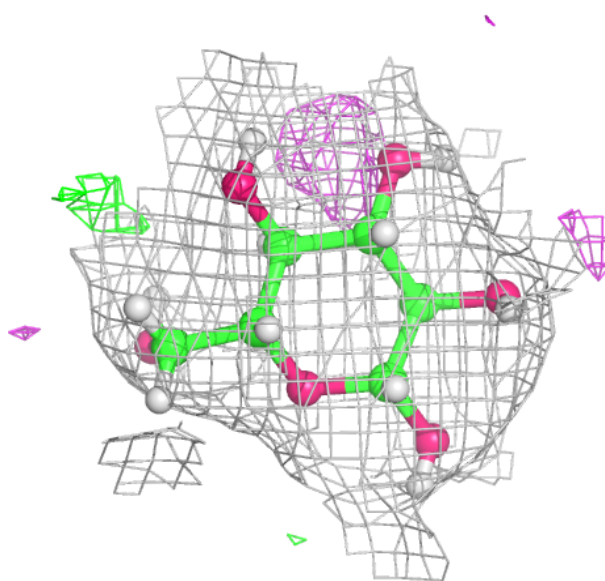
**Electron density around BGC C 501:**

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



**Electron density around BGC 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)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.