



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:07 AM UTC

PDB ID : 6KWA / pdb_00006kwa
Title : AtDAO1(dioxygenase for auxin oxidation 1 from Arabidopsis thaliana)
Authors : Rhee, S.; Jin, S.; Lee, H.
Deposited on : 2019-09-06
Resolution : 2.09 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 types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

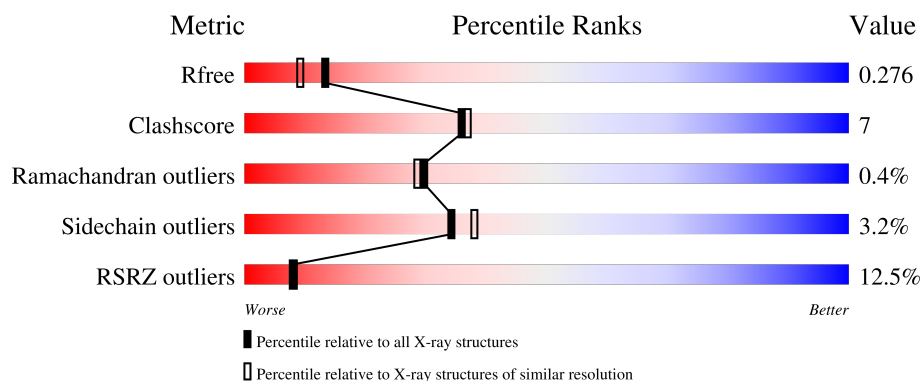
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 $\leq 5\%$. 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	270	9% 83% 14% ..
1	B	270	8% 77% 20% ..
1	C	270	18% 61% 20% • 16%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6083 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 2-oxoglutarate (2OG) and Fe(II)-dependent oxygenase super-family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			2114	1340	357	404	13			
1	B	266	Total	C	N	O	S	0	0	0
			2106	1334	356	403	13			
1	C	226	Total	C	N	O	S	0	0	0
			1805	1146	302	345	12			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	MET	-	initiating methionine	UNP Q9XI75
B	8	MET	-	initiating methionine	UNP Q9XI75
C	8	MET	-	initiating methionine	UNP Q9XI75

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	24	Total	O	0	0
			24	24		
3	B	28	Total	O	0	0
			28	28		

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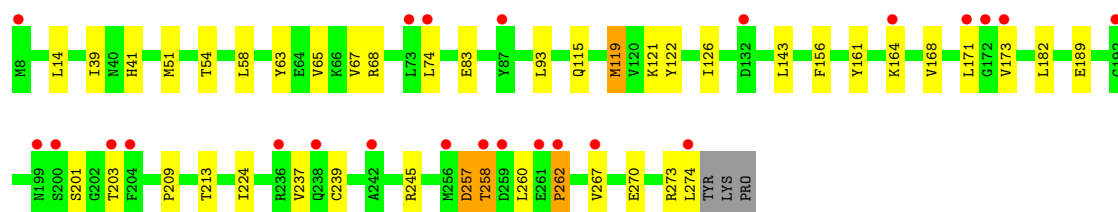
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	4	Total	O	0	0
			4	4		

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.

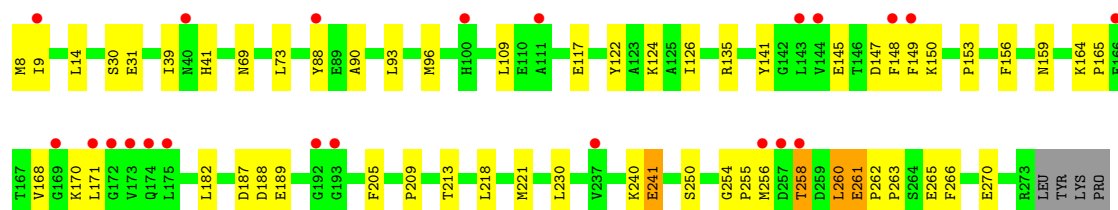
- Molecule 1: 2-oxoglutarate (2OG) and Fe(II)-dependent oxygenase superfamily protein

Chain A: 



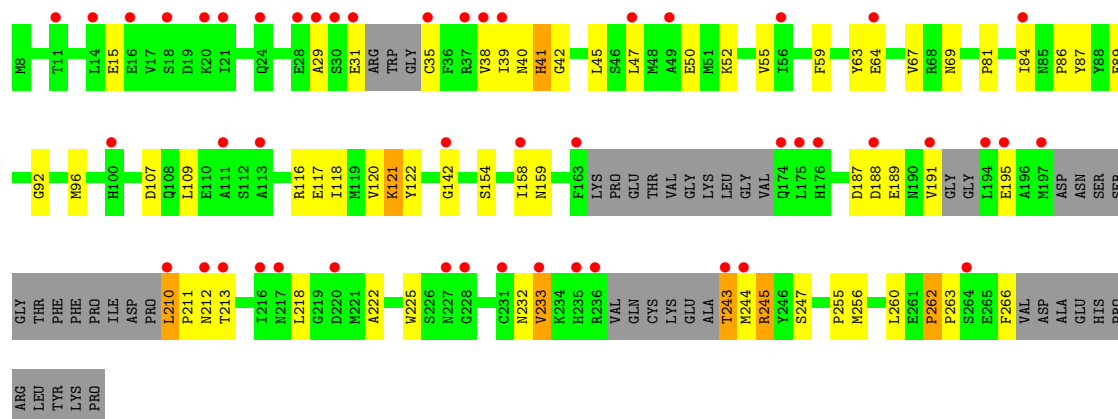
- Molecule 1: 2-oxoglutarate (2OG) and Fe(II)-dependent oxygenase superfamily protein

Chain B: 



- Molecule 1: 2-oxoglutarate (2OG) and Fe(II)-dependent oxygenase superfamily protein

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.63Å 75.55Å 165.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.88 – 2.09 33.88 – 2.09	Depositor EDS
% Data completeness (in resolution range)	99.2 (33.88-2.09) 99.1 (33.88-2.09)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.242 , 0.282 (Not available) , 0.276	Depositor DCC
R_{free} test set	2002 reflections (3.49%)	wwPDB-VP
Wilson B-factor (Å ²)	50.2	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6083	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.63	0/2160	0.92	2/2927 (0.1%)
1	B	0.60	0/2152	0.92	5/2916 (0.2%)
1	C	0.50	0/1837	0.99	7/2480 (0.3%)
All	All	0.58	0/6149	0.94	14/8323 (0.2%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	210	LEU	CA-C-N	7.51	127.47	120.03
1	C	210	LEU	C-N-CA	7.51	127.47	120.03
1	A	262	PRO	N-CA-C	6.56	117.60	110.58
1	C	42	GLY	N-CA-C	-5.97	107.42	115.36
1	C	92	GLY	N-CA-C	5.97	119.31	110.42
1	C	262	PRO	N-CA-C	5.73	116.72	110.58
1	B	261	GLU	CA-C-N	-5.56	114.66	120.38
1	B	261	GLU	C-N-CA	-5.56	114.66	120.38
1	C	38	VAL	N-CA-C	5.51	116.42	108.48
1	A	260	LEU	N-CA-C	5.50	117.67	109.59
1	B	258	THR	CA-C-N	-5.39	115.46	123.11
1	B	258	THR	C-N-CA	-5.39	115.46	123.11
1	C	233	VAL	N-CA-C	5.32	117.30	109.63
1	B	73	LEU	N-CA-C	-5.12	102.95	110.52

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	2114	0	2077	23	0
1	B	2106	0	2066	32	0
1	C	1805	0	1774	36	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	24	0	0	0	0
3	B	28	0	0	0	0
3	C	4	0	0	0	0
All	All	6083	0	5917	87	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 (87) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:GLU:OE2	1:B:88:TYR:OH	2.04	0.74
1:B:39:ILE:HG22	1:B:213:THR:HG22	1.72	0.70
1:C:210:LEU:HD12	1:C:211:PRO:HD2	1.75	0.69
1:C:187:ASP:OD1	1:C:188:ASP:N	2.25	0.66
1:C:69:ASN:HB2	1:C:109:LEU:HD11	1.78	0.65
1:A:201:SER:OG	1:A:203:THR:HG22	1.97	0.65
1:B:8:MET:SD	1:B:8:MET:N	2.71	0.64
1:C:81:PRO:HD3	1:C:89:GLU:HB3	1.77	0.63
1:A:267:VAL:HG21	1:A:273:ARG:HD3	1.82	0.61
1:C:45:LEU:HG	1:C:212:ASN:HD21	1.66	0.60
1:B:153:PRO:HD3	1:B:256:MET:HG2	1.85	0.58
1:B:221:MET:SD	1:B:260:LEU:HD11	2.44	0.58
1:B:69:ASN:HB2	1:B:109:LEU:HD11	1.87	0.57
1:A:189:GLU:OE1	1:A:189:GLU:N	2.34	0.56
1:A:39:ILE:HG22	1:A:213:THR:HG22	1.88	0.56
1:B:88:TYR:HE2	1:B:159:ASN:HB3	1.71	0.55
1:B:263:PRO:HG2	1:B:266:PHE:CD2	2.41	0.55
1:C:255:PRO:HG3	1:C:260:LEU:HB2	1.88	0.54
1:A:63:TYR:O	1:A:67:VAL:HG23	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:189:GLU:OE1	1:C:189:GLU:N	2.33	0.53
1:C:50:GLU:HB3	1:C:118:ILE:HD13	1.88	0.53
1:A:54:THR:HG23	1:A:115:GLN:HG2	1.90	0.53
1:B:31:GLU:OE2	1:B:141:TYR:OH	2.16	0.52
1:B:164:LYS:HB3	1:B:165:PRO:HD2	1.92	0.52
1:B:189:GLU:OE1	1:B:189:GLU:N	2.37	0.52
1:C:52:LYS:NZ	1:C:188:ASP:OD1	2.38	0.51
1:C:232:ASN:OD1	1:C:232:ASN:N	2.42	0.51
1:B:135:ARG:HG3	1:B:149:PHE:CZ	2.46	0.50
1:B:189:GLU:HA	1:B:209:PRO:HB2	1.94	0.50
1:B:135:ARG:HG3	1:B:149:PHE:CE2	2.48	0.49
1:A:51:MET:HE1	1:A:119:MET:SD	2.53	0.49
1:B:93:LEU:HG	1:B:156:PHE:HB3	1.93	0.49
1:C:29:ALA:C	1:C:31:GLU:H	2.21	0.49
1:C:225:TRP:O	1:C:262:PRO:HB3	2.13	0.48
1:A:65:VAL:O	1:A:68:ARG:HB2	2.13	0.48
1:C:86:PRO:HG2	1:C:87:TYR:CE2	2.48	0.48
1:A:189:GLU:HA	1:A:209:PRO:HB2	1.96	0.48
1:C:243:THR:HG22	1:C:244:MET:H	1.77	0.48
1:A:224:ILE:O	1:A:262:PRO:HG3	2.12	0.48
1:C:47:LEU:HD11	1:C:122:TYR:HA	1.95	0.48
1:B:147:ASP:HB3	1:B:150:LYS:HD3	1.96	0.48
1:B:14:LEU:HD12	1:B:41:HIS:HA	1.95	0.47
1:C:255:PRO:CG	1:C:260:LEU:HB2	2.44	0.47
1:B:240:LYS:C	1:B:241:GLU:HG3	2.40	0.47
1:C:263:PRO:HG2	1:C:266:PHE:CD2	2.50	0.47
1:A:58:LEU:HD11	1:A:119:MET:HE1	1.95	0.46
1:B:263:PRO:HG2	1:B:266:PHE:HD2	1.80	0.46
1:A:93:LEU:HG	1:A:156:PHE:HB3	1.97	0.46
1:B:30:SER:HB3	1:B:218:LEU:HD11	1.98	0.46
1:B:122:TYR:CE2	1:B:126:ILE:HG13	2.50	0.46
1:A:14:LEU:HD12	1:A:41:HIS:HA	1.96	0.46
1:C:35:CYS:HB3	1:C:233:VAL:HG11	1.97	0.46
1:A:83:GLU:HB2	1:C:142:GLY:HA3	1.99	0.45
1:C:63:TYR:O	1:C:67:VAL:HG13	2.16	0.45
1:C:45:LEU:HA	1:C:45:LEU:HD13	1.69	0.45
1:A:74:LEU:HD12	1:A:74:LEU:HA	1.76	0.45
1:A:257:ASP:HB2	1:A:258:THR:H	1.58	0.44
1:C:117:GLU:HA	1:C:120:VAL:HG22	2.00	0.44
1:C:55:VAL:HG13	1:C:158:ILE:HD12	1.99	0.44
1:B:96:MET:HE1	1:B:250:SER:OG	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:96:MET:HG2	1:C:154:SER:HB2	2.00	0.44
1:A:119:MET:HE2	1:A:119:MET:HB2	1.67	0.44
1:B:187:ASP:OD1	1:B:188:ASP:N	2.47	0.43
1:B:255:PRO:HG3	1:B:260:LEU:HB3	2.00	0.43
1:C:15:GLU:HG3	1:C:40:ASN:HB3	1.99	0.43
1:A:168:VAL:HA	1:A:239:CYS:SG	2.59	0.43
1:C:191:VAL:CG1	1:C:245:ARG:HE	2.32	0.43
1:A:122:TYR:CE2	1:A:126:ILE:HG13	2.53	0.43
1:C:41:HIS:NE2	1:C:213:THR:HA	2.34	0.43
1:A:173:VAL:HG12	1:A:237:VAL:HG12	2.01	0.42
1:C:107:ASP:OD2	1:C:116:ARG:NH1	2.41	0.42
1:B:265:GLU:HB3	1:C:84:ILE:HD12	2.01	0.42
1:C:263:PRO:HG2	1:C:266:PHE:HD2	1.84	0.42
1:C:59:PHE:CE1	1:C:89:GLU:HG3	2.55	0.42
1:A:121:LYS:HD3	1:A:121:LYS:HA	1.76	0.41
1:C:39:ILE:HG12	1:C:40:ASN:N	2.36	0.41
1:B:153:PRO:HD2	1:B:254:GLY:O	2.20	0.41
1:C:64:GLU:HA	1:C:67:VAL:HG22	2.02	0.41
1:A:161:TYR:HB2	1:A:245:ARG:HB3	2.02	0.41
1:B:30:SER:HB3	1:B:230:LEU:HD22	2.02	0.41
1:B:270:GLU:OE2	1:C:159:ASN:ND2	2.48	0.41
1:C:218:LEU:HD23	1:C:222:ALA:HB1	2.03	0.41
1:C:121:LYS:HB2	1:C:121:LYS:HE3	1.91	0.41
1:B:145:GLU:OE2	1:B:148:PHE:HA	2.21	0.41
1:B:88:TYR:CZ	1:B:90:ALA:HB2	2.57	0.40
1:B:261:GLU:HG3	1:B:262:PRO:O	2.20	0.40
1:B:9:ILE:HG23	1:B:205:PHE:CE1	2.56	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	265/270 (98%)	253 (96%)	10 (4%)	2 (1%)	16	12
1	B	264/270 (98%)	254 (96%)	10 (4%)	0	100	100
1	C	214/270 (79%)	205 (96%)	8 (4%)	1 (0%)	24	22
All	All	743/810 (92%)	712 (96%)	28 (4%)	3 (0%)	30	28

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	THR
1	C	256	MET
1	A	257	ASP

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	232/235 (99%)	226 (97%)	6 (3%)	40	46
1	B	231/235 (98%)	222 (96%)	9 (4%)	28	31
1	C	199/235 (85%)	193 (97%)	6 (3%)	36	41
All	All	662/705 (94%)	641 (97%)	21 (3%)	34	38

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	MET
1	A	143	LEU
1	A	164	LYS
1	A	171	LEU
1	A	182	LEU
1	A	274	LEU
1	B	117	GLU
1	B	124	LYS
1	B	168	VAL
1	B	170	LYS

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Mol	Chain	Res	Type
1	B	171	LEU
1	B	182	LEU
1	B	241	GLU
1	B	258	THR
1	B	260	LEU
1	C	41	HIS
1	C	121	LYS
1	C	195	GLU
1	C	243	THR
1	C	245	ARG
1	C	247	SER

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

Mol	Chain	Res	Type
1	A	85	ASN
1	A	176	HIS
1	B	85	ASN
1	B	162	HIS
1	C	212	ASN

5.3.3 RNA [i](#)

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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [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	267/270 (98%)	0.70	24 (8%) 15 16	36, 54, 87, 101	0
1	B	266/270 (98%)	0.80	22 (8%) 17 18	39, 57, 91, 104	0
1	C	226/270 (83%)	1.41	49 (21%) 2 2	51, 76, 98, 108	0
All	All	759/810 (93%)	0.95	95 (12%) 8 8	36, 62, 94, 108	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	173	VAL	6.8
1	B	173	VAL	5.8
1	C	28	GLU	5.3
1	B	149	PHE	5.2
1	A	172	GLY	5.1
1	A	171	LEU	4.7
1	A	200	SER	4.6
1	C	227	ASN	4.6
1	B	193	GLY	4.5
1	C	176	HIS	4.5
1	C	163	PHE	4.3
1	C	236	ARG	4.0
1	C	84	ILE	3.9
1	C	210	LEU	3.9
1	C	35	CYS	3.7
1	C	191	VAL	3.6
1	C	231	CYS	3.6
1	C	47	LEU	3.6
1	A	262	PRO	3.5
1	C	233	VAL	3.5
1	B	148	PHE	3.4
1	C	244	MET	3.4
1	C	30	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	11	THR	3.4
1	C	49	ALA	3.4
1	C	197	MET	3.3
1	A	242	ALA	3.3
1	C	39	ILE	3.3
1	C	37	ARG	3.3
1	B	174	GLN	3.3
1	C	235	HIS	3.2
1	C	38	VAL	3.2
1	C	217	ASN	3.2
1	C	113	ALA	3.2
1	B	237	VAL	3.1
1	B	88	TYR	3.1
1	B	256	MET	3.1
1	C	175	LEU	3.0
1	A	192	GLY	3.0
1	C	31	GLU	2.9
1	B	171	LEU	2.9
1	A	199	ASN	2.9
1	B	192	GLY	2.9
1	C	174	GLN	2.9
1	C	195	GLU	2.8
1	A	259	ASP	2.8
1	C	213	THR	2.8
1	A	258	THR	2.8
1	B	258	THR	2.7
1	B	143	LEU	2.7
1	C	111	ALA	2.7
1	A	261	GLU	2.7
1	C	188	ASP	2.6
1	C	220	ASP	2.6
1	B	111	ALA	2.5
1	A	74	LEU	2.5
1	A	238	GLN	2.4
1	B	144	VAL	2.4
1	C	20	LYS	2.4
1	B	257	ASP	2.4
1	B	100	HIS	2.4
1	C	100	HIS	2.4
1	B	172	GLY	2.3
1	C	18	SER	2.3
1	C	264	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	64	GLU	2.3
1	C	228	GLY	2.3
1	C	194	LEU	2.3
1	A	256	MET	2.3
1	C	24	GLN	2.3
1	C	16	GLU	2.2
1	C	21	ILE	2.2
1	C	216	ILE	2.2
1	A	8	MET	2.2
1	A	132	ASP	2.2
1	C	56	ILE	2.2
1	C	243	THR	2.2
1	B	169	GLY	2.2
1	A	73	LEU	2.2
1	A	236	ARG	2.2
1	A	274	LEU	2.1
1	B	40	ASN	2.1
1	C	212	ASN	2.1
1	A	87	TYR	2.1
1	C	158	ILE	2.1
1	A	267	VAL	2.1
1	C	29	ALA	2.1
1	B	175	LEU	2.1
1	B	166	GLU	2.1
1	A	204	PHE	2.1
1	C	14	LEU	2.0
1	A	164	LYS	2.0
1	B	9	ILE	2.0
1	A	203	THR	2.0
1	C	142	GLY	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

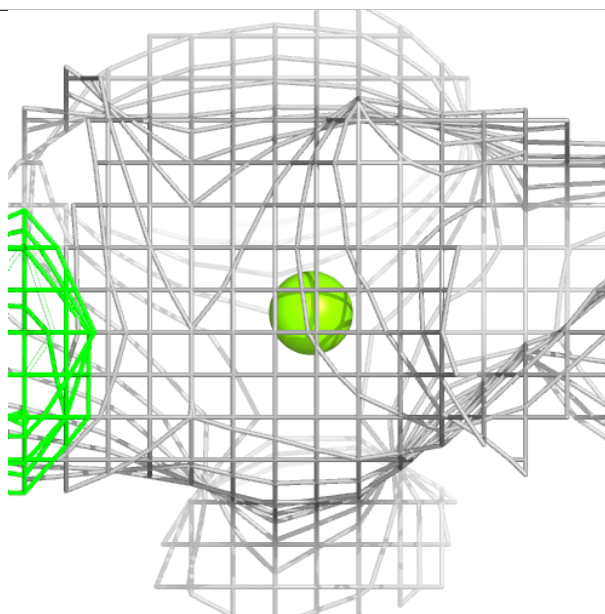
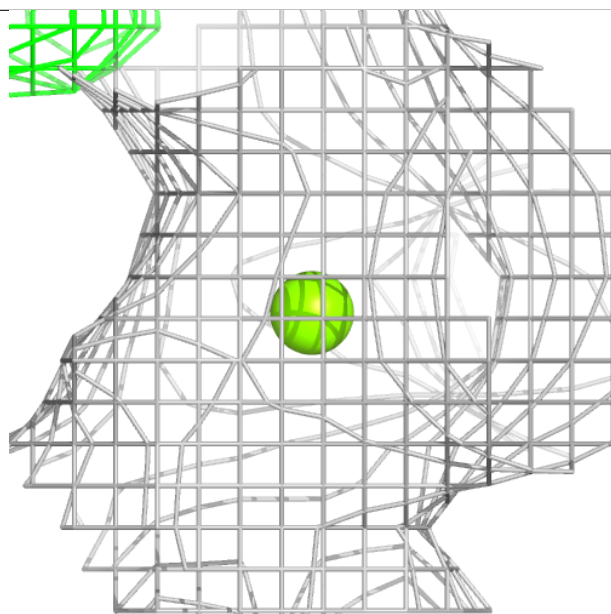
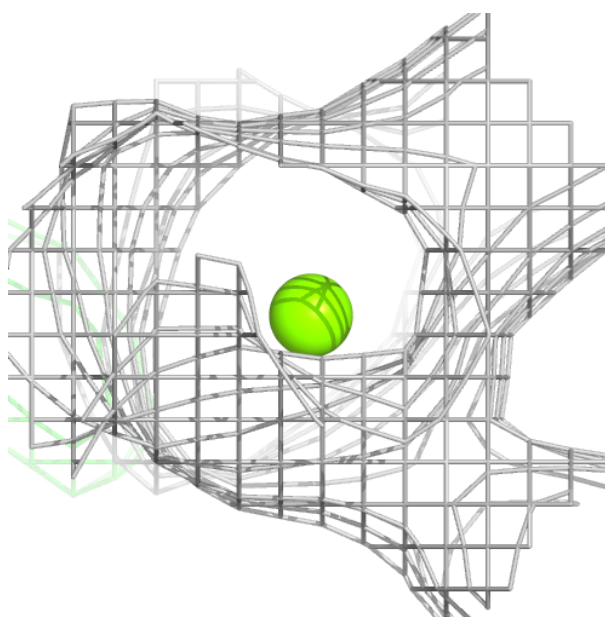
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	A	301	1/1	0.94	0.12	65,65,65,65	0
2	MG	B	301	1/1	0.96	0.07	61,61,61,61	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.

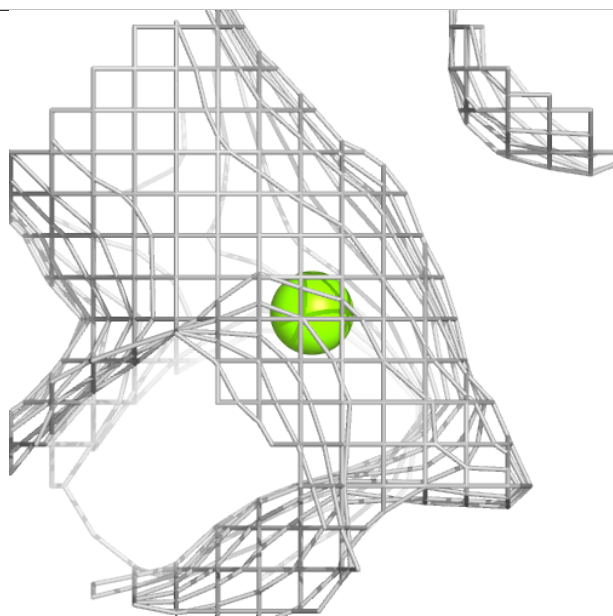
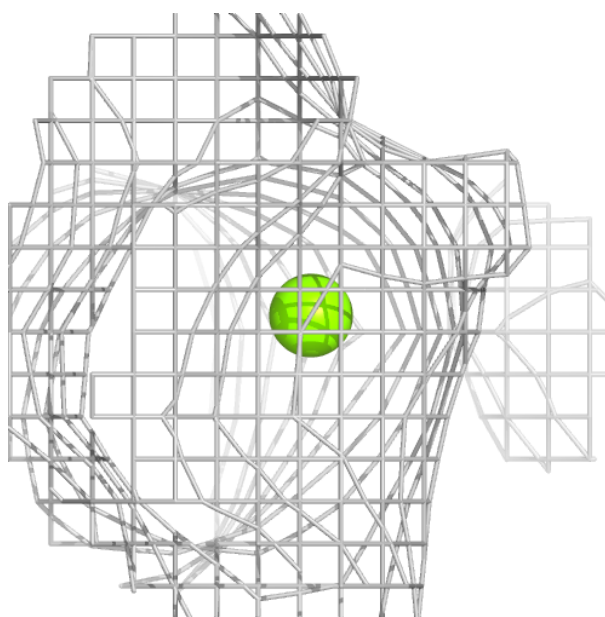
Electron density around MG A 301:

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



Electron density around MG B 301:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.