



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:36 AM UTC

PDB ID : 9L50 / pdb_00009L50
Title : The ring expansion oxygenase SpoC in complex with Mn and N-oxalylglycine (NOG)
Authors : Liu, M.; He, X.
Deposited on : 2024-12-21
Resolution : 1.85 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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
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	:	9.0.010 (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

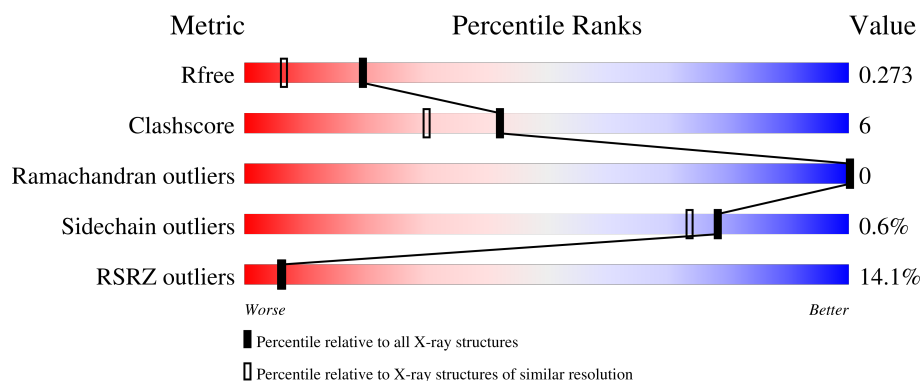
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	370	11% 73% 12% 15%
1	B	370	13% 76% 9% 15%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5592 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.

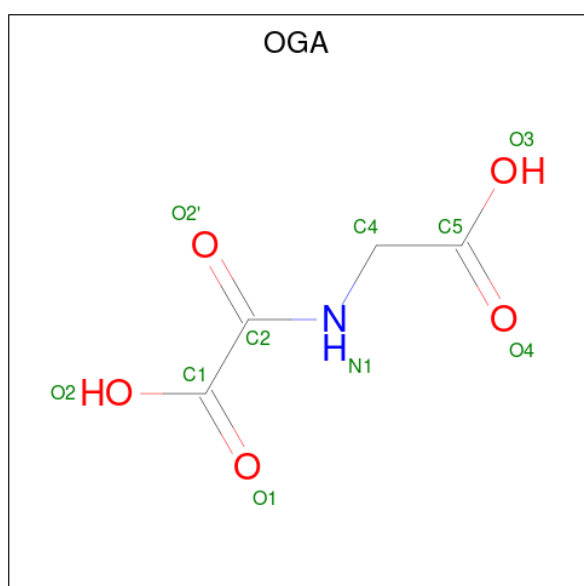
- Molecule 1 is a protein called ring expansion oxygenase SpoC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	0	4	0
			2506	1602	426	464	14			
1	B	314	Total	C	N	O	S	0	4	0
			2490	1592	420	464	14			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

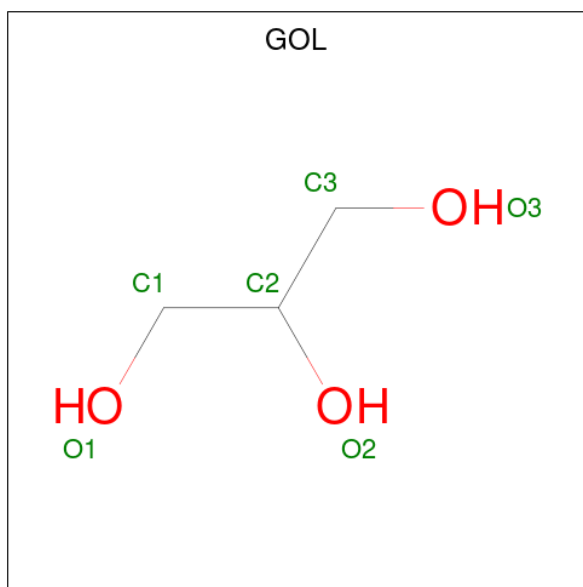
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		

- Molecule 3 is N-OXALYLGLYCINE (CCD ID: OGA) (formula: C₄H₅NO₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			10	4	1	5		
3	B	1	Total	C	N	O	0	0
			10	4	1	5		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).

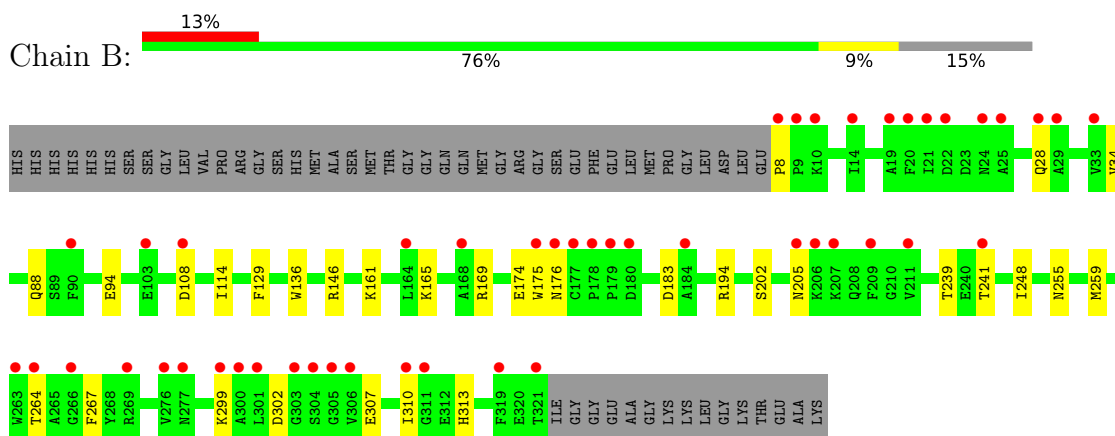


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	295	Total	O	0	0
			295	295		
5	B	261	Total	O	0	0
			261	261		

- Molecule 1: ring expansion oxygenase SpoC



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.41Å 84.30Å 104.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.07 – 1.85 39.07 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.8 (39.07-1.85) 99.9 (39.07-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 1.84Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.233 , 0.272 0.234 , 0.273	Depositor DCC
R_{free} test set	3068 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.386	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5592	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 38.54 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.6282e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: GOL, MN, OGA

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	0.32	0/2580	0.53	0/3501
1	B	0.33	0/2570	0.52	0/3490
All	All	0.32	0/5150	0.52	0/6991

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	2506	0	2467	36	0
1	B	2490	0	2448	26	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	10	0	3	1	0
3	B	10	0	3	1	0
4	A	12	0	16	1	0
4	B	6	0	8	0	0
5	A	295	0	0	15	0
5	B	261	0	0	8	0
All	All	5592	0	4945	63	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 (63) 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:11:THR:HG21	5:A:696:HOH:O	1.78	0.83
1:B:264:THR:HB	5:B:634:HOH:O	1.86	0.74
1:B:267:PHE:HB3	5:B:634:HOH:O	1.89	0.72
1:A:28:GLN:NE2	5:A:506:HOH:O	2.25	0.70
1:A:297:LYS:NZ	5:A:505:HOH:O	2.24	0.69
1:A:293:ASN:N	5:A:510:HOH:O	2.27	0.68
1:B:28:GLN:NE2	5:B:505:HOH:O	2.26	0.68
1:B:259:MET:HE3	1:B:310:ILE:HA	1.79	0.65
1:A:178:PRO:HB2	1:A:180:ASP:OD1	1.99	0.63
1:A:262:LYS:NZ	5:A:514:HOH:O	2.30	0.63
1:A:112:ALA:HB2	1:A:196:LEU:HD23	1.84	0.60
1:A:108[B]:ASP:OD1	1:A:202:SER:N	2.34	0.59
1:B:169:ARG:NH1	5:B:508:HOH:O	2.33	0.59
1:B:88:GLN:HG3	1:B:129:PHE:HA	1.84	0.59
1:A:247:VAL:HG23	5:A:520:HOH:O	2.02	0.58
1:B:299:LYS:HG2	1:B:307:GLU:HG2	1.84	0.58
1:A:94:GLU:OE1	1:A:194:ARG:NH2	2.36	0.58
1:A:310:ILE:HD12	5:A:510:HOH:O	2.04	0.57
1:A:229:THR:O	1:A:283:ARG:NH1	2.33	0.57
3:B:402:OGA:H4C1	5:B:509:HOH:O	2.06	0.55
1:B:108[A]:ASP:OD1	1:B:202:SER:N	2.39	0.55
1:B:175:TRP:CZ2	1:B:302:ASP:HB3	2.42	0.55
1:B:94[B]:GLU:OE1	1:B:194:ARG:NH2	2.40	0.55
1:A:141:PRO:HB2	1:A:144:GLU:HG3	1.91	0.53
1:B:259:MET:HE1	1:B:313:HIS:HB3	1.91	0.53
1:A:297:LYS:HE2	1:A:307:GLU:HB3	1.91	0.52
1:A:185:ALA:HA	5:A:510:HOH:O	2.09	0.52
1:B:176:ASN:ND2	5:B:517:HOH:O	2.42	0.52
1:B:114:ILE:HG13	1:B:194:ARG:HG2	1.92	0.50
1:A:211:VAL:HG22	5:A:562:HOH:O	2.13	0.49
1:B:255:ASN:OD1	5:B:501:HOH:O	2.20	0.49
1:A:204:GLU:HG2	1:A:208:GLN:HE21	1.79	0.48
1:B:34:VAL:HG21	1:B:169:ARG:HD2	1.95	0.48
1:A:18:SER:HB3	1:A:22:ASP:OD2	2.13	0.48
1:B:174:GLU:H	1:B:174:GLU:CD	2.20	0.48
1:A:60:GLN:OE1	5:A:502:HOH:O	2.20	0.47
1:A:226:GLU:O	1:A:247:VAL:HG21	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:MET:HE2	1:B:313:HIS:ND1	2.29	0.47
3:A:402:OGA:O2	4:A:403:GOL:H31	2.14	0.46
1:A:108[A]:ASP:OD1	5:A:501:HOH:O	2.20	0.46
1:A:90:PHE:HD1	1:A:114:ILE:HD13	1.80	0.46
1:B:259:MET:CE	1:B:310:ILE:HA	2.46	0.45
1:A:287:PRO:HB3	5:A:584:HOH:O	2.16	0.45
1:A:158:GLU:HG2	5:A:743:HOH:O	2.16	0.45
1:A:180:ASP:OD1	1:A:180:ASP:N	2.46	0.44
1:A:31:ASP:OD2	1:A:35:LYS:HE3	2.18	0.44
1:B:165:LYS:HE2	1:B:183:ASP:OD2	2.18	0.44
1:B:94[A]:GLU:OE1	1:B:194:ARG:NH2	2.51	0.43
1:A:204:GLU:HG2	1:A:208:GLN:NE2	2.33	0.43
1:B:161:LYS:NZ	5:B:522:HOH:O	2.51	0.43
1:A:211:VAL:HG23	1:A:275:VAL:HB	2.02	0.42
1:A:92:GLY:O	1:A:114:ILE:HG22	2.20	0.42
1:B:8:PRO:HA	1:B:239:THR:HG21	2.01	0.42
1:B:136:TRP:CZ2	1:B:146:ARG:HG3	2.55	0.41
1:B:239:THR:O	1:B:241:THR:HG23	2.20	0.41
1:A:11:THR:CG2	5:A:696:HOH:O	2.53	0.41
1:A:142:ASP:HB3	1:A:147:ILE:HG13	2.02	0.41
1:B:114:ILE:N	1:B:114:ILE:HD12	2.35	0.41
1:A:184:ALA:O	1:A:187:VAL:HG22	2.20	0.41
1:A:50:VAL:HG21	1:B:248:ILE:HG23	2.03	0.40
1:A:74:MET:HG2	5:A:556:HOH:O	2.21	0.40
1:A:209:PHE:CE2	1:A:274:ARG:HD2	2.57	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	318/370 (86%)	308 (97%)	10 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	316/370 (85%)	309 (98%)	7 (2%)	0	100	100
All	All	634/740 (86%)	617 (97%)	17 (3%)	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	275/313 (88%)	273 (99%)	2 (1%)	76	70
1	B	275/313 (88%)	274 (100%)	1 (0%)	84	82
All	All	550/626 (88%)	547 (100%)	3 (0%)	78	77

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	207	LYS
1	A	211	VAL
1	B	205	ASN

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	60	GLN
1	A	88	GLN
1	A	101	HIS
1	A	219	ASN
1	B	101	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](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	GOL	A	404	-	5,5,5	0.89	0	5,5,5	1.05	0
4	GOL	B	403	-	5,5,5	0.91	0	5,5,5	0.97	0
4	GOL	A	403	-	5,5,5	0.83	0	5,5,5	0.90	0
3	OGA	A	402	2	9,9,9	1.77	2 (22%)	8,11,11	2.03	1 (12%)
3	OGA	B	402	2	9,9,9	1.61	2 (22%)	8,11,11	2.15	3 (37%)

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	GOL	A	404	-	-	2/4/4/4	-
4	GOL	B	403	-	-	2/4/4/4	-
4	GOL	A	403	-	-	2/4/4/4	-
3	OGA	A	402	2	-	0/8/9/9	-
3	OGA	B	402	2	-	2/8/9/9	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	OGA	C2-C1	3.09	1.57	1.54
3	A	402	OGA	C4-C5	2.84	1.57	1.50
3	B	402	OGA	O2-C1	-2.50	1.23	1.30
3	A	402	OGA	C2-C1	2.40	1.57	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	OGA	C4-N1-C2	4.58	128.36	121.29
3	B	402	OGA	O2'-C2-C1	3.95	126.55	121.36
3	B	402	OGA	C4-N1-C2	2.92	125.80	121.29
3	B	402	OGA	C5-C4-N1	2.07	119.49	113.06

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	404	GOL	O1-C1-C2-O2
4	A	404	GOL	O1-C1-C2-C3
4	A	403	GOL	C1-C2-C3-O3
4	B	403	GOL	C1-C2-C3-O3
4	B	403	GOL	O2-C2-C3-O3
3	B	402	OGA	N1-C4-C5-O3
3	B	402	OGA	N1-C4-C5-O4
4	A	403	GOL	O2-C2-C3-O3

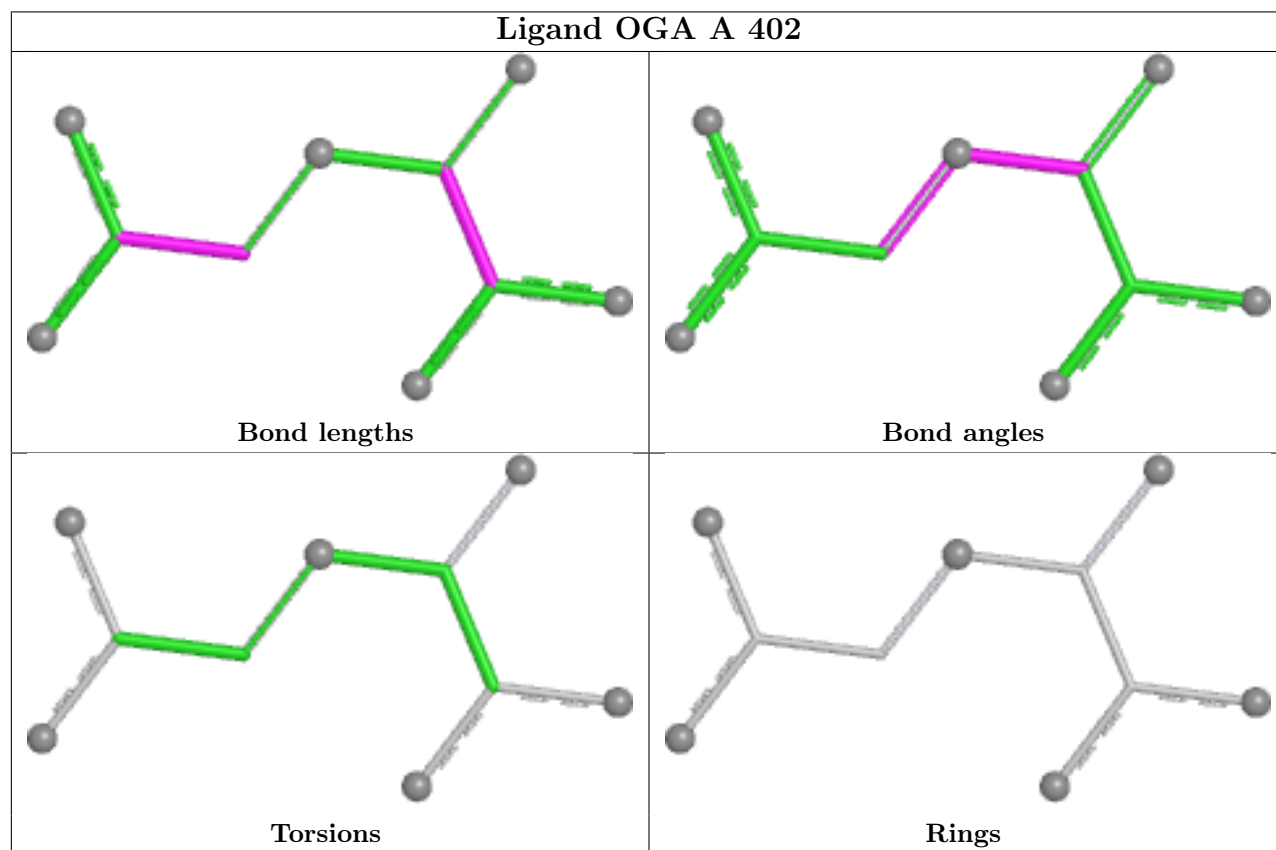
There are no ring outliers.

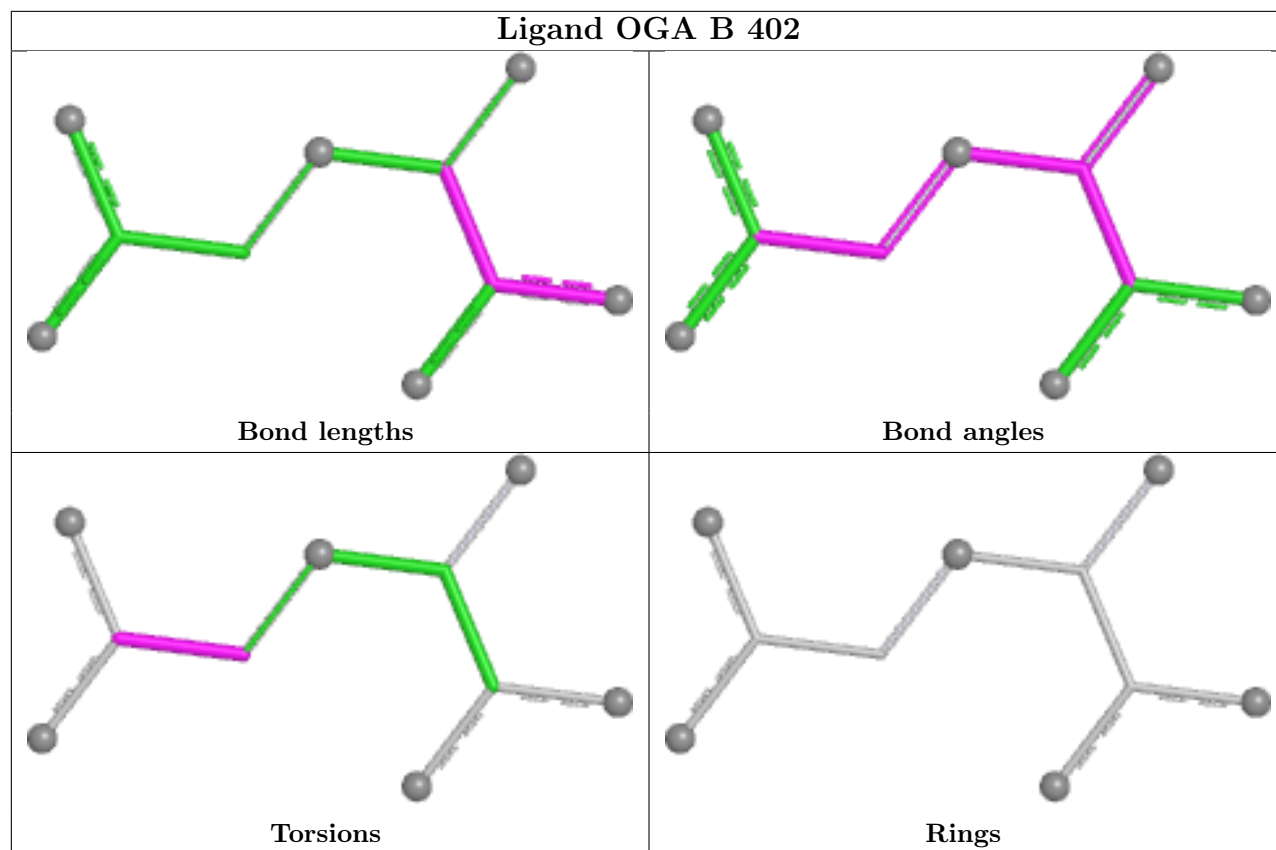
3 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	403	GOL	1	0
3	A	402	OGA	1	0
3	B	402	OGA	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	316/370 (85%)	0.96	41 (12%) 7 7	14, 31, 50, 73	4 (1%)
1	B	314/370 (84%)	1.09	48 (15%) 5 5	19, 31, 56, 72	4 (1%)
All	All	630/740 (85%)	1.03	89 (14%) 6 6	14, 31, 54, 73	8 (1%)

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	184	ALA	4.9
1	A	301	LEU	4.8
1	B	20	PHE	4.4
1	A	322	ILE	4.4
1	B	21	ILE	4.2
1	B	276	VAL	4.0
1	A	319	PHE	3.9
1	B	311	GLY	3.8
1	A	90	PHE	3.7
1	B	177	CYS	3.7
1	A	210	GLY	3.7
1	B	319	PHE	3.5
1	A	277[A]	ASN	3.5
1	B	303	GLY	3.5
1	B	300	ALA	3.4
1	B	28	GLN	3.4
1	B	211	VAL	3.3
1	B	306	VAL	3.3
1	A	205	ASN	3.2
1	B	266	GLY	3.2
1	B	321	THR	3.1
1	B	9	PRO	3.1
1	B	90	PHE	3.1
1	B	206	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	178	PRO	3.0
1	B	25	ALA	3.0
1	B	310	ILE	3.0
1	A	211	VAL	2.9
1	B	24	ASN	2.9
1	A	206	LYS	2.9
1	A	31	ASP	2.9
1	B	207	LYS	2.9
1	A	229	THR	2.8
1	A	82	VAL	2.8
1	B	209	PHE	2.8
1	A	323	GLY	2.8
1	A	119	VAL	2.8
1	B	175	TRP	2.8
1	A	23	ASP	2.7
1	A	294	ILE	2.7
1	A	276	VAL	2.7
1	B	205	ASN	2.7
1	A	27	ALA	2.6
1	B	305	GLY	2.6
1	A	321	THR	2.6
1	A	24	ASN	2.6
1	B	164	LEU	2.6
1	B	304	SER	2.6
1	B	22[A]	ASP	2.6
1	B	299	LYS	2.6
1	B	168	ALA	2.5
1	B	19	ALA	2.5
1	A	228	GLY	2.5
1	B	33	VAL	2.5
1	A	227	GLU	2.4
1	B	14	ILE	2.4
1	B	29	ALA	2.4
1	A	187	VAL	2.4
1	B	277	ASN	2.4
1	A	177	CYS	2.4
1	A	129	PHE	2.3
1	A	275	VAL	2.3
1	A	179	PRO	2.2
1	A	183	ASP	2.2
1	A	172	PRO	2.2
1	A	309	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	301	LEU	2.2
1	A	8	PRO	2.2
1	A	178	PRO	2.2
1	B	179	PRO	2.2
1	A	287	PRO	2.2
1	A	88	GLN	2.1
1	B	263	TRP	2.1
1	A	180	ASP	2.1
1	A	278	HIS	2.1
1	B	176	ASN	2.1
1	A	212[A]	GLY	2.1
1	B	269	ARG	2.1
1	B	10	LYS	2.1
1	B	180	ASP	2.1
1	B	103	GLU	2.1
1	B	8	PRO	2.1
1	A	207	LYS	2.0
1	B	264	THR	2.0
1	A	293	ASN	2.0
1	B	108[A]	ASP	2.0
1	A	139	SER	2.0
1	A	305	GLY	2.0
1	B	241	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

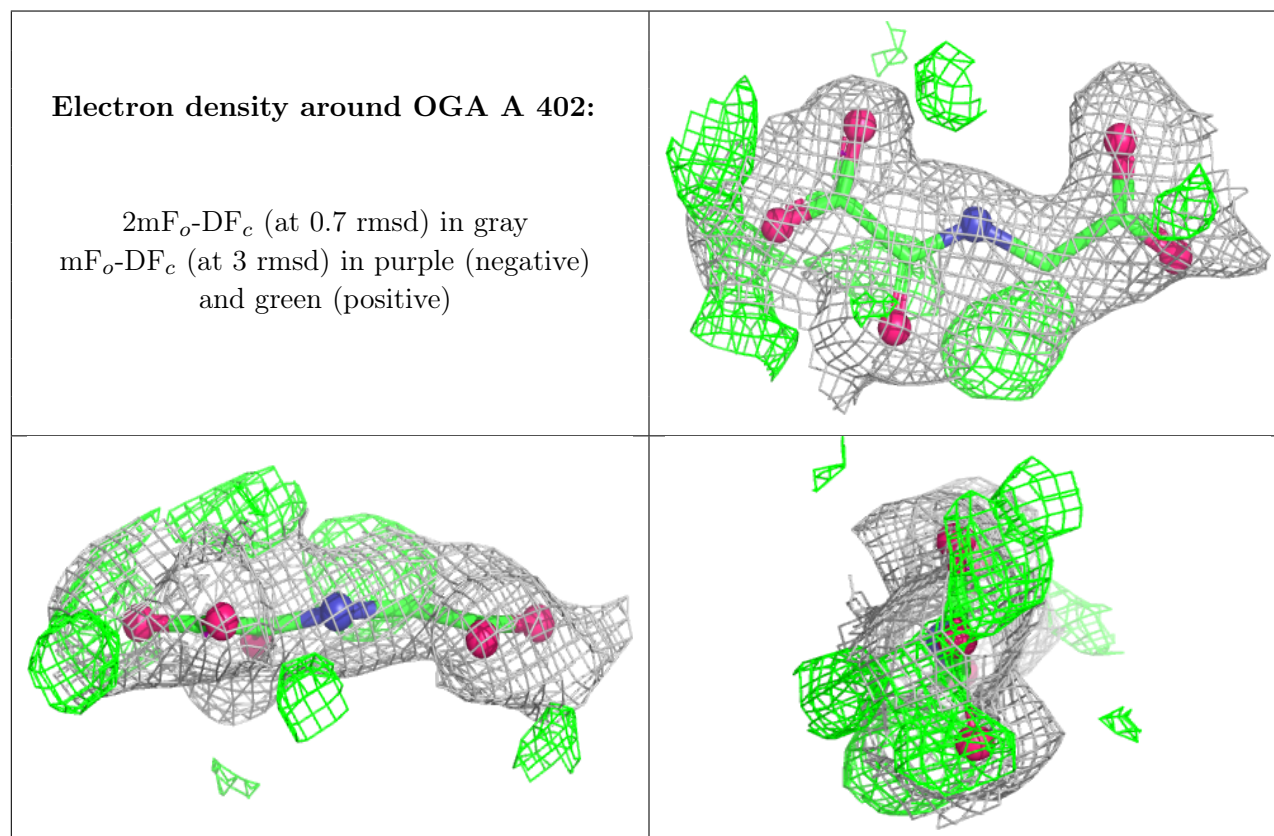
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

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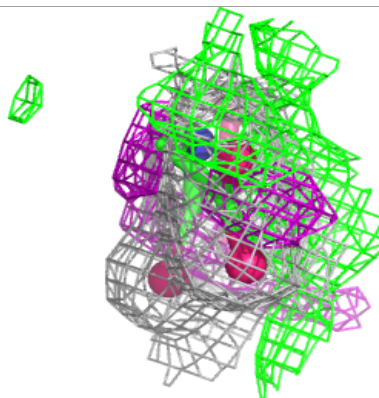
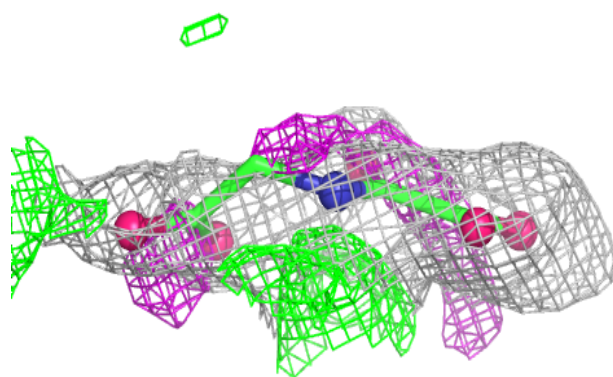
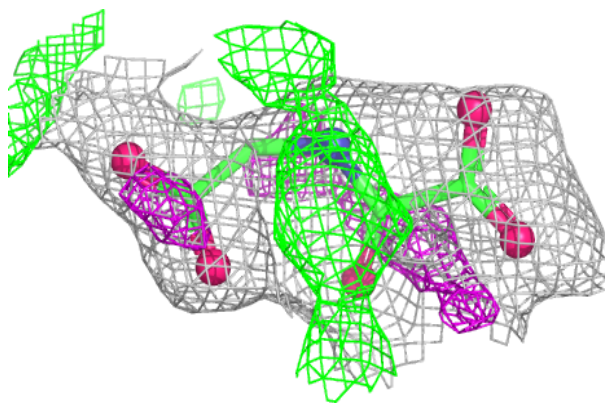
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	OGA	A	402	10/10	0.70	0.15	26,29,38,41	0
4	GOL	B	403	6/6	0.72	0.13	56,65,68,69	0
4	GOL	A	404	6/6	0.78	0.14	45,48,49,56	0
3	OGA	B	402	10/10	0.79	0.14	22,33,40,41	0
4	GOL	A	403	6/6	0.82	0.12	34,40,46,46	0
2	MN	B	401	1/1	0.96	0.04	25,25,25,25	0
2	MN	A	401	1/1	0.98	0.04	23,23,23,23	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.



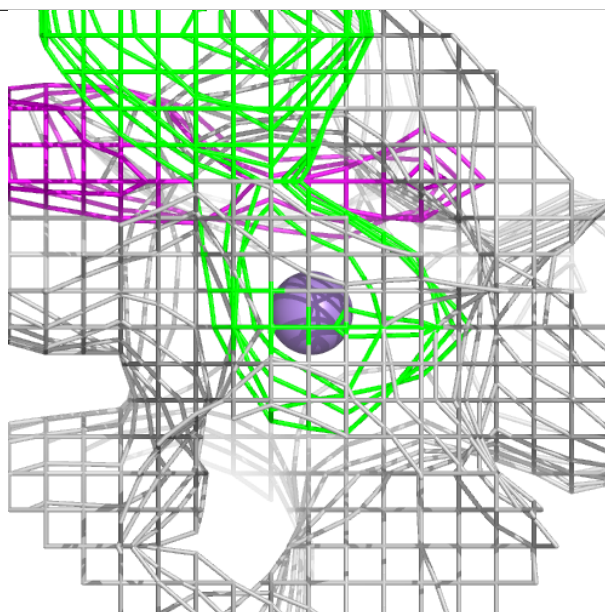
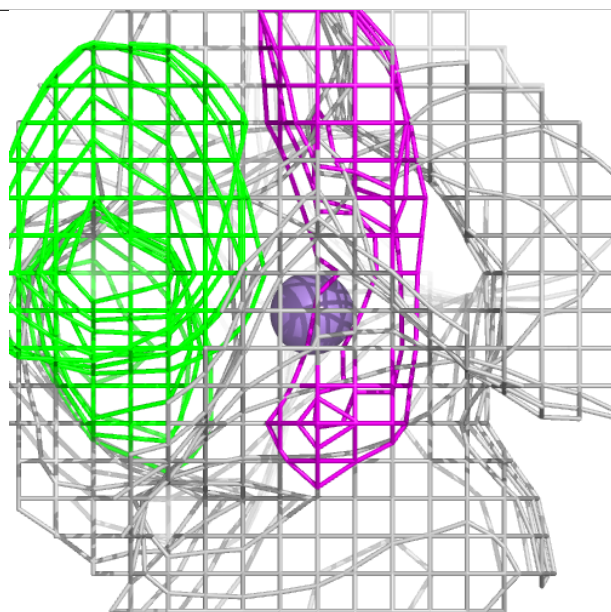
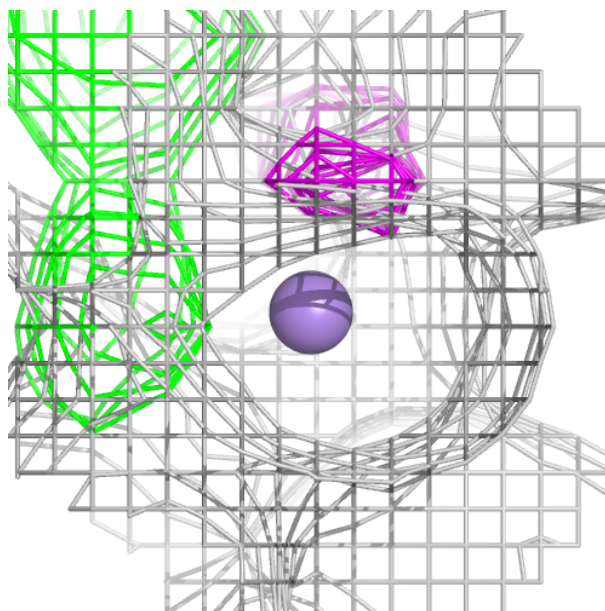
Electron density around OGA B 402:

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



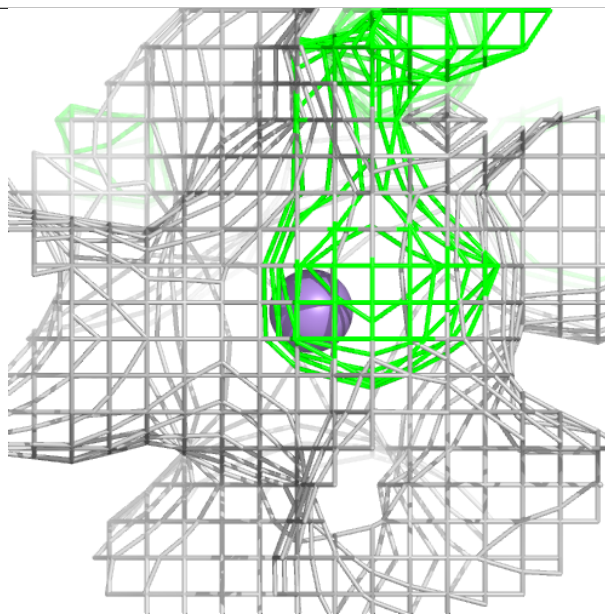
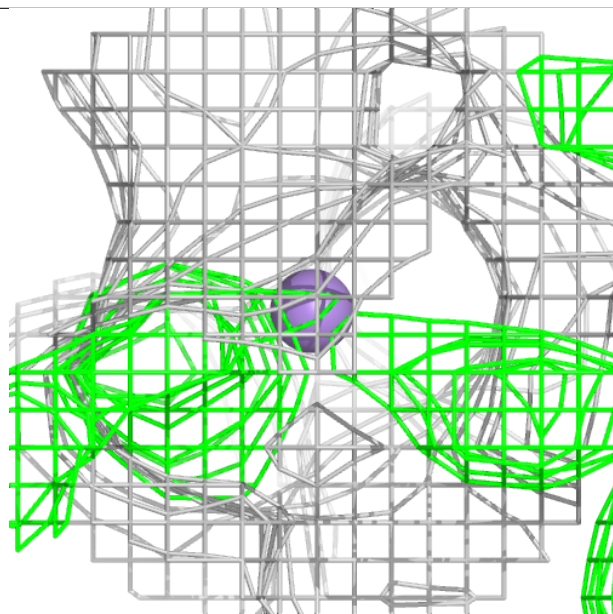
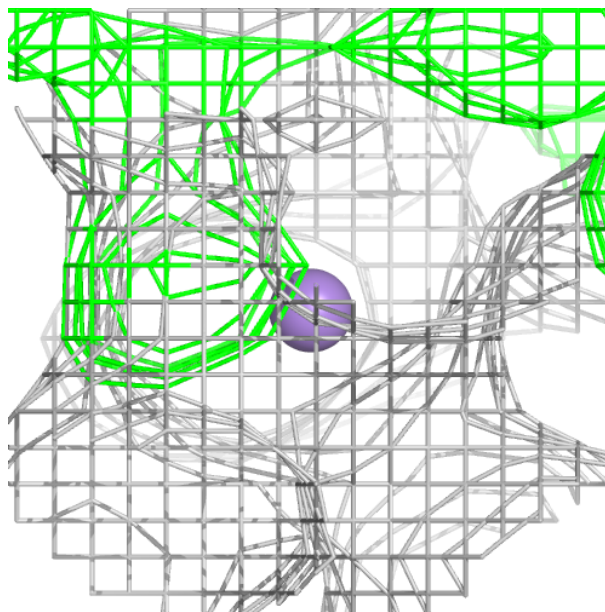
Electron density around MN B 401:

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



Electron density around MN A 401:

$2mF_o-DF_c$ (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.