wwPDB X-ray Structure Validation Summary Report (i)

May 25, 2022 – 01:22 am BST

PDB ID	:	7QWT
Title	:	Rieske non-heme iron monooxygenase for guaiacol O-demethylation
Authors	:	Hinchen, D.J.; Zahn, M.; Bleem, A.; Beckham, G.T.; McGeehan, J.E.
Deposited on	:	2022-01-25
Resolution	:	3.71 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

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

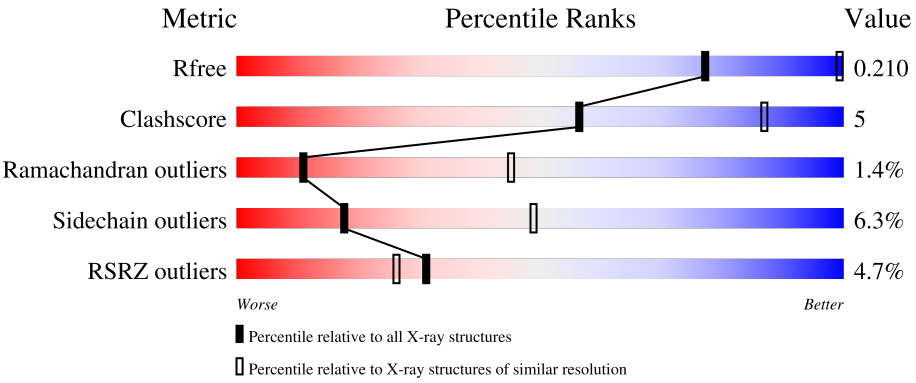
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.28.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.28.1

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 3.71 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}	130704	1089 (3.90-3.54)
Clashscore	141614	1012 (3.88-3.56)
Ramachandran outliers	138981	1114 (3.90-3.54)
Sidechain outliers	138945	1110 (3.90-3.54)
RSRZ outliers	127900	1020 (3.90-3.54)

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 <=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	373	80% 14% • 6%
1	B	373	% 77% 16% • 6%
1	C	373	78% 14% • 6%
1	D	373	8% 79% 14% • 6%
1	E	373	7% 81% 12% • 6%

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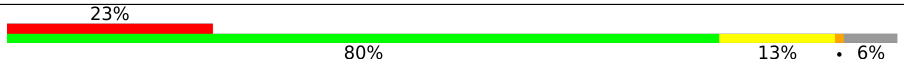
Mol	Chain	Length	Quality of chain
1	F	373	
1	G	373	
1	H	373	
1	I	373	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	FES	C	401	-	-	X	-
2	FES	F	401	-	-	X	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 49487 atoms, of which 24162 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 Rieske (2Fe-2S) domain protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	352	Total	C	H	N	O	S	84	0	0
			5499	1766	2687	505	523	18			
1	B	351	Total	C	H	N	O	S	84	0	0
			5492	1764	2684	504	522	18			
1	C	351	Total	C	H	N	O	S	84	0	0
			5492	1764	2684	504	522	18			
1	D	351	Total	C	H	N	O	S	84	0	0
			5492	1764	2684	504	522	18			
1	E	352	Total	C	H	N	O	S	84	0	0
			5499	1766	2687	505	523	18			
1	F	351	Total	C	H	N	O	S	84	0	0
			5492	1764	2684	504	522	18			
1	G	351	Total	C	H	N	O	S	84	0	0
			5492	1764	2684	504	522	18			
1	H	351	Total	C	H	N	O	S	84	0	0
			5492	1764	2684	504	522	18			
1	I	351	Total	C	H	N	O	S	84	0	0
			5492	1764	2684	504	522	18			

There are 180 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A5VBE6
A	-18	GLY	-	expression tag	UNP A5VBE6
A	-17	SER	-	expression tag	UNP A5VBE6
A	-16	SER	-	expression tag	UNP A5VBE6
A	-15	HIS	-	expression tag	UNP A5VBE6
A	-14	HIS	-	expression tag	UNP A5VBE6
A	-13	HIS	-	expression tag	UNP A5VBE6
A	-12	HIS	-	expression tag	UNP A5VBE6
A	-11	HIS	-	expression tag	UNP A5VBE6
A	-10	HIS	-	expression tag	UNP A5VBE6
A	-9	SER	-	expression tag	UNP A5VBE6

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	SER	-	expression tag	UNP A5VBE6
A	-7	GLY	-	expression tag	UNP A5VBE6
A	-6	LEU	-	expression tag	UNP A5VBE6
A	-5	VAL	-	expression tag	UNP A5VBE6
A	-4	PRO	-	expression tag	UNP A5VBE6
A	-3	ARG	-	expression tag	UNP A5VBE6
A	-2	GLY	-	expression tag	UNP A5VBE6
A	-1	SER	-	expression tag	UNP A5VBE6
A	0	HIS	-	expression tag	UNP A5VBE6
B	-19	MET	-	initiating methionine	UNP A5VBE6
B	-18	GLY	-	expression tag	UNP A5VBE6
B	-17	SER	-	expression tag	UNP A5VBE6
B	-16	SER	-	expression tag	UNP A5VBE6
B	-15	HIS	-	expression tag	UNP A5VBE6
B	-14	HIS	-	expression tag	UNP A5VBE6
B	-13	HIS	-	expression tag	UNP A5VBE6
B	-12	HIS	-	expression tag	UNP A5VBE6
B	-11	HIS	-	expression tag	UNP A5VBE6
B	-10	HIS	-	expression tag	UNP A5VBE6
B	-9	SER	-	expression tag	UNP A5VBE6
B	-8	SER	-	expression tag	UNP A5VBE6
B	-7	GLY	-	expression tag	UNP A5VBE6
B	-6	LEU	-	expression tag	UNP A5VBE6
B	-5	VAL	-	expression tag	UNP A5VBE6
B	-4	PRO	-	expression tag	UNP A5VBE6
B	-3	ARG	-	expression tag	UNP A5VBE6
B	-2	GLY	-	expression tag	UNP A5VBE6
B	-1	SER	-	expression tag	UNP A5VBE6
B	0	HIS	-	expression tag	UNP A5VBE6
C	-19	MET	-	initiating methionine	UNP A5VBE6
C	-18	GLY	-	expression tag	UNP A5VBE6
C	-17	SER	-	expression tag	UNP A5VBE6
C	-16	SER	-	expression tag	UNP A5VBE6
C	-15	HIS	-	expression tag	UNP A5VBE6
C	-14	HIS	-	expression tag	UNP A5VBE6
C	-13	HIS	-	expression tag	UNP A5VBE6
C	-12	HIS	-	expression tag	UNP A5VBE6
C	-11	HIS	-	expression tag	UNP A5VBE6
C	-10	HIS	-	expression tag	UNP A5VBE6
C	-9	SER	-	expression tag	UNP A5VBE6
C	-8	SER	-	expression tag	UNP A5VBE6
C	-7	GLY	-	expression tag	UNP A5VBE6

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	LEU	-	expression tag	UNP A5VBE6
C	-5	VAL	-	expression tag	UNP A5VBE6
C	-4	PRO	-	expression tag	UNP A5VBE6
C	-3	ARG	-	expression tag	UNP A5VBE6
C	-2	GLY	-	expression tag	UNP A5VBE6
C	-1	SER	-	expression tag	UNP A5VBE6
C	0	HIS	-	expression tag	UNP A5VBE6
D	-19	MET	-	initiating methionine	UNP A5VBE6
D	-18	GLY	-	expression tag	UNP A5VBE6
D	-17	SER	-	expression tag	UNP A5VBE6
D	-16	SER	-	expression tag	UNP A5VBE6
D	-15	HIS	-	expression tag	UNP A5VBE6
D	-14	HIS	-	expression tag	UNP A5VBE6
D	-13	HIS	-	expression tag	UNP A5VBE6
D	-12	HIS	-	expression tag	UNP A5VBE6
D	-11	HIS	-	expression tag	UNP A5VBE6
D	-10	HIS	-	expression tag	UNP A5VBE6
D	-9	SER	-	expression tag	UNP A5VBE6
D	-8	SER	-	expression tag	UNP A5VBE6
D	-7	GLY	-	expression tag	UNP A5VBE6
D	-6	LEU	-	expression tag	UNP A5VBE6
D	-5	VAL	-	expression tag	UNP A5VBE6
D	-4	PRO	-	expression tag	UNP A5VBE6
D	-3	ARG	-	expression tag	UNP A5VBE6
D	-2	GLY	-	expression tag	UNP A5VBE6
D	-1	SER	-	expression tag	UNP A5VBE6
D	0	HIS	-	expression tag	UNP A5VBE6
E	-19	MET	-	initiating methionine	UNP A5VBE6
E	-18	GLY	-	expression tag	UNP A5VBE6
E	-17	SER	-	expression tag	UNP A5VBE6
E	-16	SER	-	expression tag	UNP A5VBE6
E	-15	HIS	-	expression tag	UNP A5VBE6
E	-14	HIS	-	expression tag	UNP A5VBE6
E	-13	HIS	-	expression tag	UNP A5VBE6
E	-12	HIS	-	expression tag	UNP A5VBE6
E	-11	HIS	-	expression tag	UNP A5VBE6
E	-10	HIS	-	expression tag	UNP A5VBE6
E	-9	SER	-	expression tag	UNP A5VBE6
E	-8	SER	-	expression tag	UNP A5VBE6
E	-7	GLY	-	expression tag	UNP A5VBE6
E	-6	LEU	-	expression tag	UNP A5VBE6
E	-5	VAL	-	expression tag	UNP A5VBE6

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-4	PRO	-	expression tag	UNP A5VBE6
E	-3	ARG	-	expression tag	UNP A5VBE6
E	-2	GLY	-	expression tag	UNP A5VBE6
E	-1	SER	-	expression tag	UNP A5VBE6
E	0	HIS	-	expression tag	UNP A5VBE6
F	-19	MET	-	initiating methionine	UNP A5VBE6
F	-18	GLY	-	expression tag	UNP A5VBE6
F	-17	SER	-	expression tag	UNP A5VBE6
F	-16	SER	-	expression tag	UNP A5VBE6
F	-15	HIS	-	expression tag	UNP A5VBE6
F	-14	HIS	-	expression tag	UNP A5VBE6
F	-13	HIS	-	expression tag	UNP A5VBE6
F	-12	HIS	-	expression tag	UNP A5VBE6
F	-11	HIS	-	expression tag	UNP A5VBE6
F	-10	HIS	-	expression tag	UNP A5VBE6
F	-9	SER	-	expression tag	UNP A5VBE6
F	-8	SER	-	expression tag	UNP A5VBE6
F	-7	GLY	-	expression tag	UNP A5VBE6
F	-6	LEU	-	expression tag	UNP A5VBE6
F	-5	VAL	-	expression tag	UNP A5VBE6
F	-4	PRO	-	expression tag	UNP A5VBE6
F	-3	ARG	-	expression tag	UNP A5VBE6
F	-2	GLY	-	expression tag	UNP A5VBE6
F	-1	SER	-	expression tag	UNP A5VBE6
F	0	HIS	-	expression tag	UNP A5VBE6
G	-19	MET	-	initiating methionine	UNP A5VBE6
G	-18	GLY	-	expression tag	UNP A5VBE6
G	-17	SER	-	expression tag	UNP A5VBE6
G	-16	SER	-	expression tag	UNP A5VBE6
G	-15	HIS	-	expression tag	UNP A5VBE6
G	-14	HIS	-	expression tag	UNP A5VBE6
G	-13	HIS	-	expression tag	UNP A5VBE6
G	-12	HIS	-	expression tag	UNP A5VBE6
G	-11	HIS	-	expression tag	UNP A5VBE6
G	-10	HIS	-	expression tag	UNP A5VBE6
G	-9	SER	-	expression tag	UNP A5VBE6
G	-8	SER	-	expression tag	UNP A5VBE6
G	-7	GLY	-	expression tag	UNP A5VBE6
G	-6	LEU	-	expression tag	UNP A5VBE6
G	-5	VAL	-	expression tag	UNP A5VBE6
G	-4	PRO	-	expression tag	UNP A5VBE6
G	-3	ARG	-	expression tag	UNP A5VBE6

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-2	GLY	-	expression tag	UNP A5VBE6
G	-1	SER	-	expression tag	UNP A5VBE6
G	0	HIS	-	expression tag	UNP A5VBE6
H	-19	MET	-	initiating methionine	UNP A5VBE6
H	-18	GLY	-	expression tag	UNP A5VBE6
H	-17	SER	-	expression tag	UNP A5VBE6
H	-16	SER	-	expression tag	UNP A5VBE6
H	-15	HIS	-	expression tag	UNP A5VBE6
H	-14	HIS	-	expression tag	UNP A5VBE6
H	-13	HIS	-	expression tag	UNP A5VBE6
H	-12	HIS	-	expression tag	UNP A5VBE6
H	-11	HIS	-	expression tag	UNP A5VBE6
H	-10	HIS	-	expression tag	UNP A5VBE6
H	-9	SER	-	expression tag	UNP A5VBE6
H	-8	SER	-	expression tag	UNP A5VBE6
H	-7	GLY	-	expression tag	UNP A5VBE6
H	-6	LEU	-	expression tag	UNP A5VBE6
H	-5	VAL	-	expression tag	UNP A5VBE6
H	-4	PRO	-	expression tag	UNP A5VBE6
H	-3	ARG	-	expression tag	UNP A5VBE6
H	-2	GLY	-	expression tag	UNP A5VBE6
H	-1	SER	-	expression tag	UNP A5VBE6
H	0	HIS	-	expression tag	UNP A5VBE6
I	-19	MET	-	initiating methionine	UNP A5VBE6
I	-18	GLY	-	expression tag	UNP A5VBE6
I	-17	SER	-	expression tag	UNP A5VBE6
I	-16	SER	-	expression tag	UNP A5VBE6
I	-15	HIS	-	expression tag	UNP A5VBE6
I	-14	HIS	-	expression tag	UNP A5VBE6
I	-13	HIS	-	expression tag	UNP A5VBE6
I	-12	HIS	-	expression tag	UNP A5VBE6
I	-11	HIS	-	expression tag	UNP A5VBE6
I	-10	HIS	-	expression tag	UNP A5VBE6
I	-9	SER	-	expression tag	UNP A5VBE6
I	-8	SER	-	expression tag	UNP A5VBE6
I	-7	GLY	-	expression tag	UNP A5VBE6
I	-6	LEU	-	expression tag	UNP A5VBE6
I	-5	VAL	-	expression tag	UNP A5VBE6
I	-4	PRO	-	expression tag	UNP A5VBE6
I	-3	ARG	-	expression tag	UNP A5VBE6
I	-2	GLY	-	expression tag	UNP A5VBE6
I	-1	SER	-	expression tag	UNP A5VBE6

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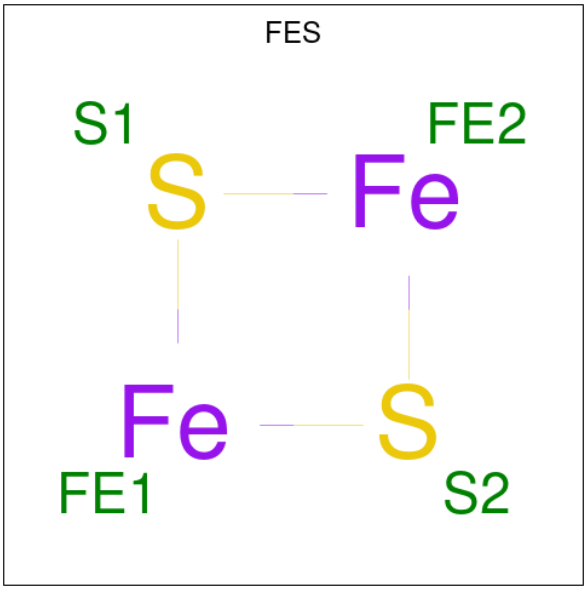
Chain	Residue	Modelled	Actual	Comment	Reference
I	0	HIS	-	expression tag	UNP A5VBE6

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe_2S_2) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe S 4 2 2	0	0
2	B	1	Total Fe S 4 2 2	0	0
2	C	1	Total Fe S 4 2 2	0	0
2	D	1	Total Fe S 4 2 2	0	0
2	E	1	Total Fe S 4 2 2	0	0
2	F	1	Total Fe S 4 2 2	0	0
2	G	1	Total Fe S 4 2 2	0	0
2	H	1	Total Fe S 4 2 2	0	0
2	I	1	Total Fe S 4 2 2	0	0

- Molecule 3 is FE (III) ION (three-letter code: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

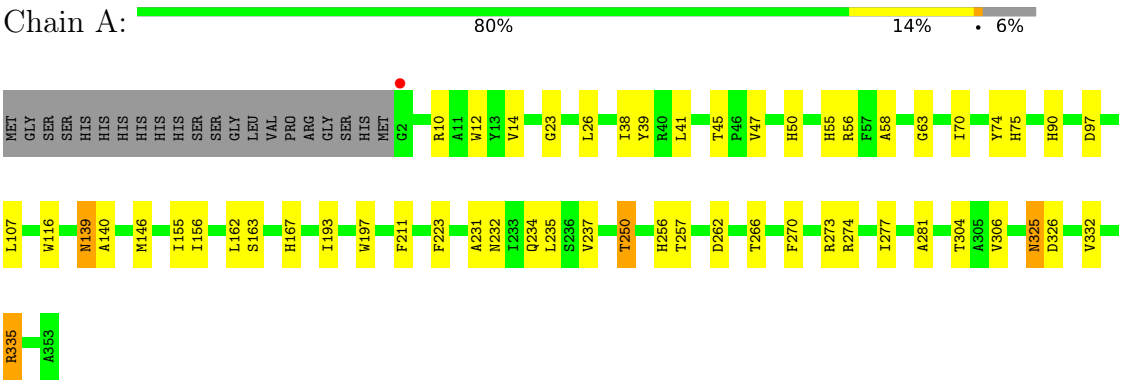
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Fe 1 1	0	0
3	B	1	Total Fe 1 1	0	0
3	C	1	Total Fe 1 1	0	0
3	D	1	Total Fe 1 1	0	0
3	E	1	Total Fe 1 1	0	0
3	F	1	Total Fe 1 1	0	0
3	G	1	Total Fe 1 1	0	0
3	H	1	Total Fe 1 1	0	0
3	I	1	Total Fe 1 1	0	0

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.

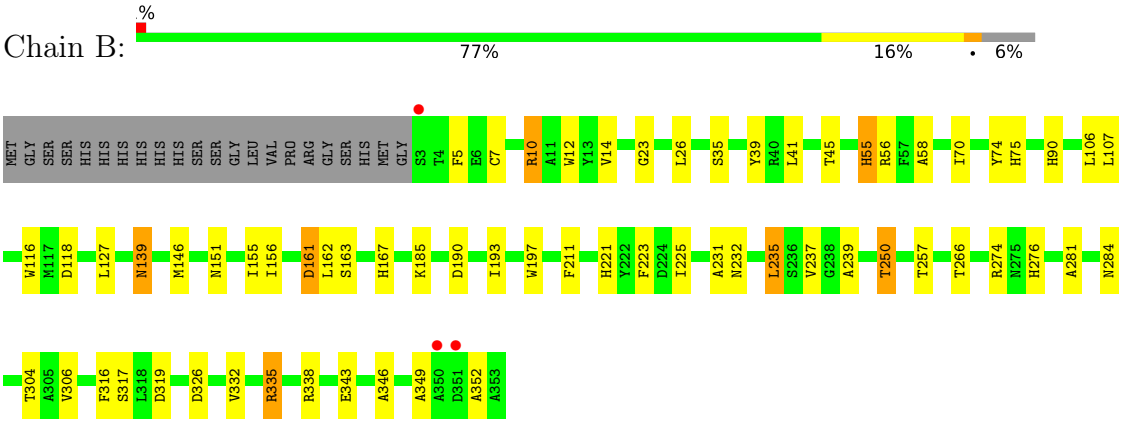
• Molecule 1: Rieske (2Fe-2S) domain protein

Chain A:



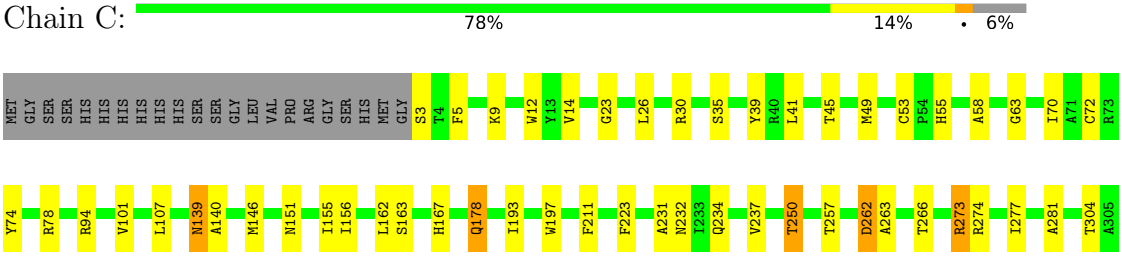
• Molecule 1: Rieske (2Fe-2S) domain protein

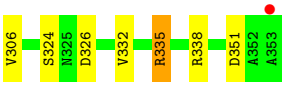
Chain B:



• Molecule 1: Rieske (2Fe-2S) domain protein

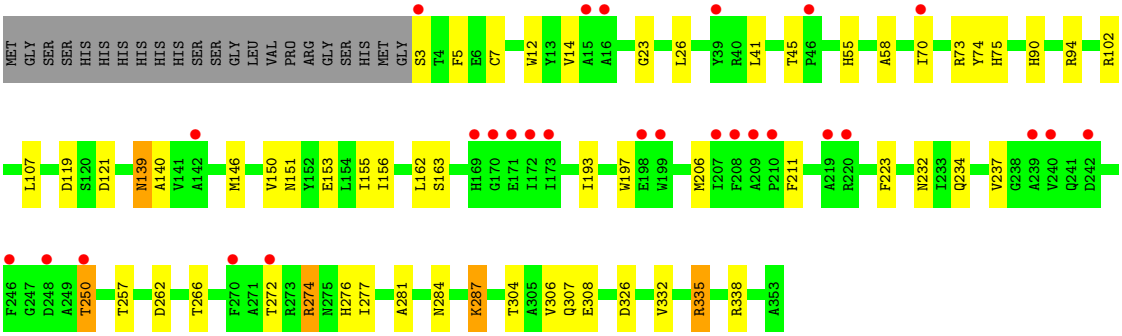
Chain C:





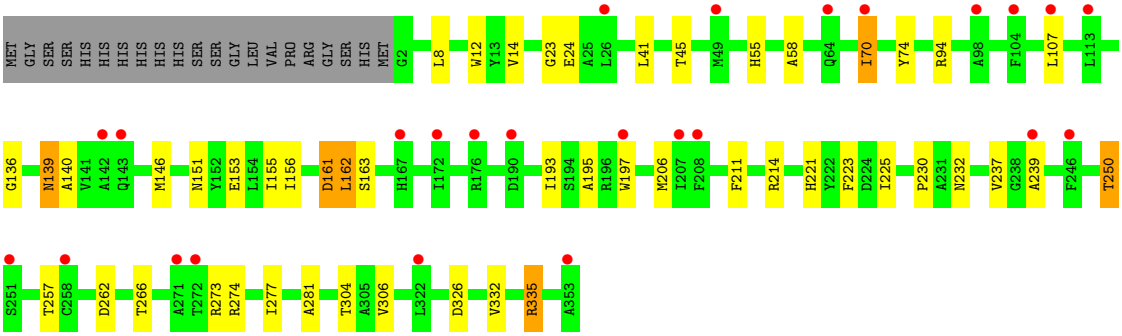
• Molecule 1: Rieske (2Fe-2S) domain protein

Chain D:



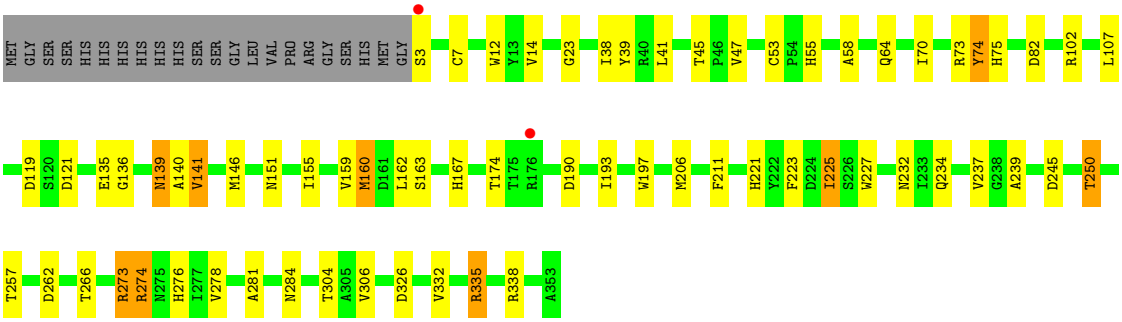
• Molecule 1: Rieske (2Fe-2S) domain protein

Chain E:



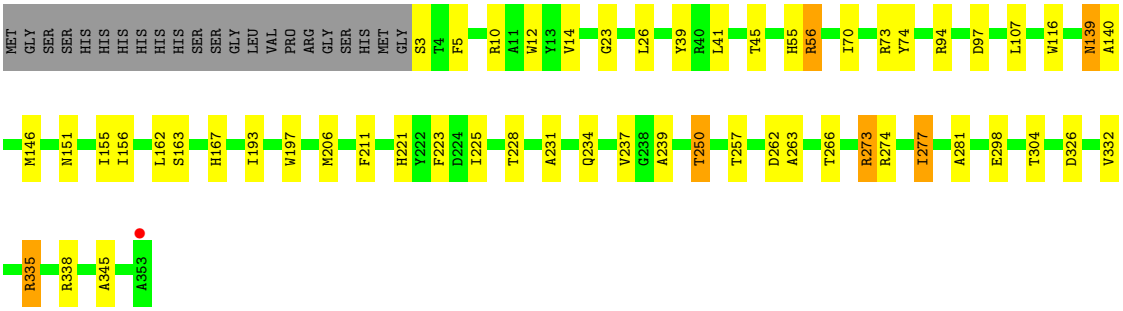
• Molecule 1: Rieske (2Fe-2S) domain protein

Chain F:



• Molecule 1: Rieske (2Fe-2S) domain protein

Chain G:

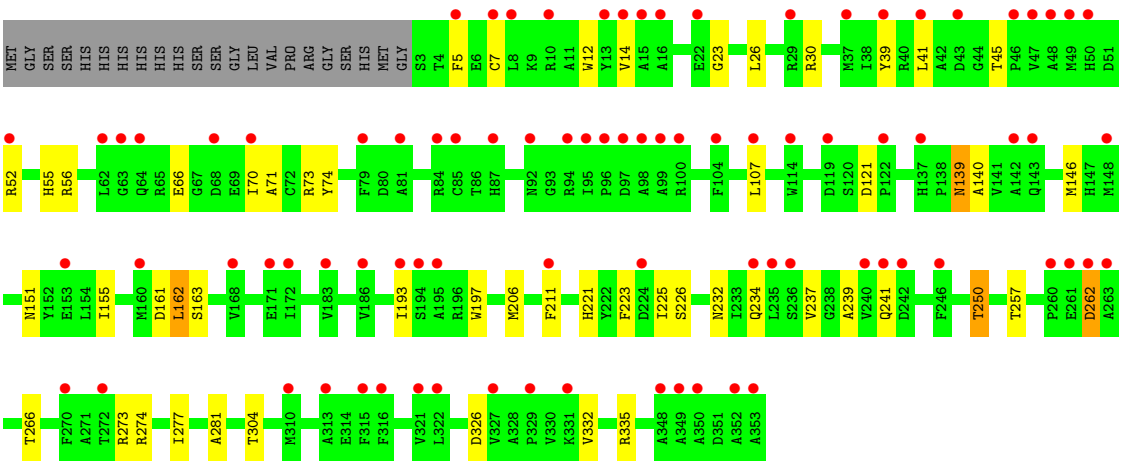


• Molecule 1: Rieske (2Fe-2S) domain protein

Chain H:

• Molecule 1: Rieske (2Fe-2S) domain protein

Chain I:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	315.58Å 315.58Å 189.96Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	157.79 – 3.71 157.79 – 3.71	Depositor EDS
% Data completeness (in resolution range)	88.0 (157.79-3.71) 88.0 (157.79-3.71)	Depositor EDS
R_{merge}	0.40	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.39 (at 3.68Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.174 , 0.216 0.177 , 0.210	Depositor DCC
R_{free} test set	3290 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	105.1	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.40$, $< L^2 > = 0.22$	Xtriage
Estimated twinning fraction	0.089 for -2/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+4/3*l,-1/3*h+1/3*k+1/3*l 0.079 for -h,1/3*h-1/3*k-4/3*l,-1/3*h-2/3*k+1/3*l 0.085 for -1/3*h+1/3*k+4/3*l,-k,2/3*h+1/3*k+1/3*l 0.089 for -h,2/3*h+1/3*k+4/3*l,1/3*h+2/3*k-1/3*l 0.096 for -1/3*h-2/3*k+4/3*l,-2/3*h-1/3*k-4/3*l,1/3*h-1/3*k-1/3*l 0.085 for 1/3*h+2/3*k-4/3*l,-k,-2/3*h-1/3*k-1/3*l 0.205 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.717 for H, K, L 0.283 for K, H, -L	Depositor
Outliers	0 of 65582 reflections	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	49487	wwPDB-VP
Average B, all atoms (Å^2)	143.0	wwPDB-VP

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% of the height of the origin peak. No significant pseudotranslation is detected.*

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FES, 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	0.76	0/2890	0.99	1/3930 (0.0%)
1	B	0.78	0/2886	1.01	1/3925 (0.0%)
1	C	0.76	0/2886	1.01	4/3925 (0.1%)
1	D	0.76	0/2886	0.99	2/3925 (0.1%)
1	E	0.76	0/2890	0.99	1/3930 (0.0%)
1	F	0.76	0/2886	1.02	6/3925 (0.2%)
1	G	0.77	0/2886	1.00	3/3925 (0.1%)
1	H	0.76	0/2886	1.00	5/3925 (0.1%)
1	I	0.75	0/2886	0.97	2/3925 (0.1%)
All	All	0.76	0/25982	1.00	25/35335 (0.1%)

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	C	0	1
1	D	0	1
1	F	0	2
1	G	0	1
1	H	0	1
All	All	0	6

There are no bond length outliers.

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	161	ASP	CB-CA-C	7.24	124.88	110.40
1	F	141	VAL	CA-CB-CG2	7.23	121.74	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	I	161	ASP	CB-CA-C	6.87	124.15	110.40
1	B	161	ASP	CA-CB-CG	6.75	128.25	113.40
1	H	245	ASP	CB-CG-OD1	6.43	124.09	118.30

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	3	SER	Peptide
1	D	3	SER	Peptide
1	F	234	GLN	Sidechain
1	F	3	SER	Peptide
1	G	3	SER	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	2812	2687	2666	27	0
1	B	2808	2684	2663	40	0
1	C	2808	2684	2665	28	0
1	D	2808	2684	2663	27	0
1	E	2812	2687	2667	28	0
1	F	2808	2684	2663	45	0
1	G	2808	2684	2663	23	0
1	H	2808	2684	2663	21	0
1	I	2808	2684	2665	20	0
2	A	4	0	0	0	0
2	B	4	0	0	1	0
2	C	4	0	0	2	0
2	D	4	0	0	0	0
2	E	4	0	0	0	0
2	F	4	0	0	5	0
2	G	4	0	0	0	0
2	H	4	0	0	1	0
2	I	4	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
3	I	1	0	0	0	0
All	All	25325	24162	23978	245	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 5.

The worst 5 of 245 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:F:74:TYR:HB3	2:F:401:FES:S2	1.97	1.03
1:F:55:HIS:HB3	2:F:401:FES:S1	2.07	0.94
1:F:160:MET:CE	1:F:227:TRP:HB2	2.04	0.88
1:E:70:ILE:N	1:E:70:ILE:HD12	2.04	0.72
1:B:316:PHE:CG	1:F:135:GLU:HG2	2.26	0.71

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	350/373 (94%)	311 (89%)	34 (10%)	5 (1%)	11 45
1	B	349/373 (94%)	309 (88%)	35 (10%)	5 (1%)	11 45
1	C	349/373 (94%)	308 (88%)	35 (10%)	6 (2%)	9 42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	349/373 (94%)	309 (88%)	35 (10%)	5 (1%)	11	45
1	E	350/373 (94%)	308 (88%)	37 (11%)	5 (1%)	11	45
1	F	349/373 (94%)	311 (89%)	35 (10%)	3 (1%)	17	53
1	G	349/373 (94%)	311 (89%)	33 (10%)	5 (1%)	11	45
1	H	349/373 (94%)	313 (90%)	31 (9%)	5 (1%)	11	45
1	I	349/373 (94%)	311 (89%)	33 (10%)	5 (1%)	11	45
All	All	3143/3357 (94%)	2791 (89%)	308 (10%)	44 (1%)	11	45

5 of 44 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	5	PHE
1	A	23	GLY
1	B	23	GLY
1	C	23	GLY
1	D	5	PHE

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	294/312 (94%)	279 (95%)	15 (5%)	24	55
1	B	294/312 (94%)	276 (94%)	18 (6%)	18	50
1	C	294/312 (94%)	277 (94%)	17 (6%)	20	51
1	D	294/312 (94%)	275 (94%)	19 (6%)	17	48
1	E	294/312 (94%)	278 (95%)	16 (5%)	22	54
1	F	294/312 (94%)	275 (94%)	19 (6%)	17	48
1	G	294/312 (94%)	273 (93%)	21 (7%)	14	45
1	H	294/312 (94%)	272 (92%)	22 (8%)	13	44
1	I	294/312 (94%)	275 (94%)	19 (6%)	17	48
All	All	2646/2808 (94%)	2480 (94%)	166 (6%)	18	49

5 of 166 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	206	MET
1	H	304	THR
1	G	262	ASP
1	H	139	ASN
1	I	56	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	178	GLN
1	D	344	GLN
1	G	325	ASN
1	E	158	ASN
1	B	232	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 18 ligands modelled in this entry, 9 are monoatomic - leaving 9 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	$\#\|Z\|>2$
2	FES	C	401	1	0,4,4	-	-	-		
2	FES	E	401	1	0,4,4	-	-	-		
2	FES	D	401	1	0,4,4	-	-	-		
2	FES	A	401	1	0,4,4	-	-	-		
2	FES	F	401	1	0,4,4	-	-	-		
2	FES	B	401	1	0,4,4	-	-	-		
2	FES	G	401	1	0,4,4	-	-	-		
2	FES	I	401	1	0,4,4	-	-	-		
2	FES	H	401	1	0,4,4	-	-	-		

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
2	FES	C	401	1	-	-	0/1/1/1
2	FES	E	401	1	-	-	0/1/1/1
2	FES	D	401	1	-	-	0/1/1/1
2	FES	A	401	1	-	-	0/1/1/1
2	FES	F	401	1	-	-	0/1/1/1
2	FES	B	401	1	-	-	0/1/1/1
2	FES	G	401	1	-	-	0/1/1/1
2	FES	I	401	1	-	-	0/1/1/1
2	FES	H	401	1	-	-	0/1/1/1

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.

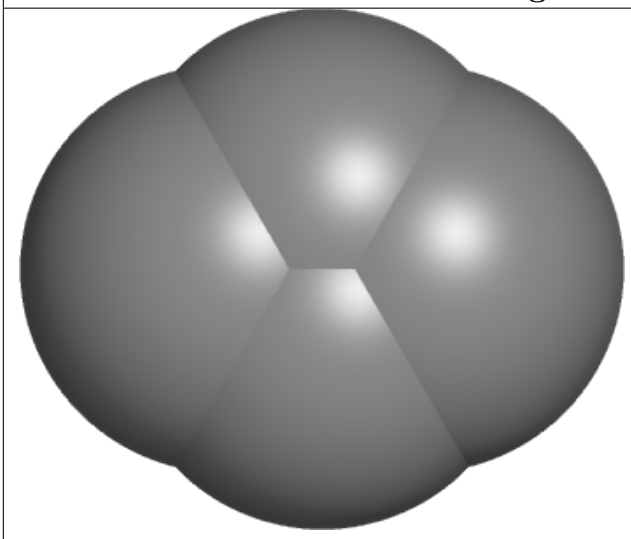
4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	FES	2	0
2	F	401	FES	5	0
2	B	401	FES	1	0
2	H	401	FES	1	0

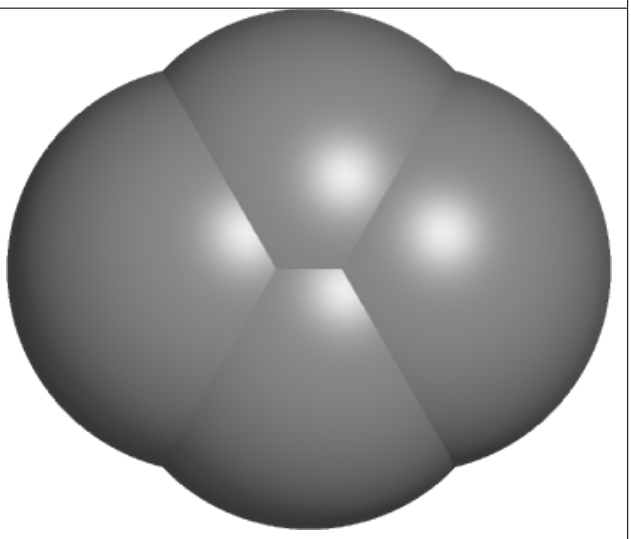
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

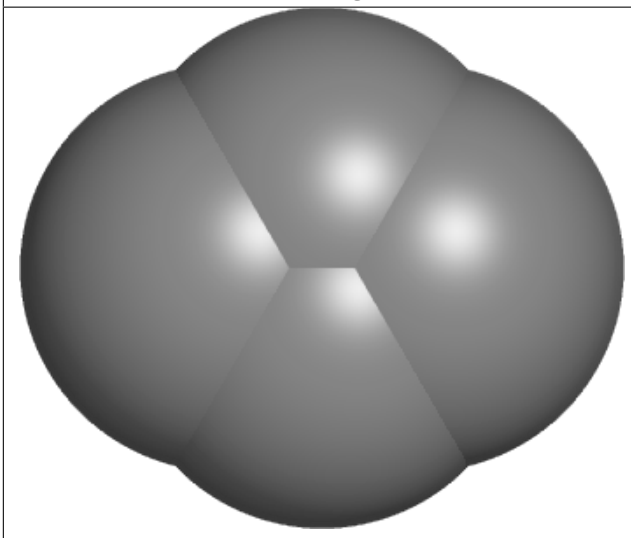
Ligand FES C 401



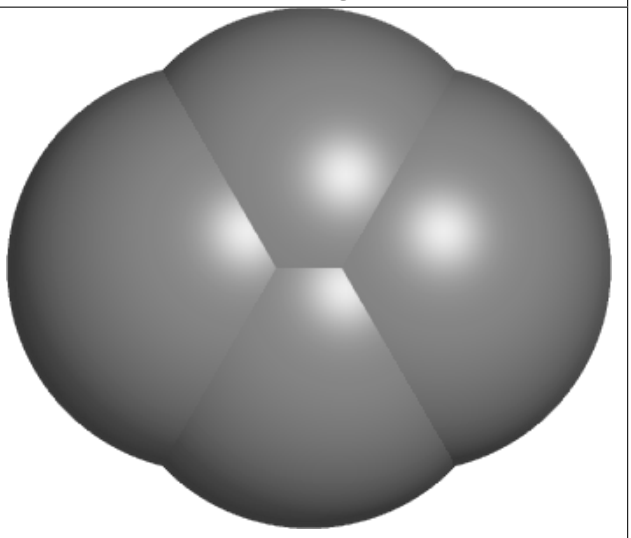
Bond lengths



Bond angles

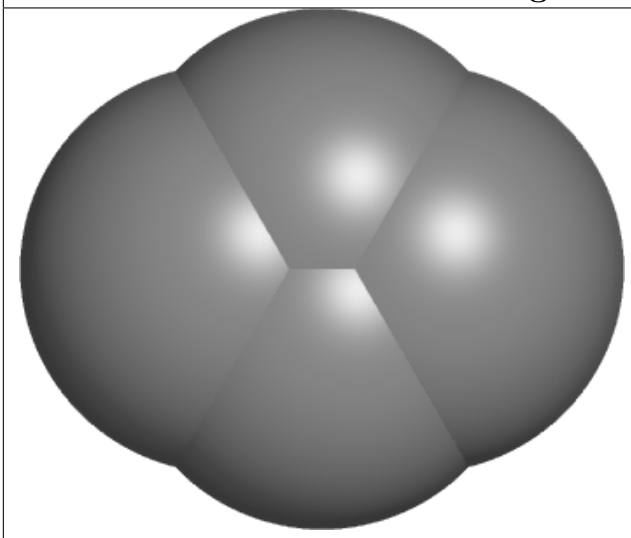


Torsions



Rings

Ligand FES E 401



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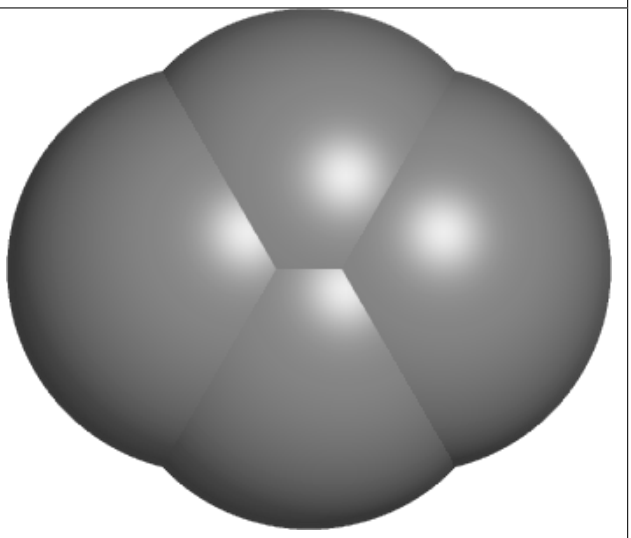

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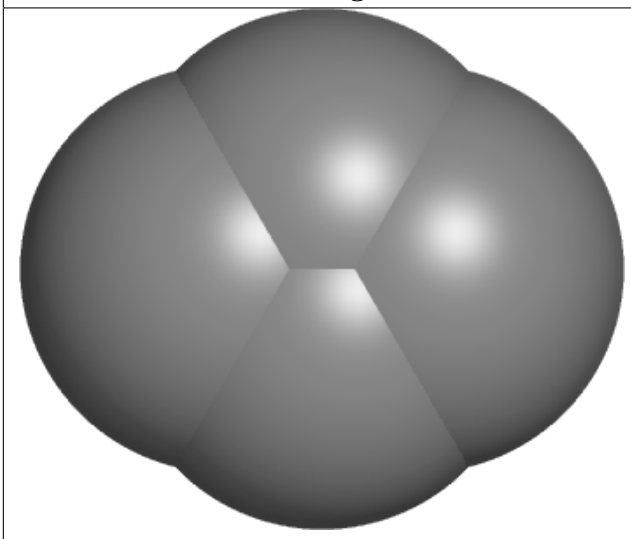

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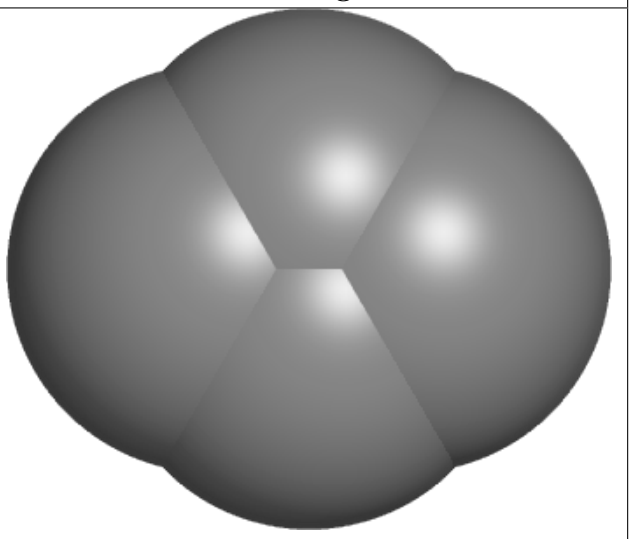

Rings

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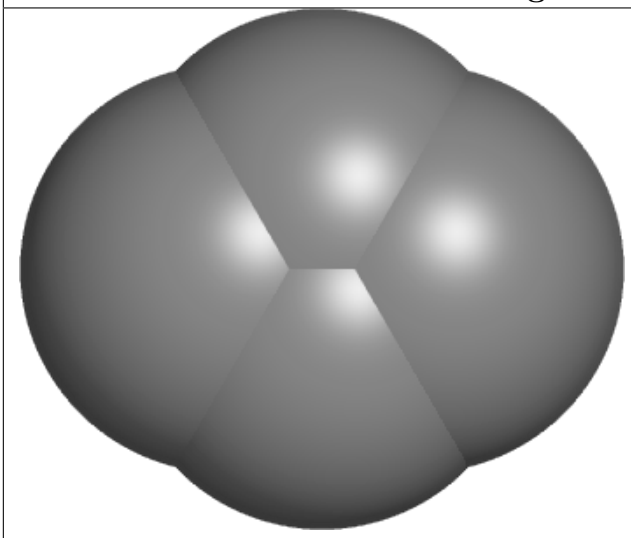

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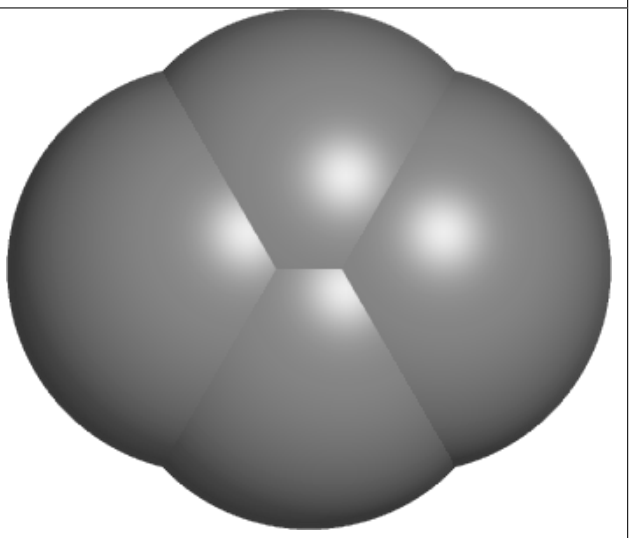

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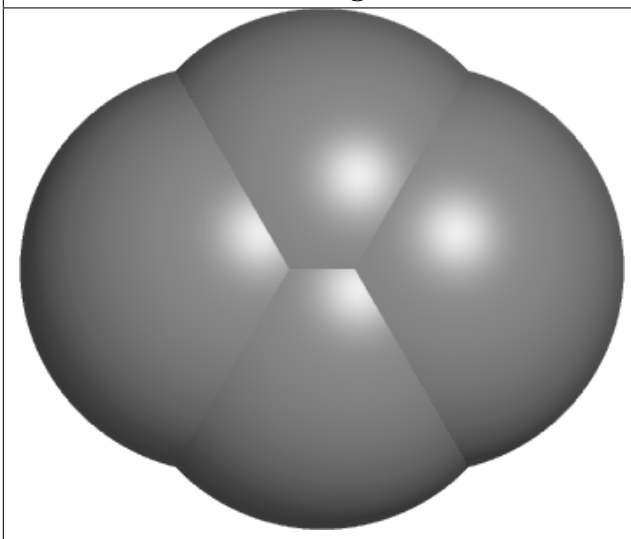

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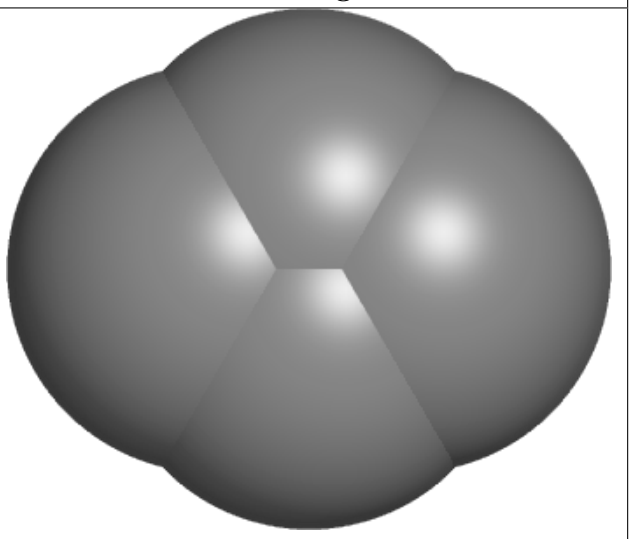
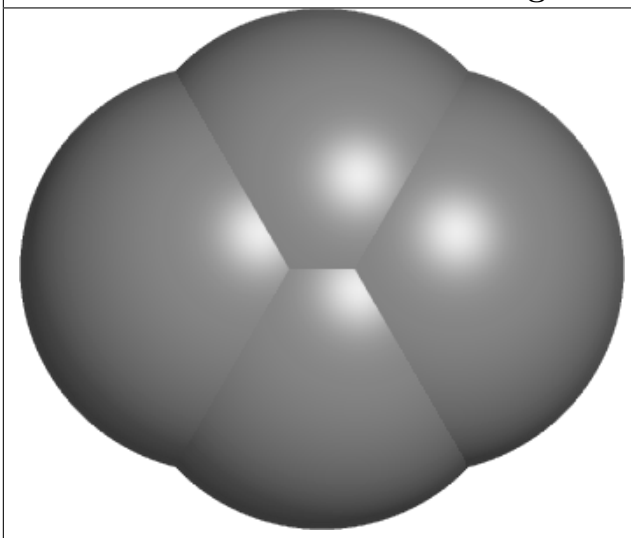


Rings

Ligand FES A 401



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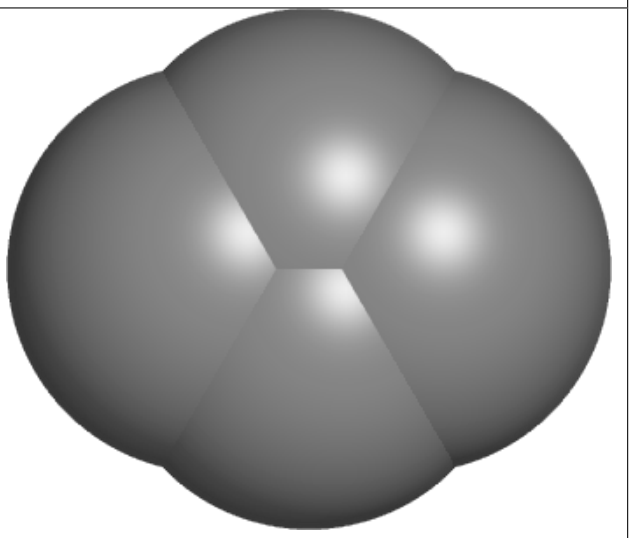
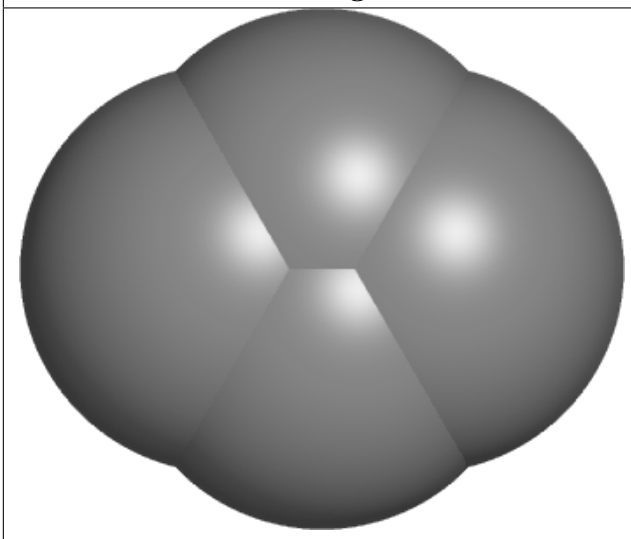


Bond angles



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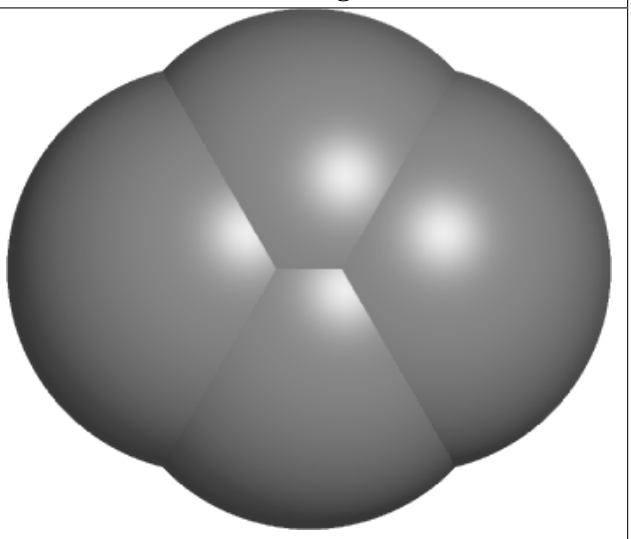
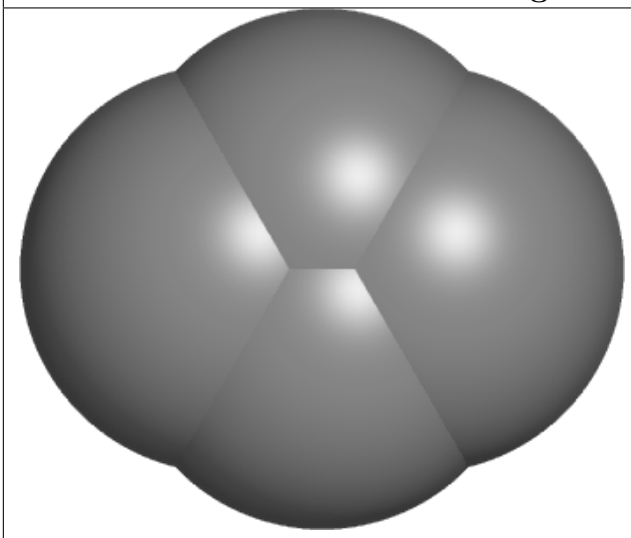
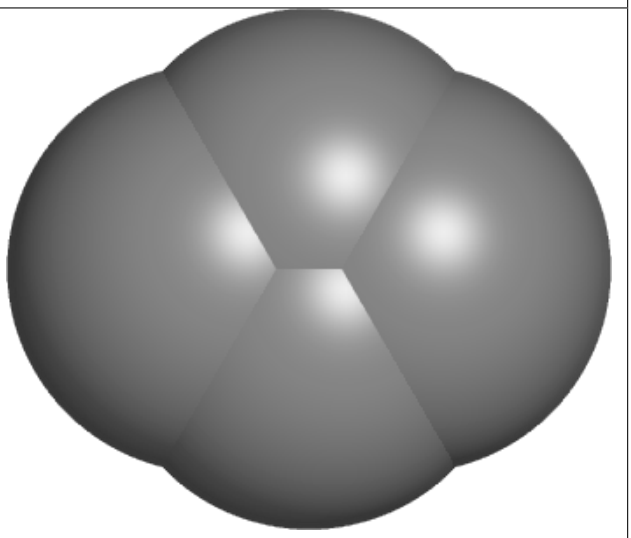
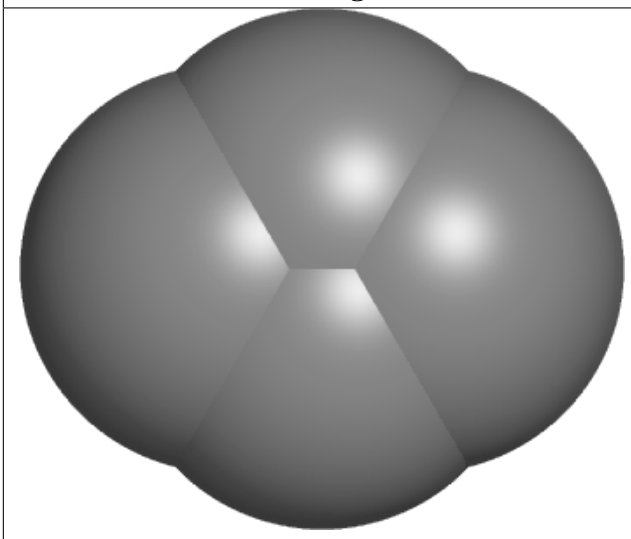
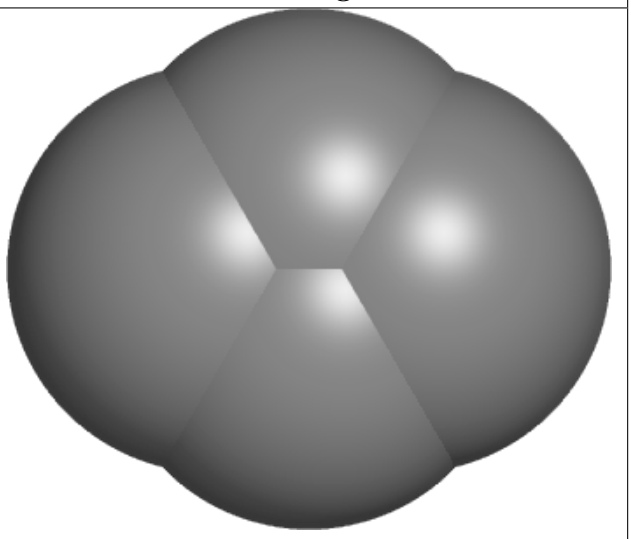
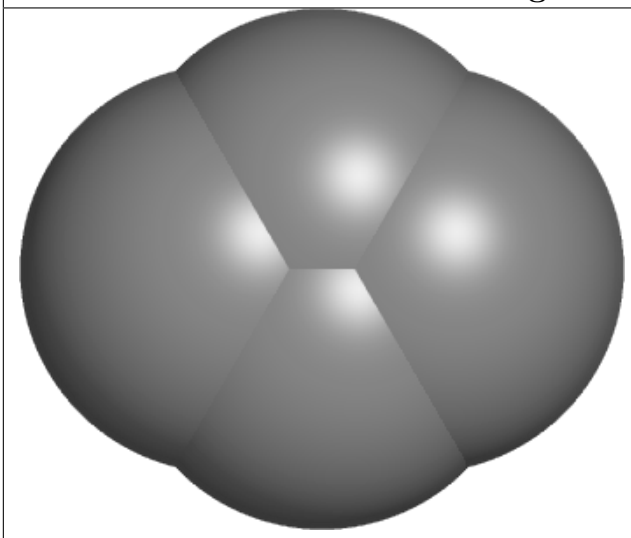


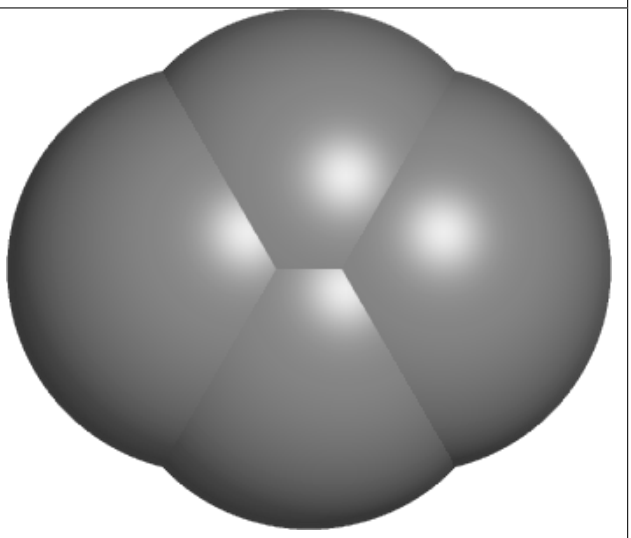
Rings

Ligand FES F 401	
	
Bond lengths	**Bond angles**
	
Torsions	**Rings**

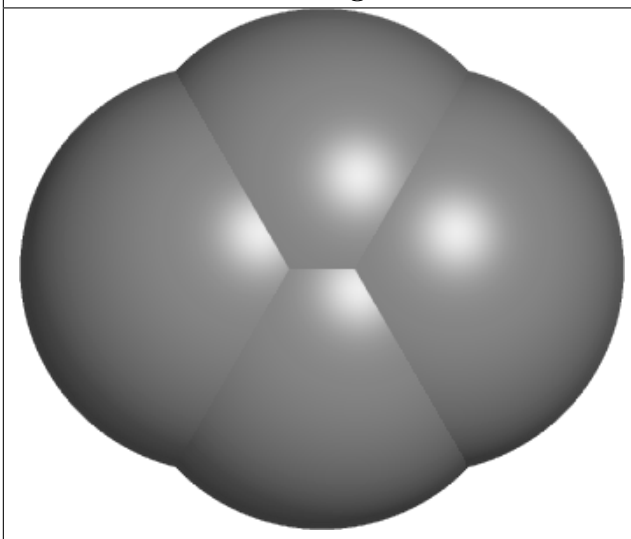
Ligand FES B 401



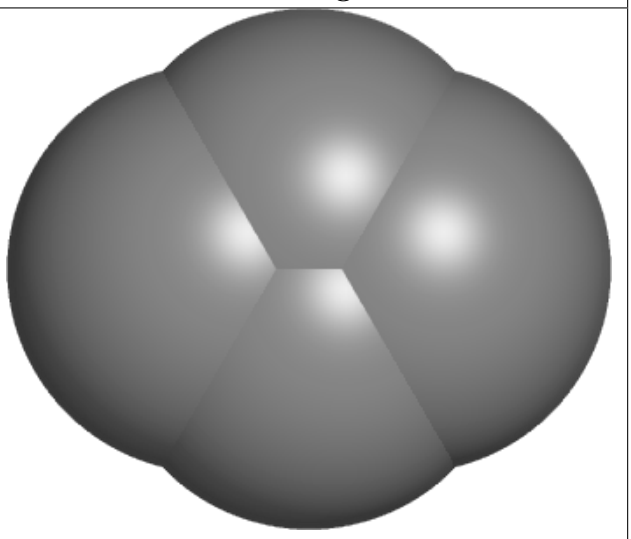
Bond lengths



Bond angles

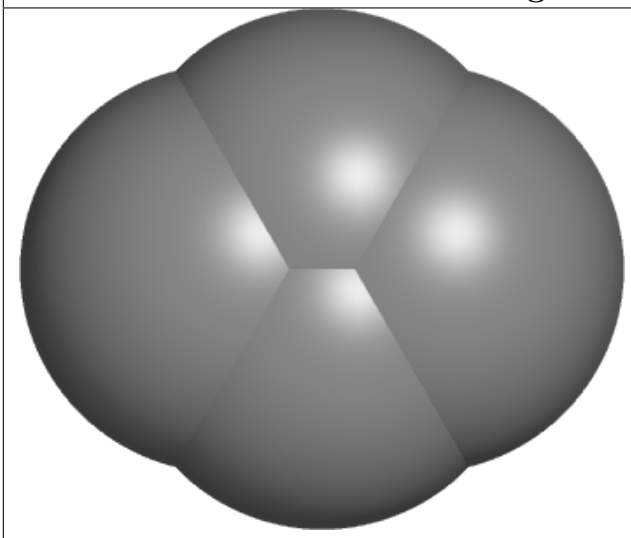


Torsions

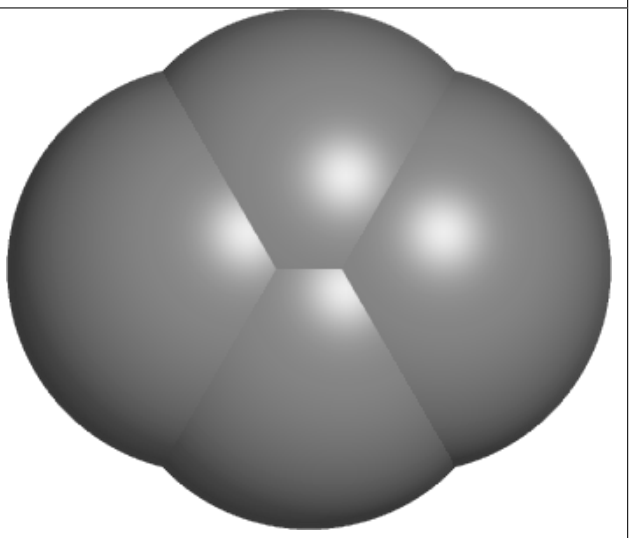


Rings

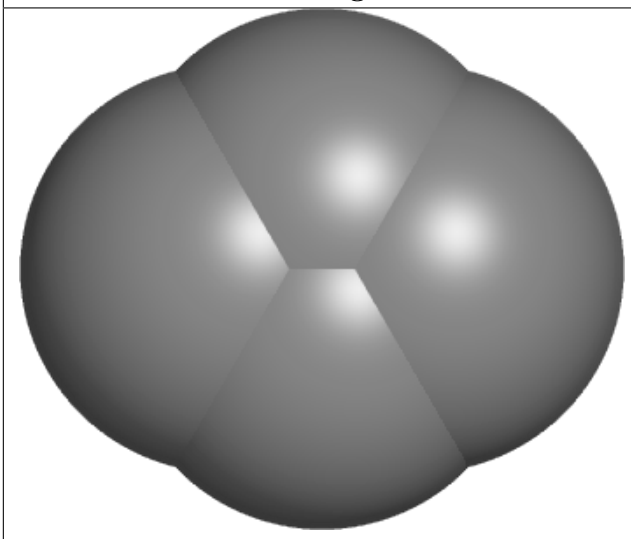
Ligand FES G 401



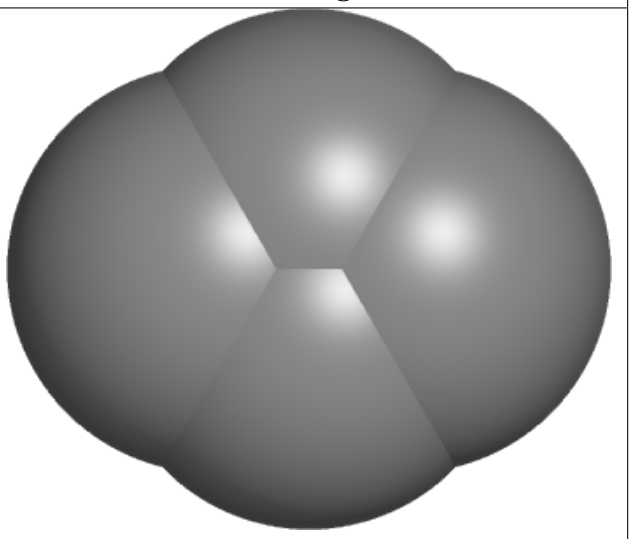
Bond lengths



Bond angles

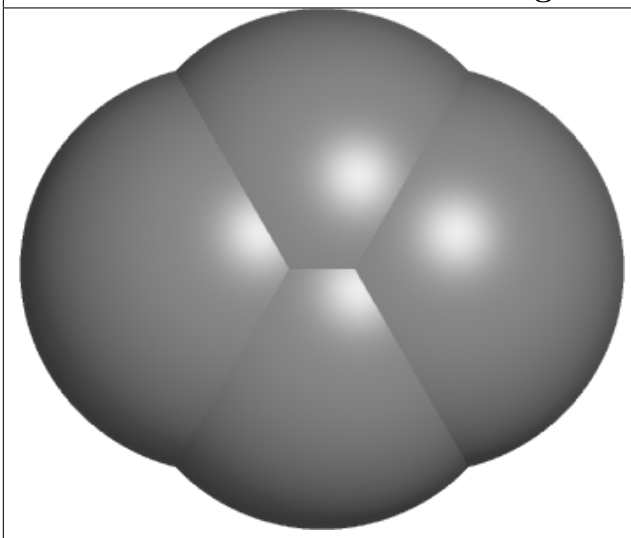


Torsions

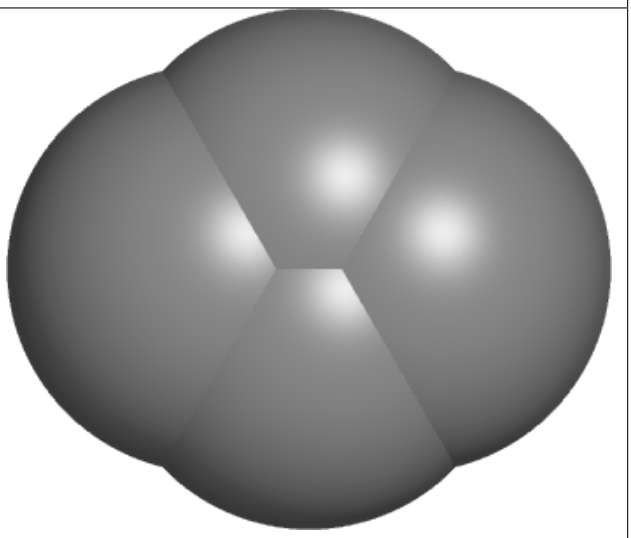


Rings

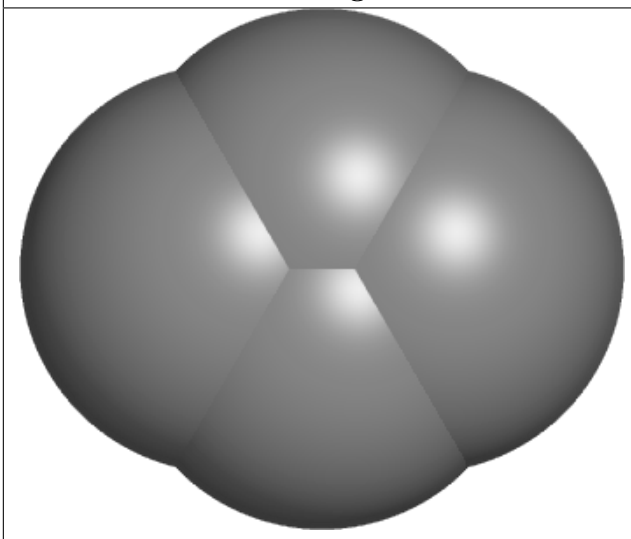
Ligand FES I 401



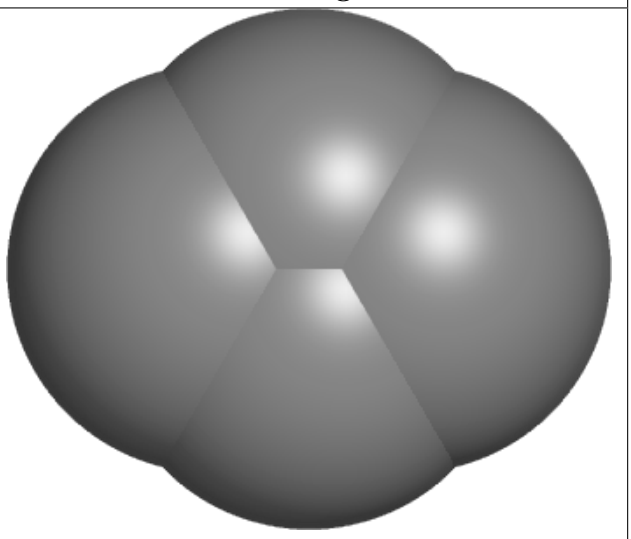
Bond lengths



Bond angles

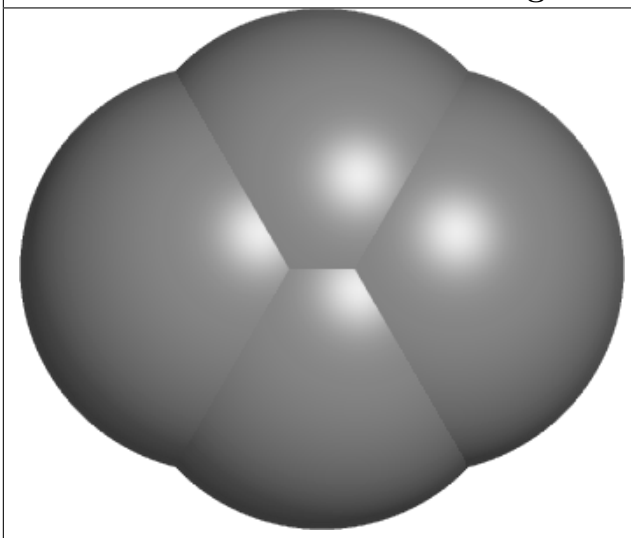


Torsions

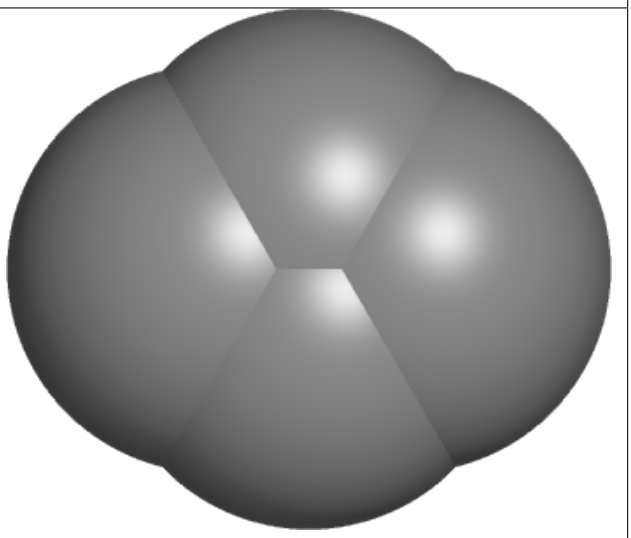


Rings

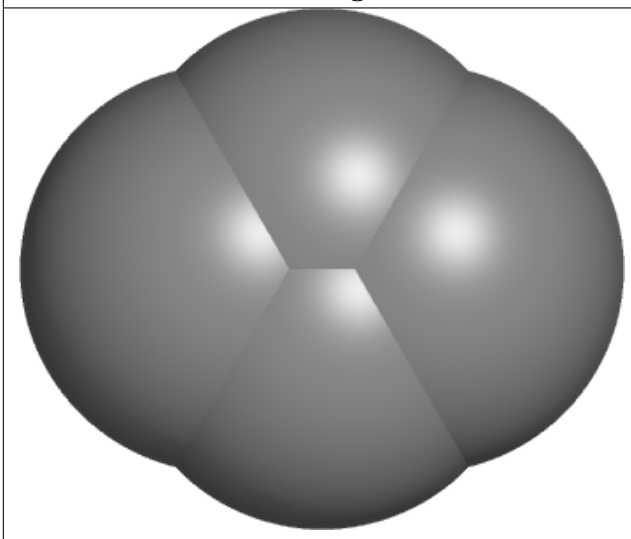
Ligand FES H 401



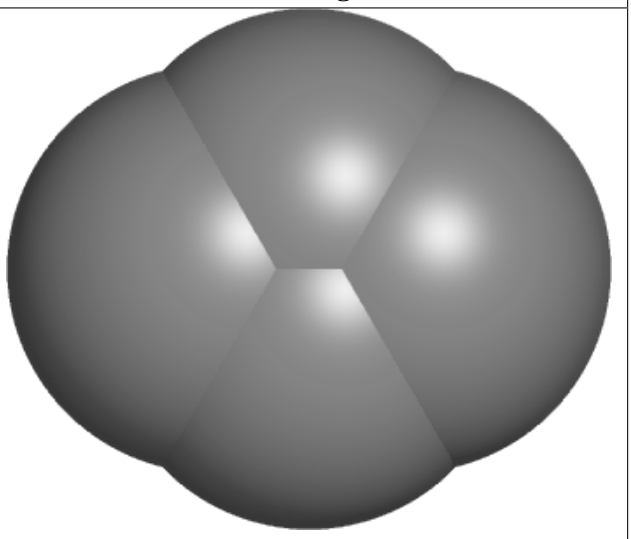
Bond lengths



Bond angles



Torsions



Rings

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, 95^{th} 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	352/373 (94%)	-0.13	1 (0%) 94 93	73, 112, 148, 191	0
1	B	351/373 (94%)	-0.19	3 (0%) 84 80	75, 106, 148, 232	0
1	C	351/373 (94%)	-0.22	1 (0%) 94 93	68, 103, 143, 179	0
1	D	351/373 (94%)	0.49	28 (7%) 12 10	117, 170, 232, 264	0
1	E	352/373 (94%)	0.48	25 (7%) 16 12	118, 180, 229, 281	0
1	F	351/373 (94%)	-0.05	2 (0%) 89 87	82, 133, 175, 200	0
1	G	351/373 (94%)	-0.23	1 (0%) 94 93	74, 105, 145, 216	0
1	H	351/373 (94%)	0.00	1 (0%) 94 93	85, 144, 174, 199	0
1	I	351/373 (94%)	1.34	86 (24%) 0 0	155, 206, 241, 271	0
All	All	3161/3357 (94%)	0.17	148 (4%) 31 26	68, 135, 219, 281	0

The worst 5 of 148 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	64	GLN	5.1
1	D	239	ALA	5.0
1	D	198	GLU	5.0
1	E	239	ALA	4.9
1	I	48	ALA	4.8

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 monosaccharides in this entry.

6.4 Ligands (i)

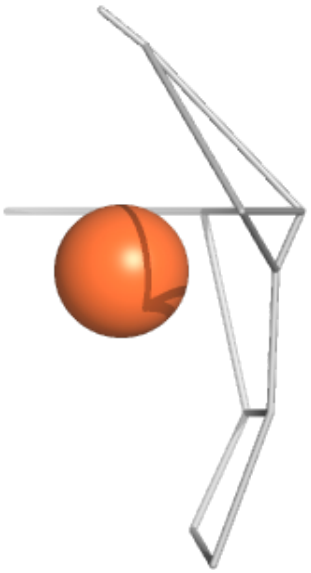
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

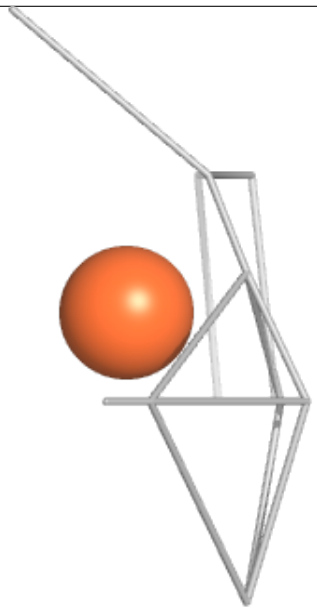
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	FE	I	402	1/1	0.89	0.13	238,238,238,238	0
3	FE	D	402	1/1	0.92	0.20	191,191,191,191	0
3	FE	E	402	1/1	0.96	0.11	202,202,202,202	0
3	FE	G	402	1/1	0.98	0.13	96,96,96,96	0
3	FE	F	402	1/1	0.98	0.12	135,135,135,135	0
3	FE	A	402	1/1	0.99	0.14	90,90,90,90	0
3	FE	B	402	1/1	0.99	0.14	90,90,90,90	0
3	FE	C	402	1/1	0.99	0.12	113,113,113,113	0
3	FE	H	402	1/1	0.99	0.11	165,165,165,165	0
2	FES	I	401	4/4	0.99	0.15	225,251,258,283	0
2	FES	C	401	4/4	1.00	0.22	66,67,69,70	0
2	FES	D	401	4/4	1.00	0.21	107,116,128,137	0
2	FES	E	401	4/4	1.00	0.19	133,163,166,186	0
2	FES	F	401	4/4	1.00	0.19	134,146,167,187	0
2	FES	G	401	4/4	1.00	0.25	66,70,76,83	0
2	FES	H	401	4/4	1.00	0.21	100,112,113,115	0
2	FES	A	401	4/4	1.00	0.24	62,69,78,81	0
2	FES	B	401	4/4	1.00	0.22	69,75,84,86	0

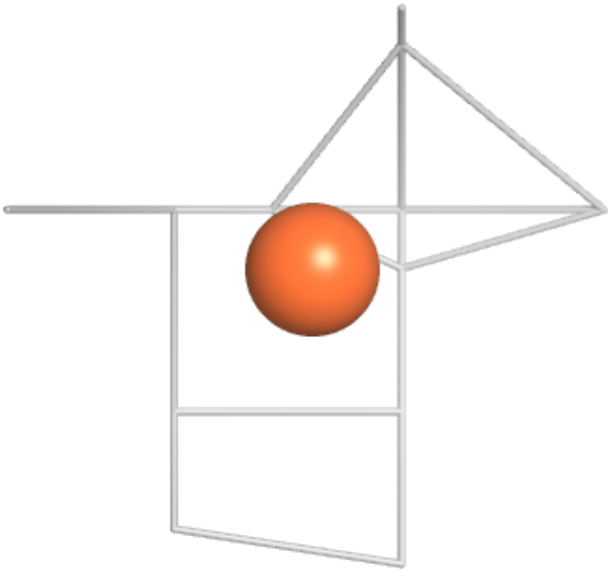
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around FE I 402:

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

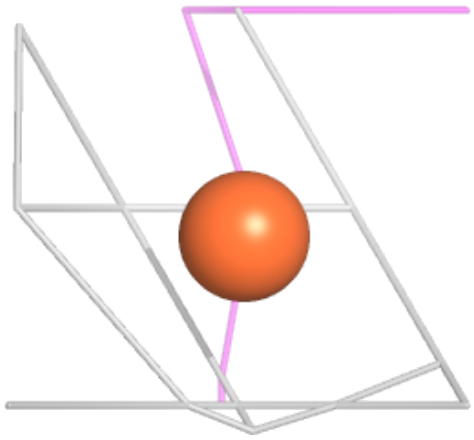


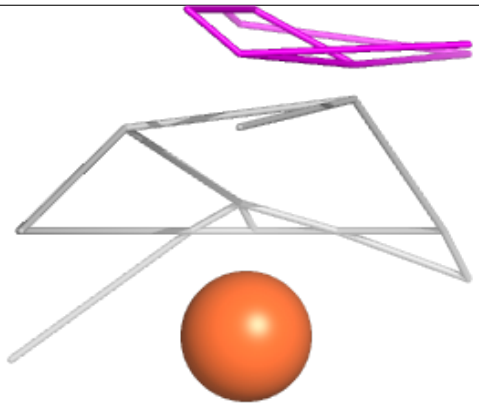


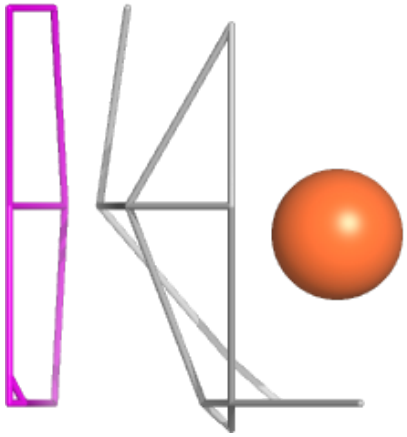


Electron density around FE D 402:

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

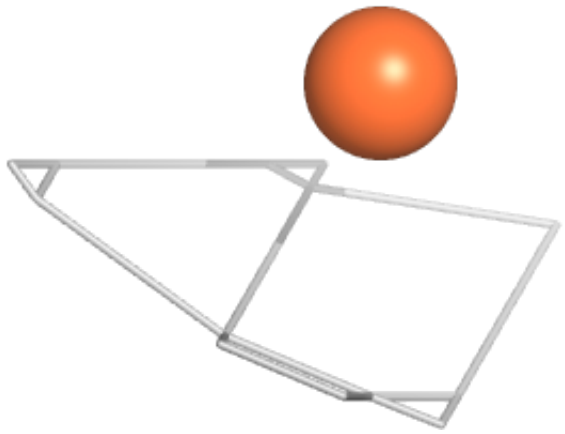


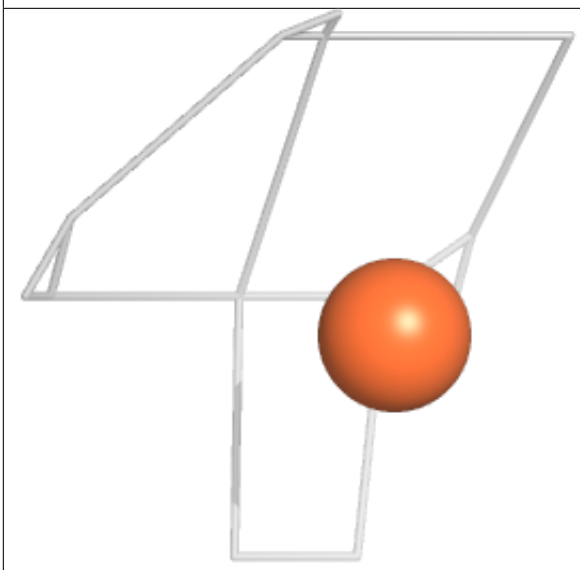


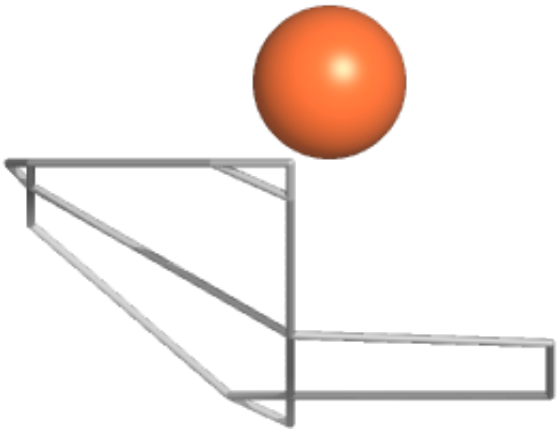


Electron density around FE E 402:

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

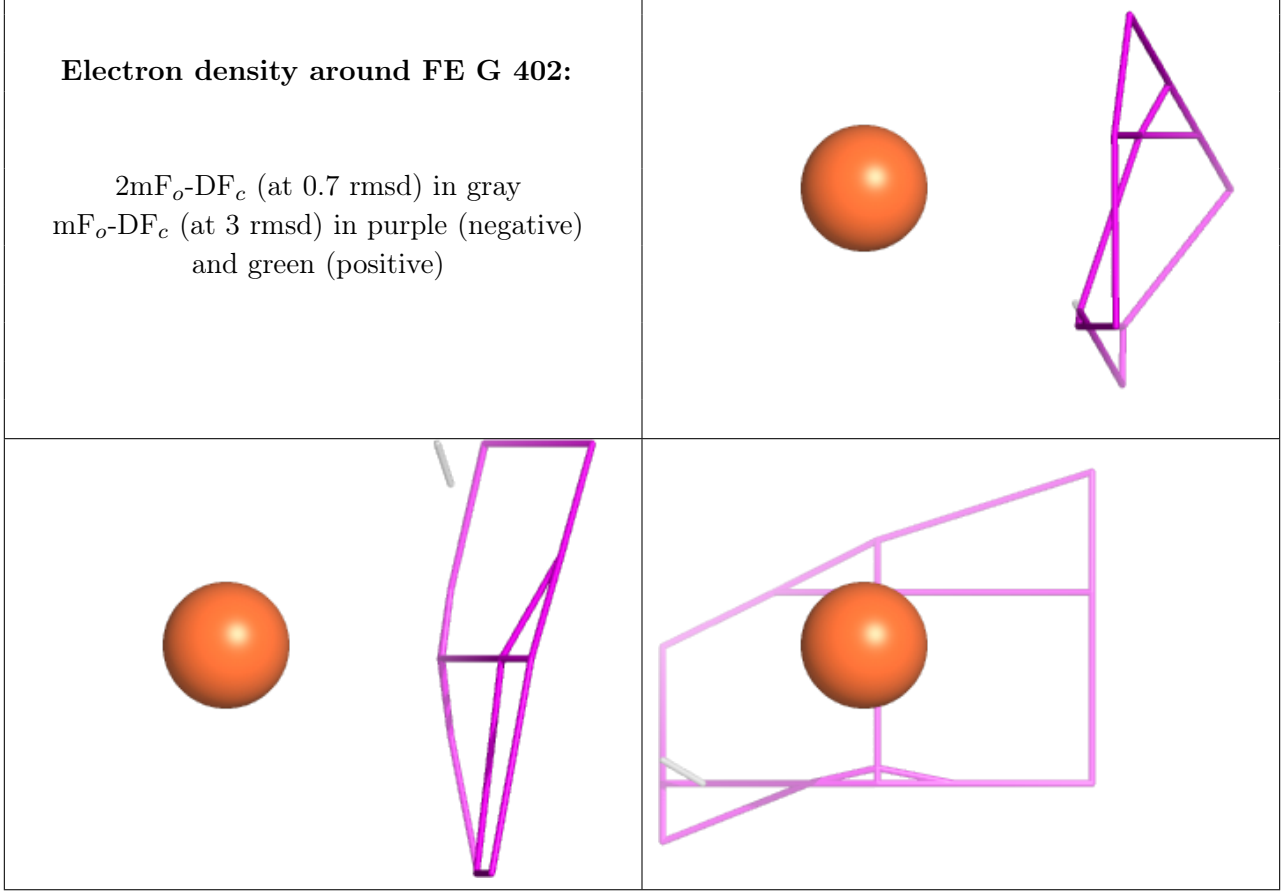






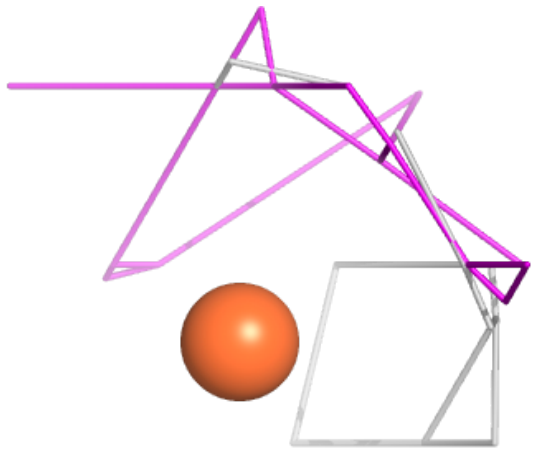
Electron density around FE G 402:

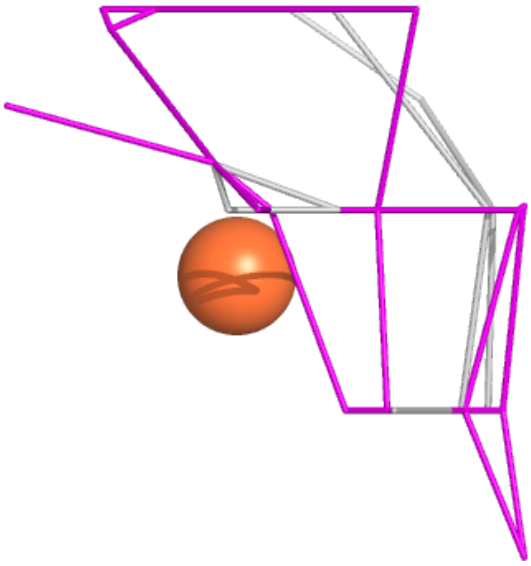
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

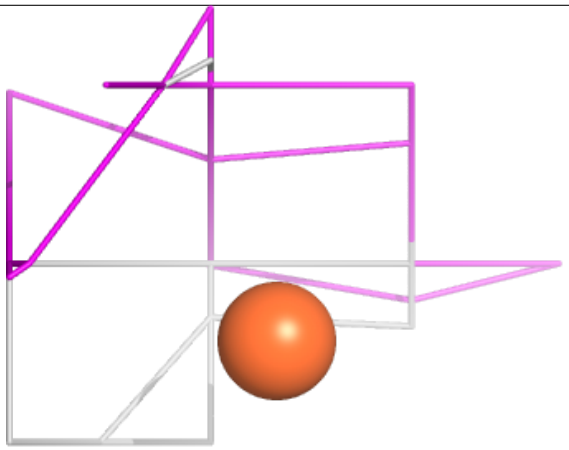


Electron density around FE F 402:

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

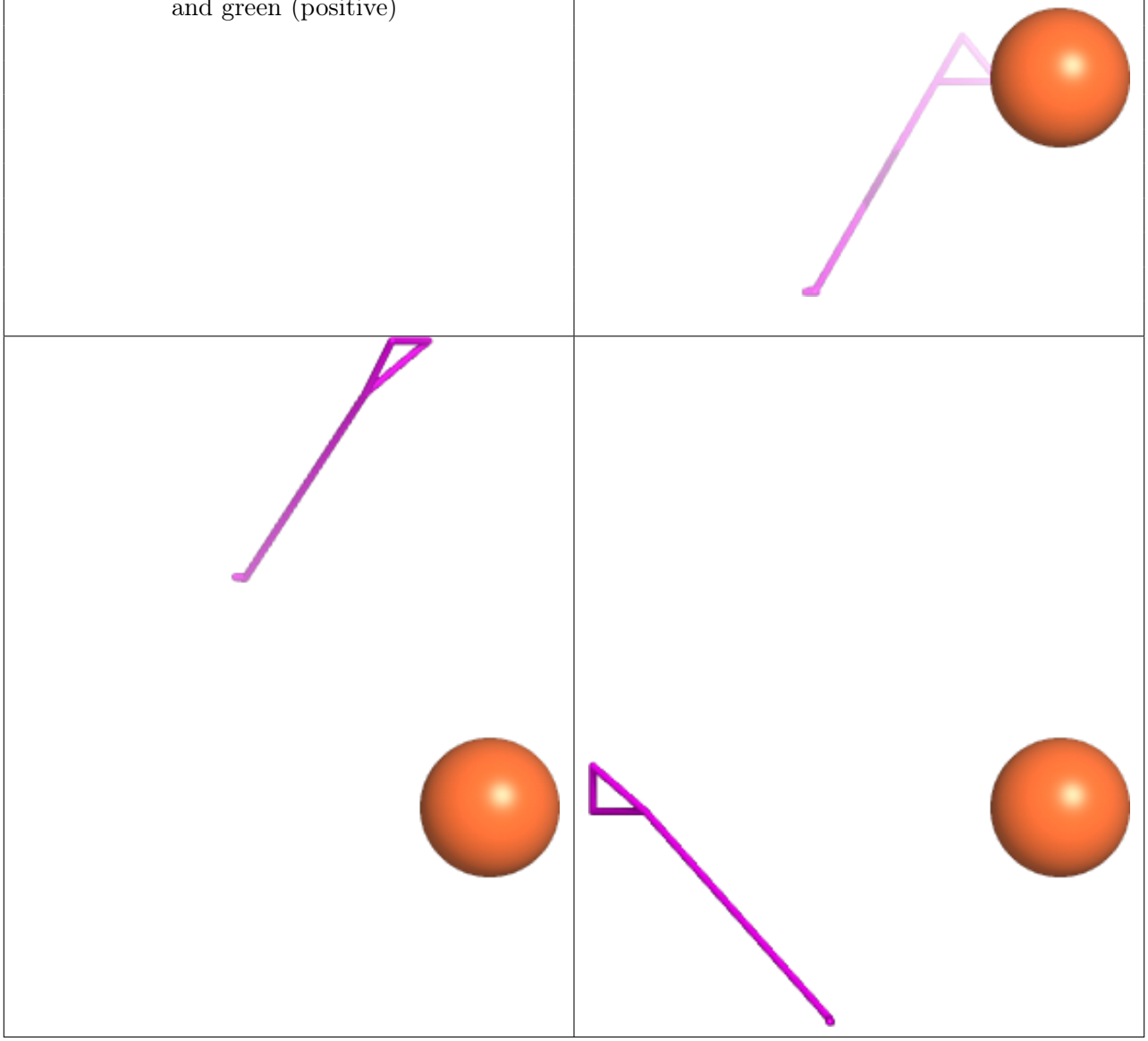






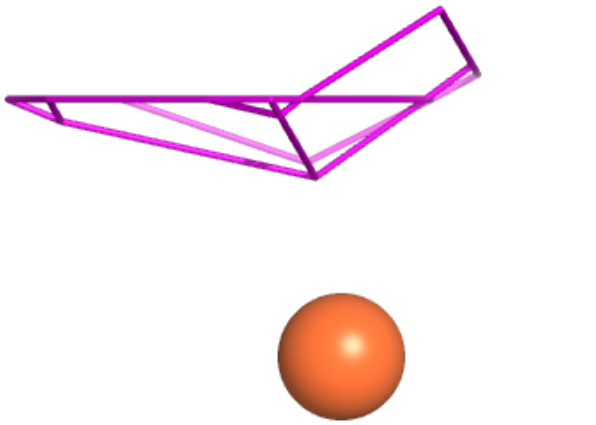
Electron density around FE A 402:

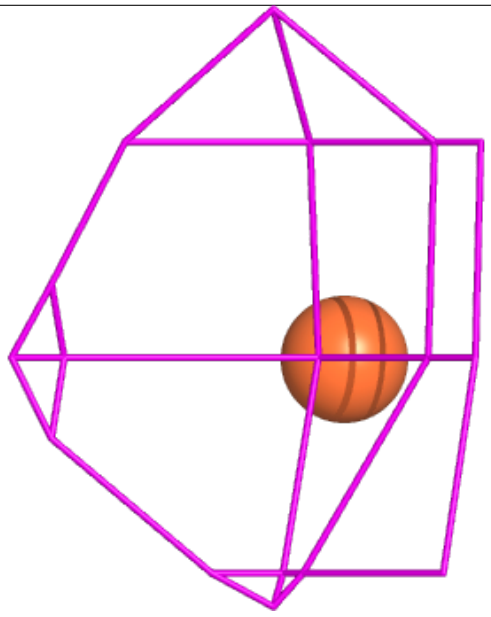
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

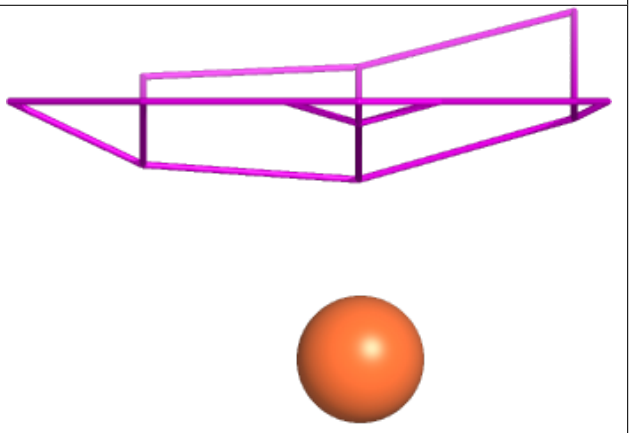


Electron density around FE B 402:

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

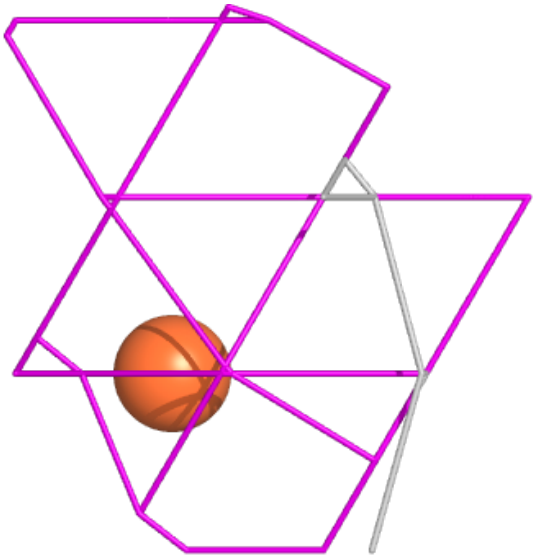


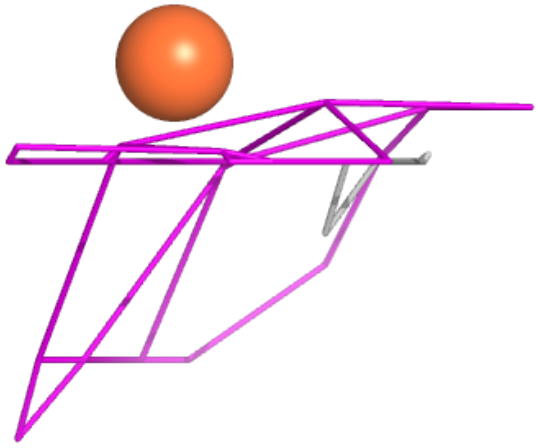


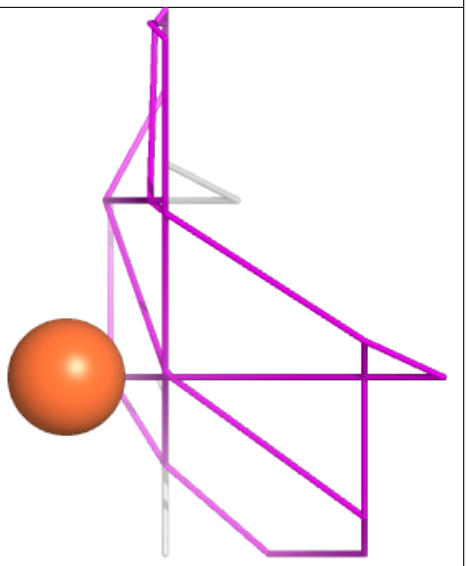


Electron density around FE C 402:

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

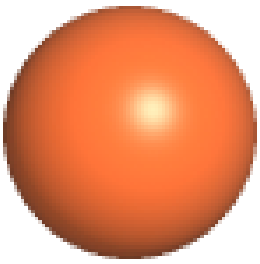


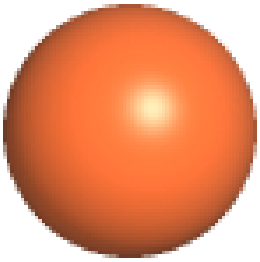


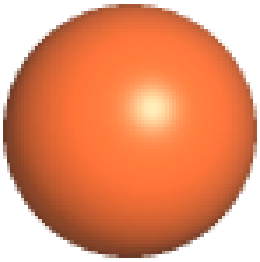


Electron density around FE H 402:

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

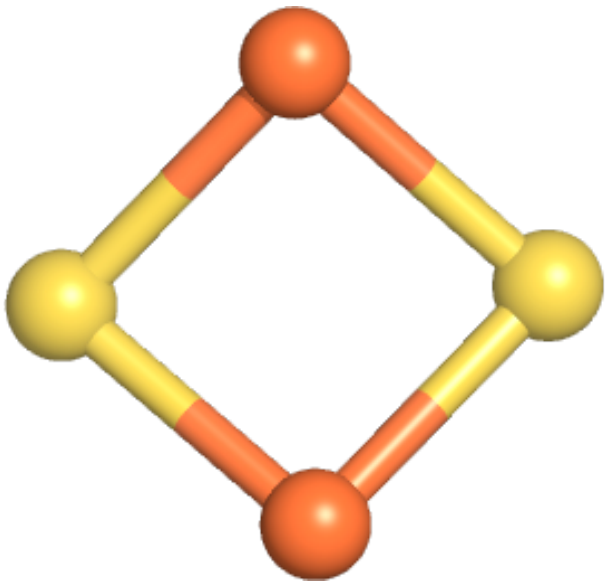


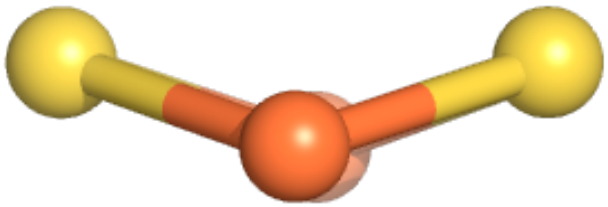




Electron density around FES I 401:

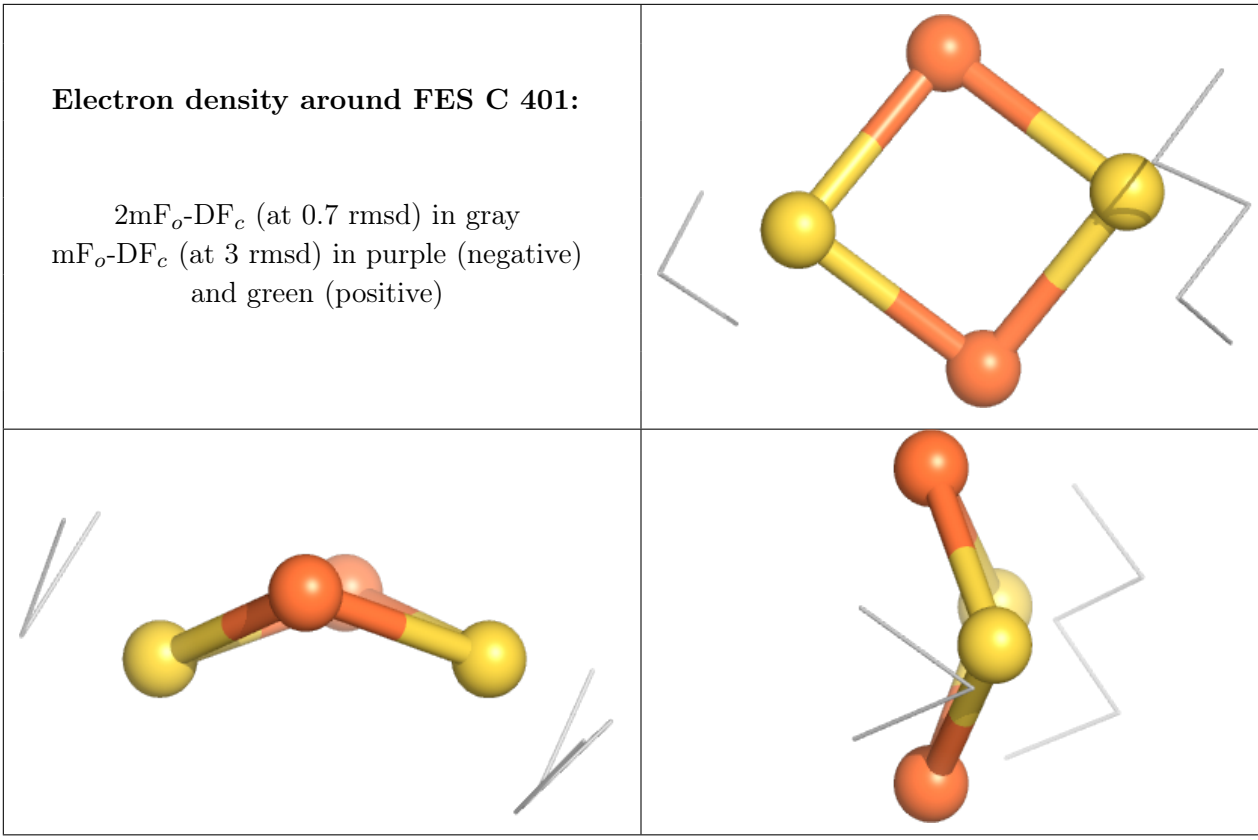
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)





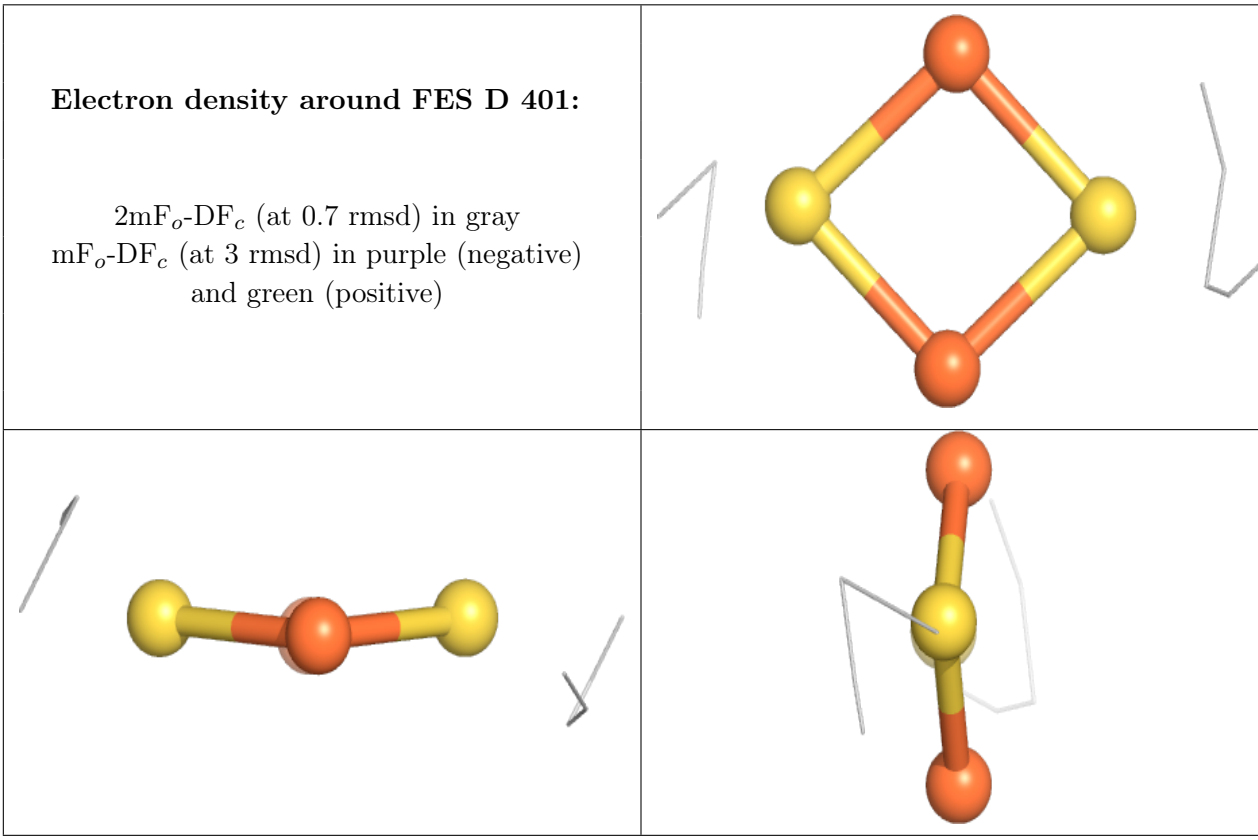
Electron density around FES C 401:

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



Electron density around FES 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 FES E 401:

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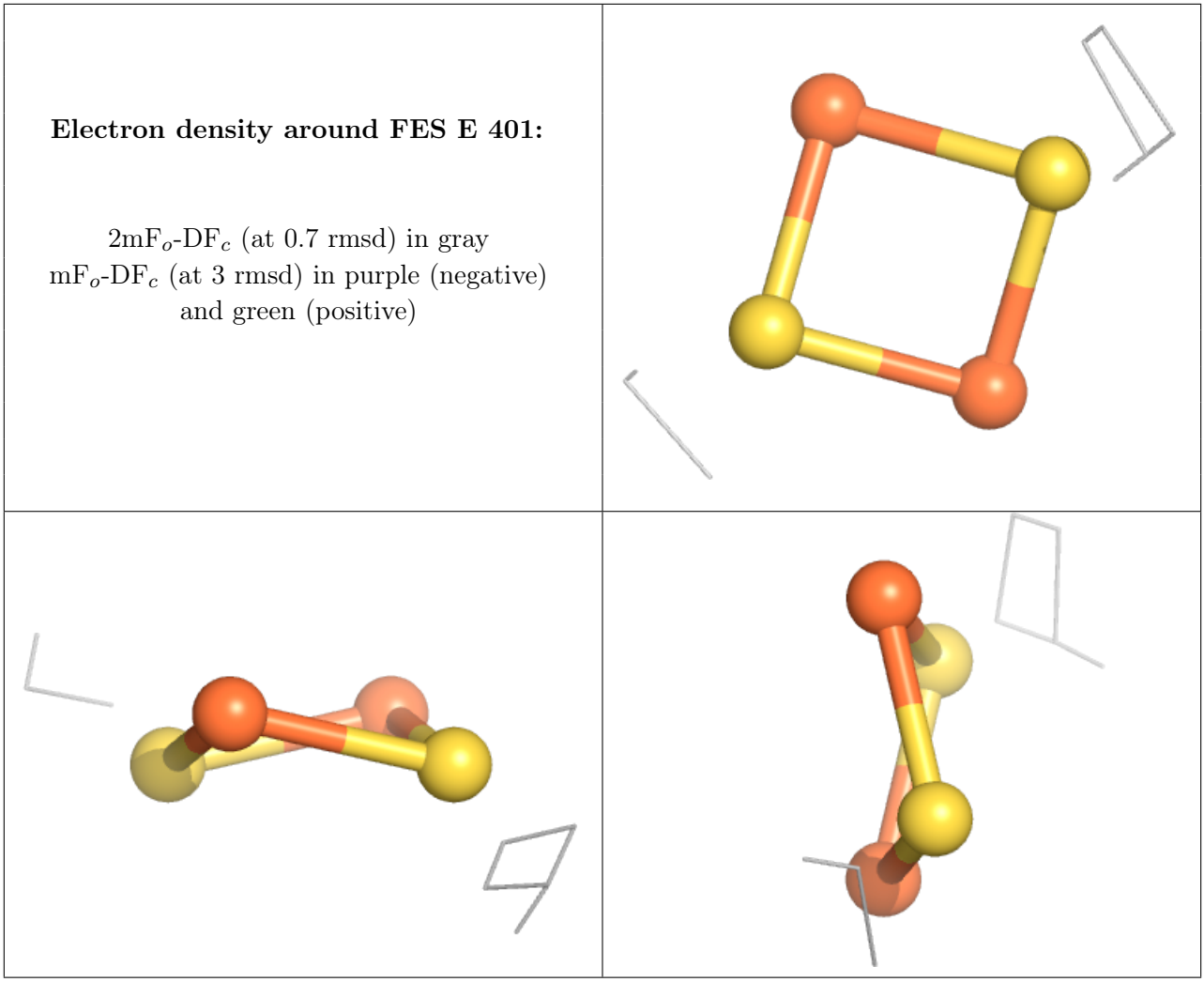
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around FES F 401:

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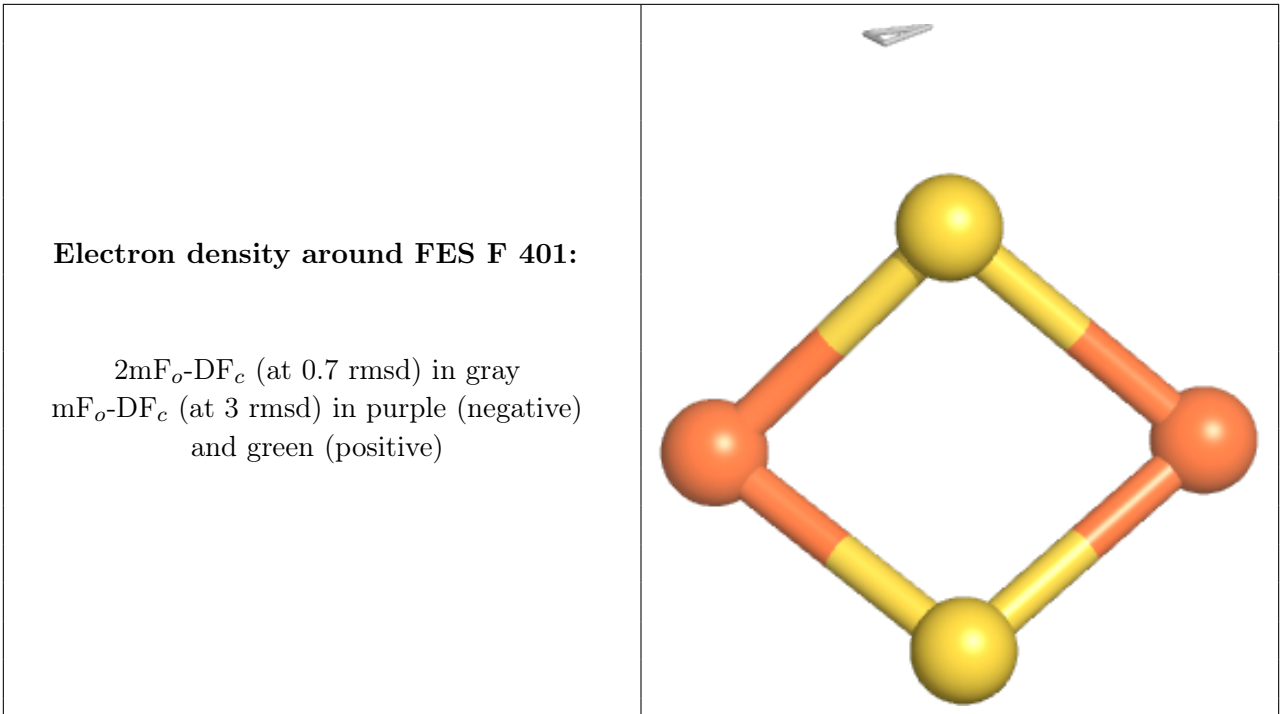
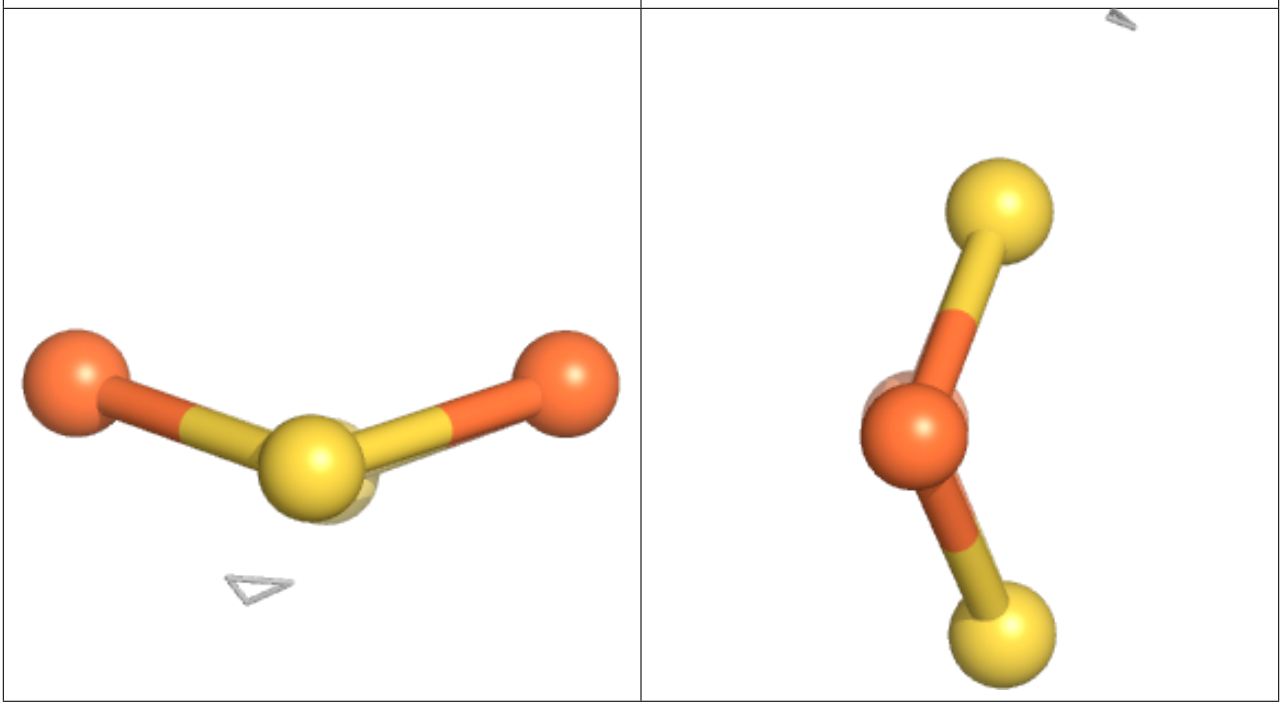
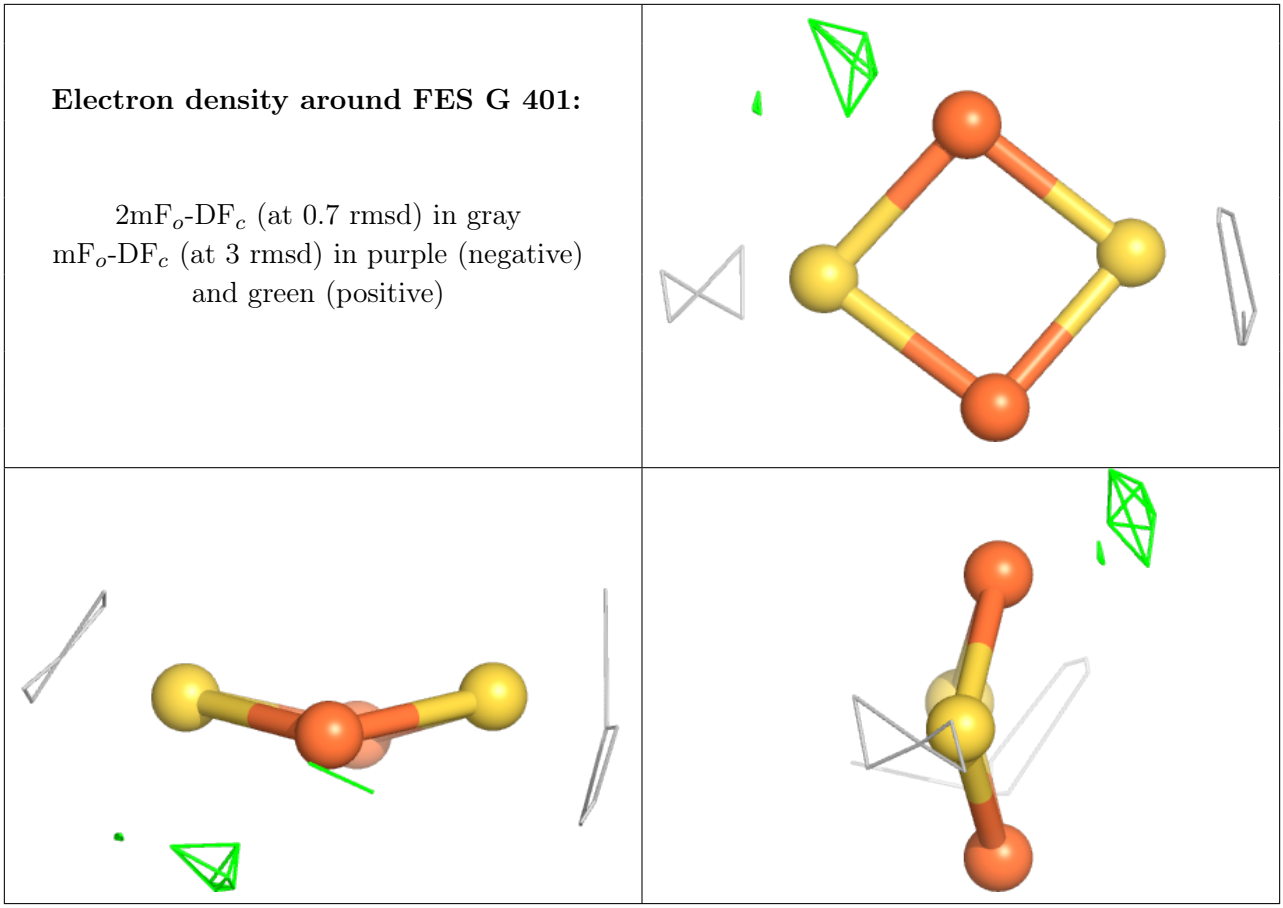
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)





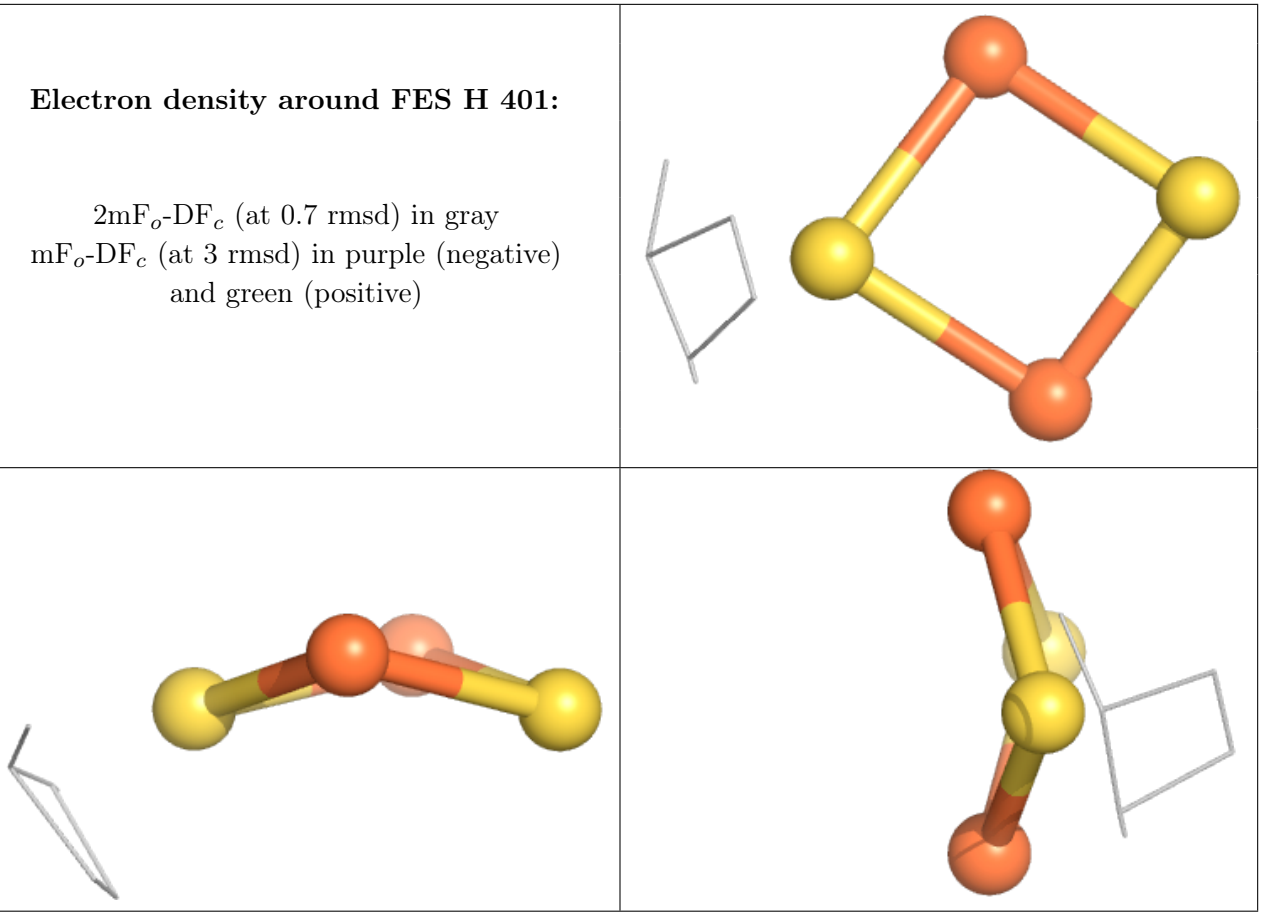
Electron density around FES G 401:

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



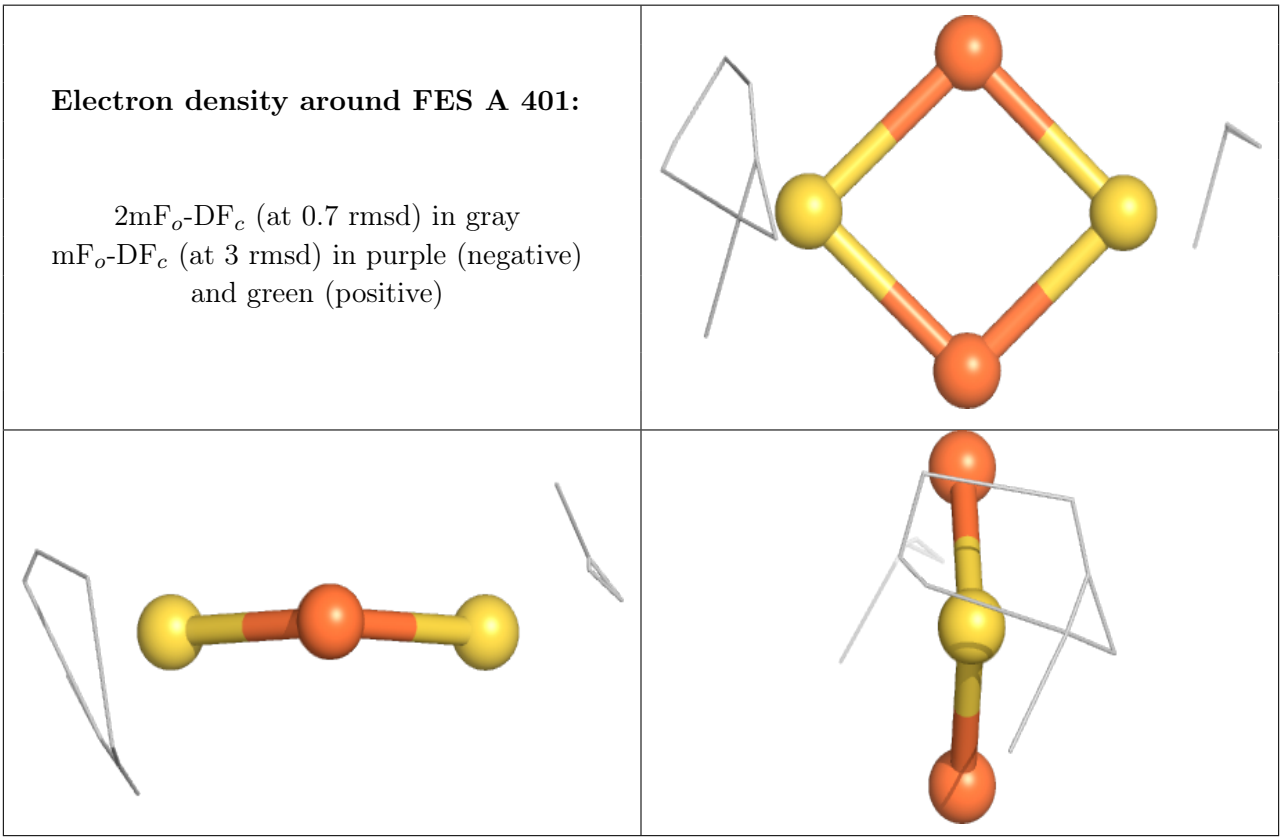
Electron density around FES H 401:

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



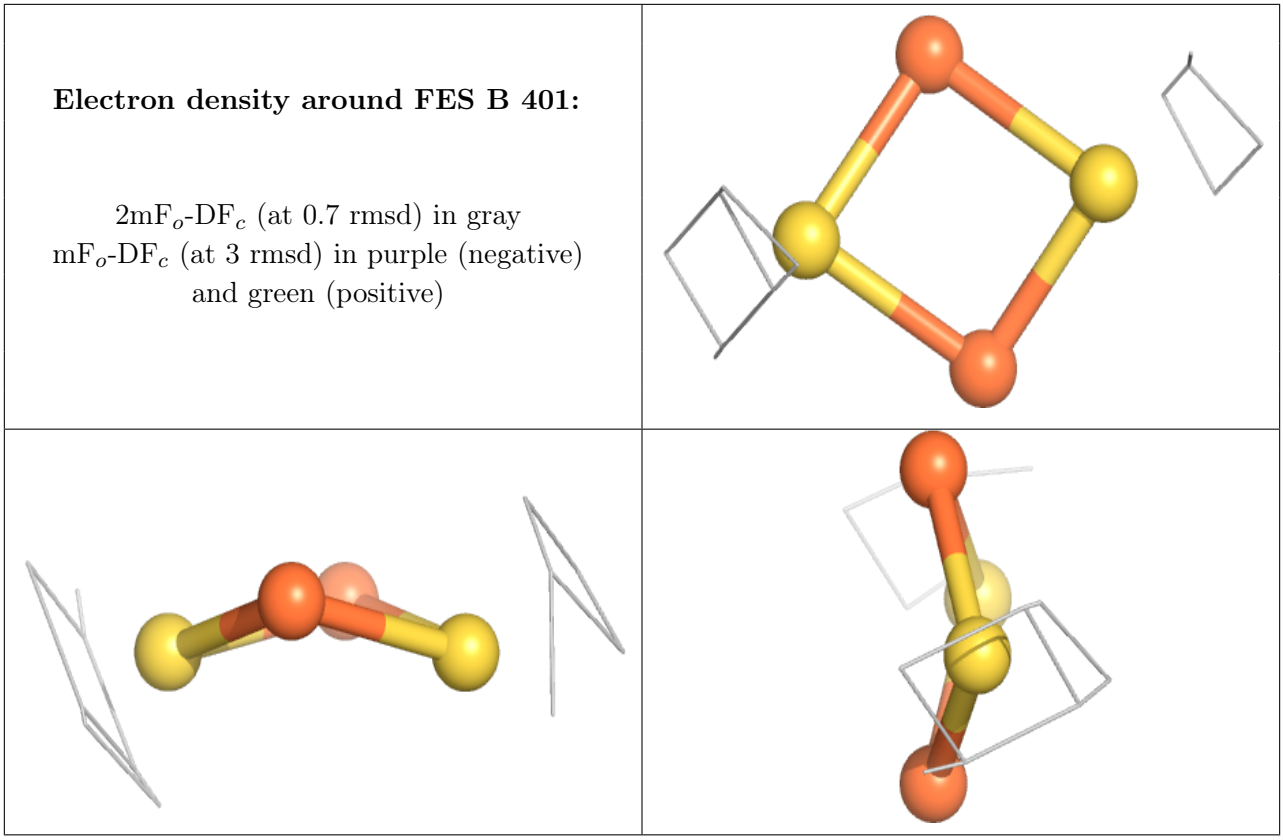
Electron density around FES A 401:

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



Electron density around FES 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)



6.5 Other polymers (i)

There are no such residues in this entry.