



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:57 PM UTC

PDB ID : 8J9L / pdb_00008j9l
Title : Crystal Structure of Human H-Ferritin variant 123F assembling in solution2
Authors : Chen, X.; Zhao, G.
Deposited on : 2023-05-04
Resolution : 2.50 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:

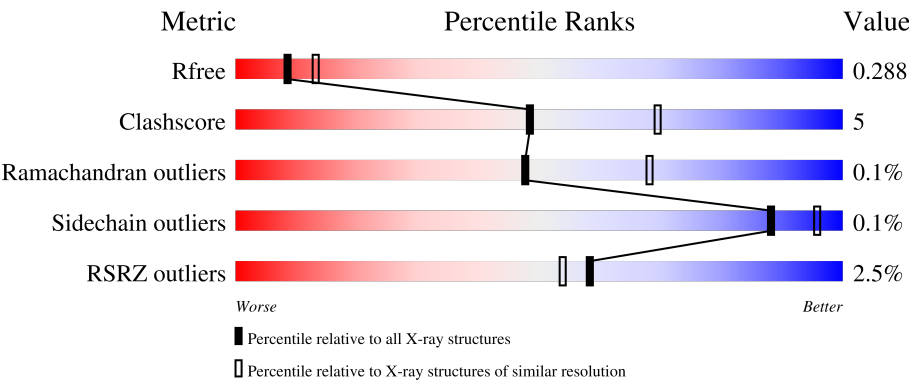
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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.



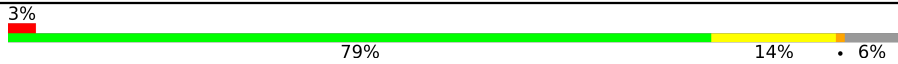
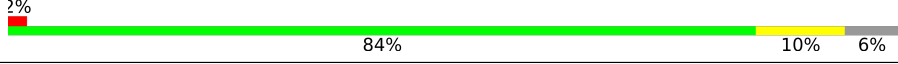
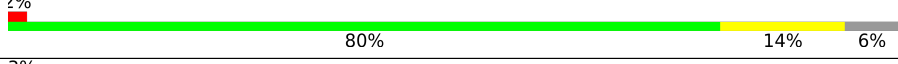
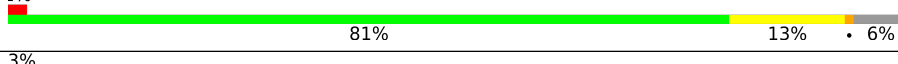
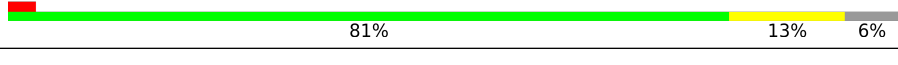
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	183	0.0% 84%10%6%
1	B	183	2.0% 80%14%6%
1	C	183	2.0% 74%20%6%
1	D	183	4.0% 85%8%6%
1	E	183	3.0% 79%15%6%

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Mol	Chain	Length	Quality of chain
1	F	183	 3% 79% 14% • 6%
1	G	183	 2% 78% 16% 6%
1	H	183	 2% 84% 10% 6%
1	I	183	 2% 80% 14% 6%
1	J	183	 3% 84% 10% 6%
1	K	183	 2% 81% 13% • 6%
1	L	183	 3% 81% 13% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17073 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 Ferritin heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	B	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	C	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	D	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	E	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	F	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	G	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	H	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	I	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	J	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	K	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			
1	L	172	Total	C	N	O	S	0	0	0
			1416	891	248	270	7			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	123	PHE	ASP	engineered mutation	UNP P02794
B	123	PHE	ASP	engineered mutation	UNP P02794
C	123	PHE	ASP	engineered mutation	UNP P02794
D	123	PHE	ASP	engineered mutation	UNP P02794
E	123	PHE	ASP	engineered mutation	UNP P02794

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Chain	Residue	Modelled	Actual	Comment	Reference
F	123	PHE	ASP	engineered mutation	UNP P02794
G	123	PHE	ASP	engineered mutation	UNP P02794
H	123	PHE	ASP	engineered mutation	UNP P02794
I	123	PHE	ASP	engineered mutation	UNP P02794
J	123	PHE	ASP	engineered mutation	UNP P02794
K	123	PHE	ASP	engineered mutation	UNP P02794
L	123	PHE	ASP	engineered mutation	UNP P02794

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe 1 1	0	0
2	B	2	Total Fe 2 2	0	0
2	C	2	Total Fe 2 2	0	0
2	D	2	Total Fe 2 2	0	0
2	E	1	Total Fe 1 1	0	0
2	F	1	Total Fe 1 1	0	0
2	G	1	Total Fe 1 1	0	0
2	H	2	Total Fe 2 2	0	0
2	I	1	Total Fe 1 1	0	0
2	J	1	Total Fe 1 1	0	0
2	K	1	Total Fe 1 1	0	0
2	L	1	Total Fe 1 1	0	0

- Molecule 3 is water.

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

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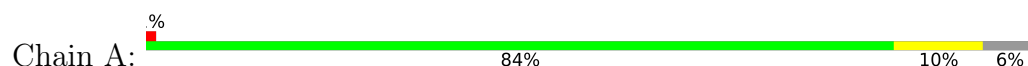
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	7	Total 7	O 7	0	0
3	C	6	Total 6	O 6	0	0
3	D	4	Total 4	O 4	0	0
3	E	5	Total 5	O 5	0	0
3	F	5	Total 5	O 5	0	0
3	G	9	Total 9	O 9	0	0
3	H	3	Total 3	O 3	0	0
3	I	3	Total 3	O 3	0	0
3	J	5	Total 5	O 5	0	0
3	K	7	Total 7	O 7	0	0
3	L	5	Total 5	O 5	0	0

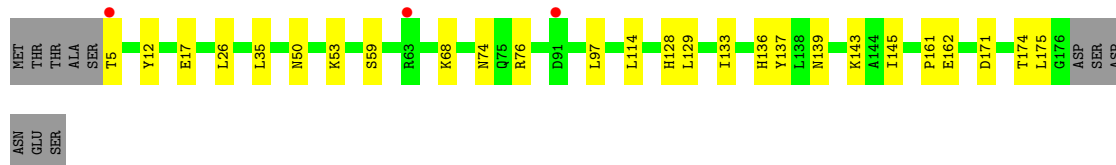
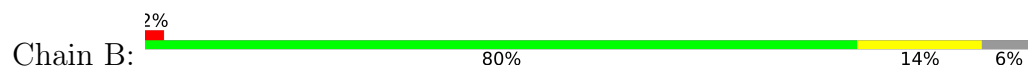
3 Residue-property plots [i](#)

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: Ferritin heavy chain



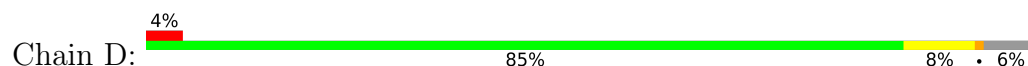
- Molecule 1: Ferritin heavy chain



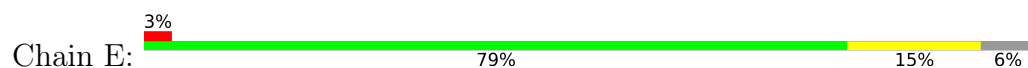
- Molecule 1: Ferritin heavy chain

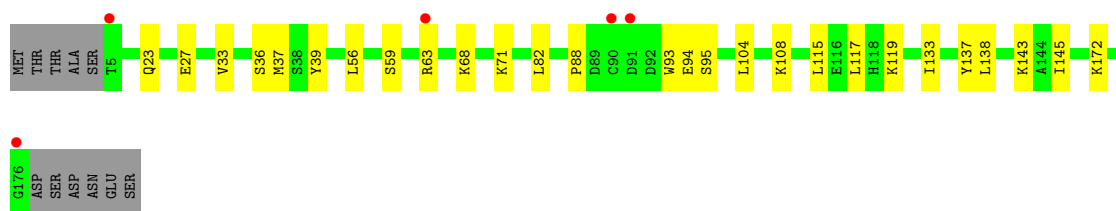


- Molecule 1: Ferritin heavy chain

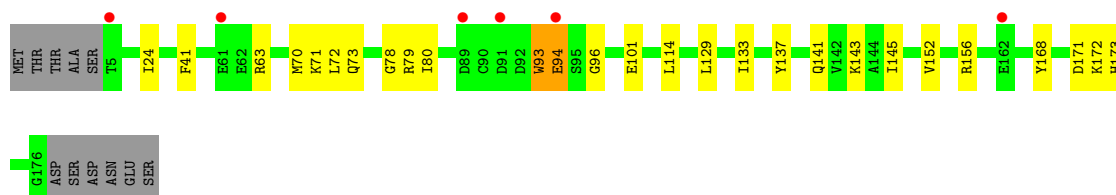
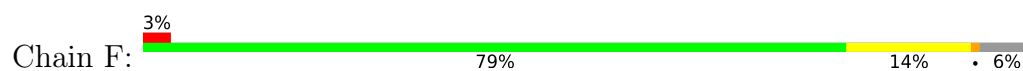


- Molecule 1: Ferritin heavy chain

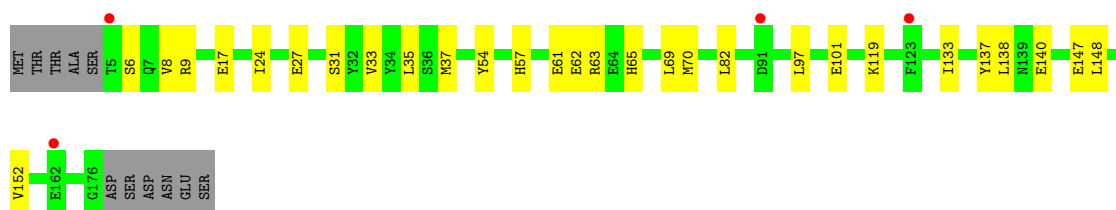
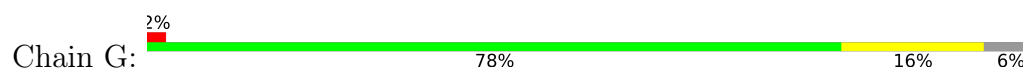




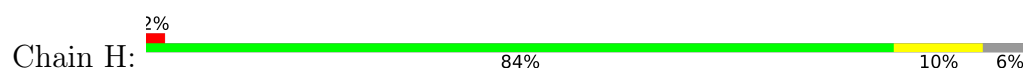
• Molecule 1: Ferritin heavy chain



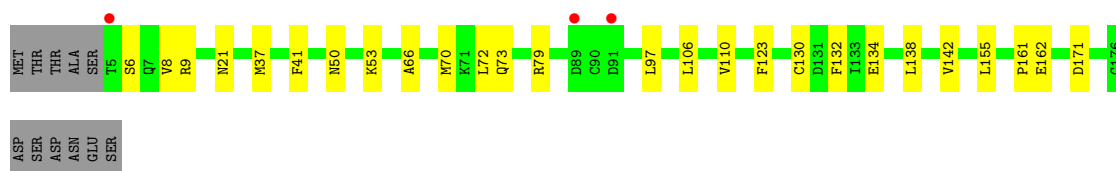
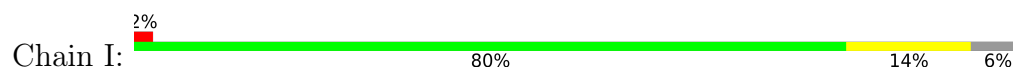
• Molecule 1: Ferritin heavy chain



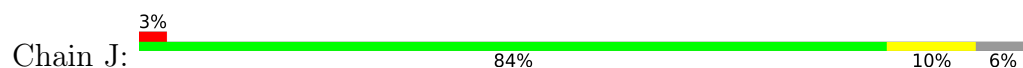
• Molecule 1: Ferritin heavy chain

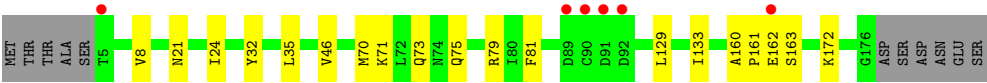


• Molecule 1: Ferritin heavy chain

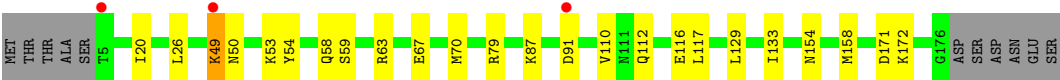
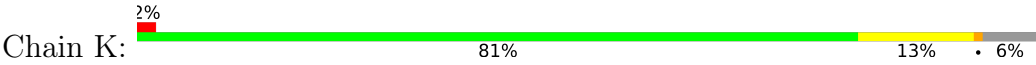


• Molecule 1: Ferritin heavy chain

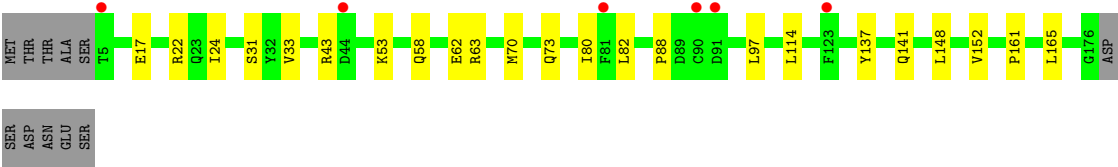
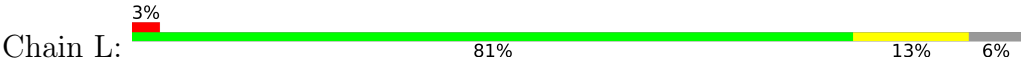




● Molecule 1: Ferritin heavy chain



● Molecule 1: Ferritin heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	261.70Å 261.70Å 320.51Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.26 – 2.50 29.26 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.3 (29.26-2.50) 97.2 (29.26-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.247 , 0.289 0.247 , 0.288	Depositor DCC
R_{free} test set	2011 reflections (1.39%)	wwPDB-VP
Wilson B-factor (Å ²)	14.0	Xtriage
Anisotropy	0.924	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -2/3*h-1/3*k+2/3*l,-1/3*h-2/3*k-2/3*l,2/3*h-2/3*k+1/3*l 0.000 for -h,1/3*h-1/3*k+2/3*l,2/3*h+4/3*k+1/3*l 0.000 for -1/3*h+1/3*k-2/3*l,-k,-4/3*h-2/3*k+1/3*l	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	17073	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.85% 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: FE

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	0.38	0/1446	0.59	0/1947
1	B	0.43	0/1446	0.59	0/1947
1	C	0.42	0/1446	0.59	0/1947
1	D	0.43	0/1446	0.68	0/1947
1	E	0.40	0/1446	0.59	0/1947
1	F	0.39	0/1446	0.65	1/1947 (0.1%)
1	G	0.42	0/1446	0.63	0/1947
1	H	0.39	0/1446	0.63	0/1947
1	I	0.40	0/1446	0.59	0/1947
1	J	0.42	1/1446 (0.1%)	0.65	0/1947
1	K	0.42	0/1446	0.66	1/1947 (0.1%)
1	L	0.42	0/1446	0.63	1/1947 (0.1%)
All	All	0.41	1/17352 (0.0%)	0.63	3/23364 (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	F	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	46	VAL	CA-CB	-5.37	1.50	1.55

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	49	LYS	CB-CG-CD	6.12	125.37	111.30
1	L	22	ARG	CB-CG-CD	6.09	125.30	111.30
1	F	94	GLU	CA-CB-CG	5.64	125.39	114.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	93	TRP	Peptide

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	1416	0	1362	16	0
1	B	1416	0	1362	20	0
1	C	1416	0	1362	25	0
1	D	1416	0	1362	12	1
1	E	1416	0	1362	18	0
1	F	1416	0	1362	16	2
1	G	1416	0	1362	20	0
1	H	1416	0	1362	12	1
1	I	1416	0	1362	19	0
1	J	1416	0	1362	14	0
1	K	1416	0	1362	17	0
1	L	1416	0	1362	17	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	2	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	6	0	0	1	0
3	B	7	0	0	2	0
3	C	6	0	0	1	0
3	D	4	0	0	0	0
3	E	5	0	0	2	0
3	F	5	0	0	0	0
3	G	9	0	0	1	0
3	H	3	0	0	0	0
3	I	3	0	0	0	0
3	J	5	0	0	0	0
3	K	7	0	0	2	0
3	L	5	0	0	0	0
All	All	17073	0	16344	183	2

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 5.

All (183) 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:K:172:LYS:NZ	3:K:301:HOH:O	1.98	0.96
1:B:59:SER:OG	1:C:63:ARG:NH2	2.13	0.80
1:K:91:ASP:OD1	3:K:302:HOH:O	2.00	0.78
1:F:78:GLY:O	1:F:79:ARG:NH1	2.18	0.76
1:H:117:LEU:HD22	1:H:133:ILE:HD11	1.74	0.69
1:J:129:LEU:O	1:J:133:ILE:HG12	1.92	0.68
1:I:50:ASN:HB2	1:I:171:ASP:OD1	1.93	0.67
1:E:117:LEU:HD22	1:E:133:ILE:HD11	1.76	0.67
1:C:114:LEU:HD13	1:C:137:TYR:HB3	1.77	0.65
1:B:5:THR:N	3:B:301:HOH:O	2.30	0.65
1:B:161:PRO:HD2	1:B:162:GLU:OE2	1.96	0.65
1:J:35:LEU:HG	1:L:70:MET:HE1	1.79	0.64
1:F:114:LEU:HD13	1:F:137:TYR:HB3	1.79	0.63
1:H:114:LEU:HD13	1:H:137:TYR:HB3	1.80	0.62
1:D:101:GLU:OE1	1:D:156:ARG:NH2	2.30	0.62
1:B:68:LYS:HE3	1:B:136:HIS:CG	2.35	0.61
1:L:53:LYS:HD2	1:L:53:LYS:C	2.26	0.61
1:G:63:ARG:HG2	1:K:63:ARG:HH21	1.66	0.60
1:A:114:LEU:HD13	1:A:137:TYR:HB3	1.84	0.60
1:C:33:VAL:HG22	1:C:88:PRO:HB3	1.83	0.60
1:B:139:ASN:ND2	1:B:143:LYS:HE3	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:21:ASN:OD1	1:J:73:GLN:NE2	2.38	0.57
1:C:161:PRO:HD2	1:C:162:GLU:OE2	2.04	0.56
1:H:97:LEU:HD23	1:H:161:PRO:HD3	1.86	0.56
1:G:54:TYR:OH	1:G:147:GLU:OE1	2.18	0.56
1:F:24:ILE:HD13	1:F:70:MET:HG2	1.89	0.55
1:C:172:LYS:HE2	3:C:305:HOH:O	2.07	0.55
1:A:60:HIS:CD2	1:E:63:ARG:NH1	2.75	0.55
1:F:152:VAL:O	1:F:156:ARG:HG2	2.07	0.54
1:G:35:LEU:HG	1:K:70:MET:HE1	1.87	0.54
1:L:70:MET:HG2	1:L:80:ILE:HD13	1.89	0.54
1:H:49:LYS:HA	1:H:49:LYS:HE2	1.89	0.54
1:I:6:SER:HB3	1:I:9:ARG:HG3	1.90	0.54
1:H:13:HIS:CD2	1:H:124:LYS:HD2	2.43	0.53
1:L:24:ILE:HD13	1:L:70:MET:HG3	1.91	0.53
1:G:6:SER:HB3	1:G:9:ARG:HB2	1.89	0.52
1:E:133:ILE:HG22	1:E:138:LEU:HG	1.91	0.52
1:A:82:LEU:HD12	1:E:36:SER:HB2	1.92	0.52
1:C:152:VAL:O	1:C:156:ARG:HG3	2.10	0.51
1:E:172:LYS:HE2	3:E:304:HOH:O	2.10	0.51
1:C:20:ILE:HD13	1:C:117:LEU:HD21	1.93	0.51
1:I:21:ASN:OD1	1:I:73:GLN:NE2	2.39	0.51
1:J:79:ARG:NH1	1:L:43:ARG:NE	2.59	0.51
1:I:50:ASN:HA	1:I:53:LYS:HG2	1.93	0.51
1:K:129:LEU:O	1:K:133:ILE:HG13	2.11	0.51
1:E:33:VAL:HG22	1:E:88:PRO:HB3	1.94	0.50
1:F:114:LEU:HD12	1:F:141:GLN:HG3	1.93	0.50
1:L:148:LEU:O	1:L:152:VAL:HG23	2.12	0.50
1:G:24:ILE:HD13	1:G:70:MET:HG2	1.94	0.50
1:E:23:GLN:O	1:E:27:GLU:HG2	2.12	0.49
1:G:63:ARG:NH2	1:K:59:SER:OG	2.44	0.49
1:I:97:LEU:HA	1:I:155:LEU:HD13	1.95	0.49
1:L:33:VAL:HG22	1:L:88:PRO:HB3	1.94	0.49
1:I:130:CYS:O	1:I:134:GLU:HG3	2.13	0.49
1:H:119:LYS:HG3	1:I:123:PHE:CE2	2.48	0.49
1:A:33:VAL:HG12	1:A:37:MET:HE3	1.93	0.49
1:F:129:LEU:O	1:F:133:ILE:HG12	2.13	0.48
1:G:57:HIS:O	1:G:61:GLU:HG2	2.13	0.48
1:K:54:TYR:O	1:K:58:GLN:HG2	2.13	0.48
1:E:93:TRP:O	1:E:95:SER:N	2.47	0.48
1:F:168:TYR:CZ	1:F:172:LYS:HE2	2.48	0.48
1:C:62:GLU:HA	1:C:62:GLU:OE1	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:LYS:NZ	3:A:301:HOH:O	2.39	0.48
1:K:20:ILE:HD13	1:K:117:LEU:HD21	1.95	0.47
1:B:97:LEU:HD23	1:B:161:PRO:HD3	1.96	0.47
1:I:138:LEU:HA	1:I:138:LEU:HD23	1.65	0.47
1:A:70:MET:HE2	1:E:39:TYR:CD2	2.49	0.47
1:C:79:ARG:HD3	1:C:79:ARG:HA	1.60	0.47
1:F:73:GLN:HG2	1:F:80:ILE:HG13	1.95	0.47
1:B:74:ASN:HD22	1:C:42:ASP:HB3	1.79	0.47
1:H:41:PHE:CZ	1:H:96:GLY:HA2	2.50	0.47
1:E:115:LEU:O	1:E:119:LYS:HD3	2.14	0.46
1:L:97:LEU:HD23	1:L:161:PRO:HD3	1.97	0.46
1:I:161:PRO:HD2	1:I:162:GLU:CD	2.41	0.46
1:A:133:ILE:HG22	1:A:138:LEU:HG	1.97	0.46
1:C:65:HIS:HB3	1:C:137:TYR:HE1	1.80	0.46
1:L:58:GLN:O	1:L:62:GLU:HG2	2.15	0.46
1:B:174:THR:HG22	1:B:175:LEU:HD23	1.98	0.46
1:D:79:ARG:HA	1:D:79:ARG:HD3	1.75	0.46
1:E:68:LYS:HB3	1:E:137:TYR:OH	2.16	0.46
1:J:35:LEU:HD13	1:L:63:ARG:HH21	1.79	0.46
1:D:72:LEU:HD22	1:D:132:PHE:CD1	2.50	0.46
1:I:50:ASN:OD1	1:I:53:LYS:HE3	2.16	0.46
1:J:32:TYR:OH	1:L:82:LEU:HB3	2.16	0.46
1:G:148:LEU:O	1:G:152:VAL:HG23	2.15	0.45
1:J:21:ASN:ND2	1:J:81:PHE:HB2	2.30	0.45
1:D:145:ILE:HG22	1:J:8:VAL:HB	1.98	0.45
1:D:159:GLY:O	1:D:162:GLU:HB2	2.16	0.45
1:G:33:VAL:HG12	1:G:37:MET:HE3	1.97	0.45
1:H:72:LEU:HD22	1:H:132:PHE:CD2	2.51	0.45
1:F:101:GLU:OE1	1:F:156:ARG:NH2	2.48	0.45
1:B:5:THR:HG22	3:B:301:HOH:O	2.15	0.45
1:C:13:HIS:CD2	1:C:124:LYS:HD2	2.51	0.45
1:C:73:GLN:HG2	1:C:80:ILE:HG13	1.99	0.45
1:K:63:ARG:O	1:K:67:GLU:HG3	2.16	0.45
1:B:162:GLU:H	1:B:162:GLU:CD	2.24	0.45
1:A:172:LYS:HG2	1:F:173:HIS:HB3	1.98	0.45
1:I:161:PRO:HD2	1:I:162:GLU:OE2	2.16	0.45
1:E:104:LEU:O	1:E:108:LYS:HG3	2.16	0.45
1:J:71:LYS:O	1:J:75:GLN:HG3	2.16	0.44
1:D:72:LEU:HD22	1:D:132:PHE:CE1	2.53	0.44
1:F:71:LYS:HE3	1:F:71:LYS:HB3	1.83	0.44
1:G:63:ARG:HG2	1:K:63:ARG:NH2	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:LEU:HD23	1:A:115:LEU:HA	1.84	0.44
1:B:50:ASN:HA	1:B:53:LYS:HB2	2.00	0.44
1:B:76:ARG:HA	1:B:76:ARG:HD2	1.77	0.44
1:C:54:TYR:O	1:C:58:GLN:HG2	2.16	0.44
1:H:72:LEU:HD22	1:H:132:PHE:CE2	2.53	0.44
1:D:123:PHE:CZ	1:G:119:LYS:HE2	2.52	0.44
1:K:49:LYS:O	1:K:53:LYS:HG2	2.17	0.44
1:B:128:HIS:CE1	1:I:142:VAL:HG21	2.52	0.44
1:G:133:ILE:HG22	1:G:138:LEU:HG	1.99	0.44
1:K:26:LEU:HD21	1:K:110:VAL:HA	1.99	0.44
1:B:12:TYR:CE2	1:B:17:GLU:HB2	2.52	0.44
1:L:114:LEU:HD12	1:L:141:GLN:HG3	2.00	0.44
1:L:165:LEU:HA	1:L:165:LEU:HD12	1.68	0.44
1:A:32:TYR:OH	1:E:82:LEU:HB3	2.18	0.44
1:C:118:HIS:HB2	1:C:133:ILE:HG21	1.99	0.44
1:K:50:ASN:HB2	1:K:171:ASP:CG	2.43	0.44
1:A:79:ARG:HD3	1:A:79:ARG:HA	1.79	0.44
1:H:53:LYS:HE2	1:H:53:LYS:HB3	1.63	0.44
1:C:101:GLU:CD	1:C:156:ARG:HH21	2.26	0.44
1:F:72:LEU:HD21	1:F:129:LEU:HD13	2.00	0.43
1:I:6:SER:OG	1:I:8:VAL:HG22	2.18	0.43
1:D:133:ILE:HG22	1:D:138:LEU:HG	2.00	0.43
1:F:93:TRP:O	1:F:94:GLU:HB3	2.12	0.43
1:I:72:LEU:HD22	1:I:132:PHE:CE2	2.54	0.43
1:A:138:LEU:HD23	1:A:138:LEU:HA	1.66	0.43
1:G:9:ARG:NH2	1:G:17:GLU:OE2	2.39	0.43
1:C:140:GLU:OE1	1:C:140:GLU:N	2.51	0.43
1:A:155:LEU:HD23	1:A:155:LEU:HA	1.88	0.43
1:C:19:ALA:HB1	1:C:117:LEU:HD13	1.99	0.43
1:E:143:LYS:NZ	3:E:301:HOH:O	1.98	0.43
1:E:33:VAL:HG12	1:E:37:MET:HE3	1.99	0.43
1:D:33:VAL:HG22	1:D:88:PRO:HB3	2.01	0.43
1:B:35:LEU:HG	1:C:70:MET:HE1	2.01	0.43
3:G:302:HOH:O	1:K:79:ARG:NH2	2.52	0.43
1:B:114:LEU:HD13	1:B:137:TYR:HB3	2.01	0.43
1:G:31:SER:HB2	1:G:62:GLU:HB2	2.01	0.42
1:K:154:ASN:O	1:K:158:MET:HG3	2.19	0.42
1:K:112:GLN:O	1:K:116:GLU:HG2	2.18	0.42
1:L:31:SER:HB2	1:L:62:GLU:HB2	2.01	0.42
1:L:114:LEU:HD13	1:L:137:TYR:HB3	2.00	0.42
1:A:114:LEU:HD12	1:A:141:GLN:HG3	1.99	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:37:MET:O	1:I:41:PHE:HD2	2.02	0.42
1:I:106:LEU:O	1:I:110:VAL:HG23	2.19	0.42
1:A:26:LEU:HD21	1:A:110:VAL:HA	2.02	0.42
1:G:82:LEU:O	1:K:87:LYS:HD3	2.19	0.42
1:J:24:ILE:HD13	1:J:70:MET:HG2	2.00	0.42
1:D:162:GLU:OE1	1:D:162:GLU:N	2.53	0.42
1:E:56:LEU:O	1:E:59:SER:HB3	2.19	0.42
1:G:27:GLU:HG3	1:G:69:LEU:HD12	2.01	0.42
1:J:172:LYS:N	1:J:172:LYS:HD2	2.34	0.42
1:C:72:LEU:HD22	1:C:132:PHE:CD2	2.55	0.42
1:D:158:MET:HG2	1:L:165:LEU:HD12	2.01	0.41
1:I:79:ARG:HA	1:I:79:ARG:HD3	1.78	0.41
1:L:17:GLU:HG3	1:L:73:GLN:OE1	2.20	0.41
1:B:171:ASP:OD1	1:B:171:ASP:C	2.63	0.41
1:H:37:MET:O	1:H:41:PHE:HD2	2.03	0.41
1:B:26:LEU:HD12	1:B:26:LEU:HA	1.74	0.41
1:F:143:LYS:HB2	1:F:143:LYS:HE3	1.74	0.41
1:C:17:GLU:OE1	1:C:79:ARG:N	2.41	0.41
1:J:160:ALA:HB1	1:J:161:PRO:HA	2.01	0.41
1:D:169:LEU:HA	1:D:169:LEU:HD23	1.78	0.41
1:F:41:PHE:CZ	1:F:96:GLY:HA2	2.56	0.41
1:A:65:HIS:HB3	1:A:137:TYR:HE2	1.84	0.41
1:C:28:LEU:HD23	1:C:28:LEU:HA	1.77	0.41
1:G:65:HIS:HB3	1:G:137:TYR:HE2	1.85	0.41
1:C:27:GLU:HB2	1:C:66:ALA:HB2	2.03	0.41
1:B:129:LEU:O	1:B:133:ILE:HG12	2.21	0.41
1:E:71:LYS:HE3	1:E:71:LYS:HB3	1.83	0.41
1:G:97:LEU:O	1:G:101:GLU:HG3	2.21	0.41
1:J:162:GLU:HG2	1:J:163:SER:H	1.86	0.41
1:C:155:LEU:HD23	1:C:155:LEU:HA	1.84	0.41
1:F:141:GLN:O	1:F:145:ILE:HD12	2.20	0.41
1:B:145:ILE:HG22	1:G:8:VAL:HB	2.04	0.40
1:C:72:LEU:HD21	1:C:129:LEU:CD1	2.52	0.40
1:H:133:ILE:HG22	1:H:138:LEU:HG	2.04	0.40
1:I:66:ALA:O	1:I:70:MET:HG3	2.21	0.40
1:E:104:LEU:HD11	1:E:145:ILE:HG23	2.02	0.40
1:G:140:GLU:OE1	1:G:140:GLU:N	2.54	0.40
1:I:162:GLU:CD	1:I:162:GLU:H	2.30	0.40
1:J:161:PRO:HD2	1:J:162:GLU:OE1	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:59:SER:OG	1:F:63:ARG:NH2[14_444]	2.16	0.04
1:F:168:TYR:OH	1:H:174:THR:OG1[9_445]	2.16	0.04

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	170/183 (93%)	165 (97%)	5 (3%)	0	100	100
1	B	170/183 (93%)	166 (98%)	4 (2%)	0	100	100
1	C	170/183 (93%)	165 (97%)	5 (3%)	0	100	100
1	D	170/183 (93%)	166 (98%)	3 (2%)	1 (1%)	21	38
1	E	170/183 (93%)	168 (99%)	1 (1%)	1 (1%)	21	38
1	F	170/183 (93%)	164 (96%)	6 (4%)	0	100	100
1	G	170/183 (93%)	164 (96%)	6 (4%)	0	100	100
1	H	170/183 (93%)	162 (95%)	8 (5%)	0	100	100
1	I	170/183 (93%)	167 (98%)	3 (2%)	0	100	100
1	J	170/183 (93%)	165 (97%)	5 (3%)	0	100	100
1	K	170/183 (93%)	167 (98%)	3 (2%)	0	100	100
1	L	170/183 (93%)	167 (98%)	3 (2%)	0	100	100
All	All	2040/2196 (93%)	1986 (97%)	52 (2%)	2 (0%)	48	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	162	GLU
1	E	94	GLU

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	153/163 (94%)	153 (100%)	0	100	100
1	B	153/163 (94%)	153 (100%)	0	100	100
1	C	153/163 (94%)	153 (100%)	0	100	100
1	D	153/163 (94%)	153 (100%)	0	100	100
1	E	153/163 (94%)	153 (100%)	0	100	100
1	F	153/163 (94%)	152 (99%)	1 (1%)	76	89
1	G	153/163 (94%)	153 (100%)	0	100	100
1	H	153/163 (94%)	153 (100%)	0	100	100
1	I	153/163 (94%)	153 (100%)	0	100	100
1	J	153/163 (94%)	153 (100%)	0	100	100
1	K	153/163 (94%)	153 (100%)	0	100	100
1	L	153/163 (94%)	153 (100%)	0	100	100
All	All	1836/1956 (94%)	1835 (100%)	1 (0%)	88	96

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

Mol	Chain	Res	Type
1	F	171	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	GLN
1	A	98	ASN
1	A	173	HIS
1	B	98	ASN
1	B	141	GLN
1	C	14	GLN
1	C	109	ASN
1	D	7	GLN

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Mol	Chain	Res	Type
1	D	118	HIS
1	D	141	GLN
1	E	111	ASN
1	E	154	ASN
1	F	111	ASN
1	F	118	HIS
1	G	109	ASN
1	G	112	GLN
1	H	25	ASN
1	H	105	HIS
1	I	151	HIS
1	I	154	ASN
1	J	65	HIS
1	J	105	HIS
1	J	109	ASN
1	J	111	ASN
1	K	65	HIS
1	K	109	ASN
1	L	109	ASN
1	L	139	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 16 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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	172/183 (93%)	0.23	2 (1%) 76 73	10, 19, 35, 46	0
1	B	172/183 (93%)	0.24	3 (1%) 69 65	11, 17, 34, 48	0
1	C	172/183 (93%)	0.17	3 (1%) 69 65	10, 18, 31, 52	0
1	D	172/183 (93%)	0.30	8 (4%) 36 32	9, 18, 33, 58	0
1	E	172/183 (93%)	0.24	5 (2%) 53 49	12, 18, 33, 50	0
1	F	172/183 (93%)	0.26	6 (3%) 47 42	11, 19, 36, 51	0
1	G	172/183 (93%)	0.20	4 (2%) 61 57	10, 17, 32, 46	0
1	H	172/183 (93%)	0.10	3 (1%) 69 65	9, 17, 33, 47	0
1	I	172/183 (93%)	0.13	3 (1%) 69 65	10, 17, 33, 45	0
1	J	172/183 (93%)	0.22	6 (3%) 47 42	11, 18, 39, 51	0
1	K	172/183 (93%)	0.17	3 (1%) 69 65	9, 18, 34, 48	0
1	L	172/183 (93%)	0.22	6 (3%) 47 42	9, 18, 38, 47	0
All	All	2064/2196 (93%)	0.21	52 (2%) 58 54	9, 18, 35, 58	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	5	THR	5.3
1	D	91	ASP	5.2
1	J	5	THR	4.9
1	C	5	THR	3.9
1	I	91	ASP	3.8
1	J	91	ASP	3.8
1	C	92	ASP	3.6
1	L	5	THR	3.4
1	K	49	LYS	3.4
1	G	5	THR	3.3
1	G	91	ASP	3.3

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Mol	Chain	Res	Type	RSRZ
1	F	5	THR	3.3
1	K	91	ASP	3.2
1	E	91	ASP	3.1
1	B	5	THR	3.0
1	A	5	THR	2.9
1	G	123	PHE	2.9
1	F	94	GLU	2.8
1	H	91	ASP	2.8
1	E	63	ARG	2.8
1	L	91	ASP	2.8
1	B	91	ASP	2.7
1	L	81	PHE	2.7
1	I	5	THR	2.6
1	L	123	PHE	2.6
1	J	90	CYS	2.6
1	D	162	GLU	2.5
1	K	5	THR	2.5
1	E	176	GLY	2.4
1	J	89	ASP	2.4
1	D	90	CYS	2.4
1	C	102	CYS	2.4
1	J	92	ASP	2.4
1	D	176	GLY	2.4
1	H	5	THR	2.3
1	H	162	GLU	2.3
1	F	162	GLU	2.3
1	D	89	ASP	2.3
1	A	91	ASP	2.2
1	F	89	ASP	2.2
1	F	61	GLU	2.2
1	L	44	ASP	2.2
1	B	63	ARG	2.2
1	D	5	THR	2.1
1	D	63	ARG	2.1
1	D	87	LYS	2.1
1	L	90	CYS	2.1
1	J	162	GLU	2.0
1	F	91	ASP	2.0
1	I	89	ASP	2.0
1	E	90	CYS	2.0
1	G	162	GLU	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](#)

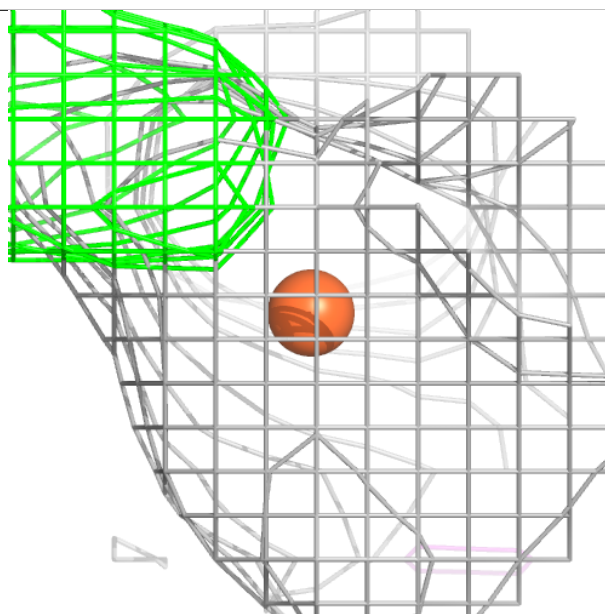
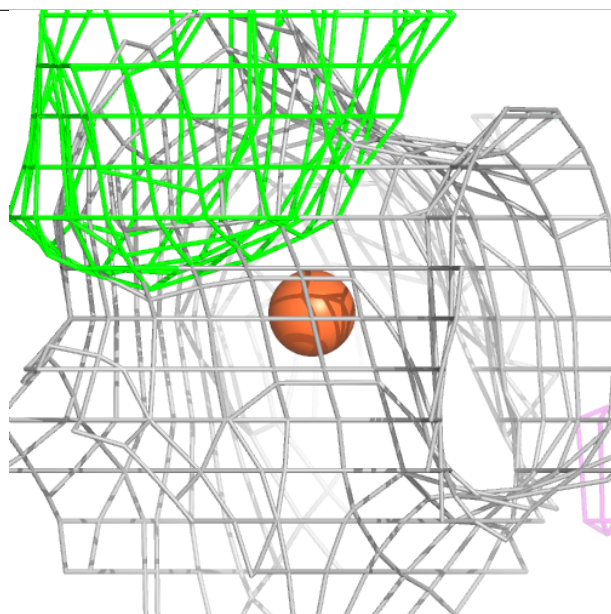
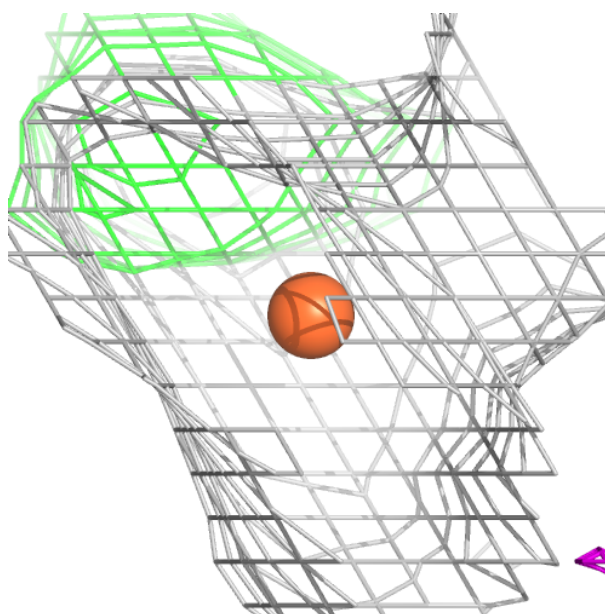
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	FE	B	202	1/1	0.88	0.09	49,49,49,49	0
2	FE	H	201	1/1	0.93	0.07	52,52,52,52	0
2	FE	A	201	1/1	0.94	0.05	49,49,49,49	0
2	FE	F	201	1/1	0.95	0.08	52,52,52,52	0
2	FE	G	201	1/1	0.95	0.04	43,43,43,43	0
2	FE	B	201	1/1	0.95	0.12	49,49,49,49	1
2	FE	E	201	1/1	0.96	0.05	43,43,43,43	0
2	FE	C	201	1/1	0.96	0.10	44,44,44,44	0
2	FE	K	201	1/1	0.96	0.04	40,40,40,40	0
2	FE	L	201	1/1	0.96	0.04	36,36,36,36	0
2	FE	C	202	1/1	0.97	0.04	55,55,55,55	0
2	FE	I	201	1/1	0.97	0.04	48,48,48,48	0
2	FE	D	201	1/1	0.98	0.03	42,42,42,42	0
2	FE	J	201	1/1	0.98	0.03	38,38,38,38	0
2	FE	D	202	1/1	0.98	0.13	50,50,50,50	1
2	FE	H	202	1/1	0.98	0.12	39,39,39,39	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.

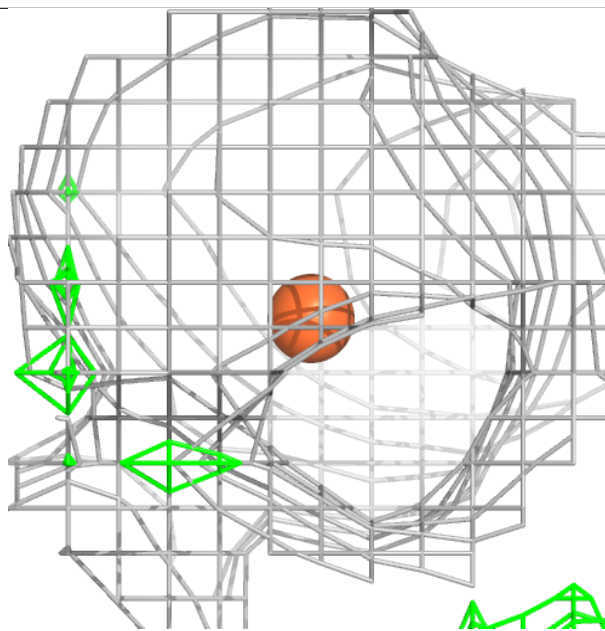
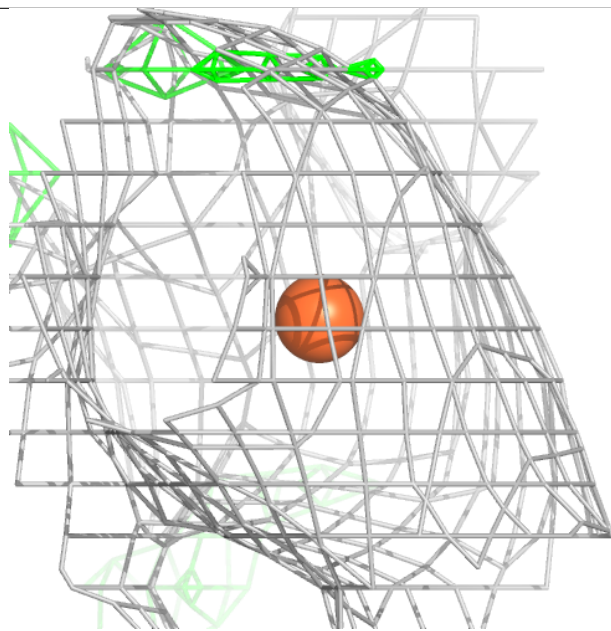
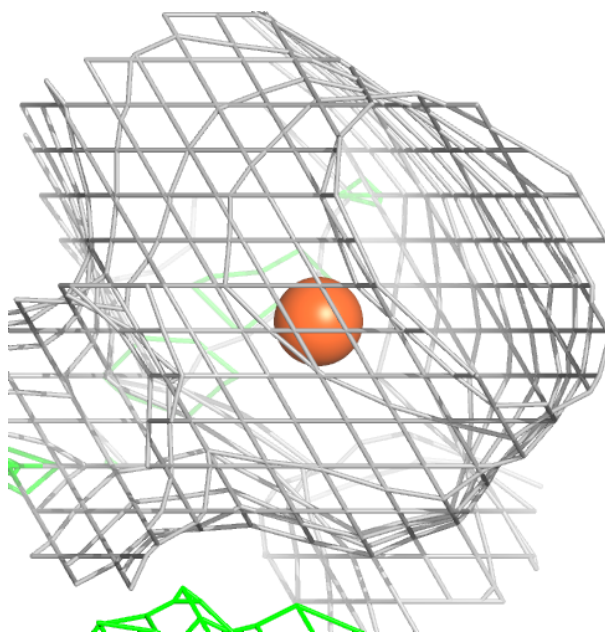
Electron density around FE B 202:

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



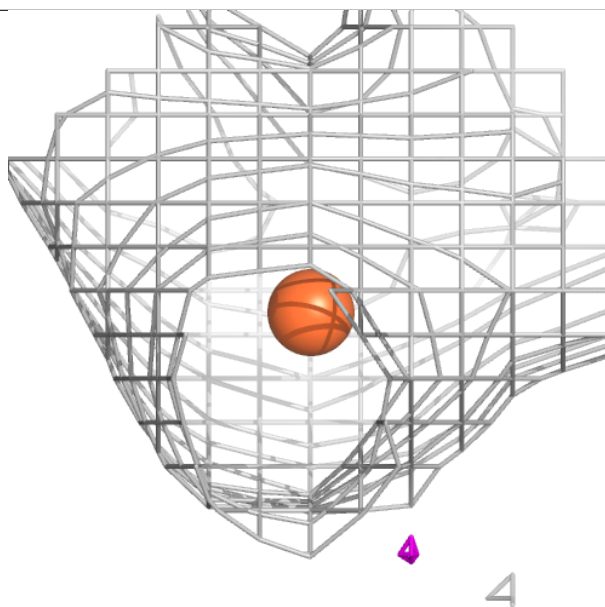
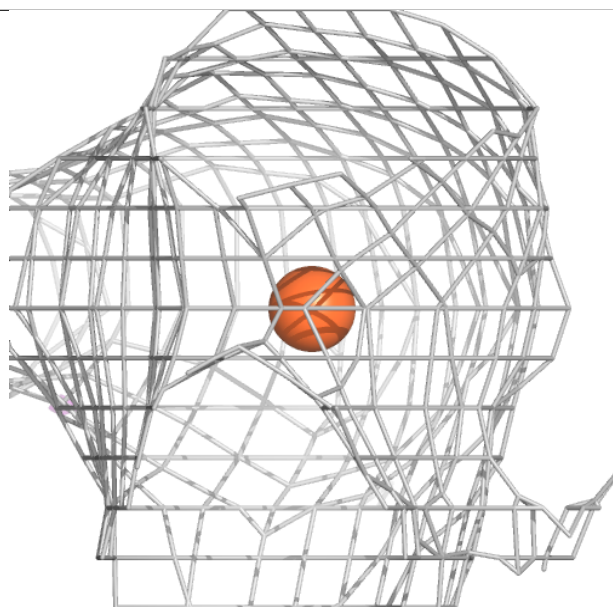
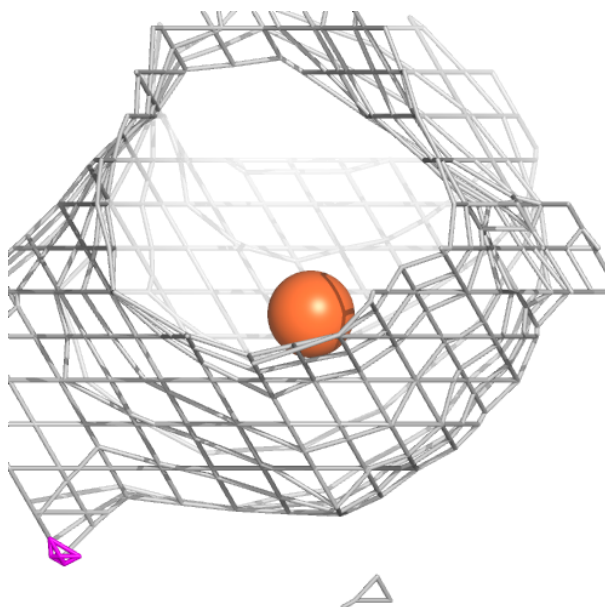
Electron density around FE H 201:

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and green (positive)



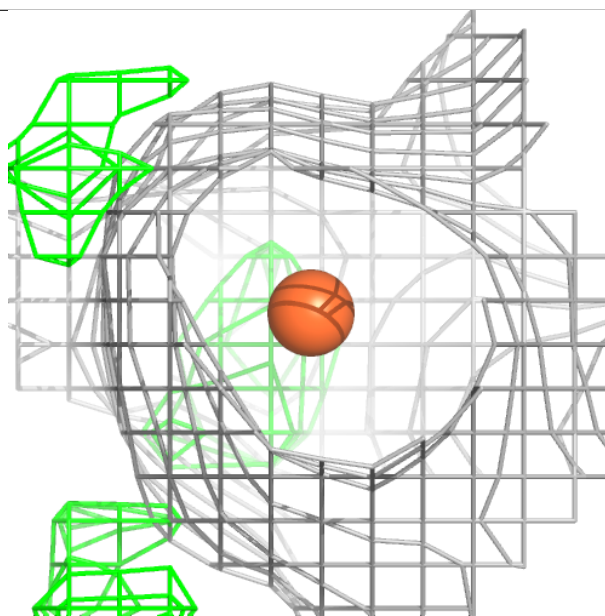
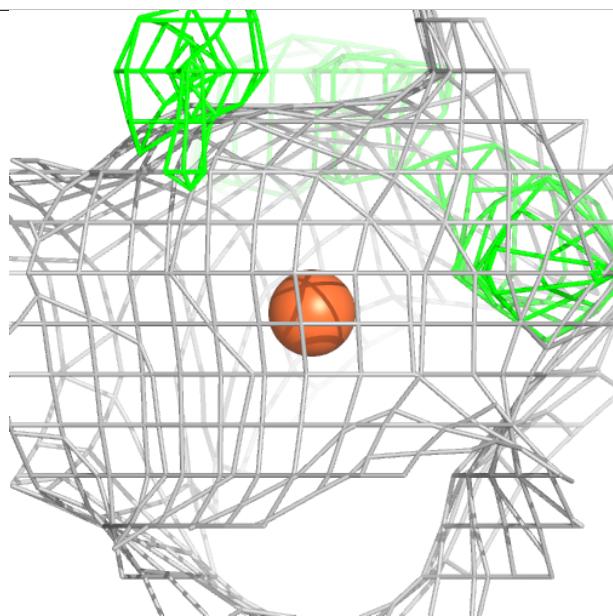
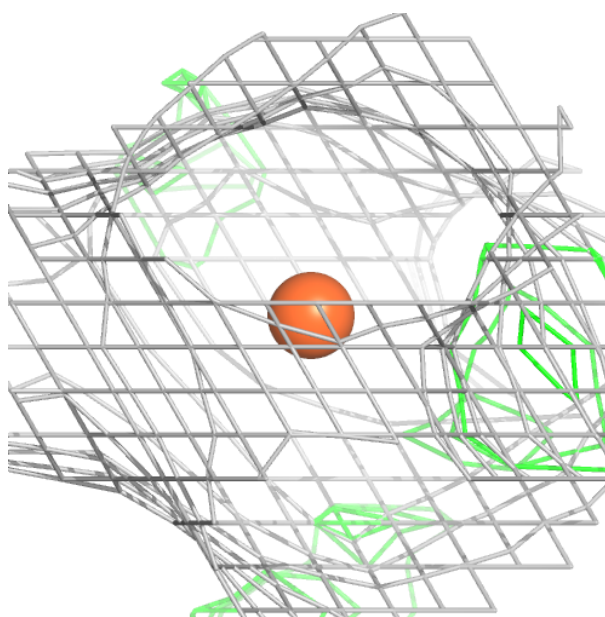
Electron density around FE A 201:

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



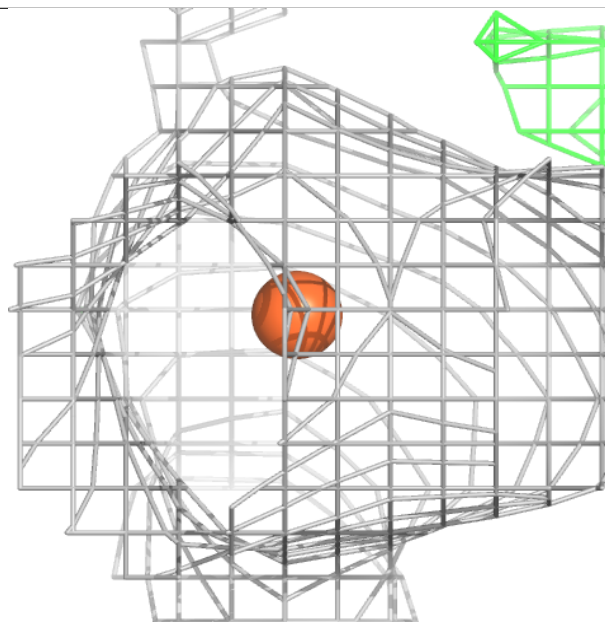
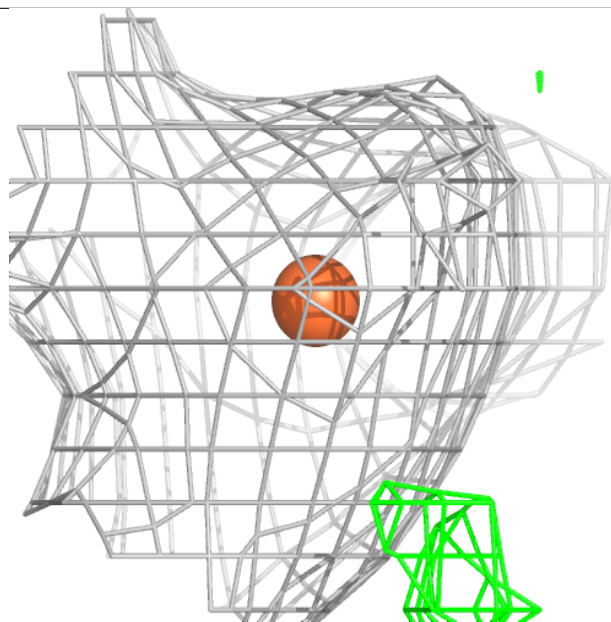
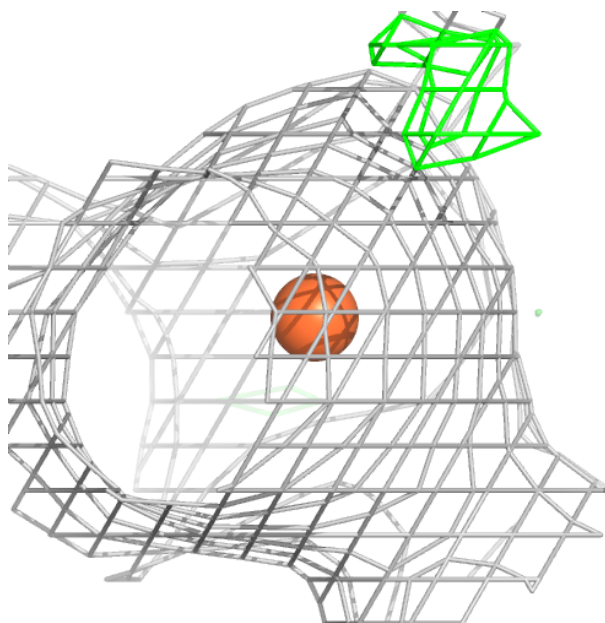
Electron density around FE F 201:

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



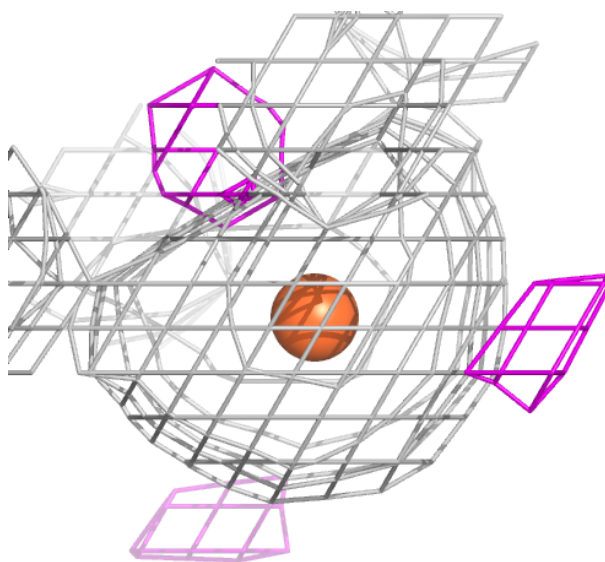
Electron density around FE G 201:

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and green (positive)

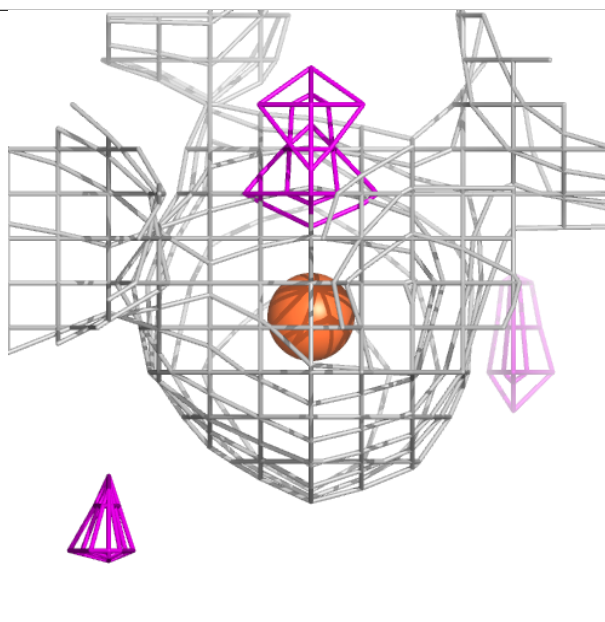
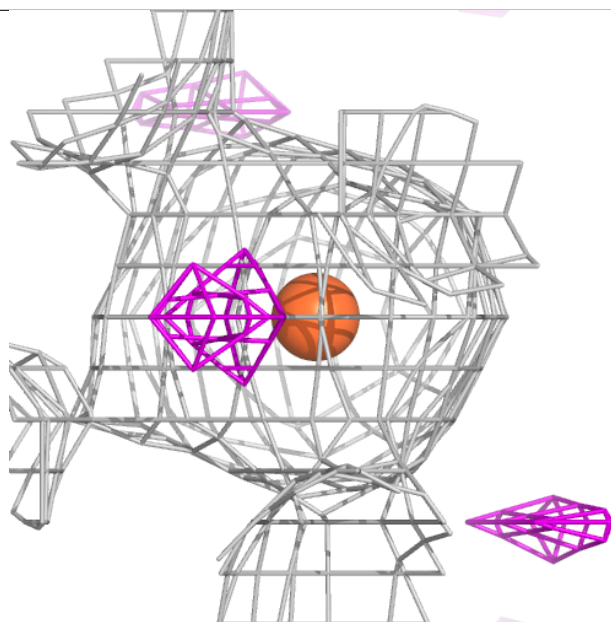


Electron density around FE B 201:

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



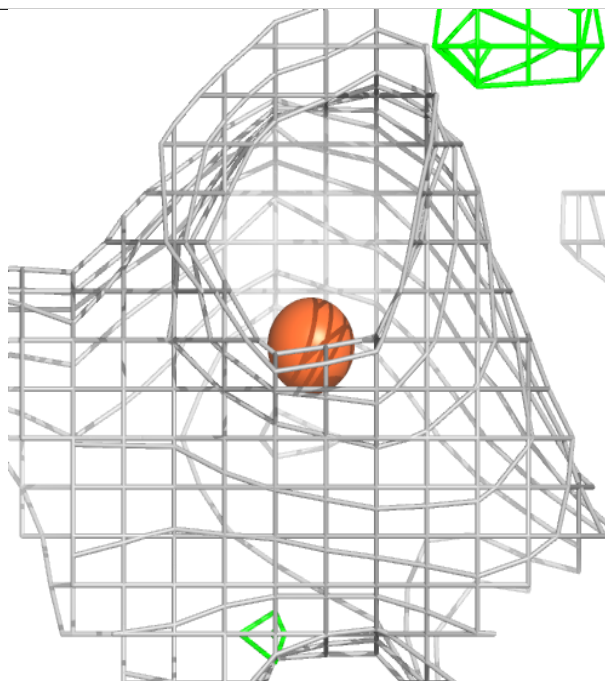
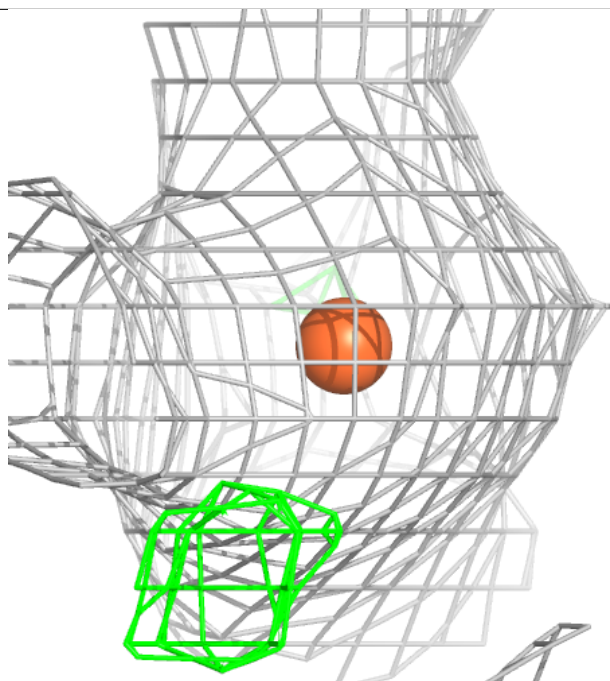
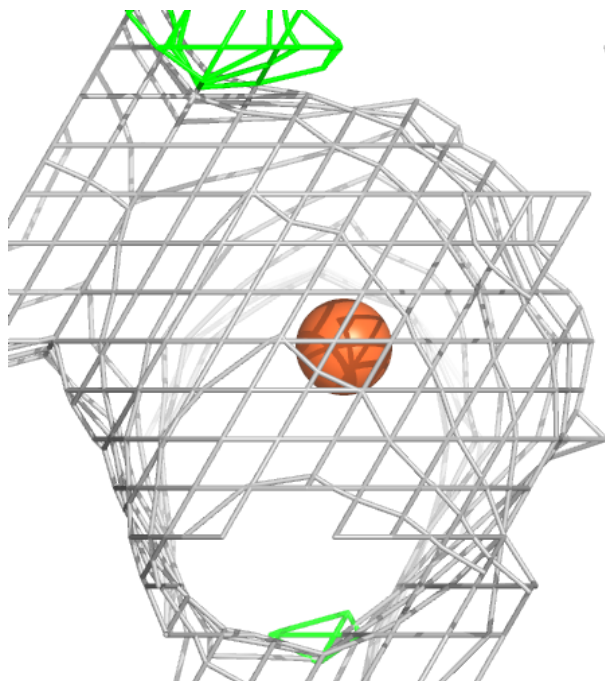
A



I

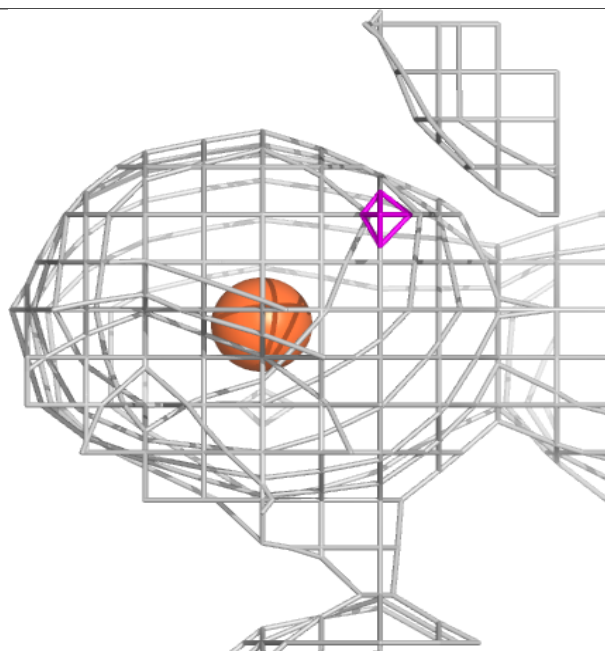
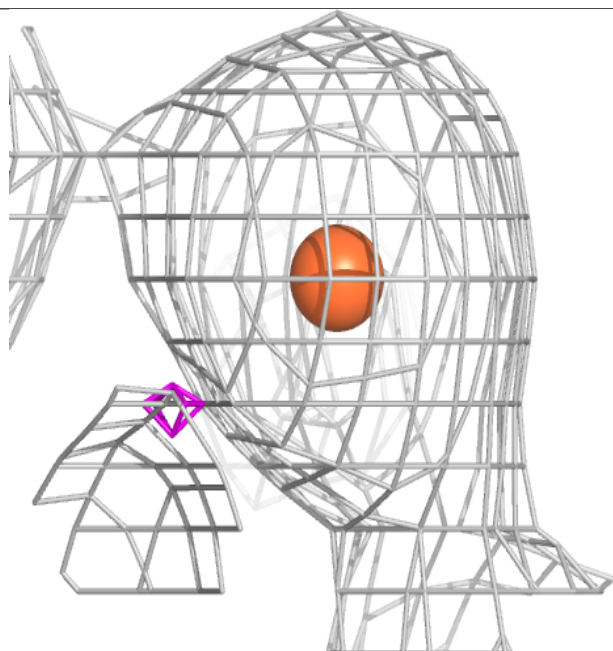
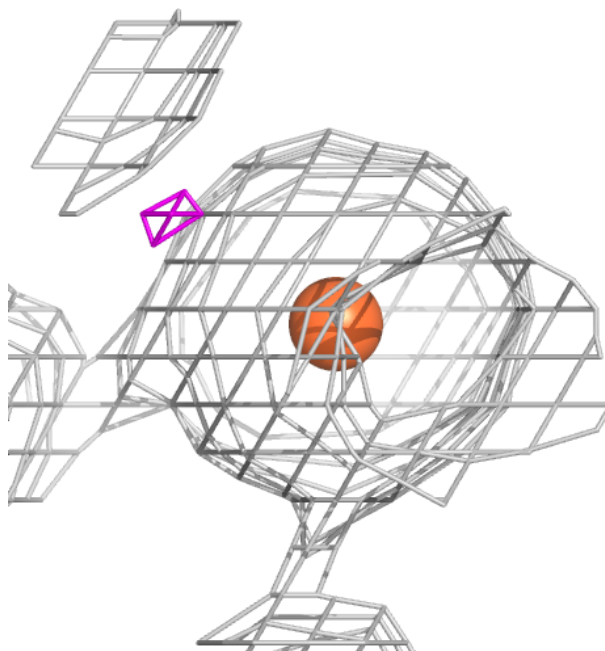
Electron density around FE E 201:

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



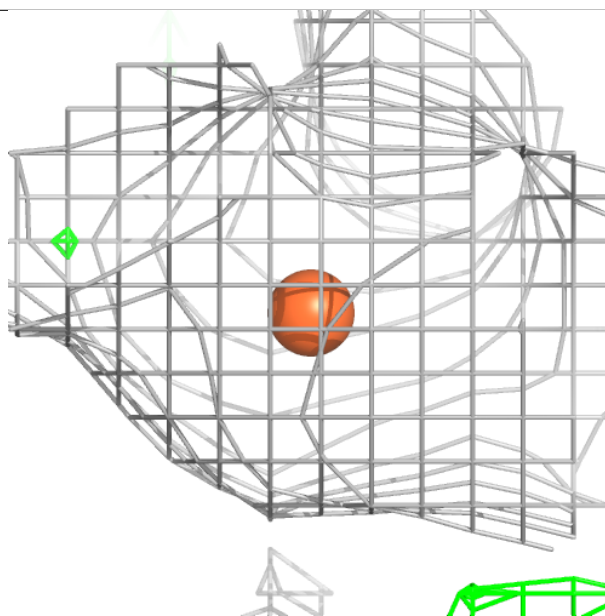
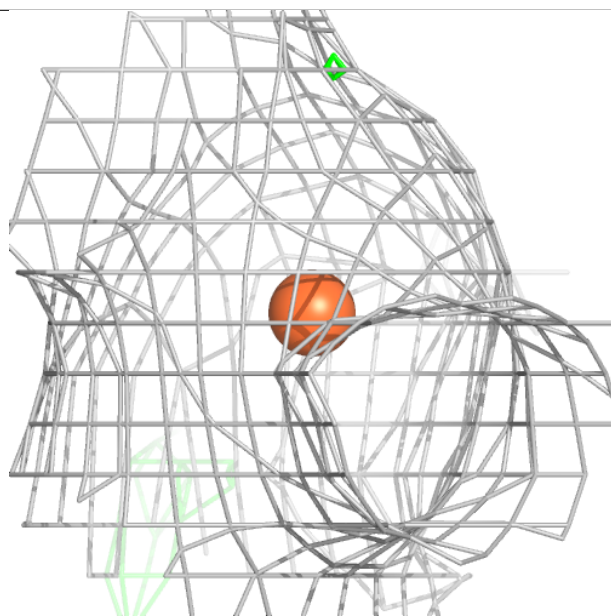
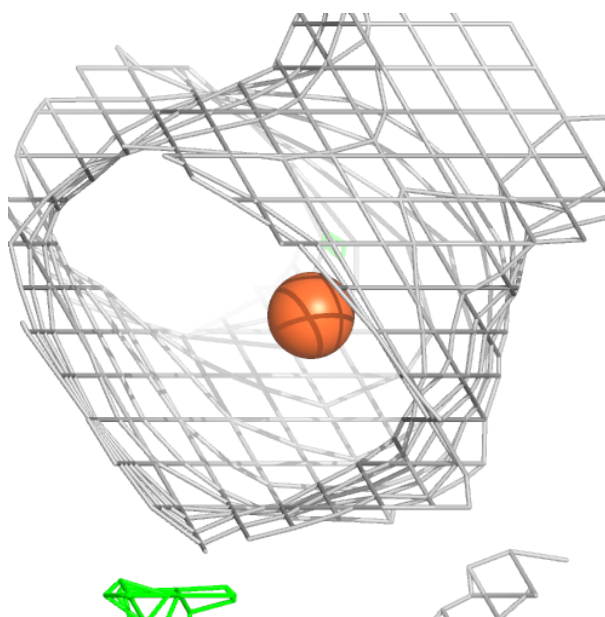
Electron density around FE C 201:

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



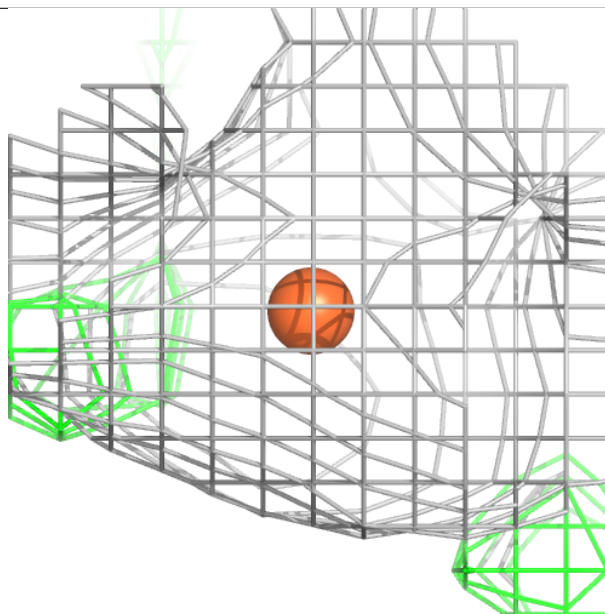
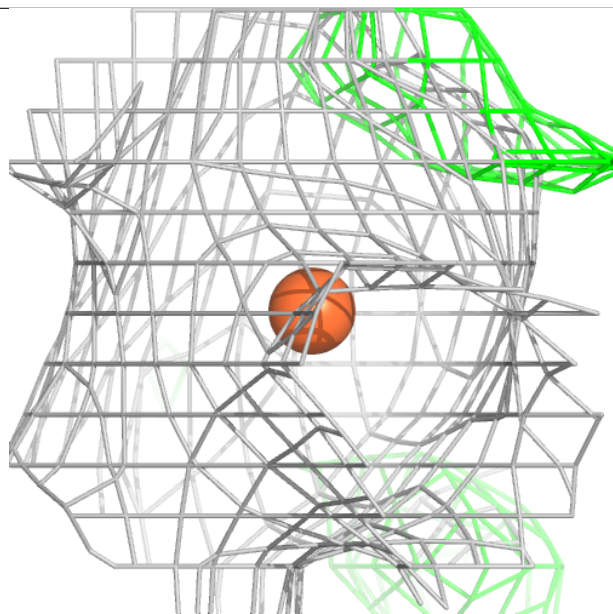
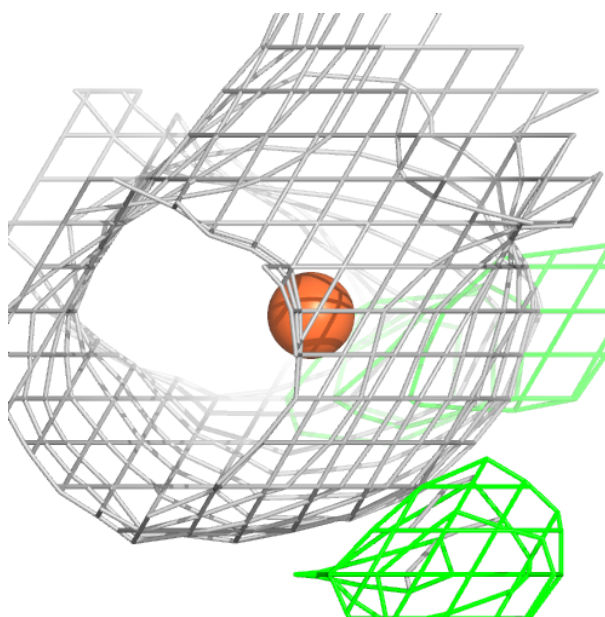
Electron density around FE K 201:

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



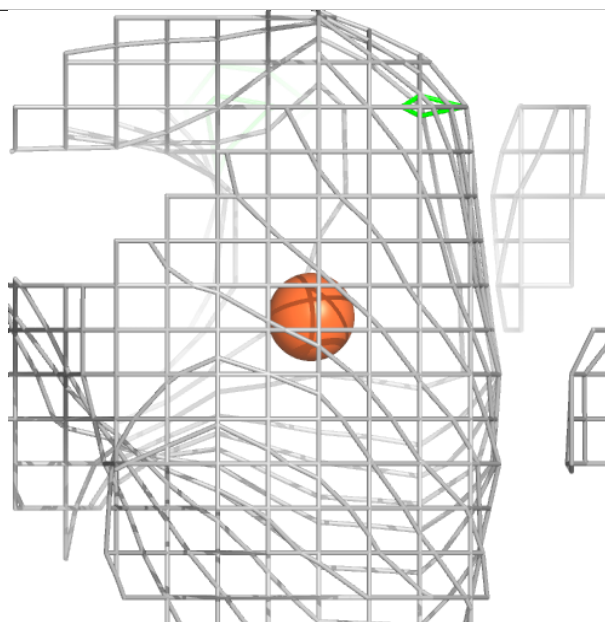
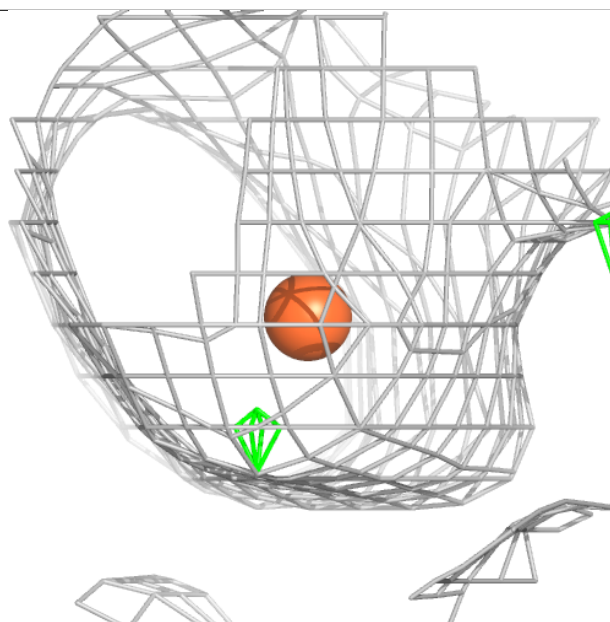
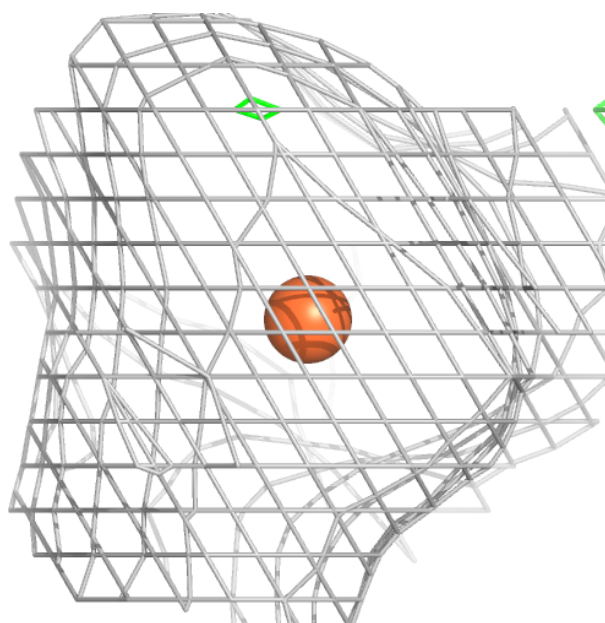
Electron density around FE L 201:

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



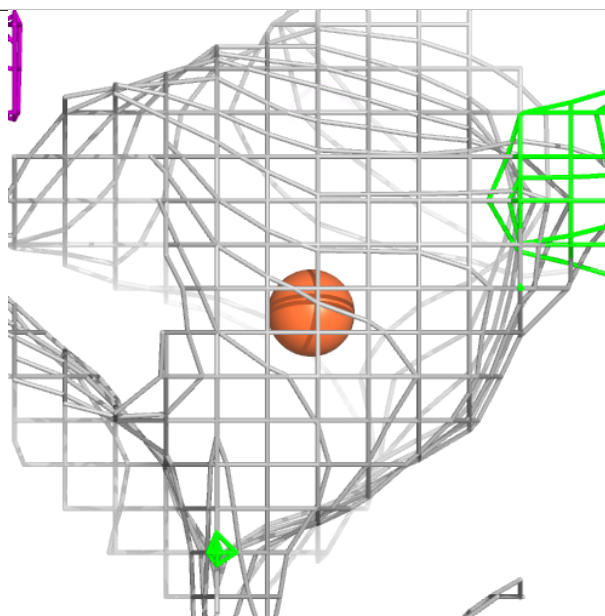
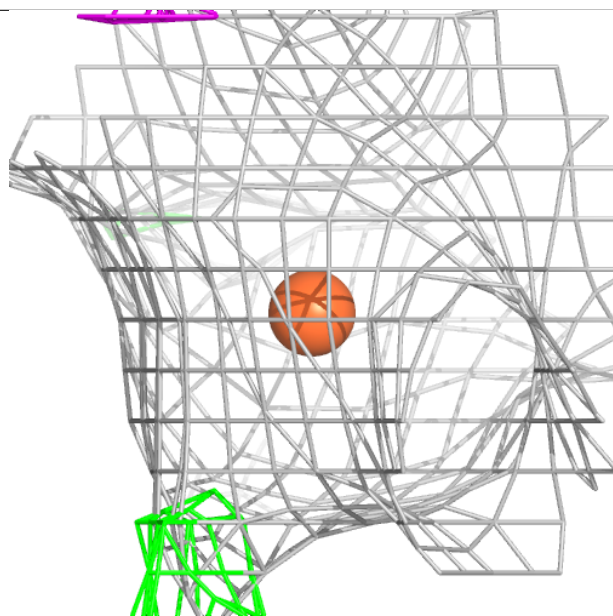
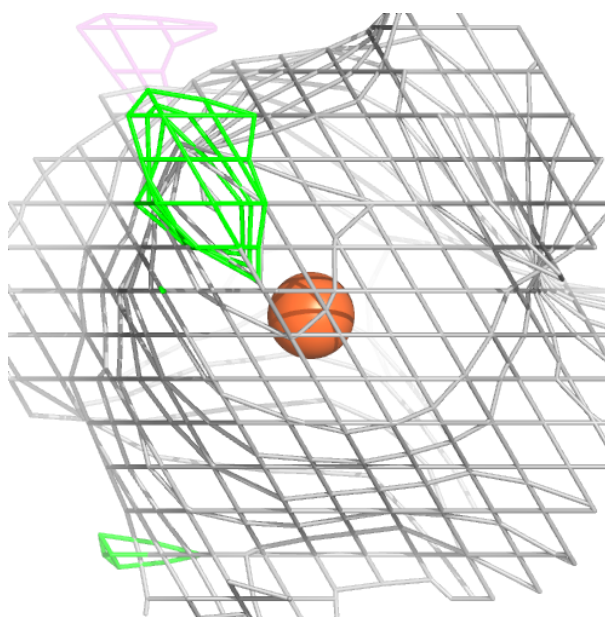
Electron density around FE C 202:

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



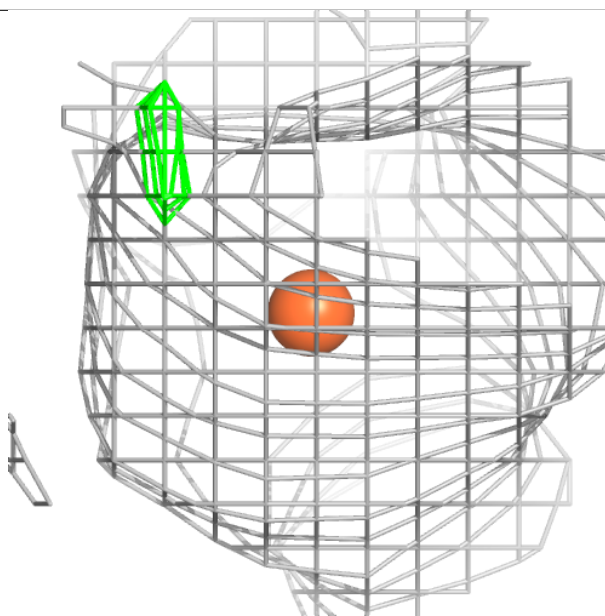
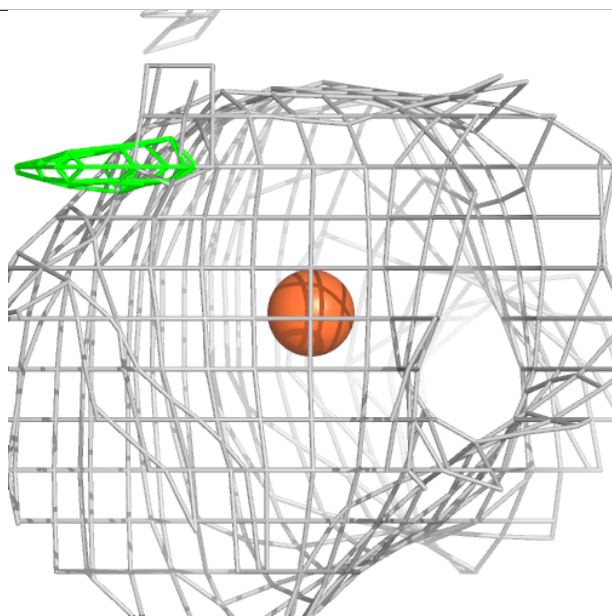
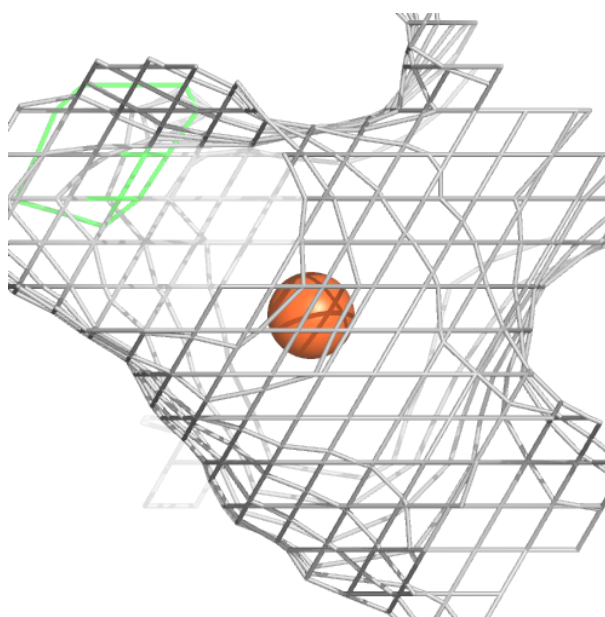
Electron density around FE I 201:

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



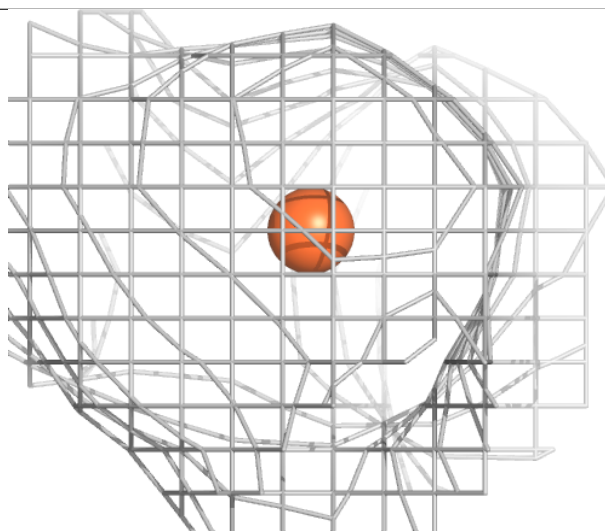
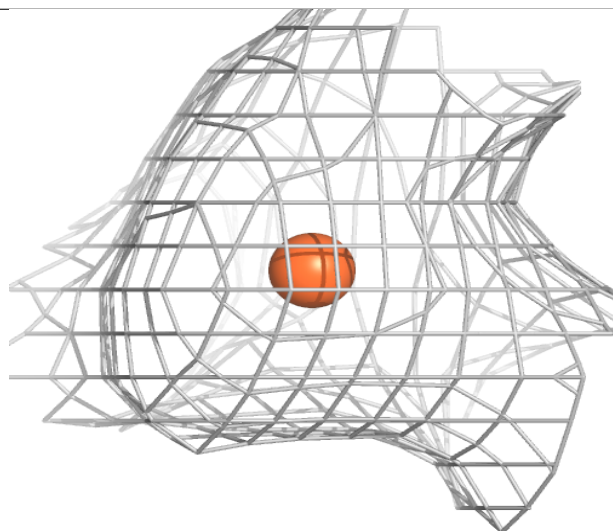
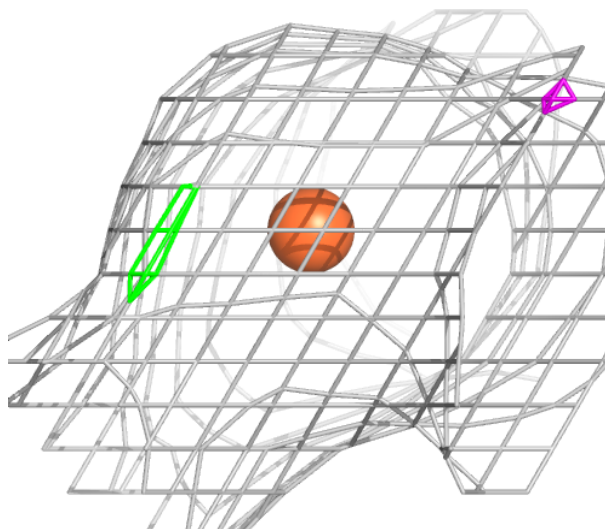
Electron density around FE D 201:

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



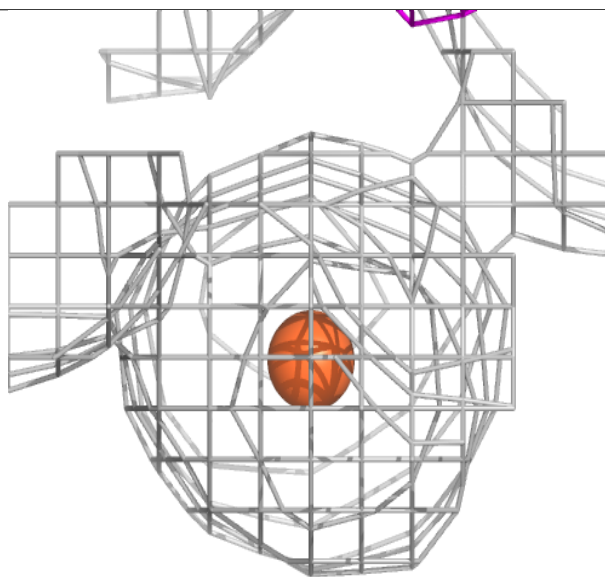
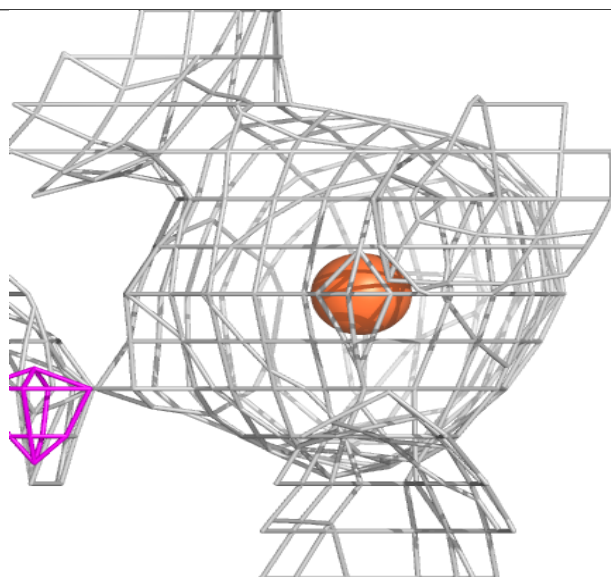
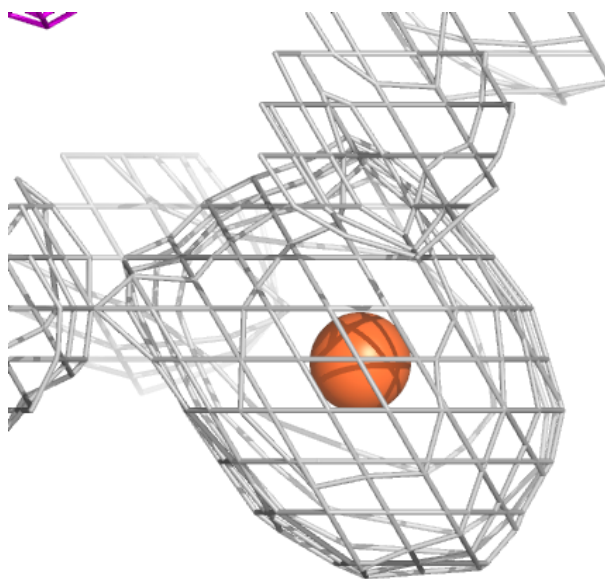
Electron density around FE J 201:

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and green (positive)



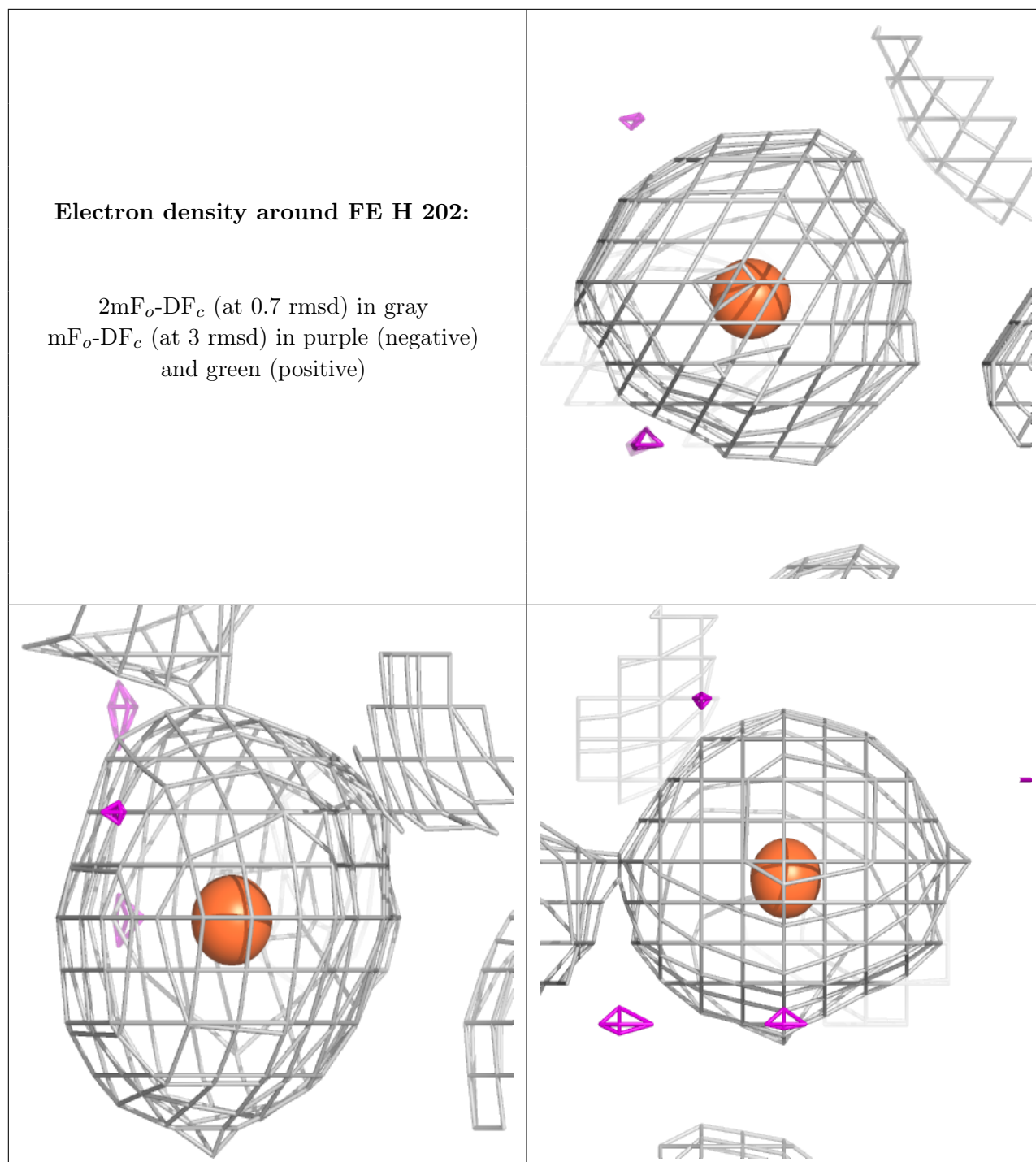
Electron density around FE D 202:

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



Electron density around FE H 202:

$2mF_o - DF_c$ (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.