



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:16 AM UTC

PDB ID : 8XV8 / pdb_00008xv8
Title : Crystal structure of PHD domain of UHRF1 in complex with hStella peptide (residues 75-121)
Authors : Du, X.; Gan, Q.; Xu, J.; Liu, J.
Deposited on : 2024-01-14
Resolution : 2.05 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

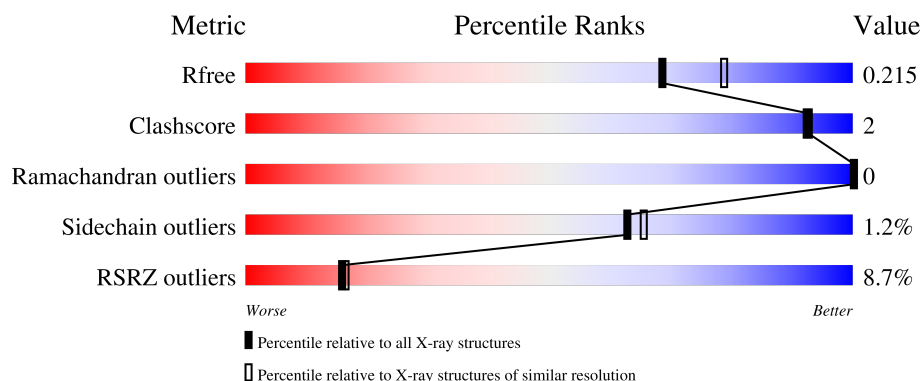
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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.



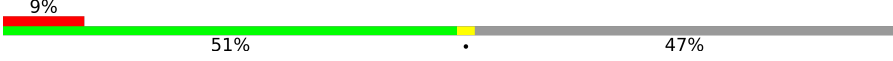
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	75	8% 92% 7%
1	C	75	3% 88% 7%
1	E	75	16% 79% 8% 13%
1	G	75	4% 81% 5% 13%
2	B	47	2% 49% 47%

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Mol	Chain	Length	Quality of chain
2	D	47	
2	F	47	
2	H	47	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3067 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 E3 ubiquitin-protein ligase UHRF1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	70	Total	C	N	O	S	0	0	0
			539	326	95	105	13			
1	C	70	Total	C	N	O	S	0	0	0
			539	326	95	105	13			
1	E	65	Total	C	N	O	S	0	0	0
			511	310	90	98	13			
1	G	65	Total	C	N	O	S	0	0	0
			511	310	90	98	13			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	293	GLY	-	expression tag	UNP Q96T88
A	294	PRO	-	expression tag	UNP Q96T88
A	295	LEU	-	expression tag	UNP Q96T88
A	296	GLY	-	expression tag	UNP Q96T88
A	297	SER	-	expression tag	UNP Q96T88
C	293	GLY	-	expression tag	UNP Q96T88
C	294	PRO	-	expression tag	UNP Q96T88
C	295	LEU	-	expression tag	UNP Q96T88
C	296	GLY	-	expression tag	UNP Q96T88
C	297	SER	-	expression tag	UNP Q96T88
E	293	GLY	-	expression tag	UNP Q96T88
E	294	PRO	-	expression tag	UNP Q96T88
E	295	LEU	-	expression tag	UNP Q96T88
E	296	GLY	-	expression tag	UNP Q96T88
E	297	SER	-	expression tag	UNP Q96T88
G	293	GLY	-	expression tag	UNP Q96T88
G	294	PRO	-	expression tag	UNP Q96T88
G	295	LEU	-	expression tag	UNP Q96T88
G	296	GLY	-	expression tag	UNP Q96T88
G	297	SER	-	expression tag	UNP Q96T88

- Molecule 2 is a protein called Developmental pluripotency-associated protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	25	Total	C	N	O	S	0	0	0
			204	128	42	32	2			
2	D	27	Total	C	N	O	S	0	0	0
			219	136	47	34	2			
2	F	25	Total	C	N	O	S	0	0	0
			201	126	42	31	2			
2	H	32	Total	C	N	O	S	0	0	0
			269	170	53	44	2			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	Zn	0	0
			4	4		
3	C	4	Total	Zn	0	0
			4	4		
3	E	3	Total	Zn	0	0
			3	3		
3	G	3	Total	Zn	0	0
			3	3		

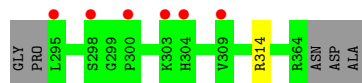
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	18	Total	O	0	0
			18	18		
4	B	2	Total	O	0	0
			2	2		
4	C	12	Total	O	0	0
			12	12		
4	D	6	Total	O	0	0
			6	6		
4	E	3	Total	O	0	0
			3	3		
4	F	1	Total	O	0	0
			1	1		
4	G	14	Total	O	0	0
			14	14		
4	H	4	Total	O	0	0
			4	4		

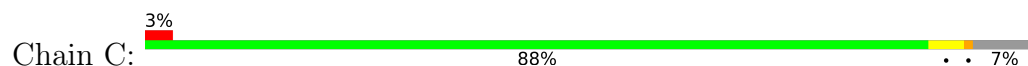
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

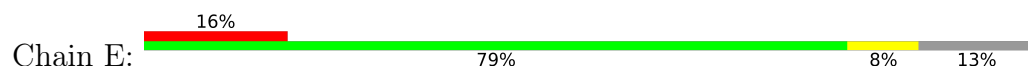
- Molecule 1: E3 ubiquitin-protein ligase UHRF1



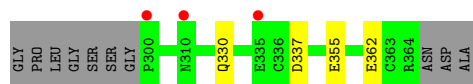
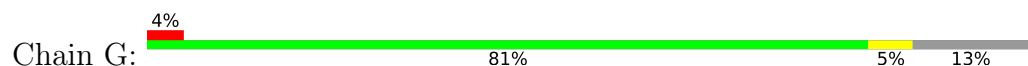
- Molecule 1: E3 ubiquitin-protein ligase UHRF1



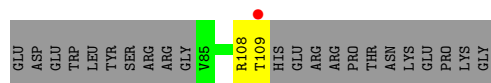
- Molecule 1: E3 ubiquitin-protein ligase UHRF1



- Molecule 1: E3 ubiquitin-protein ligase UHRF1



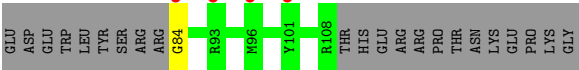
- Molecule 2: Developmental pluripotency-associated protein 3



- Molecule 2: Developmental pluripotency-associated protein 3



• Molecule 2: Developmental pluripotency-associated protein 3



• Molecule 2: Developmental pluripotency-associated protein 3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.03Å 46.02Å 82.91Å 90.00° 112.76° 90.00°	Depositor
Resolution (Å)	38.23 – 2.05 38.23 – 2.05	Depositor EDS
% Data completeness (in resolution range)	97.5 (38.23-2.05) 97.5 (38.23-2.05)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, R_{free}	0.186 , 0.207 0.194 , 0.215	Depositor DCC
R_{free} test set	1370 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	1.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3067	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.34% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.68	0/552	0.98	0/746
1	C	0.64	0/552	1.02	1/746 (0.1%)
1	E	0.58	0/524	0.98	0/708
1	G	0.64	0/524	0.96	1/708 (0.1%)
2	B	0.56	0/204	0.95	0/270
2	D	0.60	0/219	1.01	0/289
2	F	0.60	0/201	1.01	0/265
2	H	0.65	0/272	1.07	0/362
All	All	0.63	0/3048	0.99	2/4094 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	D	0	1
All	All	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	362	GLU	CB-CG-CD	5.35	121.70	112.60
1	G	362	GLU	CG-CD-OE1	5.29	130.56	118.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	314	ARG	Sidechain
2	D	100	ARG	Sidechain

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	539	0	486	0	0
1	C	539	0	486	2	0
1	E	511	0	460	3	0
1	G	511	0	460	2	0
2	B	204	0	230	1	0
2	D	219	0	246	0	0
2	F	201	0	226	1	0
2	H	269	0	284	4	0
3	A	4	0	0	0	0
3	C	4	0	0	0	0
3	E	3	0	0	0	0
3	G	3	0	0	0	0
4	A	18	0	0	0	0
4	B	2	0	0	0	0
4	C	12	0	0	0	0
4	D	6	0	0	0	0
4	E	3	0	0	0	0
4	F	1	0	0	0	0
4	G	14	0	0	0	0
4	H	4	0	0	0	0
All	All	3067	0	2878	9	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 2.

All (9) 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)
2:H:76:ASP:OD2	2:H:79:LEU:HD13	2.11	0.50
2:B:108:ARG:O	2:B:109:THR:C	2.56	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:330:GLN:O	2:H:87:THR:HA	2.15	0.45
1:E:356:ASP:HA	2:F:84:GLY:HA2	1.99	0.45
2:H:77:GLU:OE2	2:H:80:TYR:OH	2.31	0.43
1:G:355:GLU:O	2:H:85:VAL:HG23	2.19	0.42
1:C:314:ARG:NH1	1:E:362:GLU:HB3	2.34	0.42
1:C:359:TYR:O	1:C:364:ARG:NH1	2.48	0.42
1:E:347:PRO:HA	1:E:348:PRO:HD3	2.00	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	68/75 (91%)	65 (96%)	3 (4%)	0	100	100
1	C	68/75 (91%)	66 (97%)	2 (3%)	0	100	100
1	E	63/75 (84%)	60 (95%)	3 (5%)	0	100	100
1	G	63/75 (84%)	60 (95%)	3 (5%)	0	100	100
2	B	23/47 (49%)	23 (100%)	0	0	100	100
2	D	25/47 (53%)	25 (100%)	0	0	100	100
2	F	23/47 (49%)	23 (100%)	0	0	100	100
2	H	30/47 (64%)	30 (100%)	0	0	100	100
All	All	363/488 (74%)	352 (97%)	11 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	64/67 (96%)	64 (100%)	0	100	100
1	C	64/67 (96%)	63 (98%)	1 (2%)	55	56
1	E	61/67 (91%)	59 (97%)	2 (3%)	33	28
1	G	61/67 (91%)	60 (98%)	1 (2%)	55	56
2	B	22/42 (52%)	22 (100%)	0	100	100
2	D	23/42 (55%)	23 (100%)	0	100	100
2	F	21/42 (50%)	21 (100%)	0	100	100
2	H	28/42 (67%)	28 (100%)	0	100	100
All	All	344/436 (79%)	340 (99%)	4 (1%)	63	65

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	364	ARG
1	E	337	ASP
1	E	354	SER
1	G	337	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	92	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 14 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [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	70/75 (93%)	0.33	6 (8%) 16 16	24, 35, 65, 77	0
1	C	70/75 (93%)	0.30	2 (2%) 53 55	24, 38, 65, 81	0
1	E	65/75 (86%)	1.07	12 (18%) 3 2	30, 47, 83, 94	0
1	G	65/75 (86%)	0.17	3 (4%) 37 37	26, 34, 62, 73	0
2	B	25/47 (53%)	0.62	1 (4%) 42 42	29, 42, 60, 72	0
2	D	27/47 (57%)	0.54	3 (11%) 10 10	25, 40, 58, 90	0
2	F	25/47 (53%)	0.96	4 (16%) 5 4	34, 46, 70, 89	0
2	H	32/47 (68%)	0.70	2 (6%) 26 26	30, 39, 72, 83	0
All	All	379/488 (77%)	0.53	33 (8%) 16 16	24, 40, 74, 94	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	101	TYR	5.0
2	H	107	VAL	4.9
1	C	295	LEU	4.6
1	E	349	LEU	4.0
1	A	295	LEU	3.9
1	E	304	HIS	3.8
1	E	300	PRO	3.6
1	E	353	PRO	3.5
1	G	310	ASN	3.5
1	E	352	VAL	3.4
1	G	300	PRO	3.3
1	E	364	ARG	3.3
2	H	78	TRP	3.0
2	B	109	THR	3.0
2	D	109	THR	2.9
1	A	300	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
1	E	348	PRO	2.8
1	A	304	HIS	2.8
1	E	347	PRO	2.7
1	E	342	ILE	2.5
2	F	84	GLY	2.5
1	C	364	ARG	2.4
2	F	96	MET	2.4
2	D	83	ARG	2.4
2	F	93	ARG	2.3
1	G	335	GLU	2.3
1	E	351	SER	2.2
1	A	303	LYS	2.2
1	A	309	VAL	2.2
1	A	298	SER	2.1
1	E	309	VAL	2.1
1	E	346	ASP	2.0
2	D	84	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [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($\text{\AA}^2$)	Q<0.9
3	ZN	E	402	1/1	0.98	0.03	39,39,39,39	0
3	ZN	A	404	1/1	0.99	0.02	28,28,28,28	0
3	ZN	C	401	1/1	0.99	0.03	35,35,35,35	0
3	ZN	C	403	1/1	0.99	0.04	34,34,34,34	0
3	ZN	C	404	1/1	0.99	0.03	31,31,31,31	0
3	ZN	A	401	1/1	0.99	0.02	35,35,35,35	0

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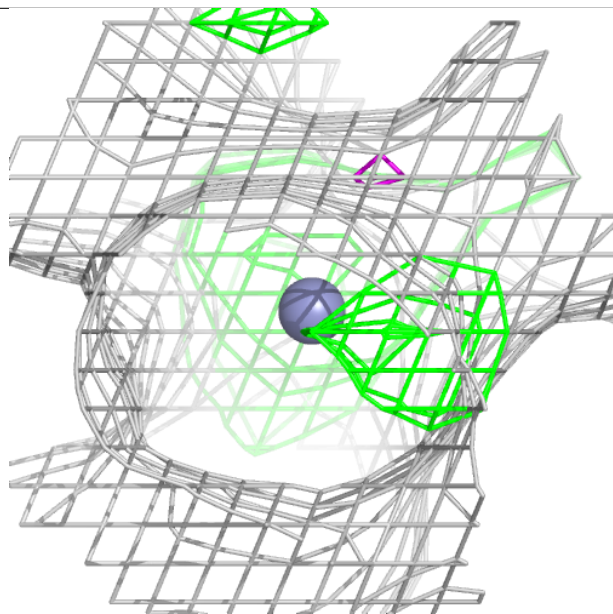
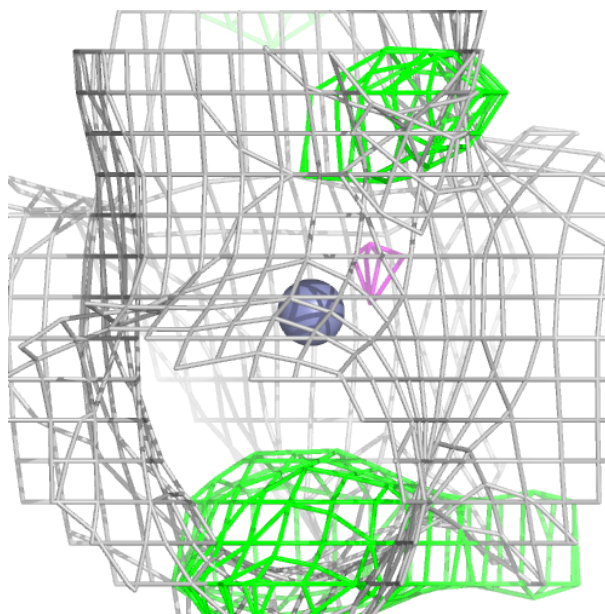
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ZN	E	403	1/1	0.99	0.03	37,37,37,37	0
3	ZN	G	401	1/1	0.99	0.03	36,36,36,36	0
3	ZN	E	401	1/1	1.00	0.01	38,38,38,38	0
3	ZN	C	402	1/1	1.00	0.01	26,26,26,26	0
3	ZN	A	402	1/1	1.00	0.01	27,27,27,27	0
3	ZN	A	403	1/1	1.00	0.02	31,31,31,31	0
3	ZN	G	402	1/1	1.00	0.02	30,30,30,30	0
3	ZN	G	403	1/1	1.00	0.02	29,29,29,29	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.

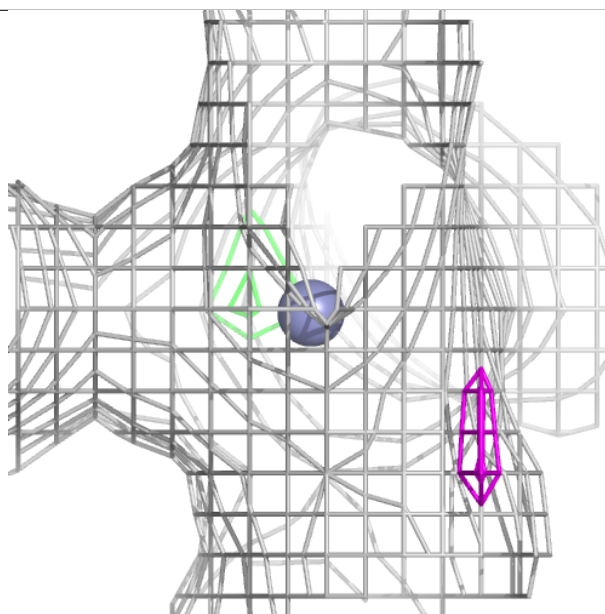
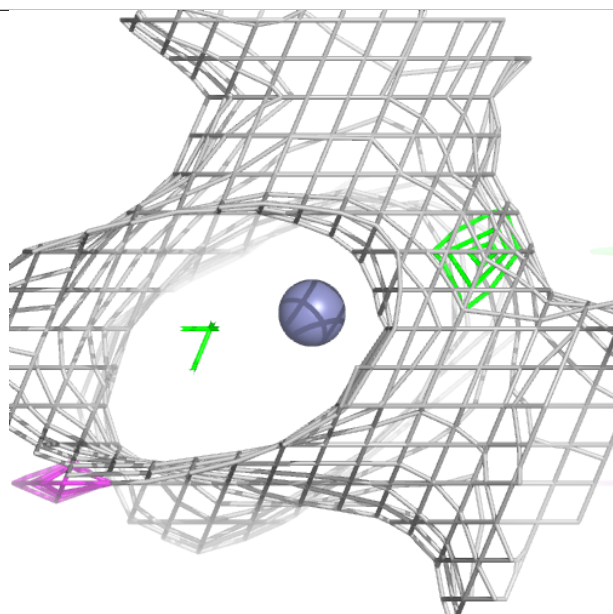
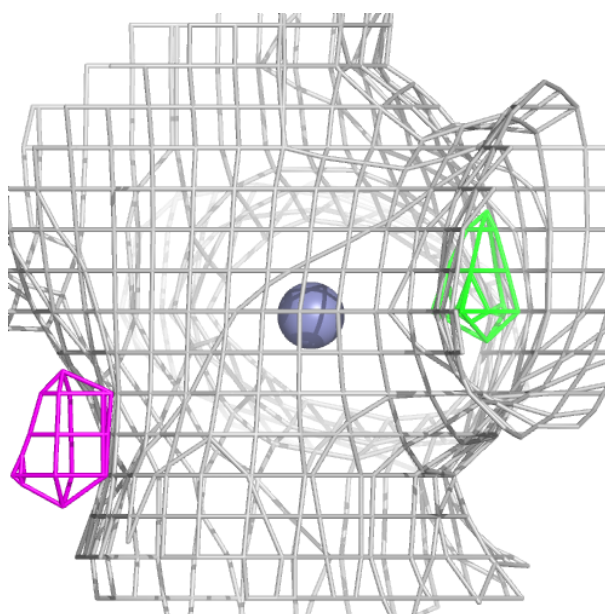
Electron density around ZN E 402:

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



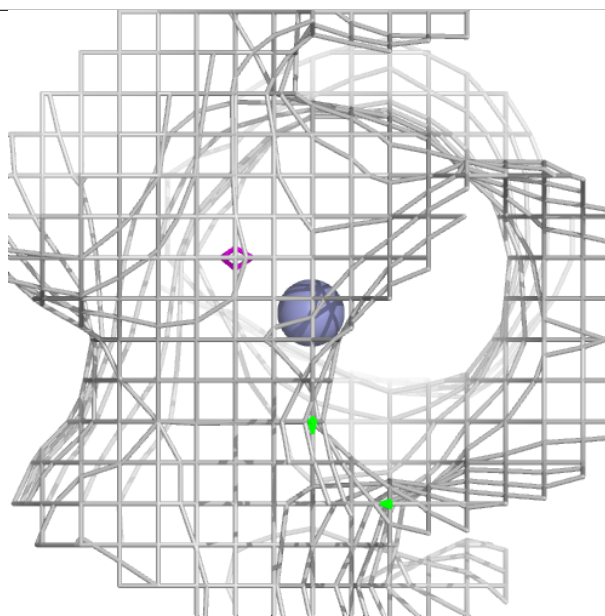
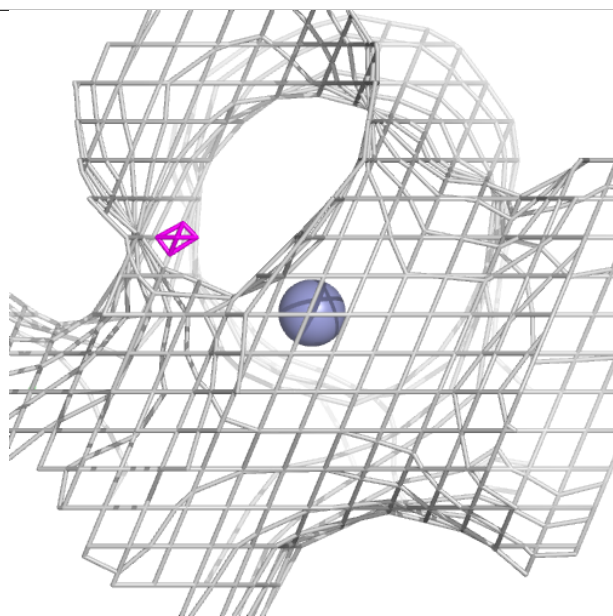
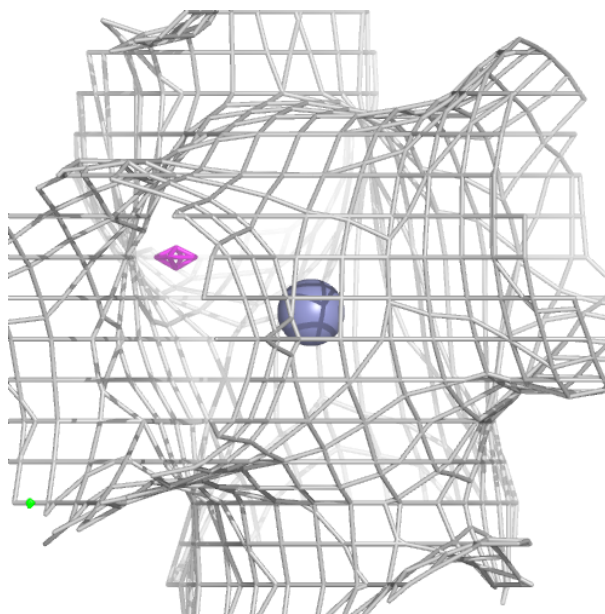
Electron density around ZN A 404:

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



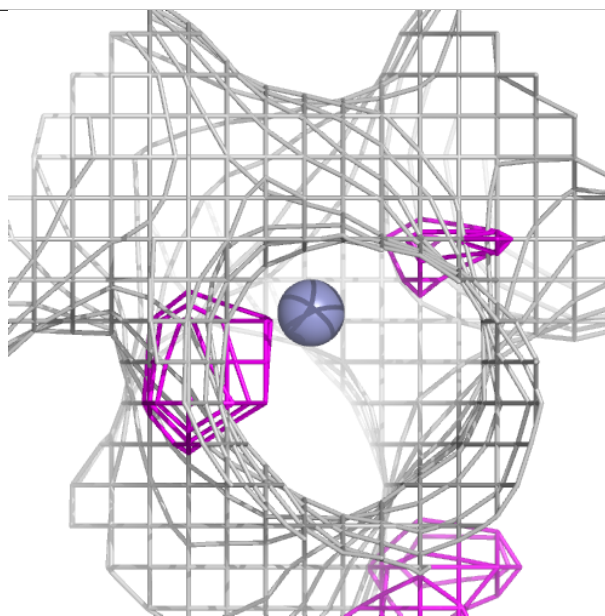
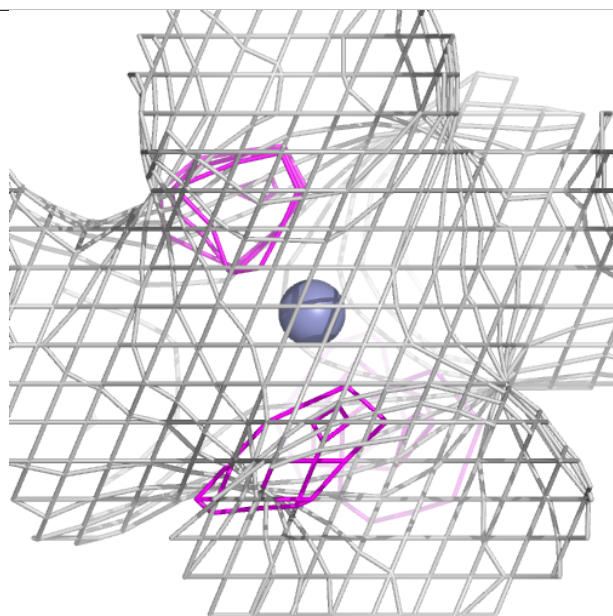
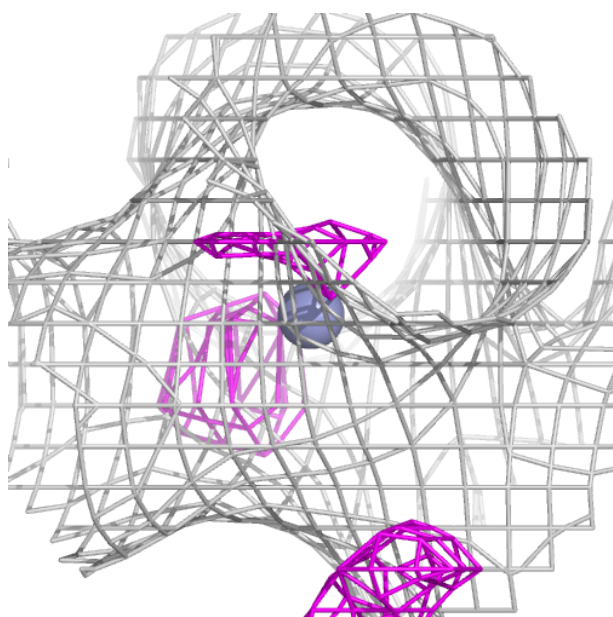
Electron density around ZN C 401:

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



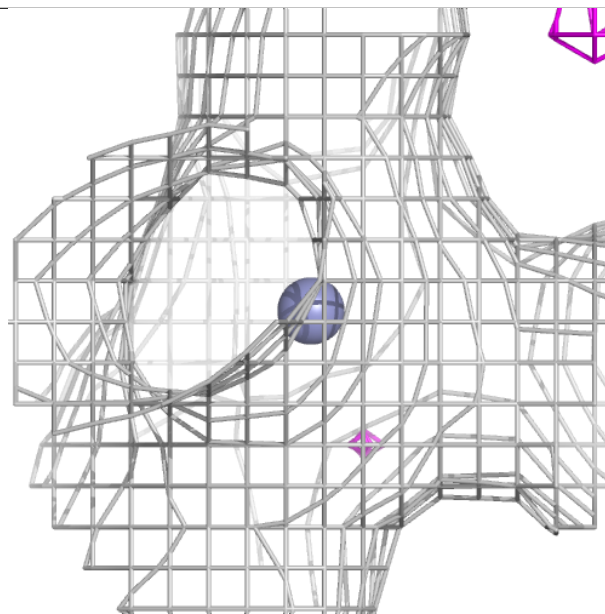
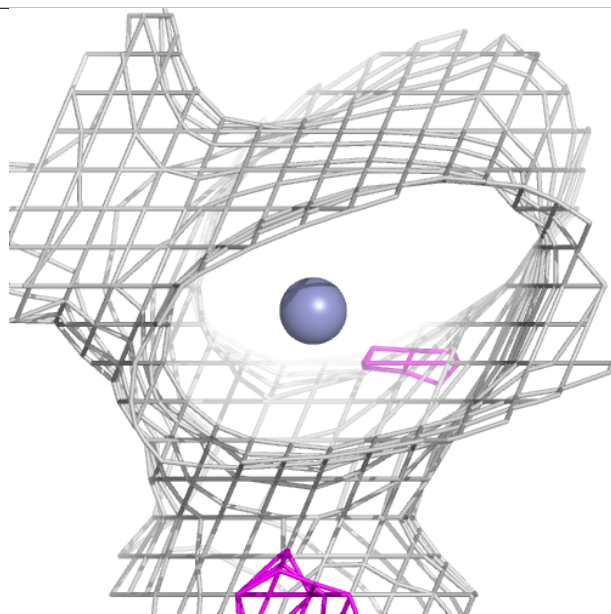
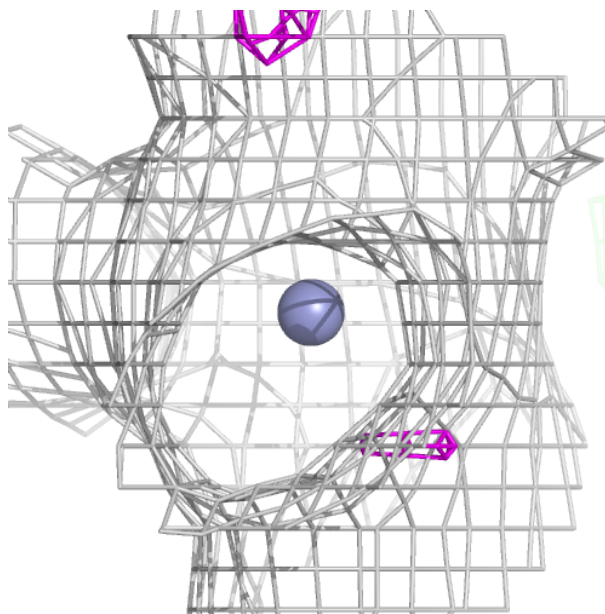
Electron density around ZN C 403:

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



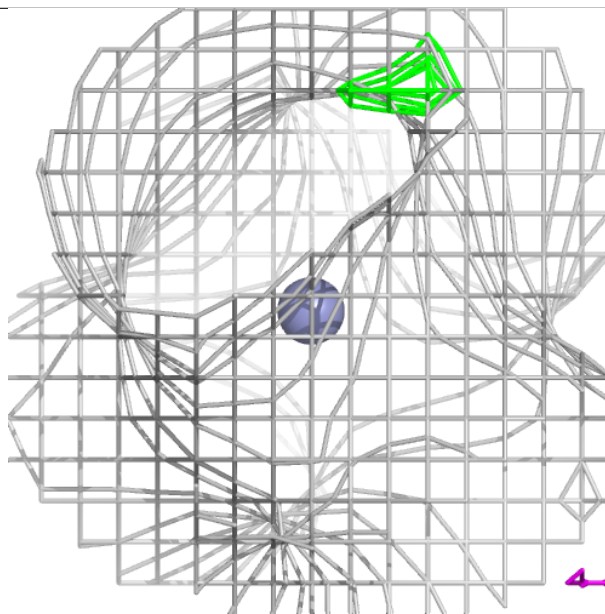
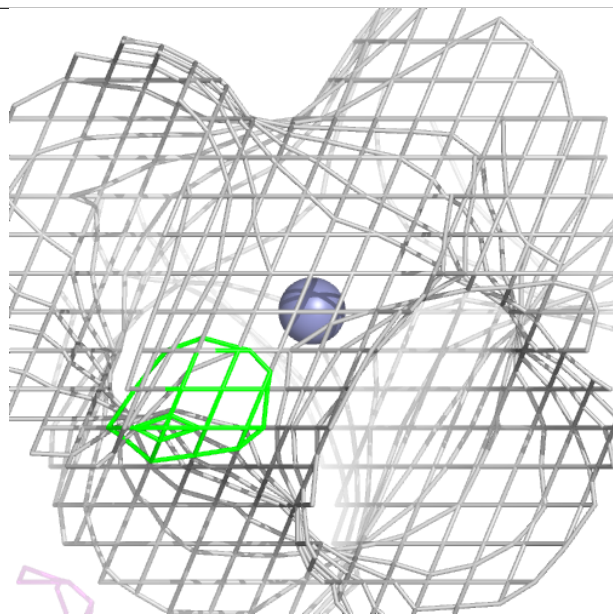
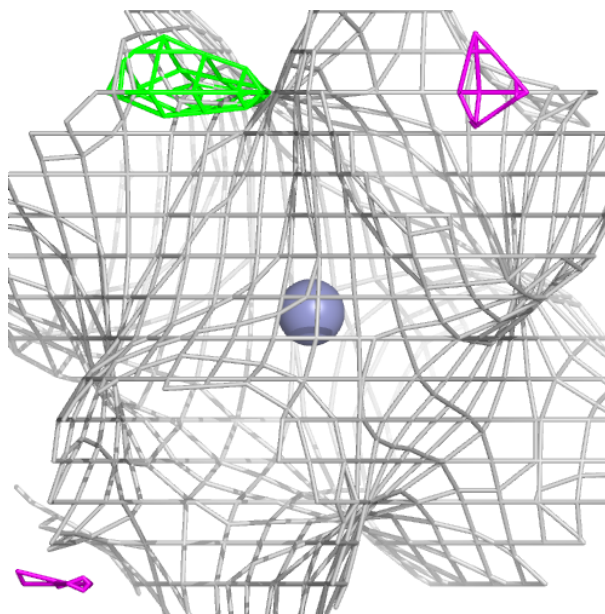
Electron density around ZN C 404:

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



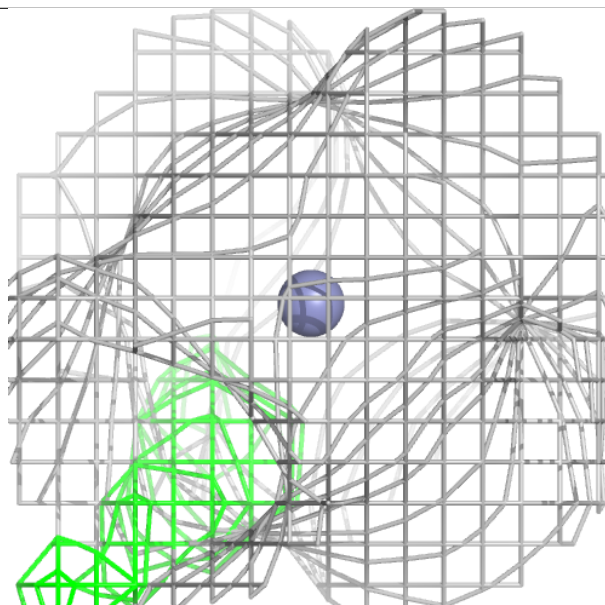
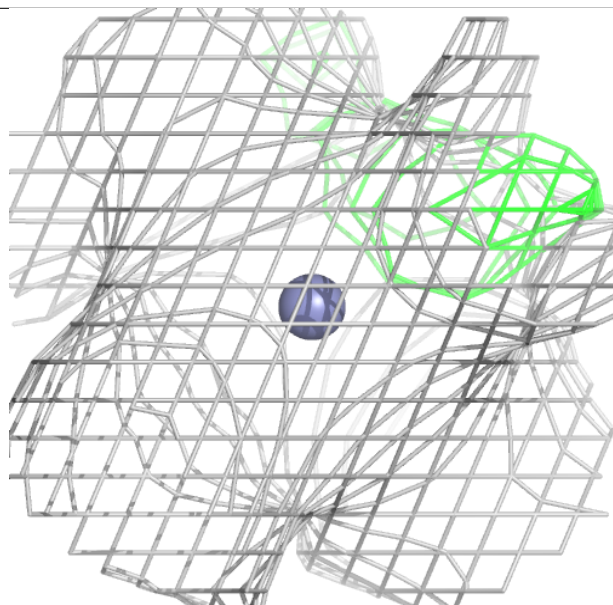
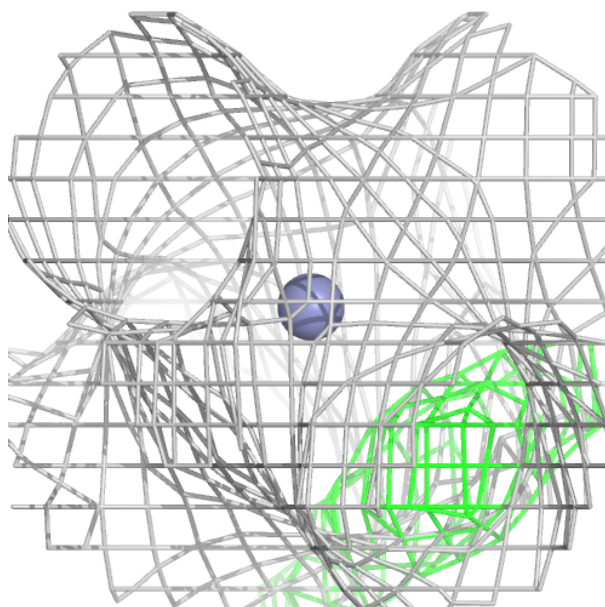
Electron density around ZN A 401:

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



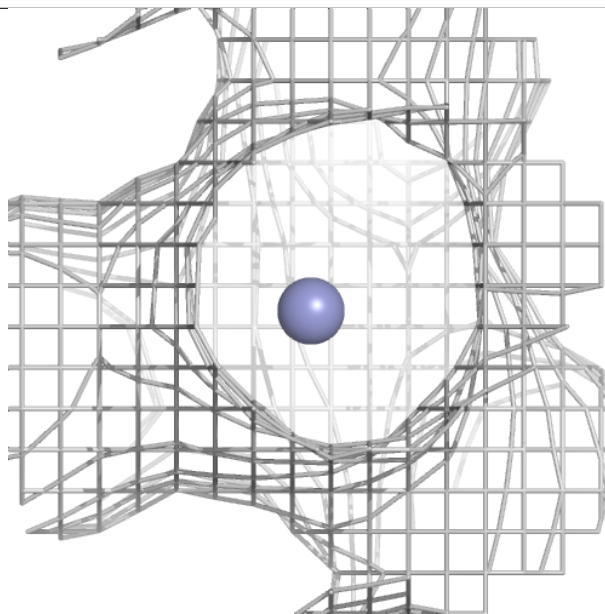
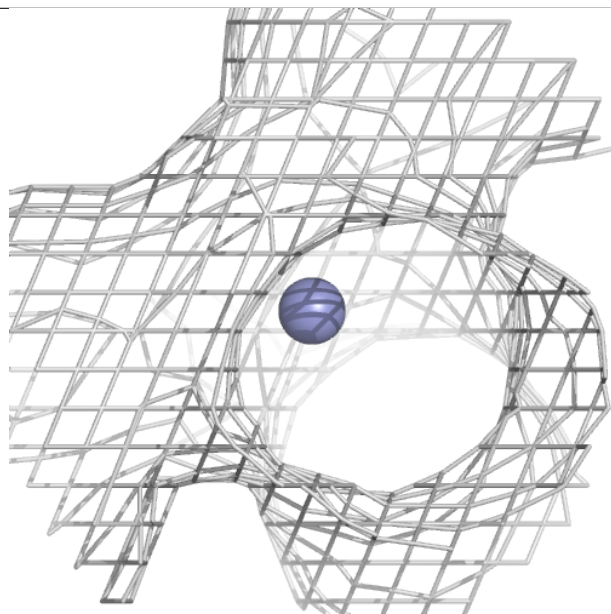
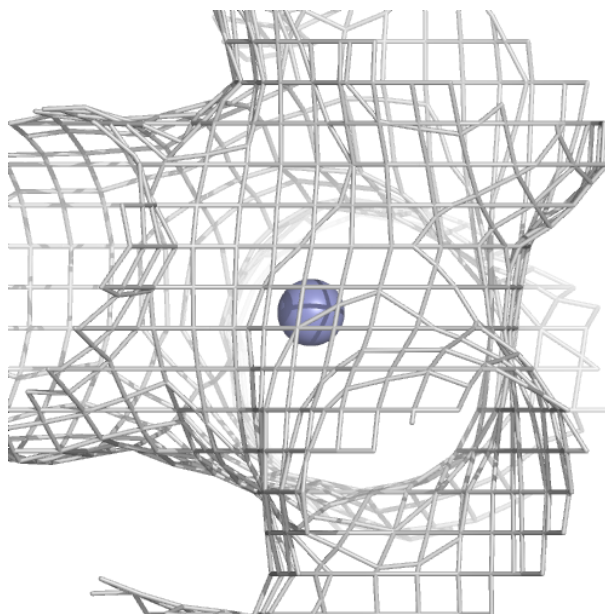
Electron density around ZN E 403:

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



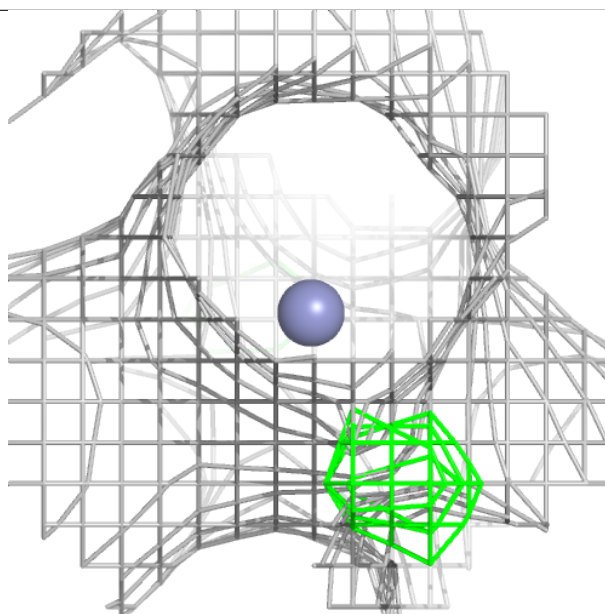
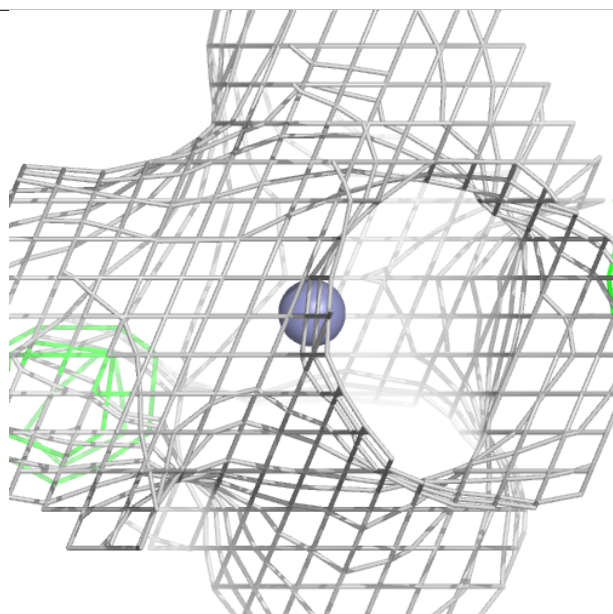
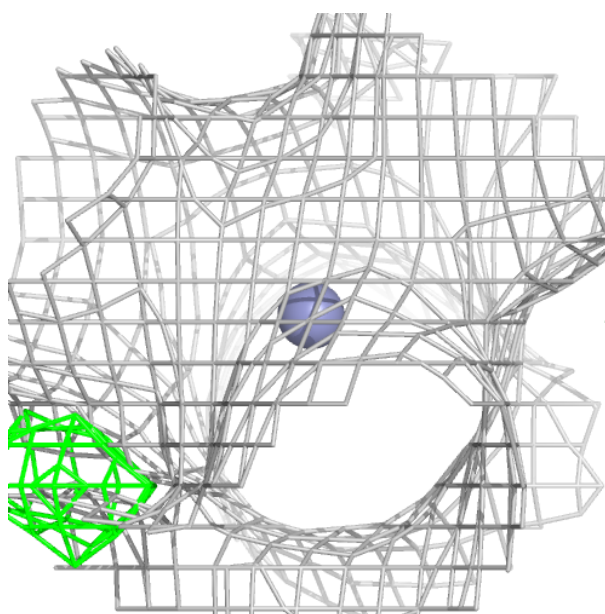
Electron density around ZN G 401:

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



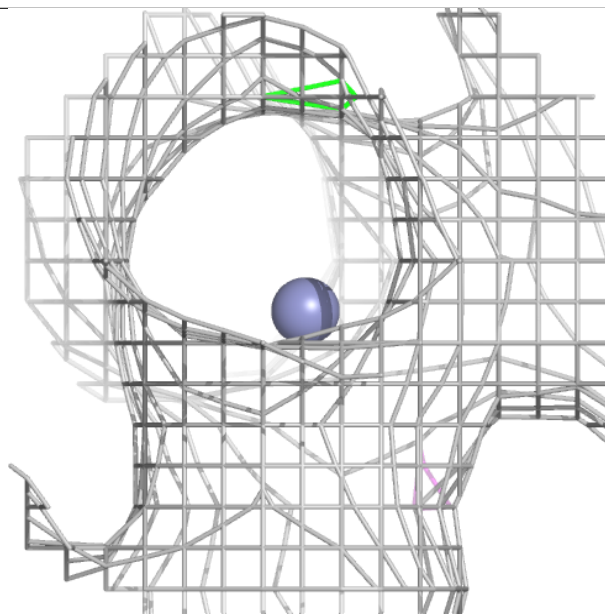
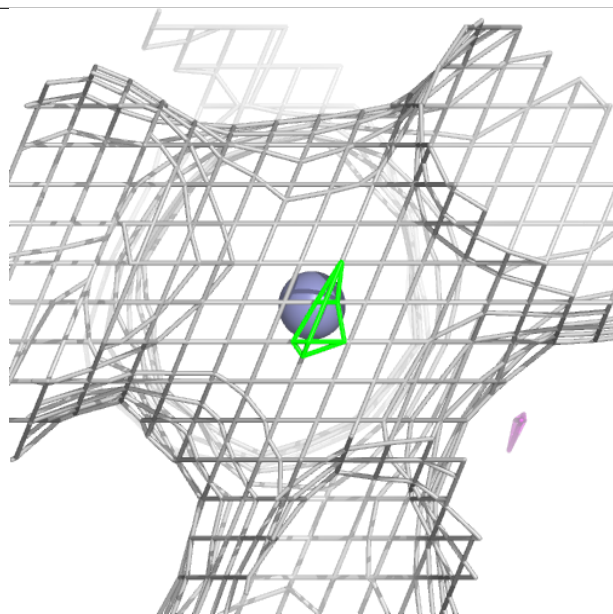
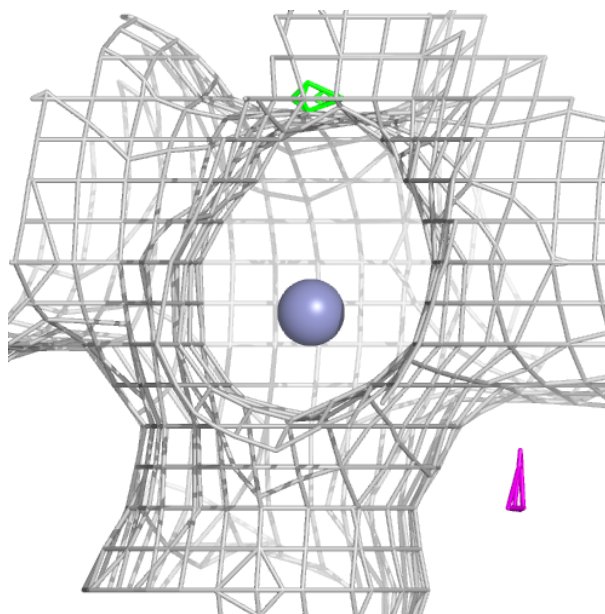
Electron density around ZN E 401:

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



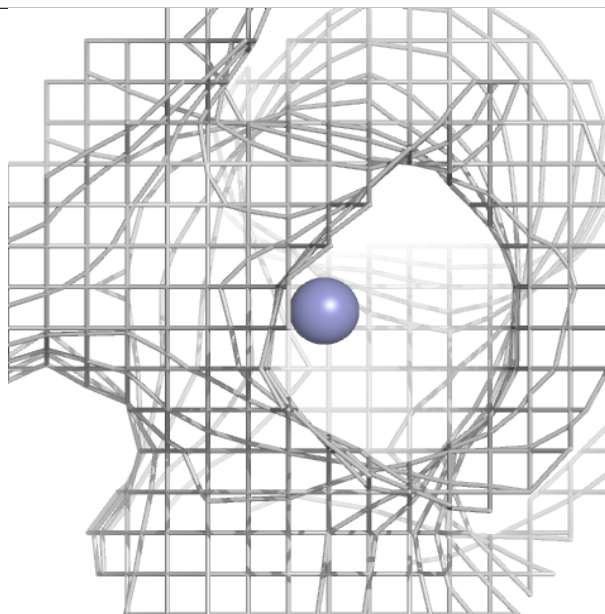
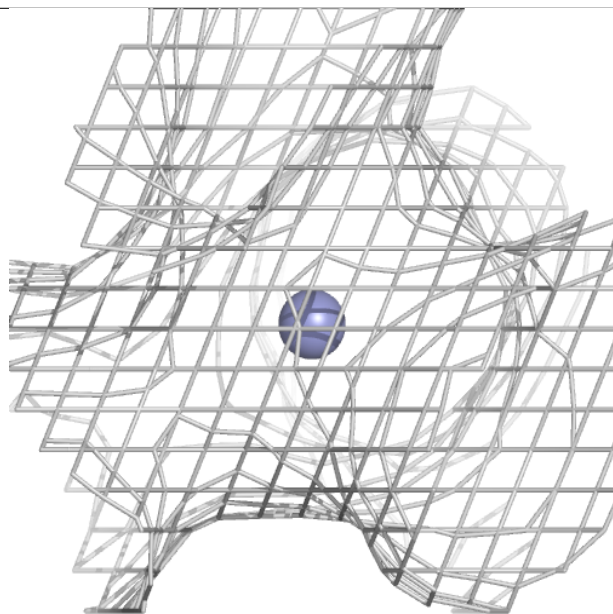
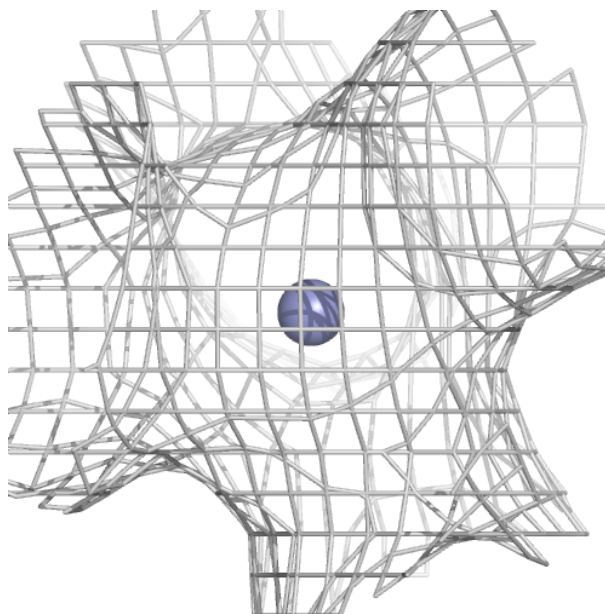
Electron density around ZN C 402:

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



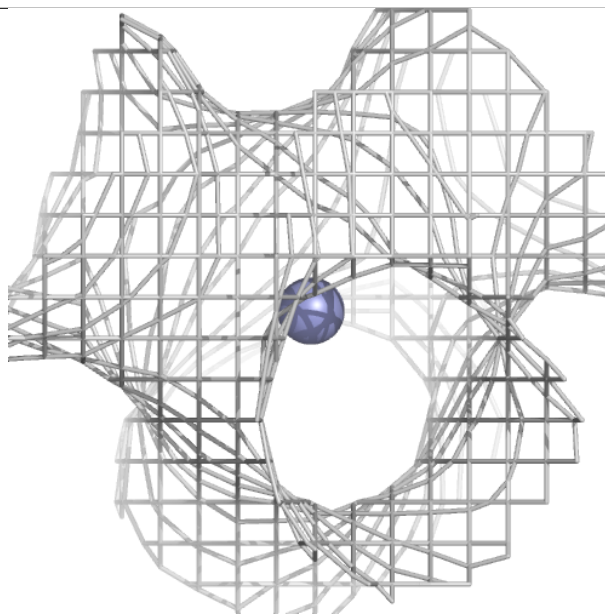
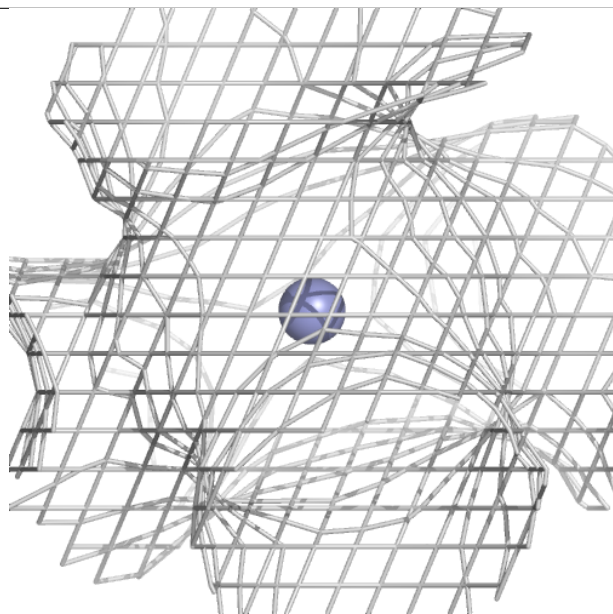
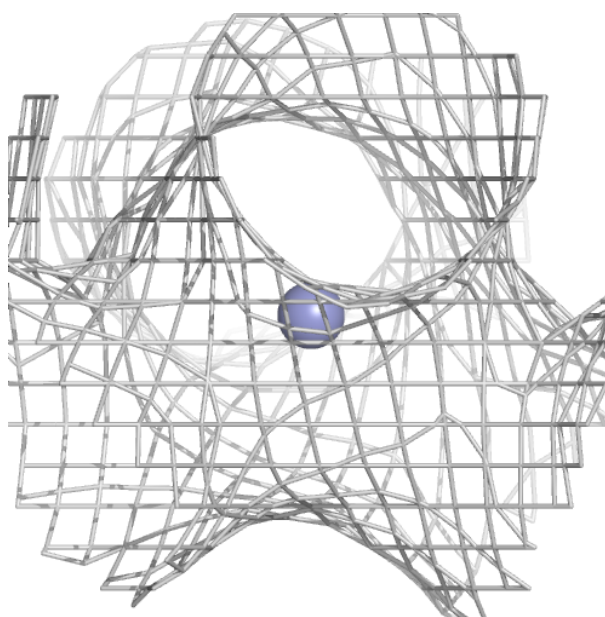
Electron density around ZN A 402:

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



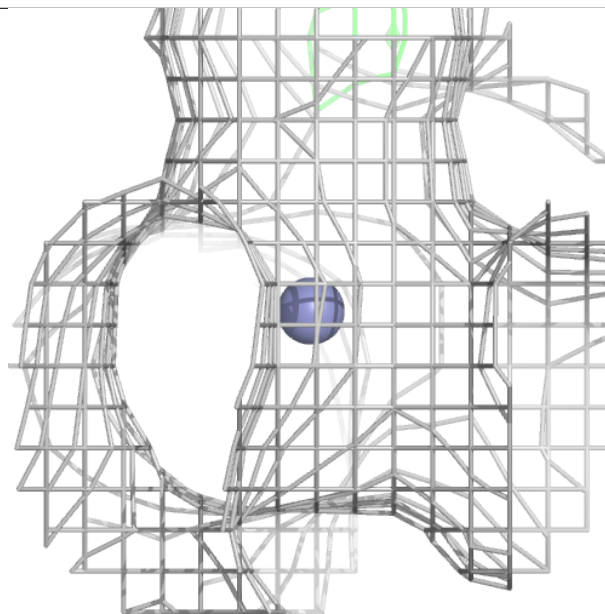
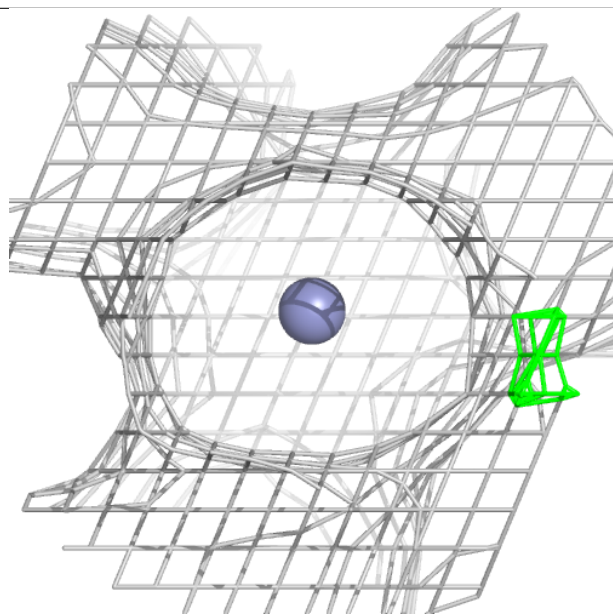
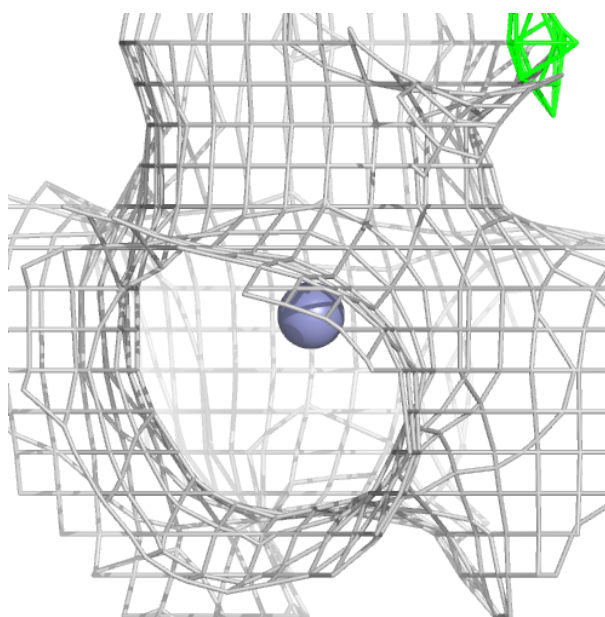
Electron density around ZN A 403:

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



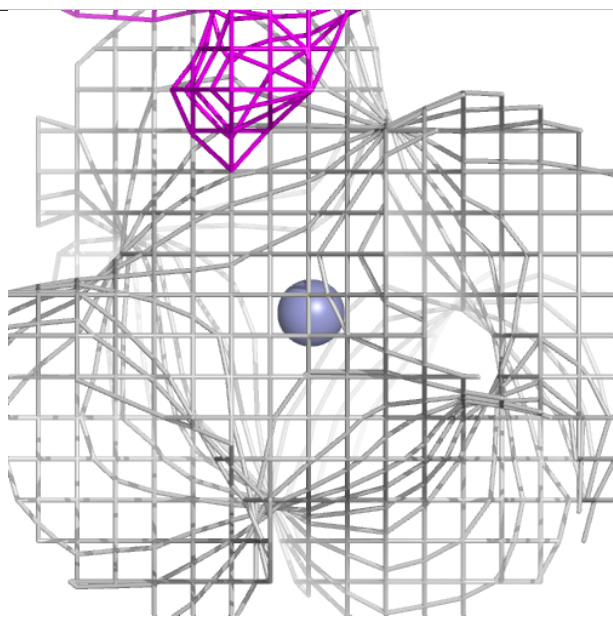
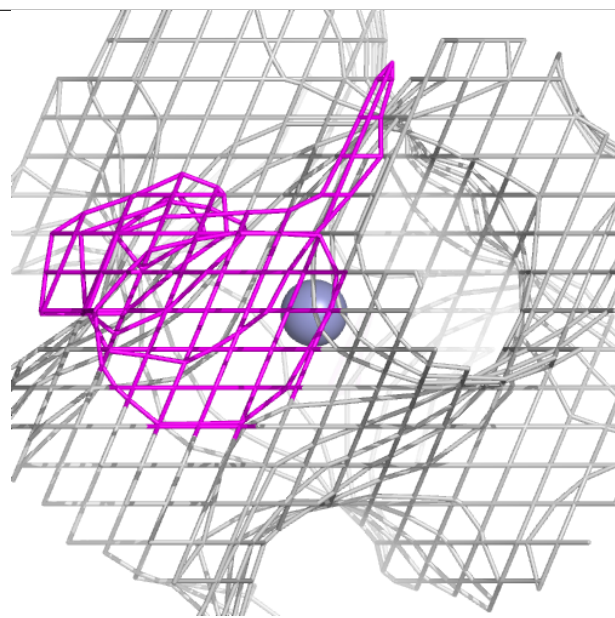
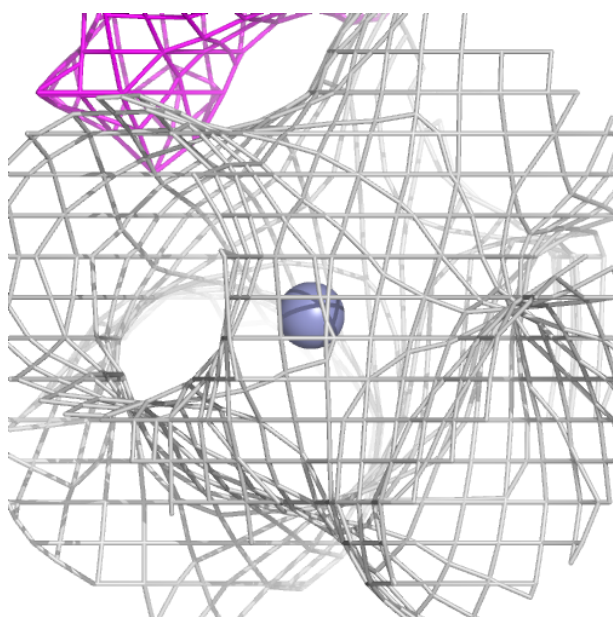
Electron density around ZN G 402:

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



Electron density around ZN G 403:

$2mF_o-DF_c$ (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.