



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 16, 2024 – 03:33 pm GMT

PDB ID : 9GOF
Title : MncA bound to nickel
Authors : Glasfeld, A.; Robinson, N.J.
Deposited on : 2024-09-05
Resolution : 1.60 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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

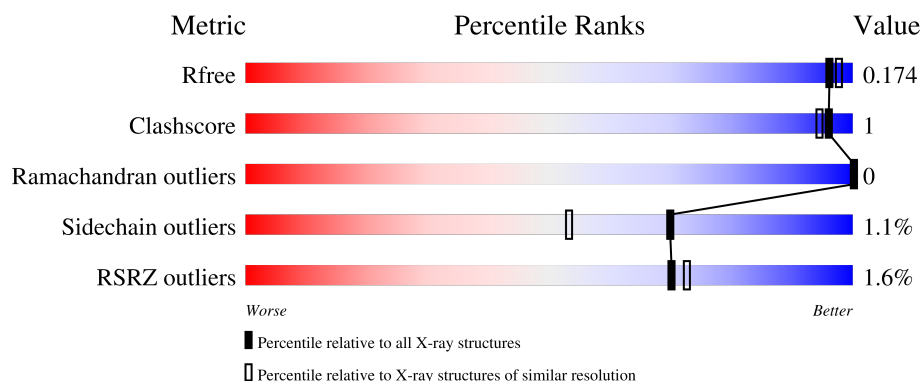
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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}	164625	4274 (1.60-1.60)
Clashscore	180529	4682 (1.60-1.60)
Ramachandran outliers	177936	4583 (1.60-1.60)
Sidechain outliers	177891	4582 (1.60-1.60)
RSRZ outliers	164620	4272 (1.60-1.60)

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	361	<div> <div>2%</div> <div> <div></div> <div>96%</div> <div>..</div> </div> </div>
1	B	361	<div> <div>%</div> <div> <div></div> <div>94%</div> <div>5% .</div> </div> </div>
1	C	361	<div> <div>2%</div> <div> <div></div> <div>96%</div> <div>..</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9298 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 Sll1358 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	356	Total	C	N	O	S	0	3	0
			2776	1783	463	524	6			
1	B	356	Total	C	N	O	S	0	1	0
			2758	1769	461	522	6			
1	C	354	Total	C	N	O	S	0	5	0
			2772	1780	462	524	6			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	MET	-	initiating methionine	UNP P73510
B	34	MET	-	initiating methionine	UNP P73510
C	34	MET	-	initiating methionine	UNP P73510

- Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni) (labeled as "Ligand of Interest" by depositor).

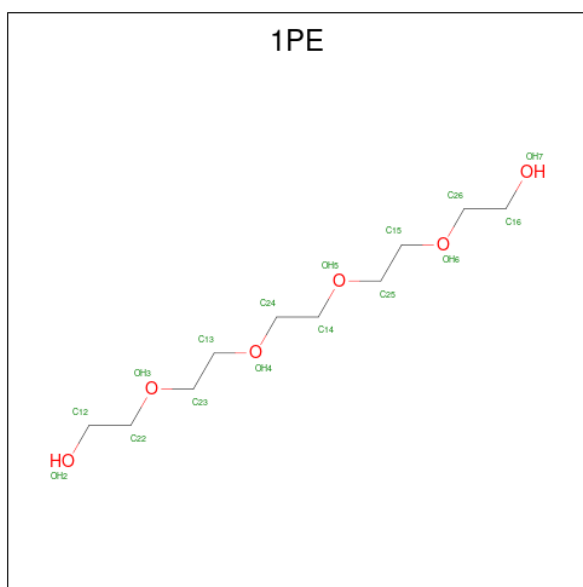
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ni	0	0
			2	2		
2	B	2	Total	Ni	0	0
			2	2		
2	C	2	Total	Ni	0	0
			2	2		

- Molecule 3 is GLYCINE (three-letter code: GLY) (formula: C₂H₅NO₂).



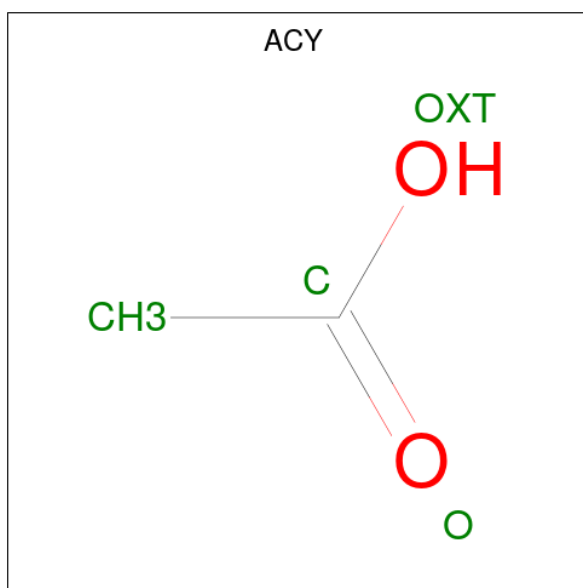
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			5	2	1	2		
3	B	1	Total	C	N	O	0	0
			5	2	1	2		
3	C	1	Total	C	N	O	0	0
			5	2	1	2		

- Molecule 4 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			16	10	6		

- Molecule 5 is ACETIC ACID (three-letter code: ACY) (formula: $C_2H_4O_2$).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	323	Total	O	0	0
			323	323		
6	B	312	Total	O	0	0
			312	312		
6	C	312	Total	O	0	0
			312	312		

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: Sll1358 protein



- Molecule 1: Sll1358 protein



- Molecule 1: Sll1358 protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	235.19Å 235.19Å 132.01Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	80.63 – 1.60 80.63 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.8 (80.63-1.60) 99.8 (80.63-1.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 1.60Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.164 , 0.177 0.167 , 0.174	Depositor DCC
R_{free} test set	2142 reflections (0.65%)	wwPDB-VP
Wilson B-factor (Å ²)	22.8	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9298	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.20% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACY, NI, 1PE

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	0.36	0/2867	0.59	0/3916
1	B	0.36	0/2844	0.58	0/3883
1	C	0.37	0/2865	0.57	0/3913
All	All	0.36	0/8576	0.58	0/11712

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 [i](#)

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	2776	0	2684	5	0
1	B	2758	0	2669	8	0
1	C	2772	0	2678	4	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	5	0	2	1	0
3	B	5	0	2	1	0
3	C	5	0	2	0	0
4	A	16	0	22	3	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	8	0	6	0	0
6	A	323	0	0	1	0
6	B	312	0	0	0	0
6	C	312	0	0	0	0
All	All	9298	0	8065	17	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 1.

All (17) 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)
4:A:1004:1PE:H141	6:A:1122:HOH:O	2.11	0.50
1:A:290:GLU:OE2	3:A:1003:GLY:N	2.48	0.46
1:B:272:ILE:HB	1:B:341:ILE:HG12	1.98	0.45
1:B:286:PRO:HG2	1:C:124:PRO:HG3	1.98	0.45
1:B:279:MET:HG3	1:B:332:ARG:HG3	1.98	0.44
1:B:109:TRP:CZ2	1:B:269:GLY:HA3	2.53	0.43
1:C:283:HIS:HB3	1:C:354:ASP:OD1	2.20	0.42
1:B:178:ASP:HB2	1:B:206:GLN:HB2	2.01	0.42
1:B:292:GLN:HG2	1:B:343:VAL:HG12	2.01	0.41
1:A:85:MET:HA	1:A:163:PHE:O	2.20	0.41
1:A:171:GLY:N	4:A:1004:1PE:H222	2.36	0.41
4:A:1004:1PE:H161	4:A:1004:1PE:H152	1.57	0.41
1:B:290:GLU:OE2	3:B:1003:GLY:N	2.53	0.41
1:B:375:PRO:O	1:B:379:LYS:HG2	2.21	0.41
1:A:124:PRO:HG3	1:C:286:PRO:HG2	2.03	0.41
1:C:109:TRP:CZ2	1:C:269:GLY:HA3	2.56	0.40
1:A:283:HIS:HB3	1:A:354:ASP:OD1	2.22	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	356/361 (99%)	351 (99%)	5 (1%)	0	100	100
1	B	355/361 (98%)	348 (98%)	7 (2%)	0	100	100
1	C	356/361 (99%)	350 (98%)	6 (2%)	0	100	100
All	All	1067/1083 (98%)	1049 (98%)	18 (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	296/298 (99%)	293 (99%)	3 (1%)	73	57
1	B	294/298 (99%)	290 (99%)	4 (1%)	62	43
1	C	296/298 (99%)	293 (99%)	3 (1%)	73	57
All	All	886/894 (99%)	876 (99%)	10 (1%)	70	53

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	HIS
1	A	191	ASN
1	A	247	SER
1	B	77	TYR
1	B	103	HIS
1	B	175	SER
1	B	191	ASN
1	C	103	HIS
1	C	175	SER
1	C	191	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	227	GLN
1	A	316	GLN
1	A	393	GLN
1	B	206	GLN
1	B	393	GLN
1	C	225	GLN

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 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 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	1PE	A	1004	-	15,15,15	0.12	0	14,14,14	0.24	0
3	GLY	B	1003	2	4,4,4	0.89	0	3,4,4	1.61	1 (33%)
3	GLY	A	1003	2	4,4,4	1.01	0	3,4,4	1.80	2 (66%)
5	ACY	B	1005	-	3,3,3	1.25	0	3,3,3	1.40	0
3	GLY	C	1003	2	4,4,4	0.99	0	3,4,4	1.63	0
5	ACY	B	1004	-	3,3,3	1.39	1 (33%)	3,3,3	1.26	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
3	GLY	A	1003	2	-	0/2/2/2	-
3	GLY	C	1003	2	-	0/2/2/2	-
3	GLY	B	1003	2	-	0/2/2/2	-
4	1PE	A	1004	-	-	5/13/13/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1004	ACY	CH3-C	2.12	1.58	1.49

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1003	GLY	OXT-C-O	-2.24	117.71	123.30
3	A	1003	GLY	OXT-C-CA	2.12	121.90	113.45
3	B	1003	GLY	OXT-C-CA	2.06	121.66	113.45

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1004	1PE	OH5-C14-C24-OH4
4	A	1004	1PE	OH2-C12-C22-OH3
4	A	1004	1PE	C16-C26-OH6-C15
4	A	1004	1PE	OH4-C13-C23-OH3
4	A	1004	1PE	OH6-C15-C25-OH5

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1004	1PE	3	0
3	B	1003	GLY	1	0
3	A	1003	GLY	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 [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	356/361 (98%)	-0.33	8 (2%) 62 65	16, 23, 33, 43	3 (0%)
1	B	356/361 (98%)	-0.36	3 (0%) 82 85	13, 23, 31, 42	1 (0%)
1	C	354/361 (98%)	-0.34	6 (1%) 69 71	12, 23, 33, 60	5 (1%)
All	All	1066/1083 (98%)	-0.35	17 (1%) 70 73	12, 23, 32, 60	9 (0%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	392	THR	6.0
1	A	39[A]	TRP	4.7
1	C	39[A]	TRP	4.7
1	B	39	TRP	3.6
1	A	40	ARG	2.9
1	A	153	GLY	2.9
1	A	247	SER	2.8
1	B	40	ARG	2.6
1	C	247[A]	SER	2.4
1	A	152	ILE	2.3
1	C	153	GLY	2.2
1	A	227	GLN	2.2
1	C	227	GLN	2.2
1	B	227	GLN	2.1
1	A	246	VAL	2.1
1	A	392	THR	2.0
1	C	40	ARG	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](#)

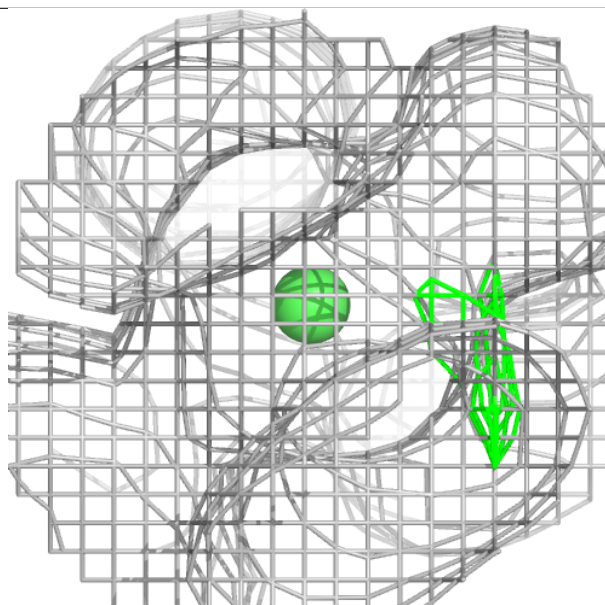
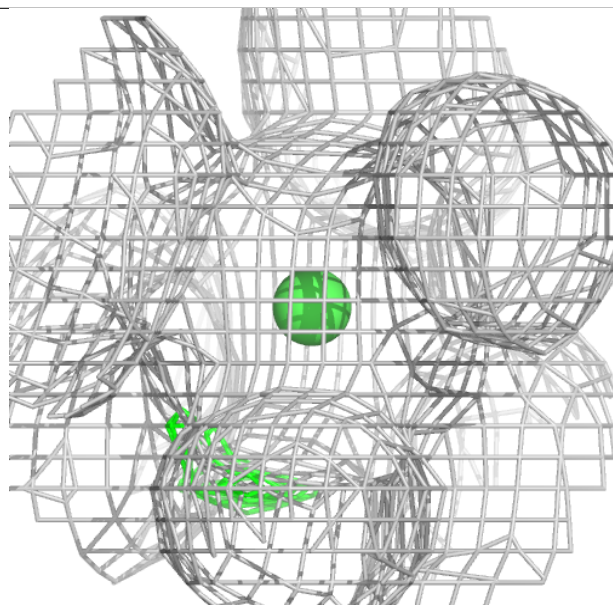
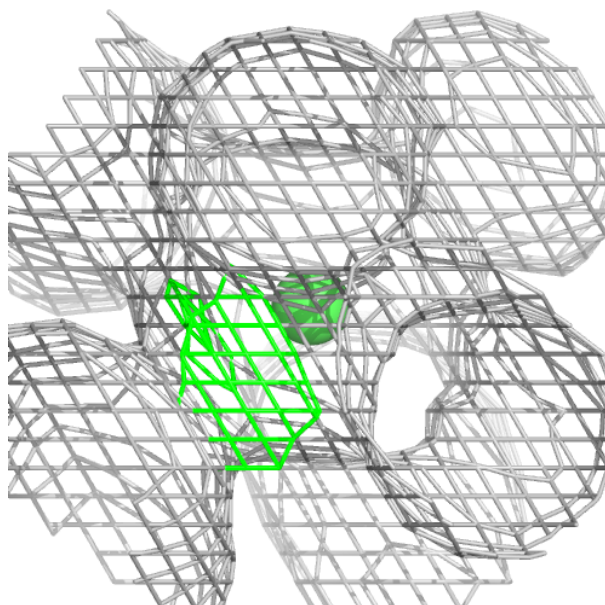
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	ACY	B	1004	4/4	0.77	0.21	32,32,36,46	0
4	1PE	A	1004	16/16	0.90	0.14	26,36,43,44	0
3	GLY	A	1003	5/5	0.96	0.08	17,21,26,29	0
3	GLY	B	1003	5/5	0.96	0.08	17,20,26,29	0
3	GLY	C	1003	5/5	0.97	0.08	17,21,25,31	0
5	ACY	B	1005	4/4	0.99	0.04	22,22,23,24	0
2	NI	A	1001	1/1	1.00	0.01	18,18,18,18	0
2	NI	A	1002	1/1	1.00	0.01	20,20,20,20	0
2	NI	B	1001	1/1	1.00	0.01	19,19,19,19	0
2	NI	B	1002	1/1	1.00	0.02	19,19,19,19	0
2	NI	C	1001	1/1	1.00	0.01	19,19,19,19	0
2	NI	C	1002	1/1	1.00	0.01	20,20,20,20	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.

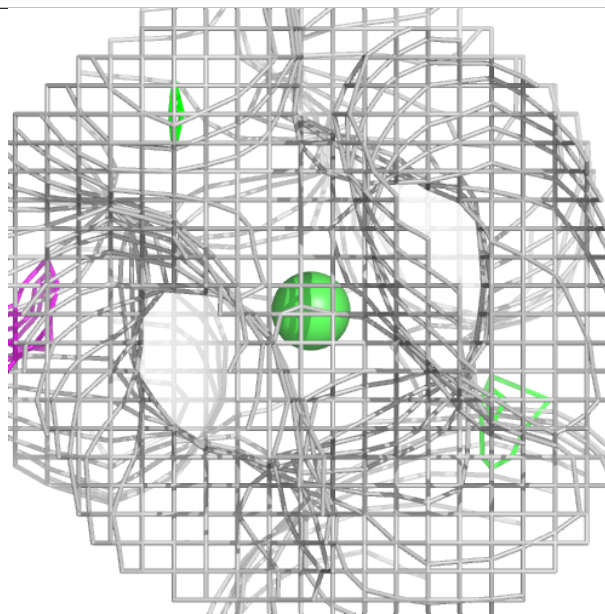
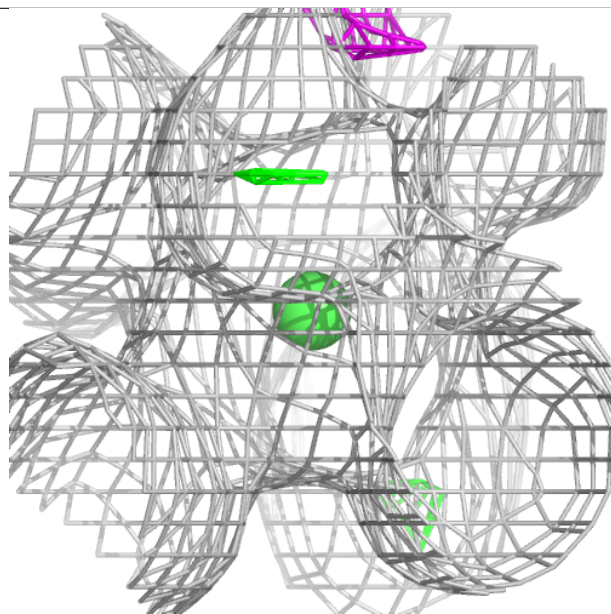
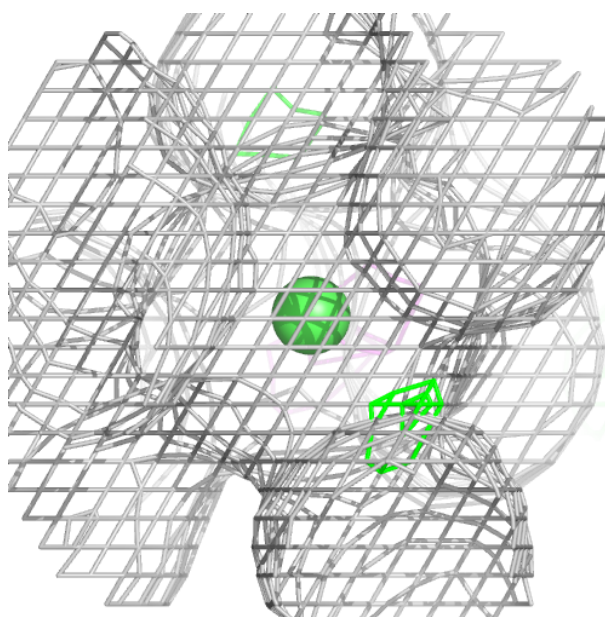
Electron density around NI A 1001:

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



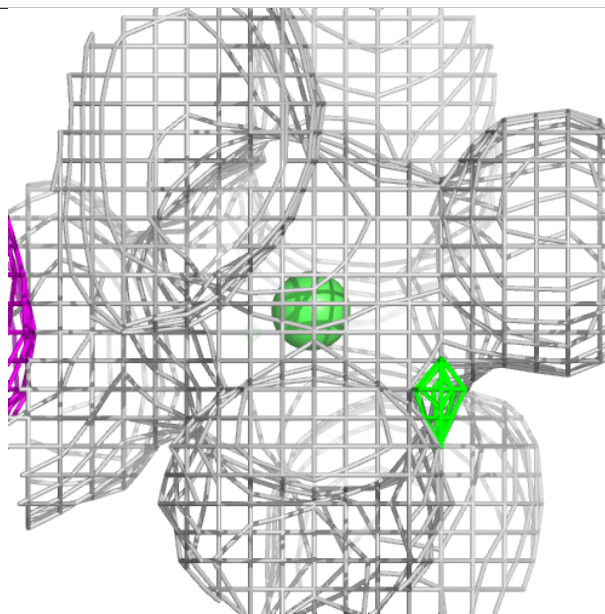
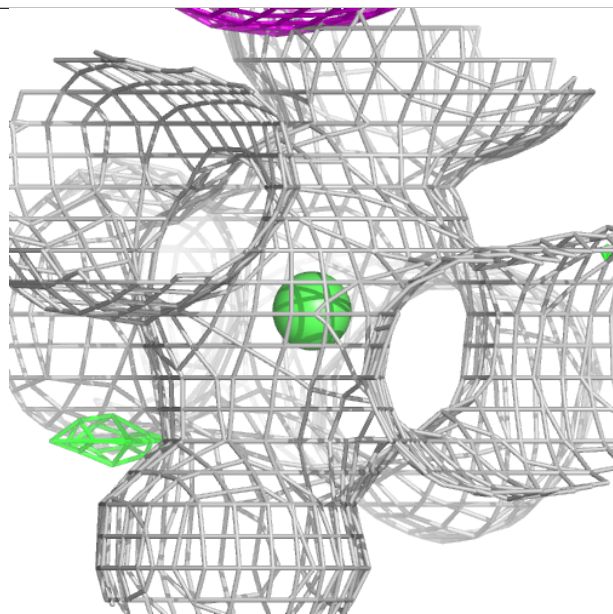
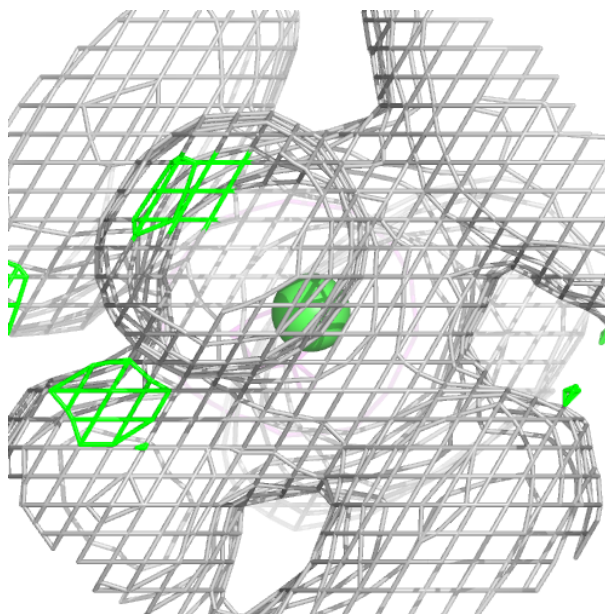
Electron density around NI A 1002:

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



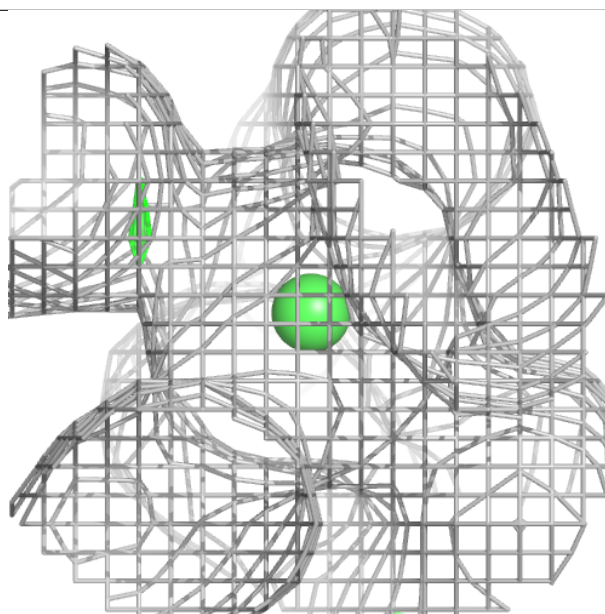
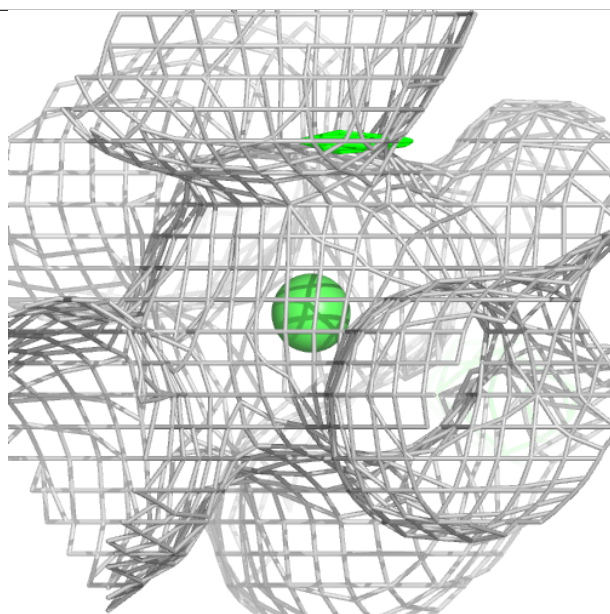
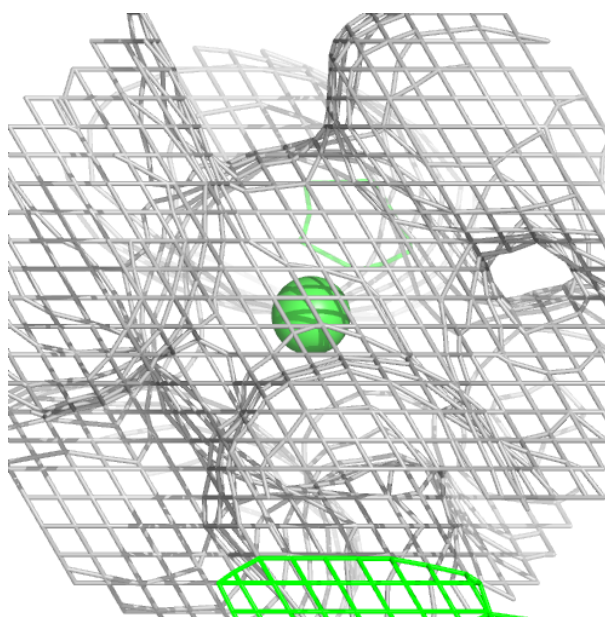
Electron density around NI B 1001:

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



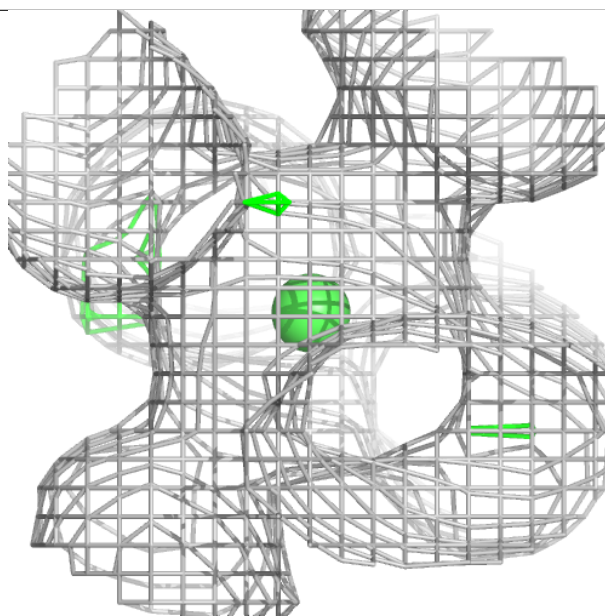
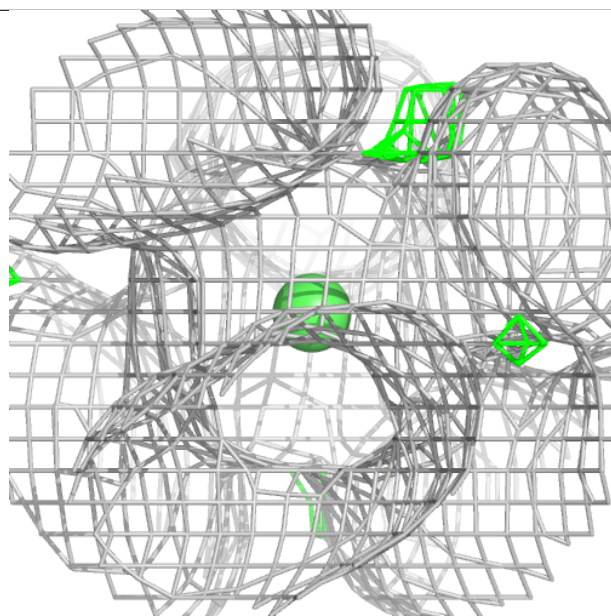
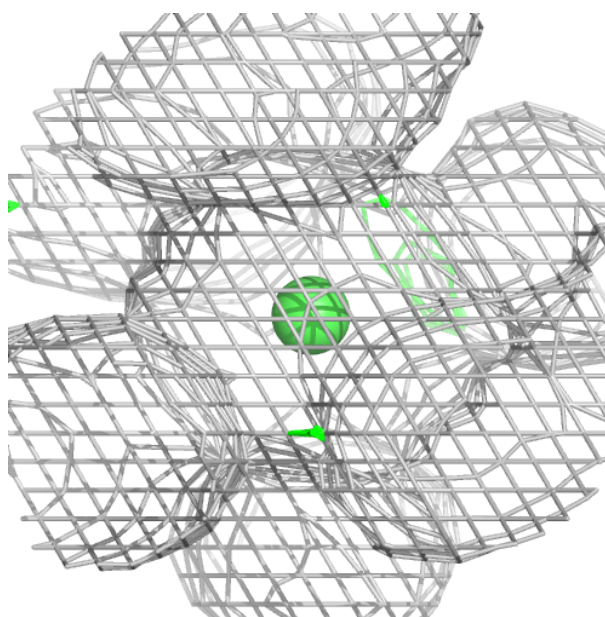
Electron density around NI B 1002:

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



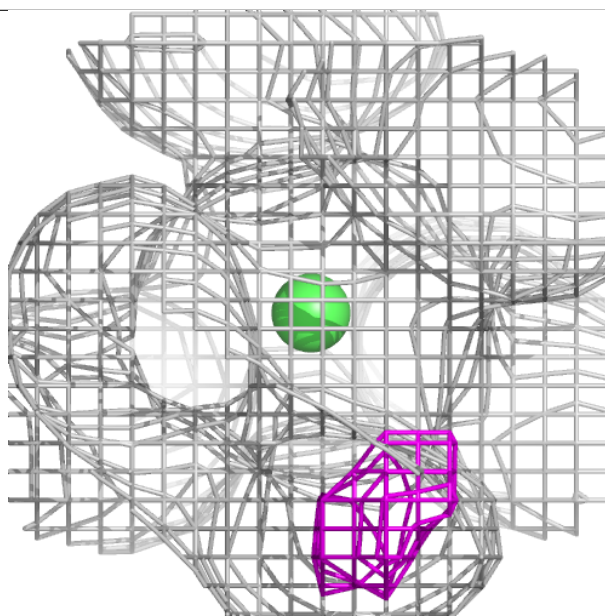
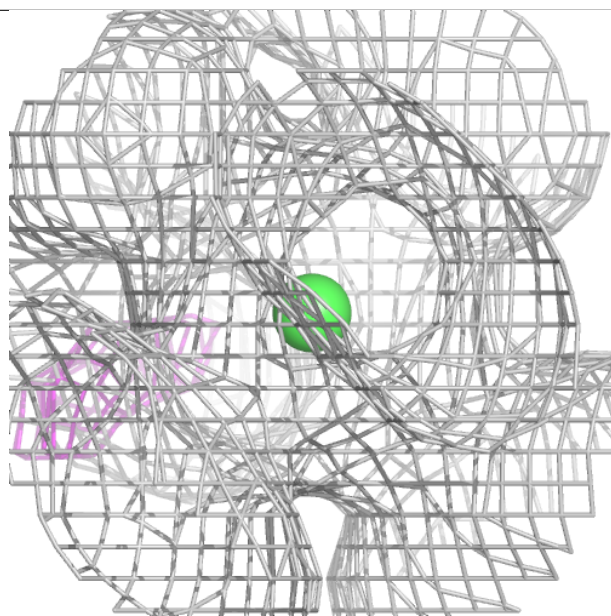
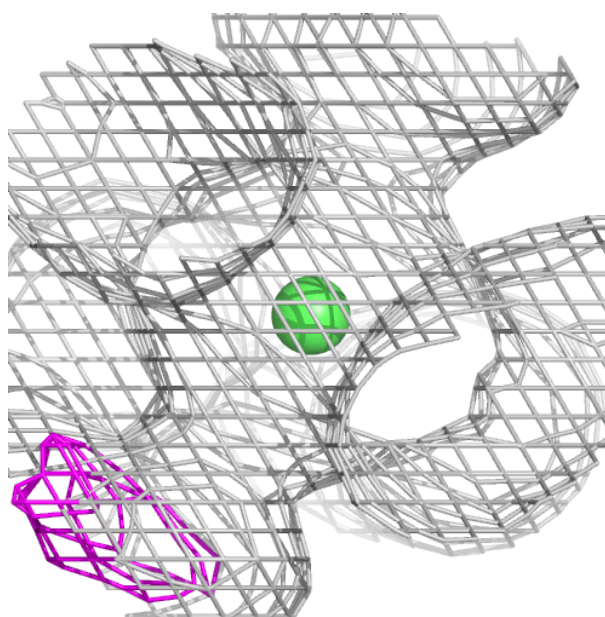
Electron density around NI C 1001:

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



Electron density around NI C 1002:

$2mF_o-DF_c$ (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.