



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 12:07 pm BST

PDB ID : 2PZI
Title : Crystal Structure of Protein kinase PknG from Mycobacterium tuberculosis
in Complex with Tetrahydrobenzothiophene AX20017
Authors : Honnappa, S.; Steinmetz, M.O.
Deposited on : 2007-05-18
Resolution : 2.40 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

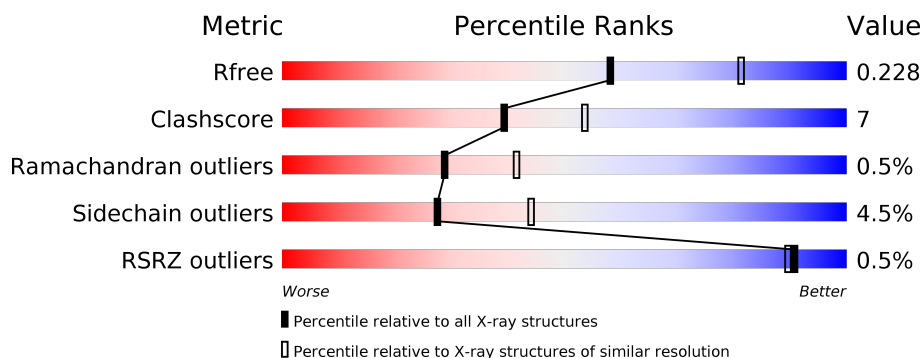
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	681	
1	B	681	

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
5	GOL	A	755	-	-	X	-
5	GOL	A	756	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10346 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 Probable serine/threonine-protein kinase pknG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	654	Total	C	N	O	S	0	0	0
			5029	3184	885	946	14			
1	B	644	Total	C	N	O	S	0	0	0
			4946	3129	872	931	14			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	69	GLY	-	CLONING ARTIFACT	UNP P65728
A	70	SER	-	CLONING ARTIFACT	UNP P65728
A	71	HIS	-	CLONING ARTIFACT	UNP P65728
A	72	MET	-	CLONING ARTIFACT	UNP P65728
B	69	GLY	-	CLONING ARTIFACT	UNP P65728
B	70	SER	-	CLONING ARTIFACT	UNP P65728
B	71	HIS	-	CLONING ARTIFACT	UNP P65728
B	72	MET	-	CLONING ARTIFACT	UNP P65728

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

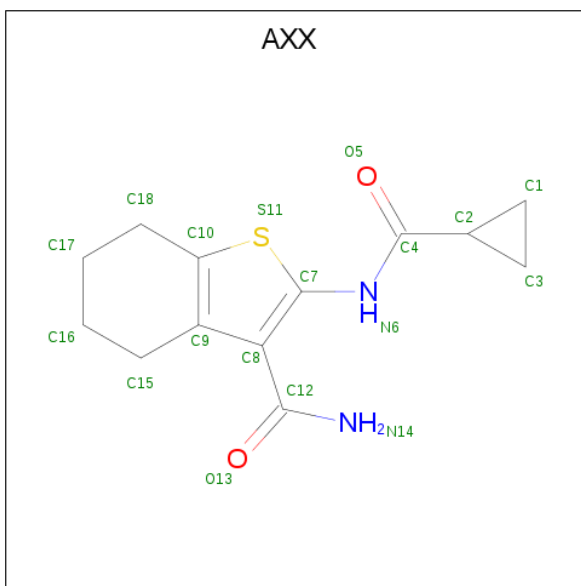
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cl	0	0
			1	1		

- Molecule 3 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cd	0	0
			1	1		
3	A	1	Total	Cd	0	0
			1	1		

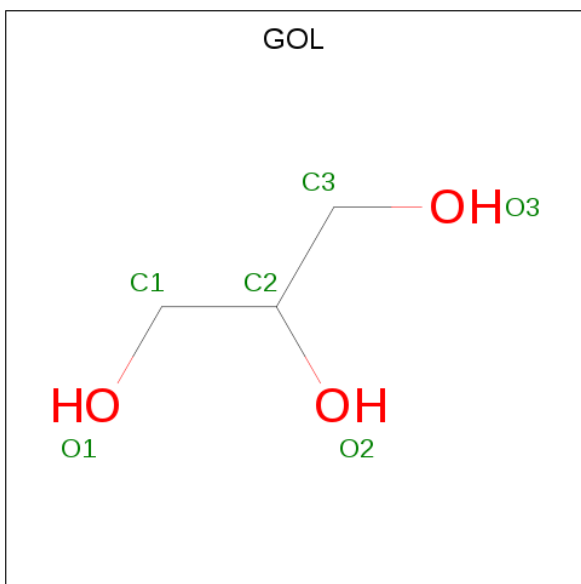
- Molecule 4 is 2-[(CYCLOPROPYLCARBONYL)AMINO]-4,5,6,7-TETRAHYDRO-1-BEN

ZOTHIOPHENE-3-CARBOXAMIDE (three-letter code: AXX) (formula: $C_{13}H_{16}N_2O_2S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			18	13	2	2	1		
4	B	1	Total	C	N	O	S	0	0
			18	13	2	2	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		


- Molecule 6 is water.

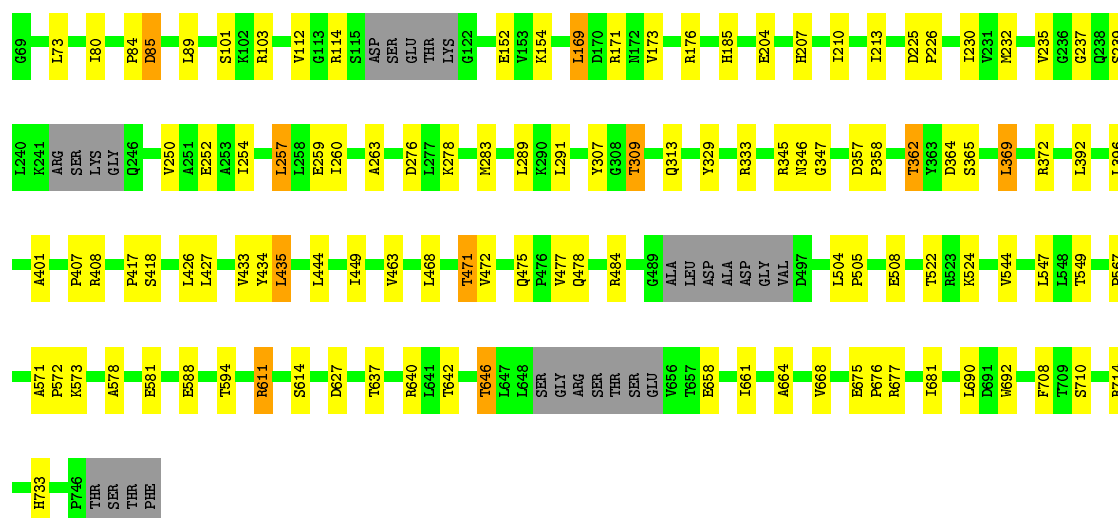
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	151	Total	O	0	0
			151	151		
6	B	145	Total	O	0	0
			145	145		

3 Residue-property plots


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.

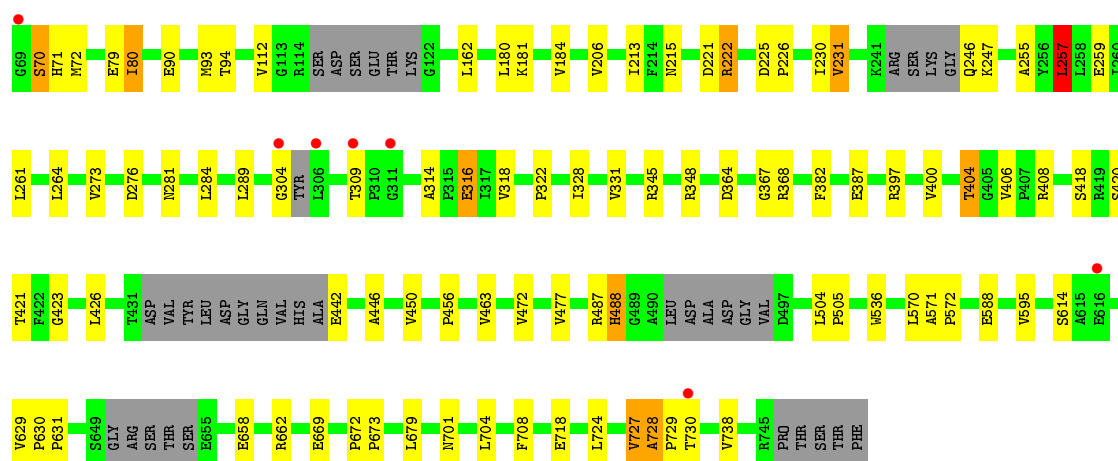
- Molecule 1: Probable serine/threonine-protein kinase pknG

Chain A: 



- Molecule 1: Probable serine/threonine-protein kinase pknG

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	122.55Å 122.55Å 243.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	106.00 – 2.40 48.92 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (106.00-2.40) 99.9 (48.92-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.84 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.183 , 0.233 0.179 , 0.228	Depositor DCC
R_{free} test set	4045 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.054 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10346	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.65% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, AXX, CL, CD

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	0.52	0/5135	0.67	4/6999 (0.1%)
1	B	0.57	2/5046 (0.0%)	0.65	3/6872 (0.0%)
All	All	0.55	2/10181 (0.0%)	0.66	7/13871 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	304	GLY	C-O	11.03	1.41	1.23
1	B	304	GLY	CA-C	5.40	1.60	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	309	THR	CA-CB-CG2	-6.97	102.64	112.40
1	B	257	LEU	CA-CB-CG	-6.95	99.32	115.30
1	A	257	LEU	CA-CB-CG	-6.23	100.96	115.30
1	B	304	GLY	CA-C-O	-6.18	109.48	120.60
1	A	611	ARG	NE-CZ-NH1	-5.91	117.34	120.30
1	A	369	LEU	CA-CB-CG	5.84	128.74	115.30
1	A	89	LEU	CA-CB-CG	5.20	127.25	115.30

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	5029	0	5023	85	0
1	B	4946	0	4947	62	0
2	A	1	0	0	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	18	0	16	2	0
4	B	18	0	16	1	0
5	A	24	0	32	28	0
5	B	12	0	16	4	0
6	A	151	0	0	6	0
6	B	145	0	0	4	0
All	All	10346	0	10050	149	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 7.

All (149) 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)
5:A:757:GOL:H2	1:B:477:VAL:HB	1.49	0.92
1:A:475:GLN:HB2	5:A:755:GOL:H31	1.54	0.89
1:B:404:THR:HG21	1:B:408:ARG:HH21	1.40	0.87
1:A:594:THR:HG21	5:A:756:GOL:H11	1.55	0.87
1:A:84:PRO:O	1:A:85:ASP:HB2	1.74	0.86
1:A:637:THR:HG21	1:A:675:GLU:HG2	1.57	0.86
1:A:573:LYS:HE2	5:A:756:GOL:H12	1.64	0.78
1:B:80:ILE:HD11	1:B:284:LEU:HD21	1.66	0.78
1:A:475:GLN:HG3	1:B:420:SER:HB2	1.66	0.78
1:A:477:VAL:HB	5:A:755:GOL:H32	1.66	0.77
1:A:477:VAL:N	5:A:755:GOL:H32	1.99	0.77
1:B:273:VAL:HG11	1:B:322:PRO:HB2	1.66	0.76
1:A:477:VAL:HB	5:A:755:GOL:H12	1.67	0.75
1:A:154:LYS:HE3	1:A:169:LEU:HD22	1.69	0.74
1:A:235:VAL:HB	1:A:283:MET:HE1	1.69	0.74
1:A:581:GLU:CD	1:A:611:ARG:HH12	1.92	0.73
1:A:103:ARG:O	1:A:114:ARG:NH1	2.21	0.73
1:A:477:VAL:H	5:A:755:GOL:H32	1.52	0.73
1:A:235:VAL:HG12	1:A:283:MET:CE	2.18	0.73
1:B:418:SER:HB2	5:B:754:GOL:H11	1.71	0.72
1:B:728:ALA:N	1:B:729:PRO:HD3	2.06	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:358:PRO:O	1:A:362:THR:HG23	1.91	0.70
1:B:669:GLU:HG3	1:B:704:LEU:HD23	1.72	0.70
1:A:417:PRO:HB3	5:A:754:GOL:H11	1.74	0.69
1:A:235:VAL:HG12	1:A:283:MET:HE3	1.73	0.69
1:A:642:THR:O	1:A:646:THR:HG23	1.92	0.69
1:A:475:GLN:CB	5:A:755:GOL:H31	2.23	0.68
1:A:478:GLN:HB2	5:A:755:GOL:H11	1.77	0.67
1:A:594:THR:HG21	5:A:756:GOL:C1	2.22	0.67
1:A:463:VAL:HG23	6:A:897:HOH:O	1.94	0.66
1:A:581:GLU:OE2	1:A:611:ARG:NH1	2.29	0.65
1:B:328:ILE:HD13	1:B:382:PHE:O	1.96	0.65
1:A:276:ASP:OD1	1:A:309:THR:HG23	1.96	0.65
1:B:504:LEU:HB3	1:B:505:PRO:HD3	1.79	0.65
1:A:235:VAL:CG1	1:A:283:MET:CE	2.75	0.64
1:A:84:PRO:O	1:A:85:ASP:CB	2.47	0.63
1:B:463:VAL:HG23	6:B:788:HOH:O	1.97	0.63
1:A:364:ASP:HB3	5:A:756:GOL:H2	1.79	0.63
1:A:477:VAL:CB	5:A:755:GOL:H12	2.30	0.62
1:A:477:VAL:CB	5:A:755:GOL:H32	2.29	0.62
1:B:364:ASP:O	1:B:368:ARG:HG3	2.00	0.62
1:B:404:THR:HG21	1:B:408:ARG:NH2	2.13	0.61
1:B:276:ASP:O	1:B:281:ASN:ND2	2.33	0.61
1:B:71:HIS:HB2	1:B:80:ILE:O	2.00	0.61
1:B:662:ARG:NH2	1:B:701:ASN:O	2.34	0.60
1:A:152:GLU:OE2	1:A:171:ARG:NH1	2.34	0.59
1:A:276:ASP:OD1	1:A:309:THR:CG2	2.51	0.59
1:A:254:ILE:HG23	1:A:392:LEU:HG	1.84	0.59
1:A:345:ARG:O	1:A:346:ASN:HB2	2.03	0.59
1:A:504:LEU:HB3	1:A:505:PRO:HD3	1.85	0.58
5:A:757:GOL:C2	1:B:477:VAL:HB	2.28	0.58
1:B:727:VAL:C	1:B:729:PRO:HD3	2.23	0.58
1:A:477:VAL:HB	5:A:755:GOL:C3	2.33	0.58
1:A:435:LEU:HD12	6:A:884:HOH:O	2.04	0.57
1:B:221:ASP:O	1:B:222:ARG:HB2	2.05	0.57
1:B:215:ASN:HB3	1:B:231:VAL:HG13	1.85	0.57
1:A:237:GLY:H	1:A:283:MET:HE3	1.71	0.56
1:A:477:VAL:CA	5:A:755:GOL:H32	2.36	0.55
1:B:418:SER:H	5:B:754:GOL:H11	1.71	0.55
1:A:484:ARG:NH2	6:A:802:HOH:O	2.40	0.55
1:B:367:GLY:HA3	6:B:844:HOH:O	2.07	0.55
1:A:477:VAL:HB	5:A:755:GOL:C1	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:404:THR:CG2	1:B:408:ARG:HH21	2.15	0.54
5:A:755:GOL:HO2	1:B:420:SER:HG	1.55	0.54
1:B:72:MET:HG3	1:B:79:GLU:OE2	2.08	0.54
1:A:449:ILE:HD13	1:A:578:ALA:HB1	1.89	0.53
1:A:80:ILE:HD11	1:A:252:GLU:HG2	1.91	0.53
1:A:664:ALA:O	1:A:668:VAL:HG23	2.09	0.52
1:A:278:LYS:NZ	1:A:309:THR:HG21	2.24	0.52
1:A:401:ALA:HB2	1:A:408:ARG:HG2	1.90	0.52
1:B:314:ALA:HB1	1:B:316:GLU:OE2	2.09	0.52
1:A:73:LEU:HD22	1:A:259:GLU:HG3	1.90	0.52
1:B:80:ILE:CD1	1:B:284:LEU:HD21	2.37	0.51
1:A:207:HIS:HB3	1:A:210:ILE:HD12	1.93	0.51
1:A:571:ALA:HB3	1:A:572:PRO:HD3	1.92	0.51
1:B:487:ARG:O	1:B:488:HIS:HB2	2.12	0.49
1:B:658:GLU:OE2	1:B:662:ARG:NH1	2.45	0.49
4:B:752:AXX:S11	4:B:752:AXX:O5	2.70	0.49
1:A:417:PRO:CB	5:A:754:GOL:H11	2.42	0.49
1:B:701:ASN:H	1:B:701:ASN:HD22	1.60	0.49
1:A:468:LEU:O	1:A:471:THR:HG22	2.12	0.49
1:B:423:GLY:HA2	1:B:426:LEU:HD13	1.95	0.48
1:A:213:ILE:HD12	1:A:230:ILE:HD11	1.95	0.48
1:A:658:GLU:HB2	1:A:692:TRP:CH2	2.49	0.48
1:A:418:SER:N	5:A:754:GOL:H12	2.29	0.48
1:A:185:HIS:HD2	6:A:907:HOH:O	1.98	0.47
1:B:571:ALA:HB3	1:B:572:PRO:HD3	1.95	0.47
1:A:235:VAL:CB	1:A:283:MET:HE1	2.42	0.47
1:A:346:ASN:HB3	1:A:347:GLY:H	1.55	0.47
1:A:477:VAL:H	5:A:755:GOL:C3	2.23	0.47
1:B:418:SER:HB2	5:B:754:GOL:C1	2.42	0.47
1:B:588:GLU:HG2	6:B:873:HOH:O	2.15	0.47
1:B:255:ALA:O	1:B:259:GLU:HG2	2.15	0.47
1:B:387:GLU:HB2	6:B:776:HOH:O	2.13	0.47
1:B:727:VAL:HG13	1:B:727:VAL:O	2.15	0.47
1:B:213:ILE:HD12	1:B:230:ILE:HD11	1.97	0.46
1:A:278:LYS:CE	1:A:309:THR:HG21	2.46	0.46
1:B:679:LEU:HB2	1:B:724:LEU:HD21	1.96	0.46
1:A:254:ILE:HG21	1:A:396:LEU:HB2	1.96	0.46
4:A:753:AXX:H141	4:A:753:AXX:C15	2.28	0.46
1:A:594:THR:CG2	5:A:756:GOL:H11	2.39	0.46
1:B:257:LEU:HD22	1:B:331:VAL:HG13	1.98	0.46
1:B:261:LEU:HD23	1:B:264:LEU:HD12	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:446:ALA:O	1:B:450:VAL:HG23	2.17	0.45
1:A:313:GLN:HB3	6:A:809:HOH:O	2.16	0.45
1:B:630:PRO:HA	1:B:631:PRO:HD3	1.88	0.45
1:B:487:ARG:O	1:B:488:HIS:CB	2.64	0.45
1:B:426:LEU:O	1:B:442:GLU:HB3	2.17	0.45
1:B:397:ARG:NH1	5:B:753:GOL:O1	2.37	0.45
1:A:260:ILE:O	1:A:263:ALA:HB3	2.16	0.45
1:A:508:GLU:OE1	1:A:524:LYS:HE2	2.17	0.45
1:B:669:GLU:HG3	1:B:704:LEU:CD2	2.45	0.45
1:B:400:VAL:O	1:B:404:THR:HB	2.17	0.44
1:A:407:PRO:C	1:A:408:ARG:HD2	2.38	0.44
1:B:225:ASP:HA	1:B:226:PRO:HD3	1.88	0.44
1:B:728:ALA:N	1:B:729:PRO:CD	2.78	0.43
1:B:246:GLN:CG	1:B:247:LYS:H	2.31	0.43
1:A:213:ILE:HD13	1:A:232:MET:HG2	1.99	0.43
1:A:676:PRO:O	1:A:677:ARG:HB2	2.19	0.43
1:B:162:LEU:HB2	1:B:181:LYS:HE2	1.99	0.43
1:A:477:VAL:HB	5:A:755:GOL:C2	2.48	0.43
1:B:456:PRO:HB3	1:B:536:TRP:CH2	2.53	0.43
1:A:426:LEU:HD11	1:A:444:LEU:HB2	2.01	0.43
1:A:594:THR:HG23	5:A:756:GOL:H32	2.01	0.43
1:A:478:GLN:N	5:A:755:GOL:C1	2.82	0.43
1:B:404:THR:HG23	1:B:406:VAL:H	1.84	0.43
1:A:235:VAL:O	4:A:753:AXX:H2	2.19	0.42
1:A:407:PRO:O	1:A:408:ARG:HD2	2.19	0.42
1:B:718:GLU:OE1	1:B:727:VAL:HG23	2.19	0.42
1:B:261:LEU:HD23	1:B:261:LEU:HA	1.84	0.42
1:A:173:VAL:HB	1:A:176:ARG:HG2	2.00	0.42
1:B:629:VAL:HA	1:B:630:PRO:HD3	1.87	0.42
1:A:365:SER:HB2	1:A:567:PRO:O	2.18	0.42
1:A:627:ASP:HA	1:A:640:ARG:HH12	1.85	0.42
1:A:434:TYR:O	1:A:733:HIS:HE1	2.02	0.42
1:B:718:GLU:HA	1:B:718:GLU:OE1	2.18	0.42
1:A:329:TYR:O	1:A:333:ARG:HG2	2.20	0.41
1:B:180:LEU:HA	1:B:230:ILE:O	2.20	0.41
1:A:690:LEU:CD1	1:A:714:ARG:HG3	2.50	0.41
1:A:658:GLU:HB2	1:A:692:TRP:CZ2	2.55	0.41
1:B:672:PRO:HA	1:B:673:PRO:HD2	1.95	0.41
1:A:478:GLN:N	5:A:755:GOL:H12	2.35	0.41
1:A:544:VAL:O	1:A:547:LEU:HB2	2.21	0.41
1:B:420:SER:OG	1:B:421:THR:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:751:CL:CL	6:A:880:HOH:O	2.60	0.40
1:A:677:ARG:O	1:A:681:ILE:HG12	2.21	0.40
1:A:225:ASP:HA	1:A:226:PRO:HD3	1.96	0.40
1:A:357:ASP:HA	1:A:358:PRO:HD3	1.86	0.40
1:B:570:LEU:HD22	1:B:595:VAL:HA	2.04	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	644/681 (95%)	624 (97%)	19 (3%)	1 (0%)	47	62
1	B	630/681 (92%)	604 (96%)	21 (3%)	5 (1%)	19	29
All	All	1274/1362 (94%)	1228 (96%)	40 (3%)	6 (0%)	29	41

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	85	ASP
1	B	488	HIS
1	B	70	SER
1	B	222	ARG
1	B	727	VAL
1	B	728	ALA

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	537/559 (96%)	510 (95%)	27 (5%)	24	40
1	B	528/559 (94%)	507 (96%)	21 (4%)	31	49
All	All	1065/1118 (95%)	1017 (96%)	48 (4%)	27	44

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	SER
1	A	112	VAL
1	A	169	LEU
1	A	204	GLU
1	A	239	SER
1	A	250	VAL
1	A	257	LEU
1	A	289	LEU
1	A	291	LEU
1	A	307	TYR
1	A	309	THR
1	A	362	THR
1	A	369	LEU
1	A	372	ARG
1	A	427	LEU
1	A	433	VAL
1	A	435	LEU
1	A	471	THR
1	A	472	VAL
1	A	522	THR
1	A	549	THR
1	A	588	GLU
1	A	614	SER
1	A	646	THR
1	A	661	ILE
1	A	708	PHE
1	A	710	SER
1	B	70	SER
1	B	80	ILE
1	B	90	GLU
1	B	93	MET
1	B	94	THR
1	B	112	VAL

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Mol	Chain	Res	Type
1	B	184	VAL
1	B	206	VAL
1	B	231	VAL
1	B	257	LEU
1	B	289	LEU
1	B	316	GLU
1	B	318	VAL
1	B	345	ARG
1	B	348	ARG
1	B	404	THR
1	B	472	VAL
1	B	614	SER
1	B	708	PHE
1	B	730	THR
1	B	738	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	733	HIS
1	B	107	ASN
1	B	172	ASN
1	B	701	ASN
1	B	711	HIS

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

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 for Mogul analysis.

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	GOL	B	754	-	5,5,5	0.41	0	5,5,5	0.47	0
5	GOL	A	756	-	5,5,5	0.47	0	5,5,5	0.54	0
4	AXX	A	753	-	18,20,20	1.02	1 (5%)	17,29,29	3.08	4 (23%)
5	GOL	A	755	-	5,5,5	1.04	1 (20%)	5,5,5	1.59	1 (20%)
4	AXX	B	752	-	18,20,20	1.07	1 (5%)	17,29,29	2.52	4 (23%)
5	GOL	A	754	-	5,5,5	0.39	0	5,5,5	0.83	0
5	GOL	A	757	-	5,5,5	0.90	0	5,5,5	0.95	0
5	GOL	B	753	-	5,5,5	0.36	0	5,5,5	0.50	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	GOL	B	754	-	-	2/4/4/4	-
5	GOL	A	756	-	-	2/4/4/4	-
4	AXX	A	753	-	-	0/6/21/21	0/3/3/3
5	GOL	A	755	-	-	3/4/4/4	-
4	AXX	B	752	-	-	0/6/21/21	0/3/3/3
5	GOL	A	754	-	-	2/4/4/4	-
5	GOL	A	757	-	-	2/4/4/4	-
5	GOL	B	753	-	-	2/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	753	AXX	C8-C7	3.03	1.47	1.41
4	B	752	AXX	C8-C7	2.74	1.47	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	755	GOL	O2-C2	-2.03	1.37	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	753	AXX	C3-C2-C4	-10.25	107.61	117.21
4	B	752	AXX	C1-C2-C4	-7.22	110.45	117.21
4	B	752	AXX	C3-C2-C4	-5.00	112.53	117.21
4	A	753	AXX	C1-C2-C4	-4.79	112.72	117.21
4	A	753	AXX	C17-C18-C10	-4.51	106.12	113.53
4	B	752	AXX	C17-C18-C10	-4.40	106.31	113.53
4	A	753	AXX	O13-C12-C8	-2.82	116.19	120.11
4	B	752	AXX	O13-C12-C8	-2.42	116.73	120.11
5	A	755	GOL	O3-C3-C2	2.08	120.17	110.20

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	754	GOL	C1-C2-C3-O3
5	B	754	GOL	O2-C2-C3-O3
5	A	756	GOL	O1-C1-C2-O2
5	A	756	GOL	O1-C1-C2-C3
5	A	754	GOL	O1-C1-C2-C3
5	B	753	GOL	O1-C1-C2-C3
5	A	757	GOL	C1-C2-C3-O3
5	A	757	GOL	O2-C2-C3-O3
5	A	755	GOL	C1-C2-C3-O3
5	A	755	GOL	O2-C2-C3-O3
5	A	754	GOL	O1-C1-C2-O2
5	B	753	GOL	O1-C1-C2-O2
5	A	755	GOL	O1-C1-C2-O2

There are no ring outliers.

8 monomers are involved in 35 short contacts:

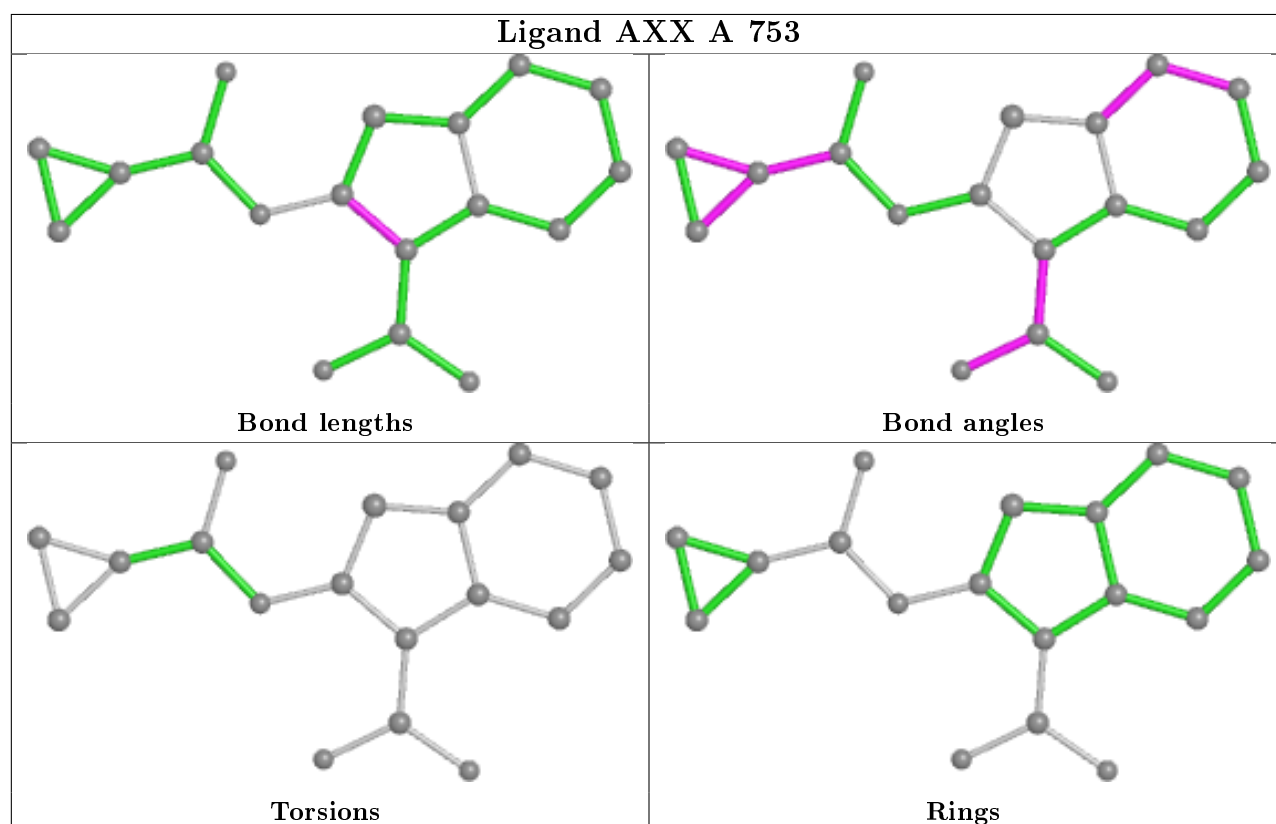
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	754	GOL	3	0
5	A	756	GOL	6	0
4	A	753	AXX	2	0
5	A	755	GOL	17	0

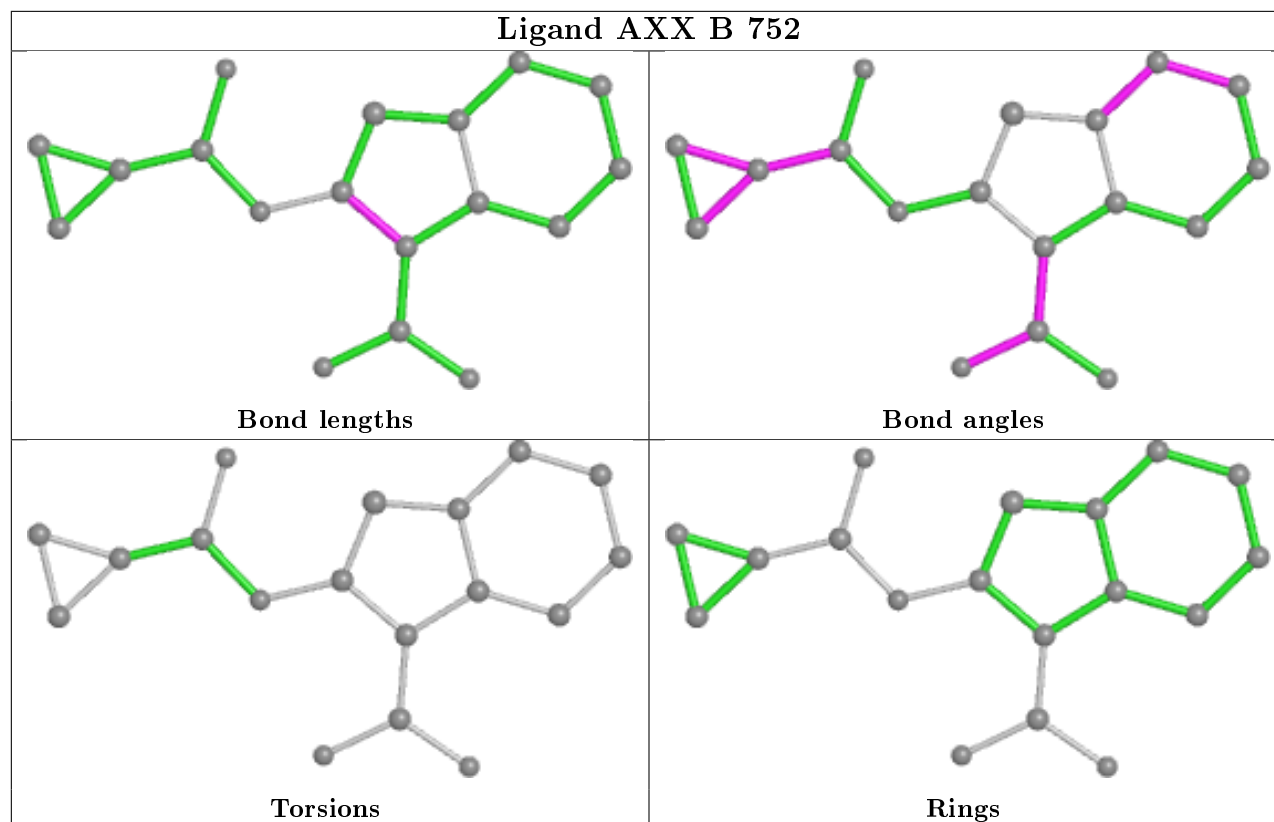
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	752	AXX	1	0
5	A	754	GOL	3	0
5	A	757	GOL	2	0
5	B	753	GOL	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	654/681 (96%)	-0.36	0 100 100	26, 44, 66, 84	0
1	B	644/681 (94%)	-0.27	7 (1%) 80 79	27, 46, 67, 91	0
All	All	1298/1362 (95%)	-0.32	7 (0%) 91 89	26, 45, 66, 91	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	69	GLY	4.9
1	B	311	GLY	2.6
1	B	306	LEU	2.4
1	B	309	THR	2.3
1	B	616	GLU	2.2
1	B	304	GLY	2.2
1	B	730	THR	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](#)

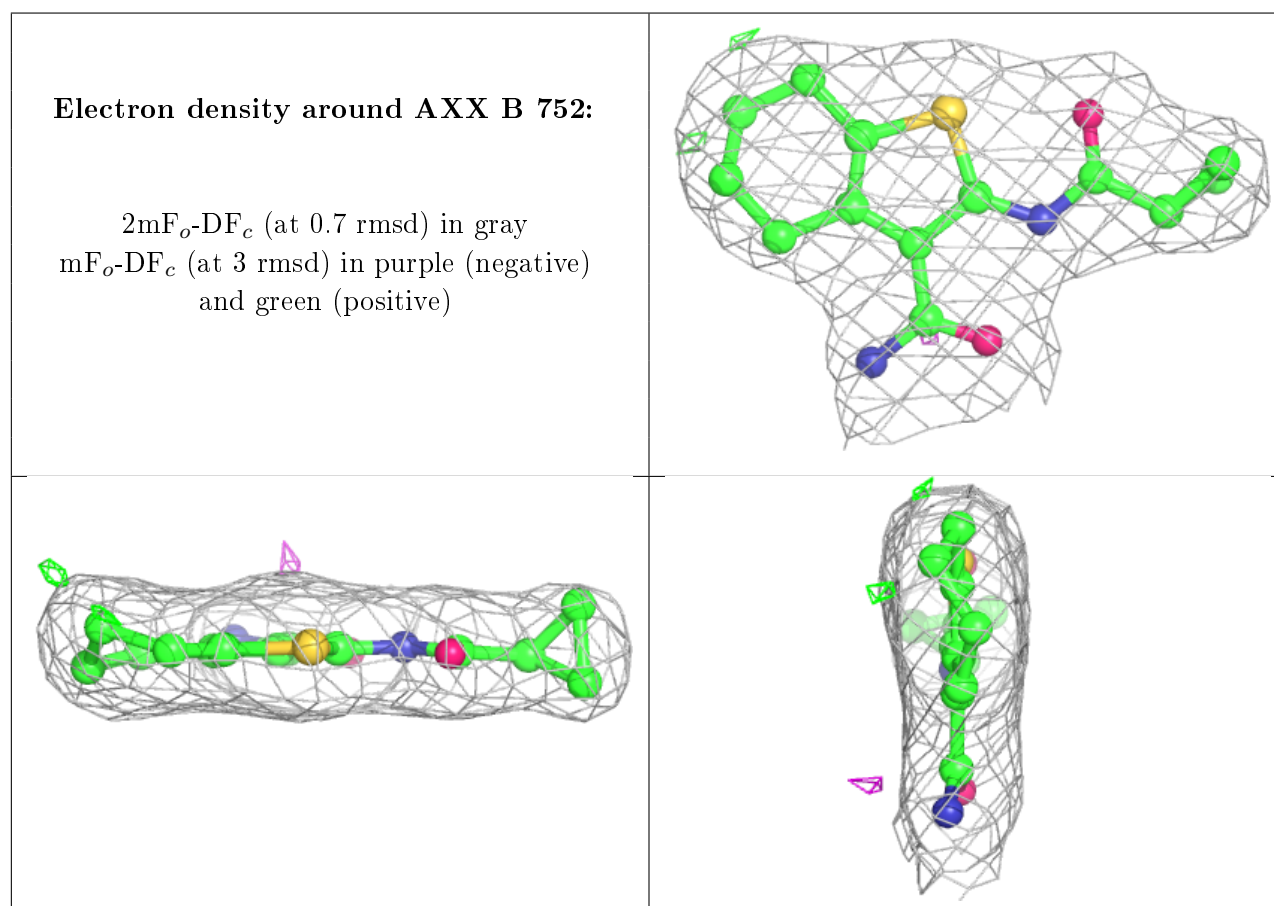
There are no carbohydrates in this entry.

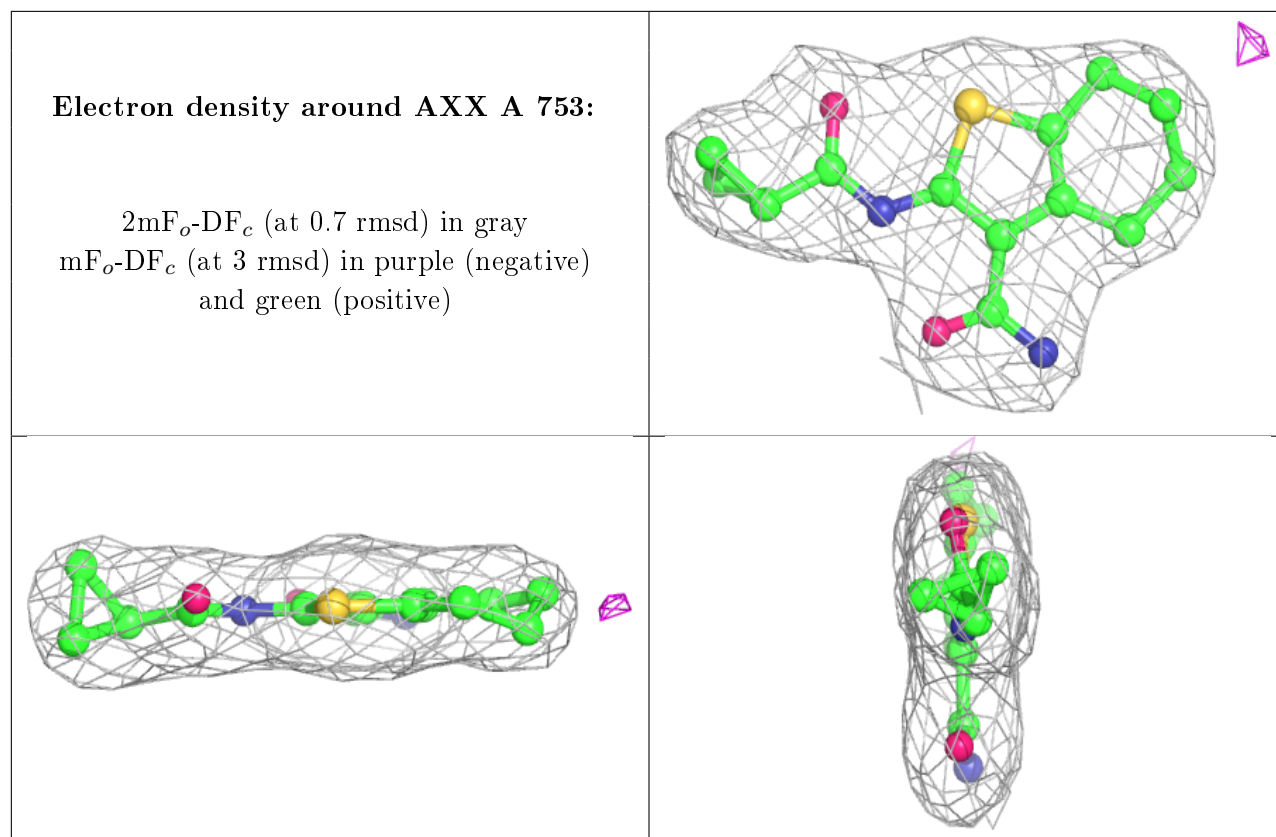
6.4 Ligands [i](#)

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
5	GOL	B	754	6/6	0.77	0.22	62,70,77,86	0
5	GOL	A	754	6/6	0.83	0.23	57,62,76,81	0
5	GOL	A	756	6/6	0.86	0.25	80,83,85,90	0
5	GOL	A	757	6/6	0.89	0.32	45,56,69,74	0
5	GOL	A	755	6/6	0.92	0.22	51,52,70,73	0
5	GOL	B	753	6/6	0.94	0.13	56,70,75,78	0
4	AXX	B	752	18/18	0.98	0.10	33,45,49,57	0
2	CL	A	751	1/1	0.98	0.13	66,66,66,66	0
3	CD	B	751	1/1	0.99	0.15	70,70,70,70	0
4	AXX	A	753	18/18	0.99	0.11	33,46,53,57	0
3	CD	A	752	1/1	0.99	0.14	53,53,53,53	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.