Full wwPDB X-ray Structure Validation Report (i)

Oct 19, 2026 – 11:41 PM JST

PDB ID	:	9X5H / pdb_00009x5h
Title	:	Magnaporthe oryzae myosin I in complex with calmodulin
Authors	:	Liu, S.T.; Zhang, F.; Zhou, Y.X.; Zhou, X.E.
Deposited on	:	2025-10-13
Resolution	:	1.82 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

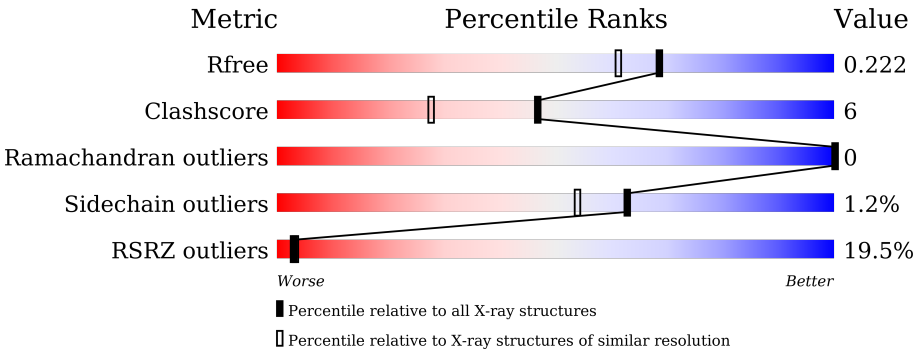
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

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 1.82 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}	180053	1112 (1.82-1.82)
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	695	7% 89% 10%
2	B	138	80% 75% 17% 7%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 7165 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 Myosin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	687	5509	3505	943	1040	21	0	0	0

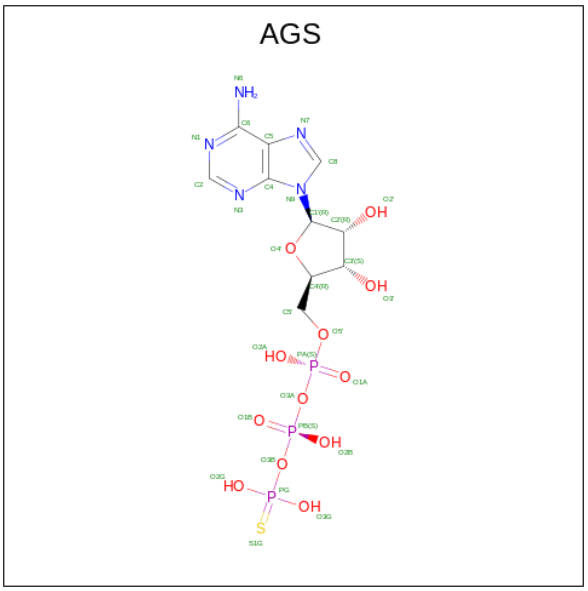
- Molecule 2 is a protein called Calmodulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	128	919	556	151	205	7	0	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
3	A	1	1	1	0	0

- Molecule 4 is PHOSPHOTHIOPHOSPHORIC ACID-ADENYLATE ESTER (CCD ID: AGS) (formula: $C_{10}H_{16}N_5O_{12}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		

- Molecule 5 is water.

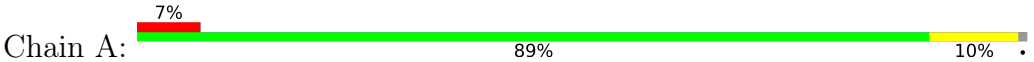
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	683	Total	O	0	0
			683	683		
5	B	22	Total	O	0	0
			22	22		

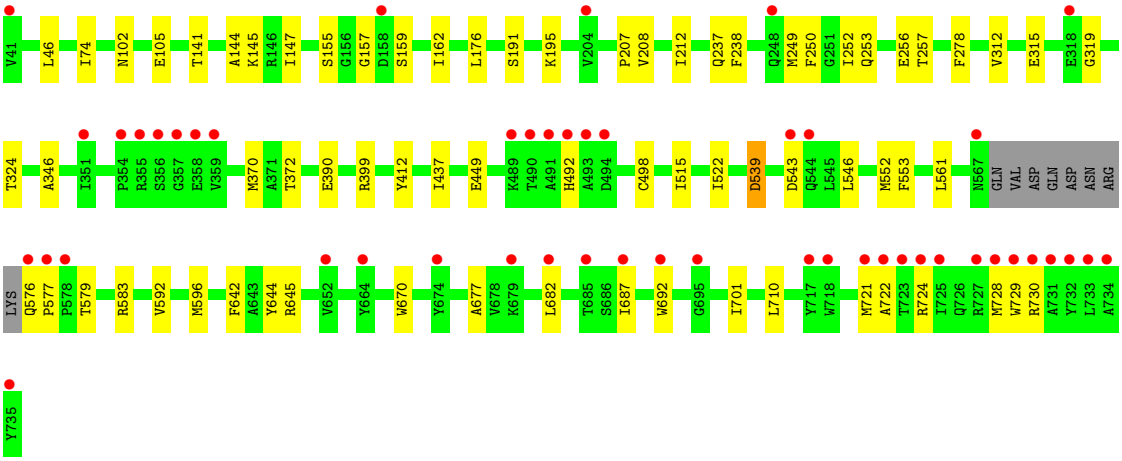
3 Residue-property plots (i)

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: Myosin-1

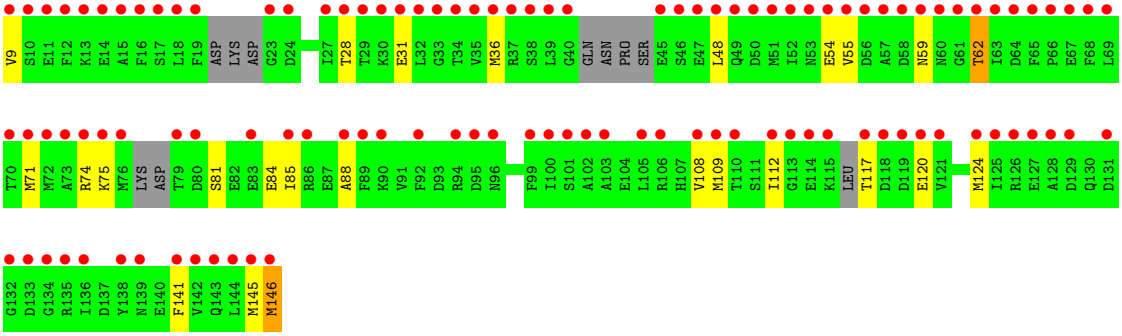
Chain A:





- Molecule 2: Calmodulin

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	58.53Å 101.86Å 167.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.00 – 1.82 49.00 – 1.82	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.00-1.82) 99.6 (49.00-1.82)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.74 (at 1.82Å)	Xtriage
Refinement program	PHENIX v1.16	Depositor
R, R_{free}	0.195 , 0.221 0.195 , 0.222	Depositor DCC
R_{free} test set	6379 reflections (7.04%)	wwPDB-VP
Wilson B-factor (Å^2)	21.2	Xtriage
Anisotropy	0.681	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 44.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7165	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: MG, AGS

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	0.39	0/5625	0.51	0/7615
2	B	0.16	0/921	0.34	0/1237
All	All	0.36	0/6546	0.49	0/8852

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	5509	0	5447	56	0
2	B	919	0	795	29	0
3	A	1	0	0	0	0
4	A	31	0	12	4	0
5	A	683	0	0	7	0
5	B	22	0	0	0	0
All	All	7165	0	6254	72	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 6.

All (72) 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:109:MET:HG3	2:B:124:MET:HE1	1.64	0.78
4:A:802:AGS:PG	5:A:901:HOH:O	2.47	0.72
1:A:721:MET:HE3	2:B:88:ALA:HA	1.73	0.71
2:B:48:LEU:HD23	2:B:75:LYS:HE3	1.74	0.69
1:A:159:SER:HB3	1:A:162:ILE:HG12	1.75	0.69
2:B:117:THR:HB	2:B:120:GLU:HG3	1.73	0.68
1:A:191:SER:CB	4:A:802:AGS:S1G	2.88	0.61
2:B:28:THR:HG23	2:B:31:GLU:H	1.67	0.60
1:A:687:ILE:HD11	1:A:710:LEU:HD21	1.83	0.59
1:A:155:SER:HB2	1:A:207:PRO:HD3	1.83	0.59
2:B:108:VAL:O	2:B:112:ILE:HG22	2.02	0.58
1:A:642:PHE:O	1:A:645:ARG:NH2	2.36	0.58
1:A:728:MET:HE2	2:B:146:MET:HG2	1.86	0.58
1:A:191:SER:N	4:A:802:AGS:S1G	2.74	0.57
1:A:191:SER:HB2	4:A:802:AGS:S1G	2.44	0.57
2:B:48:LEU:HA	2:B:75:LYS:HE3	1.87	0.56
1:A:105:GLU:HG3	2:B:74:ARG:HH21	1.70	0.56
1:A:249:MET:HE3	1:A:250:PHE:CZ	2.42	0.55
1:A:312:VAL:HA	1:A:370:MET:HE1	1.88	0.55
1:A:176:LEU:HD11	1:A:212:ILE:HD13	1.90	0.54
1:A:157:GLY:HA3	1:A:162:ILE:HG21	1.88	0.54
1:A:105:GLU:CG	2:B:74:ARG:HH21	2.21	0.54
1:A:372:THR:HG23	1:A:546:LEU:HD22	1.89	0.54
1:A:102:ASN:HD21	2:B:74:ARG:NH2	2.07	0.53
2:B:81:SER:O	2:B:85:ILE:HG12	2.08	0.53
1:A:553:PHE:CD1	1:A:561:LEU:HD21	2.44	0.53
1:A:492:HIS:CE1	1:A:498:CYS:SG	3.03	0.52
1:A:238:PHE:CZ	1:A:252:ILE:HG13	2.45	0.51
1:A:312:VAL:HG23	1:A:324:THR:OG1	2.11	0.51
1:A:390:GLU:HG3	5:A:1147:HOH:O	2.10	0.50
1:A:515:ILE:HB	1:A:522:ILE:HB	1.94	0.49
1:A:670:TRP:CD1	1:A:677:ALA:HB2	2.47	0.49
1:A:543:ASP:HB3	1:A:583:ARG:HH12	1.78	0.49
1:A:729:TRP:HA	2:B:145:MET:HE1	1.96	0.48
2:B:108:VAL:HG13	2:B:112:ILE:HG21	1.95	0.48
1:A:724:ARG:HD3	2:B:84:GLU:HB3	1.95	0.48
1:A:141:THR:HG22	1:A:145:LYS:HE2	1.95	0.47
1:A:546:LEU:H	1:A:546:LEU:HD12	1.80	0.47
2:B:36:MET:SD	2:B:48:LEU:HD22	2.55	0.46
1:A:592:VAL:O	1:A:596:MET:HG2	2.15	0.46
1:A:546:LEU:HD21	5:A:1278:HOH:O	2.16	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:721:MET:HE3	2:B:88:ALA:CA	2.43	0.46
1:A:728:MET:HG3	2:B:145:MET:CE	2.46	0.46
2:B:145:MET:HB3	2:B:145:MET:HE2	1.61	0.46
1:A:449:GLU:OE1	1:A:645:ARG:HD3	2.16	0.45
1:A:644:TYR:CD2	1:A:701:ILE:HD12	2.52	0.45
1:A:644:TYR:HB3	1:A:701:ILE:HB	2.00	0.44
1:A:437:ILE:HD11	1:A:592:VAL:HG22	1.99	0.44
1:A:724:ARG:HG2	2:B:85:ILE:HD13	1.99	0.44
1:A:346:ALA:HB2	1:A:552:MET:HB2	2.01	0.43
1:A:46:LEU:HD12	1:A:74:ILE:HG12	2.00	0.43
1:A:576:GLN:N	1:A:577:PRO:CD	2.81	0.43
1:A:682:LEU:HD13	1:A:692:TRP:CD2	2.54	0.43
2:B:48:LEU:HD23	2:B:75:LYS:CE	2.46	0.42
1:A:399:ARG:HD2	5:A:1494:HOH:O	2.19	0.42
1:A:253:GLN:HB2	1:A:257:THR:OG1	2.19	0.42
2:B:54:GLU:HG3	2:B:55:VAL:HG23	2.00	0.42
1:A:729:TRP:HA	2:B:145:MET:CE	2.50	0.42
2:B:71:MET:O	2:B:75:LYS:HG3	2.18	0.42
1:A:728:MET:HG3	2:B:145:MET:HE2	2.02	0.42
2:B:59:ASN:CB	2:B:62:THR:HG22	2.50	0.41
1:A:728:MET:C	2:B:145:MET:HE1	2.46	0.41
1:A:722:ALA:CB	2:B:112:ILE:HG23	2.51	0.41
1:A:237:GLN:HB3	1:A:278:PHE:HB2	2.02	0.41
1:A:144:ALA:HA	1:A:147:ILE:HD12	2.02	0.41
1:A:539:ASP:HB2	5:A:1112:HOH:O	2.20	0.41
1:A:315:GLU:HG2	1:A:319:GLY:HA2	2.03	0.41
1:A:645:ARG:NH1	5:A:905:HOH:O	2.54	0.41
1:A:552:MET:HE3	1:A:552:MET:HB3	1.95	0.40
2:B:141:PHE:O	2:B:145:MET:HG3	2.21	0.40
1:A:579:THR:O	1:A:583:ARG:HD3	2.21	0.40
1:A:195:LYS:HE3	5:A:1199:HOH:O	2.21	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	683/695 (98%)	673 (98%)	10 (2%)	0	100	100
2	B	118/138 (86%)	118 (100%)	0	0	100	100
All	All	801/833 (96%)	791 (99%)	10 (1%)	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	597/605 (99%)	592 (99%)	5 (1%)	73	67
2	B	90/122 (74%)	87 (97%)	3 (3%)	33	15
All	All	687/727 (94%)	679 (99%)	8 (1%)	63	54

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	208	VAL
1	A	256	GLU
1	A	412	TYR
1	A	539	ASP
1	A	730	ARG
2	B	9	VAL
2	B	62	THR
2	B	146	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	245	HIS
1	A	404	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	436	GLN
1	A	634	ASN
1	A	646	GLN
1	A	689	GLN
1	A	712	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\#\|Z\| > 2$
4	AGS	A	802	3	29,33,33	2.43	6 (20%)	39,52,52	1.81	9 (23%)

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
4	AGS	A	802	3	-	1/21/38/38	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	802	AGS	O4'-C1'	8.07	1.61	1.42
4	A	802	AGS	O4'-C4'	-5.47	1.32	1.45
4	A	802	AGS	C2'-C1'	-5.29	1.36	1.53
4	A	802	AGS	C6-N6	3.64	1.43	1.34
4	A	802	AGS	C8-N9	-2.19	1.33	1.37
4	A	802	AGS	O3'-C3'	-2.04	1.38	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	802	AGS	C5-C4-N3	-5.15	120.03	126.75
4	A	802	AGS	N3-C2-N1	-4.35	121.79	128.60
4	A	802	AGS	N3-C4-N9	3.76	133.27	127.08
4	A	802	AGS	C2-N3-C4	3.57	120.19	111.75
4	A	802	AGS	N9-C8-N7	-3.10	109.67	113.91
4	A	802	AGS	C5-N7-C8	2.85	107.55	103.51
4	A	802	AGS	C4-N9-C8	2.65	108.60	105.73
4	A	802	AGS	C4-C5-N7	-2.55	107.52	110.62
4	A	802	AGS	C3'-C2'-C1'	2.39	105.97	101.43

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	802	AGS	PA-O3A-PB-O1B

There are no ring outliers.

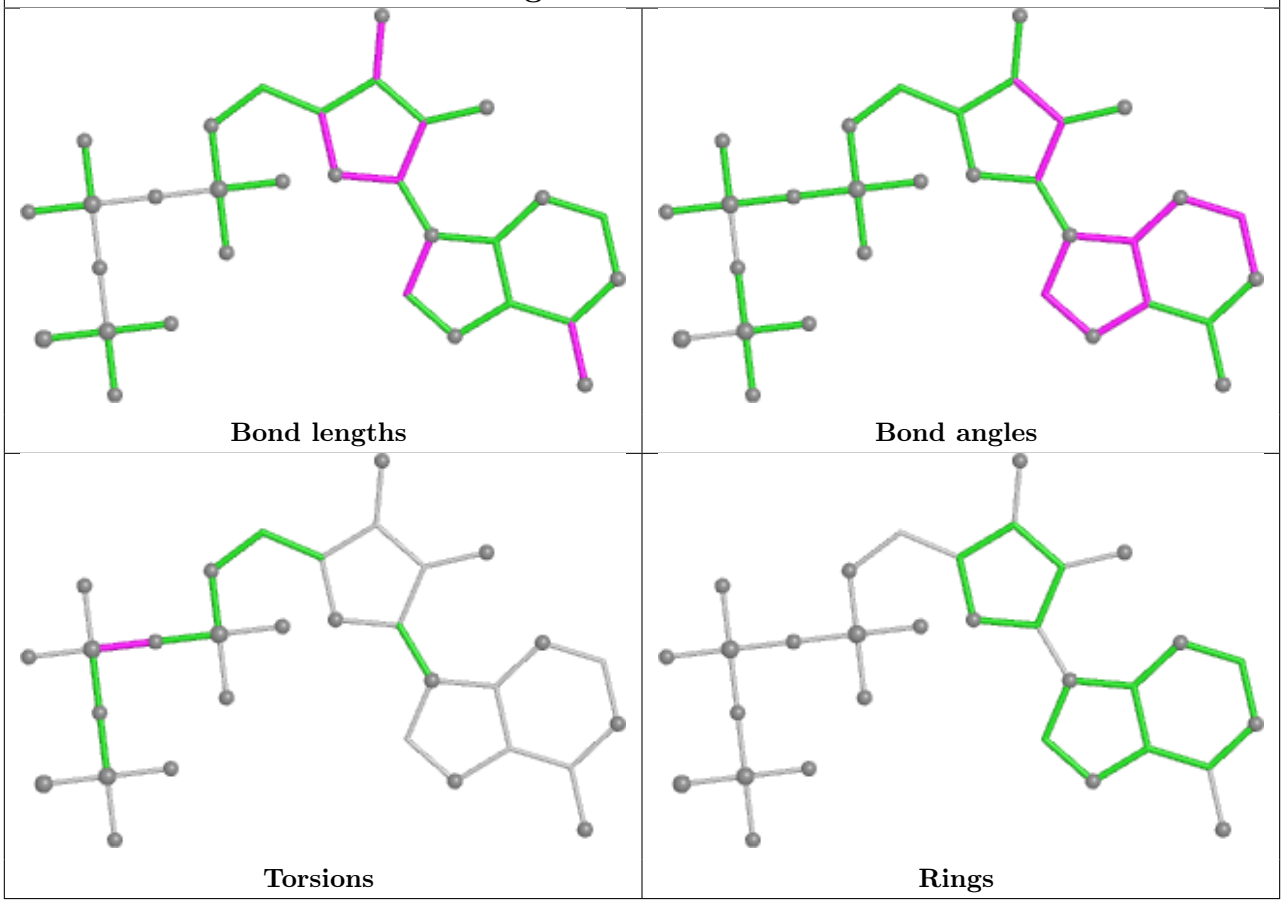
1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	802	AGS	4	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 then 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 AGS A 802



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

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^{th} 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	687/695 (98%)	0.15	49 (7%) 22 25	12, 22, 46, 62	0
2	B	128/138 (92%)	3.63	110 (85%) 0 0	42, 98, 144, 153	0
All	All	815/833 (97%)	0.70	159 (19%) 3 3	12, 24, 121, 153	0

All (159) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	112	ILE	9.3
1	A	493	ALA	8.2
2	B	55	VAL	7.4
2	B	9	VAL	7.3
2	B	73	ALA	6.9
2	B	36	MET	6.6
2	B	63	ILE	6.6
1	A	357	GLY	6.5
2	B	68	PHE	6.5
1	A	492	HIS	6.4
2	B	52	ILE	6.0
1	A	356	SER	5.9
2	B	16	PHE	5.8
2	B	132	GLY	5.7
2	B	32	LEU	5.7
2	B	12	PHE	5.6
2	B	75	LYS	5.6
2	B	125	ILE	5.5
2	B	76	MET	5.5
2	B	19	PHE	5.5
2	B	144	LEU	5.4
2	B	35	VAL	5.4
2	B	134	GLY	5.3
2	B	133	ASP	5.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	113	GLY	5.2
2	B	74	ARG	5.2
2	B	69	LEU	5.0
2	B	65	PHE	5.0
2	B	70	THR	4.8
2	B	18	LEU	4.8
2	B	15	ALA	4.7
2	B	40	GLY	4.6
2	B	117	THR	4.6
2	B	72	MET	4.6
2	B	48	LEU	4.6
2	B	128	ALA	4.5
2	B	71	MET	4.5
2	B	136	ILE	4.5
1	A	725	ILE	4.4
1	A	735	TYR	4.4
1	A	491	ALA	4.4
2	B	57	ALA	4.4
2	B	141	PHE	4.3
2	B	54	GLU	4.3
2	B	39	LEU	4.3
2	B	85	ILE	4.3
2	B	102	ALA	4.3
2	B	131	ASP	4.2
1	A	355	ARG	4.2
2	B	135	ARG	4.1
2	B	27	ILE	4.1
2	B	121	VAL	4.0
2	B	129	ASP	4.0
2	B	96	ASN	4.0
2	B	60	ASN	4.0
2	B	47	GLU	3.9
2	B	56	ASP	3.9
2	B	105	LEU	3.9
2	B	38	SER	3.9
2	B	34	THR	3.9
2	B	66	PRO	3.8
2	B	79	THR	3.8
2	B	50	ASP	3.8
2	B	124	MET	3.8
1	A	729	TRP	3.7
2	B	51	MET	3.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	67	GLU	3.7
2	B	100	ILE	3.7
2	B	142	VAL	3.6
1	A	732	TYR	3.6
1	A	354	PRO	3.6
2	B	28	THR	3.5
2	B	103	ALA	3.5
2	B	23	GLY	3.5
2	B	138	TYR	3.4
2	B	108	VAL	3.4
2	B	146	MET	3.4
2	B	95	ASP	3.4
2	B	33	GLY	3.4
2	B	62	THR	3.4
1	A	489	LYS	3.3
2	B	115	LYS	3.3
1	A	576	GLN	3.3
2	B	46	SER	3.3
1	A	577	PRO	3.3
2	B	10	SER	3.3
1	A	733	LEU	3.3
2	B	29	THR	3.3
1	A	734	ALA	3.2
2	B	139	ASN	3.2
2	B	45	GLU	3.2
2	B	30	LYS	3.1
2	B	89	PHE	3.1
1	A	567	ASN	3.1
2	B	106	ARG	3.0
1	A	351	ILE	3.0
1	A	692	TRP	3.0
1	A	718	TRP	3.0
2	B	13	LYS	3.0
2	B	119	ASP	3.0
1	A	728	MET	3.0
1	A	724	ARG	2.9
1	A	679	LYS	2.9
2	B	37	ARG	2.9
2	B	86	ARG	2.9
1	A	490	THR	2.9
2	B	110	THR	2.9
1	A	494	ASP	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	83	GLU	2.8
2	B	126	ARG	2.8
1	A	723	THR	2.8
2	B	80	ASP	2.8
1	A	721	MET	2.8
2	B	145	MET	2.7
2	B	53	ASN	2.7
2	B	11	GLU	2.7
2	B	118	ASP	2.6
1	A	358	GLU	2.6
2	B	58	ASP	2.6
1	A	687	ILE	2.6
2	B	14	GLU	2.6
1	A	359	VAL	2.6
2	B	49	GLN	2.5
2	B	31	GLU	2.5
2	B	24	ASP	2.5
1	A	685	THR	2.5
2	B	90	LYS	2.5
2	B	59	ASN	2.5
2	B	143	GLN	2.5
1	A	730	ARG	2.5
2	B	17	SER	2.5
2	B	92	PHE	2.4
2	B	61	GLY	2.4
1	A	682	LEU	2.4
2	B	99	PHE	2.4
1	A	543	ASP	2.4
2	B	64	ASP	2.4
1	A	731	ALA	2.4
2	B	114	GLU	2.4
2	B	88	ALA	2.4
2	B	120	GLU	2.4
1	A	248	GLN	2.3
1	A	204	VAL	2.3
1	A	674	TYR	2.2
1	A	717	TYR	2.2
2	B	109	MET	2.2
1	A	652	VAL	2.2
1	A	158	ASP	2.2
1	A	722	ALA	2.2
1	A	318	GLU	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	94	ARG	2.1
1	A	664	TYR	2.1
2	B	127	GLU	2.1
1	A	578	PRO	2.1
1	A	727	ARG	2.1
1	A	695	GLY	2.0
2	B	101	SER	2.0
1	A	544	GLN	2.0
1	A	41	VAL	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 oligosaccharides in this entry.

6.4 Ligands (i)

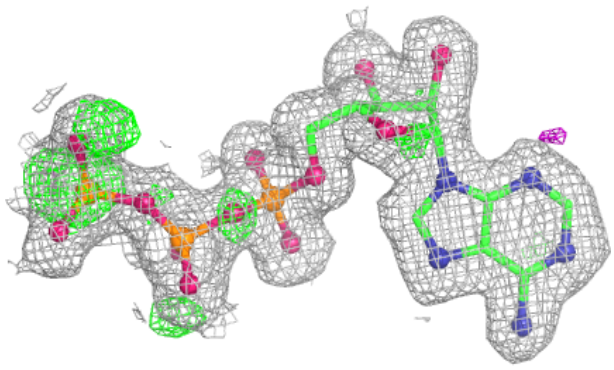
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

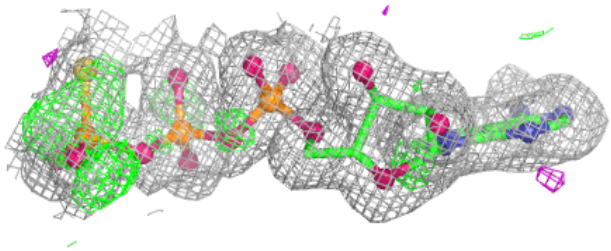
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	AGS	A	802	31/31	0.97	0.06	6,14,18,40	4
3	MG	A	801	1/1	0.99	0.02	14,14,14,14	0

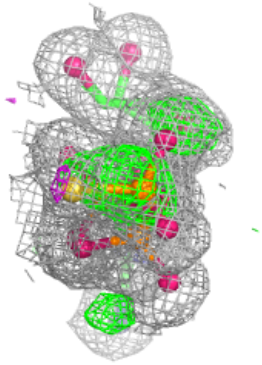
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around AGS A 802:

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

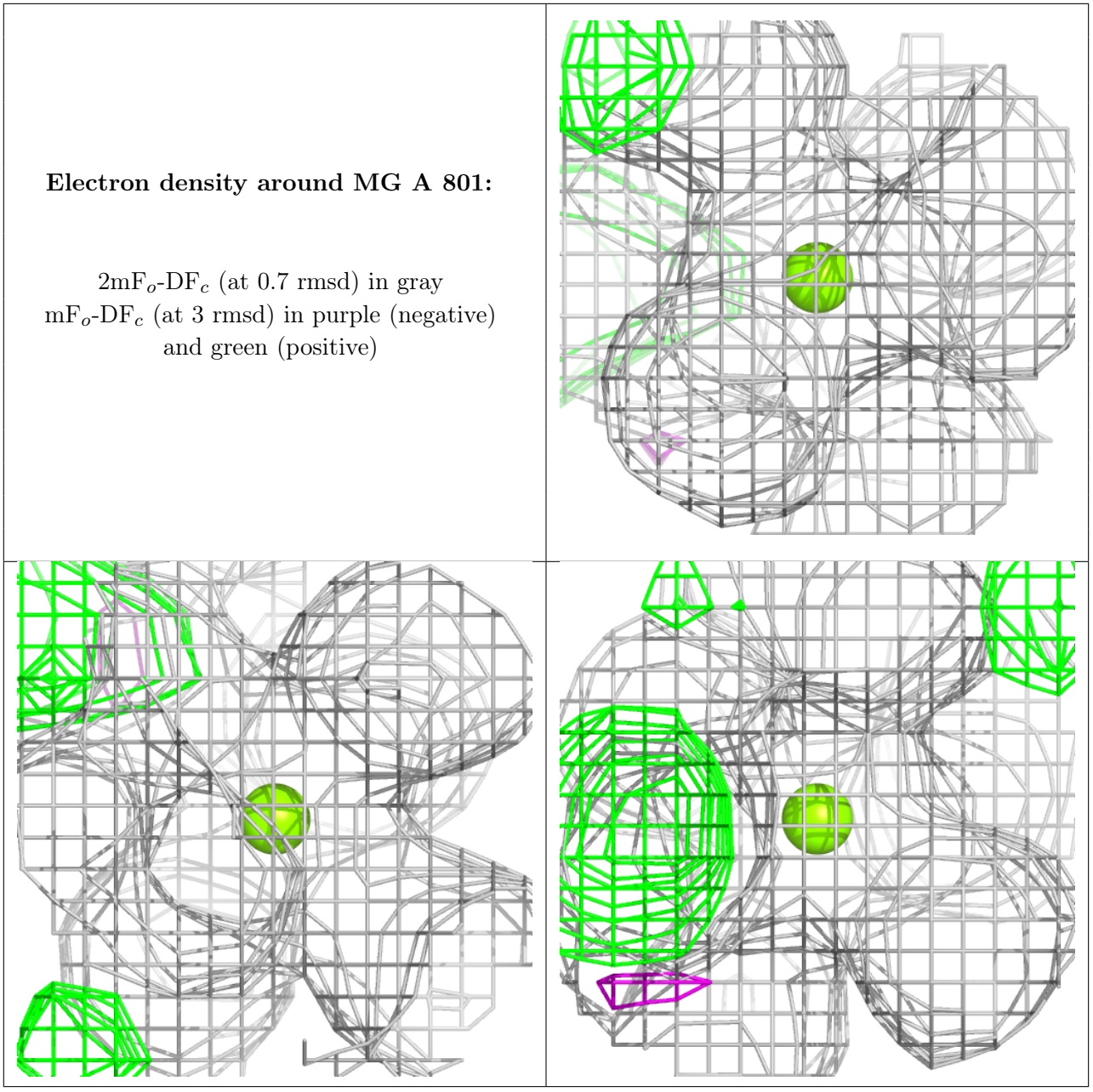






Electron density around MG A 801:

$2mF_o$-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.