



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 11:21 AM UTC

PDB ID : 10NH / pdb\_000010nh  
Title : Crystal Structure of Apurinic endonuclease (APN1) from Babesia bovis  
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)  
Deposited on : 2026-01-28  
Resolution : 2.65 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](mailto: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>.

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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  
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

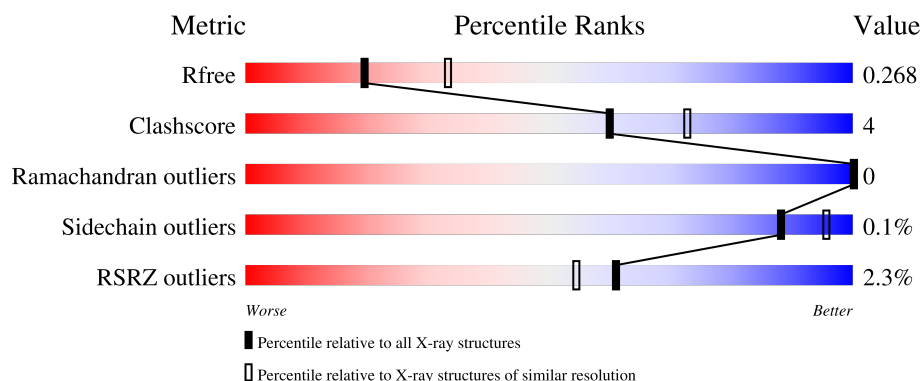
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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  $\geq 3$ , 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	323	
1	B	323	
1	C	323	
1	D	323	
1	E	323	

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Mol	Chain	Length	Quality of chain
1	F	323	<div><div></div><div>2%</div><div>87%</div><div>9%</div><div></div></div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14947 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 Apurinic endonuclease (APN1).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2460	1559	421	466	14			
1	B	313	Total	C	N	O	S	0	0	0
			2460	1559	421	466	14			
1	C	311	Total	C	N	O	S	0	0	0
			2448	1552	419	463	14			
1	D	313	Total	C	N	O	S	0	0	0
			2460	1559	421	466	14			
1	E	312	Total	C	N	O	S	0	0	0
			2453	1554	420	465	14			
1	F	312	Total	C	N	O	S	0	0	0
			2455	1557	420	464	14			

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	MET	-	initiating methionine	UNP S6B9Y3
A	46	ALA	-	expression tag	UNP S6B9Y3
A	47	HIS	-	expression tag	UNP S6B9Y3
A	48	HIS	-	expression tag	UNP S6B9Y3
A	49	HIS	-	expression tag	UNP S6B9Y3
A	50	HIS	-	expression tag	UNP S6B9Y3
A	51	HIS	-	expression tag	UNP S6B9Y3
A	52	HIS	-	expression tag	UNP S6B9Y3
A	256	VAL	GLY	conflict	UNP S6B9Y3
B	45	MET	-	initiating methionine	UNP S6B9Y3
B	46	ALA	-	expression tag	UNP S6B9Y3
B	47	HIS	-	expression tag	UNP S6B9Y3
B	48	HIS	-	expression tag	UNP S6B9Y3
B	49	HIS	-	expression tag	UNP S6B9Y3
B	50	HIS	-	expression tag	UNP S6B9Y3
B	51	HIS	-	expression tag	UNP S6B9Y3
B	52	HIS	-	expression tag	UNP S6B9Y3

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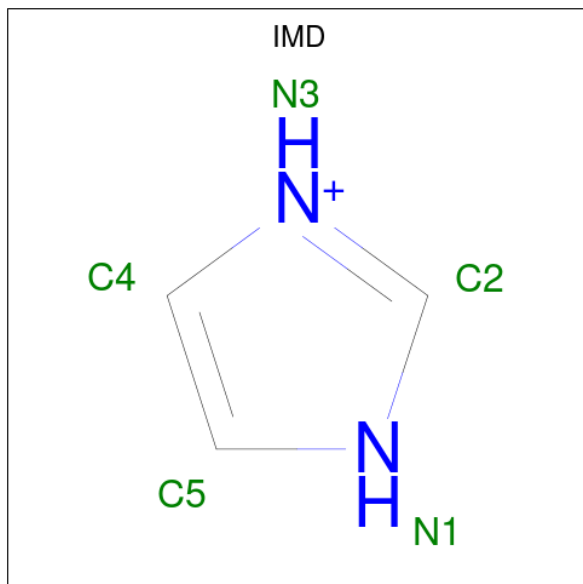
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Chain	Residue	Modelled	Actual	Comment	Reference
B	256	VAL	GLY	conflict	UNP S6B9Y3
C	45	MET	-	initiating methionine	UNP S6B9Y3
C	46	ALA	-	expression tag	UNP S6B9Y3
C	47	HIS	-	expression tag	UNP S6B9Y3
C	48	HIS	-	expression tag	UNP S6B9Y3
C	49	HIS	-	expression tag	UNP S6B9Y3
C	50	HIS	-	expression tag	UNP S6B9Y3
C	51	HIS	-	expression tag	UNP S6B9Y3
C	52	HIS	-	expression tag	UNP S6B9Y3
C	256	VAL	GLY	conflict	UNP S6B9Y3
D	45	MET	-	initiating methionine	UNP S6B9Y3
D	46	ALA	-	expression tag	UNP S6B9Y3
D	47	HIS	-	expression tag	UNP S6B9Y3
D	48	HIS	-	expression tag	UNP S6B9Y3
D	49	HIS	-	expression tag	UNP S6B9Y3
D	50	HIS	-	expression tag	UNP S6B9Y3
D	51	HIS	-	expression tag	UNP S6B9Y3
D	52	HIS	-	expression tag	UNP S6B9Y3
D	256	VAL	GLY	conflict	UNP S6B9Y3
E	45	MET	-	initiating methionine	UNP S6B9Y3
E	46	ALA	-	expression tag	UNP S6B9Y3
E	47	HIS	-	expression tag	UNP S6B9Y3
E	48	HIS	-	expression tag	UNP S6B9Y3
E	49	HIS	-	expression tag	UNP S6B9Y3
E	50	HIS	-	expression tag	UNP S6B9Y3
E	51	HIS	-	expression tag	UNP S6B9Y3
E	52	HIS	-	expression tag	UNP S6B9Y3
E	256	VAL	GLY	conflict	UNP S6B9Y3
F	45	MET	-	initiating methionine	UNP S6B9Y3
F	46	ALA	-	expression tag	UNP S6B9Y3
F	47	HIS	-	expression tag	UNP S6B9Y3
F	48	HIS	-	expression tag	UNP S6B9Y3
F	49	HIS	-	expression tag	UNP S6B9Y3
F	50	HIS	-	expression tag	UNP S6B9Y3
F	51	HIS	-	expression tag	UNP S6B9Y3
F	52	HIS	-	expression tag	UNP S6B9Y3
F	256	VAL	GLY	conflict	UNP S6B9Y3

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

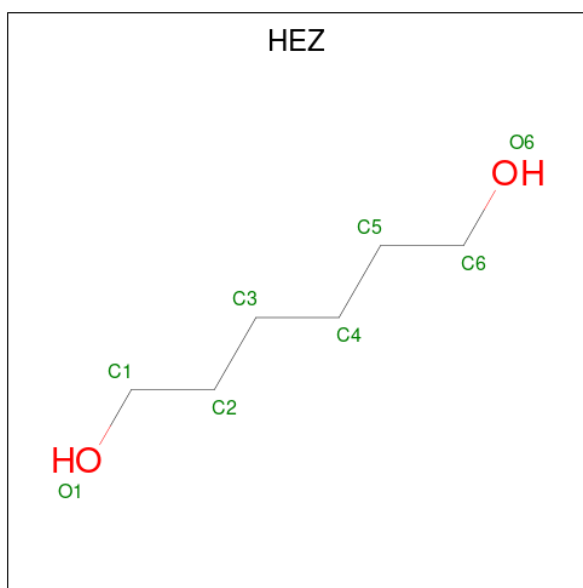
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total 3	Zn 3	0	0
2	B	3	Total 3	Zn 3	0	0
2	C	3	Total 3	Zn 3	0	0
2	D	3	Total 3	Zn 3	0	0
2	E	3	Total 3	Zn 3	0	0
2	F	3	Total 3	Zn 3	0	0

- Molecule 3 is IMIDAZOLE (CCD ID: IMD) (formula:  $C_3H_5N_2$ ).



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

- Molecule 4 is HEXANE-1,6-DIOL (CCD ID: HEZ) (formula:  $C_6H_{14}O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	F	1	Total	C	O	0	0
			8	6	2		

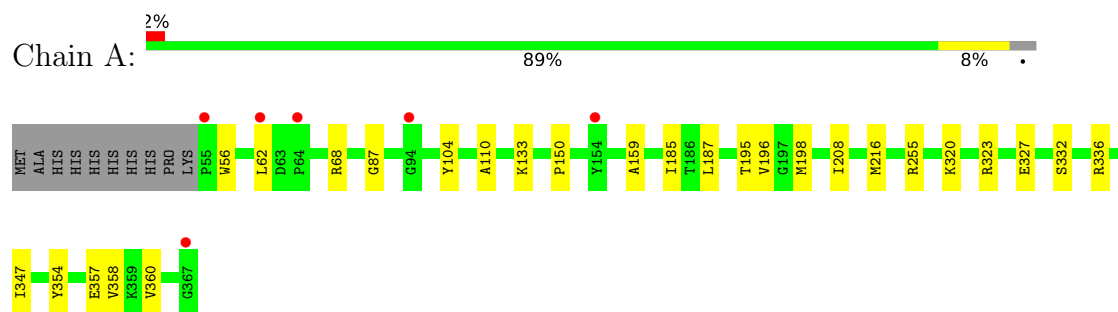
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	21	Total	O	0	0
			21	21		
5	B	25	Total	O	0	0
			25	25		
5	C	25	Total	O	0	0
			25	25		
5	D	23	Total	O	0	0
			23	23		
5	E	34	Total	O	0	0
			34	34		
5	F	27	Total	O	0	0
			27	27		

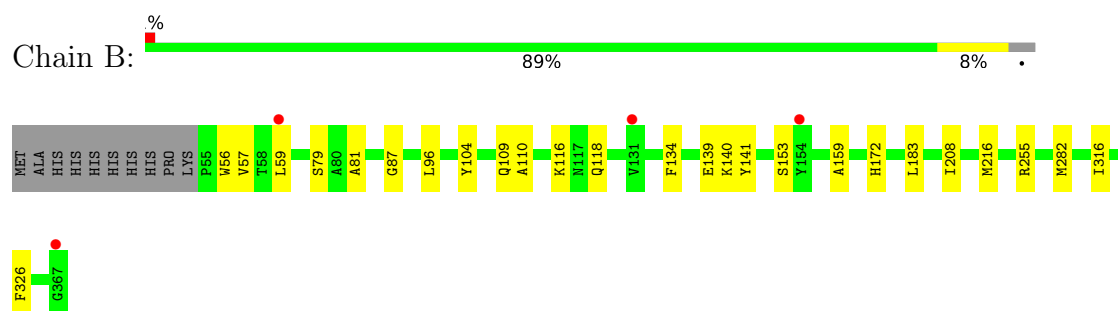
### 3 Residue-property plots [i](#)

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.

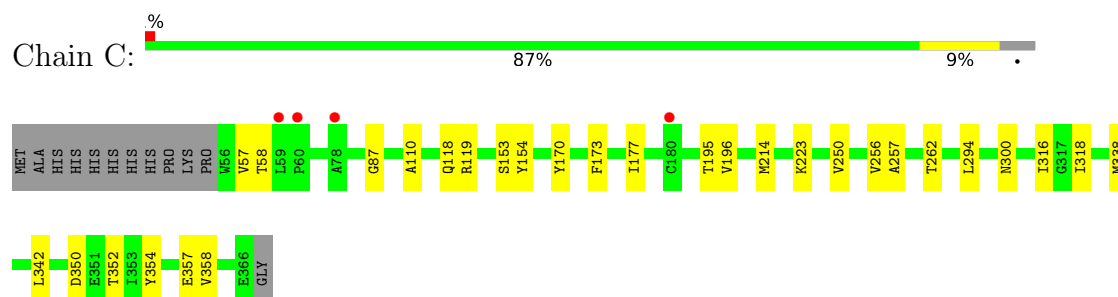
- Molecule 1: Apurinic endonuclease (APN1)



- Molecule 1: Apurinic endonuclease (APN1)



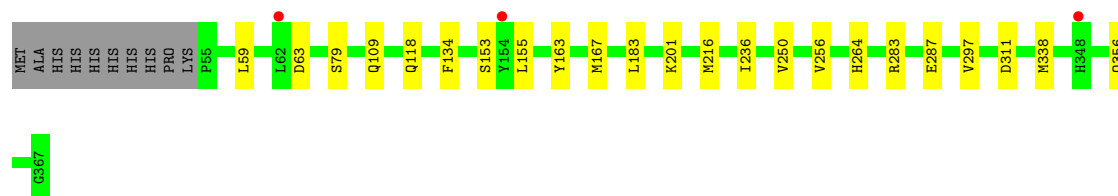
- Molecule 1: Apurinic endonuclease (APN1)



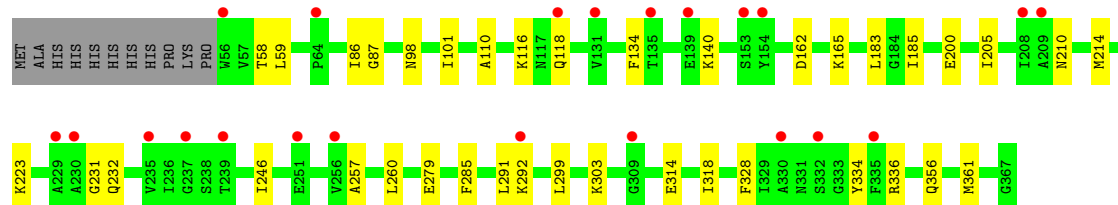
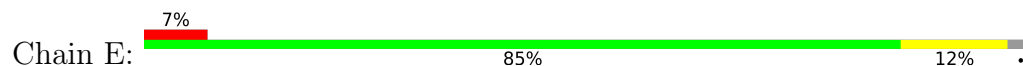
- Molecule 1: Apurinic endonuclease (APN1)



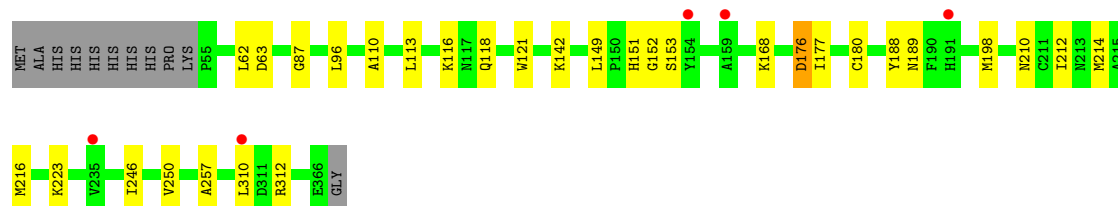
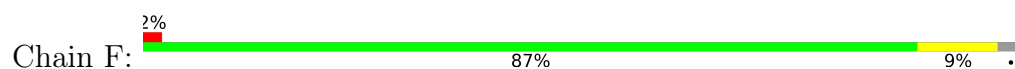




- Molecule 1: Apurinic endonuclease (APN1)



- Molecule 1: Apurinic endonuclease (APN1)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.62Å 93.09Å 102.88Å 91.66° 91.18° 91.09°	Depositor
Resolution (Å)	46.52 – 2.65 46.52 – 2.65	Depositor EDS
% Data completeness (in resolution range)	97.0 (46.52-2.65) 97.1 (46.52-2.65)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 2.65Å)	Xtriage
Refinement program	PHENIX (2.0_5936: ???)	Depositor
R, $R_{free}$	0.205 , 0.264 0.217 , 0.268	Depositor DCC
$R_{free}$ test set	3214 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	54.6	Xtriage
Anisotropy	0.342	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 51.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-l 0.015 for -h,k,-l 0.000 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	14947	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	69.0	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 48.69 % 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 8.2972e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: IMD, HEZ, 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	0.19	0/2510	0.38	0/3392
1	B	0.18	0/2510	0.36	0/3392
1	C	0.19	0/2497	0.36	0/3376
1	D	0.19	0/2510	0.35	0/3392
1	E	0.20	0/2502	0.40	0/3381
1	F	0.19	0/2505	0.36	0/3387
All	All	0.19	0/15034	0.37	0/20320

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	2460	0	2428	22	0
1	B	2460	0	2428	18	0
1	C	2448	0	2417	19	0
1	D	2460	0	2428	15	0
1	E	2453	0	2420	29	0
1	F	2455	0	2425	24	0
2	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	3	0	0	0	0
2	C	3	0	0	0	0
2	D	3	0	0	0	0
2	E	3	0	0	0	0
2	F	3	0	0	0	0
3	A	5	0	5	0	0
3	B	5	0	5	0	0
3	C	5	0	5	0	0
3	D	5	0	5	0	0
3	E	5	0	5	0	0
3	F	5	0	5	0	0
4	F	8	0	14	0	0
5	A	21	0	0	0	0
5	B	25	0	0	0	0
5	C	25	0	0	0	0
5	D	23	0	0	0	0
5	E	34	0	0	0	0
5	F	27	0	0	0	0
All	All	14947	0	14590	125	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 (125) 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:D:250:VAL:HG11	1:D:256:VAL:HG21	1.60	0.82
1:C:250:VAL:HG11	1:C:256:VAL:HG21	1.70	0.73
1:E:336:ARG:HH11	1:E:336:ARG:HG3	1.59	0.67
1:F:198:MET:HA	1:F:198:MET:HE2	1.76	0.66
1:F:310:LEU:HD22	1:F:312:ARG:HH22	1.62	0.65
1:C:316:ILE:HD11	1:C:342:LEU:HD21	1.79	0.64
1:F:310:LEU:HD22	1:F:312:ARG:NH2	2.12	0.64
1:D:297:VAL:HG23	1:D:338:MET:CE	2.27	0.64
1:D:250:VAL:HG11	1:D:256:VAL:CG2	2.28	0.63
1:C:316:ILE:HD11	1:C:342:LEU:CD2	2.29	0.62
1:D:297:VAL:HG23	1:D:338:MET:HE3	1.81	0.61
1:C:250:VAL:HG11	1:C:256:VAL:CG2	2.30	0.61
1:A:323:ARG:HD3	1:A:327:GLU:OE2	2.01	0.60
1:F:121:TRP:CE3	1:F:168:LYS:HE3	2.36	0.60
1:B:216:MET:HE2	1:B:255:ARG:CB	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:PHE:CD2	1:B:183:LEU:HD23	2.37	0.59
1:F:223:LYS:HD3	1:F:257:ALA:HB2	1.85	0.59
1:E:291:LEU:HD11	1:E:334:TYR:CE1	2.38	0.59
1:E:292:LYS:O	1:E:292:LYS:HD2	2.03	0.58
1:E:162:ASP:OD1	1:E:165:LYS:HG2	2.03	0.58
1:E:200:GLU:HA	1:E:200:GLU:OE2	2.03	0.58
1:C:294:LEU:HG	1:C:338:MET:HE1	1.86	0.57
1:E:299:LEU:HD21	1:E:361:MET:HE2	1.86	0.57
1:F:63:ASP:OD1	1:F:63:ASP:C	2.49	0.56
1:A:216:MET:HE2	1:A:255:ARG:CB	2.35	0.56
1:B:216:MET:HE2	1:B:255:ARG:HB2	1.90	0.54
1:C:154:TYR:O	1:C:154:TYR:CD2	2.61	0.54
1:C:350:ASP:OD1	1:C:352:THR:HG23	2.08	0.54
1:E:292:LYS:HD2	1:E:292:LYS:C	2.33	0.54
1:E:98:ASN:HA	1:E:101:ILE:HD12	1.90	0.53
1:B:96:LEU:HD23	1:B:183:LEU:HD21	1.91	0.52
1:E:118:GLN:N	1:E:118:GLN:CD	2.68	0.52
1:F:87:GLY:HA3	1:F:110:ALA:O	2.08	0.52
1:F:96:LEU:CD1	1:F:113:LEU:HD12	2.39	0.52
1:A:216:MET:HE2	1:A:255:ARG:HB2	1.92	0.52
1:E:303:LYS:HE2	1:E:314:GLU:OE1	2.10	0.52
1:C:223:LYS:HD3	1:C:257:ALA:HB2	1.93	0.51
1:E:86:ILE:HD13	1:E:361:MET:HB3	1.91	0.51
1:D:59:LEU:HD23	1:D:356:GLN:HB2	1.92	0.51
1:F:142:LYS:O	1:F:142:LYS:HG3	2.11	0.51
1:B:139:GLU:O	1:B:139:GLU:OE1	2.28	0.51
1:D:163:TYR:CE1	1:D:167:MET:HE3	2.45	0.51
1:F:96:LEU:HD12	1:F:96:LEU:O	2.10	0.50
1:A:68:ARG:H	1:A:68:ARG:CD	2.25	0.50
1:A:68:ARG:H	1:A:68:ARG:HD3	1.76	0.50
1:B:316:ILE:HD12	1:B:326:PHE:CE1	2.47	0.50
1:F:142:LYS:O	1:F:142:LYS:CG	2.60	0.50
1:B:216:MET:HE2	1:B:255:ARG:HB3	1.94	0.49
1:C:118:GLN:OE1	1:C:153:SER:HA	2.14	0.48
1:F:310:LEU:HD13	1:F:312:ARG:HH21	1.78	0.48
1:E:134:PHE:CD2	1:E:183:LEU:HD22	2.48	0.48
1:E:279:GLU:HA	1:E:328:PHE:CZ	2.50	0.47
1:A:347:ILE:HD12	1:A:347:ILE:H	1.78	0.47
1:C:195:THR:O	1:C:196:VAL:HG22	2.15	0.47
1:E:183:LEU:HB2	1:E:185:ILE:HD12	1.97	0.47
1:A:198:MET:HA	1:A:198:MET:HE3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:GLN:OE1	1:B:153:SER:HA	2.15	0.47
1:D:63:ASP:OD2	1:D:63:ASP:C	2.57	0.46
1:E:231:GLY:N	1:E:232:GLN:OE1	2.48	0.46
1:A:336:ARG:NH2	1:C:119:ARG:CZ	2.78	0.46
1:F:246:ILE:O	1:F:250:VAL:HG23	2.16	0.46
1:A:87:GLY:HA3	1:A:110:ALA:O	2.16	0.46
1:A:216:MET:HE2	1:A:255:ARG:HB3	1.97	0.46
1:E:87:GLY:HA3	1:E:110:ALA:O	2.16	0.46
1:F:210:ASN:O	1:F:214:MET:HG3	2.16	0.46
1:E:232:GLN:OE1	1:E:232:GLN:N	2.48	0.46
1:A:150:PRO:CG	1:A:185:ILE:HG21	2.46	0.46
1:B:59:LEU:H	1:B:59:LEU:HD12	1.81	0.45
1:E:291:LEU:HD11	1:E:334:TYR:CD1	2.51	0.45
1:B:282:MET:HE2	1:B:282:MET:HA	1.99	0.45
1:B:87:GLY:HA3	1:B:110:ALA:O	2.17	0.45
1:F:116:LYS:HE2	1:F:176:ASP:OD1	2.17	0.45
1:A:195:THR:O	1:A:196:VAL:HG22	2.17	0.44
1:B:56:TRP:CE3	1:B:57:VAL:HB	2.52	0.44
1:E:59:LEU:HB3	1:E:356:GLN:HE21	1.82	0.44
1:C:316:ILE:HD12	1:C:357:GLU:HB3	1.98	0.44
1:E:223:LYS:HD3	1:E:257:ALA:HB2	1.99	0.44
1:A:332:SER:O	1:A:336:ARG:NH1	2.51	0.44
1:C:87:GLY:HA3	1:C:110:ALA:O	2.17	0.44
1:F:176:ASP:HB3	1:F:188:TYR:CE1	2.53	0.44
1:A:320:LYS:HB3	1:A:320:LYS:HE3	1.86	0.44
1:A:159:ALA:HB2	1:A:208:ILE:HD11	1.99	0.44
1:C:318:ILE:HD12	1:C:318:ILE:N	2.32	0.44
1:E:140:LYS:N	1:E:140:LYS:HD3	2.33	0.43
1:A:150:PRO:HD2	1:A:187:LEU:O	2.17	0.43
1:F:142:LYS:O	1:F:142:LYS:HD2	2.18	0.43
1:F:149:LEU:HD11	1:F:189:ASN:HB2	2.01	0.43
1:E:291:LEU:HD11	1:E:334:TYR:HE1	1.80	0.43
1:B:159:ALA:HB2	1:B:208:ILE:HD11	2.01	0.43
1:C:262:THR:OG1	1:C:300:ASN:O	2.37	0.43
1:F:212:ILE:O	1:F:216:MET:HG3	2.19	0.43
1:A:104:TYR:CD1	1:A:104:TYR:C	2.97	0.43
1:E:205:ILE:HG23	1:E:246:ILE:HG12	2.00	0.43
1:C:57:VAL:HG22	1:C:58:THR:N	2.34	0.43
1:D:264:HIS:NE2	1:D:311:ASP:OD2	2.52	0.43
1:D:201:LYS:HG2	1:D:236:ILE:HG12	2.01	0.42
1:E:210:ASN:O	1:E:214:MET:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:LYS:HE3	1:B:172:HIS:CE1	2.54	0.42
1:D:79:SER:HB2	1:D:109:GLN:HG3	2.00	0.42
1:C:173:PHE:CZ	1:C:177:ILE:HD11	2.54	0.42
1:D:134:PHE:CD2	1:D:183:LEU:HD22	2.55	0.42
1:A:56:TRP:CE3	1:A:323:ARG:HG3	2.55	0.42
1:D:283:ARG:HH21	1:D:287:GLU:HG3	1.85	0.42
1:A:354:TYR:O	1:A:358:VAL:HG23	2.20	0.42
1:B:104:TYR:CD1	1:B:141:TYR:HD2	2.38	0.42
1:B:140:LYS:HE3	1:B:141:TYR:CE1	2.55	0.41
1:E:318:ILE:O	1:E:318:ILE:HG22	2.19	0.41
1:A:133:LYS:HD2	1:A:133:LYS:HA	1.89	0.41
1:A:62:LEU:H	1:A:62:LEU:HD23	1.85	0.41
1:B:79:SER:HB2	1:B:109:GLN:HG3	2.03	0.41
1:E:116:LYS:O	1:E:118:GLN:OE1	2.38	0.41
1:A:357:GLU:O	1:A:360:VAL:HG22	2.21	0.41
1:B:81:ALA:CB	1:D:155:LEU:HD21	2.50	0.41
1:E:58:THR:O	1:E:58:THR:HG23	2.20	0.41
1:F:62:LEU:HD13	1:F:63:ASP:N	2.35	0.41
1:F:177:ILE:O	1:F:180:CYS:HB3	2.20	0.41
1:C:354:TYR:O	1:C:358:VAL:HG23	2.20	0.41
1:E:260:LEU:HD11	1:E:285:PHE:CE2	2.56	0.41
1:D:216:MET:HE1	1:D:256:VAL:HG23	2.02	0.40
1:E:303:LYS:CE	1:E:314:GLU:OE1	2.68	0.40
1:F:118:GLN:OE1	1:F:153:SER:HA	2.21	0.40
1:F:151:HIS:HD1	1:F:152:GLY:N	2.20	0.40
1:D:118:GLN:OE1	1:D:153:SER:HA	2.21	0.40
1:C:170:TYR:OH	1:C:214:MET:HG2	2.22	0.40
1:F:168:LYS:HD2	1:F:168:LYS:O	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	311/323 (96%)	304 (98%)	7 (2%)	0	100	100
1	B	311/323 (96%)	301 (97%)	10 (3%)	0	100	100
1	C	309/323 (96%)	299 (97%)	10 (3%)	0	100	100
1	D	311/323 (96%)	306 (98%)	5 (2%)	0	100	100
1	E	310/323 (96%)	301 (97%)	9 (3%)	0	100	100
1	F	310/323 (96%)	303 (98%)	7 (2%)	0	100	100
All	All	1862/1938 (96%)	1814 (97%)	48 (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	263/272 (97%)	263 (100%)	0	100	100
1	B	263/272 (97%)	263 (100%)	0	100	100
1	C	262/272 (96%)	262 (100%)	0	100	100
1	D	263/272 (97%)	263 (100%)	0	100	100
1	E	262/272 (96%)	262 (100%)	0	100	100
1	F	263/272 (97%)	262 (100%)	1 (0%)	84	93
All	All	1576/1632 (97%)	1575 (100%)	1 (0%)	88	95

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	176	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	GLN

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Mol	Chain	Res	Type
1	A	337	ASN
1	B	172	HIS
1	B	356	GLN
1	C	234	ASN
1	C	315	ASN
1	D	89	HIS
1	D	102	ASN
1	D	234	ASN
1	D	315	ASN
1	D	337	ASN
1	E	117	ASN
1	E	118	GLN
1	E	137	ASN
1	E	298	HIS
1	E	356	GLN
1	F	89	HIS

### 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 25 ligands modelled in this entry, 18 are monoatomic - leaving 7 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
3	IMD	E	404	2	5,5,5	0.40	0	5,5,5	0.89	0
3	IMD	A	404	2	5,5,5	0.34	0	5,5,5	0.91	0
3	IMD	F	405	2	5,5,5	0.40	0	5,5,5	0.96	0
3	IMD	B	404	2	5,5,5	0.38	0	5,5,5	0.90	0
3	IMD	C	404	2	5,5,5	0.42	0	5,5,5	0.92	0
3	IMD	D	404	2	5,5,5	0.46	0	5,5,5	0.93	0
4	HEZ	F	404	-	7,7,7	0.16	0	6,6,6	0.37	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	IMD	E	404	2	-	-	0/1/1/1
3	IMD	A	404	2	-	-	0/1/1/1
3	IMD	F	405	2	-	-	0/1/1/1
3	IMD	B	404	2	-	-	0/1/1/1
3	IMD	C	404	2	-	-	0/1/1/1
3	IMD	D	404	2	-	-	0/1/1/1
4	HEZ	F	404	-	-	2/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	404	HEZ	C2-C3-C4-C5
4	F	404	HEZ	C3-C4-C5-C6

There are no ring outliers.

No monomer is involved in short contacts.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	313/323 (96%)	0.07	6 (1%)	66 61	40, 57, 92, 140	0
1	B	313/323 (96%)	0.10	4 (1%)	75 71	46, 65, 98, 142	0
1	C	311/323 (96%)	0.27	4 (1%)	75 71	44, 65, 99, 159	0
1	D	313/323 (96%)	0.19	3 (0%)	79 76	42, 63, 96, 181	0
1	E	312/323 (96%)	0.76	22 (7%)	22 16	49, 77, 110, 183	0
1	F	312/323 (96%)	0.25	5 (1%)	70 66	44, 66, 94, 145	0
All	All	1874/1938 (96%)	0.27	44 (2%)	61 54	40, 65, 101, 183	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	154	TYR	4.1
1	B	154	TYR	3.6
1	A	367	GLY	3.5
1	D	154	TYR	3.5
1	A	62	LEU	3.4
1	A	55	PRO	3.3
1	C	59	LEU	3.2
1	E	230	ALA	3.2
1	A	154	TYR	3.1
1	E	153	SER	3.0
1	E	239	THR	2.9
1	A	94	GLY	2.9
1	E	309	GLY	2.9
1	E	131	VAL	2.8
1	E	256	VAL	2.7
1	E	237	GLY	2.7
1	E	209	ALA	2.7
1	F	154	TYR	2.6
1	F	191	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	E	56	TRP	2.5
1	E	208	ILE	2.5
1	C	60	PRO	2.5
1	A	64	PRO	2.4
1	E	229	ALA	2.4
1	E	118	GLN	2.4
1	E	292	LYS	2.3
1	F	310	LEU	2.3
1	D	62	LEU	2.3
1	E	139	GLU	2.2
1	C	78	ALA	2.2
1	E	330	ALA	2.2
1	C	180	CYS	2.2
1	B	131	VAL	2.2
1	E	135	THR	2.2
1	E	64	PRO	2.1
1	E	251	GLU	2.1
1	B	59	LEU	2.1
1	B	367	GLY	2.1
1	E	335	PHE	2.1
1	D	348	HIS	2.0
1	F	235	VAL	2.0
1	F	159	ALA	2.0
1	E	332	SER	2.0
1	E	235	VAL	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

## 6.4 Ligands ⓘ

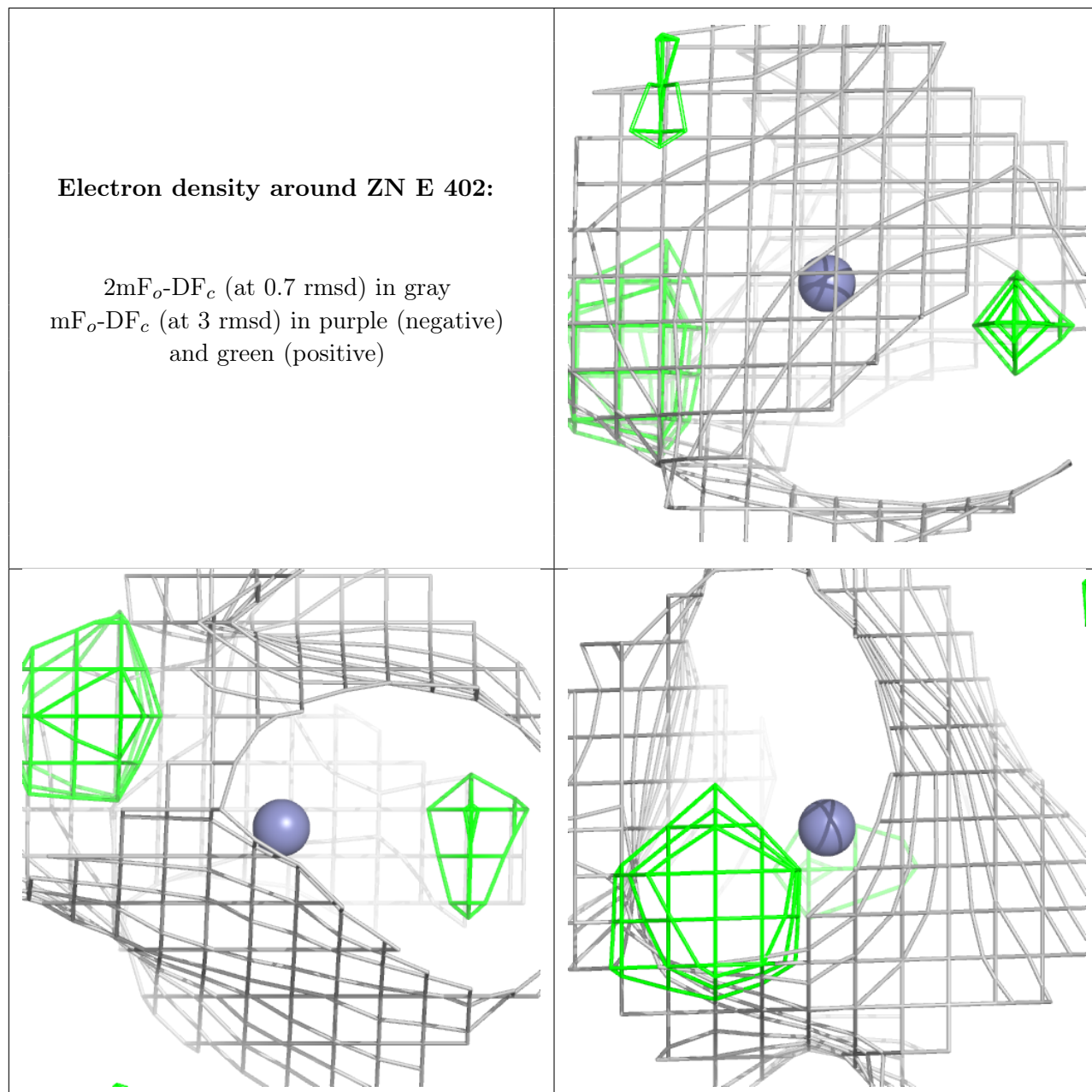
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
4	HEZ	F	404	8/8	0.73	0.17	68,76,82,89	0
3	IMD	E	404	5/5	0.77	0.15	57,65,86,87	0
3	IMD	F	405	5/5	0.84	0.17	50,50,68,85	0
3	IMD	D	404	5/5	0.84	0.13	37,48,51,78	0
3	IMD	A	404	5/5	0.94	0.12	50,57,71,76	0
3	IMD	B	404	5/5	0.94	0.11	43,53,74,86	0
2	ZN	E	402	1/1	0.94	0.06	86,86,86,86	0
2	ZN	E	401	1/1	0.95	0.06	83,83,83,83	0
3	IMD	C	404	5/5	0.95	0.12	39,43,60,67	0
2	ZN	A	402	1/1	0.96	0.09	73,73,73,73	0
2	ZN	F	402	1/1	0.97	0.05	65,65,65,65	0
2	ZN	E	403	1/1	0.97	0.05	74,74,74,74	0
2	ZN	C	402	1/1	0.98	0.04	64,64,64,64	0
2	ZN	D	401	1/1	0.98	0.03	69,69,69,69	0
2	ZN	D	402	1/1	0.98	0.04	54,54,54,54	0
2	ZN	D	403	1/1	0.98	0.05	66,66,66,66	0
2	ZN	A	401	1/1	0.98	0.04	60,60,60,60	0
2	ZN	B	401	1/1	0.98	0.06	65,65,65,65	0
2	ZN	C	401	1/1	0.98	0.03	69,69,69,69	0
2	ZN	F	401	1/1	0.98	0.05	63,63,63,63	0
2	ZN	B	402	1/1	0.99	0.03	57,57,57,57	0
2	ZN	C	403	1/1	0.99	0.02	55,55,55,55	0
2	ZN	F	403	1/1	0.99	0.03	57,57,57,57	0
2	ZN	B	403	1/1	0.99	0.03	59,59,59,59	0
2	ZN	A	403	1/1	0.99	0.03	58,58,58,58	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.

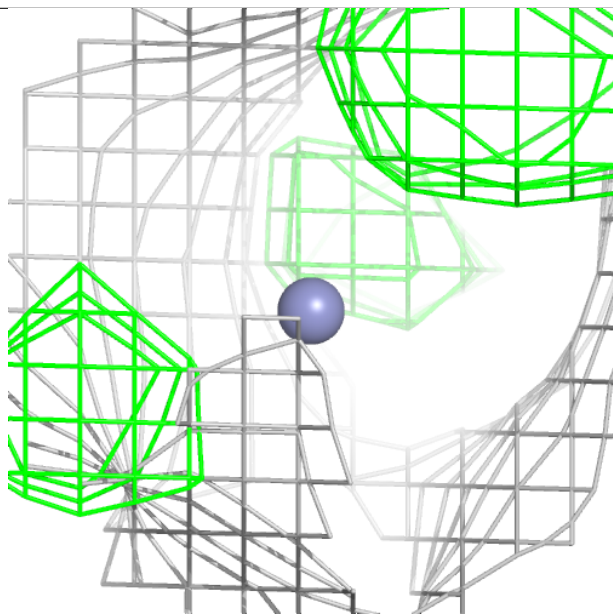
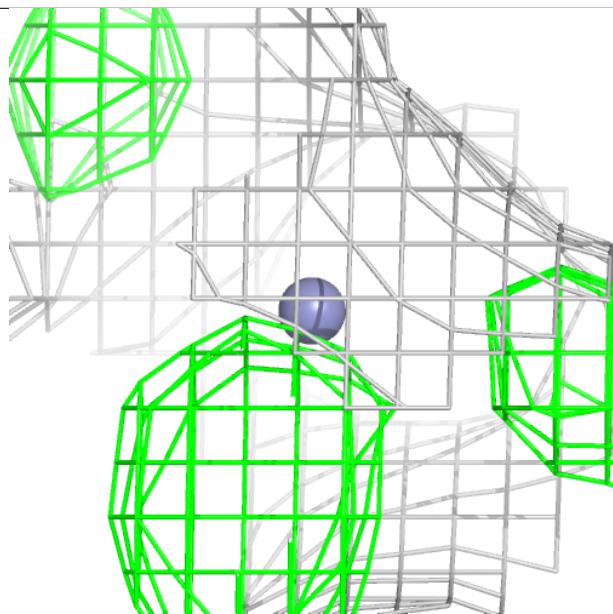
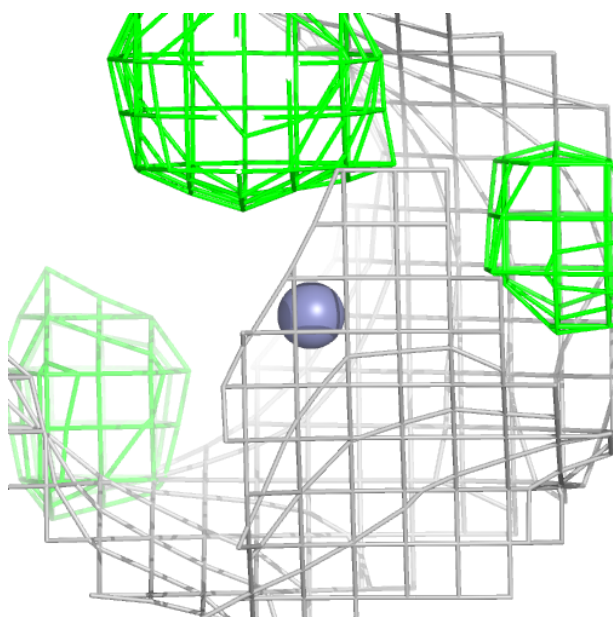
**Electron density around ZN E 402:**

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



**Electron density around ZN E 401:**

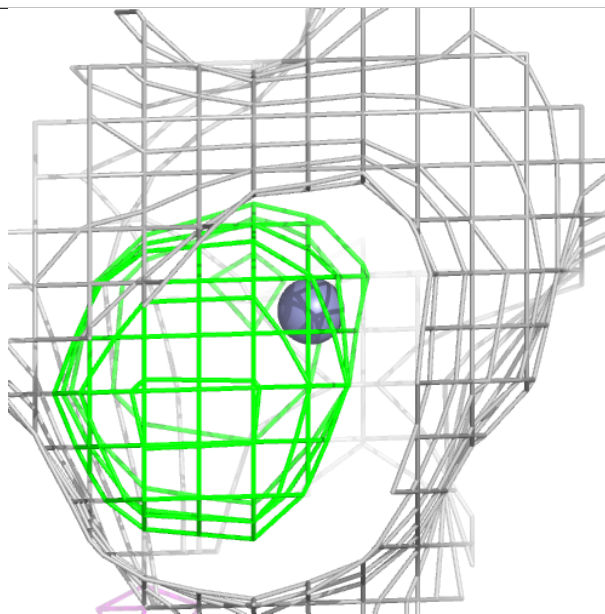
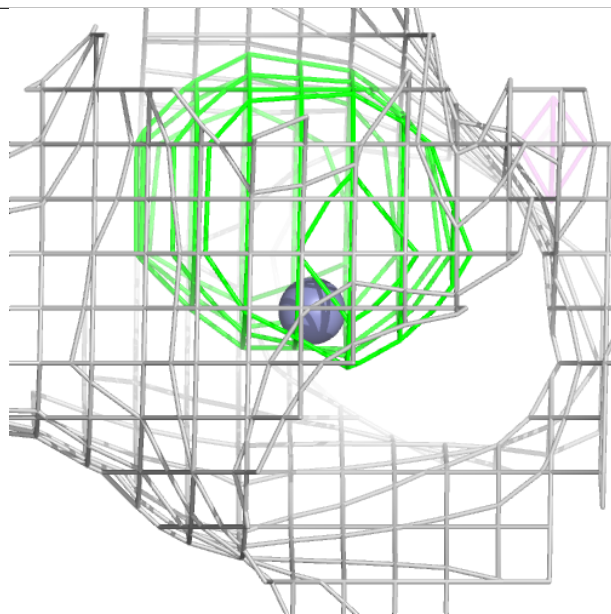
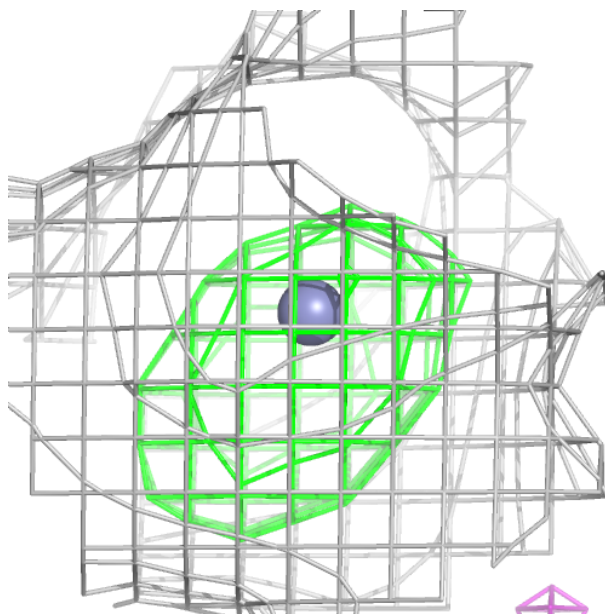
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





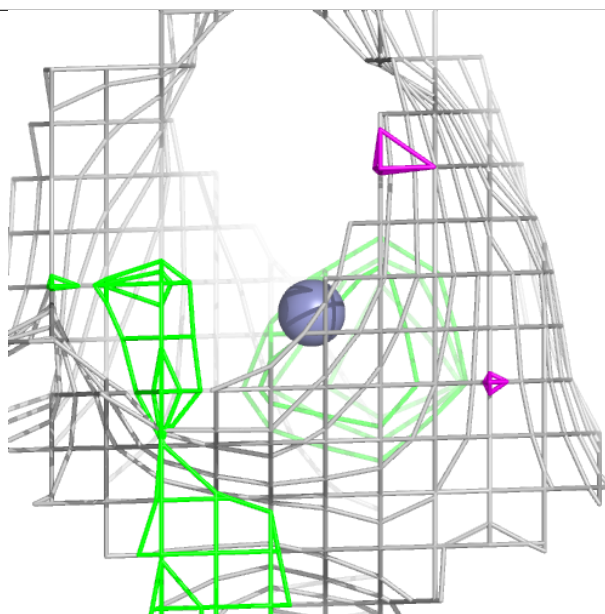
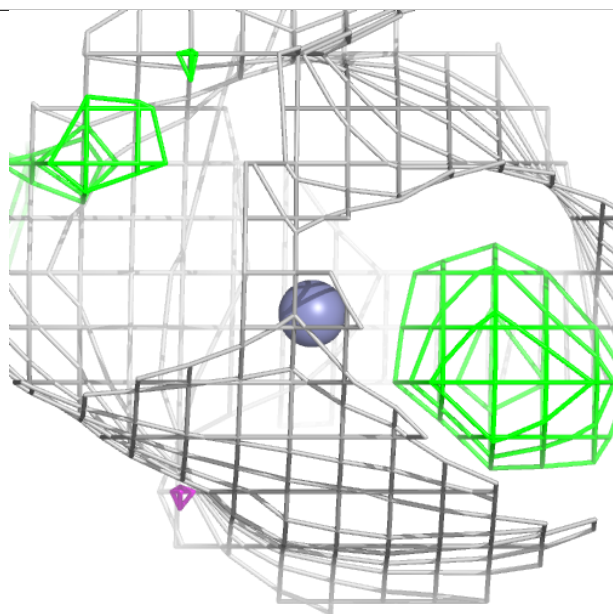
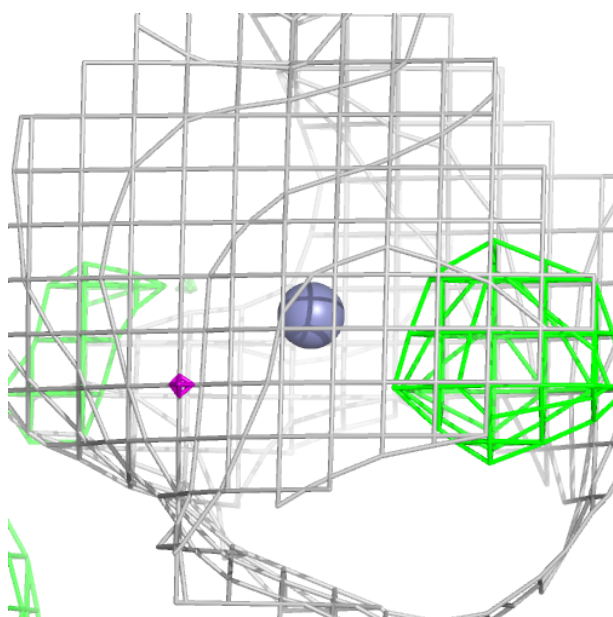
**Electron density around ZN A 402:**

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



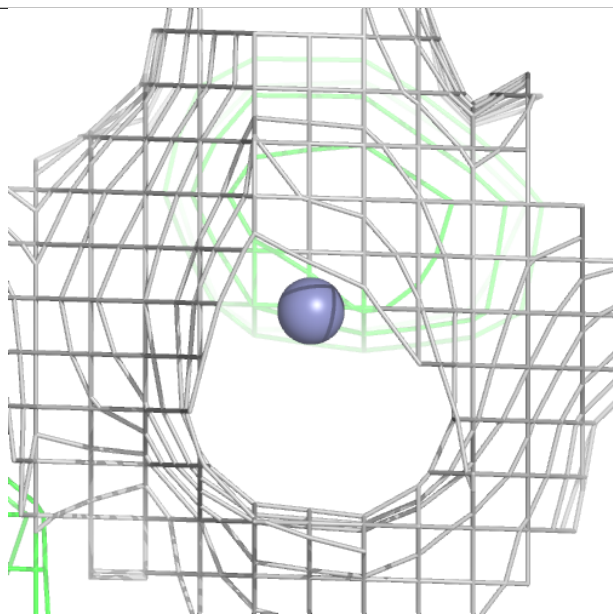
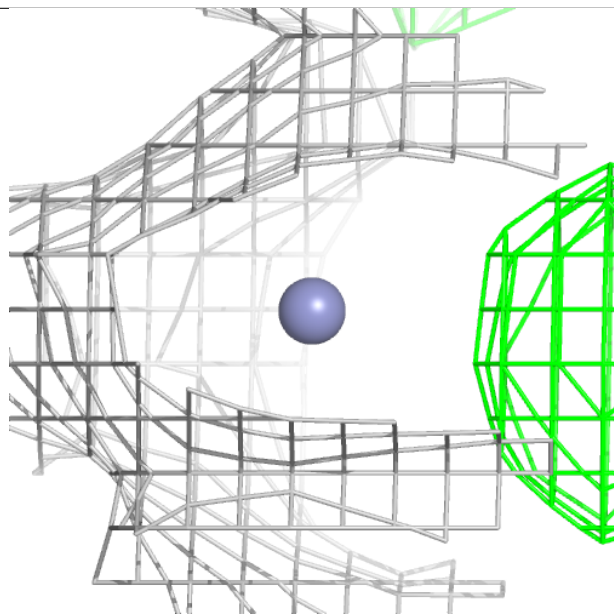
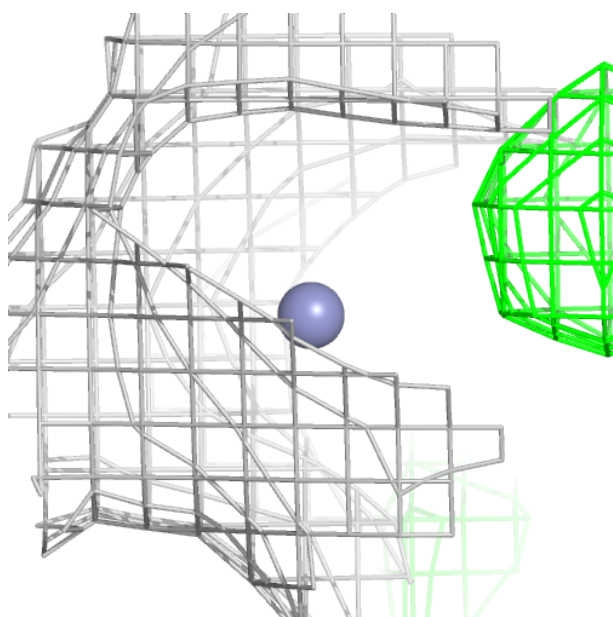
**Electron density around ZN F 402:**

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



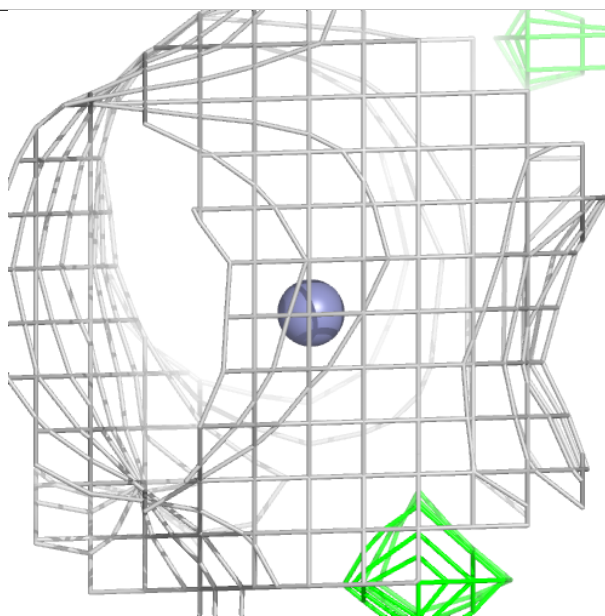
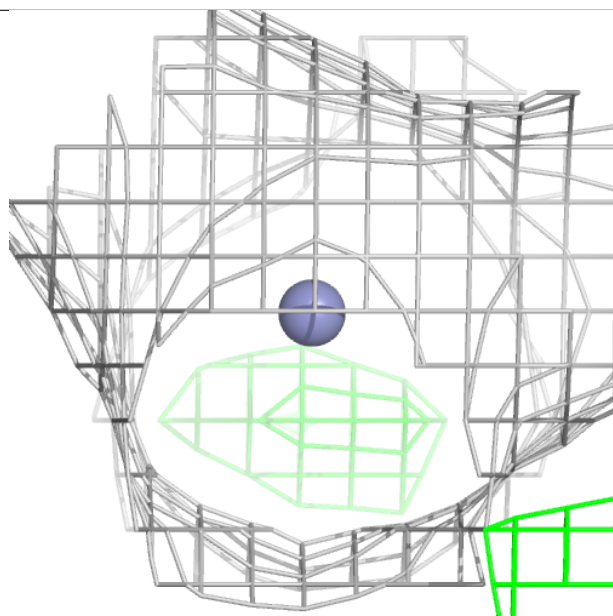
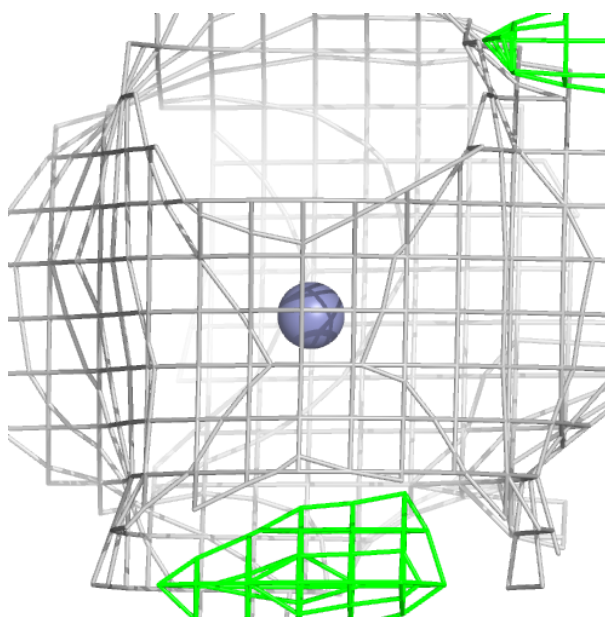
**Electron density around ZN E 403:**

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



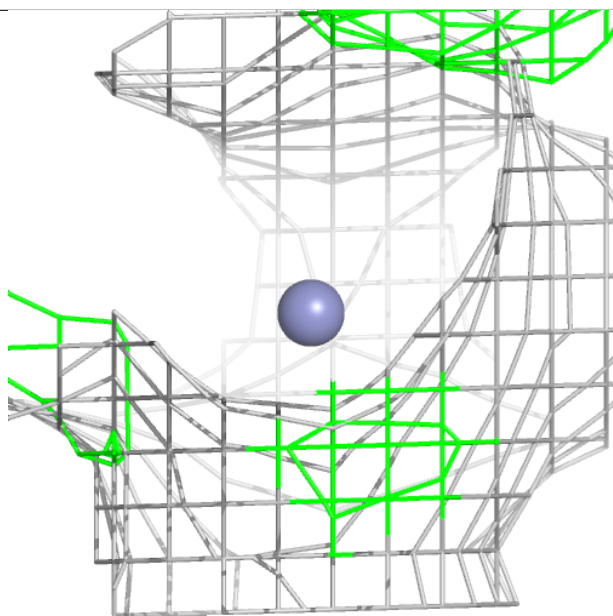
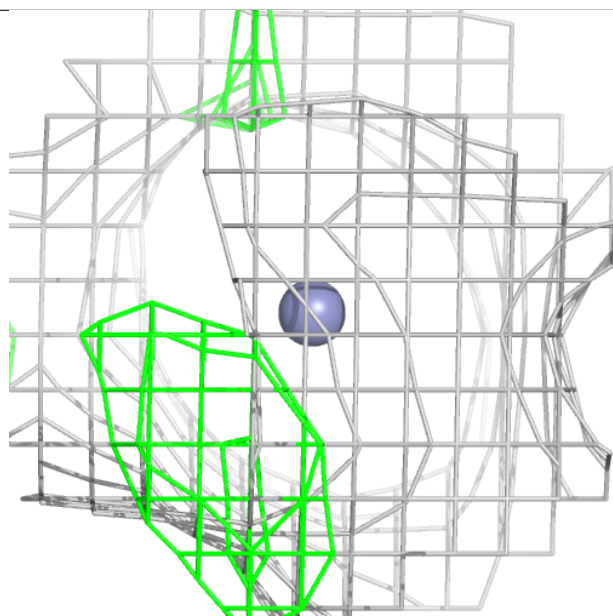
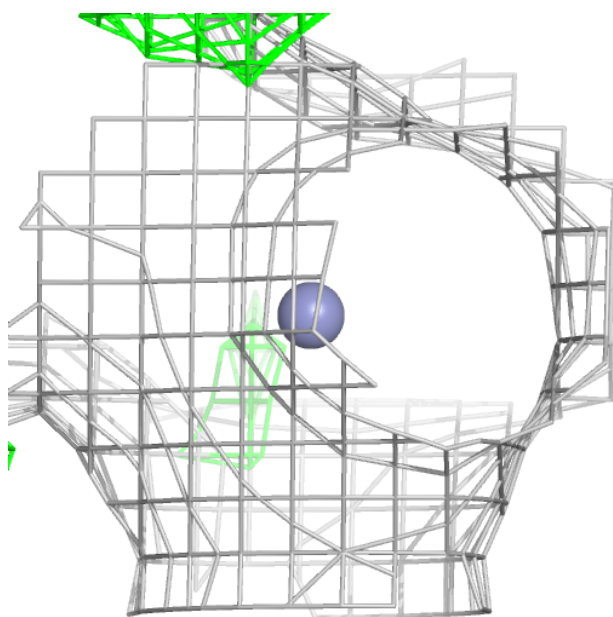
**Electron density around ZN C 402:**

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



**Electron density around ZN D 401:**

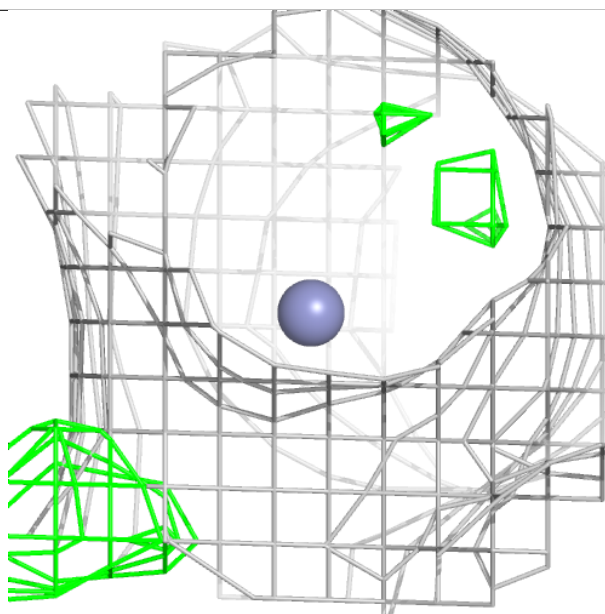
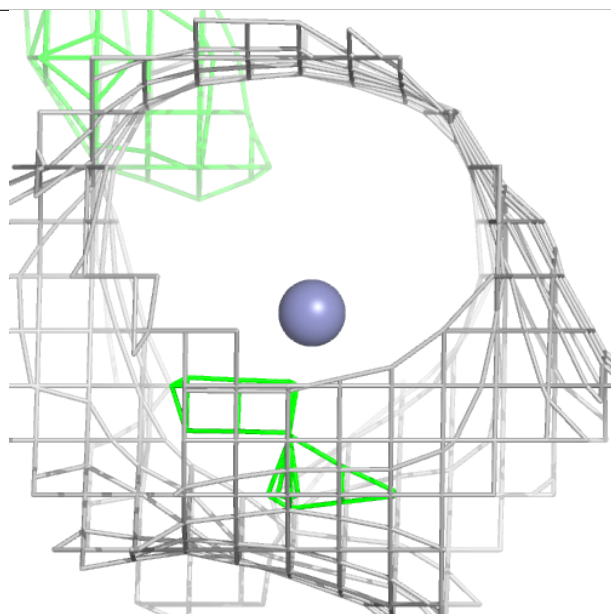
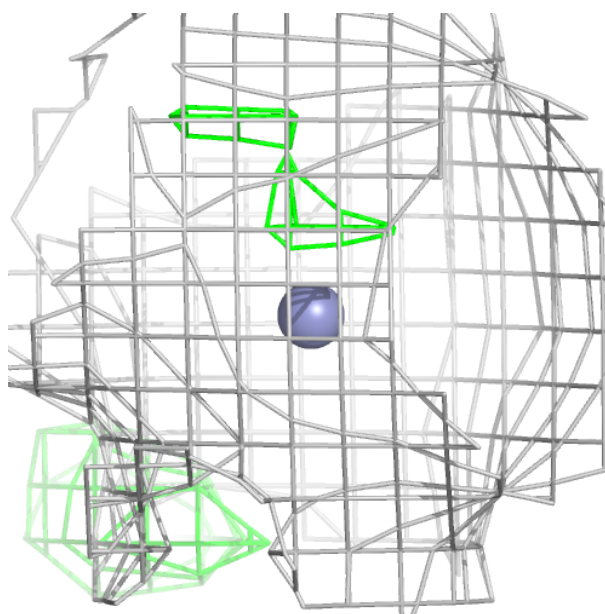
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





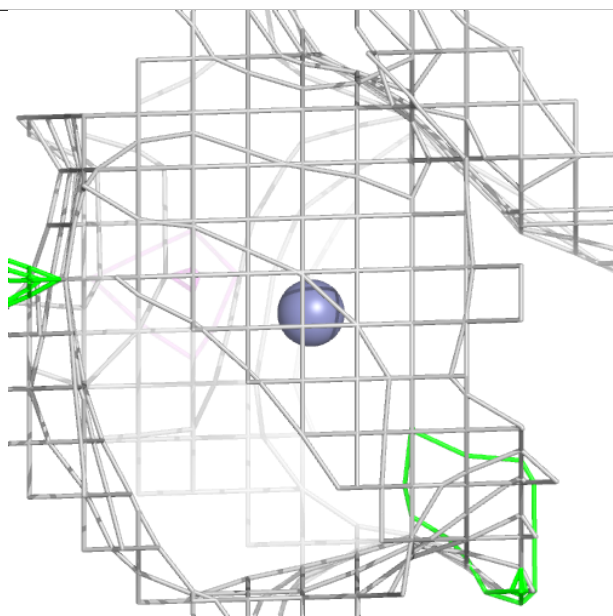
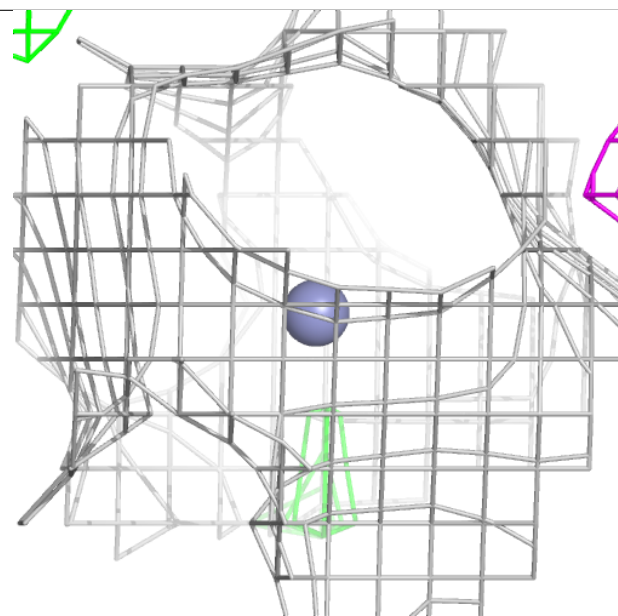
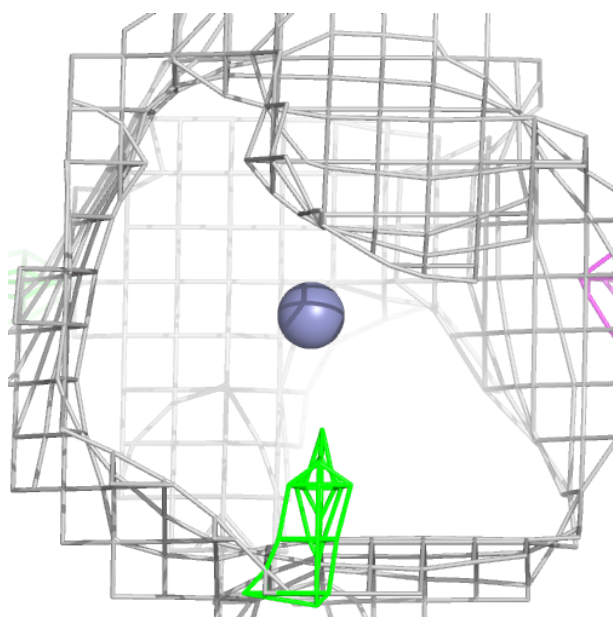
**Electron density around ZN D 402:**

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



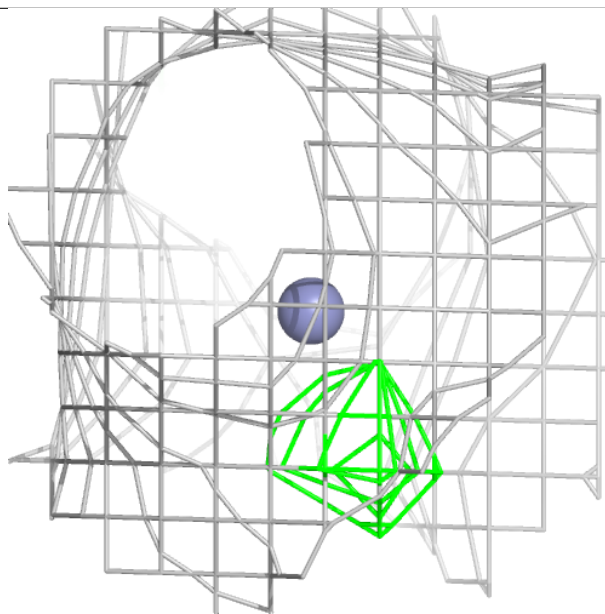
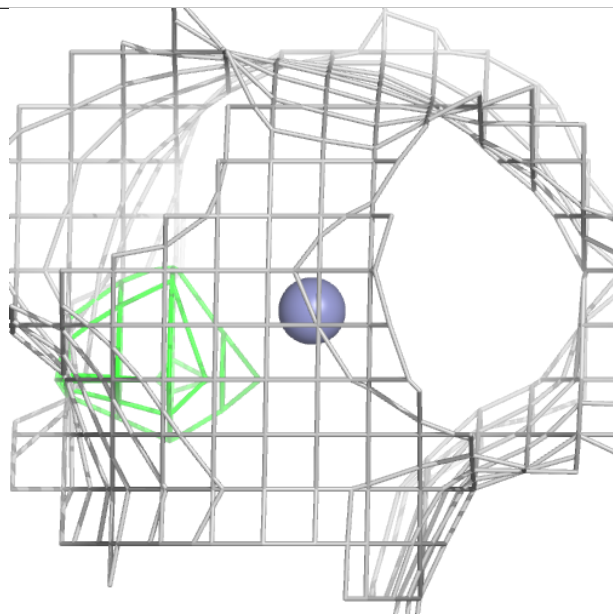
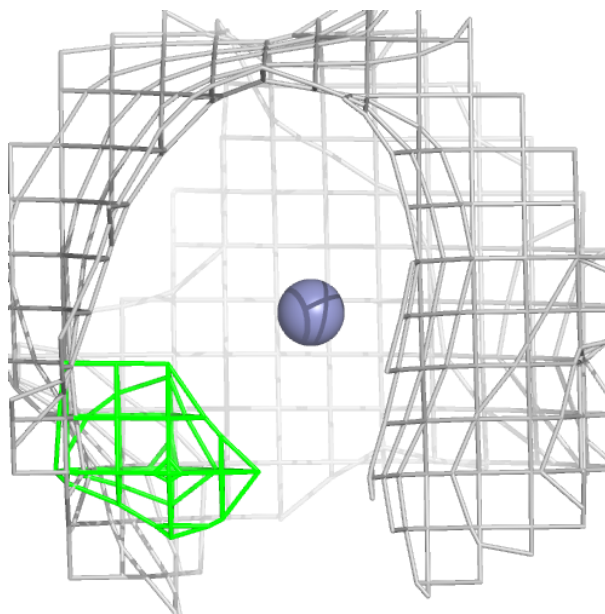
**Electron density around ZN D 403:**

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



**Electron density around ZN A 401:**

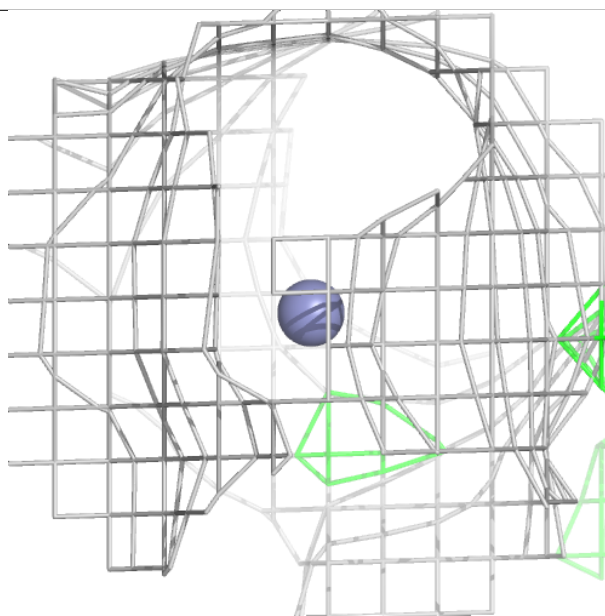
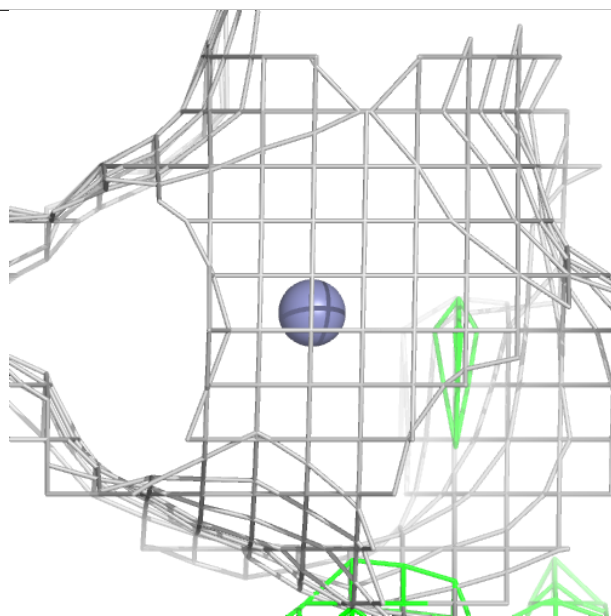
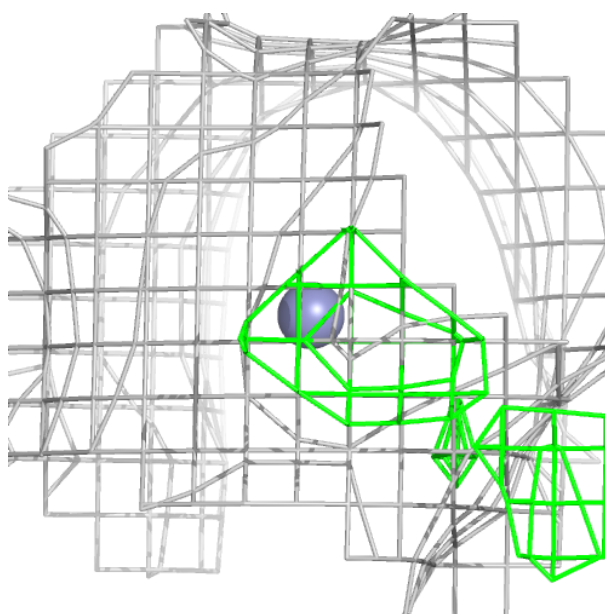
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





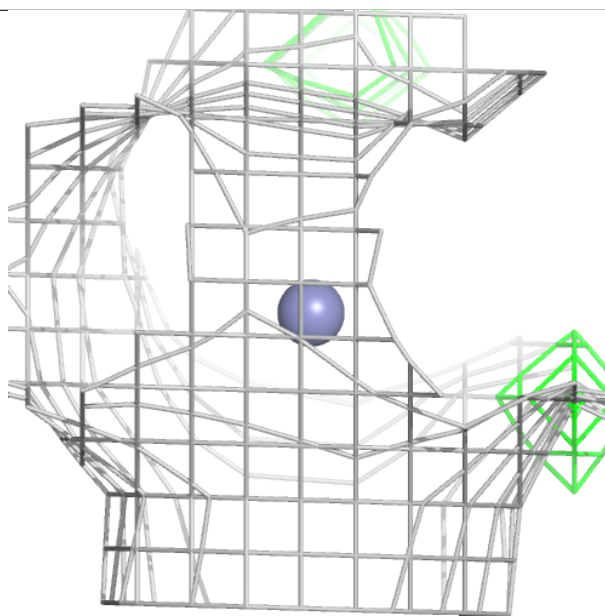
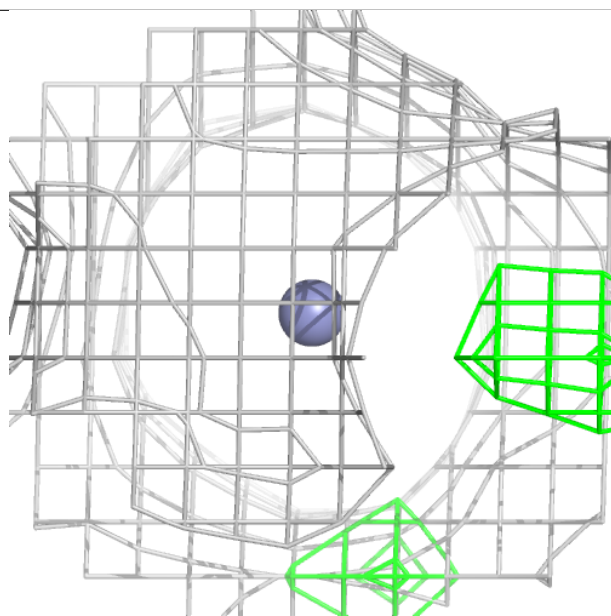
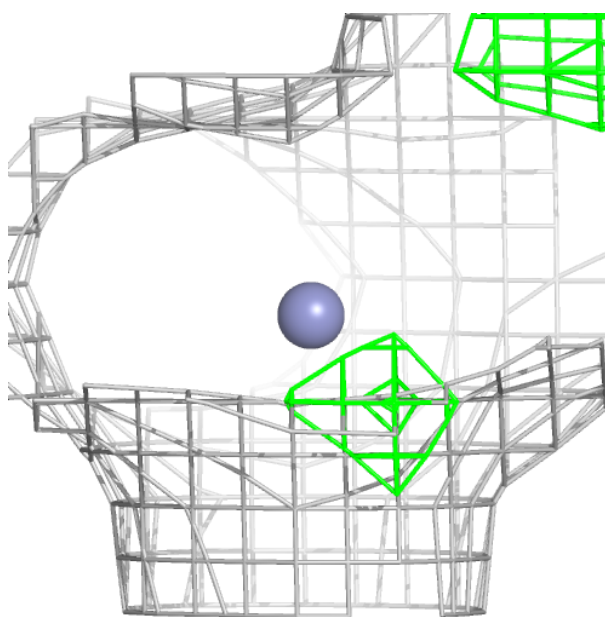
**Electron density around ZN B 401:**

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



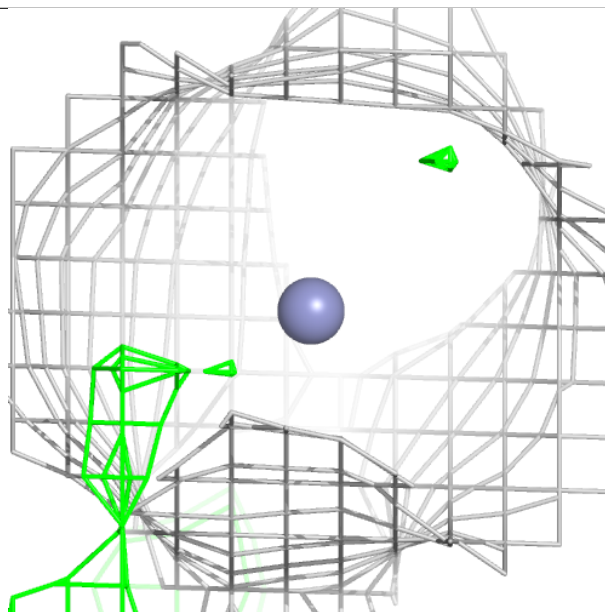
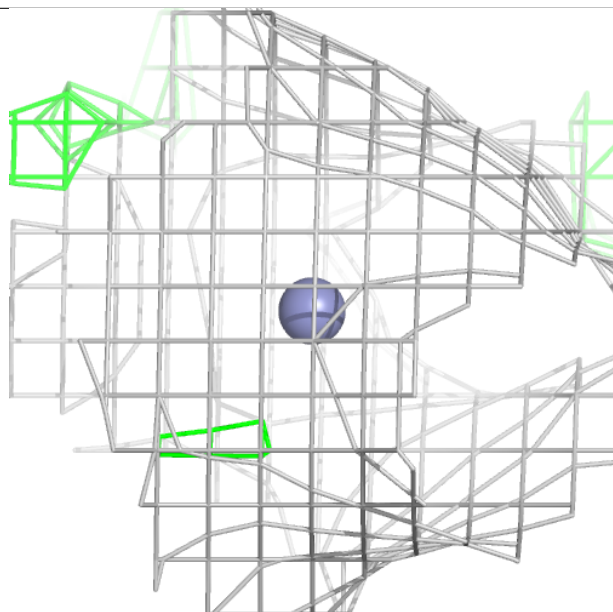
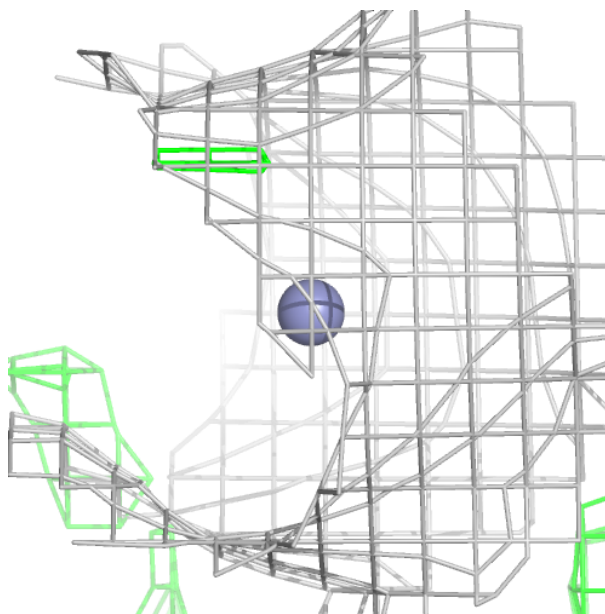
**Electron density around ZN C 401:**

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



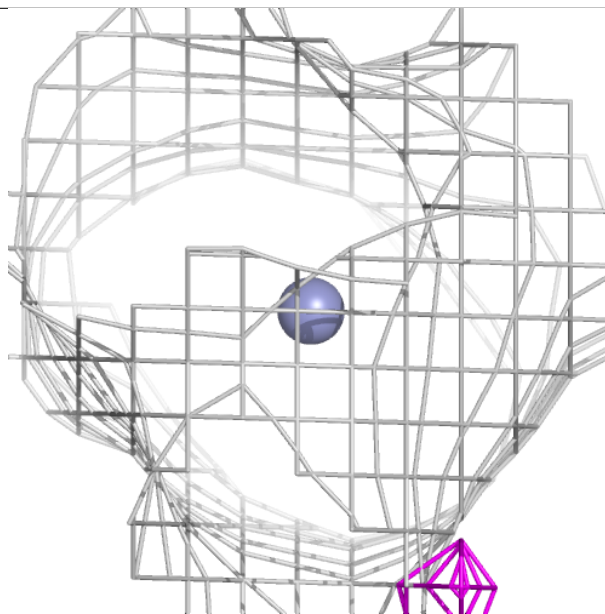
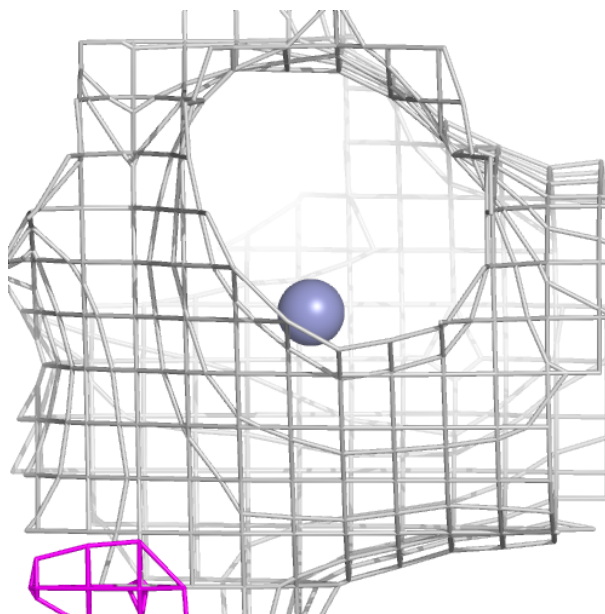
**Electron density around ZN F 401:**

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



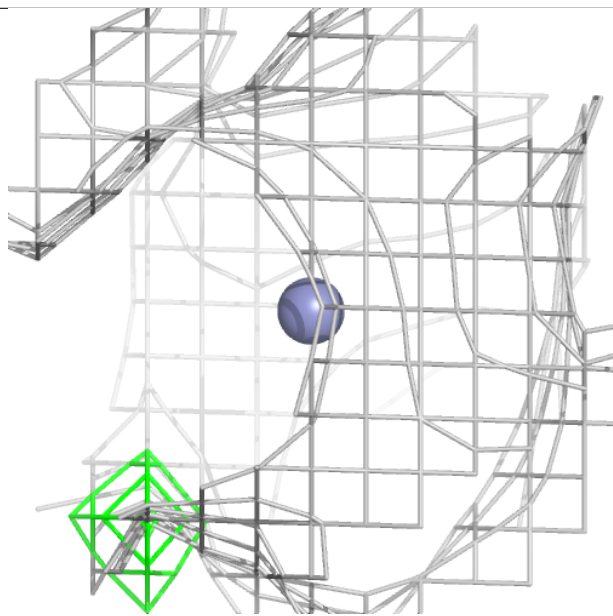
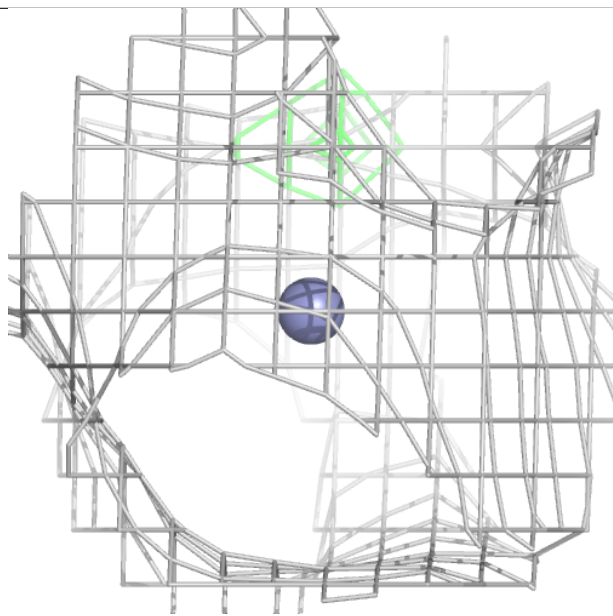
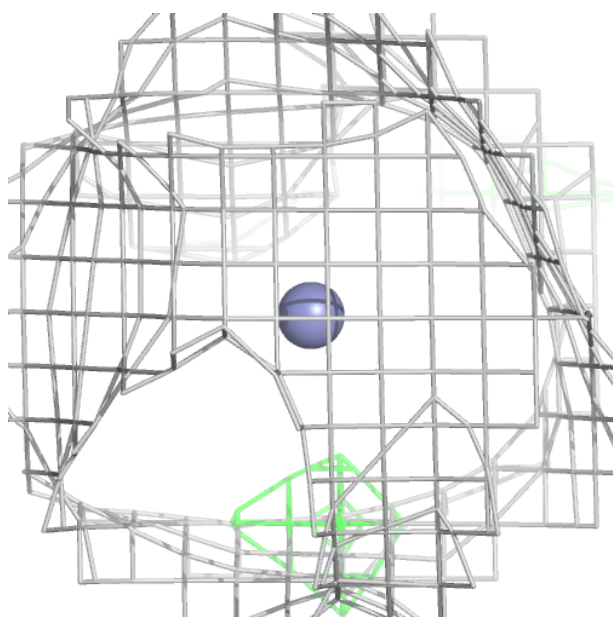
**Electron density around ZN B 402:**

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



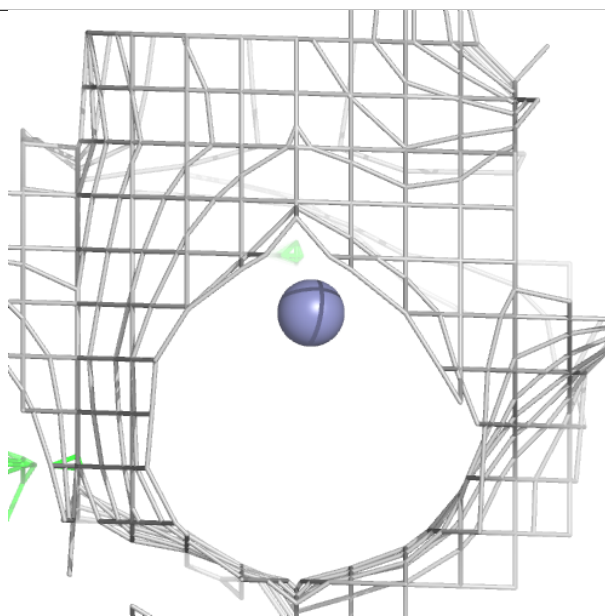
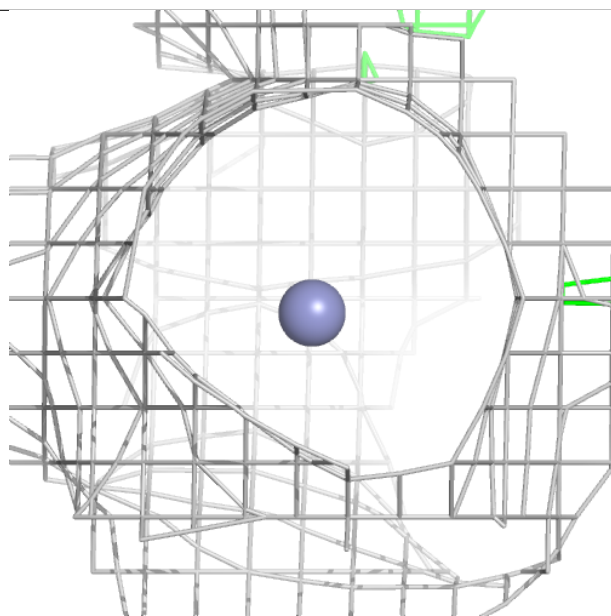
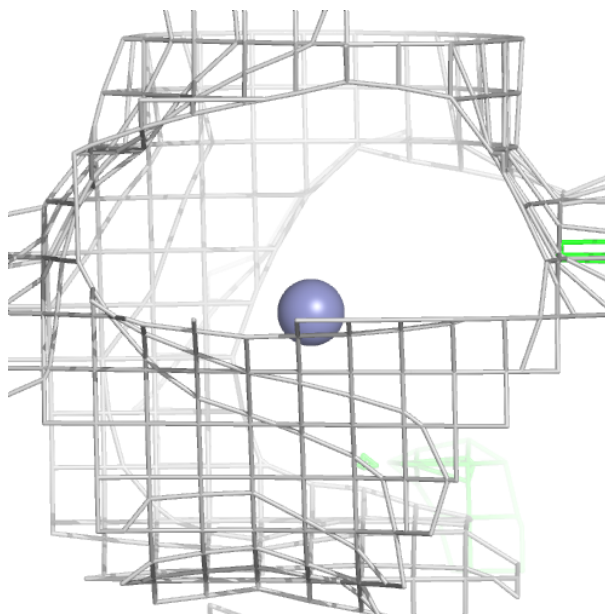
**Electron density around ZN C 403:**

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



**Electron density around ZN F 403:**

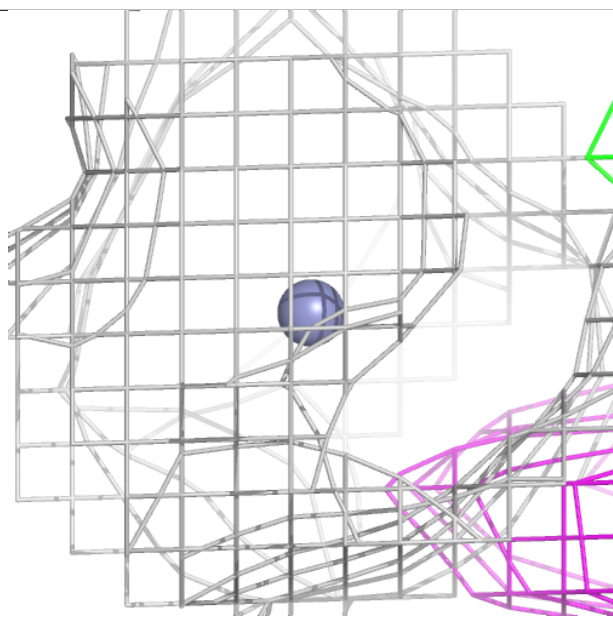
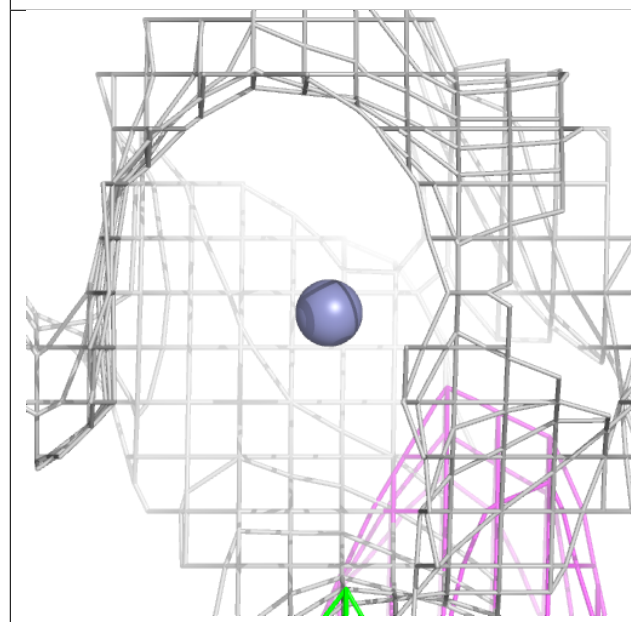
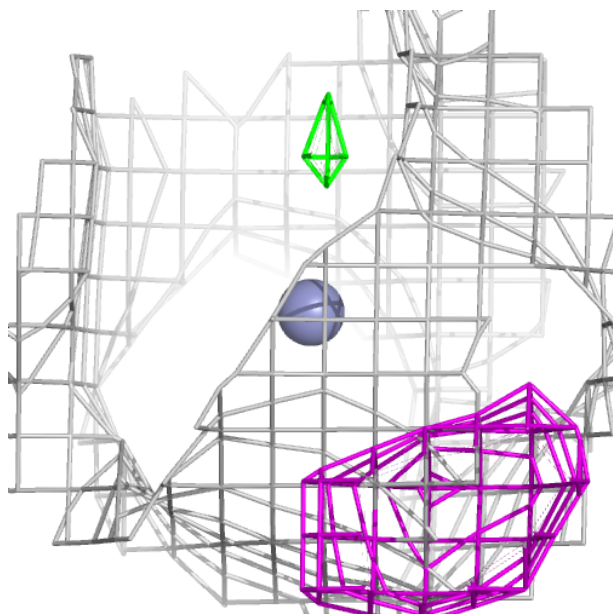
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





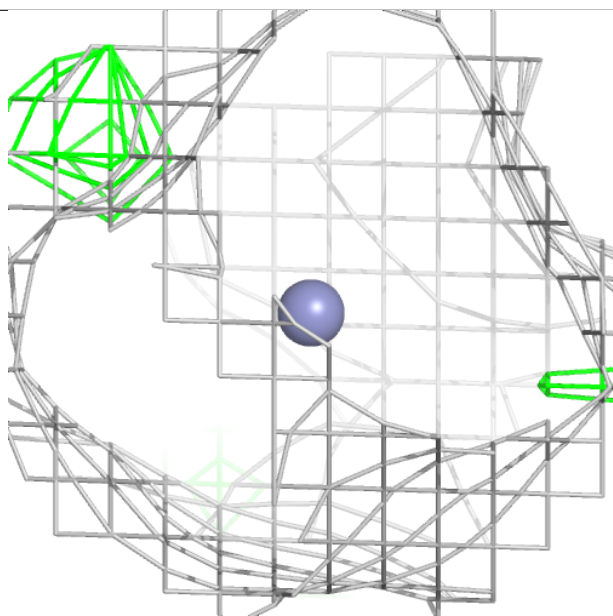
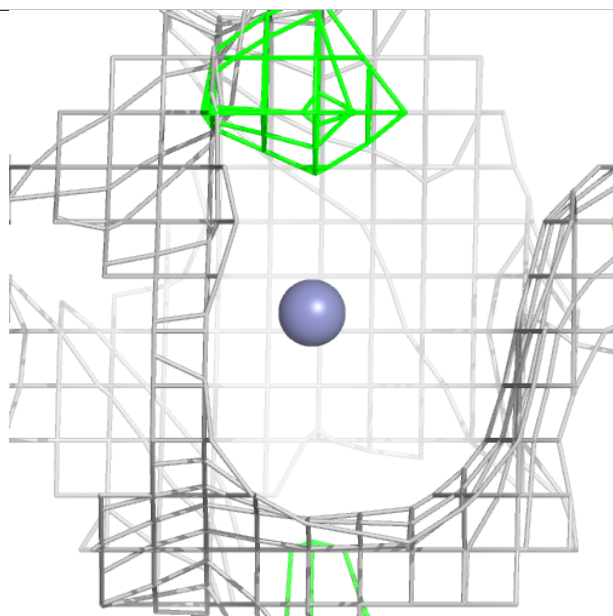
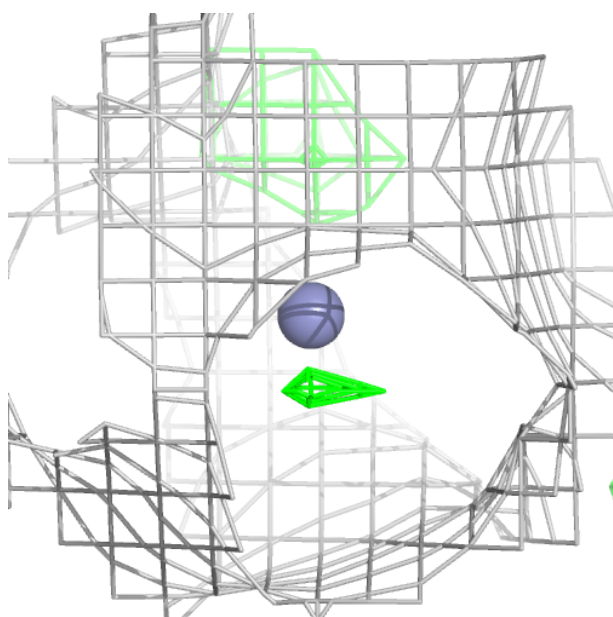
**Electron density around ZN B 403:**

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



**Electron density around ZN 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)



## 6.5 Other polymers ⓘ

There are no such residues in this entry.