



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 06:57 PM BST

PDB ID : 6VCZ
Title : Crystal structure of Arabidopsis thaliana S-adenosylmethionine Synthase 2 (AtMAT2)
Authors : Sekula, B.; Ruszkowski, M.; Dauter, Z.
Deposited on : 2019-12-23
Resolution : 1.52 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 : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

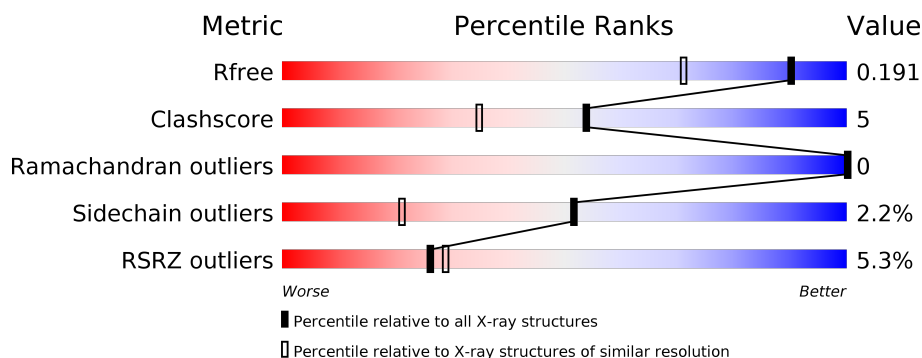
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.52 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	4009 (1.54-1.50)
Clashscore	141614	4249 (1.54-1.50)
Ramachandran outliers	138981	4148 (1.54-1.50)
Sidechain outliers	138945	4146 (1.54-1.50)
RSRZ outliers	127900	3943 (1.54-1.50)

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	402	<div> <div>5%</div> <div> <div></div> <div>87%</div> <div>7%</div> <div>6%</div> </div> </div>
1	B	402	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>8%</div> <div>5%</div> </div> </div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7107 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 S-adenosylmethionine synthase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	378	Total	C	N	O	S	0	10	0
			2976	1883	509	566	18			
1	B	380	Total	C	N	O	S	0	14	0
			3012	1908	513	573	18			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP P17562
A	-1	ASN	-	expression tag	UNP P17562
A	0	ALA	-	expression tag	UNP P17562
A	393	LEU	-	expression tag	UNP P17562
A	394	ASN	-	expression tag	UNP P17562
A	395	ASN	-	expression tag	UNP P17562
A	396	ILE	-	expression tag	UNP P17562
A	397	GLY	-	expression tag	UNP P17562
A	398	SER	-	expression tag	UNP P17562
A	399	GLY	-	expression tag	UNP P17562
B	-2	SER	-	expression tag	UNP P17562
B	-1	ASN	-	expression tag	UNP P17562
B	0	ALA	-	expression tag	UNP P17562
B	393	LEU	-	expression tag	UNP P17562
B	394	ASN	-	expression tag	UNP P17562
B	395	ASN	-	expression tag	UNP P17562
B	396	ILE	-	expression tag	UNP P17562
B	397	GLY	-	expression tag	UNP P17562
B	398	SER	-	expression tag	UNP P17562
B	399	GLY	-	expression tag	UNP P17562

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).

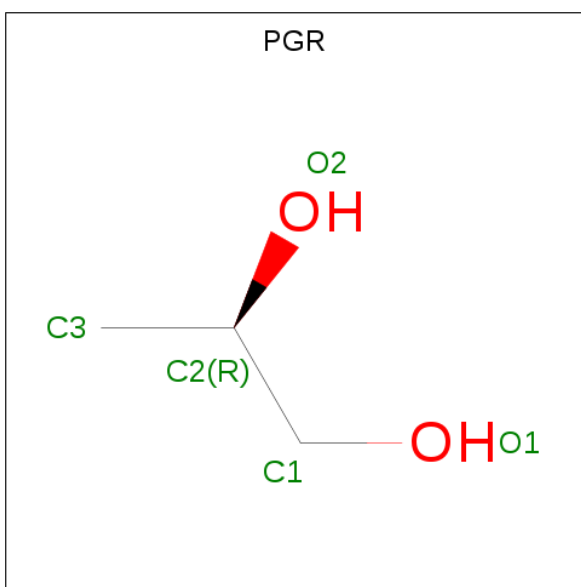


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by author).

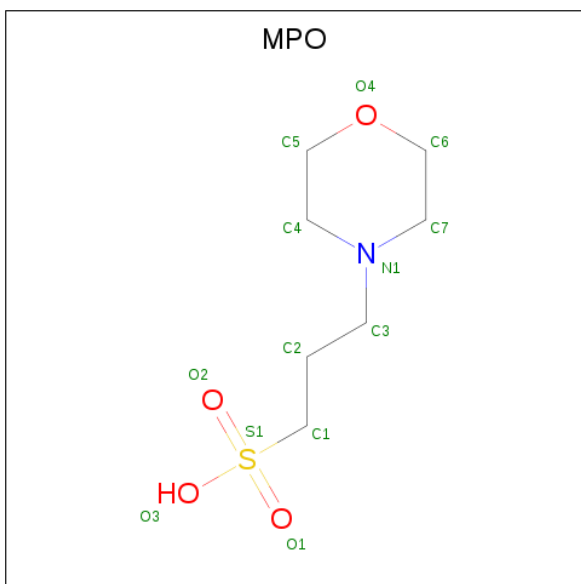
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total	Mg	0	0
			4	4		
3	A	3	Total	Mg	0	0
			3	3		

- Molecule 4 is R-1,2-PROPANEDIOL (three-letter code: PGR) (formula: C₃H₈O₂).



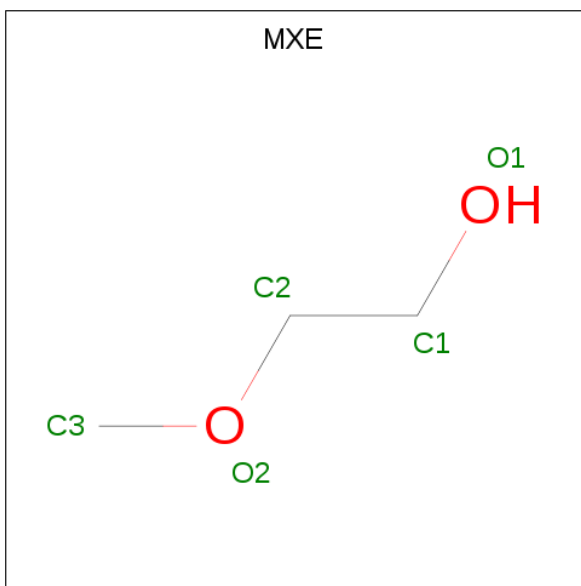
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			5	3	2		
4	B	1	Total	C	O	0	0
			5	3	2		

- Molecule 5 is 3[N-MORPHOLINO]PROPANE SULFONIC ACID (three-letter code: MPO) (formula: $C_7H_{15}NO_4S$).



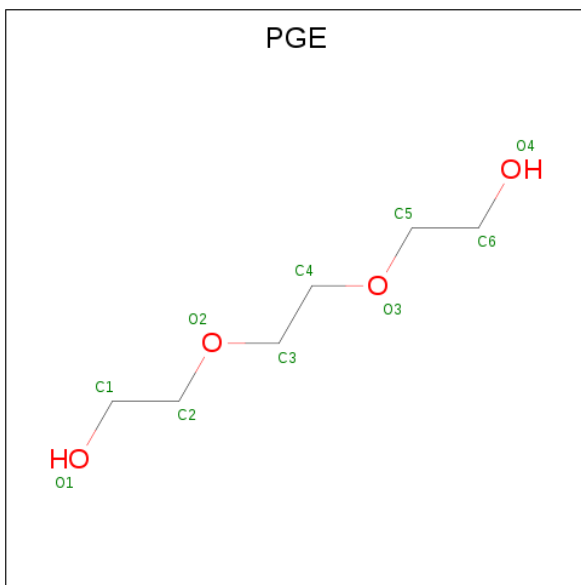
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			13	7	1	4	1		

- Molecule 6 is 2-METHOXYETHANOL (three-letter code: MXE) (formula: $C_3H_8O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			5	3	2		

- Molecule 7 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			10	6	4		
7	B	1	Total	C	O	0	0
			10	6	4		

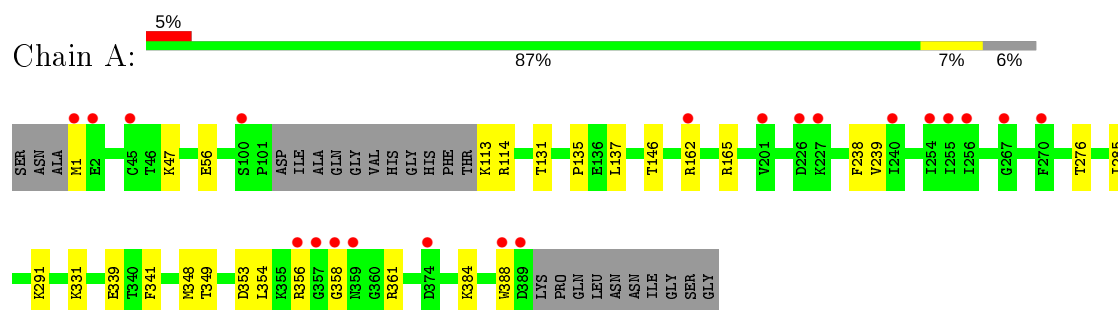
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	498	Total 501	O 501	0	3
8	B	516	Total 523	O 523	0	7

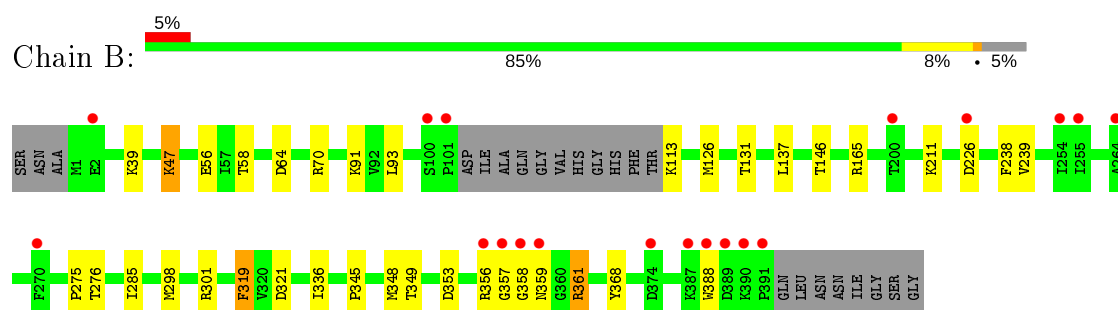
3 Residue-property plots [i](#)

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: S-adenosylmethionine synthase 2



• Molecule 1: S-adenosylmethionine synthase 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.80 Å 99.28 Å 86.93 Å 90.00° 109.74° 90.00°	Depositor
Resolution (Å)	42.44 – 1.52 42.44 – 1.52	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.44-1.52) 99.8 (42.44-1.52)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.52 Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.160 , 0.180 0.172 , 0.191	Depositor DCC
R_{free} test set	1542 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7107	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.26% 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, PGE, MPO, PO4, PGR, MXE

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	0.87	0/3066	0.96	0/4149
1	B	0.87	0/3115	0.97	0/4214
All	All	0.87	0/6181	0.97	0/8363

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	2976	0	2987	27	0
1	B	3012	0	3035	36	0
2	A	20	0	0	1	0
2	B	20	0	0	1	0
3	A	3	0	0	0	0
3	B	4	0	0	0	0
4	A	5	0	8	1	0
4	B	5	0	8	0	0
5	B	13	0	14	4	0
6	B	5	0	8	0	0
7	B	20	0	28	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	501	0	0	12	0
8	B	523	0	0	19	0
All	All	7107	0	6088	65	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (65) 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:B:361:ARG:HB2	8:B:505:HOH:O	1.39	1.20
1:A:146[B]:THR:HG21	8:A:665:HOH:O	1.72	0.88
1:A:56:GLU:HG2	8:A:517:HOH:O	1.80	0.79
1:B:91[A]:LYS:HE2	1:B:93:LEU:HD21	1.64	0.77
1:B:276:THR:HG23	8:B:697:HOH:O	1.83	0.76
1:B:146[B]:THR:HG21	8:B:678:HOH:O	1.86	0.74
1:A:388:TRP:HA	8:A:823:HOH:O	1.88	0.73
1:B:359:ASN:ND2	8:B:503:HOH:O	2.23	0.72
1:A:353:ASP:OD2	1:A:356:ARG:NE	2.24	0.70
1:B:358:GLY:N	8:B:505:HOH:O	2.26	0.69
1:B:56:GLU:HG3	8:B:511:HOH:O	1.95	0.67
1:A:146[B]:THR:CG2	8:A:665:HOH:O	2.36	0.66
1:B:146[B]:THR:CG2	8:B:678:HOH:O	2.44	0.64
1:A:56:GLU:HG3	8:A:702:HOH:O	1.97	0.64
1:B:358:GLY:CA	8:B:508:HOH:O	2.47	0.61
1:A:276:THR:HG22	1:A:349[B]:THR:HG21	1.84	0.60
1:B:64:ASP:OD2	5:B:409:MPO:N1	2.36	0.59
1:B:357:GLY:C	8:B:508:HOH:O	2.41	0.59
1:A:114:ARG:HG3	1:A:114:ARG:HH21	1.67	0.58
1:B:358:GLY:HA2	8:B:508:HOH:O	2.03	0.58
1:B:358:GLY:N	8:B:508:HOH:O	2.36	0.58
1:A:358:GLY:N	8:A:502:HOH:O	2.22	0.58
1:B:113:LYS:HE2	8:B:829:HOH:O	2.03	0.57
1:B:211[B]:LYS:HG3	8:B:774:HOH:O	2.03	0.57
1:A:165:ARG:NH1	1:A:239[A]:VAL:HG13	2.20	0.57
1:B:56:GLU:HG3	8:B:704:HOH:O	2.09	0.52
1:B:275:PRO:O	1:B:349[B]:THR:HG22	2.10	0.52
1:A:135:PRO:HD2	4:A:408:PGR:H33	1.90	0.52
1:A:291:LYS:HZ1	2:A:404:PO4:P	2.33	0.51
1:A:276:THR:HG23	8:A:604:HOH:O	2.09	0.51
1:B:131:THR:O	1:B:137:LEU:HA	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239[A]:VAL:CG1	8:A:756:HOH:O	2.59	0.49
1:B:359:ASN:CG	8:B:503:HOH:O	2.51	0.49
7:B:413:PGE:O1	7:B:413:PGE:H3	2.13	0.49
1:B:276:THR:HA	1:B:349[B]:THR:HG21	1.95	0.48
1:B:319[B]:PHE:CZ	1:B:321[B]:ASP:HB2	2.49	0.48
1:B:91[A]:LYS:CE	1:B:93:LEU:HD21	2.37	0.48
1:A:276:THR:CG2	1:A:349[B]:THR:HG21	2.43	0.48
1:A:131:THR:O	1:A:137:LEU:HA	2.13	0.47
1:B:64:ASP:OD2	5:B:409:MPO:H31	2.14	0.47
1:B:285:ILE:HG21	1:B:348:MET:SD	2.55	0.47
1:B:39[B]:LYS:HB2	1:B:58:THR:HB	1.98	0.46
1:A:353:ASP:OD2	1:A:356:ARG:NH1	2.48	0.45
1:A:349[A]:THR:HG23	1:A:354:LEU:HB2	1.97	0.45
1:B:298:MET:HE2	1:B:388:TRP:HB3	1.98	0.45
1:A:276:THR:HG22	8:A:785:HOH:O	2.16	0.44
1:A:1:MET:HA	8:A:544:HOH:O	2.16	0.44
1:B:64:ASP:OD2	5:B:409:MPO:C3	2.65	0.44
1:B:165:ARG:NH1	1:B:239[B]:VAL:HG13	2.33	0.43
2:B:404:PO4:O2	8:B:501:HOH:O	2.21	0.42
1:A:276:THR:OG1	8:A:501:HOH:O	2.21	0.42
1:B:70:ARG:HD2	5:B:409:MPO:O2	2.19	0.42
1:A:239[A]:VAL:HG11	8:A:756:HOH:O	2.19	0.42
1:A:341:PHE:CE1	1:A:384:LYS:HD3	2.54	0.42
1:A:331:LYS:HB2	1:A:331:LYS:HE2	1.60	0.42
1:B:368:TYR:HA	8:B:543:HOH:O	2.20	0.42
1:A:165:ARG:HH12	1:A:239[A]:VAL:HG13	1.85	0.41
1:B:113:LYS:N	8:B:531:HOH:O	2.53	0.41
1:B:298:MET:HE1	1:B:336:ILE:HG21	2.02	0.41
1:B:47[A]:LYS:HB2	1:B:47[A]:LYS:HE3	1.92	0.41
1:B:353:ASP:CG	1:B:356:ARG:HD3	2.42	0.41
1:A:388:TRP:CG	1:A:388:TRP:O	2.72	0.40
1:A:285:ILE:HG21	1:A:348:MET:SD	2.62	0.40
1:B:56:GLU:CG	8:B:511:HOH:O	2.62	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	384/402 (96%)	377 (98%)	7 (2%)	0	100	100
1	B	390/402 (97%)	382 (98%)	8 (2%)	0	100	100
All	All	774/804 (96%)	759 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	329/337 (98%)	322 (98%)	7 (2%)	53	22
1	B	335/337 (99%)	325 (97%)	10 (3%)	41	12
All	All	664/674 (98%)	647 (97%)	17 (3%)	52	16

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47[A]	LYS
1	A	47[B]	LYS
1	A	113	LYS
1	A	162	ARG
1	A	238	PHE
1	A	339	GLU
1	A	361	ARG
1	B	47[A]	LYS

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Mol	Chain	Res	Type
1	B	47[B]	LYS
1	B	126	MET
1	B	226	ASP
1	B	238	PHE
1	B	301	ARG
1	B	319[A]	PHE
1	B	319[B]	PHE
1	B	345	PRO
1	B	361	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 21 ligands modelled in this entry, 7 are monoatomic - leaving 14 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	$\# Z > 2$
2	PO4	B	402	3	4,4,4	0.70	0	6,6,6	0.92	0
2	PO4	A	404	-	4,4,4	1.02	0	6,6,6	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PGR	B	411	-	3,4,4	0.27	0	1,4,4	0.33	0
4	PGR	A	408	-	3,4,4	0.23	0	1,4,4	0.08	0
2	PO4	A	401	3	4,4,4	1.16	0	6,6,6	0.47	0
2	PO4	B	401	3	4,4,4	1.19	0	6,6,6	0.53	0
7	PGE	B	412	-	9,9,9	0.23	0	8,8,8	0.26	0
6	MXE	B	410	-	4,4,4	0.06	0	3,3,3	0.27	0
2	PO4	B	404	3	4,4,4	0.56	0	6,6,6	0.37	0
2	PO4	A	402	3	4,4,4	1.03	0	6,6,6	0.73	0
2	PO4	B	403	-	4,4,4	0.91	0	6,6,6	0.70	0
2	PO4	A	403	-	4,4,4	0.69	0	6,6,6	0.46	0
5	MPO	B	409	-	13,13,13	0.82	1 (7%)	17,17,17	1.10	2 (11%)
7	PGE	B	413	-	9,9,9	0.25	0	8,8,8	0.23	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
4	PGR	B	411	-	-	1/2/2/2	-
4	PGR	A	408	-	-	0/2/2/2	-
7	PGE	B	412	-	-	4/7/7/7	-
6	MXE	B	410	-	-	1/2/2/2	-
5	MPO	B	409	-	-	4/7/15/15	0/1/1/1
7	PGE	B	413	-	-	4/7/7/7	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	409	MPO	O2-S1	2.51	1.52	1.45

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	409	MPO	O3-S1-O1	2.85	118.24	111.27
5	B	409	MPO	C2-C1-S1	-2.15	109.96	113.25

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	411	PGR	O1-C1-C2-O2
5	B	409	MPO	C2-C3-N1-C7
7	B	412	PGE	O1-C1-C2-O2
7	B	412	PGE	O3-C5-C6-O4
7	B	413	PGE	O2-C3-C4-O3
5	B	409	MPO	C2-C3-N1-C4
7	B	412	PGE	O2-C3-C4-O3
7	B	413	PGE	O1-C1-C2-O2
7	B	413	PGE	O3-C5-C6-O4
5	B	409	MPO	C1-C2-C3-N1
7	B	412	PGE	C3-C4-O3-C5
7	B	413	PGE	C3-C4-O3-C5
5	B	409	MPO	S1-C1-C2-C3
6	B	410	MXE	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	404	PO4	1	0
4	A	408	PGR	1	0
2	B	404	PO4	1	0
5	B	409	MPO	4	0
7	B	413	PGE	1	0

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, 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	378/402 (94%)	0.16	21 (5%) 24 26	11, 18, 34, 60	0
1	B	380/402 (94%)	0.18	19 (5%) 28 31	11, 19, 36, 59	0
All	All	758/804 (94%)	0.17	40 (5%) 26 29	11, 19, 35, 60	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	389	ASP	5.7
1	B	358	GLY	5.1
1	B	356	ARG	4.8
1	B	100	SER	4.4
1	B	391	PRO	4.2
1	B	357	GLY	4.0
1	A	357	GLY	3.9
1	A	359	ASN	3.8
1	A	255	ILE	3.7
1	B	390	LYS	3.6
1	B	254	ILE	3.6
1	A	1	MET	3.4
1	B	359	ASN	3.3
1	A	374	ASP	3.3
1	A	356	ARG	3.2
1	B	374	ASP	3.2
1	A	254	ILE	3.2
1	A	388	TRP	3.2
1	B	226	ASP	3.1
1	B	387	LYS	3.1
1	A	226	ASP	3.0
1	B	255	ILE	3.0
1	B	389	ASP	2.9
1	A	2	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	101	PRO	2.8
1	A	256	ILE	2.5
1	A	358	GLY	2.5
1	B	270	PHE	2.3
1	B	388	TRP	2.3
1	A	100	SER	2.2
1	A	270	PHE	2.2
1	B	264	ALA	2.2
1	A	201	VAL	2.2
1	A	227	LYS	2.2
1	A	240	ILE	2.2
1	A	162	ARG	2.1
1	A	45[A]	CYS	2.1
1	B	2	GLU	2.1
1	A	267	GLY	2.0
1	B	200	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 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, 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
4	PGR	B	411	5/5	0.57	0.16	44,48,51,51	0
4	PGR	A	408	5/5	0.64	0.15	46,48,52,55	0
2	PO4	A	404	5/5	0.76	0.22	29,29,34,35	5
7	PGE	B	413	10/10	0.81	0.16	42,50,54,56	0
6	MXE	B	410	5/5	0.83	0.25	47,48,51,52	0
7	PGE	B	412	10/10	0.88	0.17	23,52,53,55	0
2	PO4	A	403	5/5	0.92	0.15	27,32,36,39	5

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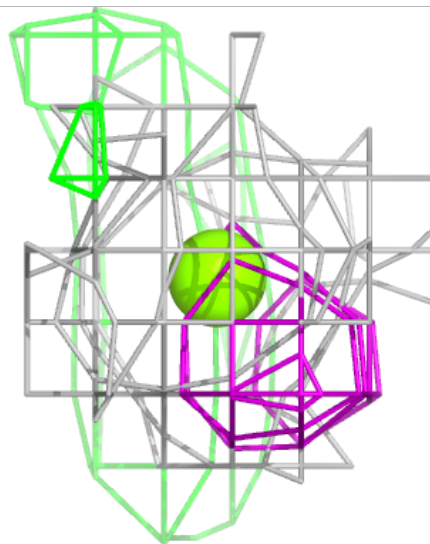
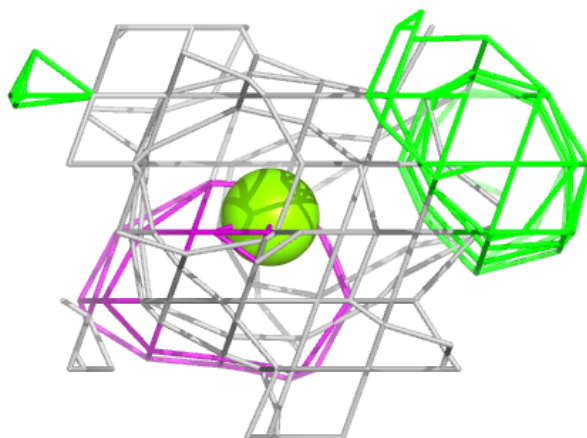
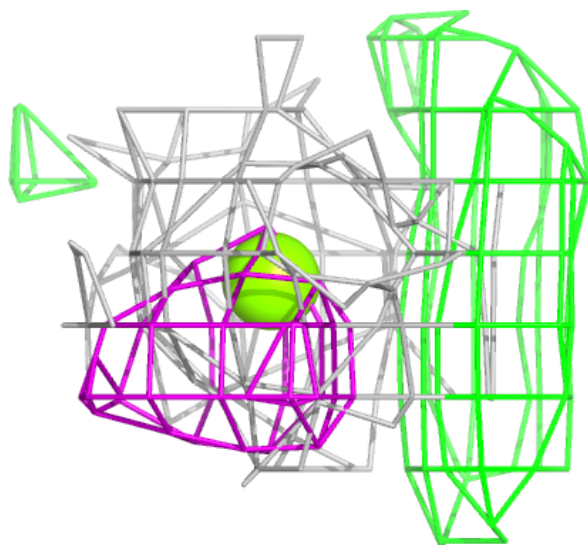
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	MPO	B	409	13/13	0.92	0.16	25,30,33,34	13
2	PO4	B	404	5/5	0.93	0.16	19,27,34,35	5
3	MG	B	408	1/1	0.96	0.06	19,19,19,19	1
2	PO4	B	403	5/5	0.96	0.09	26,26,32,36	5
3	MG	A	407	1/1	0.97	0.06	16,16,16,16	1
3	MG	B	407	1/1	0.98	0.06	18,18,18,18	0
3	MG	A	406	1/1	0.99	0.07	20,20,20,20	0
2	PO4	B	402	5/5	0.99	0.07	14,15,16,18	0
2	PO4	A	401	5/5	0.99	0.06	15,16,17,17	0
3	MG	B	406	1/1	0.99	0.08	17,17,17,17	0
2	PO4	B	401	5/5	0.99	0.07	13,13,15,15	0
2	PO4	A	402	5/5	0.99	0.07	14,16,18,21	0
3	MG	B	405	1/1	1.00	0.05	12,12,12,12	0
3	MG	A	405	1/1	1.00	0.06	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.

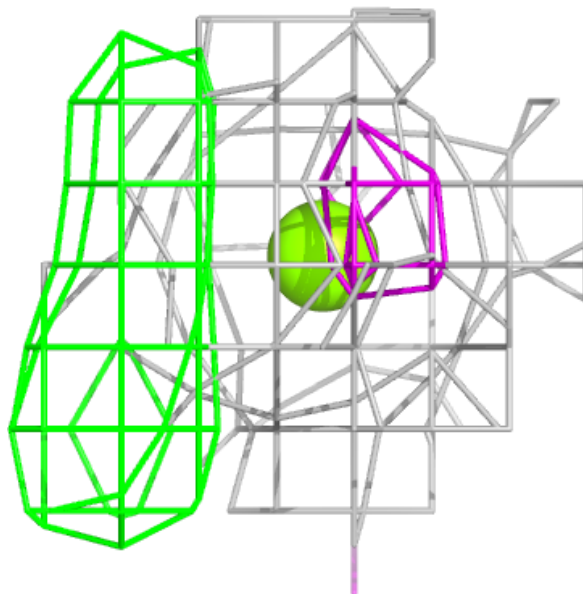
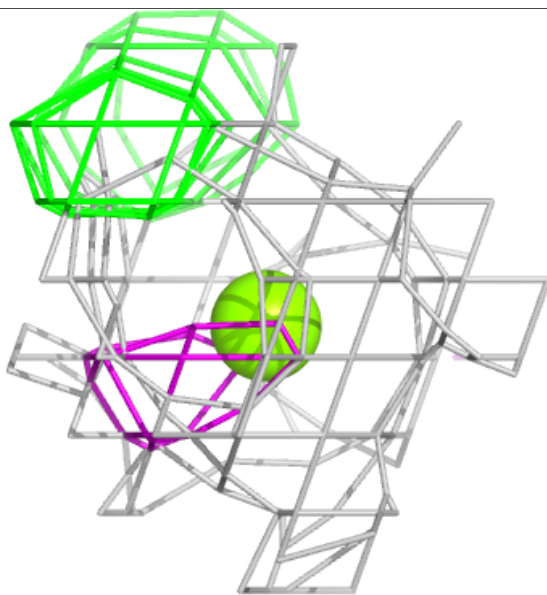
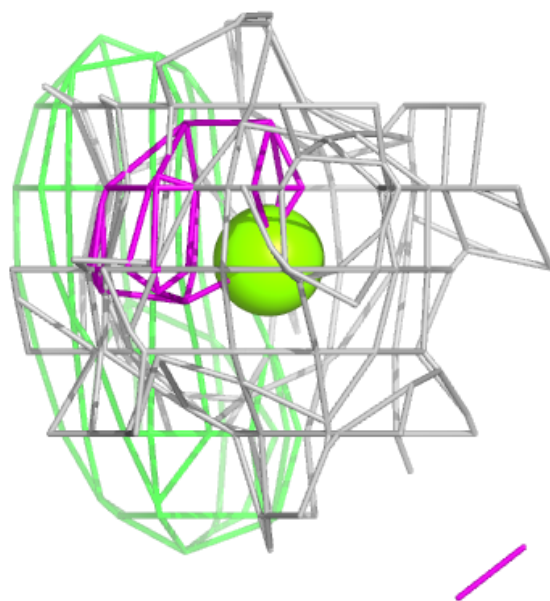
Electron density around MG B 408:

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and green (positive)



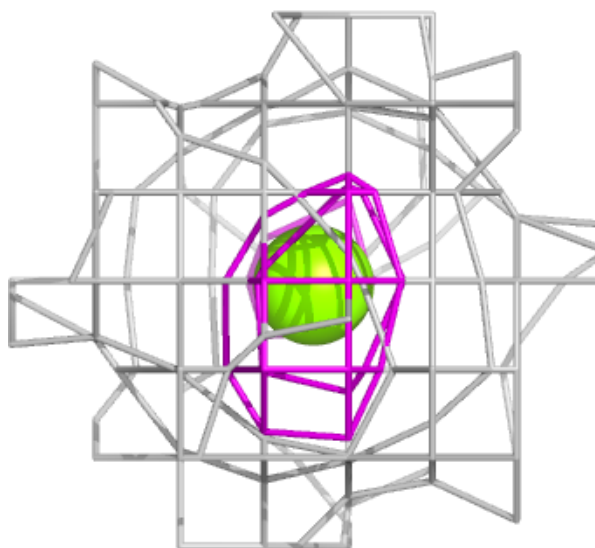
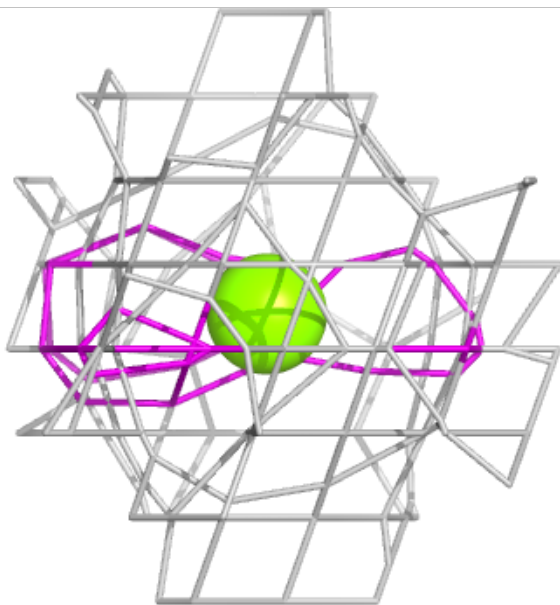
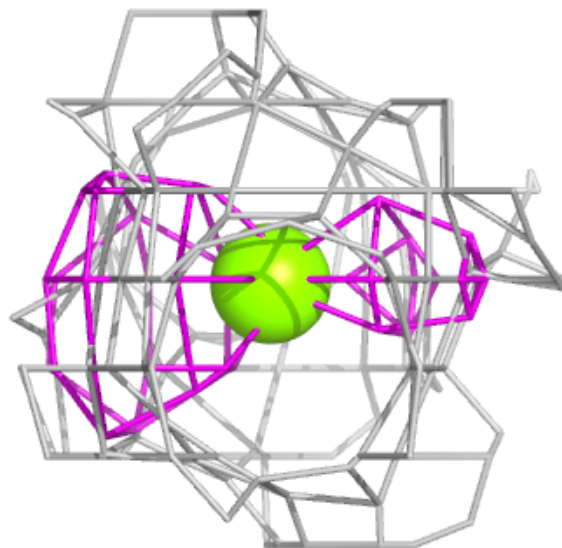
Electron density around MG A 407:

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and green (positive)



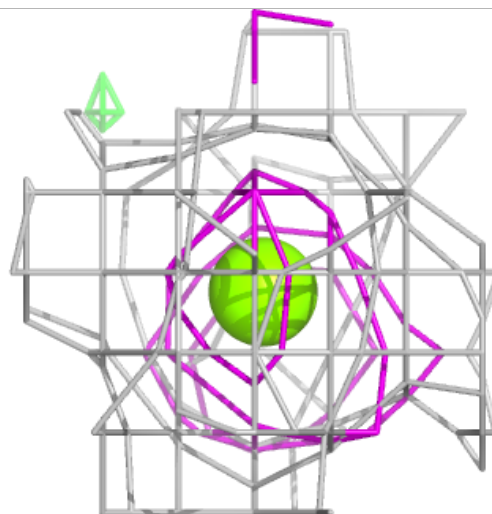
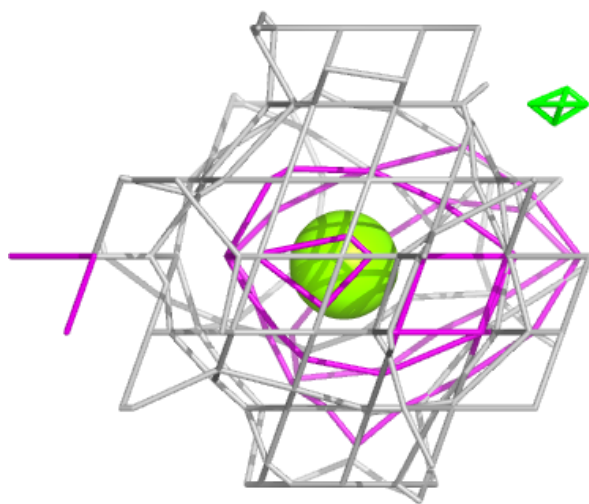
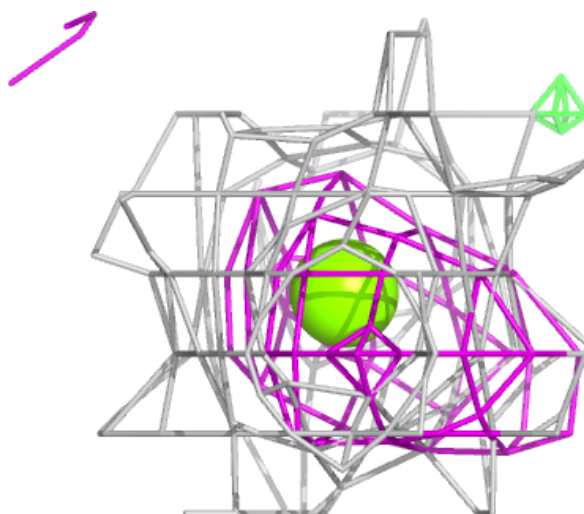
Electron density around MG B 407:

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and green (positive)



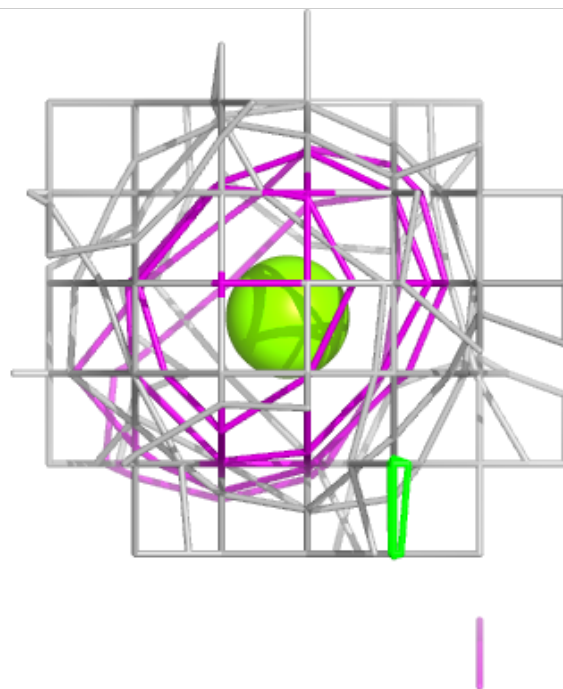
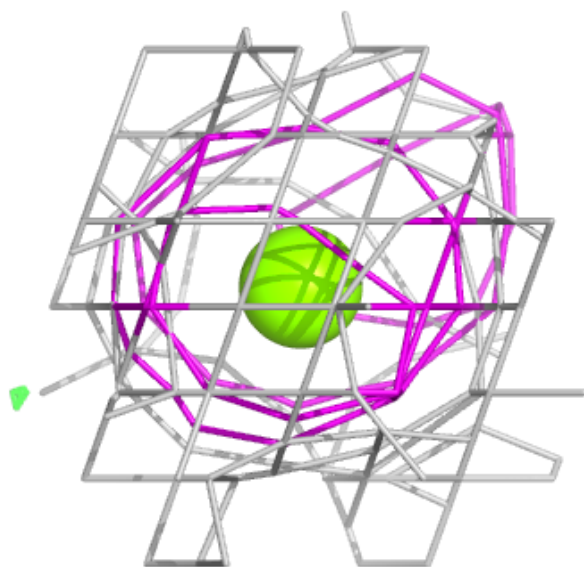
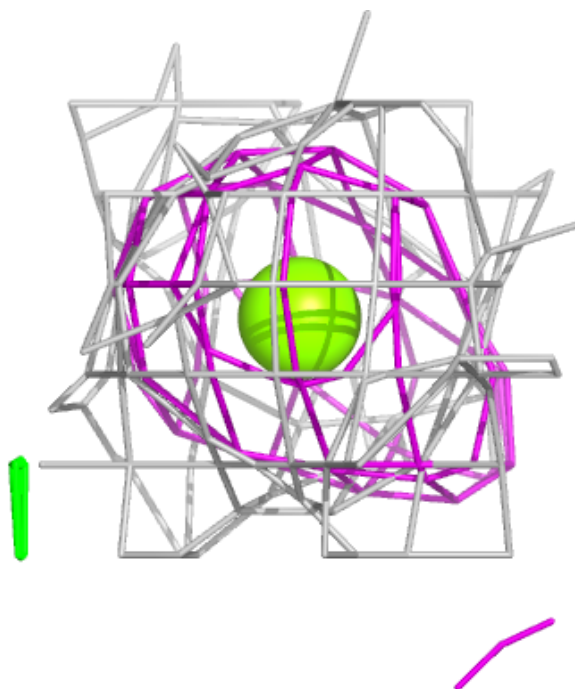
Electron density around MG A 406:

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and green (positive)



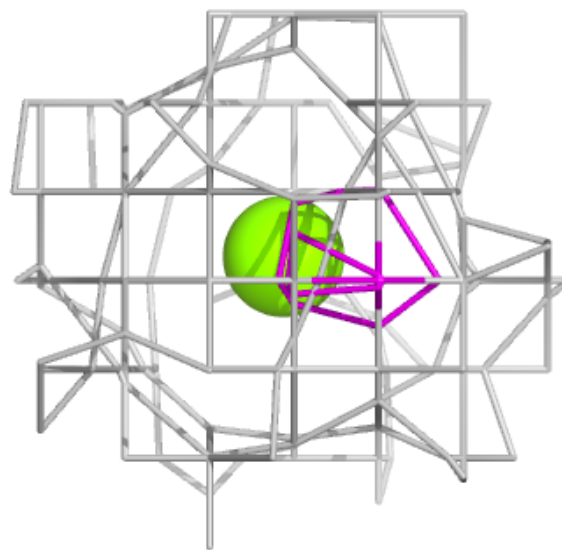
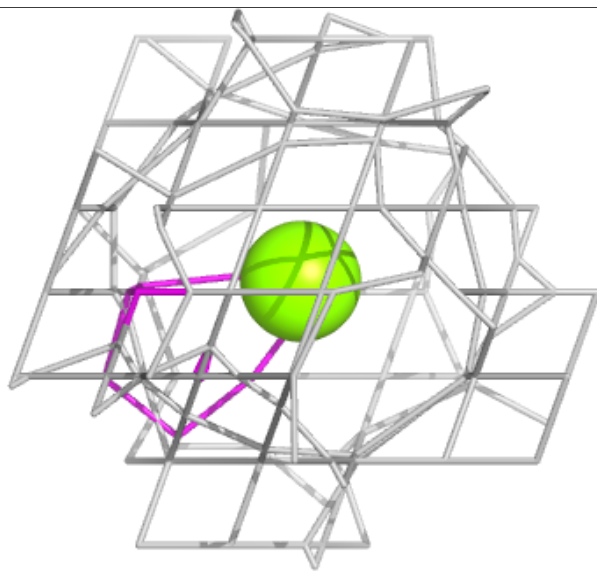
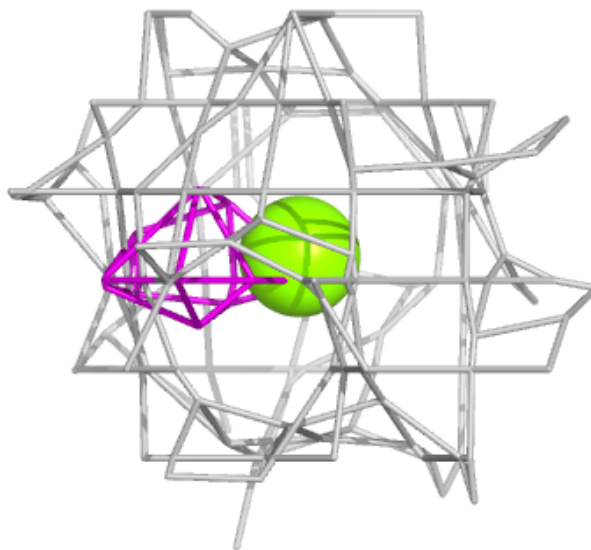
Electron density around MG B 406:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



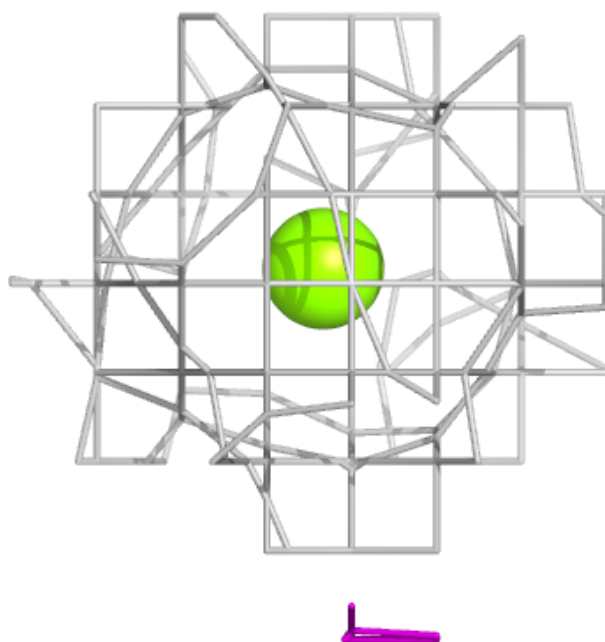
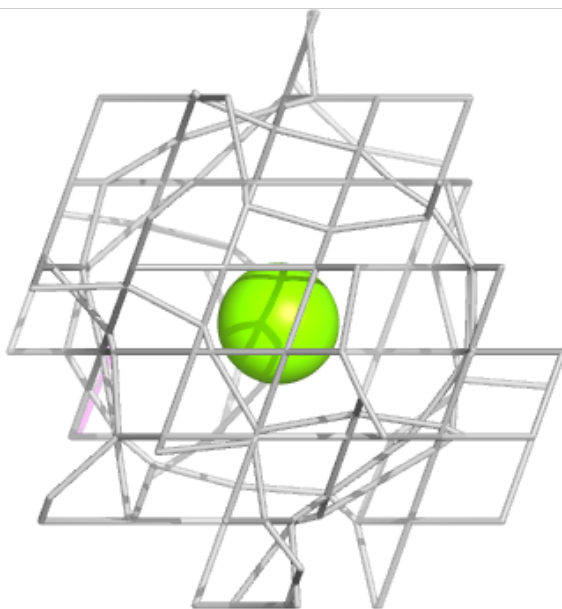
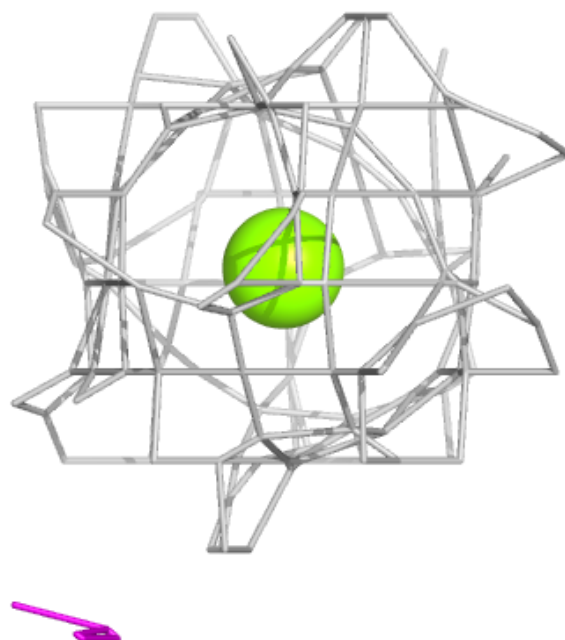
Electron density around MG B 405:

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



Electron density around MG A 405:

$2mF_o-DF_c$ (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.