



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 26, 2025 – 02:08 PM EDT

PDB ID : 9MK2 / pdb_00009mk2
Title : Crystal structure of Neisseria meningitidis ClpP protease complex with non-covalent activator, ACP1-01
Authors : Mabanglo, M.F.; Houry, W.A.
Deposited on : 2024-12-16
Resolution : 1.95 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 Phenix2.0rc1
Mogul : 2022.3.0, CSD as543be (2022)
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 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

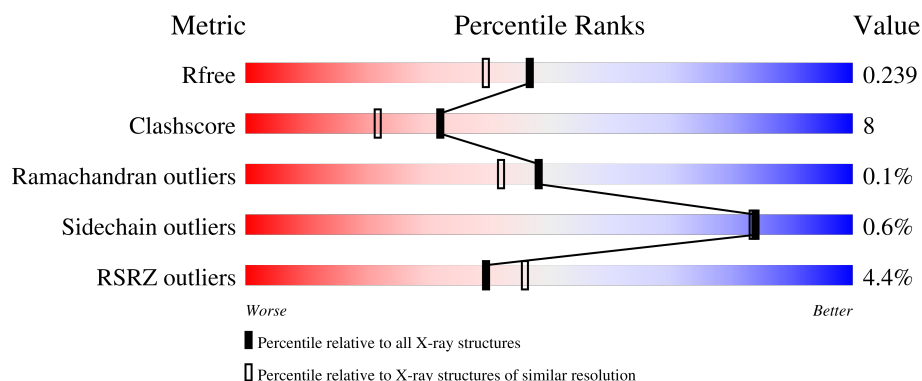
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3187 (1.96-1.96)
Clashscore	180529	3412 (1.96-1.96)
Ramachandran outliers	177936	3390 (1.96-1.96)
Sidechain outliers	177891	3390 (1.96-1.96)
RSRZ outliers	164620	3186 (1.96-1.96)

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	218	<div> <div>2%</div> <div>71% 11% 19%</div> </div>
1	B	218	<div> <div>3%</div> <div>68% 12% 18%</div> </div>
1	C	218	<div> <div>4%</div> <div>74% 9% 17%</div> </div>
1	D	218	<div> <div>2%</div> <div>71% 11% 18%</div> </div>
1	E	218	<div> <div>5%</div> <div>70% 12% 17%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	218	
1	G	218	

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
2	A1BLZ	B	301	-	-	X	-
2	A1BLZ	C	301	-	-	X	-
2	A1BLZ	E	301	-	-	X	-
2	A1BLZ	F	301	-	-	X	-
2	A1BLZ	G	301	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10464 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 ATP-dependent Clp protease proteolytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	177	Total	C	N	O	S	0	0	0
			1384	873	236	267	8			
1	B	178	Total	C	N	O	S	0	0	0
			1399	887	237	267	8			
1	C	181	Total	C	N	O	S	0	0	0
			1414	895	240	271	8			
1	D	178	Total	C	N	O	S	0	0	0
			1392	879	237	268	8			
1	E	180	Total	C	N	O	S	0	0	0
			1410	893	239	270	8			
1	F	183	Total	C	N	O	S	0	0	0
			1420	896	242	274	8			
1	G	183	Total	C	N	O	S	0	0	0
			1436	908	245	275	8			

There are 98 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	HIS	-	expression tag	UNP Q9JZ38
A	-12	HIS	-	expression tag	UNP Q9JZ38
A	-11	HIS	-	expression tag	UNP Q9JZ38
A	-10	HIS	-	expression tag	UNP Q9JZ38
A	-9	HIS	-	expression tag	UNP Q9JZ38
A	-8	HIS	-	expression tag	UNP Q9JZ38
A	-7	GLU	-	expression tag	UNP Q9JZ38
A	-6	ASN	-	expression tag	UNP Q9JZ38
A	-5	LEU	-	expression tag	UNP Q9JZ38
A	-4	TYR	-	expression tag	UNP Q9JZ38
A	-3	PHE	-	expression tag	UNP Q9JZ38
A	-2	GLN	-	expression tag	UNP Q9JZ38
A	-1	SER	-	expression tag	UNP Q9JZ38
A	0	ASN	-	expression tag	UNP Q9JZ38
B	-13	HIS	-	expression tag	UNP Q9JZ38

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	HIS	-	expression tag	UNP Q9JZ38
B	-11	HIS	-	expression tag	UNP Q9JZ38
B	-10	HIS	-	expression tag	UNP Q9JZ38
B	-9	HIS	-	expression tag	UNP Q9JZ38
B	-8	HIS	-	expression tag	UNP Q9JZ38
B	-7	GLU	-	expression tag	UNP Q9JZ38
B	-6	ASN	-	expression tag	UNP Q9JZ38
B	-5	LEU	-	expression tag	UNP Q9JZ38
B	-4	TYR	-	expression tag	UNP Q9JZ38
B	-3	PHE	-	expression tag	UNP Q9JZ38
B	-2	GLN	-	expression tag	UNP Q9JZ38
B	-1	SER	-	expression tag	UNP Q9JZ38
B	0	ASN	-	expression tag	UNP Q9JZ38
C	-13	HIS	-	expression tag	UNP Q9JZ38
C	-12	HIS	-	expression tag	UNP Q9JZ38
C	-11	HIS	-	expression tag	UNP Q9JZ38
C	-10	HIS	-	expression tag	UNP Q9JZ38
C	-9	HIS	-	expression tag	UNP Q9JZ38
C	-8	HIS	-	expression tag	UNP Q9JZ38
C	-7	GLU	-	expression tag	UNP Q9JZ38
C	-6	ASN	-	expression tag	UNP Q9JZ38
C	-5	LEU	-	expression tag	UNP Q9JZ38
C	-4	TYR	-	expression tag	UNP Q9JZ38
C	-3	PHE	-	expression tag	UNP Q9JZ38
C	-2	GLN	-	expression tag	UNP Q9JZ38
C	-1	SER	-	expression tag	UNP Q9JZ38
C	0	ASN	-	expression tag	UNP Q9JZ38
D	-13	HIS	-	expression tag	UNP Q9JZ38
D	-12	HIS	-	expression tag	UNP Q9JZ38
D	-11	HIS	-	expression tag	UNP Q9JZ38
D	-10	HIS	-	expression tag	UNP Q9JZ38
D	-9	HIS	-	expression tag	UNP Q9JZ38
D	-8	HIS	-	expression tag	UNP Q9JZ38
D	-7	GLU	-	expression tag	UNP Q9JZ38
D	-6	ASN	-	expression tag	UNP Q9JZ38
D	-5	LEU	-	expression tag	UNP Q9JZ38
D	-4	TYR	-	expression tag	UNP Q9JZ38
D	-3	PHE	-	expression tag	UNP Q9JZ38
D	-2	GLN	-	expression tag	UNP Q9JZ38
D	-1	SER	-	expression tag	UNP Q9JZ38
D	0	ASN	-	expression tag	UNP Q9JZ38
E	-13	HIS	-	expression tag	UNP Q9JZ38

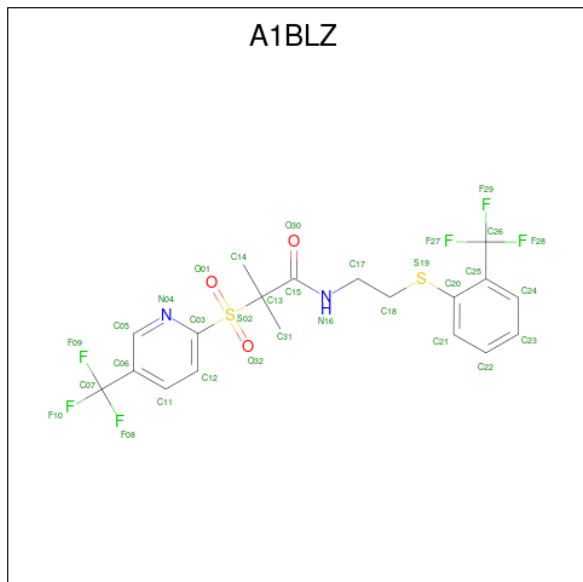
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Chain	Residue	Modelled	Actual	Comment	Reference
E	-12	HIS	-	expression tag	UNP Q9JZ38
E	-11	HIS	-	expression tag	UNP Q9JZ38
E	-10	HIS	-	expression tag	UNP Q9JZ38
E	-9	HIS	-	expression tag	UNP Q9JZ38
E	-8	HIS	-	expression tag	UNP Q9JZ38
E	-7	GLU	-	expression tag	UNP Q9JZ38
E	-6	ASN	-	expression tag	UNP Q9JZ38
E	-5	LEU	-	expression tag	UNP Q9JZ38
E	-4	TYR	-	expression tag	UNP Q9JZ38
E	-3	PHE	-	expression tag	UNP Q9JZ38
E	-2	GLN	-	expression tag	UNP Q9JZ38
E	-1	SER	-	expression tag	UNP Q9JZ38
E	0	ASN	-	expression tag	UNP Q9JZ38
F	-13	HIS	-	expression tag	UNP Q9JZ38
F	-12	HIS	-	expression tag	UNP Q9JZ38
F	-11	HIS	-	expression tag	UNP Q9JZ38
F	-10	HIS	-	expression tag	UNP Q9JZ38
F	-9	HIS	-	expression tag	UNP Q9JZ38
F	-8	HIS	-	expression tag	UNP Q9JZ38
F	-7	GLU	-	expression tag	UNP Q9JZ38
F	-6	ASN	-	expression tag	UNP Q9JZ38
F	-5	LEU	-	expression tag	UNP Q9JZ38
F	-4	TYR	-	expression tag	UNP Q9JZ38
F	-3	PHE	-	expression tag	UNP Q9JZ38
F	-2	GLN	-	expression tag	UNP Q9JZ38
F	-1	SER	-	expression tag	UNP Q9JZ38
F	0	ASN	-	expression tag	UNP Q9JZ38
G	-13	HIS	-	expression tag	UNP Q9JZ38
G	-12	HIS	-	expression tag	UNP Q9JZ38
G	-11	HIS	-	expression tag	UNP Q9JZ38
G	-10	HIS	-	expression tag	UNP Q9JZ38
G	-9	HIS	-	expression tag	UNP Q9JZ38
G	-8	HIS	-	expression tag	UNP Q9JZ38
G	-7	GLU	-	expression tag	UNP Q9JZ38
G	-6	ASN	-	expression tag	UNP Q9JZ38
G	-5	LEU	-	expression tag	UNP Q9JZ38
G	-4	TYR	-	expression tag	UNP Q9JZ38
G	-3	PHE	-	expression tag	UNP Q9JZ38
G	-2	GLN	-	expression tag	UNP Q9JZ38
G	-1	SER	-	expression tag	UNP Q9JZ38
G	0	ASN	-	expression tag	UNP Q9JZ38

- Molecule 2 is 2-methyl-N-(2-{[2-(trifluoromethyl)phenyl]sulfanyl}ethyl)-2-[5-(trifluorometh

yl)pyridine-2-sulfonyl]propanamide (CCD ID: A1BLZ) (formula: C₁₉H₁₈F₆N₂O₃S₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 32	C 19	F 6	N 2	O 3	S 2	0	0
2	B	1	Total 32	C 19	F 6	N 2	O 3	S 2	0	0
2	C	1	Total 32	C 19	F 6	N 2	O 3	S 2	0	0
2	D	1	Total 32	C 19	F 6	N 2	O 3	S 2	0	0
2	E	1	Total 32	C 19	F 6	N 2	O 3	S 2	0	0
2	F	1	Total 32	C 19	F 6	N 2	O 3	S 2	0	0
2	G	1	Total 32	C 19	F 6	N 2	O 3	S 2	0	0

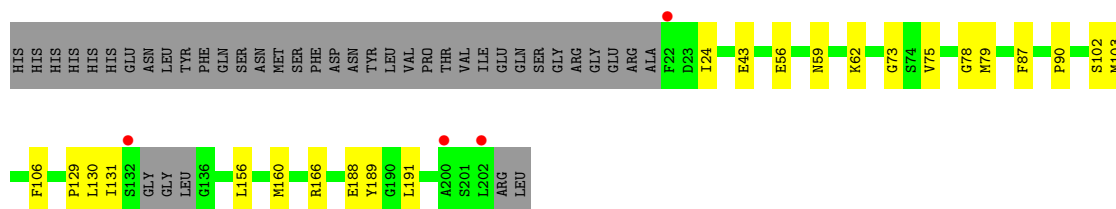
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	55	Total	O	0	0
			55	55		
3	B	57	Total	O	0	0
			57	57		
3	C	57	Total	O	0	0
			57	57		

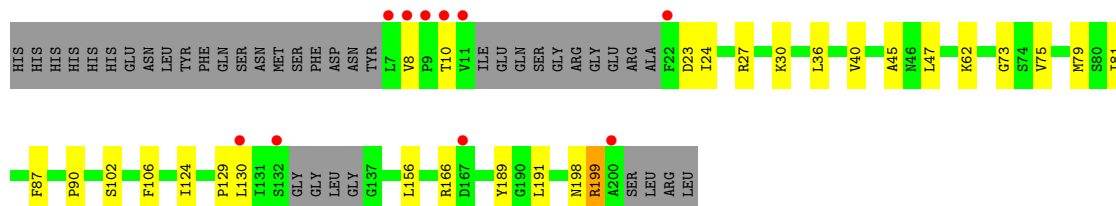
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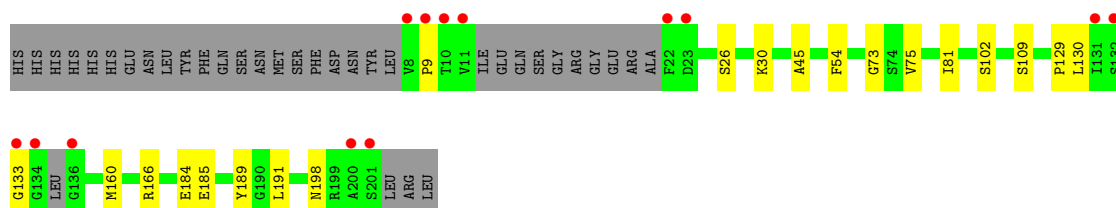
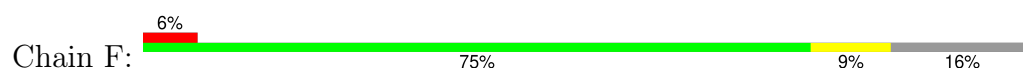
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	56	Total 56	O 56	0	0
3	E	54	Total 54	O 54	0	0
3	F	52	Total 52	O 52	0	0
3	G	54	Total 54	O 54	0	0



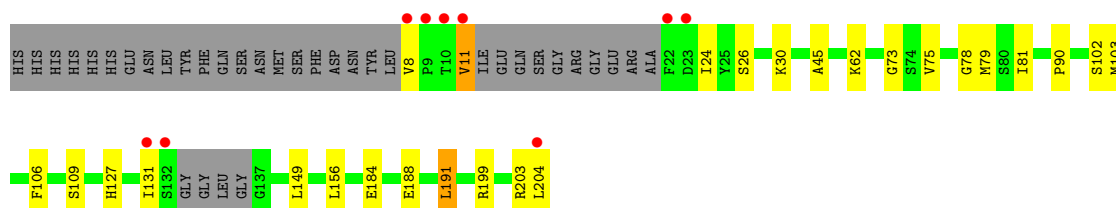
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	97.35Å 118.70Å 127.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	75.27 – 1.95 75.27 – 1.95	Depositor EDS
% Data completeness (in resolution range)	100.0 (75.27-1.95) 99.9 (75.27-1.95)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.194 , 0.239 0.194 , 0.239	Depositor DCC
R_{free} test set	5294 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	19.5	Xtriage
Anisotropy	0.736	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10464	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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: A1BLZ

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	0/1404	0.56	0/1888
1	B	0.67	3/1419 (0.2%)	0.71	2/1910 (0.1%)
1	C	0.39	0/1434	0.57	0/1930
1	D	0.38	0/1412	0.52	0/1899
1	E	0.37	0/1430	0.51	0/1925
1	F	0.36	0/1440	0.53	0/1937
1	G	0.44	0/1456	0.55	0/1958
All	All	0.44	3/9995 (0.0%)	0.57	2/13447 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	144	ILE	N-CA	-6.47	1.39	1.46
1	B	146	ALA	C-O	-5.55	1.17	1.24
1	B	152	ILE	C-O	-5.25	1.18	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	144	ILE	CA-C-O	-7.61	113.10	121.17
1	B	148	GLU	N-CA-CB	6.37	119.49	110.12

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	1384	0	1385	20	0
1	B	1399	0	1409	29	0
1	C	1414	0	1422	20	0
1	D	1392	0	1396	18	0
1	E	1410	0	1419	23	0
1	F	1420	0	1422	19	0
1	G	1436	0	1448	30	0
2	A	32	0	0	7	0
2	B	32	0	0	14	0
2	C	32	0	0	13	0
2	D	32	0	0	7	0
2	E	32	0	0	9	0
2	F	32	0	0	11	0
2	G	32	0	0	17	0
3	A	55	0	0	2	0
3	B	57	0	0	0	0
3	C	57	0	0	0	0
3	D	56	0	0	1	0
3	E	54	0	0	0	0
3	F	52	0	0	2	0
3	G	54	0	0	0	0
All	All	10464	0	9901	158	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

The worst 5 of 158 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:75:VAL:HG23	2:C:301:A1BLZ:C14	1.95	0.96
3:F:419:HOH:O	1:G:199:ARG:HD2	1.65	0.96
1:C:129:PRO:HA	2:C:301:A1BLZ:O32	1.71	0.91
1:E:75:VAL:HG23	2:E:301:A1BLZ:C14	2.02	0.89
2:G:301:A1BLZ:F29	2:G:301:A1BLZ:S19	2.21	0.88

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	173/218 (79%)	171 (99%)	2 (1%)	0	100	100
1	B	172/218 (79%)	169 (98%)	3 (2%)	0	100	100
1	C	175/218 (80%)	173 (99%)	2 (1%)	0	100	100
1	D	174/218 (80%)	172 (99%)	2 (1%)	0	100	100
1	E	174/218 (80%)	170 (98%)	4 (2%)	0	100	100
1	F	177/218 (81%)	173 (98%)	3 (2%)	1 (1%)	22	13
1	G	177/218 (81%)	172 (97%)	5 (3%)	0	100	100
All	All	1222/1526 (80%)	1200 (98%)	21 (2%)	1 (0%)	48	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	133	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	151/187 (81%)	151 (100%)	0	100	100
1	B	154/187 (82%)	151 (98%)	3 (2%)	52	47
1	C	155/187 (83%)	155 (100%)	0	100	100
1	D	152/187 (81%)	152 (100%)	0	100	100
1	E	155/187 (83%)	154 (99%)	1 (1%)	84	83

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	155/187 (83%)	155 (100%)	0	100	100
1	G	158/187 (84%)	156 (99%)	2 (1%)	65	62
All	All	1080/1309 (82%)	1074 (99%)	6 (1%)	84	83

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	199	ARG
1	G	11	VAL
1	G	191	LEU
1	B	22	PHE
1	B	11	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	198	ASN
1	F	121	ASN
1	F	194	GLN
1	D	59	ASN
1	D	46	ASN

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](#)

7 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	A1BLZ	B	301	-	32,33,33	2.79	14 (43%)	37,51,51	5.85	17 (45%)
2	A1BLZ	F	301	-	32,33,33	3.02	13 (40%)	37,51,51	4.94	20 (54%)
2	A1BLZ	G	301	-	32,33,33	2.94	15 (46%)	37,51,51	5.18	18 (48%)
2	A1BLZ	C	301	-	32,33,33	2.79	15 (46%)	37,51,51	5.86	17 (45%)
2	A1BLZ	A	301	-	32,33,33	2.66	8 (25%)	37,51,51	4.91	20 (54%)
2	A1BLZ	E	301	-	32,33,33	2.62	12 (37%)	37,51,51	5.36	16 (43%)
2	A1BLZ	D	301	-	32,33,33	2.79	15 (46%)	37,51,51	5.86	17 (45%)

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	A1BLZ	B	301	-	-	18/37/40/40	0/2/2/2
2	A1BLZ	F	301	-	-	15/37/40/40	0/2/2/2
2	A1BLZ	G	301	-	-	13/37/40/40	0/2/2/2
2	A1BLZ	C	301	-	-	18/37/40/40	0/2/2/2
2	A1BLZ	A	301	-	-	10/37/40/40	0/2/2/2
2	A1BLZ	E	301	-	-	14/37/40/40	0/2/2/2
2	A1BLZ	D	301	-	-	18/37/40/40	0/2/2/2

The worst 5 of 92 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	301	A1BLZ	O32-S02	-7.95	1.36	1.44
2	F	301	A1BLZ	O32-S02	-7.88	1.36	1.44
2	E	301	A1BLZ	O32-S02	-7.81	1.36	1.44
2	A	301	A1BLZ	O32-S02	-7.78	1.36	1.44
2	B	301	A1BLZ	O32-S02	-6.68	1.37	1.44

The worst 5 of 125 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	A1BLZ	O32-S02-O01	-23.42	89.66	119.06
2	B	301	A1BLZ	O32-S02-O01	-23.41	89.68	119.06
2	C	301	A1BLZ	O32-S02-O01	-23.40	89.68	119.06
2	E	301	A1BLZ	O32-S02-O01	-17.81	96.71	119.06
2	F	301	A1BLZ	O32-S02-O01	-16.52	98.32	119.06

There are no chirality outliers.

5 of 106 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	A1BLZ	C15-C13-S02-O01
2	A	301	A1BLZ	C15-C13-S02-O32
2	A	301	A1BLZ	C14-C13-S02-C03
2	A	301	A1BLZ	C14-C13-S02-O01
2	A	301	A1BLZ	C14-C13-S02-O32

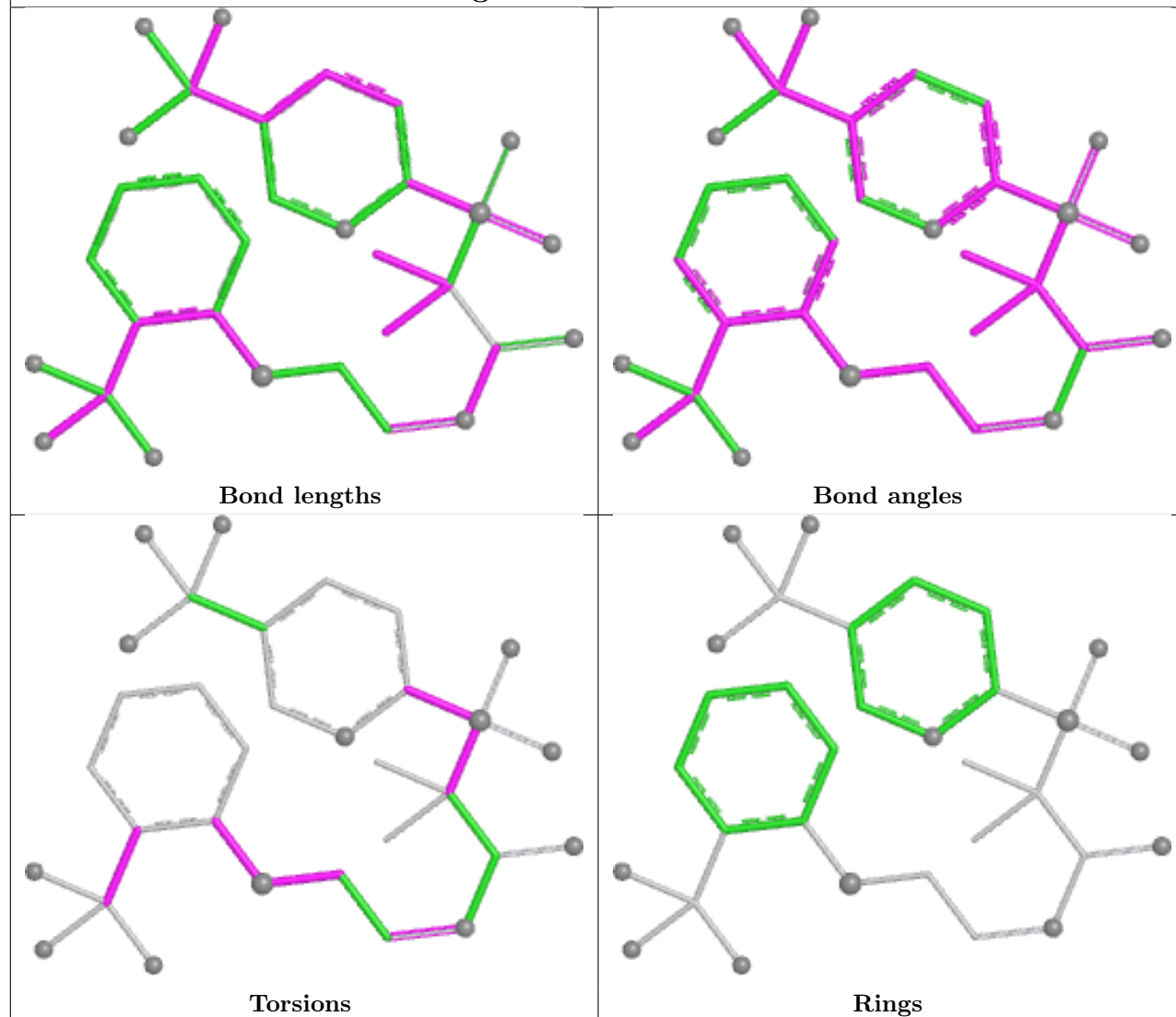
There are no ring outliers.

7 monomers are involved in 78 short contacts:

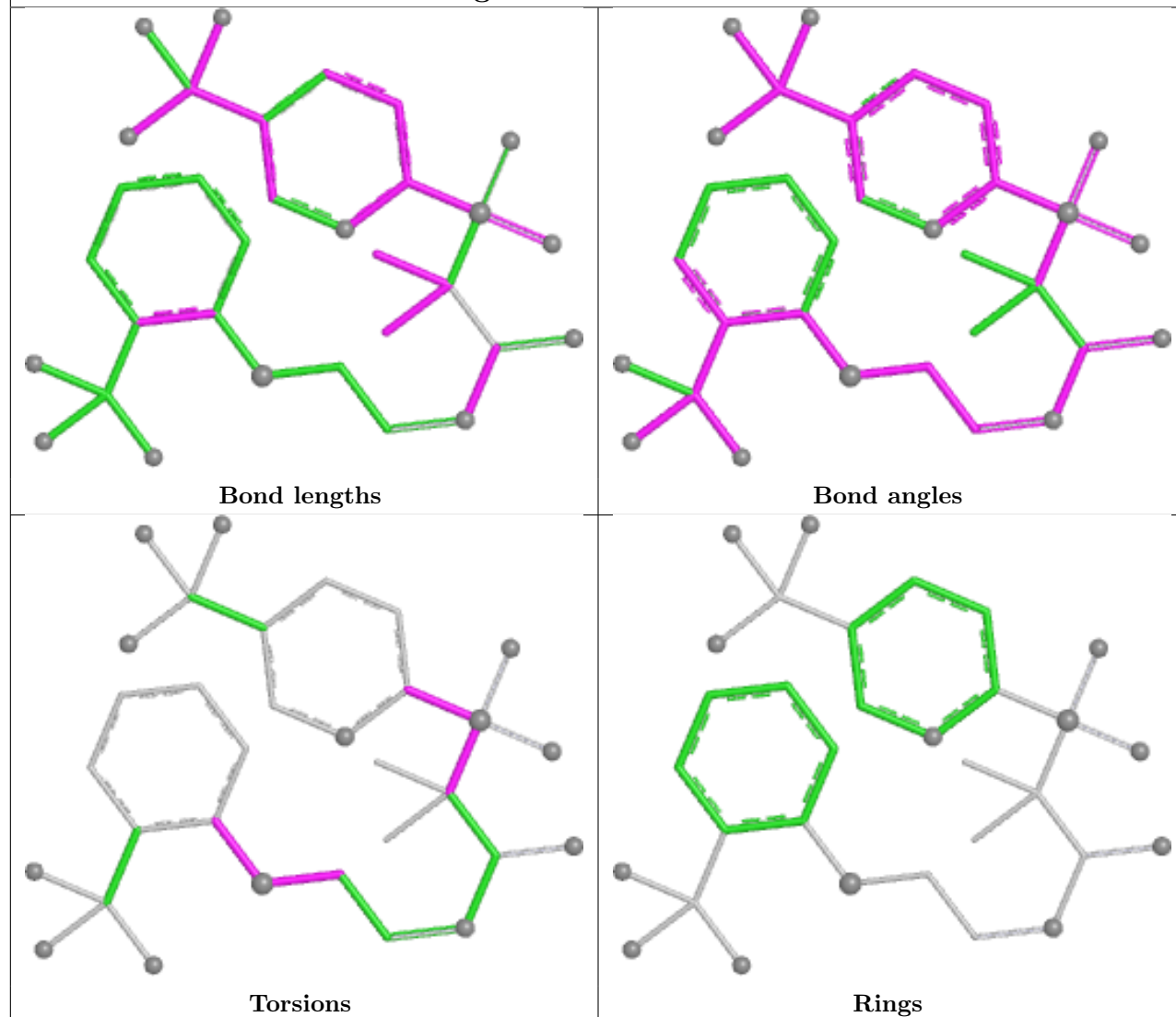
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	A1BLZ	14	0
2	F	301	A1BLZ	11	0
2	G	301	A1BLZ	17	0
2	C	301	A1BLZ	13	0
2	A	301	A1BLZ	7	0
2	E	301	A1BLZ	9	0
2	D	301	A1BLZ	7	0

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.

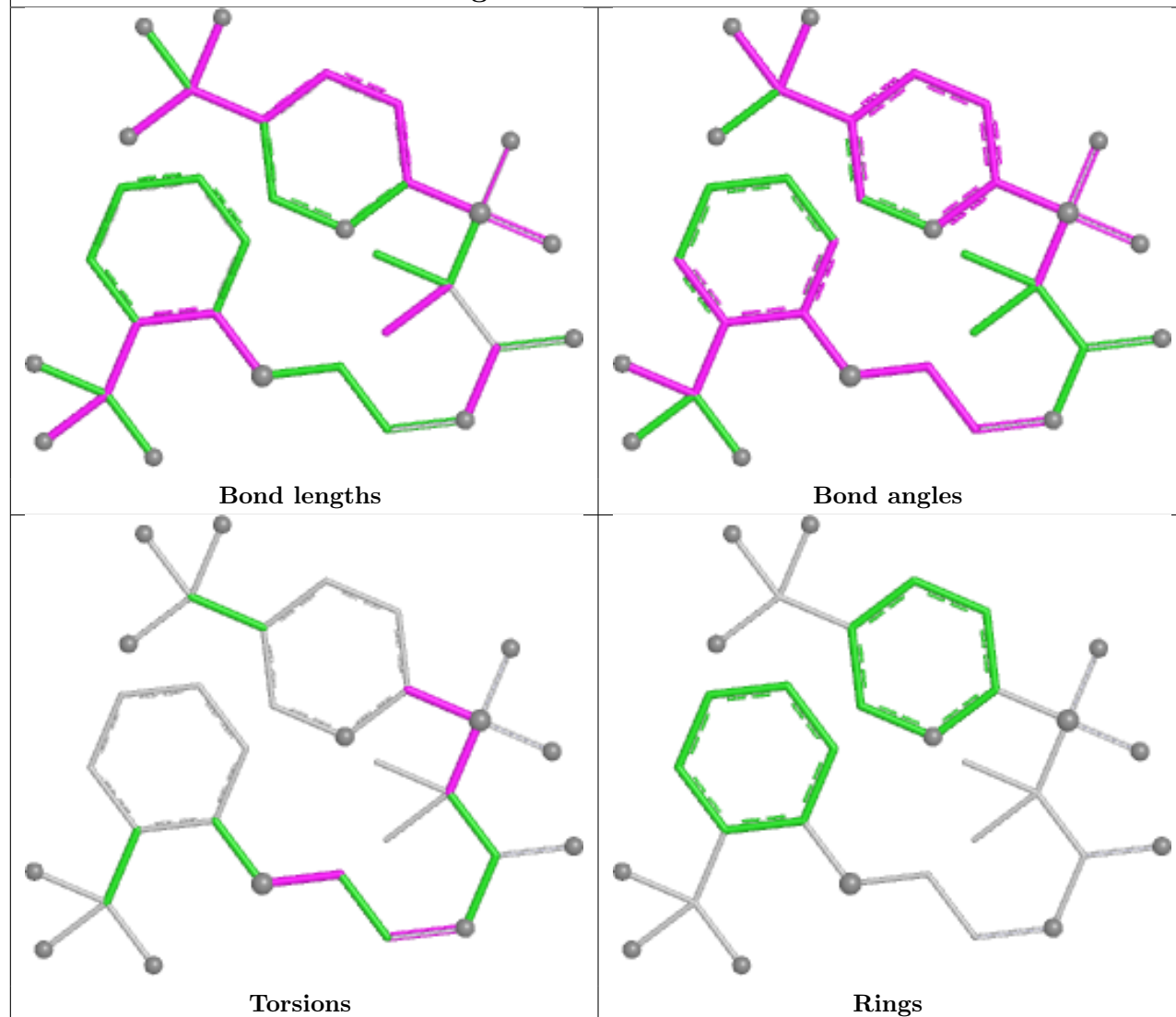
Ligand A1BLZ B 301



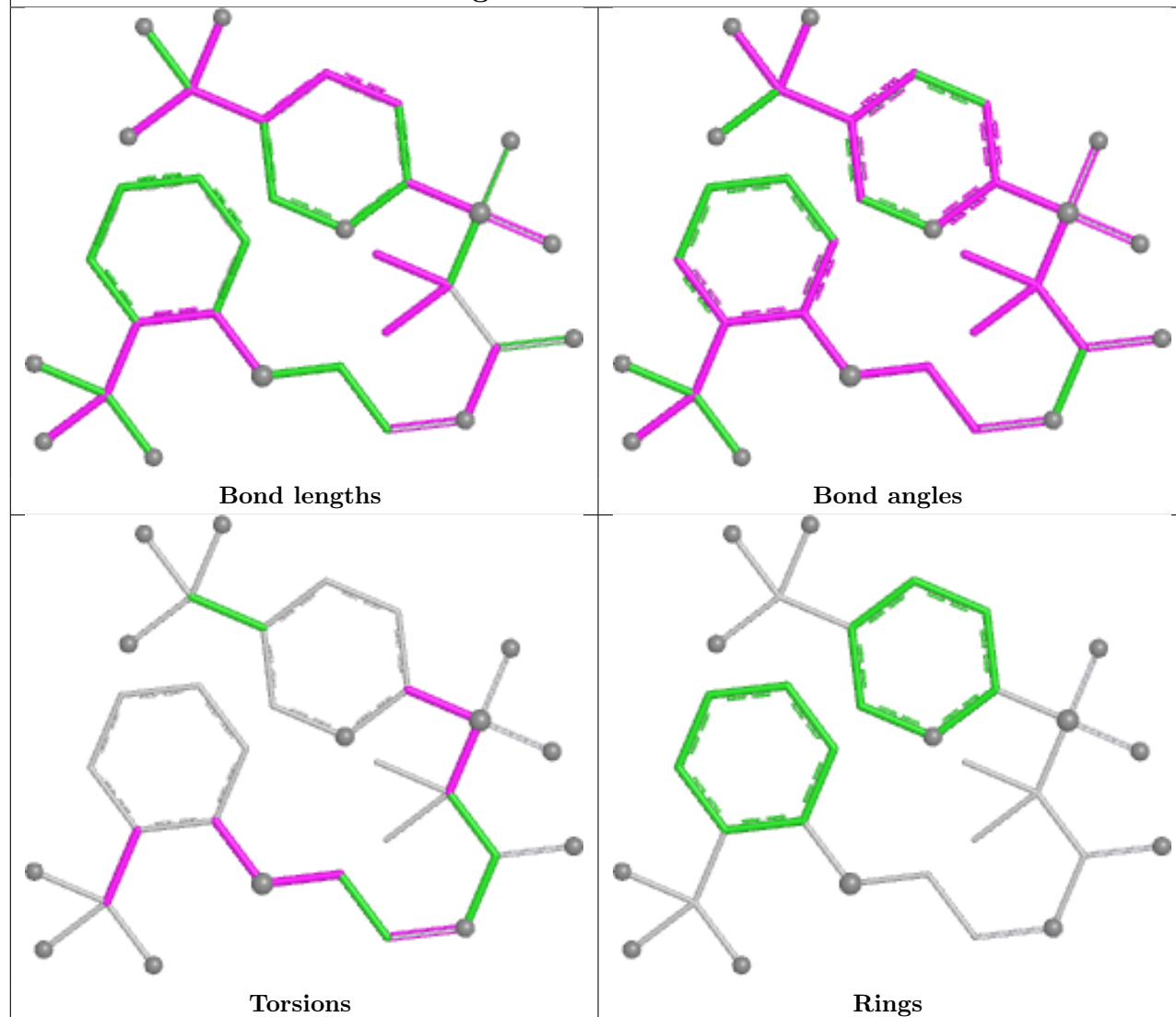
Ligand A1BLZ F 301



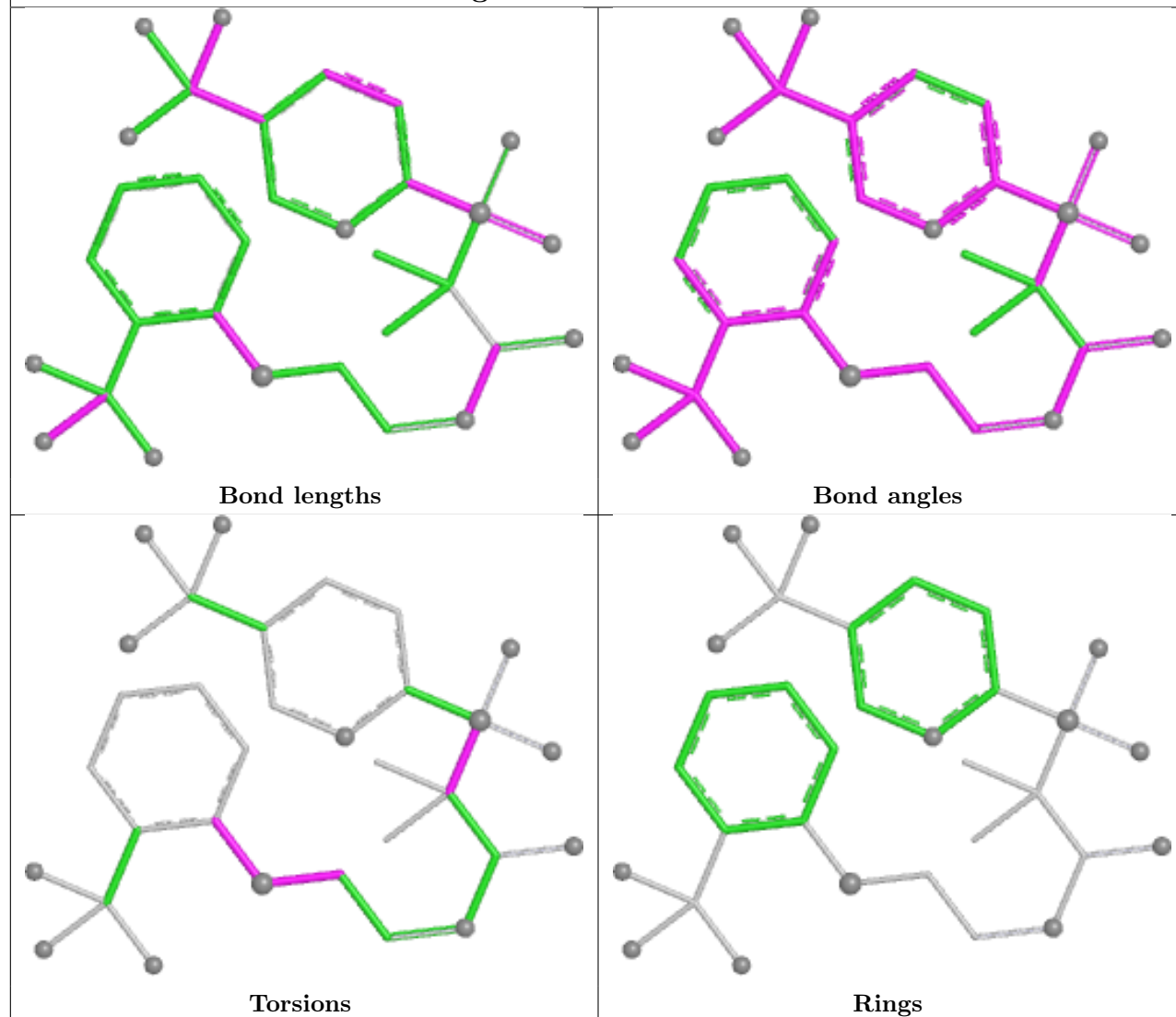
Ligand A1BLZ G 301



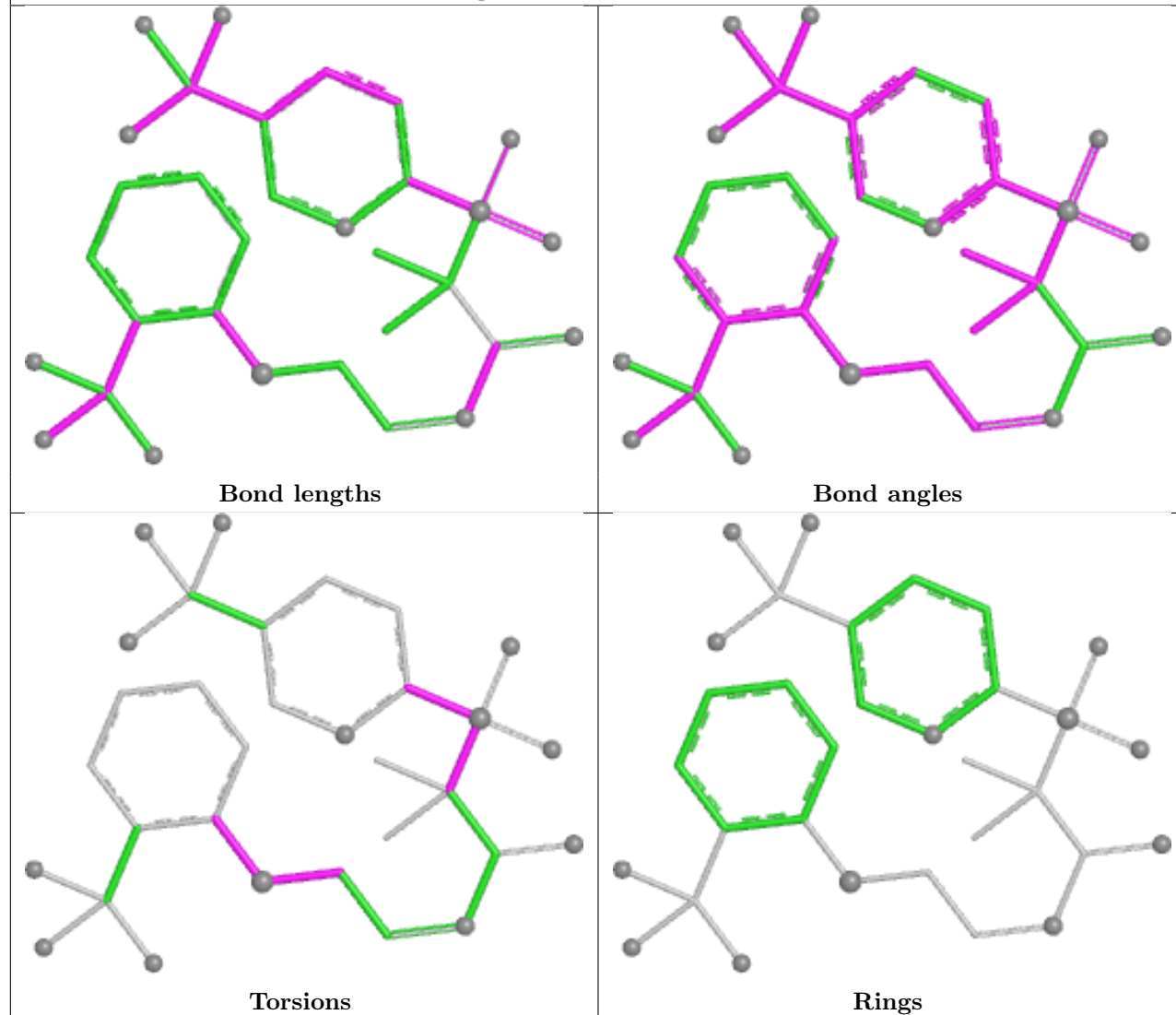
Ligand A1BLZ C 301

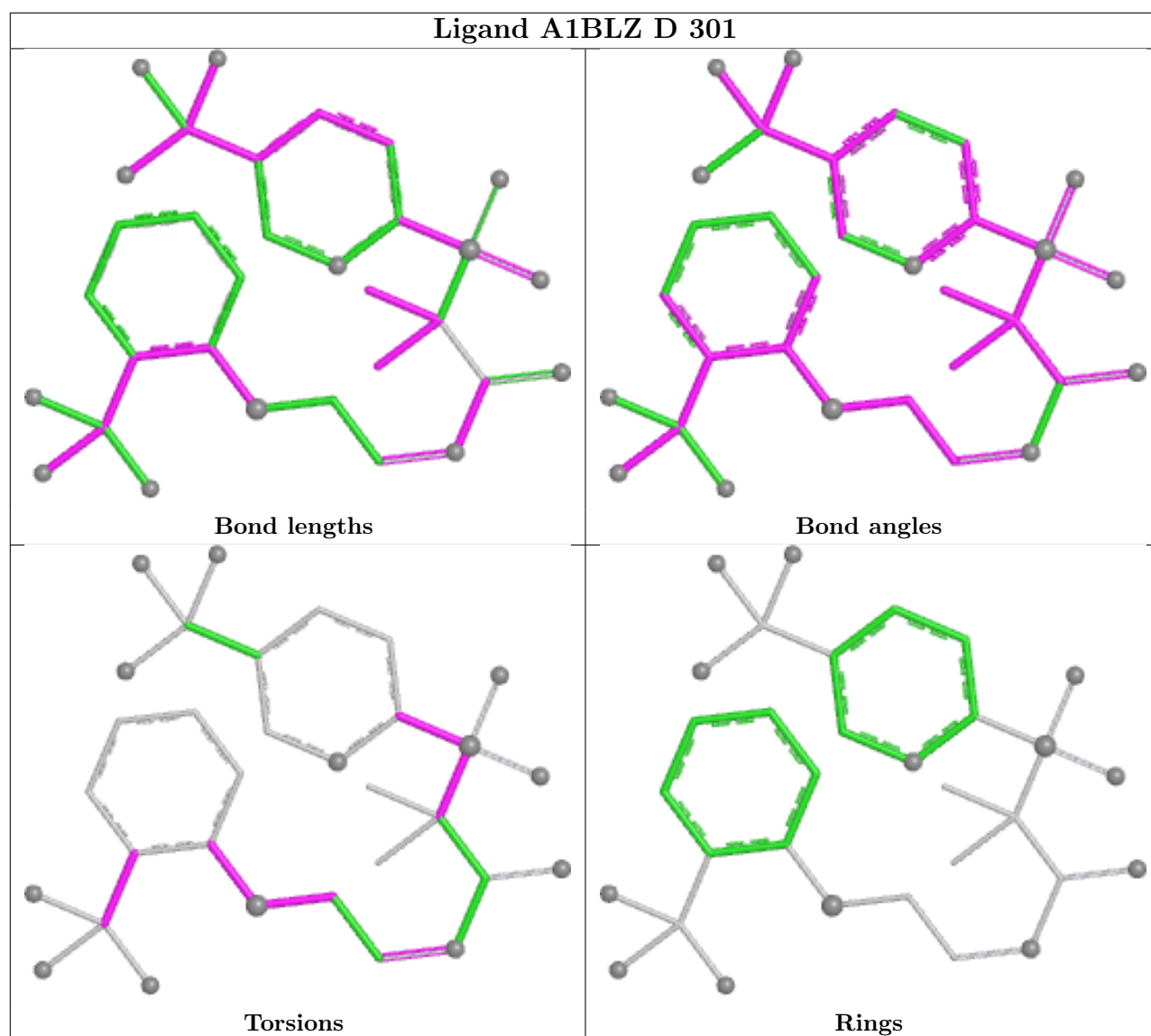


Ligand A1BLZ A 301



Ligand A1BLZ E 301





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	177/218 (81%)	-0.20	5 (2%) 55 61	12, 19, 35, 66	0
1	B	178/218 (81%)	-0.01	7 (3%) 44 50	14, 20, 38, 51	0
1	C	181/218 (83%)	-0.09	8 (4%) 39 46	14, 19, 35, 73	0
1	D	178/218 (81%)	-0.14	4 (2%) 62 68	14, 20, 36, 71	0
1	E	180/218 (82%)	-0.01	10 (5%) 31 38	14, 20, 44, 67	0
1	F	183/218 (83%)	0.11	13 (7%) 23 28	13, 21, 44, 77	0
1	G	183/218 (83%)	-0.06	9 (4%) 36 42	14, 19, 39, 81	0
All	All	1260/1526 (82%)	-0.06	56 (4%) 39 46	12, 20, 39, 81	0

The worst 5 of 56 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	22	PHE	7.7
1	G	22	PHE	7.1
1	F	22	PHE	6.9
1	F	10	THR	6.4
1	D	22	PHE	6.1

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](#)

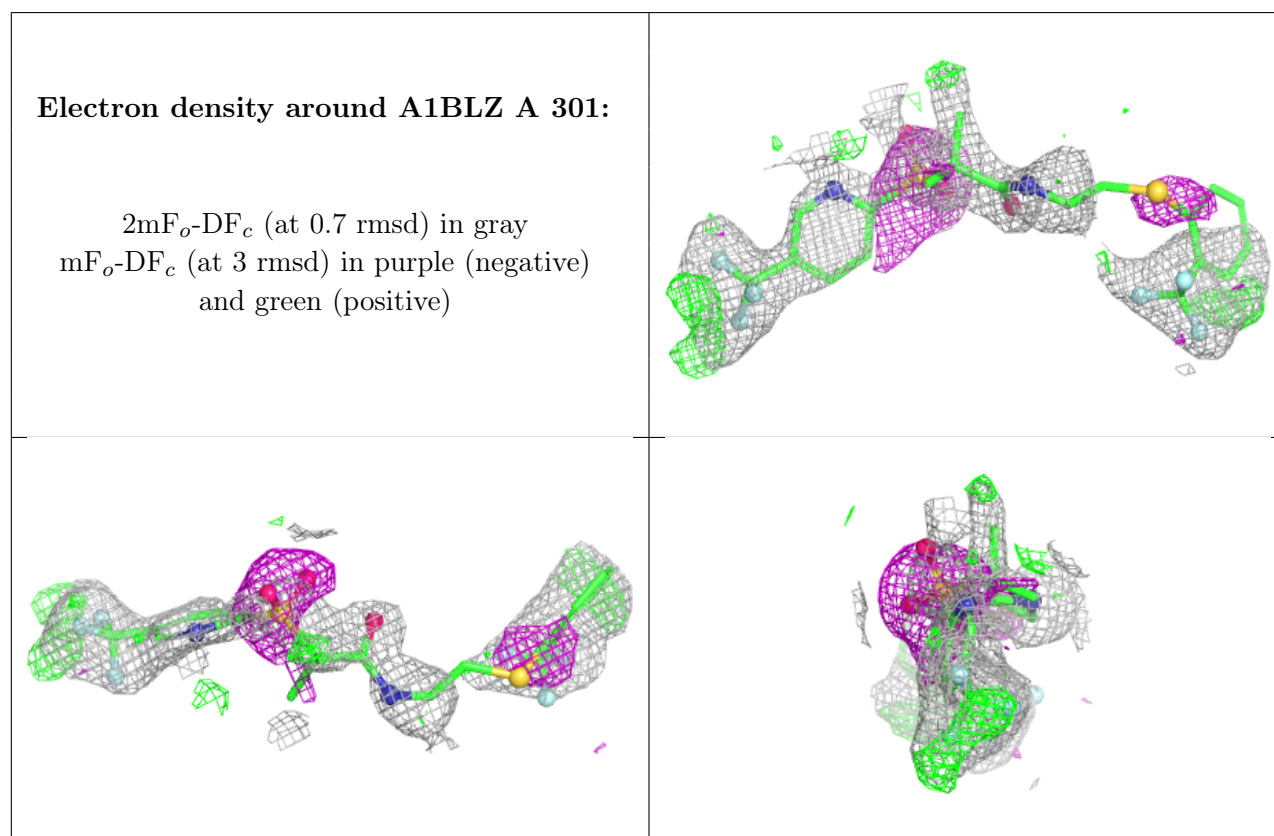
There are no oligosaccharides in this entry.

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, 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.

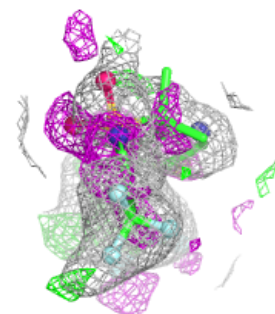
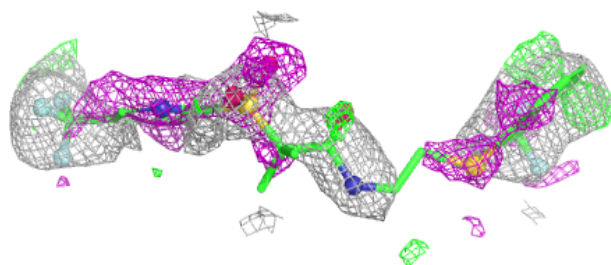
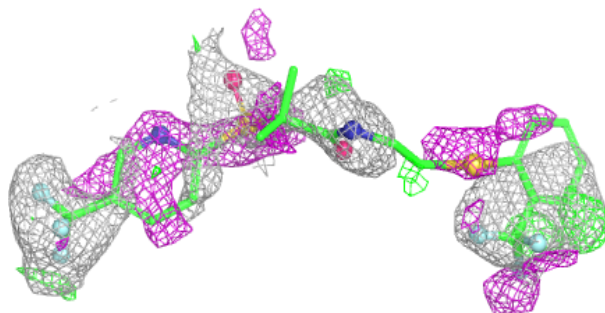
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	A1BLZ	A	301	32/32	0.68	0.22	24,42,59,86	32
2	A1BLZ	B	301	32/32	0.68	0.27	20,47,62,83	32
2	A1BLZ	C	301	32/32	0.69	0.23	23,44,59,96	32
2	A1BLZ	E	301	32/32	0.72	0.23	26,43,57,79	32
2	A1BLZ	D	301	32/32	0.74	0.21	23,41,51,78	32
2	A1BLZ	G	301	32/32	0.74	0.22	22,43,55,66	32
2	A1BLZ	F	301	32/32	0.78	0.18	28,43,60,69	32

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.

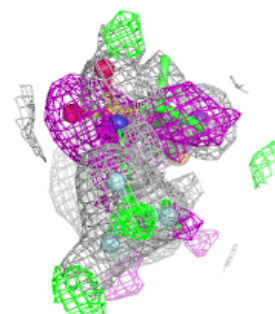
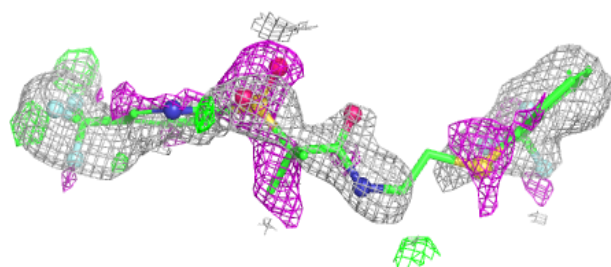
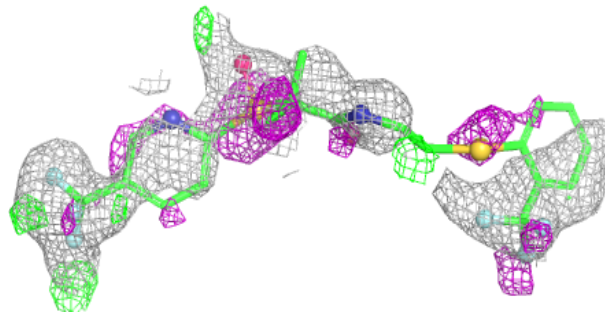


Electron density around A1BLZ B 301:

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

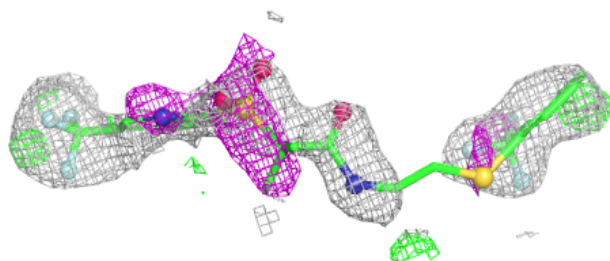
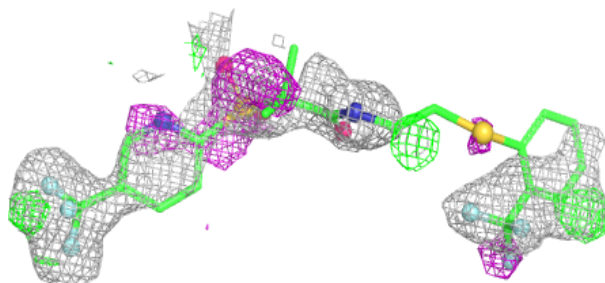
**Electron density around A1BLZ C 301:**

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

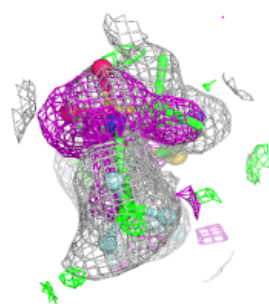
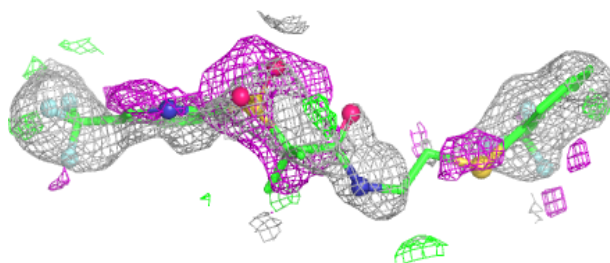
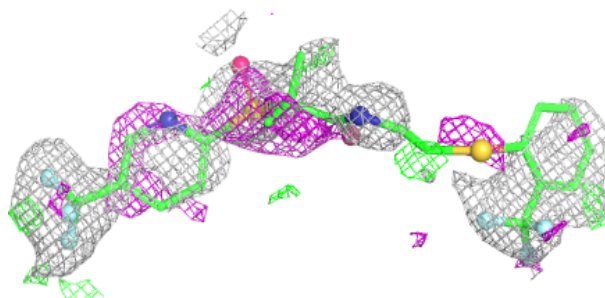


Electron density around A1BLZ E 301:

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

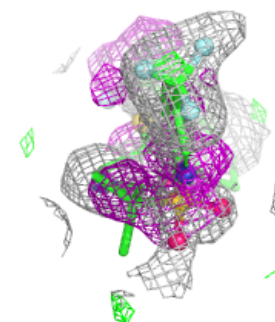
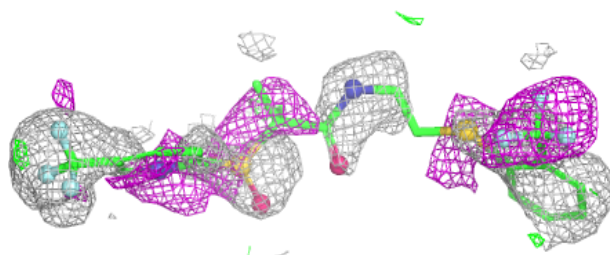
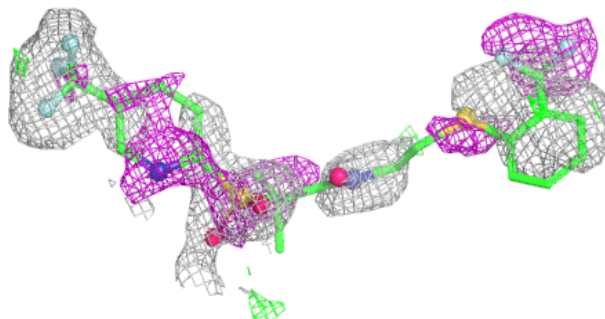
**Electron density around A1BLZ D 301:**

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

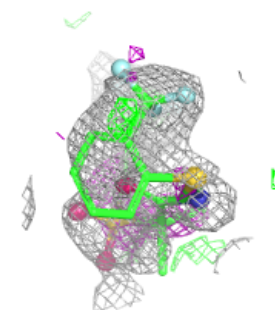
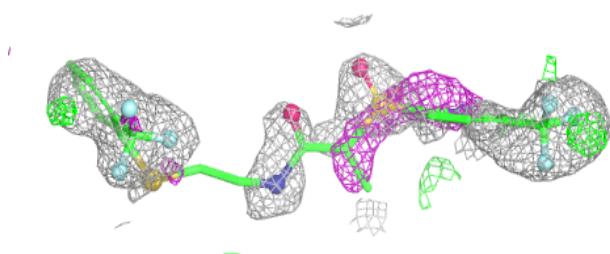
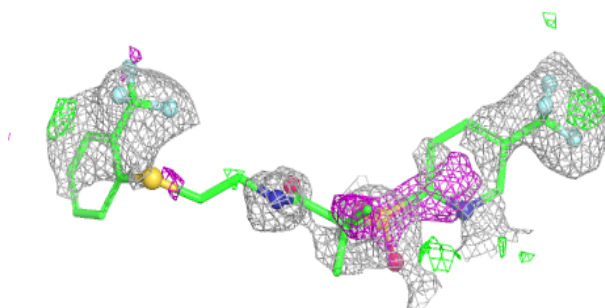


Electron density around A1BLZ G 301:

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

**Electron density around A1BLZ F 301:**

$2mF_o-DF_c$ (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.