



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 7, 2020 – 03:22 am BST

PDB ID : 6JN4
Title : Serine Beta-Lactamase KPC-2 in Complex with Dual MBL/SBL Inhibitor WL-001
Authors : Li, G.-B.; Liu, S.
Deposited on : 2019-03-13
Resolution : 1.90 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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

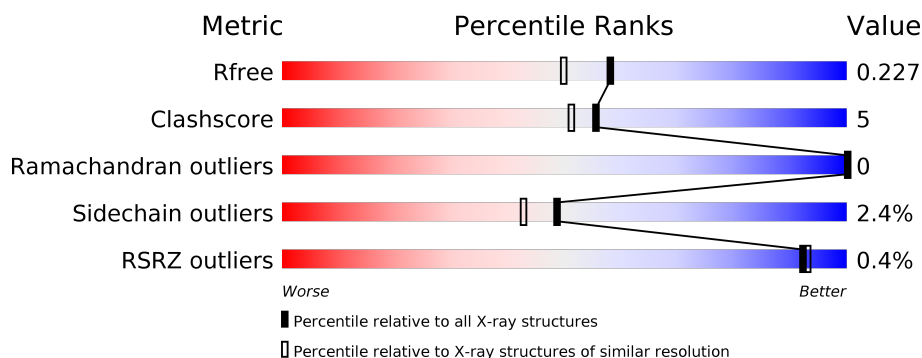
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 1.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	265	 93% 5% •
1	B	265	 91% 9% •
1	C	265	 91% 6% • •
1	D	265	 91% 9%

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	ACT	A	302	-	-	X	-
3	ACT	D	302	-	-	X	-
4	ACY	A	304	-	-	X	-
4	ACY	B	303	-	-	X	-
4	ACY	C	303	-	-	X	-
4	ACY	D	303	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8652 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.

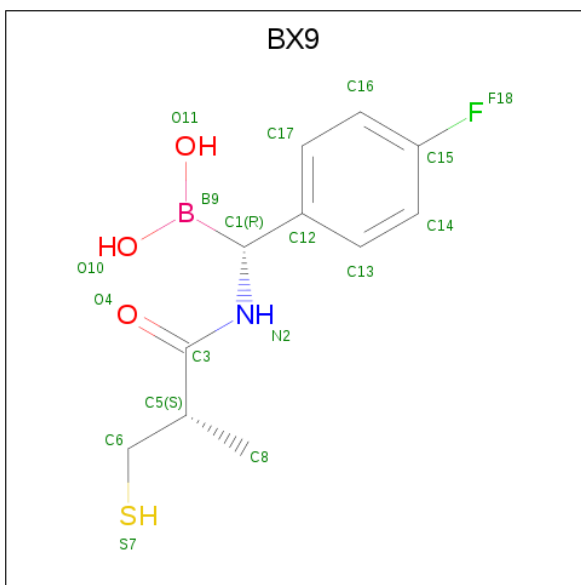
- Molecule 1 is a protein called Serine Beta-Lactamase KPC-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	1
			1971	1240	346	380	5			
1	B	263	Total	C	N	O	S	0	0	1
			1961	1235	345	376	5			
1	C	258	Total	C	N	O	S	0	0	1
			1924	1212	339	368	5			
1	D	265	Total	C	N	O	S	0	0	1
			1972	1241	346	380	5			

There are 4 discrepancies between the modelled and reference sequences:

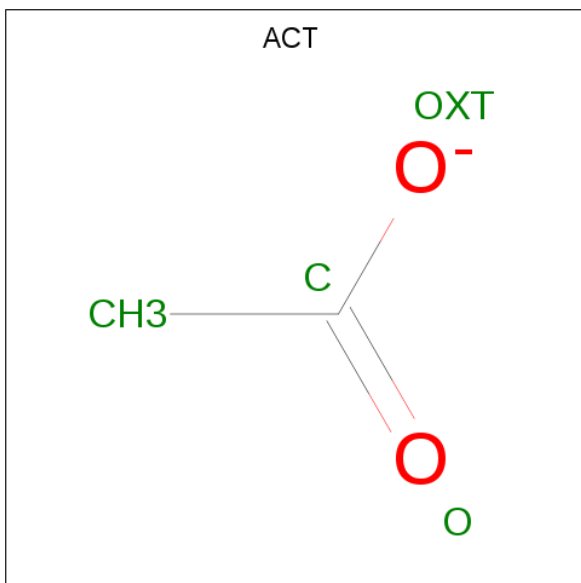
Chain	Residue	Modelled	Actual	Comment	Reference
A	25	SER	-	expression tag	UNP Q93LQ9
B	25	SER	-	expression tag	UNP Q93LQ9
C	25	SER	-	expression tag	UNP Q93LQ9
D	25	SER	-	expression tag	UNP Q93LQ9

- Molecule 2 is [(R)-(4-fluorophenyl)-[[(2S)-2-methyl-3-sulfanyl-propanoyl]amino]methyl]boronic acid (three-letter code: BX9) (formula: C₁₁H₁₅BFNO₃S).



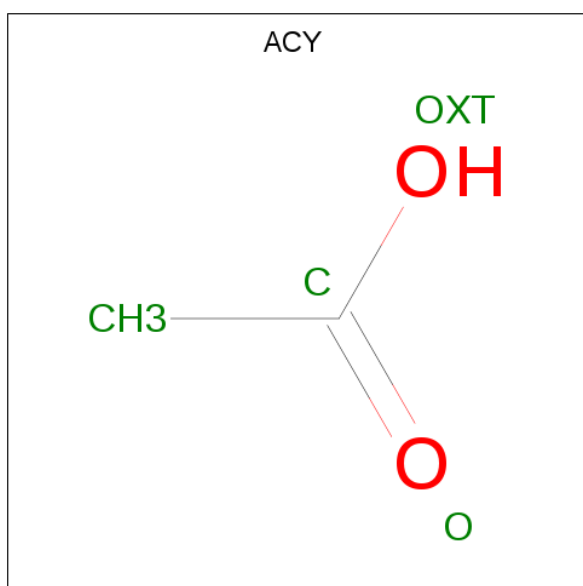
Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	B	C	F	N	O	S		
2	A	1	Total 18	1	11	1	1	3	1	0	0
2	B	1	Total 18	1	11	1	1	3	1	0	0
2	C	1	Total 18	1	11	1	1	3	1	0	0
2	D	1	Total 18	1	11	1	1	3	1	0	0

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



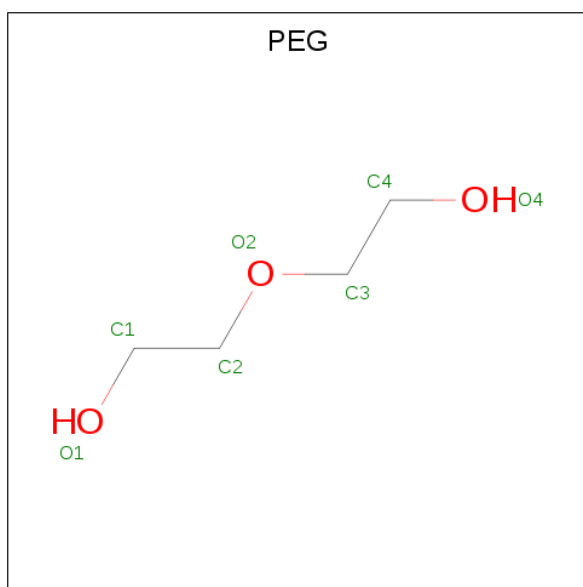
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 4 2 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0
3	D	1	Total C O 4 2 2	0	0

- Molecule 4 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).



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

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

- Molecule 6 is water.

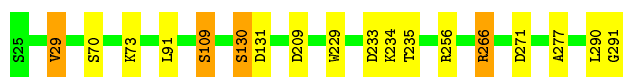
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	200	Total	O	0	0
			200	200		
6	B	213	Total	O	0	0
			213	213		
6	C	135	Total	O	0	1
			136	136		
6	D	109	Total	O	0	2
			111	111		

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: Serine Beta-Lactamase KPC-2

Chain A: 

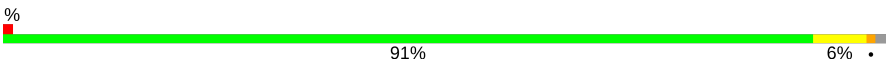


• Molecule 1: Serine Beta-Lactamase KPC-2

Chain B: 



• Molecule 1: Serine Beta-Lactamase KPC-2

Chain C: 



• Molecule 1: Serine Beta-Lactamase KPC-2

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	164.78 Å 164.78 Å 94.68 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.57 – 1.90 47.57 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.57-1.90) 97.3 (47.57-1.90)	Depositor EDS
R_{merge}	0.30	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 1.90 Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.180 , 0.221 0.189 , 0.227	Depositor DCC
R_{free} test set	1990 reflections (1.73%)	wwPDB-VP
Wilson B-factor (Å ²)	18.0	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 29.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.136 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8652	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BX9, ACY, PEG, ACT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.57	0/2011	0.72	2/2741 (0.1%)
1	B	0.61	0/2001	0.68	1/2726 (0.0%)
1	C	0.53	0/1964	0.67	0/2674
1	D	0.49	0/2012	0.62	0/2742
All	All	0.55	0/7988	0.67	3/10883 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	266	ARG	NE-CZ-NH2	-7.15	116.72	120.30
1	A	266	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	B	178	ARG	NE-CZ-NH2	-5.34	117.63	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1971	0	1944	22	0
1	B	1961	0	1941	17	0
1	C	1924	0	1906	17	0
1	D	1972	0	1947	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	18	0	0	3	0
2	B	18	0	0	4	0
2	C	18	0	0	2	0
2	D	18	0	0	5	0
3	A	4	0	3	7	0
3	B	4	0	3	1	0
3	C	4	0	3	1	0
3	D	4	0	3	2	0
4	A	8	0	6	5	0
4	B	4	0	3	5	0
4	C	4	0	3	5	0
4	D	4	0	3	8	0
5	A	14	0	20	1	0
5	B	28	0	40	1	0
5	C	14	0	20	4	0
6	A	200	0	0	4	2
6	B	213	0	0	3	2
6	C	136	0	0	1	0
6	D	111	0	0	2	0
All	All	8652	0	7845	81	2

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 5.

All (81) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:291:GLY:N	3:A:302:ACT:OXT	1.86	1.08
2:B:301:BX9:S7	6:B:575:HOH:O	2.15	1.05
1:C:235:THR:OG1	4:C:303:ACY:O	1.79	1.01
1:A:235:THR:OG1	4:A:304:ACY:O	1.79	1.00
1:B:235:THR:OG1	4:B:303:ACY:OXT	1.86	0.92
1:B:284:ARG:NH1	6:B:401:HOH:O	2.11	0.84
2:D:301:BX9:C13	4:D:303:ACY:H2	2.08	0.84
1:C:70:SER:HB2	4:C:303:ACY:H2	1.62	0.79
1:A:70:SER:HB2	4:A:304:ACY:H2	1.68	0.75
1:D:204:ARG:NH2	6:D:402:HOH:O	2.21	0.74
2:D:301:BX9:C14	4:D:303:ACY:H2	2.22	0.69
1:D:70:SER:HB2	4:D:303:ACY:H1	1.76	0.68
1:A:271:ASP:OD1	6:A:401:HOH:O	2.11	0.67
1:A:70:SER:HB2	4:A:304:ACY:CH3	2.30	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:SER:OG	1:A:131:ASP:OD1	2.18	0.61
1:D:235:THR:OG1	4:D:303:ACY:OXT	2.14	0.61
1:B:70:SER:HB2	4:B:303:ACY:H1	1.84	0.59
1:C:85:GLN:HE21	5:C:304:PEG:H41	1.68	0.58
1:A:256:ARG:HH12	3:A:302:ACT:H2	1.70	0.56
2:B:301:BX9:C13	4:B:303:ACY:H2	2.36	0.56
1:B:191:GLN:OE1	5:B:304:PEG:H11	2.06	0.56
1:D:104:PRO:HB2	1:D:105:TRP:CE3	2.41	0.55
1:B:31:GLU:HG2	6:B:444:HOH:O	2.05	0.55
1:D:237:THR:H	4:D:303:ACY:H3	1.73	0.53
1:A:130:SER:OG	4:A:304:ACY:H3	2.09	0.53
1:B:70:SER:OG	2:B:301:BX9:C1	2.57	0.52
1:D:70:SER:CB	4:D:303:ACY:H1	2.39	0.52
1:B:70:SER:HB2	4:B:303:ACY:CH3	2.39	0.52
1:D:226:PRO:HD2	1:D:229:TRP:CD2	2.46	0.51
1:D:237:THR:N	4:D:303:ACY:H3	2.25	0.51
1:B:49:MET:HG2	1:B:191:GLN:HE22	1.76	0.51
1:D:64:GLU:OE1	6:D:401:HOH:O	2.20	0.51
1:C:237:THR:OG1	4:C:303:ACY:OXT	2.21	0.48
1:A:291:GLY:N	3:A:302:ACT:C	2.70	0.48
1:A:29:VAL:O	1:A:29:VAL:HG22	2.14	0.48
1:C:99:LYS:HD2	1:C:102:LEU:HD12	1.96	0.48
1:A:291:GLY:N	3:A:302:ACT:H1	2.29	0.47
5:A:305:PEG:O1	5:A:306:PEG:O1	2.24	0.47
1:D:166:GLU:HB3	1:D:167:LEU:HD12	1.97	0.47
1:A:70:SER:OG	2:A:301:BX9:O11	2.32	0.47
3:A:302:ACT:H3	6:A:525:HOH:O	2.14	0.47
1:B:252:PRO:HB2	1:B:256:ARG:HB2	1.97	0.47
1:A:290:LEU:C	3:A:302:ACT:H1	2.35	0.46
1:A:229:TRP:HZ2	1:C:168:GLU:HG2	1.80	0.46
1:D:276:GLU:H	3:D:302:ACT:CH3	2.28	0.46
1:A:266:ARG:HB2	1:A:266:ARG:CZ	2.46	0.46
1:C:85:GLN:HE21	5:C:304:PEG:C4	2.28	0.46
1:D:106:SER:O	1:D:110:GLU:HG3	2.16	0.46
1:A:277:ALA:H	4:A:303:ACY:H3	1.79	0.46
1:C:70:SER:OG	2:C:301:BX9:O10	2.34	0.45
1:A:29:VAL:HG22	6:A:439:HOH:O	2.16	0.45
1:D:29:VAL:O	1:D:32:PRO:HD2	2.17	0.45
1:A:73:LYS:HB2	1:A:234:LYS:HE3	1.98	0.45
1:A:256:ARG:HH12	3:A:302:ACT:CH3	2.30	0.45
1:B:104:PRO:HB2	1:B:105:TRP:CE3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:70:SER:OG	2:D:301:BX9:O10	2.29	0.45
1:C:277:ALA:H	3:C:302:ACT:CH3	2.30	0.44
1:A:70:SER:OG	2:A:301:BX9:C1	2.65	0.44
1:C:291:GLY:N	6:C:405:HOH:O	2.50	0.44
1:D:225:VAL:HG13	1:D:229:TRP:CE3	2.52	0.44
1:A:70:SER:OG	2:A:301:BX9:O10	2.36	0.44
1:B:70:SER:OG	2:B:301:BX9:O11	2.35	0.44
1:D:202:PRO:HG2	1:D:203:GLN:OE1	2.18	0.43
1:B:225:VAL:CG1	1:B:229:TRP:HB2	2.49	0.43
1:C:235:THR:HG1	4:C:303:ACY:C	2.18	0.43
1:B:27:ASN:HB2	1:B:28:LEU:H	1.68	0.42
2:D:301:BX9:O10	4:D:303:ACY:H3	2.19	0.42
1:B:79:ALA:HB3	1:B:142:LEU:HD21	2.00	0.42
1:B:229:TRP:HE3	1:B:250:VAL:CG1	2.33	0.42
1:B:276:GLU:H	3:B:302:ACT:CH3	2.32	0.42
1:D:70:SER:OG	2:D:301:BX9:O11	2.36	0.42
1:C:49:MET:HG3	5:C:305:PEG:H22	2.02	0.42
1:C:35:LYS:HG3	1:C:36:LEU:N	2.34	0.42
1:C:70:SER:CB	4:C:303:ACY:H2	2.43	0.41
1:D:118:THR:OG1	1:D:121:GLU:HG3	2.20	0.41
1:C:70:SER:OG	2:C:301:BX9:C1	2.68	0.41
1:B:237:THR:OG1	4:B:303:ACY:O	2.25	0.41
1:A:209:ASP:OD2	6:A:402:HOH:O	2.22	0.41
1:C:49:MET:HG3	5:C:305:PEG:C2	2.50	0.41
1:D:276:GLU:H	3:D:302:ACT:H1	1.86	0.41
1:C:252:PRO:HB2	1:C:256:ARG:HB3	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:508:HOH:O	6:B:550:HOH:O[4_648]	1.67	0.53
6:A:545:HOH:O	6:B:554:HOH:O[4_648]	2.18	0.02

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	263/265 (99%)	257 (98%)	6 (2%)	0	100	100
1	B	261/265 (98%)	259 (99%)	2 (1%)	0	100	100
1	C	256/265 (97%)	253 (99%)	3 (1%)	0	100	100
1	D	263/265 (99%)	258 (98%)	5 (2%)	0	100	100
All	All	1043/1060 (98%)	1027 (98%)	16 (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	199/200 (100%)	194 (98%)	5 (2%)	47	41
1	B	198/200 (99%)	195 (98%)	3 (2%)	65	62
1	C	194/200 (97%)	187 (96%)	7 (4%)	35	26
1	D	199/200 (100%)	195 (98%)	4 (2%)	55	51
All	All	790/800 (99%)	771 (98%)	19 (2%)	49	43

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	VAL
1	A	91	LEU
1	A	109	SER
1	A	130	SER
1	A	233	ASP
1	B	130	SER
1	B	140	LYS
1	B	273	LYS

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Mol	Chain	Res	Type
1	C	35	LYS
1	C	38	GLN
1	C	49	MET
1	C	130	SER
1	C	142	LEU
1	C	168	GLU
1	C	225	VAL
1	D	130	SER
1	D	167	LEU
1	D	168	GLU
1	D	256	ARG

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

21 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
3	ACT	D	302	-	1,3,3	0.65	0	0,3,3	0.00	-
2	BX9	A	301	-	15,18,18	1.55	2 (13%)	15,24,24	1.87	6 (40%)
3	ACT	B	302	-	1,3,3	0.97	0	0,3,3	0.00	-
5	PEG	B	305	-	6,6,6	0.46	0	5,5,5	0.46	0
4	ACY	C	303	-	1,3,3	10.26	1 (100%)	0,3,3	0.00	-
4	ACY	A	304	-	1,3,3	12.98	1 (100%)	0,3,3	0.00	-
4	ACY	A	303	-	1,3,3	0.24	0	0,3,3	0.00	-
4	ACY	D	303	-	1,3,3	5.00	1 (100%)	0,3,3	0.00	-
5	PEG	A	306	-	6,6,6	0.55	0	5,5,5	0.42	0
4	ACY	B	303	-	1,3,3	15.49	1 (100%)	0,3,3	0.00	-
2	BX9	B	301	-	15,18,18	1.51	3 (20%)	15,24,24	1.92	3 (20%)
3	ACT	C	302	-	1,3,3	0.65	0	0,3,3	0.00	-
5	PEG	A	305	-	6,6,6	0.55	0	5,5,5	0.49	0
3	ACT	A	302	-	1,3,3	1.35	0	0,3,3	0.00	-
5	PEG	C	305	-	6,6,6	0.46	0	5,5,5	0.29	0
2	BX9	D	301	-	15,18,18	1.62	2 (13%)	15,24,24	1.71	3 (20%)
5	PEG	B	307	-	6,6,6	0.46	0	5,5,5	0.37	0
5	PEG	C	304	-	6,6,6	0.49	0	5,5,5	0.38	0
5	PEG	B	304	-	6,6,6	0.65	0	5,5,5	1.07	0
5	PEG	B	306	-	6,6,6	0.50	0	5,5,5	0.33	0
2	BX9	C	301	-	15,18,18	1.53	2 (13%)	15,24,24	1.55	5 (33%)

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	BX9	A	301	-	-	2/13/18/18	0/1/1/1
5	PEG	B	305	-	-	0/4/4/4	-
5	PEG	B	307	-	-	3/4/4/4	-
5	PEG	B	306	-	-	2/4/4/4	-
5	PEG	A	306	-	-	3/4/4/4	-
2	BX9	B	301	-	-	1/13/18/18	0/1/1/1
5	PEG	A	305	-	-	2/4/4/4	-
5	PEG	C	305	-	-	0/4/4/4	-
2	BX9	D	301	-	-	2/13/18/18	0/1/1/1
5	PEG	C	304	-	-	2/4/4/4	-
5	PEG	B	304	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BX9	C	301	-	-	2/13/18/18	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	303	ACY	CH3-C	-15.49	1.29	1.48
4	A	304	ACY	CH3-C	-12.98	1.32	1.48
4	C	303	ACY	CH3-C	-10.26	1.35	1.48
4	D	303	ACY	CH3-C	-5.00	1.42	1.48
2	D	301	BX9	C3-N2	4.28	1.43	1.34
2	A	301	BX9	C3-N2	4.07	1.43	1.34
2	C	301	BX9	C1-N2	-3.83	1.41	1.47
2	C	301	BX9	C3-N2	3.66	1.42	1.34
2	D	301	BX9	C1-N2	-3.59	1.42	1.47
2	B	301	BX9	C3-N2	3.47	1.41	1.34
2	B	301	BX9	O4-C3	-2.89	1.17	1.23
2	B	301	BX9	C1-N2	-2.77	1.43	1.47
2	A	301	BX9	C6-C5	2.64	1.56	1.53

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	BX9	C8-C5-C3	-3.89	103.16	109.42
2	B	301	BX9	C5-C3-N2	-3.84	112.30	116.42
2	B	301	BX9	C17-C16-C15	3.58	122.06	118.36
2	A	301	BX9	C5-C6-S7	3.48	118.69	114.05
2	A	301	BX9	C16-C15-C14	-3.16	118.63	122.83
2	B	301	BX9	C16-C15-C14	-3.07	118.75	122.83
2	A	301	BX9	C17-C16-C15	2.79	121.25	118.36
2	D	301	BX9	C16-C15-C14	-2.44	119.58	122.83
2	D	301	BX9	C17-C16-C15	2.35	120.79	118.36
2	C	301	BX9	C5-C3-N2	-2.31	113.94	116.42
2	C	301	BX9	O4-C3-C5	2.22	125.59	121.33
2	C	301	BX9	C5-C6-S7	-2.22	111.09	114.05
2	C	301	BX9	C17-C16-C15	2.19	120.63	118.36
2	A	301	BX9	O4-C3-C5	2.16	125.46	121.33
2	C	301	BX9	C16-C15-C14	-2.15	119.97	122.83
2	A	301	BX9	F18-C15-C16	2.13	122.16	118.54
2	A	301	BX9	O4-C3-N2	-2.09	119.06	122.93

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	BX9	C3-C5-C6-S7
2	A	301	BX9	C8-C5-C6-S7
2	D	301	BX9	C3-C5-C6-S7
2	D	301	BX9	C8-C5-C6-S7
2	C	301	BX9	C3-C5-C6-S7
2	C	301	BX9	C8-C5-C6-S7
5	B	304	PEG	O2-C3-C4-O4
5	B	307	PEG	O1-C1-C2-O2
5	B	306	PEG	O2-C3-C4-O4
5	A	305	PEG	O2-C3-C4-O4
5	C	304	PEG	O1-C1-C2-O2
5	A	306	PEG	O1-C1-C2-O2
5	A	306	PEG	O2-C3-C4-O4
5	A	306	PEG	C4-C3-O2-C2
5	C	304	PEG	C1-C2-O2-C3
5	B	306	PEG	C1-C2-O2-C3
5	B	307	PEG	C1-C2-O2-C3
5	A	305	PEG	C1-C2-O2-C3
2	B	301	BX9	C3-C5-C6-S7
5	B	307	PEG	C4-C3-O2-C2
5	B	304	PEG	C4-C3-O2-C2

There are no ring outliers.

18 monomers are involved in 50 short contacts:

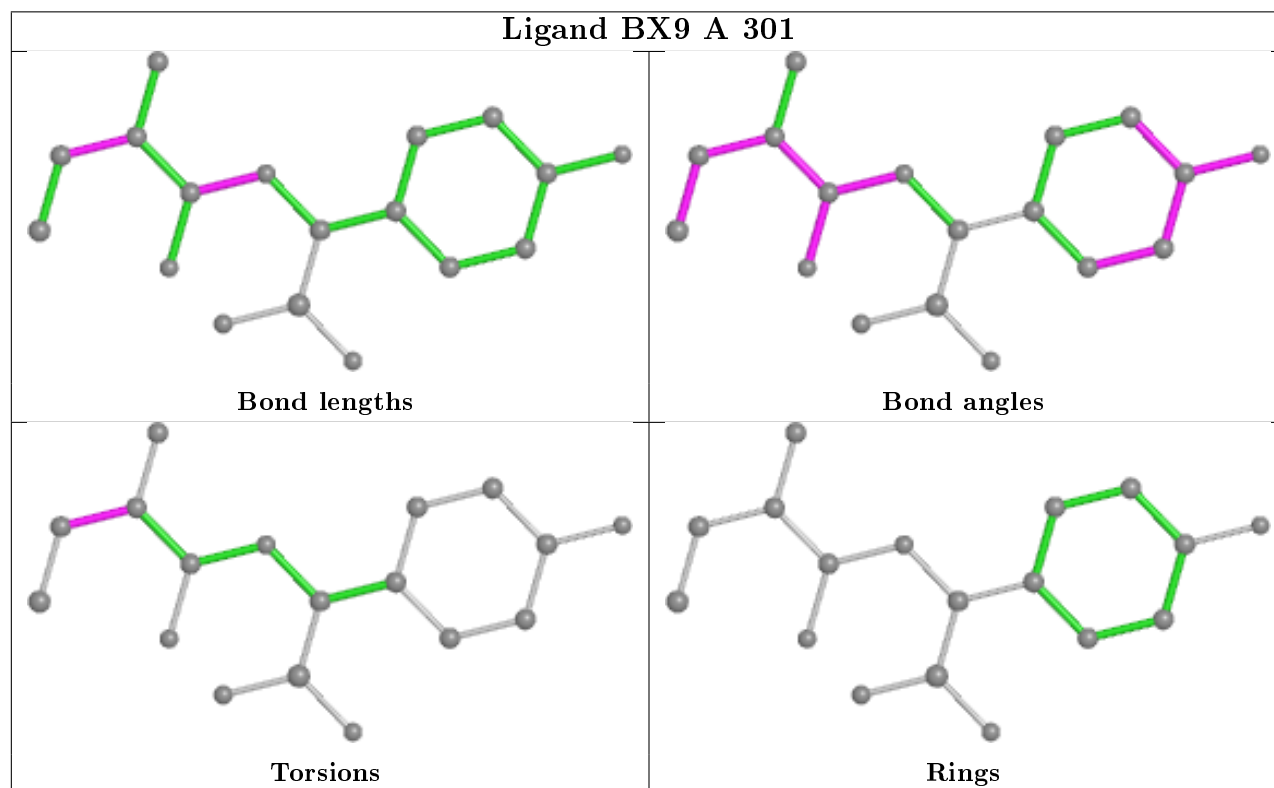
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	302	ACT	2	0
2	A	301	BX9	3	0
3	B	302	ACT	1	0
4	C	303	ACY	5	0
4	A	304	ACY	4	0
4	A	303	ACY	1	0
4	D	303	ACY	8	0
5	A	306	PEG	1	0
4	B	303	ACY	5	0
2	B	301	BX9	4	0
3	C	302	ACT	1	0
5	A	305	PEG	1	0
3	A	302	ACT	7	0
5	C	305	PEG	2	0
2	D	301	BX9	5	0

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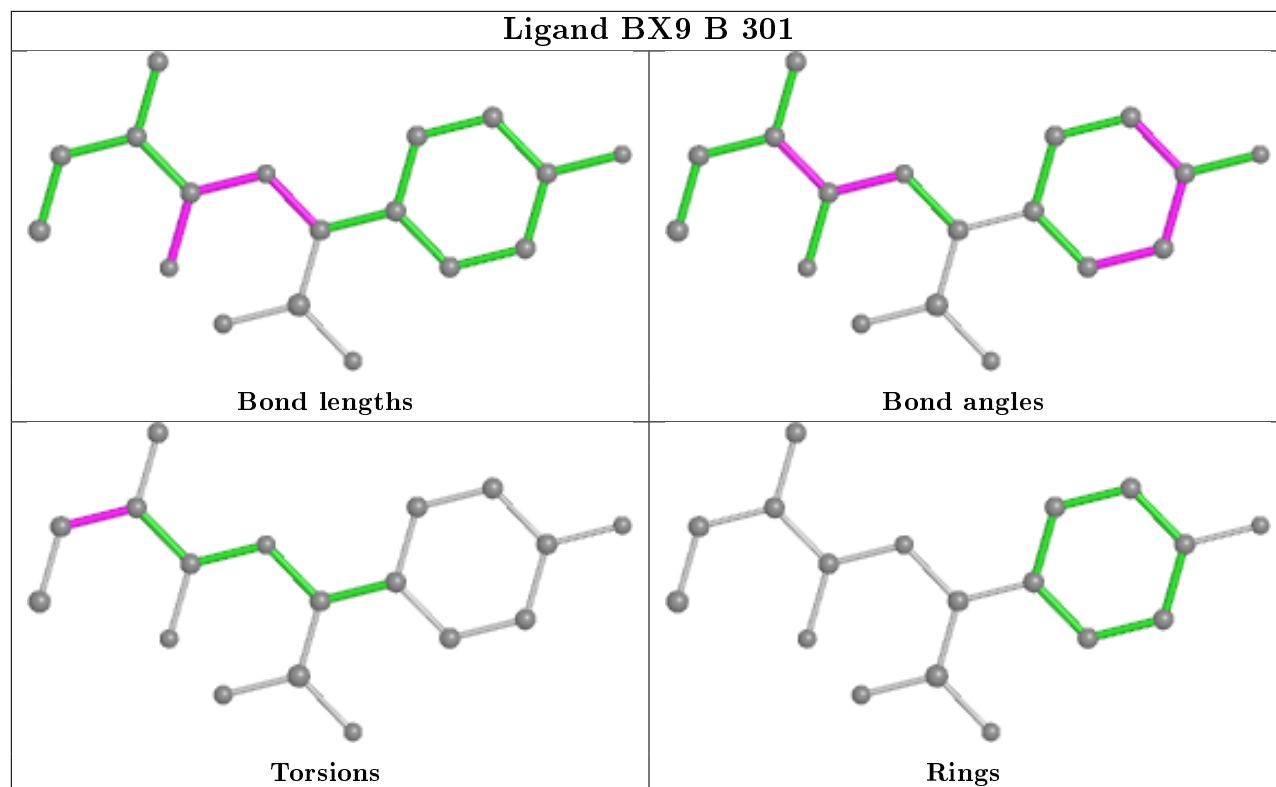
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	304	PEG	2	0
5	B	304	PEG	1	0
2	C	301	BX9	2	0

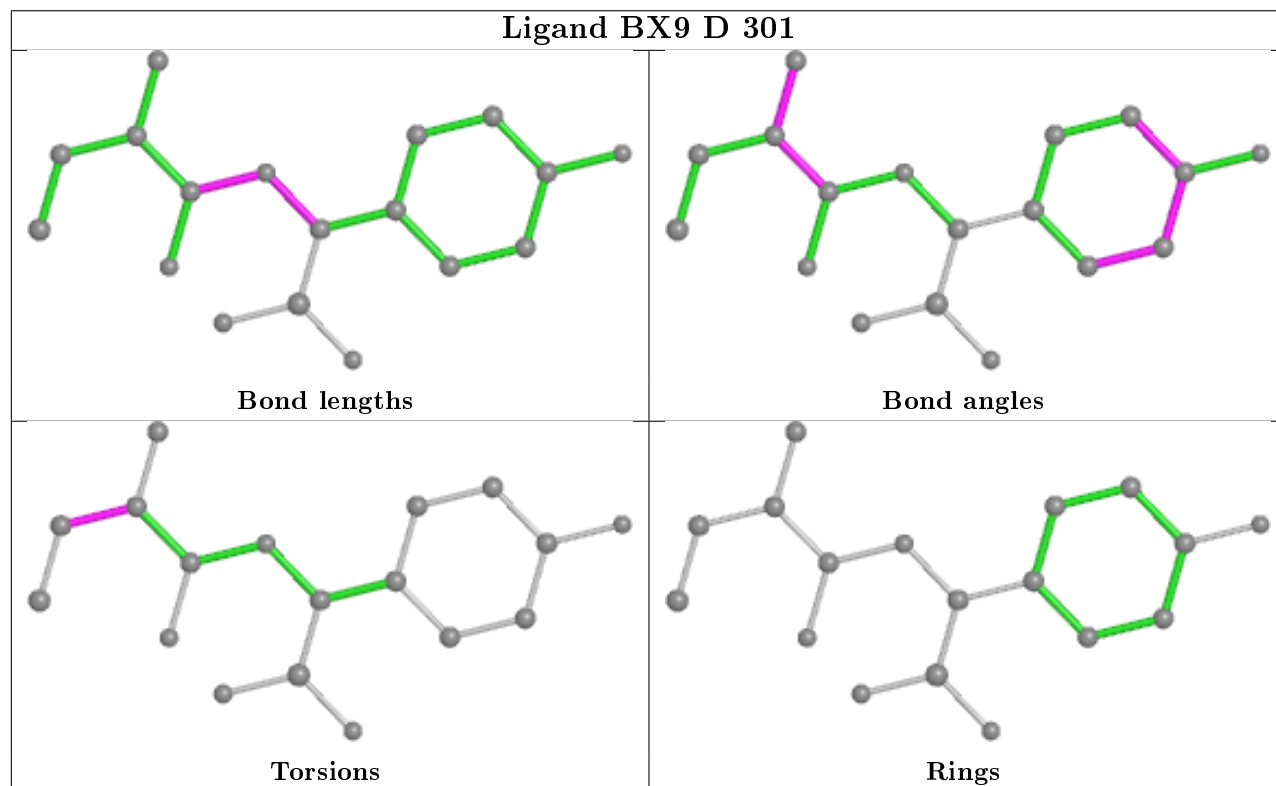
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

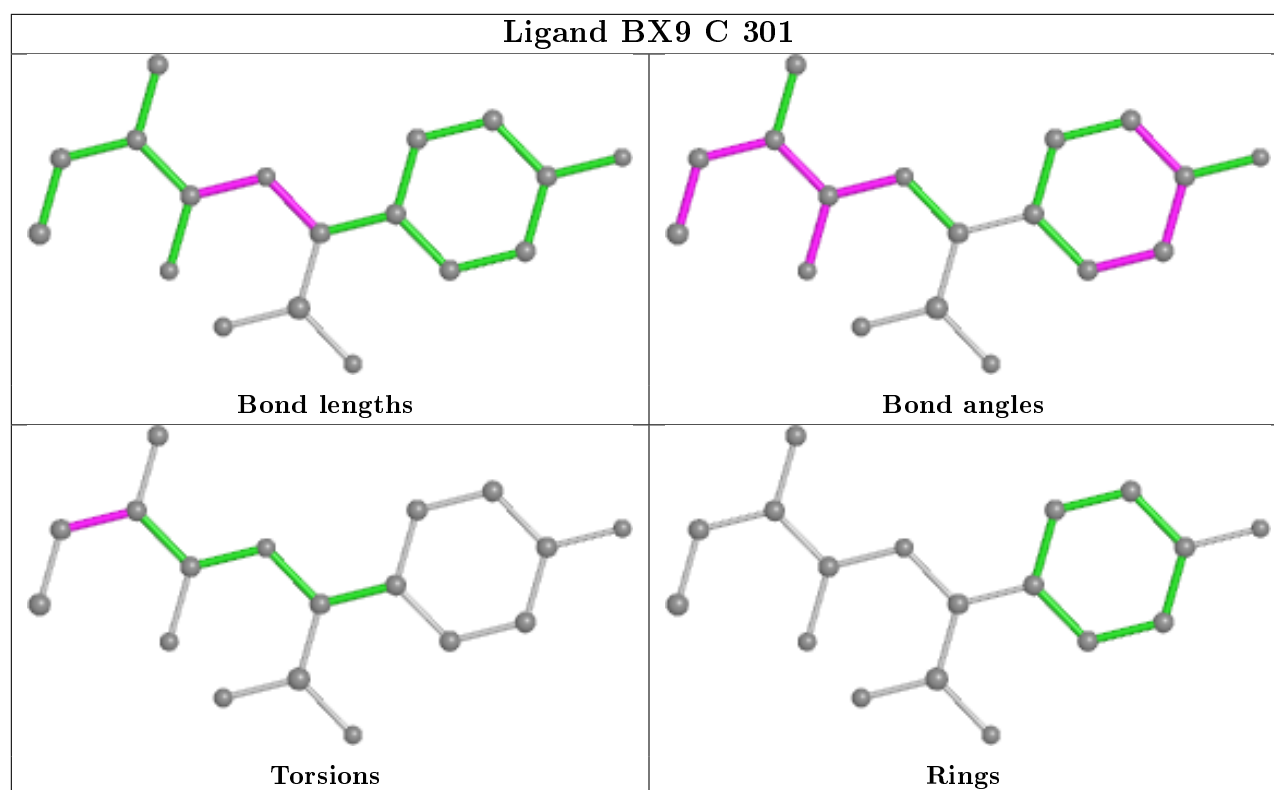


Ligand BX9 B 301



Ligand BX9 D 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	265/265 (100%)	-0.66	0	100 100	10, 17, 35, 53	0
1	B	263/265 (99%)	-0.67	0	100 100	7, 13, 32, 55	0
1	C	258/265 (97%)	-0.52	3 (1%)	79 81	10, 18, 48, 77	0
1	D	265/265 (100%)	-0.39	1 (0%)	92 93	16, 26, 52, 81	0
All	All	1051/1060 (99%)	-0.56	4 (0%)	92 93	7, 19, 44, 81	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	105	TRP	3.6
1	C	38	GLN	2.9
1	C	100	ASN	2.8
1	D	256	ARG	2.2

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 carbohydrates in this entry.

6.4 Ligands [i](#)

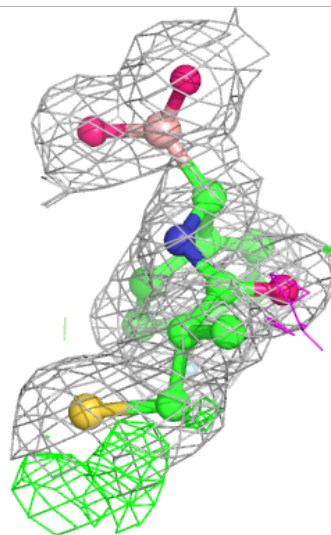
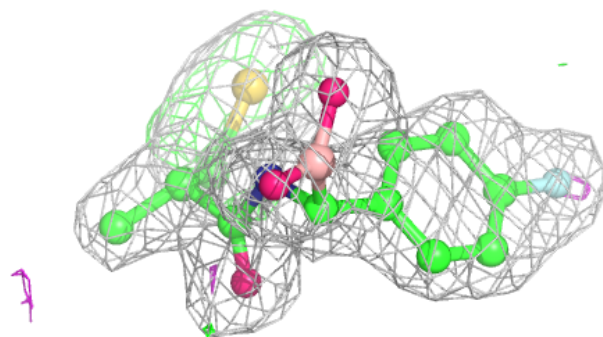
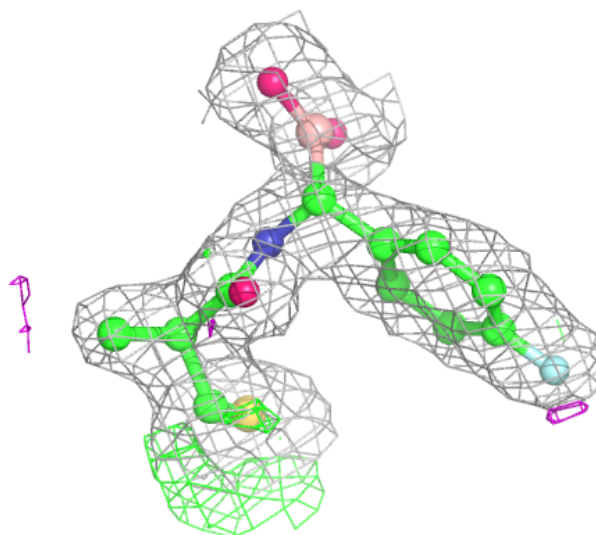
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	ACY	A	303	4/4	0.73	0.20	19,24,31,48	0
5	PEG	B	306	7/7	0.84	0.16	39,42,45,47	0
5	PEG	C	304	7/7	0.88	0.16	44,45,50,54	0
5	PEG	B	307	7/7	0.88	0.11	36,42,47,49	0
5	PEG	C	305	7/7	0.90	0.11	32,35,40,47	0
5	PEG	A	306	7/7	0.91	0.23	37,41,56,61	0
5	PEG	B	304	7/7	0.91	0.14	12,23,32,52	0
3	ACT	C	302	4/4	0.91	0.14	26,32,38,55	0
5	PEG	A	305	7/7	0.92	0.10	32,35,38,40	0
4	ACY	C	303	4/4	0.92	0.26	9,15,35,40	0
2	BX9	C	301	18/18	0.93	0.10	12,25,32,45	0
2	BX9	D	301	18/18	0.94	0.08	21,29,37,45	0
4	ACY	A	304	4/4	0.94	0.30	3,13,35,41	0
5	PEG	B	305	7/7	0.94	0.08	17,18,29,31	0
3	ACT	B	302	4/4	0.95	0.11	22,24,28,35	0
3	ACT	D	302	4/4	0.95	0.12	20,21,24,31	0
2	BX9	A	301	18/18	0.95	0.09	11,18,26,43	0
4	ACY	B	303	4/4	0.96	0.24	1,11,32,38	0
3	ACT	A	302	4/4	0.96	0.07	21,22,23,30	0
4	ACY	D	303	4/4	0.96	0.25	8,27,35,36	0
2	BX9	B	301	18/18	0.97	0.07	7,11,18,21	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.

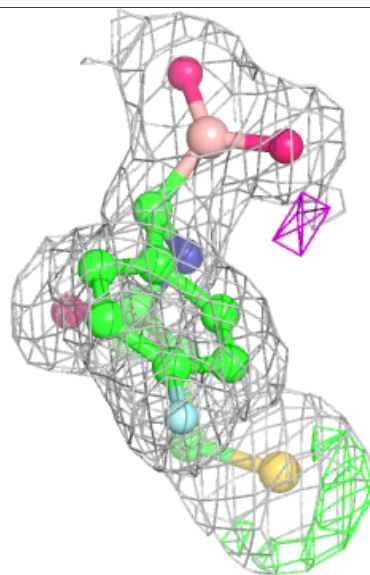
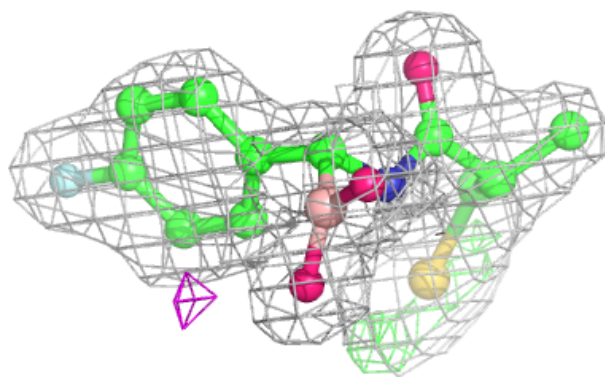
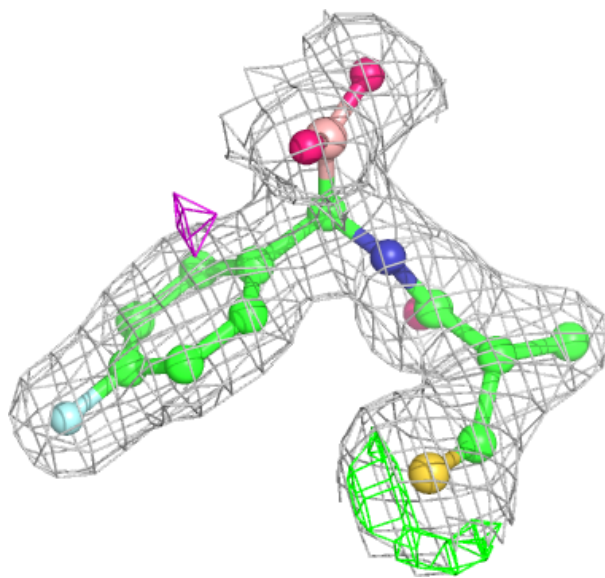
Electron density around BX9 C 301:

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



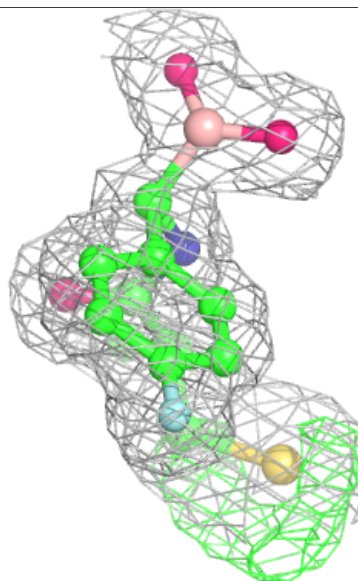
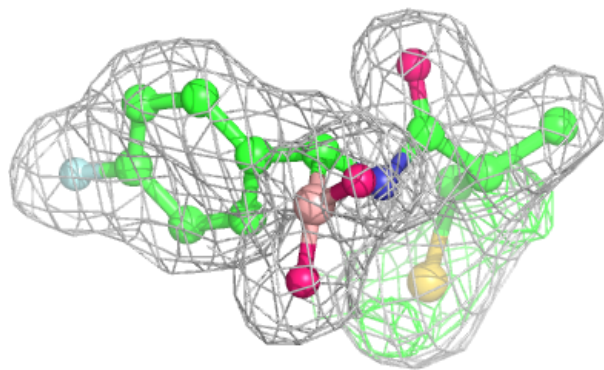
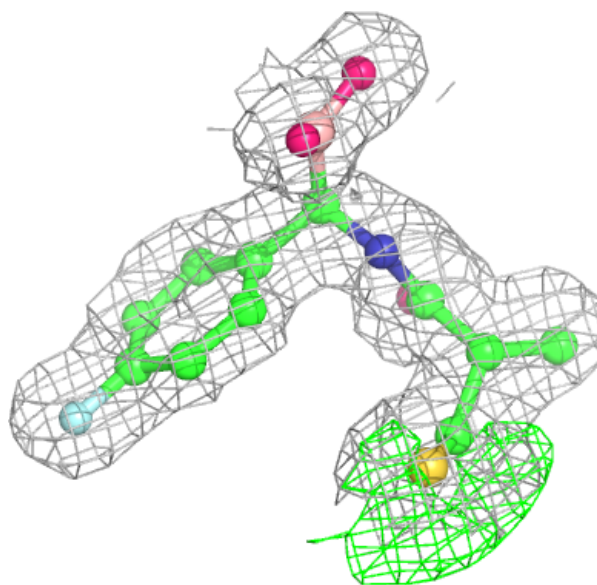
Electron density around BX9 D 301:

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



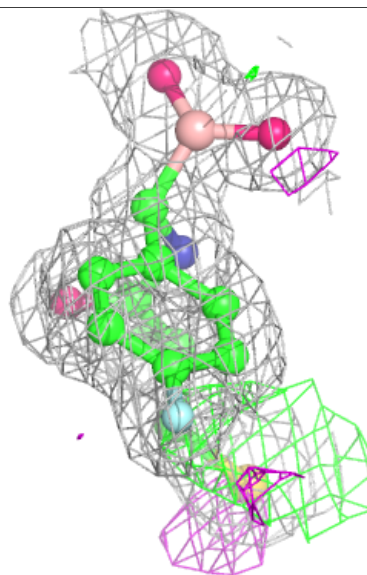
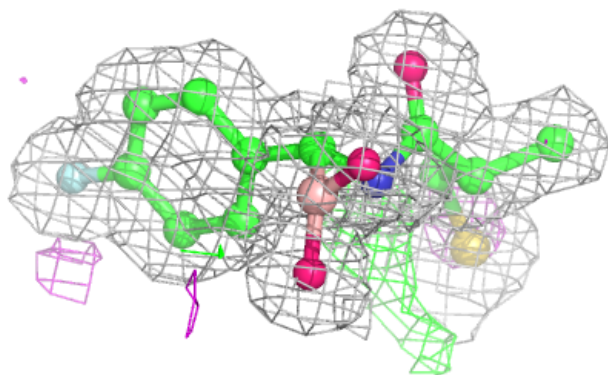
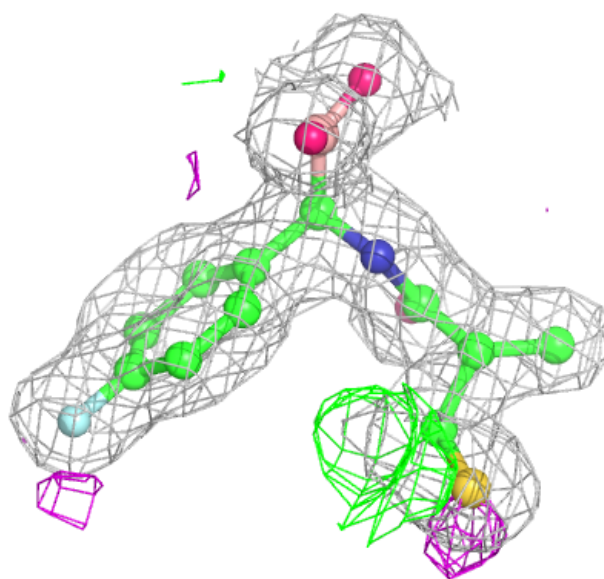
Electron density around BX9 A 301:

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



Electron density around BX9 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.