



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2025 – 04:54 PM EST

PDB ID : 9MGL  
Title : Crystal structure of PRMT5:MEP50 in complex with MTA and GSK3326595  
Authors : Whittington, D.A.  
Deposited on : 2024-12-11  
Resolution : 2.25 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

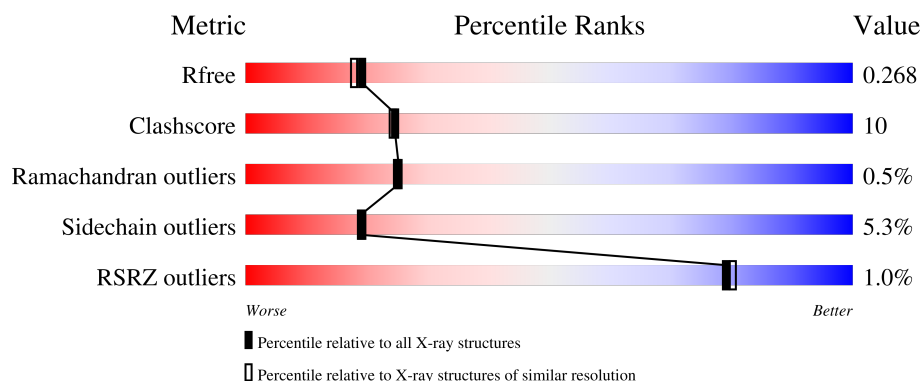
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.25 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	164625	1763 (2.26-2.26)
Clashscore	180529	1919 (2.26-2.26)
Ramachandran outliers	177936	1884 (2.26-2.26)
Sidechain outliers	177891	1885 (2.26-2.26)
RSRZ outliers	164620	1763 (2.26-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	645	<div> <div style="display: flex; align-items: center;"> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="margin-left: 10px;"> <div style="width: 76%; height: 10px; background-color: red;"></div> <div style="width: 19%; height: 10px; background-color: orange;"></div> <div style="width: 5%; height: 10px; background-color: yellow;"></div> <div style="width: 5%; height: 10px; background-color: green;"></div> <div style="width: 5%; height: 10px; background-color: grey;"></div> </div> </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>76%</span> <span>19%</span> <span>• •</span> </div> </div>
2	B	350	<div> <div style="display: flex; align-items: center;"> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="margin-left: 10px;"> <div style="width: 60%; height: 10px; background-color: red;"></div> <div style="width: 24%; height: 10px; background-color: orange;"></div> <div style="width: 1%; height: 10px; background-color: yellow;"></div> <div style="width: 1%; height: 10px; background-color: green;"></div> <div style="width: 10%; height: 10px; background-color: grey;"></div> </div> </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>60%</span> <span>24%</span> <span>•</span> <span>13%</span> </div> </div>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 14927 atoms, of which 7282 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 Protein arginine N-methyltransferase 5.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	625	10013	3237	4952	869	931	24	0	0	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP O14744
A	-6	ASP	-	expression tag	UNP O14744
A	-5	TYR	-	expression tag	UNP O14744
A	-4	LYS	-	expression tag	UNP O14744
A	-3	ASP	-	expression tag	UNP O14744
A	-2	ASP	-	expression tag	UNP O14744
A	-1	ASP	-	expression tag	UNP O14744
A	0	ASP	-	expression tag	UNP O14744
A	1	LYS	-	expression tag	UNP O14744

- Molecule 2 is a protein called Methylosome protein 50.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	304	4510	1443	2213	393	448	13	0	0	0

There are 9 discrepancies between the modelled and reference sequences:

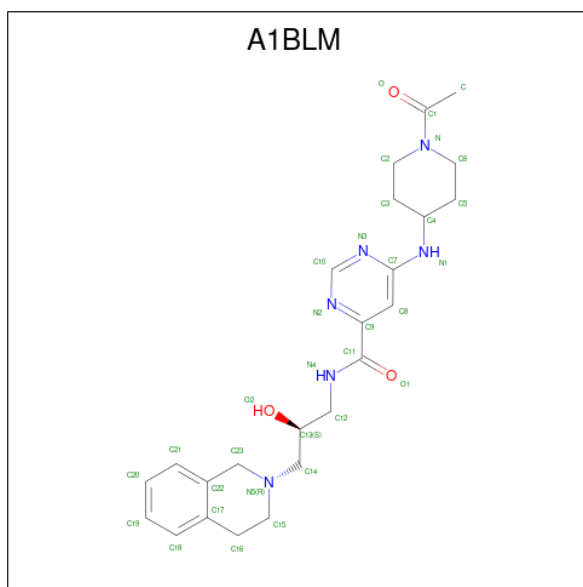
Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	MET	-	initiating methionine	UNP Q9BQA1
B	-6	HIS	-	expression tag	UNP Q9BQA1
B	-5	HIS	-	expression tag	UNP Q9BQA1
B	-4	HIS	-	expression tag	UNP Q9BQA1
B	-3	HIS	-	expression tag	UNP Q9BQA1
B	-2	HIS	-	expression tag	UNP Q9BQA1
B	-1	HIS	-	expression tag	UNP Q9BQA1

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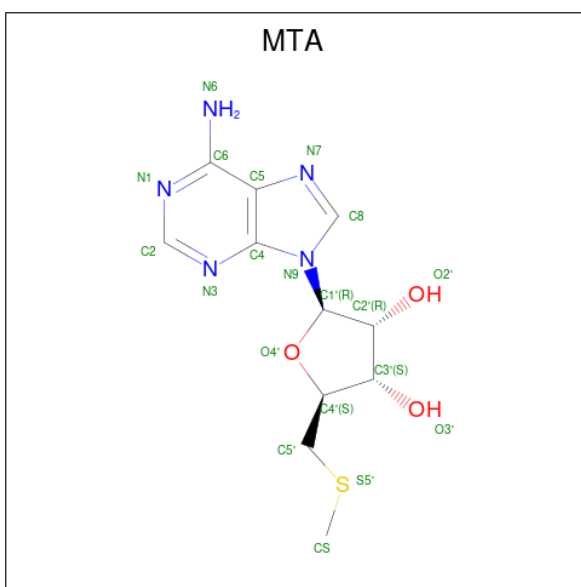
Chain	Residue	Modelled	Actual	Comment	Reference
B	0	HIS	-	expression tag	UNP Q9BQA1
B	1	HIS	-	expression tag	UNP Q9BQA1

- Molecule 3 is pemrametostat (three-letter code: A1BLM) (formula:  $C_{24}H_{32}N_6O_3$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			65	24	32	6	3		

- Molecule 4 is 5'-DEOXY-5'-METHYLTHIOADENOSINE (three-letter code: MTA) (formula:  $C_{11}H_{15}N_5O_3S$ ).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	S	0	0
			35	11	15	5	3	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	B	1	Total	C	H	O	0	0
			10	2	6	2		
6	B	1	Total	C	H	O	0	0
			10	2	6	2		
6	B	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	H	O	0	0
			17	4	10	3		

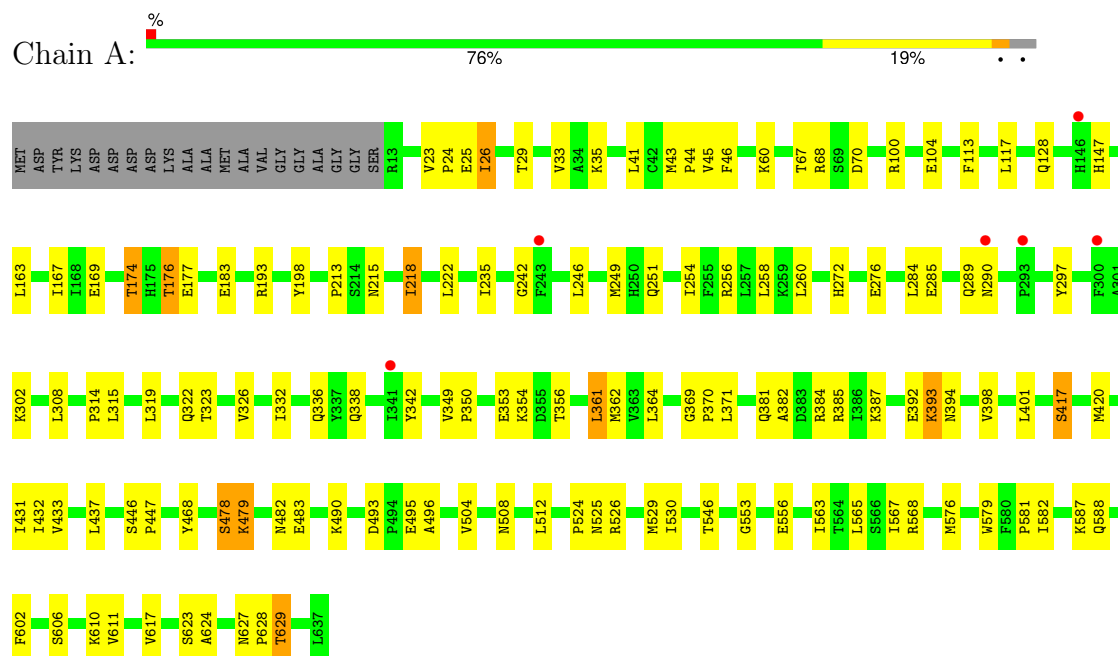
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	124	Total	O	0	0
			124	124		
8	B	48	Total	O	0	0
			48	48		

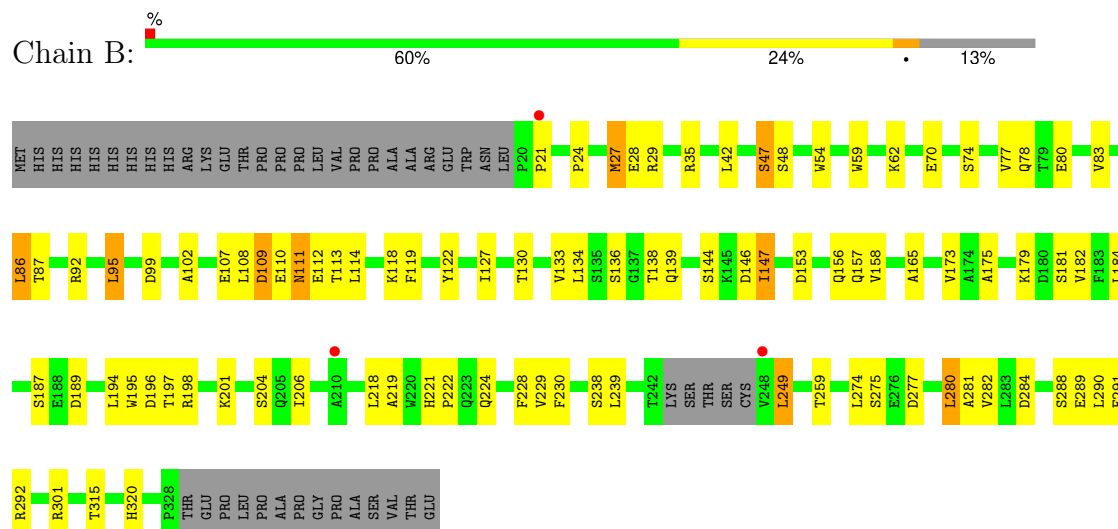
### 3 Residue-property plots [i](#)

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: Protein arginine N-methyltransferase 5



- Molecule 2: Methylosome protein 50





## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	99.74Å 136.97Å 178.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.62 – 2.25 34.62 – 2.25	Depositor EDS
% Data completeness (in resolution range)	42.6 (34.62-2.25) 42.6 (34.62-2.25)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.73 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.192 , 0.268 0.192 , 0.268	Depositor DCC
$R_{free}$ test set	34661 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.2	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 33.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	14927	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, A1BLM, SO4, EDO, MTA

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	0.69	0/5203	0.75	0/7078
2	B	0.55	0/2353	0.73	0/3215
All	All	0.65	0/7556	0.74	0/10293

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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	5061	4952	4952	82	0
2	B	2297	2213	2212	63	0
3	A	33	32	0	0	0
4	A	20	15	15	1	0
5	A	15	0	0	0	0
6	A	28	42	42	3	0
6	B	12	18	18	3	0
7	A	7	10	10	0	0
8	A	124	0	0	3	0
8	B	48	0	0	1	0
All	All	7645	7282	7249	142	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 10.

All (142) 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)
2:B:130:THR:HG21	2:B:173:VAL:H	1.41	0.84
1:A:349:VAL:HG23	1:A:384:ARG:HE	1.42	0.84
1:A:45:VAL:HG23	1:A:46:PHE:HD1	1.45	0.81
2:B:109:ASP:OD2	2:B:111:ASN:ND2	2.15	0.80
1:A:26:ILE:HD12	1:A:43:MET:HE1	1.63	0.78
1:A:254:ILE:O	1:A:258:LEU:HD12	1.86	0.76
1:A:254:ILE:HG22	1:A:258:LEU:HD11	1.69	0.74
1:A:163:LEU:HD23	1:A:163:LEU:O	1.90	0.72
2:B:284:ASP:HB3	2:B:290:LEU:HD11	1.74	0.70
2:B:194:LEU:HB2	2:B:206:ILE:HD11	1.74	0.69
2:B:102:ALA:HB2	2:B:122:TYR:CD1	2.28	0.69
2:B:99:ASP:HA	2:B:127:ILE:HG23	1.75	0.69
2:B:28:GLU:HG2	2:B:48:SER:HB2	1.74	0.68
1:A:349:VAL:HG23	1:A:384:ARG:NE	2.10	0.67
1:A:100:ARG:O	1:A:104:GLU:HG3	1.95	0.66
1:A:314:PRO:HB3	1:A:319:LEU:HD11	1.77	0.66
1:A:174:THR:HG22	1:A:176:THR:HG23	1.80	0.64
1:A:512:LEU:HD13	1:A:546:THR:HG21	1.81	0.62
2:B:221:HIS:CD2	2:B:222:PRO:HD2	2.34	0.62
1:A:24:PRO:HA	2:B:54:TRP:HB2	1.82	0.62
1:A:350:PRO:HD2	1:A:353:GLU:HG3	1.82	0.61
2:B:182:VAL:HA	2:B:195:TRP:O	1.99	0.61
1:A:437:LEU:HD21	1:A:468:TYR:CD2	2.36	0.61
2:B:282:VAL:HB	2:B:291:PHE:HB3	1.83	0.61
1:A:169:GLU:HA	2:B:201:LYS:HD3	1.83	0.61
2:B:47:SER:HB3	2:B:83:VAL:H	1.67	0.59
2:B:62:LYS:HG3	2:B:74:SER:OG	2.04	0.58
2:B:70:GLU:OE1	2:B:70:GLU:N	2.34	0.58
1:A:45:VAL:HG23	1:A:46:PHE:CD1	2.35	0.58
2:B:35:ARG:NH2	2:B:86:LEU:O	2.37	0.58
2:B:156:GLN:HB3	2:B:158:VAL:HG22	1.86	0.57
2:B:181:SER:HB2	2:B:197:THR:HG22	1.85	0.57
1:A:25:GLU:O	1:A:29:THR:HG23	2.04	0.57
1:A:174:THR:HG22	1:A:176:THR:CG2	2.34	0.57
2:B:184:LEU:HD11	2:B:228:PHE:CD2	2.40	0.57
1:A:623:SER:HB2	6:A:712:EDO:H22	1.87	0.57
1:A:493:ASP:HB3	1:A:496:ALA:HB2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:364:LEU:HB3	1:A:420:MET:CE	2.35	0.56
2:B:221:HIS:ND1	2:B:224:GLN:HG3	2.21	0.56
1:A:198:TYR:O	1:A:482:ASN:ND2	2.37	0.55
2:B:144:SER:HB3	2:B:146:ASP:OD1	2.07	0.55
2:B:259:THR:HG21	2:B:301:ARG:NH2	2.22	0.54
1:A:26:ILE:CD1	1:A:43:MET:HE1	2.37	0.54
1:A:628:PRO:O	1:A:629:THR:HB	2.08	0.54
1:A:183:GLU:OE1	8:A:801:HOH:O	2.19	0.54
1:A:322:GLN:O	1:A:326:VAL:HG13	2.07	0.54
1:A:342:TYR:HD1	1:A:382:ALA:HB2	1.73	0.54
2:B:111:ASN:ND2	2:B:111:ASN:H	2.06	0.53
2:B:134:LEU:HD12	2:B:139:GLN:HB3	1.91	0.53
2:B:274:LEU:HD12	2:B:274:LEU:N	2.24	0.53
2:B:87:THR:HG21	2:B:133:VAL:HG23	1.90	0.53
1:A:246:LEU:O	1:A:251:GLN:NE2	2.41	0.53
1:A:371:LEU:HD13	1:A:433:VAL:HG12	1.91	0.53
1:A:362:MET:HB2	1:A:432:ILE:HG22	1.91	0.52
1:A:606:SER:HB3	1:A:611:VAL:HG22	1.90	0.52
2:B:153:ASP:OD2	2:B:156:GLN:HB2	2.10	0.52
2:B:184:LEU:HB3	2:B:218:LEU:HD13	1.93	0.51
1:A:553:GLY:HA3	1:A:582:ILE:HG22	1.93	0.51
2:B:113:THR:O	2:B:114:LEU:HD12	2.10	0.51
2:B:181:SER:CB	2:B:197:THR:HG22	2.41	0.51
1:A:213:PRO:HD2	1:A:218:ILE:HD11	1.93	0.50
1:A:308:LEU:HD23	1:A:504:VAL:HB	1.92	0.50
2:B:24:PRO:HD2	2:B:59:TRP:CH2	2.47	0.50
1:A:29:THR:O	1:A:33:VAL:HG13	2.12	0.50
1:A:361:LEU:HD11	1:A:431:ILE:HD12	1.93	0.49
1:A:606:SER:CB	1:A:611:VAL:HG22	2.43	0.49
1:A:113:PHE:CE2	1:A:117:LEU:HD21	2.48	0.49
2:B:134:LEU:HD23	2:B:175:ALA:HB1	1.94	0.49
1:A:44:PRO:HB2	1:A:46:PHE:O	2.13	0.49
1:A:393:LYS:HG3	1:A:394:ASN:N	2.26	0.49
2:B:229:VAL:CG1	2:B:239:LEU:HD12	2.42	0.49
1:A:364:LEU:HB3	1:A:420:MET:HE1	1.93	0.49
2:B:107:GLU:HB2	2:B:118:LYS:HG3	1.95	0.49
2:B:146:ASP:HA	6:B:401:EDO:C2	2.43	0.49
2:B:289:GLU:OE2	2:B:292:ARG:HB2	2.13	0.48
1:A:319:LEU:HD23	1:A:323:THR:HG21	1.95	0.48
1:A:222:LEU:HD23	1:A:260:LEU:HD22	1.95	0.48
1:A:70:ASP:N	1:A:70:ASP:OD1	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:GLU:HA	1:A:356:THR:HG22	1.96	0.48
2:B:147:ILE:O	2:B:147:ILE:HG22	2.14	0.47
1:A:128:GLN:NE2	8:A:803:HOH:O	2.37	0.47
1:A:392:GLU:OE2	4:A:702:MTA:O3'	2.26	0.47
2:B:119:PHE:CD2	2:B:157:GLN:HG2	2.49	0.47
1:A:495:GLU:OE1	1:A:587:LYS:HE2	2.14	0.47
1:A:629:THR:CG2	1:A:629:THR:O	2.61	0.47
1:A:41:LEU:CD2	1:A:43:MET:HE2	2.45	0.47
1:A:617:VAL:H	1:A:623:SER:HG	1.61	0.47
2:B:92:ARG:NH1	2:B:110:GLU:HA	2.31	0.46
1:A:588:GLN:HB2	6:A:712:EDO:O1	2.15	0.46
1:A:556:GLU:OE2	1:A:568:ARG:HD3	2.15	0.46
1:A:624:ALA:H	6:A:712:EDO:H22	1.79	0.46
1:A:398:VAL:HG21	1:A:417:SER:HB2	1.98	0.46
2:B:77:VAL:HG22	2:B:78:GLN:H	1.81	0.46
2:B:108:LEU:HD22	2:B:112:GLU:HA	1.98	0.46
2:B:113:THR:C	2:B:114:LEU:HD12	2.36	0.46
1:A:369:GLY:H	1:A:370:PRO:HD3	1.82	0.45
1:A:446:SER:HB3	1:A:447:PRO:HD3	1.98	0.45
2:B:280:LEU:HG	2:B:281:ALA:N	2.31	0.45
1:A:167:ILE:HD11	2:B:165:ALA:HB2	1.98	0.45
1:A:369:GLY:N	1:A:370:PRO:HD3	2.32	0.45
1:A:504:VAL:O	1:A:581:PRO:HD2	2.17	0.44
2:B:239:LEU:HB3	2:B:249:LEU:HD23	2.00	0.44
2:B:275:SER:HB3	2:B:277:ASP:OD1	2.17	0.44
1:A:371:LEU:HD13	1:A:433:VAL:CG1	2.48	0.44
2:B:130:THR:HB	2:B:173:VAL:HG22	2.00	0.44
2:B:187:SER:HB3	2:B:189:ASP:OD1	2.17	0.44
1:A:67:THR:HG23	1:A:68:ARG:O	2.17	0.44
2:B:92:ARG:HH12	2:B:110:GLU:HA	1.82	0.44
2:B:27:MET:HG2	2:B:320:HIS:CE1	2.53	0.44
1:A:524:PRO:HG3	1:A:526:ARG:HH21	1.82	0.44
1:A:338:GLN:NE2	1:A:381:GLN:OE1	2.51	0.43
1:A:193:ARG:HH11	1:A:478:SER:HB3	1.83	0.43
1:A:627:ASN:N	1:A:628:PRO:CD	2.80	0.43
1:A:565:LEU:HD22	1:A:576:MET:SD	2.58	0.43
2:B:146:ASP:HA	6:B:401:EDO:H21	2.01	0.43
1:A:23:VAL:O	1:A:23:VAL:HG13	2.18	0.43
1:A:525:ASN:OD1	1:A:530:ILE:HG23	2.18	0.43
1:A:23:VAL:HG23	1:A:29:THR:OG1	2.18	0.43
2:B:77:VAL:HG22	2:B:78:GLN:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:GLU:O	1:A:289:GLN:HG2	2.19	0.42
2:B:86:LEU:C	2:B:86:LEU:HD12	2.39	0.42
1:A:26:ILE:HG22	8:B:501:HOH:O	2.19	0.42
1:A:369:GLY:N	1:A:370:PRO:CD	2.83	0.42
2:B:95:LEU:HD23	2:B:95:LEU:HA	1.98	0.42
2:B:27:MET:CG	2:B:320:HIS:HE1	2.32	0.42
1:A:393:LYS:HE2	1:A:393:LYS:HB2	1.69	0.42
1:A:290:ASN:O	1:A:290:ASN:CG	2.58	0.42
1:A:479:LYS:HB2	1:A:508:ASN:HD22	1.85	0.42
1:A:193:ARG:NH1	1:A:478:SER:HB3	2.35	0.41
1:A:315:LEU:HD23	1:A:315:LEU:HA	1.87	0.41
2:B:21:PRO:HG2	2:B:114:LEU:HD11	2.02	0.41
2:B:86:LEU:O	2:B:86:LEU:HD12	2.21	0.41
1:A:297:TYR:HB2	8:A:853:HOH:O	2.20	0.41
1:A:336:GLN:OE1	1:A:563:ILE:HG23	2.20	0.41
2:B:230:PHE:CE1	2:B:238:SER:HB2	2.55	0.41
1:A:567:ILE:HG21	1:A:579:TRP:HB2	2.02	0.41
2:B:42:LEU:C	2:B:42:LEU:HD23	2.42	0.41
2:B:146:ASP:HA	6:B:401:EDO:H22	2.02	0.41
1:A:479:LYS:O	1:A:483:GLU:HG3	2.22	0.40
2:B:196:ASP:OD1	2:B:198:ARG:HB2	2.22	0.40
2:B:219:ALA:HB3	2:B:229:VAL:CG2	2.51	0.40
2:B:280:LEU:HD13	2:B:315:THR:HG21	2.03	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	623/645 (97%)	584 (94%)	37 (6%)	2 (0%)	37	41
2	B	300/350 (86%)	281 (94%)	16 (5%)	3 (1%)	13	10

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	923/995 (93%)	865 (94%)	53 (6%)	5 (0%)	25	25

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	147	ILE
1	A	276	GLU
2	B	249	LEU
2	B	109	ASP
1	A	242	GLY

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	558/570 (98%)	528 (95%)	30 (5%)	18	18
2	B	257/298 (86%)	244 (95%)	13 (5%)	20	21
All	All	815/868 (94%)	772 (95%)	43 (5%)	19	19

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ILE
1	A	35	LYS
1	A	60	LYS
1	A	147	HIS
1	A	174	THR
1	A	176	THR
1	A	177	GLU
1	A	215	ASN
1	A	218	ILE
1	A	235	ILE
1	A	249	MET
1	A	256	ARG
1	A	272	HIS

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Mol	Chain	Res	Type
1	A	284	LEU
1	A	302	LYS
1	A	332	ILE
1	A	354	LYS
1	A	361	LEU
1	A	385	ARG
1	A	387	LYS
1	A	393	LYS
1	A	401	LEU
1	A	417	SER
1	A	478	SER
1	A	479	LYS
1	A	490	LYS
1	A	529	MET
1	A	602	PHE
1	A	610	LYS
1	A	629	THR
2	B	27	MET
2	B	29	ARG
2	B	47	SER
2	B	80	GLU
2	B	86	LEU
2	B	95	LEU
2	B	111	ASN
2	B	136	SER
2	B	138	THR
2	B	179	LYS
2	B	204	SER
2	B	280	LEU
2	B	288	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	508	ASN
2	B	224	GLN
2	B	320	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.



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

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

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

16 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	SO4	A	703	-	4,4,4	0.34	0	6,6,6	0.42	0
7	PEG	A	713	-	6,6,6	0.23	0	5,5,5	0.13	0
6	EDO	A	707	-	3,3,3	0.43	0	2,2,2	0.56	0
3	A1BLM	A	701	-	36,36,36	0.30	0	47,49,49	0.96	2 (4%)
5	SO4	A	704	-	4,4,4	0.32	0	6,6,6	0.35	0
6	EDO	A	708	-	3,3,3	0.52	0	2,2,2	0.13	0
6	EDO	A	709	-	3,3,3	0.47	0	2,2,2	0.29	0
6	EDO	A	710	-	3,3,3	0.53	0	2,2,2	0.34	0
6	EDO	A	712	-	3,3,3	0.74	0	2,2,2	0.23	0
6	EDO	B	402	-	3,3,3	0.49	0	2,2,2	0.18	0
6	EDO	B	403	-	3,3,3	0.45	0	2,2,2	0.40	0
5	SO4	A	705	-	4,4,4	0.35	0	6,6,6	0.60	0
6	EDO	A	711	-	3,3,3	0.69	0	2,2,2	0.14	0
6	EDO	A	706	-	3,3,3	0.50	0	2,2,2	0.30	0
4	MTA	A	702	-	18,22,22	1.40	3 (16%)	18,32,32	1.32	3 (16%)
6	EDO	B	401	-	3,3,3	0.50	0	2,2,2	0.35	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
7	PEG	A	713	-	-	1/4/4/4	-
6	EDO	A	707	-	-	0/1/1/1	-
3	A1BLM	A	701	-	-	4/21/40/40	0/4/4/4
6	EDO	A	708	-	-	1/1/1/1	-
6	EDO	A	712	-	-	0/1/1/1	-
6	EDO	A	709	-	-	1/1/1/1	-
6	EDO	A	710	-	-	1/1/1/1	-
6	EDO	B	403	-	-	0/1/1/1	-
6	EDO	B	402	-	-	0/1/1/1	-
6	EDO	A	711	-	-	1/1/1/1	-
6	EDO	A	706	-	-	0/1/1/1	-
4	MTA	A	702	-	-	3/3/23/23	0/3/3/3
6	EDO	B	401	-	-	1/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	702	MTA	C1'-N9	-3.18	1.42	1.49
4	A	702	MTA	C4-N3	-2.86	1.31	1.35
4	A	702	MTA	C8-N7	-2.81	1.29	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	A1BLM	C22-C23-N5	4.30	117.08	112.14
4	A	702	MTA	C5-C6-N6	2.65	124.34	120.31
4	A	702	MTA	O2'-C2'-C3'	2.59	120.13	111.82
4	A	702	MTA	O4'-C4'-C5'	2.28	114.69	108.83
3	A	701	A1BLM	C8-C7-N1	-2.21	117.29	122.07

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	702	MTA	C4'-C5'-S5'-CS
6	A	710	EDO	O1-C1-C2-O2
3	A	701	A1BLM	C3-C4-N1-C7
3	A	701	A1BLM	N4-C11-C9-N2
3	A	701	A1BLM	N4-C11-C9-C8
4	A	702	MTA	O4'-C4'-C5'-S5'
6	B	401	EDO	O1-C1-C2-O2
3	A	701	A1BLM	C5-C4-N1-C7

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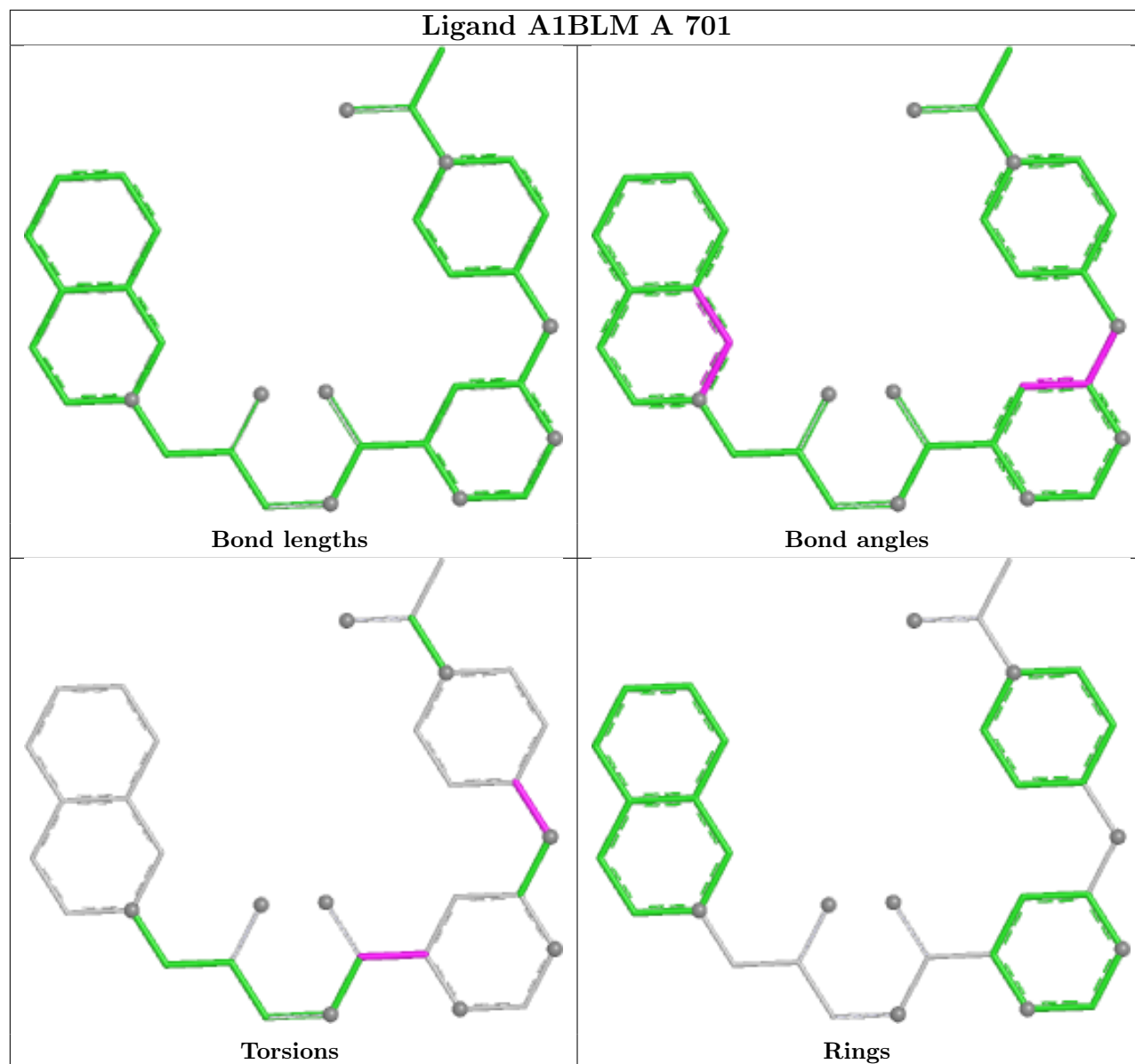
Mol	Chain	Res	Type	Atoms
6	A	708	EDO	O1-C1-C2-O2
4	A	702	MTA	C3'-C4'-C5'-S5'
6	A	709	EDO	O1-C1-C2-O2
6	A	711	EDO	O1-C1-C2-O2
7	A	713	PEG	O2-C3-C4-O4

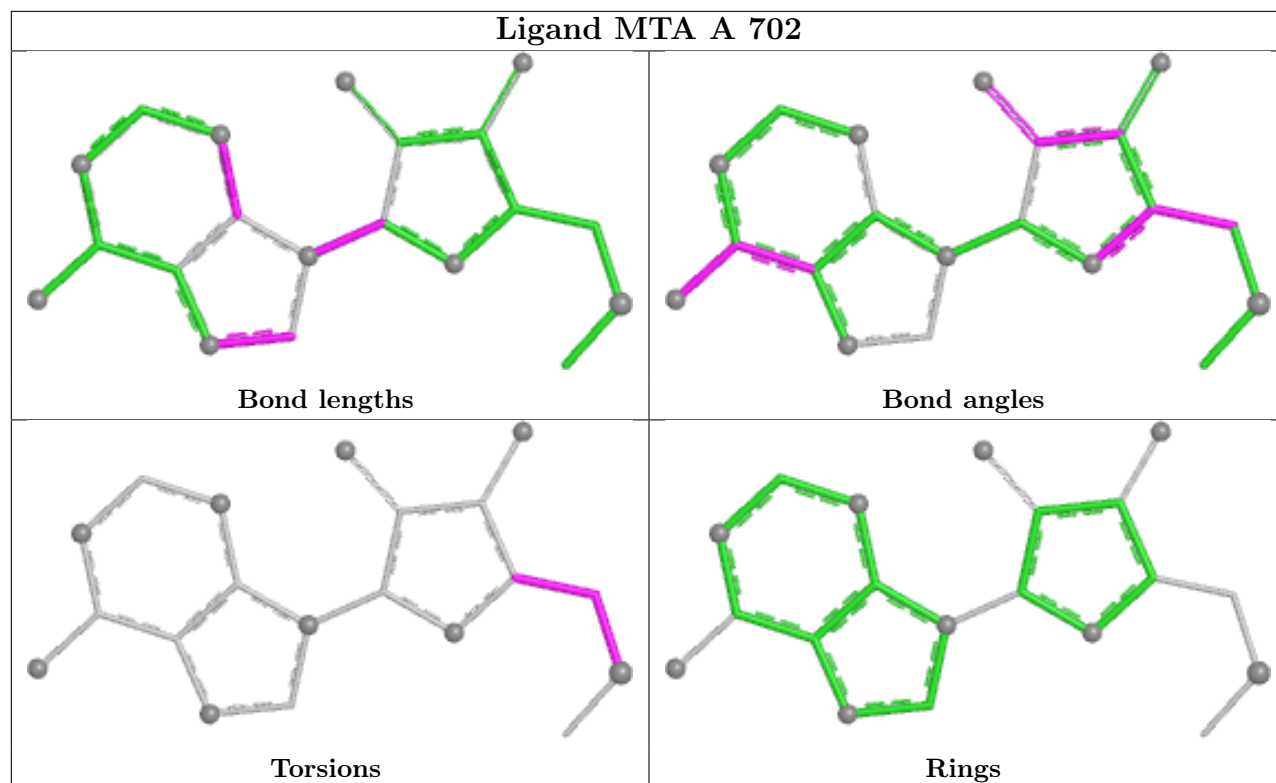
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	712	EDO	3	0
4	A	702	MTA	1	0
6	B	401	EDO	3	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, 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	625/645 (96%)	-0.20	6 (0%) 79 80	13, 37, 72, 96	0
2	B	304/350 (86%)	0.21	3 (0%) 79 80	30, 50, 76, 109	0
All	All	929/995 (93%)	-0.07	9 (0%) 79 80	13, 43, 74, 109	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	21	PRO	4.0
2	B	248	VAL	2.9
1	A	300	PHE	2.8
1	A	146	HIS	2.8
2	B	210	ALA	2.7
1	A	290	ASN	2.4
1	A	341	ILE	2.1
1	A	293	PRO	2.1
1	A	243	PHE	2.0

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

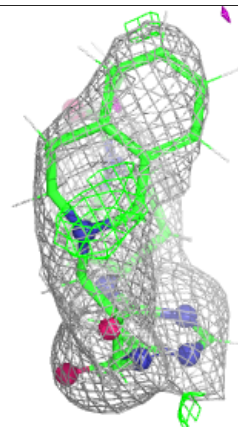
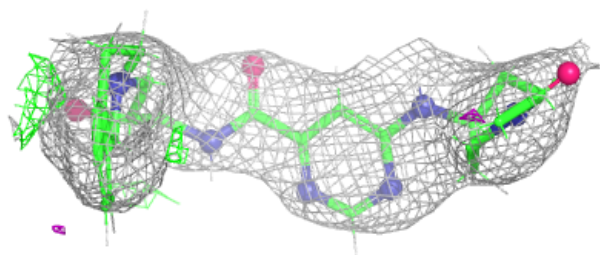
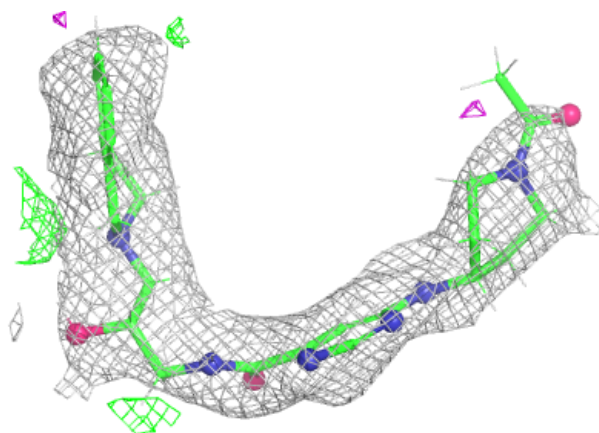
median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
7	PEG	A	713	7/7	0.58	0.16	55,67,84,84	0
6	EDO	A	708	4/4	0.69	0.13	52,62,67,73	0
5	SO4	A	704	5/5	0.73	0.11	70,72,79,83	0
6	EDO	B	401	4/4	0.76	0.20	40,53,63,65	0
6	EDO	A	711	4/4	0.85	0.18	38,46,53,56	0
5	SO4	A	705	5/5	0.86	0.10	53,55,70,87	0
5	SO4	A	703	5/5	0.86	0.17	59,61,67,89	0
6	EDO	A	709	4/4	0.86	0.09	48,57,65,65	0
3	A1BLM	A	701	33/33	0.90	0.11	36,56,77,88	0
6	EDO	B	402	4/4	0.91	0.07	45,54,61,61	0
6	EDO	A	712	4/4	0.92	0.18	25,30,35,40	0
6	EDO	A	706	4/4	0.92	0.08	30,36,39,44	0
6	EDO	B	403	4/4	0.94	0.07	48,57,59,65	0
6	EDO	A	710	4/4	0.94	0.09	44,53,64,64	0
6	EDO	A	707	4/4	0.95	0.07	34,41,47,50	0
4	MTA	A	702	20/20	0.95	0.07	19,27,46,46	0

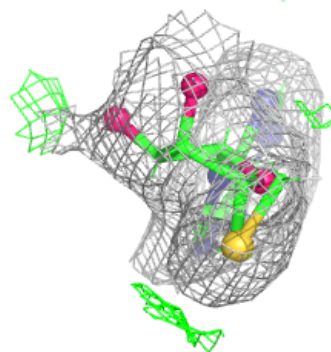
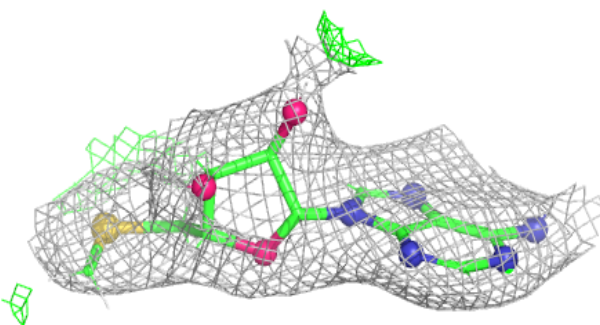
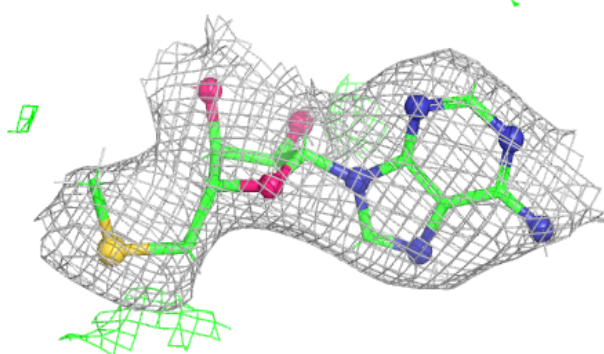
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around A1BLM A 701:**

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

**Electron density around MTA A 702:**

$2mF_o-DF_c$  (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.