



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 25, 2020 – 02:24 PM BST

PDB ID : 4LOP  
Title : Structural basis of autoactivation of p38 alpha induced by TAB1 (Tetragonal crystal form)  
Authors : Chaikuad, A.; DeNicola, G.F.; Krojer, T.; Allerston, C.K.; von Delft, F.; Arrowsmith, C.H.; Edwards, A.M.; Bountra, C.; Marber, M.S.; Knapp, S.; Structural Genomics Consortium (SGC)  
Deposited on : 2013-07-13  
Resolution : 2.05 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.

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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 : 2.13  
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) : 2.13

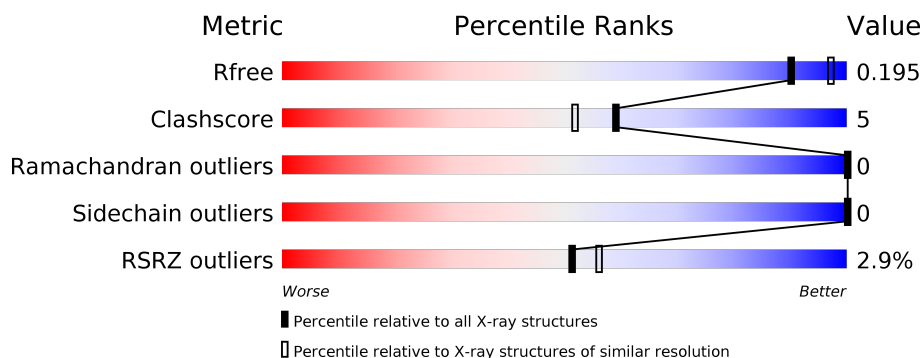
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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  $\geq 3$ , 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	361	<div> <div>2%</div> <div> <div></div> <div>88%</div> <div>9%</div> <div>•</div> </div> </div>
1	B	361	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>9%</div> <div>•</div> </div> </div>
1	C	361	<div> <div>4%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>6%</div> </div> </div>
1	D	361	<div> <div>•</div> <div> <div></div> <div>87%</div> <div>10%</div> <div>••</div> </div> </div>
2	K	29	<div> <div>3%</div> <div> <div></div> <div>55%</div> <div>17%</div> <div>28%</div> </div> </div>
2	L	29	<div> <div>7%</div> <div> <div></div> <div>72%</div> <div>28%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	M	29	<div><div></div><div>10%</div><div>41%</div><div>21%</div><div>38%</div></div>
2	N	29	<div><div></div><div>7%</div><div>55%</div><div>10%</div><div>34%</div></div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12902 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 Mitogen-activated protein kinase 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	4	0
			2844	1821	486	523	14			
1	B	350	Total	C	N	O	S	0	8	0
			2854	1826	488	527	13			
1	C	341	Total	C	N	O	S	0	5	0
			2708	1738	457	500	13			
1	D	351	Total	C	N	O	S	0	7	0
			2856	1829	486	527	14			

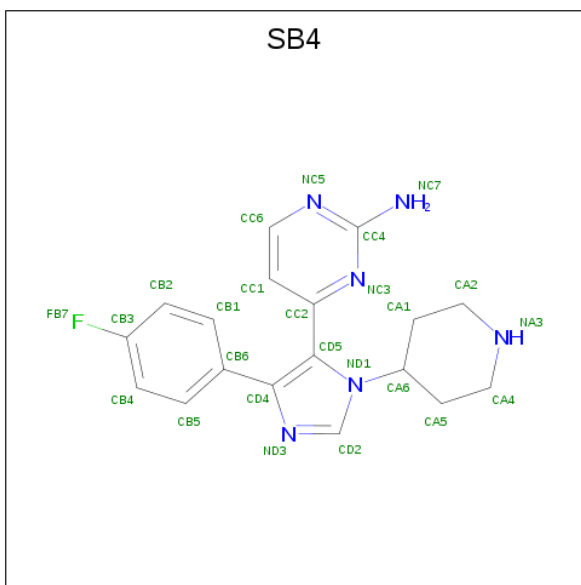
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	EXPRESSION TAG	UNP P47811
B	0	GLY	-	EXPRESSION TAG	UNP P47811
C	0	GLY	-	EXPRESSION TAG	UNP P47811
D	0	GLY	-	EXPRESSION TAG	UNP P47811

- Molecule 2 is a protein called TGF-beta-activated kinase 1 and MAP3K7-binding protein 1.

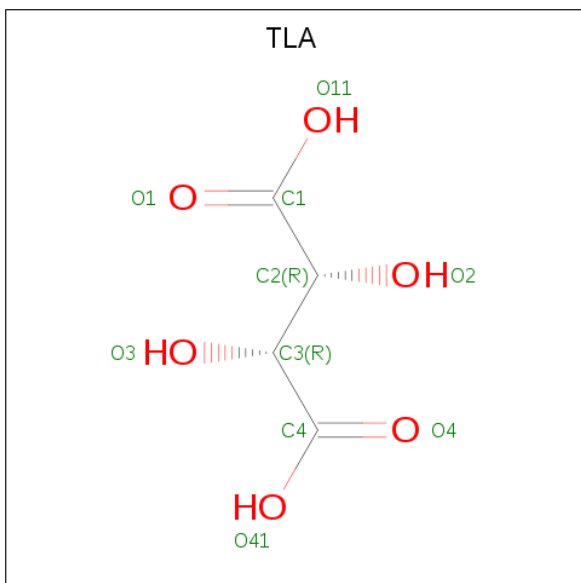
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	21	Total	C	N	O	S	0	0	0
			160	103	25	31	1			
2	L	21	Total	C	N	O	S	0	0	0
			158	101	25	31	1			
2	M	18	Total	C	N	O	S	0	0	0
			134	89	18	26	1			
2	N	19	Total	C	N	O	S	0	1	0
			143	96	19	27	1			

- Molecule 3 is 4-(4-FLUOROPHENYL)-1-(4-PIPERIDINYL)-5-(2-AMINO-4-PYRIMIDINYL)-IMIDAZOLE (three-letter code: SB4) (formula: C<sub>18</sub>H<sub>19</sub>FN<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 25	C 18	F 1	N 6	0	0
3	B	1	Total 25	C 18	F 1	N 6	0	0
3	C	1	Total 25	C 18	F 1	N 6	0	0
3	D	1	Total 25	C 18	F 1	N 6	0	0

- Molecule 4 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula:  $\text{C}_4\text{H}_6\text{O}_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	4	6		
4	D	1	Total	C	O	0	0
			10	4	6		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

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

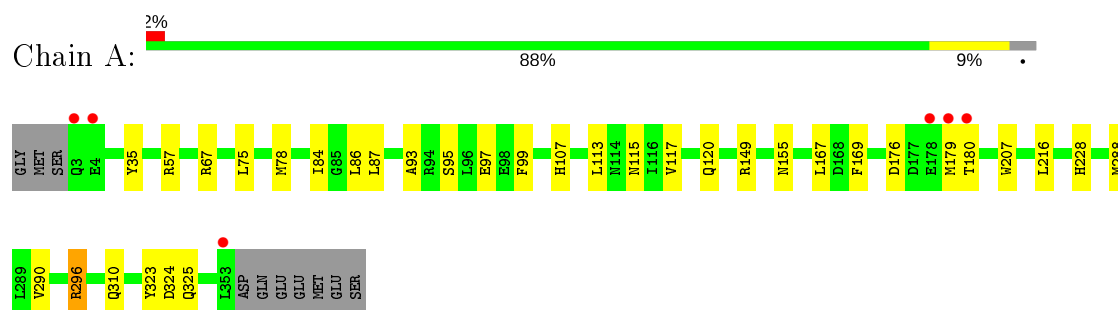
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	237	Total O 237 237	0	0
6	B	184	Total O 184 184	0	0
6	C	180	Total O 180 180	0	0
6	D	226	Total O 226 226	0	0
6	K	11	Total O 11 11	0	0
6	L	15	Total O 15 15	0	0
6	M	6	Total O 6 6	0	0
6	N	6	Total O 6 6	0	0

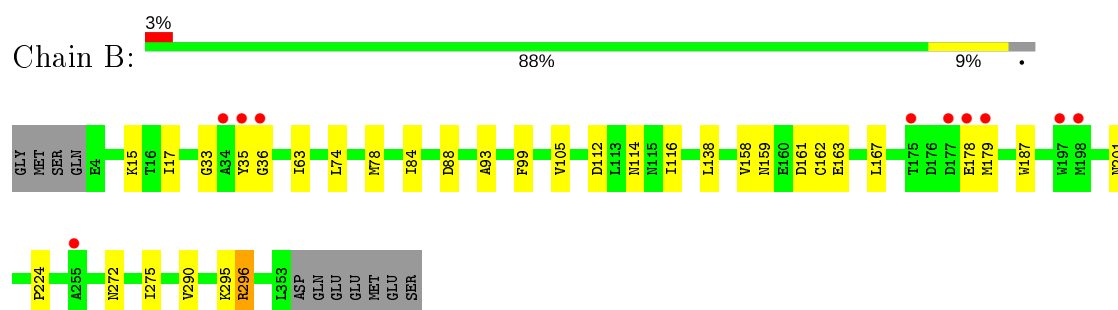
### 3 Residue-property plots [i](#)

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

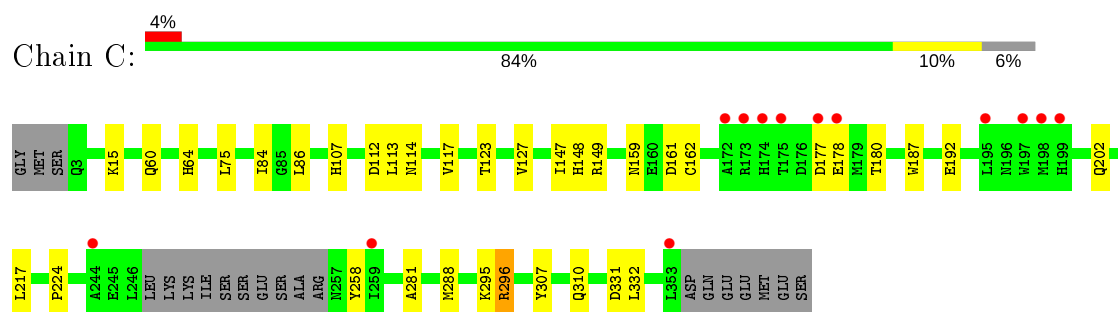
- Molecule 1: Mitogen-activated protein kinase 14



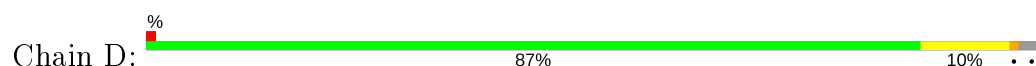
- Molecule 1: Mitogen-activated protein kinase 14



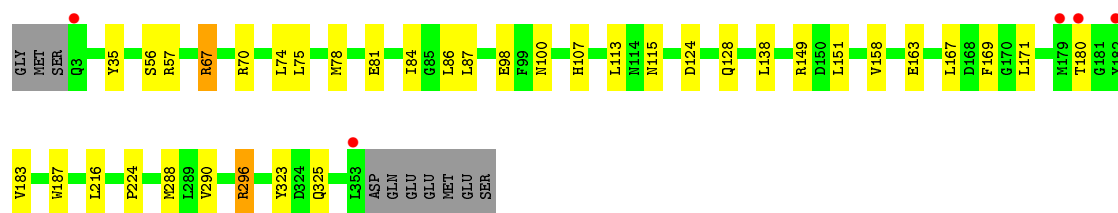
- Molecule 1: Mitogen-activated protein kinase 14



- Molecule 1: Mitogen-activated protein kinase 14



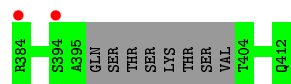




- Molecule 2: TGF-beta-activated kinase 1 and MAP3K7-binding protein 1



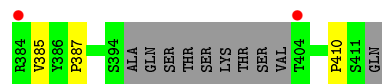
- Molecule 2: TGF-beta-activated kinase 1 and MAP3K7-binding protein 1



- Molecule 2: TGF-beta-activated kinase 1 and MAP3K7-binding protein 1



- Molecule 2: TGF-beta-activated kinase 1 and MAP3K7-binding protein 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.54Å 86.54Å 226.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	86.55 – 2.05 68.81 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.3 (86.55-2.05) 99.9 (68.81-2.05)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.76 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.148 , 0.188 0.161 , 0.195	Depositor DCC
$R_{free}$ test set	5308 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.9	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 47.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.022 for h,-k,-l	Xtriage
Reported twinning fraction	0.783 for H, K, L 0.217 for K, H, -L	Depositor
Outliers	0 of 103979 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	12902	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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: TLA, EDO, SB4

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	0.86	1/2923 (0.0%)	0.86	4/3969 (0.1%)
1	B	0.76	0/2928	0.77	2/3977 (0.1%)
1	C	0.78	0/2786	0.77	1/3791 (0.0%)
1	D	0.84	0/2944	0.89	7/3998 (0.2%)
2	K	0.71	0/163	0.84	0/221
2	L	0.74	0/161	0.74	0/218
2	M	0.67	0/137	0.67	0/188
2	N	0.52	0/149	0.70	0/205
All	All	0.80	1/12191 (0.0%)	0.82	14/16567 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	207	TRP	CD2-CE2	5.05	1.47	1.41

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	296	ARG	NE-CZ-NH2	-15.76	112.42	120.30
1	A	296	ARG	NE-CZ-NH2	-13.35	113.63	120.30
1	D	296	ARG	NE-CZ-NH1	12.69	126.65	120.30
1	A	296	ARG	NE-CZ-NH1	11.10	125.85	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	296	ARG	NE-CZ-NH2	-8.65	115.97	120.30
1	B	296	ARG	NE-CZ-NH1	7.40	124.00	120.30
1	D	149	ARG	NE-CZ-NH1	6.97	123.79	120.30
1	A	149	ARG	NE-CZ-NH2	6.22	123.41	120.30
1	D	70	ARG	NE-CZ-NH1	6.19	123.40	120.30
1	C	296	ARG	NE-CZ-NH2	-5.83	117.39	120.30
1	D	296	ARG	CD-NE-CZ	5.62	131.46	123.60
1	D	67	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	D	67	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	324	ASP	CB-CG-OD2	5.38	123.14	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	258	TYR	Peptide

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2844	0	2830	28	0
1	B	2854	0	2822	24	0
1	C	2708	0	2639	26	0
1	D	2856	0	2840	31	0
2	K	160	0	162	4	0
2	L	158	0	155	0	0
2	M	134	0	136	4	0
2	N	143	0	147	2	0
3	A	25	0	19	2	0
3	B	25	0	19	2	0
3	C	25	0	19	1	0
3	D	25	0	19	2	0
4	A	10	0	4	0	0
4	D	10	0	4	0	0
5	A	12	0	18	0	0
5	B	16	0	24	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	16	0	24	0	0
5	D	16	0	24	0	0
6	A	237	0	0	1	0
6	B	184	0	0	4	0
6	C	180	0	0	5	0
6	D	226	0	0	2	0
6	K	11	0	0	0	0
6	L	15	0	0	0	0
6	M	6	0	0	0	0
6	N	6	0	0	0	0
All	All	12902	0	11905	112	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 5.

All (112) 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:D:84:ILE:HG13	1:D:167:LEU:HB3	1.71	0.73
1:B:15:LYS:O	1:D:180[A]:THR:HG22	1.91	0.70
3:A:401:SB4:H6	3:A:401:SB4:NC3	2.05	0.70
1:A:180[B]:THR:HG22	1:C:15:LYS:HB3	1.71	0.69
1:A:176:ASP:HB2	1:A:179[A]:MET:SD	2.32	0.69
1:A:35:TYR:CE2	1:A:67:ARG:HD3	2.28	0.67
1:B:178:GLU:HB2	1:D:57:ARG:HE	1.61	0.66
1:D:288:MET:O	1:D:296:ARG:HD2	1.97	0.64
1:B:33[A]:GLY:N	1:B:36[A]:GLY:O	2.28	0.63
1:A:115:ASN:HB3	2:M:410:PRO:HG3	1.80	0.63
1:C:107:HIS:ND1	6:C:616:HOH:O	2.31	0.62
1:A:120:GLN:NE2	2:M:408:VAL:HG22	2.15	0.61
1:C:159[A]:ASN:ND2	6:C:556:HOH:O	2.33	0.61
1:D:87[A]:LEU:HD21	1:D:107:HIS:CE1	2.35	0.61
1:D:323:TYR:CE2	1:D:325:GLN:HG2	2.35	0.61
1:A:117:VAL:CG1	1:A:216:LEU:HD23	2.31	0.61
1:B:35[B]:TYR:N	1:B:35[B]:TYR:CD2	2.68	0.60
1:B:84:ILE:HG13	1:B:167:LEU:HB3	1.84	0.59
1:A:78:MET:HG3	1:A:169:PHE:CZ	2.38	0.58
1:B:161:ASP:O	1:B:162[A]:CYS:HB2	2.03	0.58
1:D:98:GLU:O	1:D:100:ASN:ND2	2.37	0.57
1:C:161:ASP:O	1:C:162[A]:CYS:HB2	2.05	0.57
1:B:63:ILE:N	1:B:63:ILE:HD13	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:ILE:HG13	1:B:158:VAL:HG21	1.86	0.56
1:D:81:GLU:HG3	6:D:710:HOH:O	2.05	0.56
1:C:112:ASP:OD1	1:C:114:ASN:HB3	2.06	0.56
1:D:35:TYR:CE2	1:D:67:ARG:HD3	2.41	0.55
1:A:180[B]:THR:HG22	1:C:15:LYS:O	2.06	0.55
1:A:113:LEU:O	1:A:117:VAL:HG22	2.07	0.55
1:A:290:VAL:O	1:A:296:ARG:HD3	2.07	0.55
1:C:60:GLN:OE1	1:C:64:HIS:CD2	2.60	0.55
1:A:323:TYR:CE2	1:A:325:GLN:HG2	2.42	0.55
1:B:33[A]:GLY:O	1:B:36[A]:GLY:O	2.25	0.55
1:B:33[A]:GLY:HA2	6:B:661:HOH:O	2.07	0.54
1:A:57:ARG:HD2	1:C:178:GLU:HB2	1.88	0.54
6:C:671:HOH:O	2:K:395:ALA:HB3	2.08	0.54
1:A:35:TYR:CZ	1:A:67:ARG:HD3	2.43	0.53
1:A:167:LEU:HD11	3:A:401:SB4:H52	1.91	0.53
1:A:95:SER:OG	1:A:97[A]:GLU:HG2	2.09	0.53
1:B:272:ASN:O	1:B:275:ILE:HD11	2.09	0.52
2:N:385:VAL:HG12	2:N:387:PRO:HD3	1.90	0.52
1:B:15:LYS:O	1:D:180[A]:THR:CG2	2.57	0.52
1:D:124:ASP:OD2	1:D:128:GLN:NE2	2.42	0.52
1:B:178:GLU:O	1:B:179:MET:CB	2.57	0.51
3:C:401:SB4:NC3	3:C:401:SB4:H6	2.25	0.51
2:K:385:VAL:HG12	2:K:387:PRO:HD3	1.93	0.51
1:D:187:TRP:CD1	1:D:224:PRO:HA	2.45	0.51
1:D:290:VAL:O	1:D:296:ARG:HD3	2.11	0.51
1:D:78:MET:HG3	1:D:169:PHE:CZ	2.46	0.51
1:A:288:MET:O	1:A:296:ARG:HD2	2.11	0.50
2:M:385:VAL:HG12	2:M:387:PRO:HD3	1.93	0.50
1:C:177:ASP:HB2	1:C:180:THR:CG2	2.43	0.49
1:C:281:ALA:HB2	1:C:307:TYR:CE1	2.48	0.49
1:D:56:SER:O	1:D:57:ARG:C	2.51	0.49
1:D:78:MET:HG3	1:D:169:PHE:CE1	2.48	0.49
1:D:87[A]:LEU:HD11	1:D:107:HIS:CE1	2.47	0.49
2:M:393:SER:O	2:M:394:SER:C	2.51	0.49
1:C:177:ASP:HB2	1:C:180:THR:HG21	1.96	0.48
3:B:401:SB4:H2	6:B:661:HOH:O	2.12	0.48
1:A:180[B]:THR:CG2	1:C:15:LYS:O	2.62	0.47
1:A:87:LEU:HD21	1:A:107:HIS:CE1	2.50	0.47
1:B:93:ALA:HB2	1:B:99:PHE:HA	1.96	0.47
1:B:112:ASP:OD1	1:B:114:ASN:HB3	2.15	0.47
1:D:323:TYR:CD2	1:D:325:GLN:HG2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:84:ILE:CG1	1:D:167:LEU:HB3	2.44	0.47
1:A:75:LEU:HB3	1:A:86:LEU:HG	1.97	0.47
3:B:401:SB4:NC3	3:B:401:SB4:H6	2.30	0.47
1:C:147:ILE:HG12	1:C:202:GLN:HG2	1.97	0.46
1:B:290:VAL:O	1:B:296:ARG:HD3	2.15	0.46
1:B:295:LYS:NZ	6:B:666:HOH:O	2.28	0.46
1:A:228:HIS:HD2	6:A:697:HOH:O	1.99	0.46
1:C:127:VAL:HG21	1:C:217:LEU:HD23	1.98	0.46
1:C:295:LYS:O	6:C:579:HOH:O	2.20	0.46
1:C:331:ASP:O	1:C:332:LEU:HD23	2.16	0.46
1:C:159[A]:ASN:OD1	1:C:161:ASP:OD1	2.35	0.45
1:C:187:TRP:CD1	1:C:224:PRO:HA	2.52	0.45
1:D:67:ARG:HD2	6:D:582:HOH:O	2.17	0.45
1:B:201:ASN:C	1:B:201:ASN:OD1	2.55	0.44
1:B:74:LEU:O	1:B:78:MET:HG2	2.16	0.44
1:C:148:HIS:O	1:C:149:ARG:HB2	2.17	0.44
1:D:74:LEU:O	1:D:78:MET:HG2	2.17	0.44
1:A:155:ASN:O	1:A:167:LEU:HD23	2.17	0.44
1:C:123:THR:HG22	2:K:390:VAL:HG11	1.98	0.44
1:A:117:VAL:HG12	1:A:216:LEU:HD23	2.00	0.43
1:D:87[A]:LEU:HA	1:D:87[A]:LEU:HD23	1.81	0.43
1:D:74:LEU:HD12	1:D:171:LEU:HD23	2.00	0.43
1:D:75:LEU:HB3	1:D:86:LEU:HG	1.99	0.43
1:A:84:ILE:HG13	1:A:167:LEU:HB3	2.00	0.43
1:C:288:MET:O	1:C:296:ARG:HD2	2.18	0.43
2:K:405:LEU:HD23	2:K:405:LEU:HA	1.70	0.42
1:D:167:LEU:HD21	3:D:401:SB4:H52	2.01	0.42
1:B:159[A]:ASN:OD1	1:B:163:GLU:HB3	2.20	0.42
1:A:93:ALA:HB2	1:A:99:PHE:HA	2.00	0.42
1:D:158:VAL:HA	1:D:163:GLU:O	2.20	0.42
1:A:93:ALA:CB	1:A:99:PHE:HA	2.50	0.42
1:D:115:ASN:HB3	2:N:410:PRO:HG2	2.02	0.42
1:D:138:LEU:HD21	1:D:151:LEU:HD11	2.01	0.42
1:C:310:GLN:O	1:C:310:GLN:CG	2.68	0.41
1:A:78:MET:HG3	1:A:169:PHE:CE2	2.55	0.41
1:D:113:LEU:HD13	1:D:216:LEU:HD21	2.02	0.41
1:A:84:ILE:CG1	1:A:167:LEU:HB3	2.50	0.41
1:C:84:ILE:CD1	6:C:535:HOH:O	2.68	0.41
6:B:595:HOH:O	3:D:401:SB4:H22	2.20	0.41
1:C:75:LEU:HB3	1:C:86:LEU:HG	2.03	0.41
1:C:113:LEU:O	1:C:117[A]:VAL:HG22	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:192:GLU:OE2	1:C:296:ARG:NH2	2.49	0.41
1:B:17:ILE:HG13	1:D:183:VAL:HG11	2.02	0.41
1:A:310:GLN:O	1:A:310:GLN:HG2	2.21	0.41
1:B:138:LEU:HA	1:B:138:LEU:HD12	1.77	0.40
1:B:88:ASP:HB3	1:B:105:VAL:HB	2.04	0.40
1:D:113:LEU:HA	1:D:113:LEU:HD23	1.96	0.40
1:B:187:TRP:CD1	1:B:224:PRO:HA	2.57	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	353/361 (98%)	343 (97%)	10 (3%)	0	100	100
1	B	356/361 (99%)	344 (97%)	12 (3%)	0	100	100
1	C	342/361 (95%)	333 (97%)	9 (3%)	0	100	100
1	D	356/361 (99%)	349 (98%)	7 (2%)	0	100	100
2	K	17/29 (59%)	16 (94%)	1 (6%)	0	100	100
2	L	17/29 (59%)	17 (100%)	0	0	100	100
2	M	14/29 (48%)	14 (100%)	0	0	100	100
2	N	16/29 (55%)	15 (94%)	1 (6%)	0	100	100
All	All	1471/1560 (94%)	1431 (97%)	40 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	311/319 (98%)	311 (100%)	0	100	100
1	B	309/319 (97%)	309 (100%)	0	100	100
1	C	287/319 (90%)	287 (100%)	0	100	100
1	D	313/319 (98%)	313 (100%)	0	100	100
2	K	20/28 (71%)	20 (100%)	0	100	100
2	L	19/28 (68%)	19 (100%)	0	100	100
2	M	18/28 (64%)	18 (100%)	0	100	100
2	N	19/28 (68%)	19 (100%)	0	100	100
All	All	1296/1388 (93%)	1296 (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. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	100	ASN

### 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 monosaccharides 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$
5	EDO	A	405	-	3,3,3	1.02	0	2,2,2	0.07	0
5	EDO	D	403	-	3,3,3	0.26	0	2,2,2	0.43	0
3	SB4	A	401	-	24,28,28	3.10	5 (20%)	29,39,39	3.21	12 (41%)
3	SB4	C	401	-	24,28,28	3.25	6 (25%)	29,39,39	4.03	12 (41%)
5	EDO	B	404	-	3,3,3	0.48	0	2,2,2	0.16	0
5	EDO	C	404	-	3,3,3	0.63	0	2,2,2	0.19	0
5	EDO	B	402	-	3,3,3	0.40	0	2,2,2	0.55	0
5	EDO	D	404	-	3,3,3	0.39	0	2,2,2	0.46	0
5	EDO	A	403	-	3,3,3	0.23	0	2,2,2	0.97	0
5	EDO	D	405	-	3,3,3	0.61	0	2,2,2	0.15	0
5	EDO	B	405	-	3,3,3	0.39	0	2,2,2	0.45	0
3	SB4	B	401	-	24,28,28	3.09	5 (20%)	29,39,39	3.20	15 (51%)
5	EDO	A	404	-	3,3,3	0.53	0	2,2,2	0.94	0
5	EDO	C	403	-	3,3,3	0.38	0	2,2,2	1.07	0
4	TLA	A	402	-	3,9,9	0.96	0	6,12,12	1.17	0
5	EDO	C	402	-	3,3,3	0.76	0	2,2,2	0.25	0
5	EDO	D	406	-	3,3,3	1.03	0	2,2,2	0.32	0
3	SB4	D	401	-	24,28,28	3.04	7 (29%)	29,39,39	5.16	18 (62%)
4	TLA	D	402	-	3,9,9	0.50	0	6,12,12	1.62	1 (16%)
5	EDO	C	405	-	3,3,3	0.50	0	2,2,2	0.52	0
5	EDO	B	403	-	3,3,3	0.57	0	2,2,2	0.74	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
5	EDO	A	405	-	-	1/1/1/1	-
5	EDO	D	403	-	-	1/1/1/1	-
3	SB4	A	401	-	-	2/5/20/20	0/4/4/4
3	SB4	C	401	-	-	2/5/20/20	0/4/4/4
5	EDO	B	404	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	C	404	-	-	1/1/1/1	-
5	EDO	B	402	-	-	0/1/1/1	-
5	EDO	D	404	-	-	1/1/1/1	-
5	EDO	A	403	-	-	1/1/1/1	-
5	EDO	D	405	-	-	1/1/1/1	-
5	EDO	B	405	-	-	1/1/1/1	-
3	SB4	B	401	-	-	1/5/20/20	0/4/4/4
5	EDO	A	404	-	-	1/1/1/1	-
5	EDO	C	403	-	-	1/1/1/1	-
4	TLA	A	402	-	-	0/4/12/12	-
5	EDO	C	402	-	-	0/1/1/1	-
5	EDO	D	406	-	-	1/1/1/1	-
3	SB4	D	401	-	-	2/5/20/20	0/4/4/4
4	TLA	D	402	-	-	0/4/12/12	-
5	EDO	C	405	-	-	0/1/1/1	-
5	EDO	B	403	-	-	1/1/1/1	-

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	401	SB4	CB6-CD4	-13.06	1.34	1.49
3	A	401	SB4	CB6-CD4	-12.41	1.35	1.49
3	B	401	SB4	CB6-CD4	-11.49	1.36	1.49
3	D	401	SB4	CB6-CD4	-11.04	1.36	1.49
3	B	401	SB4	CD5-CC2	-7.27	1.37	1.49
3	A	401	SB4	CD5-CC2	-6.75	1.38	1.49
3	C	401	SB4	CD5-CC2	-6.01	1.39	1.49
3	D	401	SB4	CD5-CC2	-5.69	1.39	1.49
3	D	401	SB4	CC6-NC5	4.99	1.45	1.34
3	B	401	SB4	CD5-CD4	-3.95	1.34	1.43
3	C	401	SB4	CD5-CD4	-3.43	1.35	1.43
3	D	401	SB4	CC2-NC3	-2.95	1.30	1.34
3	D	401	SB4	CD5-CD4	-2.95	1.36	1.43
3	A	401	SB4	CD5-CD4	-2.79	1.37	1.43
3	C	401	SB4	CC6-NC5	2.49	1.39	1.34
3	D	401	SB4	CA5-CA4	2.41	1.60	1.52
3	C	401	SB4	CA5-CA4	2.38	1.60	1.52
3	B	401	SB4	CD2-ND3	-2.36	1.31	1.35
3	B	401	SB4	CC6-NC5	2.22	1.39	1.34
3	A	401	SB4	CB4-CB3	2.21	1.41	1.37
3	D	401	SB4	CA5-CA6	-2.17	1.47	1.52
3	A	401	SB4	CC6-NC5	2.14	1.39	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	401	SB4	CA1-CA2	-2.07	1.45	1.52

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	401	SB4	CC2-NC3-CC4	17.50	126.32	116.34
3	C	401	SB4	CC2-NC3-CC4	16.69	125.86	116.34
3	B	401	SB4	CC2-NC3-CC4	12.49	123.46	116.34
3	A	401	SB4	CC2-NC3-CC4	10.67	122.43	116.34
3	D	401	SB4	CA4-CA5-CA6	9.96	118.23	110.44
3	D	401	SB4	CC6-CC1-CC2	7.94	124.71	117.22
3	D	401	SB4	CA2-CA1-CA6	-7.82	104.32	110.44
3	A	401	SB4	CA2-CA1-CA6	-7.48	104.59	110.44
3	D	401	SB4	CD4-CD5-ND1	-6.97	104.52	108.14
3	D	401	SB4	CC1-CC2-NC3	-6.34	113.75	121.97
3	C	401	SB4	CD4-CD5-ND1	-5.91	105.07	108.14
3	C	401	SB4	CC1-CC2-NC3	-5.73	114.55	121.97
3	D	401	SB4	CC1-CC6-NC5	-5.55	117.07	123.96
3	A	401	SB4	CD4-CD5-ND1	-4.89	105.60	108.14
3	D	401	SB4	CA1-CA2-NA3	-4.75	101.03	110.64
3	D	401	SB4	NC7-CC4-NC5	4.74	122.25	117.44
3	C	401	SB4	CA2-CA1-CA6	-4.65	106.80	110.44
3	B	401	SB4	NC5-CC4-NC3	-4.44	120.29	125.70
3	A	401	SB4	CD5-CC2-NC3	3.94	122.35	115.62
3	C	401	SB4	NC5-CC4-NC3	-3.85	121.01	125.70
3	C	401	SB4	CA4-CA5-CA6	-3.73	107.52	110.44
3	A	401	SB4	CC1-CC2-NC3	-3.72	117.15	121.97
3	D	401	SB4	CD5-CC2-NC3	3.66	121.87	115.62
3	A	401	SB4	CC6-CC1-CC2	3.56	120.58	117.22
3	B	401	SB4	CC1-CC2-NC3	-3.55	117.37	121.97
3	B	401	SB4	CC1-CC6-NC5	-3.54	119.56	123.96
3	C	401	SB4	CA1-CA2-NA3	-3.53	103.49	110.64
3	A	401	SB4	NC5-CC4-NC3	-3.53	121.40	125.70
3	A	401	SB4	CC1-CC6-NC5	-3.52	119.58	123.96
4	D	402	TLA	C1-C2-C3	-3.41	105.77	113.11
3	B	401	SB4	CC6-NC5-CC4	3.40	119.72	116.24
3	C	401	SB4	CC6-CC1-CC2	3.31	120.34	117.22
3	C	401	SB4	CD5-CC2-NC3	3.29	121.24	115.62
3	B	401	SB4	CD4-CD5-ND1	-3.00	106.58	108.14
3	B	401	SB4	CB1-CB2-CB3	2.99	121.46	118.36
3	C	401	SB4	CB5-CB4-CB3	2.98	121.44	118.36
3	B	401	SB4	CD5-CC2-NC3	2.96	120.69	115.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	401	SB4	CB5-CB4-CB3	2.96	121.42	118.36
3	B	401	SB4	CA2-CA1-CA6	2.94	112.74	110.44
3	D	401	SB4	CA5-CA6-CA1	-2.88	104.64	110.16
3	C	401	SB4	NC7-CC4-NC5	2.83	120.31	117.44
3	B	401	SB4	CB4-CB3-CB2	-2.80	119.11	122.83
3	B	401	SB4	NC7-CC4-NC5	2.79	120.27	117.44
3	D	401	SB4	CB1-CB2-CB3	2.54	120.99	118.36
3	A	401	SB4	CC6-NC5-CC4	2.52	118.82	116.24
3	D	401	SB4	CB4-CB3-CB2	-2.52	119.47	122.83
3	B	401	SB4	CD2-ND3-CD4	2.50	107.75	102.99
3	D	401	SB4	CA5-CA4-NA3	2.49	115.68	110.64
3	D	401	SB4	NC5-CC4-NC3	-2.42	122.75	125.70
3	B	401	SB4	CC6-CC1-CC2	2.40	119.49	117.22
3	A	401	SB4	CB4-CB3-CB2	-2.38	119.66	122.83
3	A	401	SB4	FB7-CB3-CB2	2.36	122.55	118.54
3	B	401	SB4	CA5-CA4-NA3	-2.34	105.91	110.64
3	C	401	SB4	CB4-CB3-CB2	-2.32	119.74	122.83
3	D	401	SB4	CD5-ND1-CA6	2.24	131.17	126.72
3	D	401	SB4	CD2-ND3-CD4	2.24	107.25	102.99
3	A	401	SB4	NC7-CC4-NC3	2.16	120.62	117.25
3	B	401	SB4	CD5-ND1-CA6	2.03	130.74	126.72

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	404	EDO	O1-C1-C2-O2
5	A	405	EDO	O1-C1-C2-O2
5	D	403	EDO	O1-C1-C2-O2
5	D	406	EDO	O1-C1-C2-O2
3	A	401	SB4	CB1-CB6-CD4-ND3
5	D	404	EDO	O1-C1-C2-O2
5	B	403	EDO	O1-C1-C2-O2
5	C	403	EDO	O1-C1-C2-O2
3	C	401	SB4	CB1-CB6-CD4-ND3
3	D	401	SB4	CB1-CB6-CD4-ND3
5	A	403	EDO	O1-C1-C2-O2
3	B	401	SB4	CB1-CB6-CD4-ND3
5	B	405	EDO	O1-C1-C2-O2
5	C	404	EDO	O1-C1-C2-O2
3	A	401	SB4	CB5-CB6-CD4-ND3
3	C	401	SB4	CB5-CB6-CD4-ND3

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Mol	Chain	Res	Type	Atoms
3	D	401	SB4	CB5-CB6-CD4-ND3
5	D	405	EDO	O1-C1-C2-O2

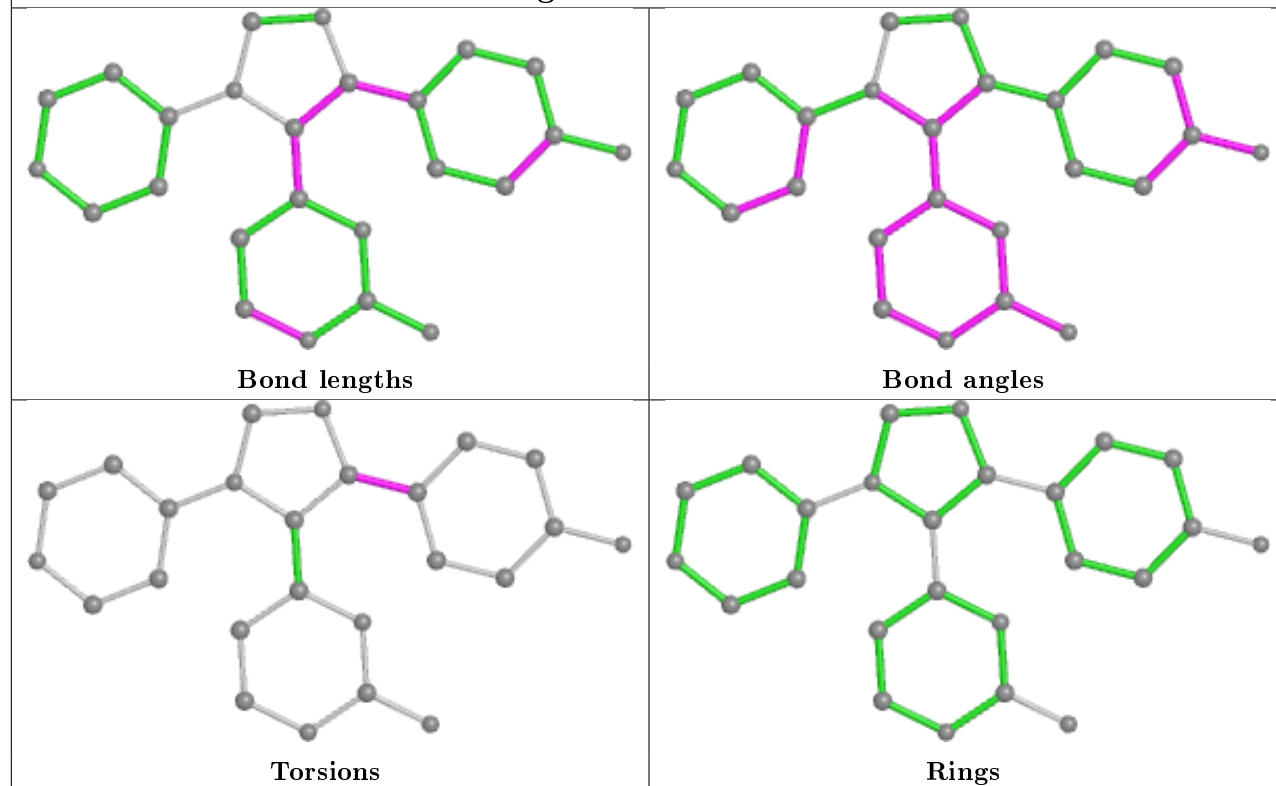
There are no ring outliers.

4 monomers are involved in 7 short contacts:

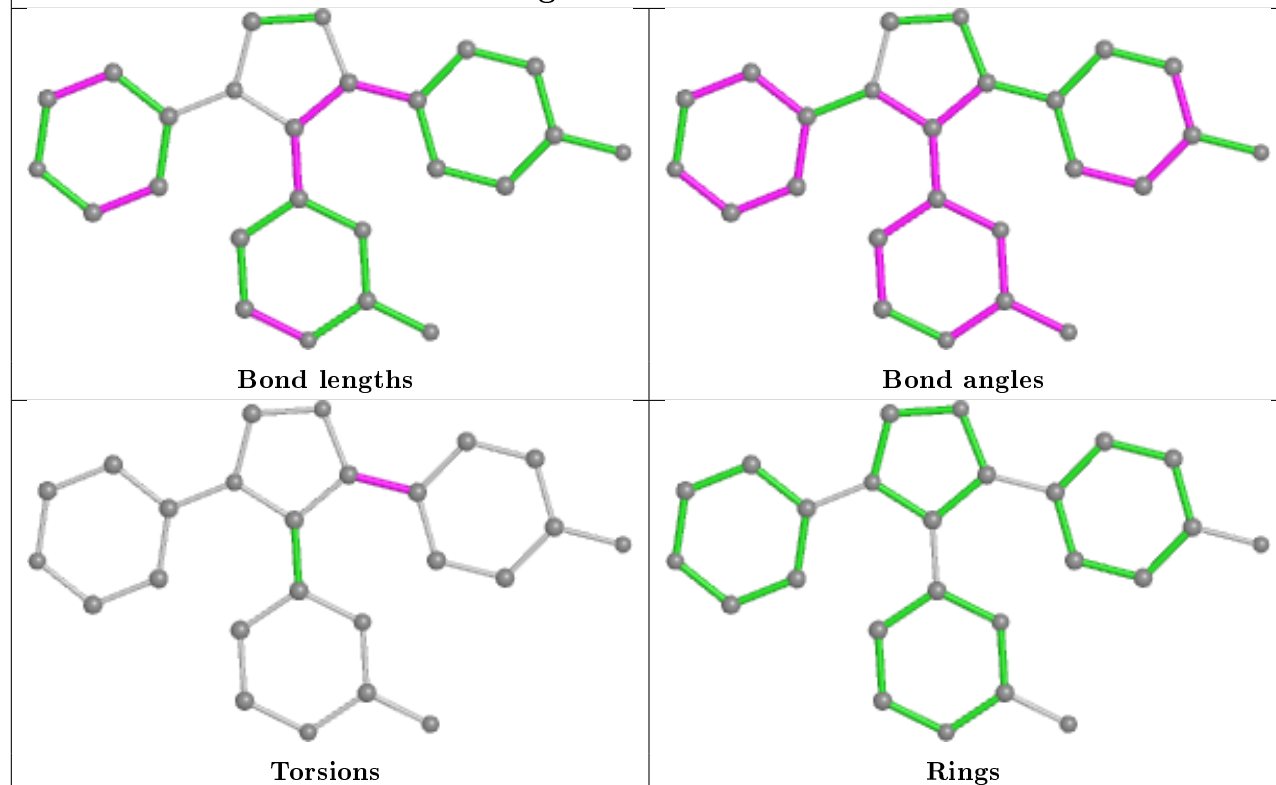
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	SB4	2	0
3	C	401	SB4	1	0
3	B	401	SB4	2	0
3	D	401	SB4	2	0

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.

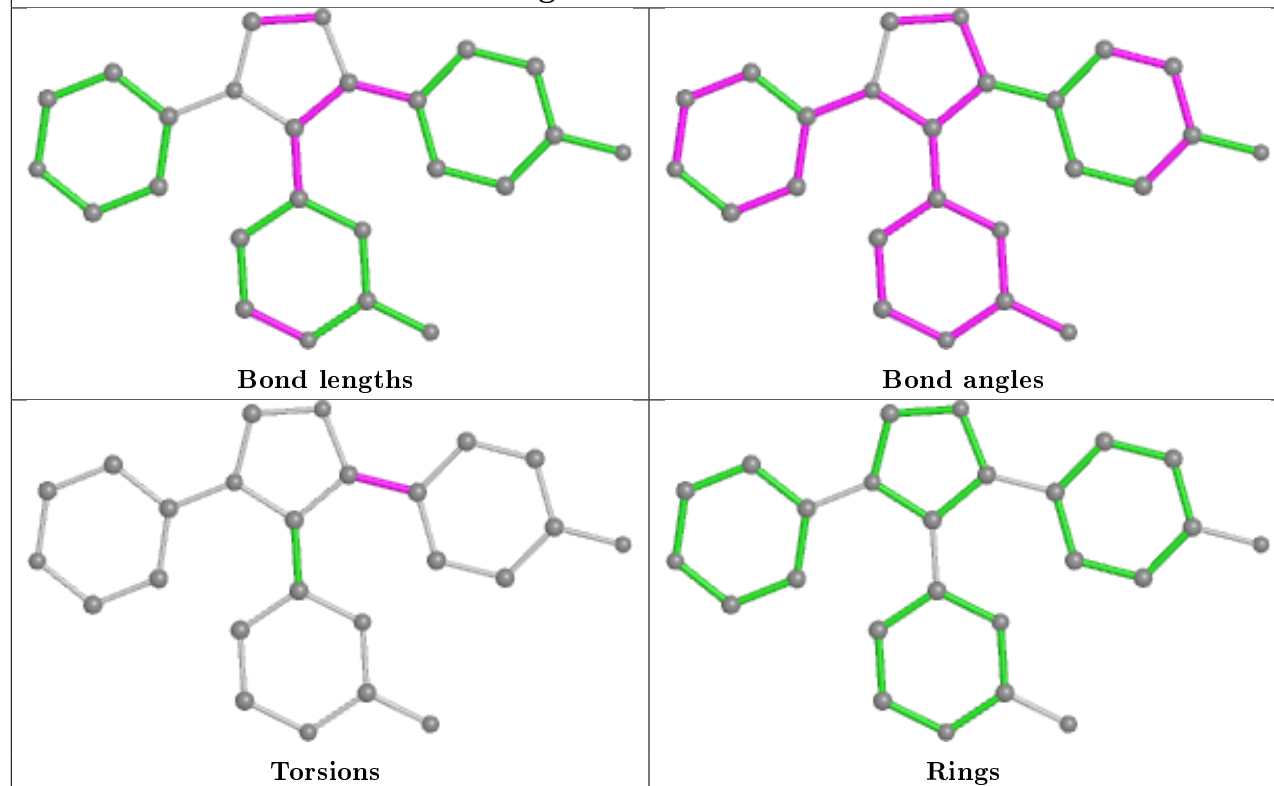
## Ligand SB4 A 401



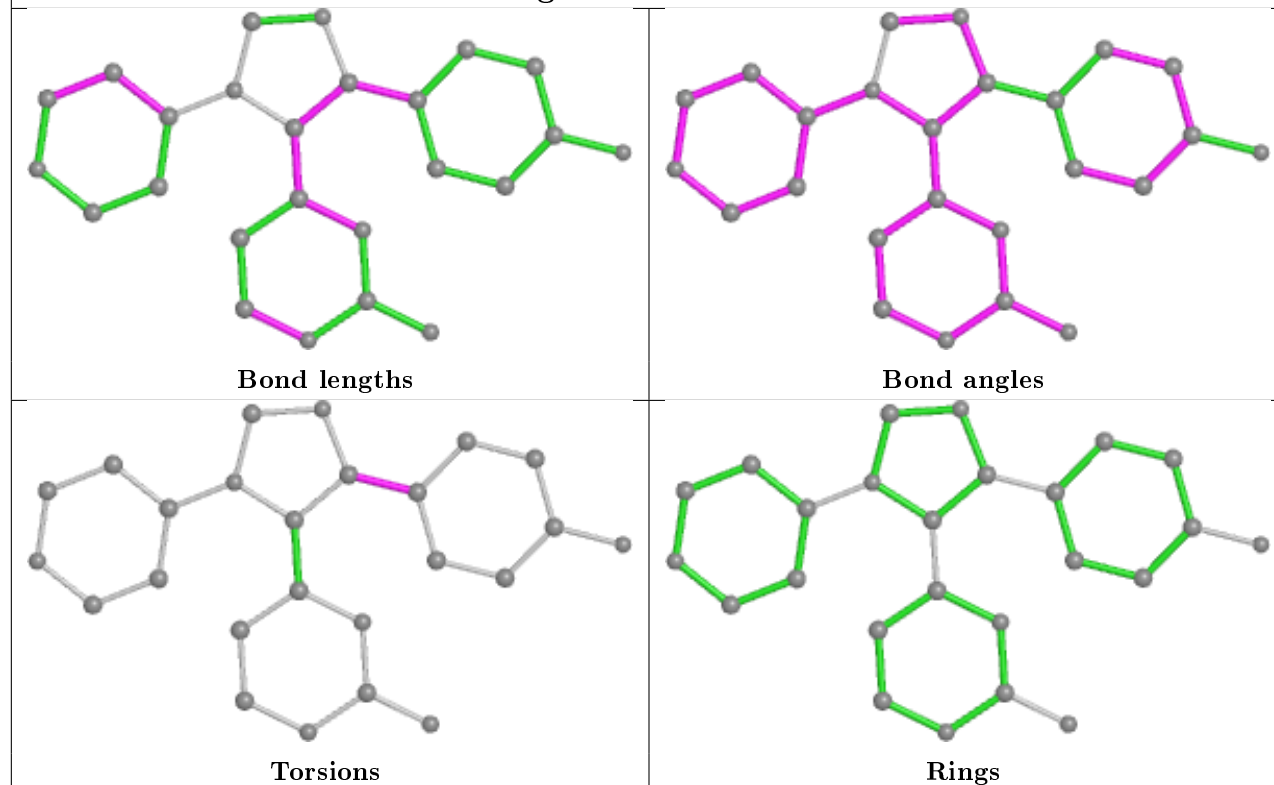
## Ligand SB4 C 401



## Ligand SB4 B 401



## Ligand SB4 D 401





## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	351/361 (97%)	-0.12	6 (1%) 70 73	16, 27, 42, 69	0
1	B	350/361 (96%)	0.04	10 (2%) 51 56	22, 33, 54, 71	0
1	C	341/361 (94%)	0.06	13 (3%) 40 44	19, 31, 62, 79	5 (1%)
1	D	351/361 (97%)	-0.12	5 (1%) 75 78	18, 27, 43, 63	0
2	K	21/29 (72%)	0.14	1 (4%) 30 33	24, 32, 55, 62	0
2	L	21/29 (72%)	0.31	2 (9%) 8 9	25, 33, 62, 73	0
2	M	18/29 (62%)	0.52	3 (16%) 1 1	30, 39, 59, 62	0
2	N	19/29 (65%)	0.61	2 (10%) 6 6	34, 44, 72, 84	0
All	All	1472/1560 (94%)	-0.02	42 (2%) 51 56	16, 30, 54, 84	5 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	198	MET	7.1
1	C	195	LEU	6.3
1	C	197	TRP	6.1
1	C	178	GLU	6.0
1	D	3	GLN	5.7
1	C	175	THR	4.7
1	D	353	LEU	4.4
1	C	174	HIS	4.3
1	B	178	GLU	4.1
1	B	35[A]	TYR	4.0
1	C	177	ASP	4.0
1	C	199	HIS	3.9
1	A	353	LEU	3.4
2	L	384	ARG	3.3
1	B	198	MET	3.3
1	D	179	MET	3.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	180[A]	THR	3.1
1	D	180[A]	THR	3.0
1	B	34[A]	ALA	3.0
1	B	255	ALA	2.9
1	C	172	ALA	2.9
1	B	179	MET	2.9
1	A	178	GLU	2.5
1	C	259	ILE	2.4
2	M	386	TYR	2.4
1	C	173	ARG	2.4
1	B	177	ASP	2.4
1	D	182	TYR	2.3
2	L	394	SER	2.3
1	A	179[A]	MET	2.3
2	N	404	THR	2.3
1	B	175	THR	2.2
2	N	384	ARG	2.2
2	M	411	SER	2.1
1	A	3	GLN	2.1
1	A	4	GLU	2.1
1	B	36[A]	GLY	2.1
2	M	404	THR	2.1
2	K	395	ALA	2.1
1	C	353	LEU	2.0
1	B	197	TRP	2.0
1	C	244	ALA	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](#)

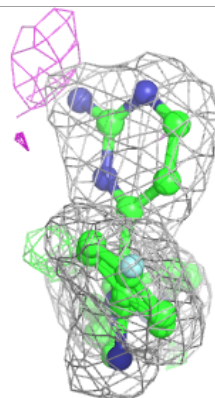
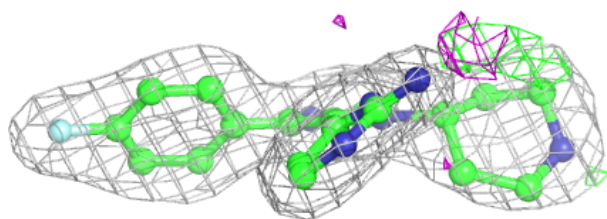
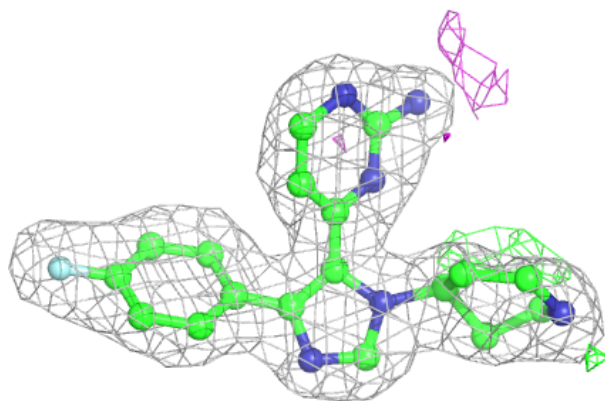
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	EDO	C	404	4/4	0.83	0.27	38,46,49,49	0
5	EDO	B	404	4/4	0.86	0.30	35,36,39,40	0
5	EDO	D	406	4/4	0.87	0.13	34,35,36,39	0
5	EDO	D	404	4/4	0.88	0.14	34,37,39,41	0
5	EDO	A	405	4/4	0.89	0.18	35,38,38,41	0
5	EDO	B	403	4/4	0.89	0.20	30,33,34,39	0
5	EDO	D	405	4/4	0.90	0.18	35,36,41,42	0
5	EDO	C	405	4/4	0.91	0.15	41,44,47,47	0
5	EDO	C	402	4/4	0.91	0.10	36,39,42,43	0
5	EDO	A	404	4/4	0.92	0.19	38,41,42,46	0
5	EDO	B	405	4/4	0.93	0.24	33,33,34,38	0
5	EDO	D	403	4/4	0.95	0.14	33,33,39,41	0
5	EDO	C	403	4/4	0.95	0.11	34,38,38,38	0
5	EDO	B	402	4/4	0.95	0.11	31,32,33,35	0
3	SB4	D	401	25/25	0.96	0.13	18,23,32,38	0
4	TLA	D	402	10/10	0.96	0.14	29,30,33,34	0
5	EDO	A	403	4/4	0.96	0.10	30,33,35,40	0
4	TLA	A	402	10/10	0.96	0.16	25,32,34,36	0
3	SB4	C	401	25/25	0.97	0.12	19,23,33,34	0
3	SB4	A	401	25/25	0.97	0.12	20,24,34,40	0
3	SB4	B	401	25/25	0.97	0.11	22,26,33,35	0

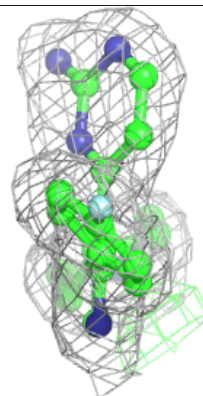
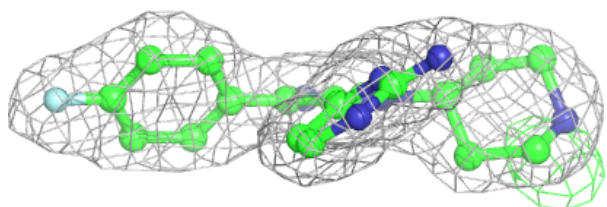
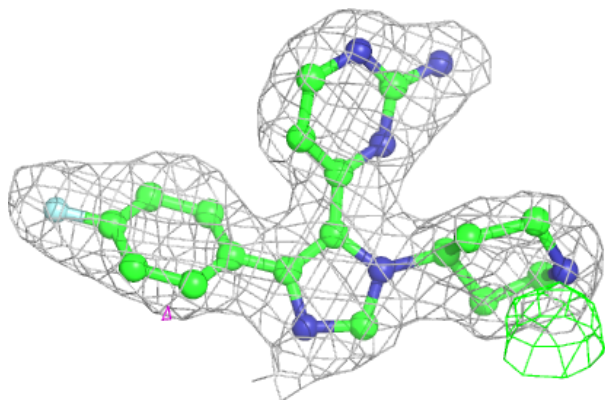
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around SB4 D 401:**

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

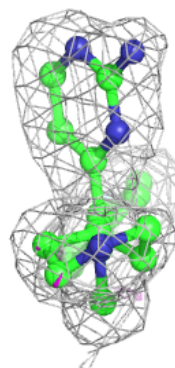
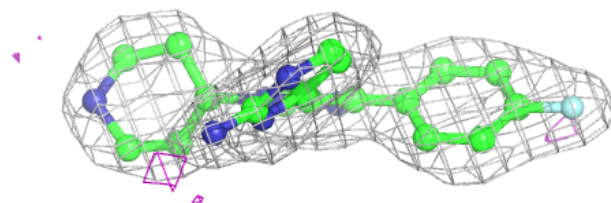
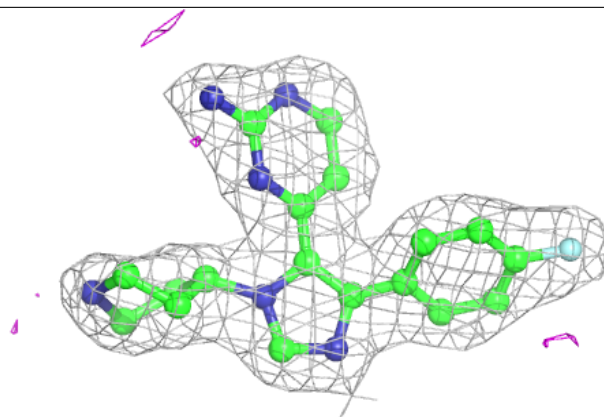
**Electron density around SB4 C 401:**

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

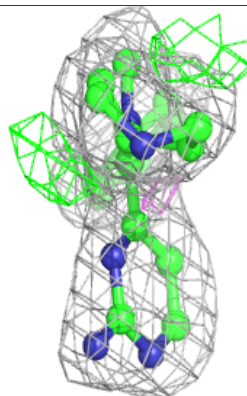
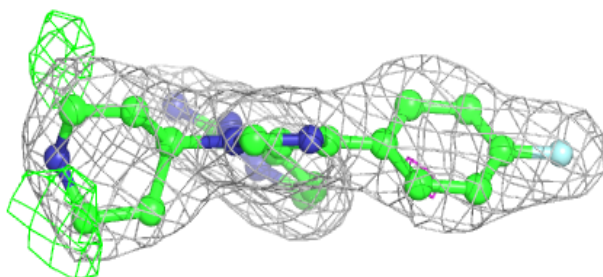
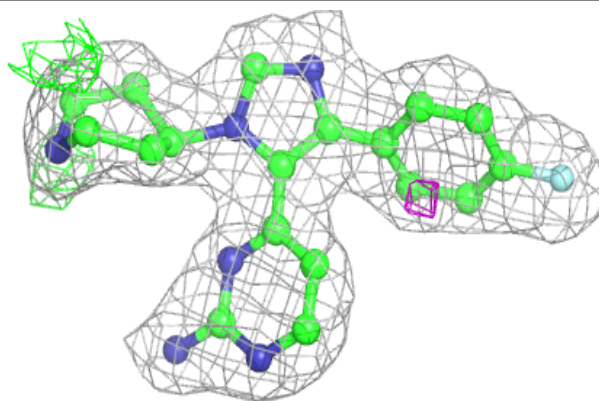


**Electron density around SB4 A 401:**

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

**Electron density around SB4 B 401:**

$2mF_o-DF_c$  (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.