



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 10:09 PM BST

PDB ID : 6OZN
Title : Crystal structure of Mus musculus (Mm) Endonuclease V in complex with a 23mer RNA oligo containing an inosine after a 15 min soak in 10 mM Mn²⁺
Authors : Samara, N.L.; Yang, W.
Deposited on : 2019-05-15
Resolution : 1.90 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
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

i

X-RAY DIFFRACTION

A.



R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

that have poor fit to the electron density. The numeric value is given above the bar.

Iteration	Category	Count	Confidence
1	A	253	<div> <div></div> <div>6%</div> <div>88%</div> <div>8%</div> <div></div> </div>
1	B	253	<div> <div></div> <div>20%</div> <div>86%</div> <div>10%</div> <div></div> </div>
2	C	23	<div> <div></div> <div>4%</div> <div>48%</div> <div>13%</div> <div></div> <div>35%</div> </div>
2	D	23	<div> <div></div> <div>4%</div> <div>48%</div> <div>13%</div> <div></div> <div>35%</div> </div>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 4798 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 Endonuclease V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	S	0	3	0
			1916	1215	346	347	8			
1	B	245	Total	C	N	O	S	0	5	0
			1938	1232	352	346	8			

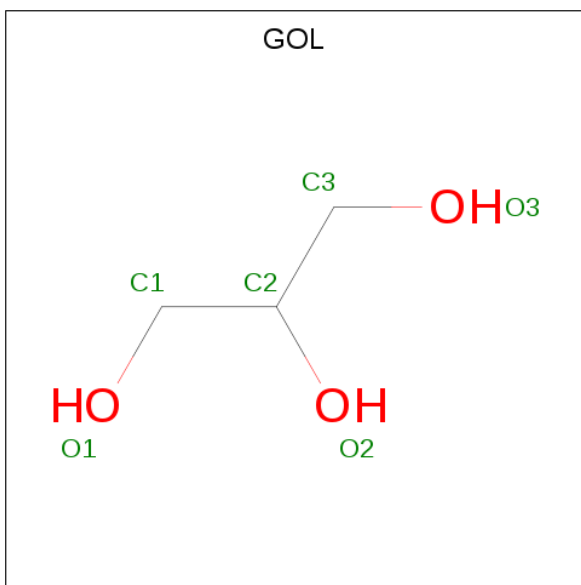
- Molecule 2 is DNA/RNA hybrid called DNA/RNA (5'-R(P*CP*GP*GP*UP*AP*AP*CP*CP*C)-D(P*I)-R(P*AP*UP*AP*UP*GP*CP*AP*UP*GP*CP*AP*UP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	15	Total	C	N	O	P	0	2	0
			328	147	53	112	16			
2	D	15	Total	C	N	O	P	0	2	0
			328	147	53	112	16			

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by author).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Mn	0	0
			3	3		
3	A	3	Total	Mn	0	0
			3	3		
3	D	3	Total	Mn	0	0
			3	3		
3	C	1	Total	Mn	0	0
			1	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 5 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



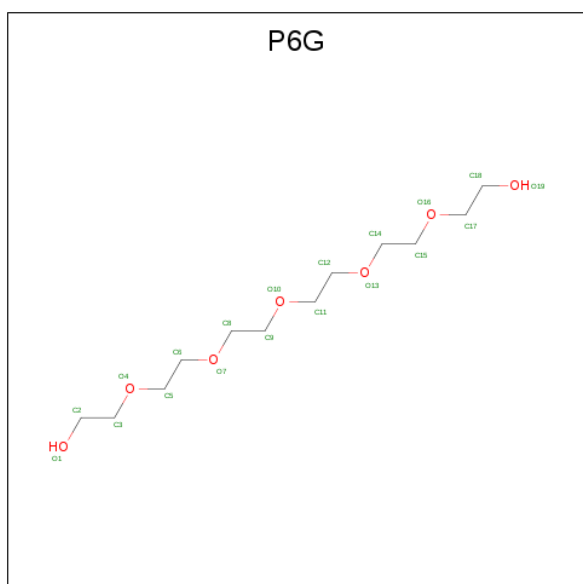
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		
5	B	1	Total	C	O	0	0
			10	6	4		

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



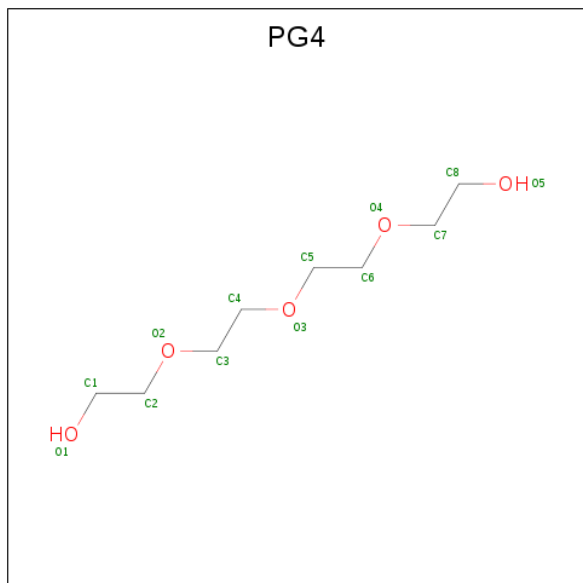
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: $C_{12}H_{26}O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			19	12	7		

- Molecule 8 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			13	8	5		
8	A	1	Total	C	O	0	0
			13	8	5		

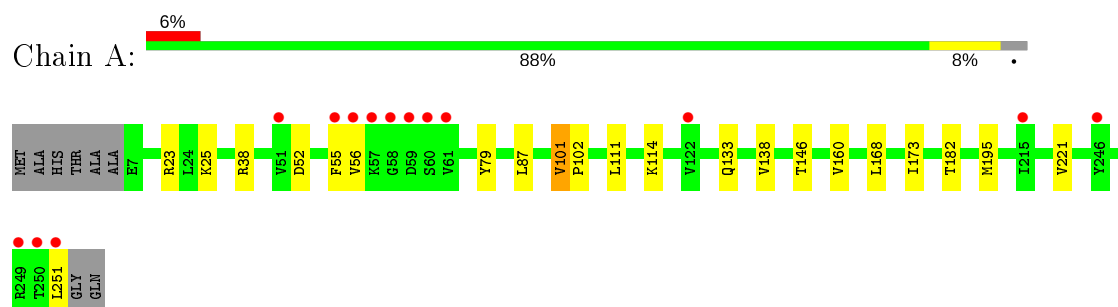
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	93	Total	O	0	0
			93	93		
9	B	60	Total	O	0	0
			60	60		
9	C	15	Total	O	0	0
			15	15		
9	D	17	Total	O	0	0
			17	17		

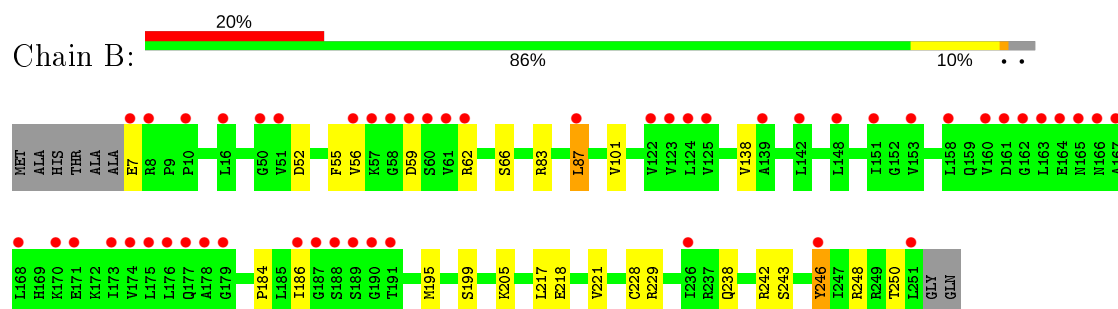
3 Residue-property plots [i](#)

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.

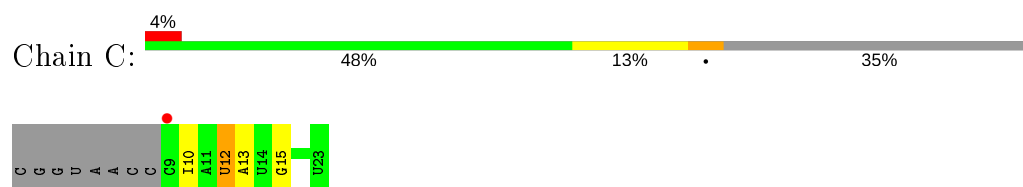
• Molecule 1: Endonuclease V



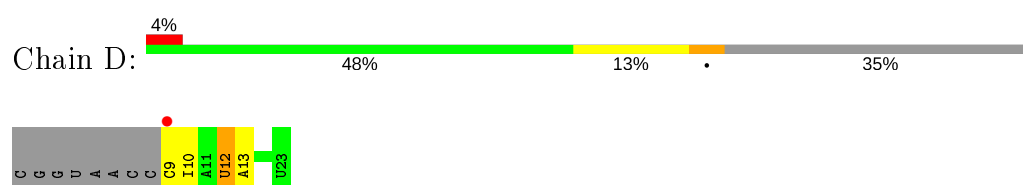
• Molecule 1: Endonuclease V



• Molecule 2: DNA/RNA (5'-R(P*CP*GP*GP*UP*AP*AP*CP*CP*C)-D(P*I)-R(P*AP*UP*A P*UP*GP*CP*AP*UP*GP*CP*AP*UP*U)-3')



• Molecule 2: DNA/RNA (5'-R(P*CP*GP*GP*UP*AP*AP*CP*CP*C)-D(P*I)-R(P*AP*UP*A P*UP*GP*CP*AP*UP*GP*CP*AP*UP*U)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.40Å 72.94Å 155.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.50 – 1.90 41.92 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.5 (35.50-1.90) 97.7 (41.92-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.17 (at 1.89Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.192 , 0.221 0.192 , 0.222	Depositor DCC
R_{free} test set	3189 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	29.5	Xtriage
Anisotropy	0.797	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 56.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for k,h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4798	wwPDB-VP
Average B, all atoms (Å ²)	52.0	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 25.88 % 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 2.9202e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PGE, MN, EDO, PG4, P6G

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	0.42	0/1962	0.60	2/2659 (0.1%)
1	B	0.37	0/1991	0.57	2/2694 (0.1%)
2	C	5.32	2/350 (0.6%)	0.88	0/536
2	D	5.23	2/350 (0.6%)	1.07	3/536 (0.6%)
All	All	2.08	4/4653 (0.1%)	0.67	7/6425 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	12[B]	U	OP3-P	70.04	2.45	1.61
2	C	12[A]	U	OP3-P	70.04	2.45	1.61
2	D	12[B]	U	OP3-P	68.89	2.43	1.61
2	D	12[A]	U	OP3-P	68.89	2.43	1.61

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	52	ASP	CB-CG-OD1	7.38	124.94	118.30
2	D	9	C	C2-N1-C1'	6.93	126.42	118.80
1	A	52	ASP	CB-CG-OD1	5.92	123.62	118.30
1	A	52	ASP	CB-CG-OD2	-5.63	113.23	118.30
1	B	52	ASP	CB-CG-OD2	-5.60	113.26	118.30
2	D	9	C	N1-C2-O2	5.23	122.04	118.90
2	D	9	C	C6-N1-C1'	-5.12	114.66	120.80

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	1916	0	1954	12	0
1	B	1938	0	1992	15	0
2	C	328	0	164	5	0
2	D	328	0	164	4	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	1	0	0	0	0
3	D	3	0	0	0	0
4	A	6	0	8	1	0
4	C	6	0	8	0	0
5	A	10	0	14	0	0
5	B	10	0	14	0	0
6	A	16	0	24	1	0
7	A	19	0	26	2	0
8	A	26	0	36	1	0
9	A	93	0	0	1	0
9	B	60	0	0	0	0
9	C	15	0	0	1	0
9	D	17	0	0	0	0
All	All	4798	0	4404	32	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 4.

All (32) 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:23:ARG:NH1	9:A:401:HOH:O	2.20	0.74
1:A:173:ILE:HG22	6:A:309:EDO:H21	1.79	0.65
1:A:195:MET:HG3	1:A:221:VAL:HG21	1.78	0.64
1:B:199:SER:HB2	1:B:228:CYS:SG	2.41	0.61
1:A:138:VAL:H	2:C:10:DI:H1	1.52	0.57
1:B:248[A]:ARG:CZ	2:C:15:G:H5"	2.35	0.57
1:B:195:MET:HG3	1:B:221:VAL:HG21	1.86	0.57
1:B:56:VAL:HB	1:B:59:ASP:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:55[A]:PHE:CD2	2:D:12[A]:U:H4'	2.41	0.56
1:B:195:MET:HG2	1:B:217:LEU:HD21	1.88	0.56
1:A:114:LYS:HE2	4:A:304:GOL:H11	1.89	0.55
1:A:133:GLN:HA	1:A:160:VAL:HG21	1.89	0.54
1:B:238:GLN:OE1	1:B:242[A]:ARG:NH1	2.43	0.52
1:B:184:PRO:HB2	1:B:186:ILE:HD11	1.92	0.52
1:B:138:VAL:H	2:D:10:DI:H1	1.58	0.51
1:A:182:THR:HG21	1:A:221:VAL:HG11	1.95	0.48
2:C:12[B]:U:OP3	9:C:201:HOH:O	0.51	0.47
1:A:25:LYS:HE2	8:A:312:PG4:H12	1.97	0.46
1:B:55[A]:PHE:HD1	1:B:87:LEU:HD21	1.80	0.46
1:B:55[A]:PHE:HD2	2:D:12[A]:U:H4'	1.79	0.46
7:A:310:P6G:H111	7:A:310:P6G:H82	1.60	0.45
1:B:205:LYS:HE2	1:B:205:LYS:HB3	1.75	0.44
2:C:12[B]:U:H2'	2:C:13:A:C8	2.52	0.44
2:C:12[A]:U:H2'	2:C:13:A:C8	2.53	0.43
1:B:218:GLU:H	1:B:218:GLU:CD	2.22	0.43
1:B:246:TYR:HE1	1:B:250:THR:HG21	1.85	0.42
1:A:133:GLN:HA	1:A:160:VAL:CG2	2.50	0.42
2:D:12[B]:U:H2'	2:D:13:A:C8	2.55	0.42
1:A:79:TYR:CZ	1:A:111:LEU:HD13	2.55	0.41
1:A:101:VAL:HG12	1:A:102:PRO:HD3	2.02	0.41
1:B:66[A]:SER:OG	1:B:243:SER:HB3	2.20	0.41
1:A:146:THR:O	7:A:310:P6G:H61	2.20	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	246/253 (97%)	237 (96%)	9 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	248/253 (98%)	241 (97%)	7 (3%)	0	100	100
All	All	494/506 (98%)	478 (97%)	16 (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	211/215 (98%)	204 (97%)	7 (3%)	38	29
1	B	214/215 (100%)	207 (97%)	7 (3%)	38	29
All	All	425/430 (99%)	411 (97%)	14 (3%)	37	29

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	ARG
1	A	55	PHE
1	A	56	VAL
1	A	87	LEU
1	A	101	VAL
1	A	168	LEU
1	A	251	LEU
1	B	7	GLU
1	B	62	ARG
1	B	83	ARG
1	B	87	LEU
1	B	101	VAL
1	B	229	ARG
1	B	246	TYR

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

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 21 ligands modelled in this entry, 10 are monoatomic - leaving 11 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	$\# Z > 2$
6	EDO	A	306	-	3,3,3	0.45	0	2,2,2	0.33	0
5	PGE	B	304	-	9,9,9	0.32	0	8,8,8	0.31	0
6	EDO	A	307	-	3,3,3	0.42	0	2,2,2	0.39	0
6	EDO	A	309	-	3,3,3	0.45	0	2,2,2	0.24	0
4	GOL	C	102	-	5,5,5	0.36	0	5,5,5	0.51	0
8	PG4	A	311	-	12,12,12	0.52	0	11,11,11	0.25	0
7	P6G	A	310	-	18,18,18	0.53	0	17,17,17	0.29	0
5	PGE	A	305	-	9,9,9	0.32	0	8,8,8	0.24	0
8	PG4	A	312	-	12,12,12	0.51	0	11,11,11	0.30	0
6	EDO	A	308	-	3,3,3	0.50	0	2,2,2	0.31	0
4	GOL	A	304	-	5,5,5	0.41	0	5,5,5	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
6	EDO	A	306	-	-	1/1/1/1	-
5	PGE	B	304	-	-	3/7/7/7	-
6	EDO	A	307	-	-	0/1/1/1	-
6	EDO	A	309	-	-	0/1/1/1	-
4	GOL	C	102	-	-	0/4/4/4	-
8	PG4	A	311	-	-	6/10/10/10	-
7	P6G	A	310	-	-	10/16/16/16	-
5	PGE	A	305	-	-	4/7/7/7	-
8	PG4	A	312	-	-	6/10/10/10	-
6	EDO	A	308	-	-	1/1/1/1	-
4	GOL	A	304	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	304	GOL	C1-C2-C3-O3
7	A	310	P6G	O4-C5-C6-O7
4	A	304	GOL	O2-C2-C3-O3
5	A	305	PGE	O3-C5-C6-O4
8	A	311	PG4	O4-C7-C8-O5
7	A	310	P6G	O13-C14-C15-O16
4	A	304	GOL	O1-C1-C2-C3
4	A	304	GOL	O1-C1-C2-O2
7	A	310	P6G	C8-C9-O10-C11
7	A	310	P6G	O16-C17-C18-O19
6	A	306	EDO	O1-C1-C2-O2
5	B	304	PGE	C1-C2-O2-C3
7	A	310	P6G	C5-C6-O7-C8
5	A	305	PGE	O1-C1-C2-O2
8	A	312	PG4	C6-C5-O3-C4
8	A	311	PG4	C3-C4-O3-C5
8	A	312	PG4	C1-C2-O2-C3
5	A	305	PGE	C1-C2-O2-C3
7	A	310	P6G	C2-C3-O4-C5
5	A	305	PGE	O2-C3-C4-O3
8	A	311	PG4	C4-C3-O2-C2
5	B	304	PGE	C6-C5-O3-C4

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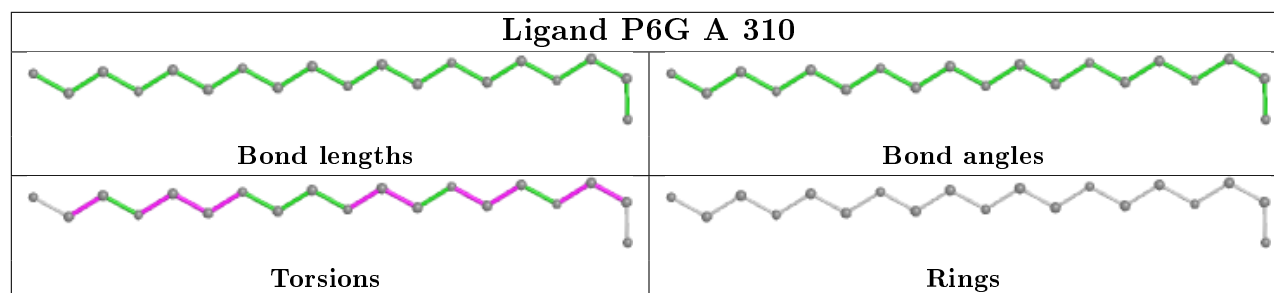
Mol	Chain	Res	Type	Atoms
7	A	310	P6G	C14-C15-O16-C17
8	A	311	PG4	C5-C6-O4-C7
8	A	312	PG4	O4-C7-C8-O5
8	A	312	PG4	C5-C6-O4-C7
8	A	312	PG4	O1-C1-C2-O2
6	A	308	EDO	O1-C1-C2-O2
8	A	312	PG4	O3-C5-C6-O4
8	A	311	PG4	C1-C2-O2-C3
7	A	310	P6G	C15-C14-O13-C12
5	B	304	PGE	O2-C3-C4-O3
8	A	311	PG4	O3-C5-C6-O4
7	A	310	P6G	O1-C2-C3-O4
7	A	310	P6G	O7-C8-C9-O10

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	309	EDO	1	0
7	A	310	P6G	2	0
8	A	312	PG4	1	0
4	A	304	GOL	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

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 ⓘ

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	245/253 (96%)	0.39	14 (5%) 23 26	26, 39, 77, 130	0
1	B	245/253 (96%)	0.91	51 (20%) 1 1	31, 51, 91, 124	0
2	C	14/23 (60%)	-0.04	1 (7%) 16 17	39, 53, 89, 109	0
2	D	14/23 (60%)	0.03	1 (7%) 16 17	39, 58, 97, 146	0
All	All	518/552 (93%)	0.62	67 (12%) 3 3	26, 44, 90, 146	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	58	GLY	10.8
1	B	61	VAL	8.5
1	B	58	GLY	6.7
1	A	250	THR	6.5
1	B	59	ASP	6.4
1	B	62	ARG	5.5
1	B	60	SER	5.4
1	A	56	VAL	5.4
1	B	7	GLU	5.3
1	B	167	ALA	5.2
1	A	251	LEU	4.7
1	A	249	ARG	4.6
1	B	56	VAL	4.5
1	B	177	GLN	4.5
1	B	175	LEU	4.4
1	B	174	VAL	4.2
1	B	189	SER	4.2
2	D	9	C	4.1
1	B	188	SER	4.0
1	B	168	LEU	4.0
1	B	178	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	190	GLY	3.7
1	A	57	LYS	3.7
1	B	164	GLU	3.5
1	B	161	ASP	3.3
1	B	191	THR	3.3
1	A	59	ASP	3.3
1	A	60	SER	3.2
1	B	16	LEU	3.1
1	B	148	LEU	3.1
1	B	51	VAL	3.1
1	B	8	ARG	3.0
1	B	186	ILE	3.0
1	B	163	LEU	3.0
1	B	187	GLY	2.9
1	B	125	VAL	2.9
1	B	151	ILE	2.8
1	A	55	PHE	2.7
1	B	165	ASN	2.7
1	A	61	VAL	2.6
1	B	87	LEU	2.6
1	B	122	VAL	2.5
1	B	57	LYS	2.5
1	B	171	GLU	2.5
1	B	124	LEU	2.5
2	C	9	C	2.5
1	B	160	VAL	2.4
1	A	51	VAL	2.4
1	B	50	GLY	2.3
1	B	123	VAL	2.3
1	B	162	GLY	2.3
1	B	246	TYR	2.3
1	A	246	TYR	2.3
1	B	236	ILE	2.3
1	B	158	LEU	2.3
1	B	173	ILE	2.2
1	B	251	LEU	2.2
1	B	176	LEU	2.2
1	B	10	PRO	2.1
1	B	166	ASN	2.1
1	B	179	GLY	2.1
1	B	170	LYS	2.1
1	A	122	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	153	VAL	2.1
1	B	142	LEU	2.1
1	B	139	ALA	2.0
1	A	215	ILE	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
3	MN	B	303	1/1	0.69	0.38	51,51,51,51	1
6	EDO	A	306	4/4	0.75	0.28	81,82,85,90	0
6	EDO	A	308	4/4	0.75	0.18	55,71,79,81	0
3	MN	D	102	1/1	0.79	0.16	38,38,38,38	1
8	PG4	A	312	13/13	0.80	0.22	69,78,91,92	0
4	GOL	C	102	6/6	0.80	0.23	51,53,61,67	0
5	PGE	B	304	10/10	0.83	0.15	73,81,89,90	0
6	EDO	A	307	4/4	0.86	0.11	73,74,79,85	0
6	EDO	A	309	4/4	0.86	0.50	80,81,81,86	0
4	GOL	A	304	6/6	0.87	0.22	72,73,75,79	0
8	PG4	A	311	13/13	0.87	0.21	59,67,74,74	0
5	PGE	A	305	10/10	0.88	0.12	63,69,78,82	0
3	MN	B	302	1/1	0.90	0.06	52,52,52,52	1
3	MN	D	103	1/1	0.92	0.22	50,50,50,50	1
3	MN	C	101	1/1	0.92	0.07	39,39,39,39	1
7	P6G	A	310	19/19	0.92	0.19	51,67,86,87	0
3	MN	A	303	1/1	0.93	0.06	52,52,52,52	1
3	MN	D	101	1/1	0.97	0.07	37,37,37,37	0
3	MN	A	302	1/1	0.98	0.08	30,30,30,30	0

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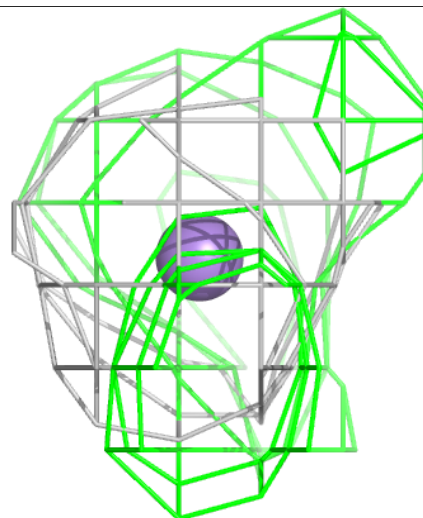
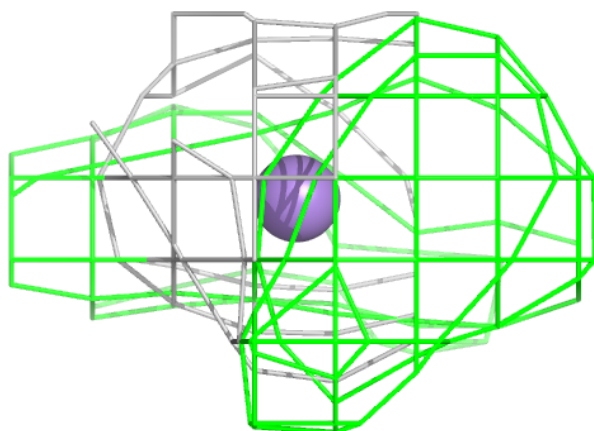
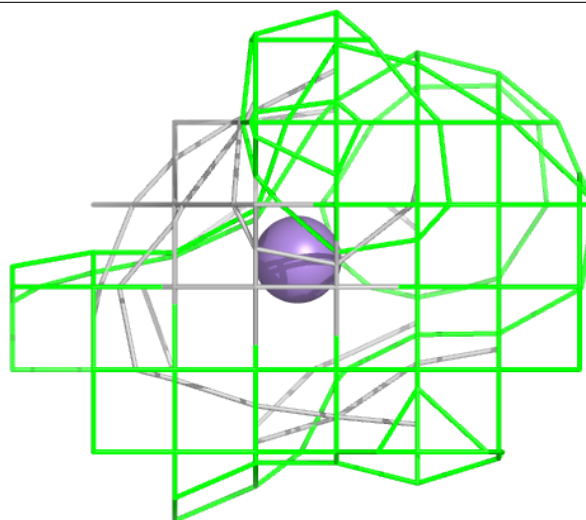
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MN	B	301	1/1	0.98	0.06	34,34,34,34	1
3	MN	A	301	1/1	0.99	0.09	36,36,36,36	0

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.

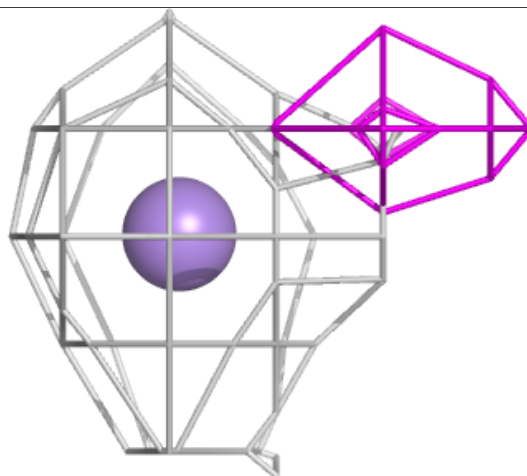
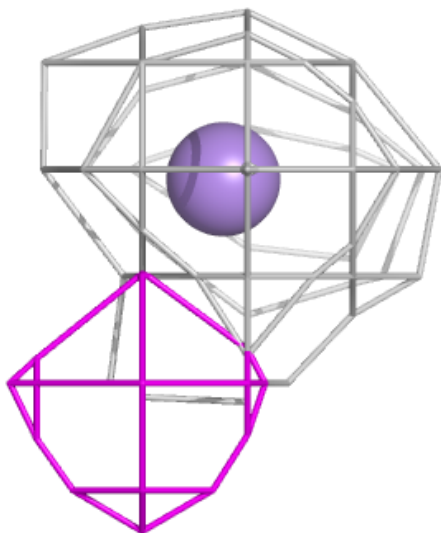
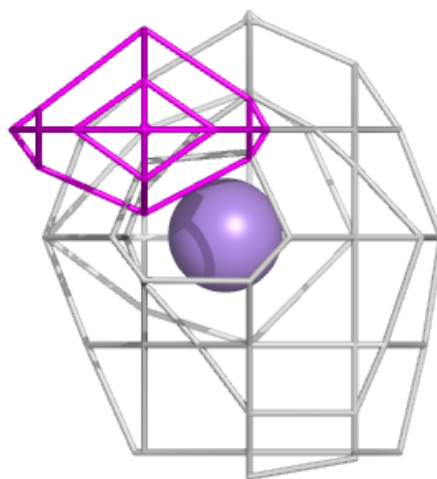
Electron density around MN B 303:

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



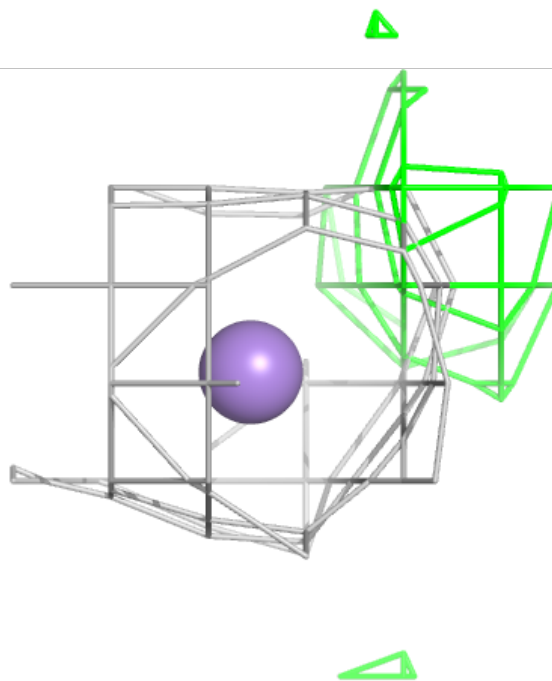
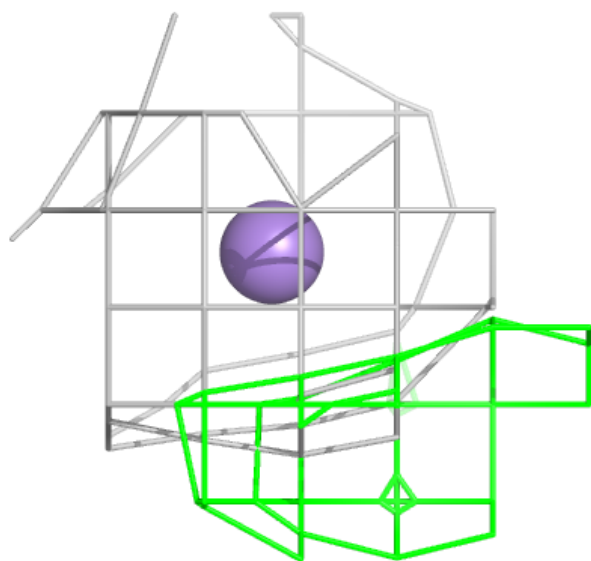
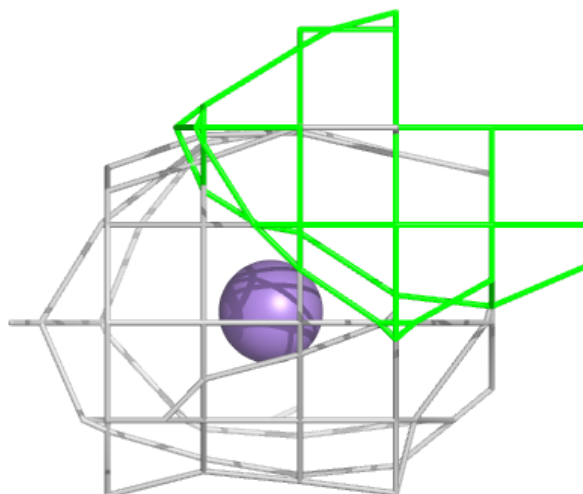
Electron density around MN D 102:

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



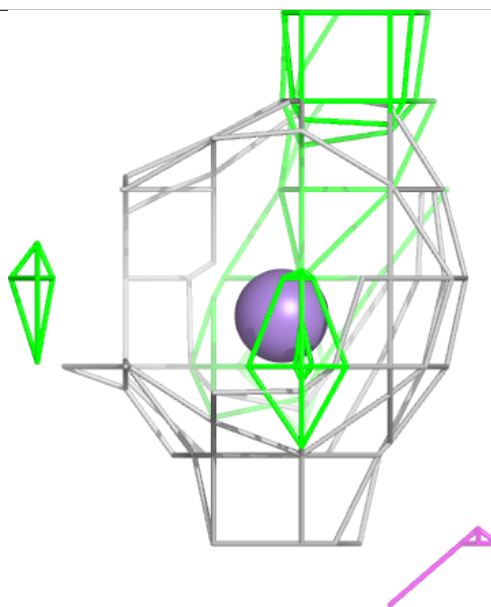
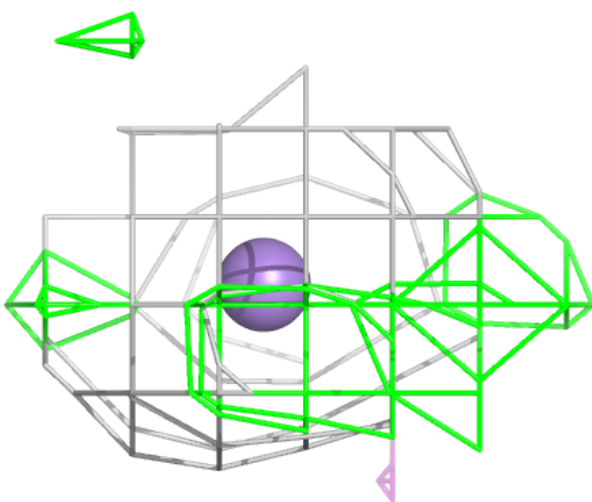
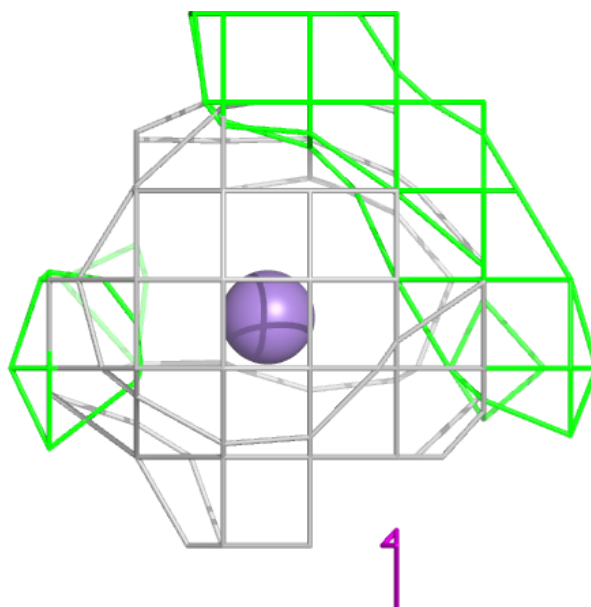
Electron density around MN B 302:

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



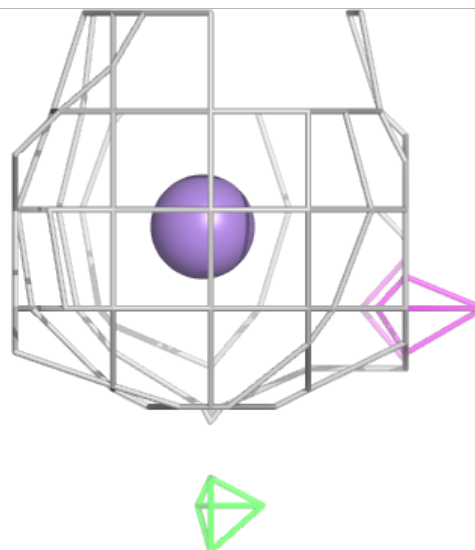
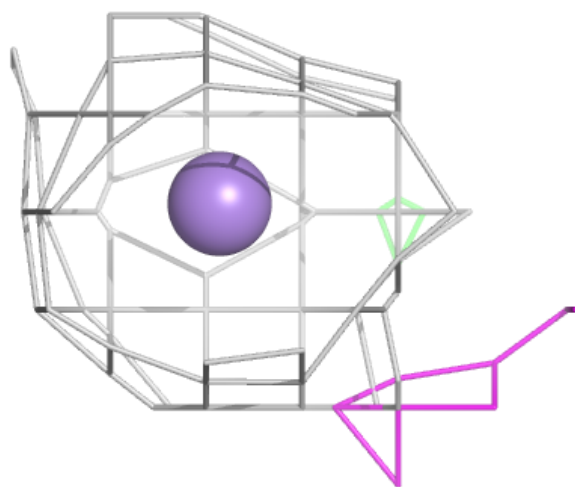
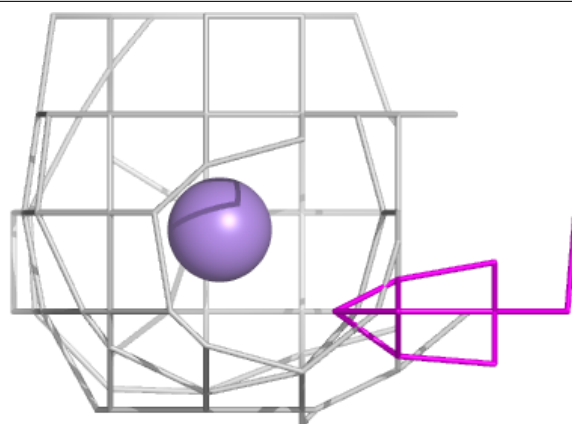
Electron density around MN D 103:

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



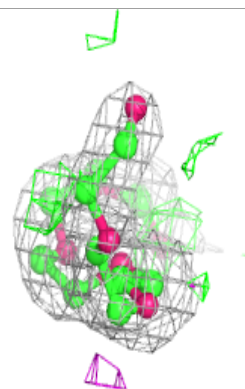
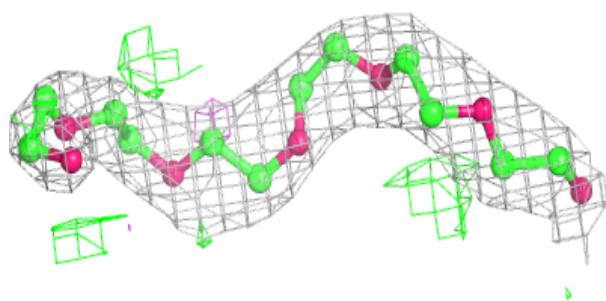
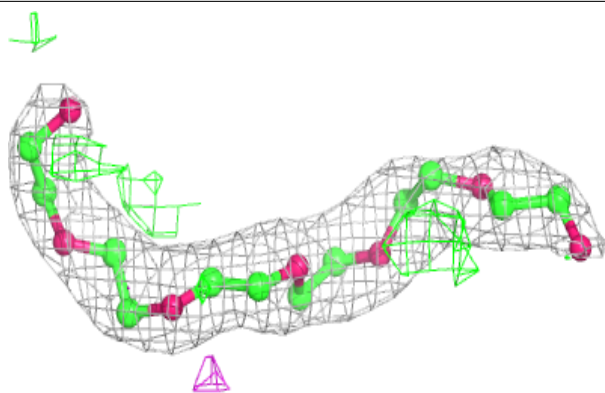
Electron density around MN C 101:

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



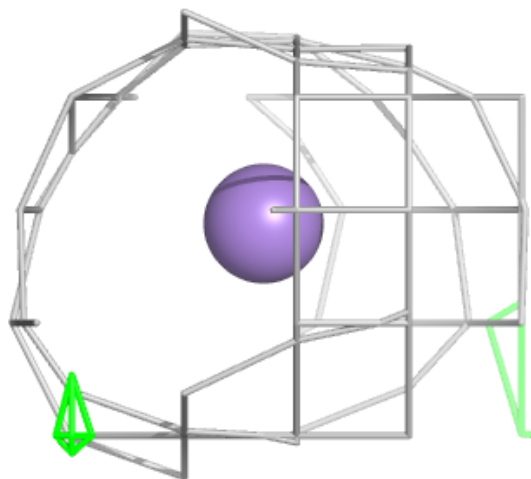
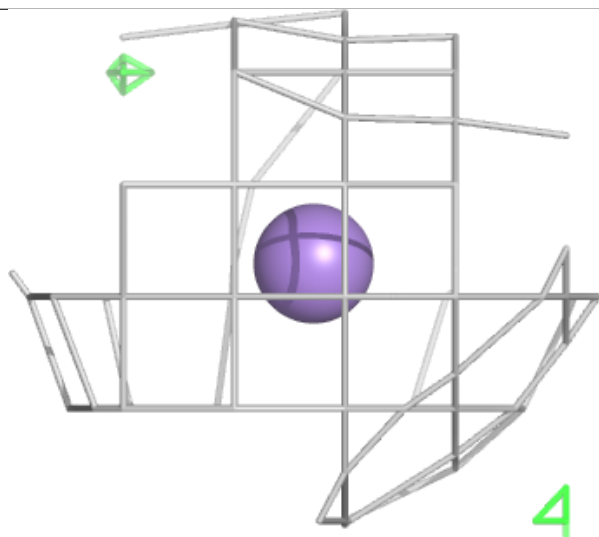
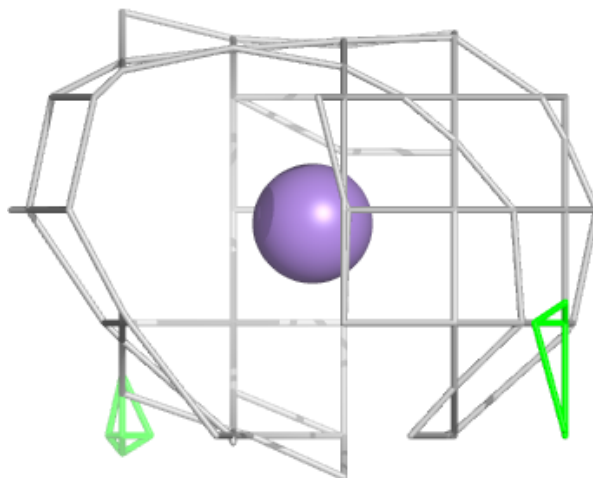
Electron density around P6G A 310:

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



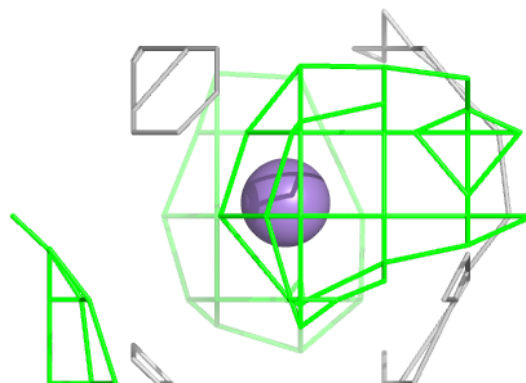
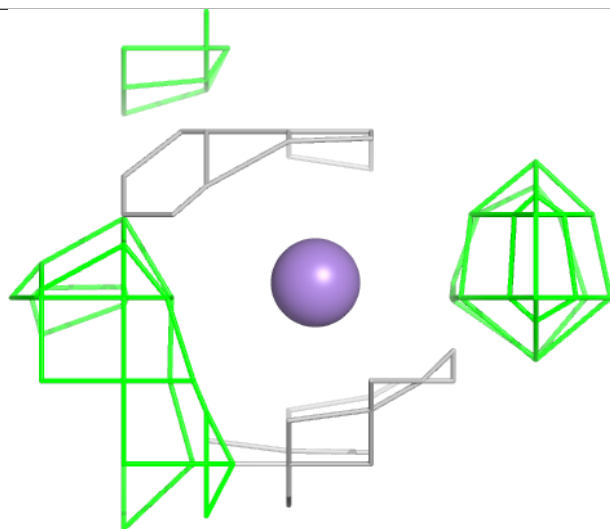
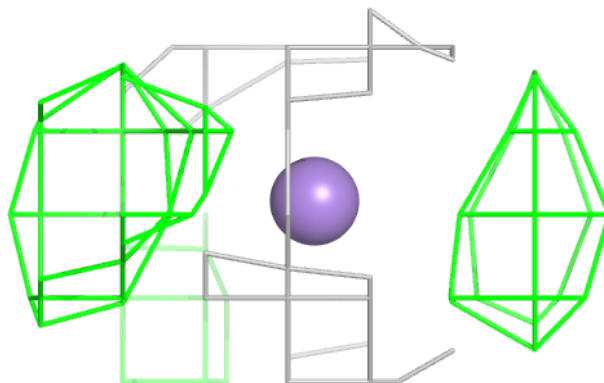
Electron density around MN A 303:

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



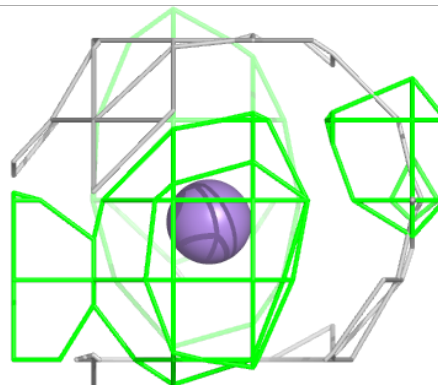
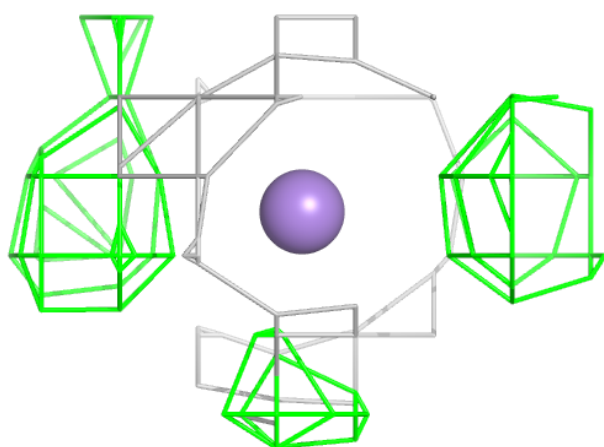
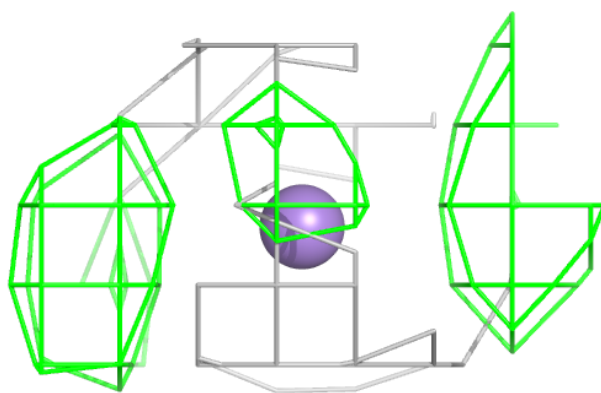
Electron density around MN D 101:

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



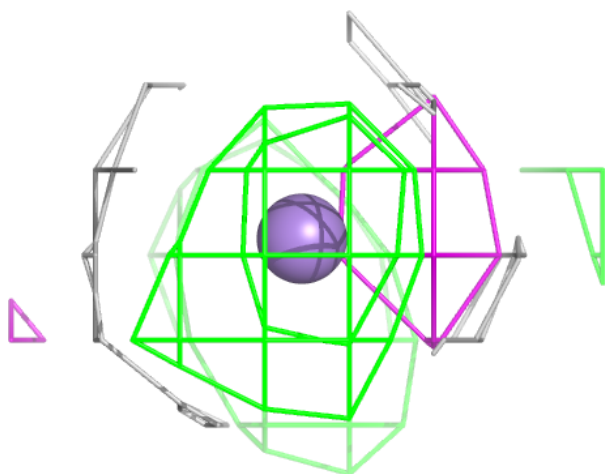
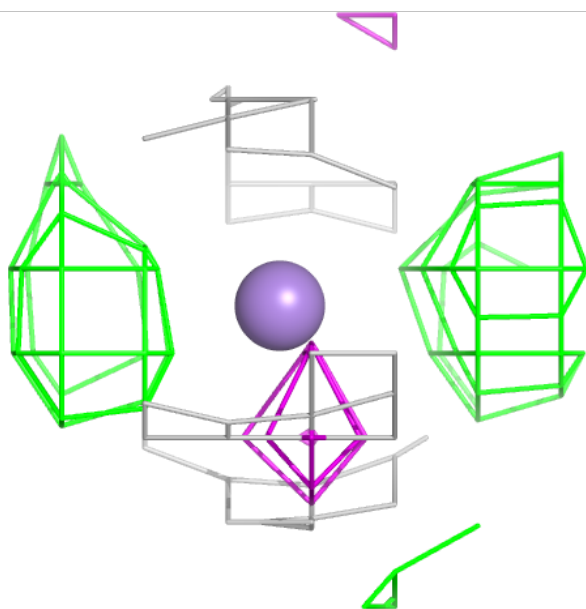
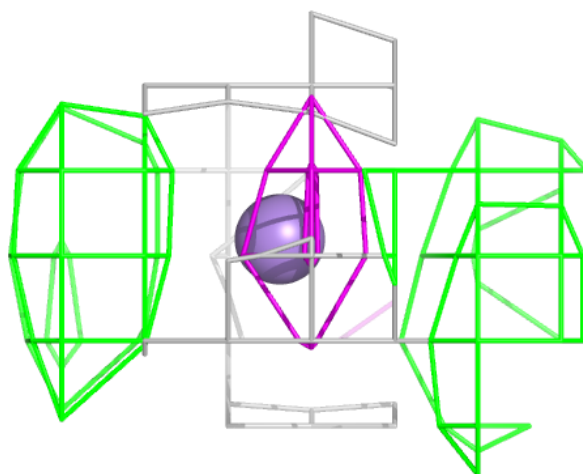
Electron density around MN A 302:

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



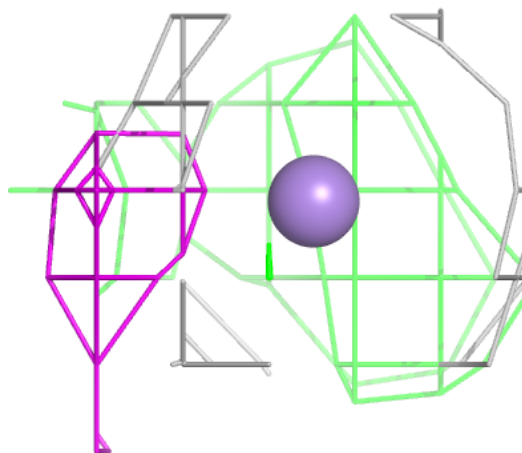
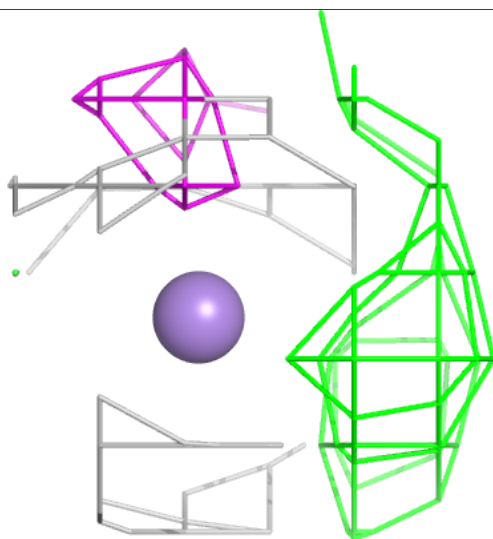
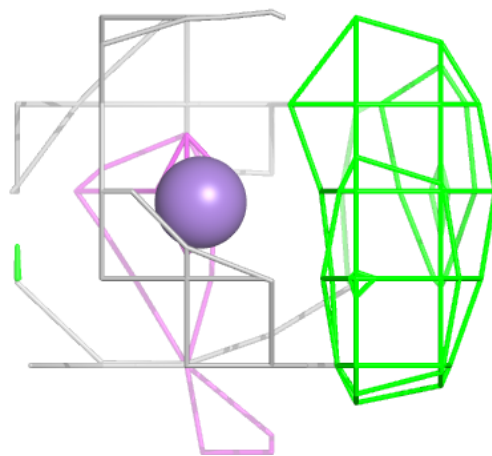
Electron density around MN B 301:

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



Electron density around MN 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)



6.5 Other polymers ⓘ

There are no such residues in this entry.