



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 02:54 AM BST

PDB ID : 6WN3
Title : Structure of the Rieske non-heme iron oxygenase SxtT
Authors : Bridwell-Rabb, J.; Liu, J.
Deposited on : 2020-04-22
Resolution : 1.86 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.

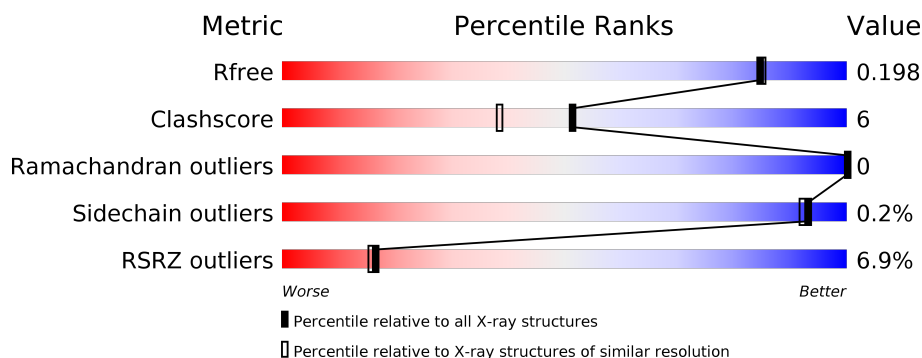
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

i

X-RAY DIFFRACTION

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	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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 on 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	334	<div> <div></div> <div>6%</div> <div>88%</div> <div>9%</div> <div></div> </div>	•
1	B	334	<div> <div></div> <div>5%</div> <div>91%</div> <div>7%</div> <div></div> </div>	•
1	C	334	<div> <div></div> <div>9%</div> <div>87%</div> <div>8%</div> <div>5%</div> </div>	

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
3	GOL	A	503	-	-	X	-

2 Entry composition [i](#)

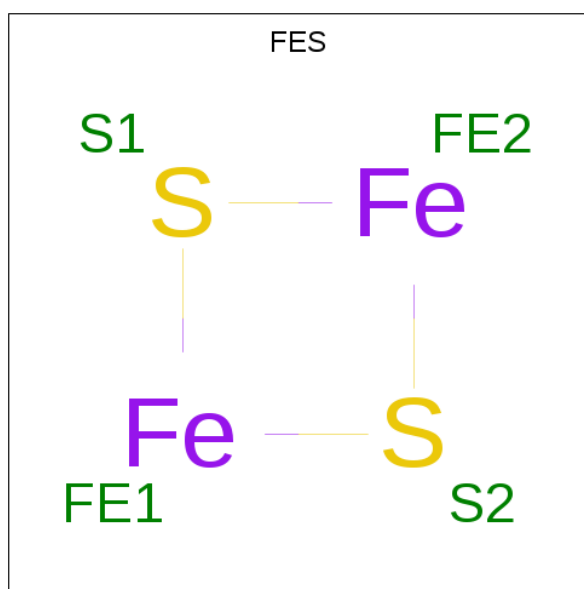
There are 6 unique types of molecules in this entry. The entry contains 9106 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SxtT.

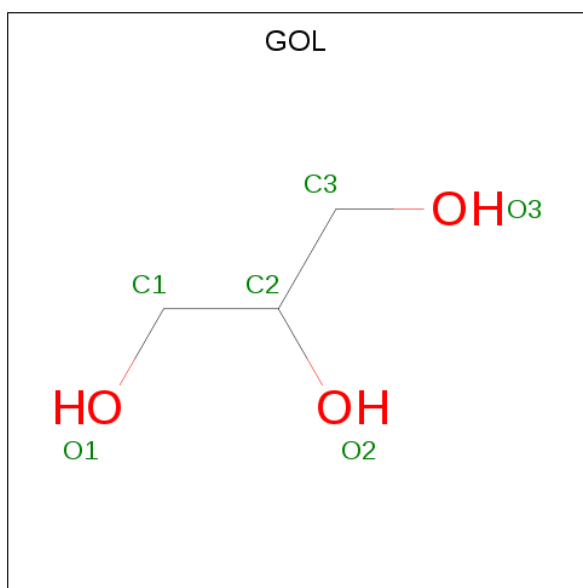
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	12	0
			2694	1708	460	500	26			
1	B	328	Total	C	N	O	S	0	21	0
			2800	1770	480	521	29			
1	C	316	Total	C	N	O	S	0	7	0
			2600	1657	443	477	23			

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe₂S₂) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			4	2	2		
2	B	1	Total	Fe	S	0	0
			4	2	2		
2	C	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe) (labeled as "Ligand of Interest" by author).

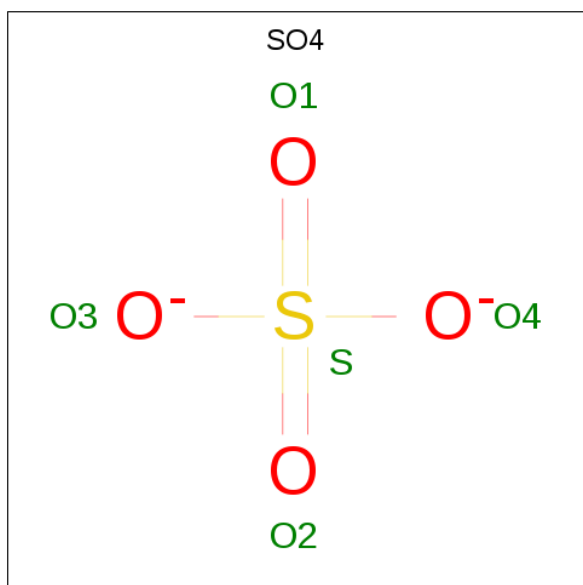
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Fe	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Fe	0	0
			1	1		
4	C	1	Total	Fe	0	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

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

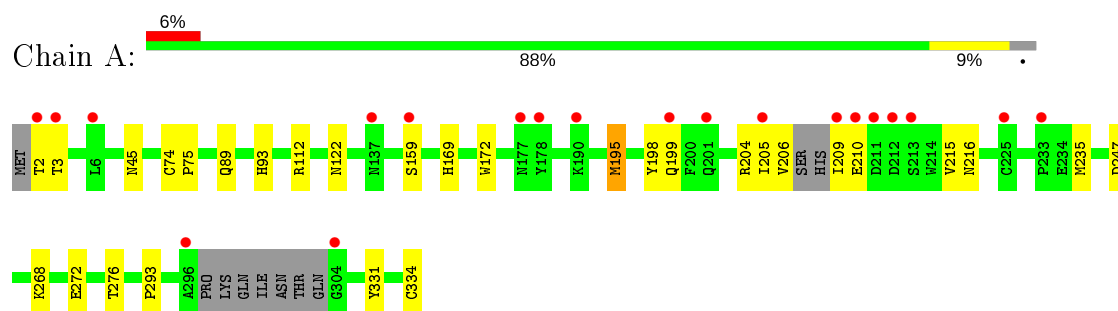
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	309	Total	O	0	0
			309	309		
6	B	368	Total	O	0	2
			368	368		
6	C	206	Total	O	0	0
			206	206		

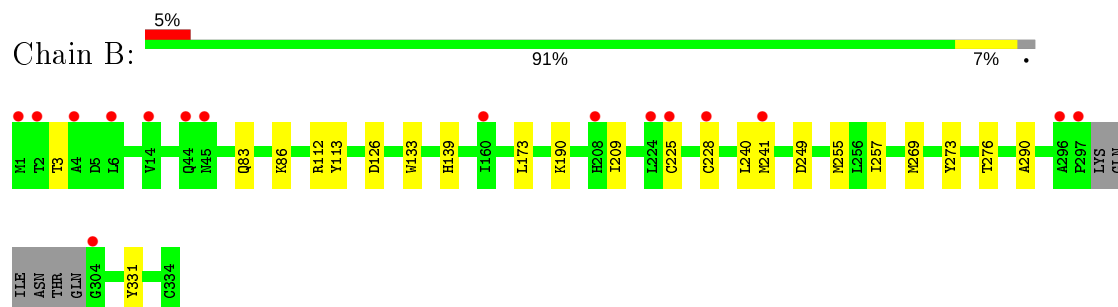
3 Residue-property plots [i](#)

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.

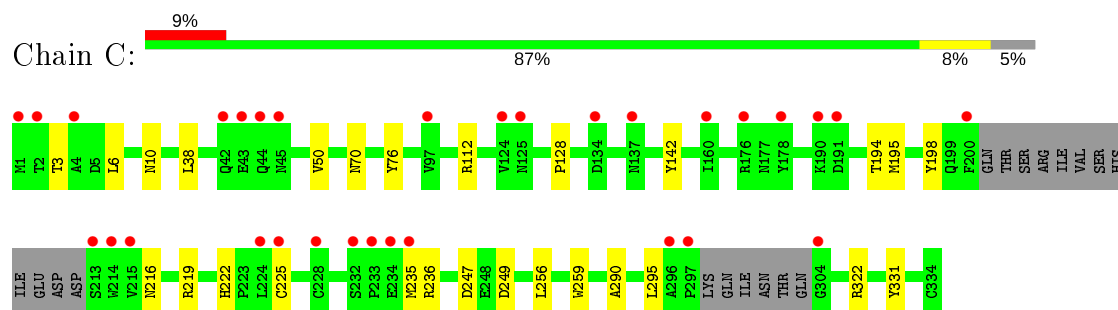
• Molecule 1: SxtT



• Molecule 1: SxtT



• Molecule 1: SxtT



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	152.28Å 159.58Å 115.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.92 – 1.86 29.92 – 1.86	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.92-1.86) 100.0 (29.92-1.86)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 1.87Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.170 , 0.198 0.170 , 0.198	Depositor DCC
R_{free} test set	5895 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.507	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9106	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4, 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	0.40	0/2767	0.58	0/3772
1	B	0.41	0/2876	0.57	0/3920
1	C	0.35	0/2674	0.54	0/3646
All	All	0.39	0/8317	0.56	0/11338

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2694	0	2590	39	2
1	B	2800	0	2686	24	1
1	C	2600	0	2509	24	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
3	A	24	0	32	8	0
3	B	18	0	24	6	0
3	C	12	0	16	0	0
4	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	A	25	0	0	0	0
5	B	20	0	0	0	0
5	C	15	0	0	1	0
6	A	309	0	0	6	2
6	B	368	0	0	7	0
6	C	206	0	0	5	0
All	All	9106	0	7857	91	5

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 6.

All (91) 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:A:172:TRP:CZ3	1:A:205:ILE:CD1	2.21	1.22
1:A:172:TRP:CZ3	1:A:205:ILE:HD12	1.74	1.19
1:A:172:TRP:HZ3	1:A:205:ILE:HD12	0.98	1.14
1:A:172:TRP:HZ3	1:A:205:ILE:CD1	1.68	1.00
1:A:172:TRP:CZ3	1:A:205:ILE:HD11	1.93	0.99
1:A:206:VAL:HG12	1:A:235:MET:HE1	1.42	0.97
1:A:2:THR:HG22	1:A:3:THR:N	1.78	0.97
1:B:228[B]:CYS:HB3	6:B:909[B]:HOH:O	1.70	0.92
1:B:133:TRP:HE1	3:B:502:GOL:H31	1.34	0.91
1:B:257:ILE:HG21	1:B:269[B]:MET:HE2	1.54	0.89
1:C:235:MET:HE1	1:C:259:TRP:HB2	1.57	0.85
1:A:2:THR:HG22	1:A:3:THR:H	1.40	0.84
1:A:172:TRP:CH2	1:A:205:ILE:HD11	2.13	0.84
1:A:206:VAL:CG1	1:A:235:MET:HE1	2.08	0.82
1:C:112:ARG:NH1	6:C:601:HOH:O	2.18	0.77
1:A:45:ASN:H	3:A:502:GOL:H12	1.51	0.76
1:B:190:LYS:HD3	1:B:190:LYS:H	1.51	0.75
1:B:257:ILE:HG21	1:B:269[B]:MET:CE	2.17	0.74
3:A:503:GOL:H11	6:A:744:HOH:O	1.90	0.71
3:A:503:GOL:C1	6:A:744:HOH:O	2.38	0.70
1:B:290:ALA:HB3	3:B:503:GOL:H11	1.74	0.69
1:B:3:THR:HB	1:B:331:TYR:CD1	2.28	0.68
1:A:2:THR:CG2	1:A:3:THR:N	2.50	0.68
1:C:3:THR:HB	1:C:331:TYR:CD1	2.29	0.67
1:C:194:THR:HG22	1:C:219:ARG:HB2	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:236:ARG:NH2	5:C:506:SO4:O3	2.28	0.66
1:C:195:MET:CE	1:C:198:TYR:CD1	2.82	0.62
1:C:70:ASN:ND2	6:C:605:HOH:O	2.32	0.62
1:B:190:LYS:H	1:B:190:LYS:CD	2.13	0.61
3:B:502:GOL:H32	6:B:759:HOH:O	2.00	0.61
1:A:195:MET:HE3	1:A:198:TYR:HD1	1.66	0.61
1:A:195:MET:HE3	1:A:198:TYR:CD1	2.36	0.60
1:B:190:LYS:HD3	1:B:190:LYS:N	2.17	0.59
1:C:195:MET:HE3	1:C:198:TYR:CD1	2.39	0.57
1:A:199:GLN:HG2	1:A:215[A]:VAL:HG12	1.86	0.57
1:C:219:ARG:NH2	6:C:608:HOH:O	2.37	0.57
1:A:195:MET:HE1	1:A:198:TYR:CE1	2.40	0.56
1:A:112:ARG:HH12	3:A:503:GOL:H32	1.71	0.56
3:B:502:GOL:O2	6:B:601:HOH:O	2.16	0.56
1:A:45:ASN:H	3:A:502:GOL:C1	2.18	0.56
1:B:241[A]:MET:HG3	1:B:255[A]:MET:HG2	1.88	0.55
1:A:172:TRP:CH2	1:A:205:ILE:CD1	2.77	0.55
1:C:195:MET:HE2	1:C:198:TYR:CD1	2.41	0.55
1:B:290:ALA:H	3:B:503:GOL:H32	1.72	0.54
1:B:173[A]:LEU:HD13	1:B:276:THR:HG21	1.89	0.54
1:C:295:LEU:HD12	1:C:322:ARG:HG3	1.91	0.53
1:A:195:MET:CE	1:A:198:TYR:CD1	2.92	0.53
1:A:45:ASN:ND2	6:A:612:HOH:O	2.40	0.53
1:A:293:PRO:HA	1:A:334[B]:CYS:HB2	1.90	0.52
1:A:112:ARG:HH22	3:A:503:GOL:H31	1.75	0.52
1:A:93:HIS:HB3	3:A:505:GOL:H2	1.92	0.51
1:A:159[B]:SER:OG	6:A:601:HOH:O	2.18	0.51
1:B:139:HIS:HD2	6:B:920:HOH:O	1.92	0.51
1:C:38:LEU:HG	1:C:50:VAL:HG22	1.93	0.50
1:C:198:TYR:CE1	1:C:216:ASN:HB2	2.47	0.49
1:A:209:ILE:HG13	1:A:210:GLU:N	2.27	0.49
1:A:89:GLN:NE2	6:A:615:HOH:O	2.45	0.48
1:A:74[B]:CYS:SG	1:A:75:PRO:HD2	2.53	0.48
1:C:198:TYR:CZ	1:C:216:ASN:HB2	2.49	0.48
1:C:194:THR:HG22	1:C:219:ARG:CB	2.42	0.48
1:B:209:ILE:H	1:B:209:ILE:HD12	1.79	0.48
1:A:205:ILE:HG22	1:A:268:LYS:HB3	1.95	0.47
1:A:169:HIS:CE1	1:A:276[B]:THR:HG23	2.50	0.47
1:A:122:ASN:HB3	6:A:756:HOH:O	2.14	0.47
1:B:249:ASP:HB3	6:B:687:HOH:O	2.14	0.47
1:C:6:LEU:HD22	1:C:10:ASN:ND2	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:MET:HE2	1:A:195:MET:HB3	1.65	0.46
1:B:112[B]:ARG:HD3	1:B:113:TYR:CZ	2.52	0.45
1:A:2:THR:HG22	1:A:3:THR:OG1	2.17	0.45
1:B:209:ILE:HD12	1:B:209:ILE:N	2.32	0.44
1:C:290:ALA:O	6:C:603:HOH:O	2.21	0.44
1:A:112:ARG:HH12	3:A:503:GOL:C3	2.30	0.44
1:A:247:ASP:HA	1:A:331:TYR:OH	2.17	0.44
1:B:269[B]:MET:HE2	1:B:269[B]:MET:HB3	1.87	0.44
1:A:2:THR:CG2	1:A:3:THR:H	2.08	0.43
1:B:83:GLN:HB2	3:B:504:GOL:H31	1.99	0.43
1:A:204:ARG:HB3	1:A:272:GLU:OE2	2.17	0.43
1:C:235:MET:CE	1:C:259:TRP:HB2	2.40	0.43
1:B:225[B]:CYS:SG	1:B:240:LEU:HD11	2.58	0.43
1:C:235:MET:HE3	1:C:235:MET:HB2	1.82	0.42
1:B:86:LYS:HE3	6:B:705:HOH:O	2.20	0.42
1:C:249:ASP:HB3	6:C:649:HOH:O	2.19	0.42
1:C:128:PRO:HG2	1:C:225[B]:CYS:SG	2.60	0.42
1:B:83:GLN:NE2	6:B:614:HOH:O	2.50	0.42
1:C:142:TYR:CE2	1:C:256:LEU:HD13	2.55	0.42
1:C:222:HIS:HB3	1:C:225[B]:CYS:HB3	2.02	0.41
1:B:269[B]:MET:HE1	1:B:273:TYR:HE1	1.84	0.41
1:A:198:TYR:CE1	1:A:216:ASN:HB2	2.56	0.41
1:C:247:ASP:HA	1:C:331:TYR:OH	2.20	0.41
1:A:209:ILE:HG13	1:A:210:GLU:H	1.86	0.40
1:B:269[B]:MET:CE	1:B:273:TYR:CE1	3.04	0.40

All (5) 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:A:206:VAL:O	1:A:209:ILE:CG2[6_445]	1.84	0.36
1:A:205:ILE:CG1	1:A:205:ILE:CG2[6_445]	2.06	0.14
6:A:744:HOH:O	6:A:778:HOH:O[3_556]	2.10	0.10
6:A:808:HOH:O	6:A:889:HOH:O[3_556]	2.16	0.04
1:B:126:ASP:OD2	1:B:190:LYS:NZ[3_555]	2.19	0.01

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	329/334 (98%)	316 (96%)	13 (4%)	0	100	100
1	B	345/334 (103%)	327 (95%)	18 (5%)	0	100	100
1	C	316/334 (95%)	305 (96%)	11 (4%)	0	100	100
All	All	990/1002 (99%)	948 (96%)	42 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	305/303 (101%)	304 (100%)	1 (0%)	92	91
1	B	318/303 (105%)	318 (100%)	0	100	100
1	C	292/303 (96%)	291 (100%)	1 (0%)	92	91
All	All	915/909 (101%)	913 (100%)	2 (0%)	93	92

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	195	MET
1	C	76	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 27 ligands modelled in this entry, 3 are monoatomic - leaving 24 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	$\# Z > 2$
5	SO4	C	507	-	4,4,4	0.16	0	6,6,6	0.10	0
5	SO4	A	511	-	4,4,4	0.39	0	6,6,6	0.04	0
5	SO4	B	507	-	4,4,4	0.16	0	6,6,6	0.05	0
3	GOL	A	503	-	5,5,5	1.15	1 (20%)	5,5,5	0.90	0
3	GOL	A	504	-	5,5,5	0.84	0	5,5,5	1.03	0
2	FES	C	501	1	0,4,4	0.00	-	-		
2	FES	B	501	1	0,4,4	0.00	-	-		
5	SO4	A	508	-	4,4,4	0.17	0	6,6,6	0.17	0
3	GOL	C	503	-	5,5,5	1.07	0	5,5,5	0.88	0
3	GOL	B	503	-	5,5,5	0.89	0	5,5,5	0.85	0
5	SO4	A	509	-	4,4,4	0.14	0	6,6,6	0.05	0
5	SO4	A	507	-	4,4,4	0.14	0	6,6,6	0.07	0
5	SO4	B	508	-	4,4,4	0.15	0	6,6,6	0.19	0
5	SO4	C	506	-	4,4,4	0.15	0	6,6,6	0.08	0
2	FES	A	501	1	0,4,4	0.00	-	-		
5	SO4	B	506	-	4,4,4	0.17	0	6,6,6	0.12	0
3	GOL	B	504	-	5,5,5	0.83	0	5,5,5	0.97	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	C	502	-	5,5,5	0.91	0	5,5,5	0.95	0
3	GOL	A	505	-	5,5,5	0.98	0	5,5,5	0.98	0
3	GOL	B	502	-	5,5,5	0.93	0	5,5,5	0.98	0
5	SO4	B	509	-	4,4,4	0.15	0	6,6,6	0.11	0
5	SO4	C	505	-	4,4,4	0.17	0	6,6,6	0.11	0
3	GOL	A	502	-	5,5,5	0.85	0	5,5,5	1.01	0
5	SO4	A	510	-	4,4,4	0.13	0	6,6,6	0.13	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	505	-	-	4/4/4/4	-
3	GOL	B	502	-	-	3/4/4/4	-
3	GOL	A	503	-	-	4/4/4/4	-
3	GOL	A	504	-	-	4/4/4/4	-
3	GOL	C	503	-	-	0/4/4/4	-
2	FES	C	501	1	-	-	0/1/1/1
3	GOL	B	503	-	-	2/4/4/4	-
2	FES	A	501	1	-	-	0/1/1/1
2	FES	B	501	1	-	-	0/1/1/1
3	GOL	A	502	-	-	2/4/4/4	-
3	GOL	B	504	-	-	2/4/4/4	-
3	GOL	C	502	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	503	GOL	O2-C2	-2.09	1.37	1.43

There are no bond angle outliers.

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	503	GOL	O1-C1-C2-C3
3	A	504	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	C	502	GOL	C1-C2-C3-O3
3	A	505	GOL	O1-C1-C2-C3
3	A	505	GOL	O1-C1-C2-O2
3	A	503	GOL	C1-C2-C3-O3
3	A	504	GOL	C1-C2-C3-O3
3	B	503	GOL	O1-C1-C2-C3
3	B	504	GOL	C1-C2-C3-O3
3	B	502	GOL	O1-C1-C2-C3
3	A	503	GOL	O2-C2-C3-O3
3	A	504	GOL	O1-C1-C2-O2
3	A	504	GOL	O2-C2-C3-O3
3	C	502	GOL	O2-C2-C3-O3
3	A	503	GOL	O1-C1-C2-O2
3	B	502	GOL	O1-C1-C2-O2
3	B	502	GOL	O2-C2-C3-O3
3	B	503	GOL	O1-C1-C2-O2
3	B	504	GOL	O2-C2-C3-O3
3	A	502	GOL	O2-C2-C3-O3
3	A	502	GOL	C1-C2-C3-O3
3	A	505	GOL	O2-C2-C3-O3
3	A	505	GOL	C1-C2-C3-O3

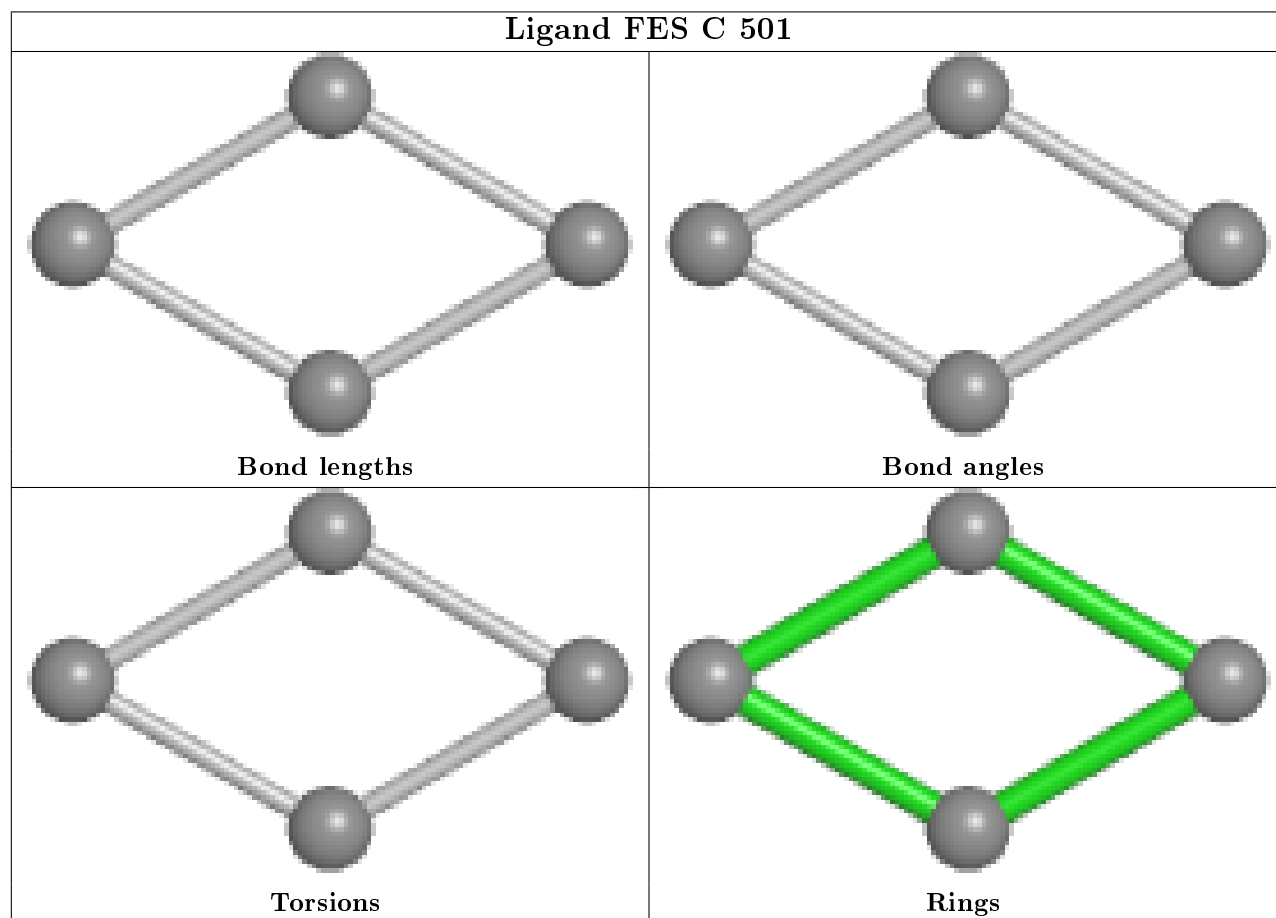
There are no ring outliers.

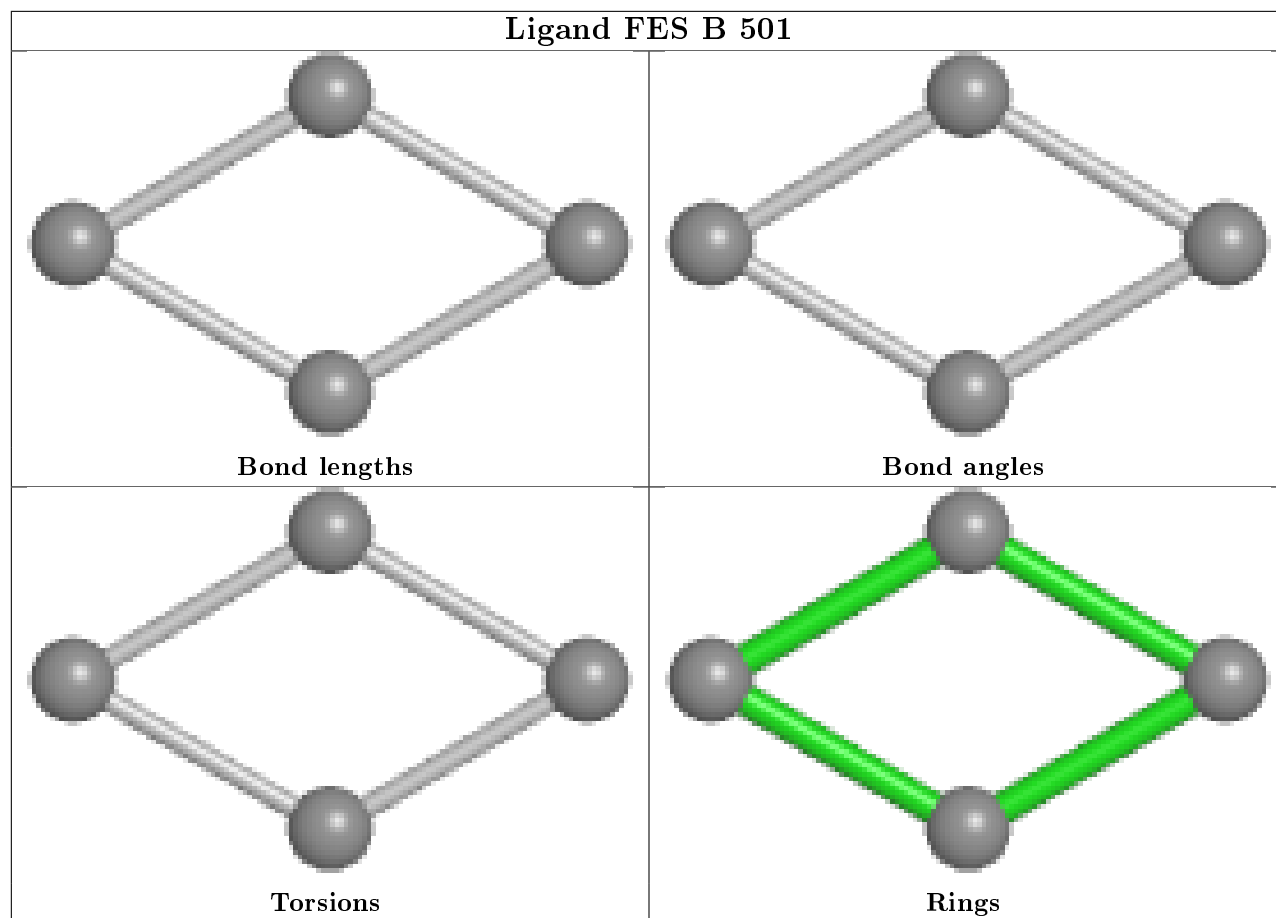
7 monomers are involved in 15 short contacts:

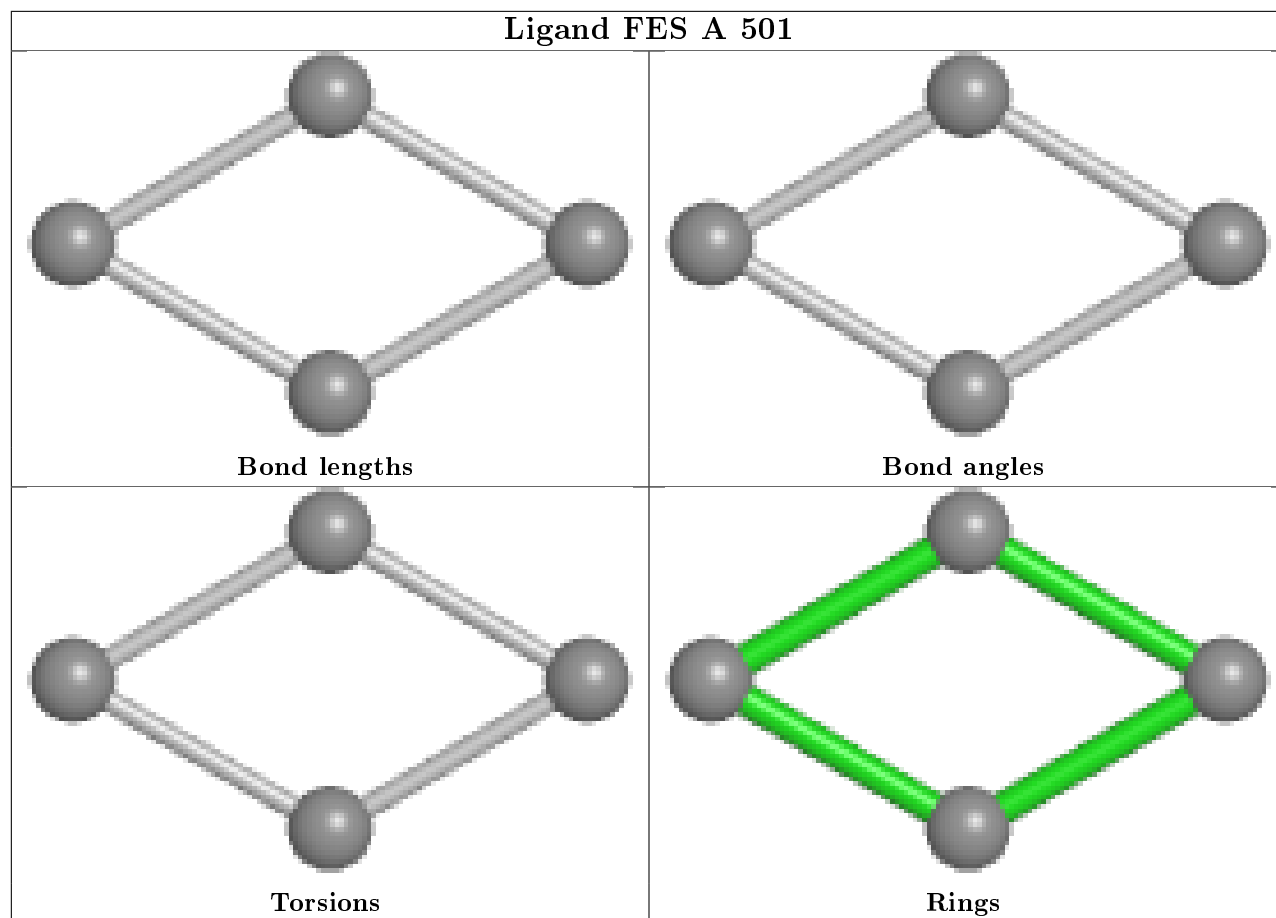
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	GOL	5	0
3	B	503	GOL	2	0
5	C	506	SO4	1	0
3	B	504	GOL	1	0
3	A	505	GOL	1	0
3	B	502	GOL	3	0
3	A	502	GOL	2	0

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 than 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.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	324/334 (97%)	0.25	20 (6%)	20 20	22, 36, 71, 103	0
1	B	328/334 (98%)	0.10	16 (4%)	29 28	23, 31, 56, 93	0
1	C	316/334 (94%)	0.40	31 (9%)	7 7	27, 42, 73, 103	0
All	All	968/1002 (96%)	0.25	67 (6%)	16 16	22, 36, 68, 103	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	233	PRO	7.1
1	C	297	PRO	6.8
1	B	297	PRO	6.7
1	A	3	THR	6.6
1	A	2	THR	6.2
1	C	200[A]	PHE	6.0
1	C	1	MET	5.9
1	C	213	SER	5.8
1	A	205	ILE	5.3
1	A	178	TYR	5.2
1	A	209	ILE	5.1
1	B	1	MET	5.0
1	A	213	SER	4.7
1	C	234	GLU	4.5
1	B	44	GLN	4.4
1	A	190	LYS	4.2
1	C	214	TRP	4.1
1	A	6	LEU	4.1
1	C	4	ALA	3.9
1	C	190	LYS	3.9
1	C	178	TYR	3.8
1	C	232	SER	3.8
1	C	304	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	235	MET	3.2
1	A	210	GLU	3.2
1	B	4	ALA	3.2
1	C	134	ASP	3.1
1	B	225[A]	CYS	3.1
1	C	43	GLU	3.0
1	C	2	THR	3.0
1	C	45	ASN	3.0
1	C	160	ILE	2.9
1	B	208	HIS	2.9
1	A	212	ASP	2.8
1	B	241[A]	MET	2.7
1	A	296	ALA	2.7
1	B	296	ALA	2.7
1	B	14	VAL	2.6
1	B	224	LEU	2.6
1	C	191	ASP	2.6
1	A	233	PRO	2.6
1	B	6	LEU	2.6
1	C	137	ASN	2.6
1	B	228[A]	CYS	2.5
1	C	224	LEU	2.4
1	A	159[A]	SER	2.4
1	C	42	GLN	2.3
1	A	304	GLY	2.3
1	A	211	ASP	2.3
1	A	177	ASN	2.3
1	C	124	VAL	2.3
1	C	225[A]	CYS	2.3
1	A	201	GLN	2.2
1	C	296	ALA	2.2
1	C	97	VAL	2.2
1	B	2	THR	2.2
1	A	199	GLN	2.2
1	B	160	ILE	2.1
1	B	304	GLY	2.1
1	C	125	ASN	2.1
1	C	44	GLN	2.1
1	C	215	VAL	2.1
1	C	228[A]	CYS	2.1
1	B	45	ASN	2.1
1	C	176	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	137	ASN	2.0
1	A	225[A]	CYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	SO4	A	509	5/5	0.70	0.34	92,95,115,132	0
5	SO4	A	510	5/5	0.74	0.37	74,75,92,103	0
3	GOL	C	503	6/6	0.75	0.33	59,66,71,79	0
5	SO4	B	509	5/5	0.81	0.30	80,81,113,120	0
5	SO4	C	505	5/5	0.81	0.39	57,66,93,105	0
3	GOL	C	502	6/6	0.81	0.20	50,62,65,66	0
5	SO4	A	511	5/5	0.83	0.42	89,91,113,116	0
3	GOL	A	503	6/6	0.85	0.16	52,57,62,69	0
5	SO4	C	506	5/5	0.85	0.32	94,98,107,119	0
5	SO4	B	507	5/5	0.86	0.35	74,84,107,110	0
5	SO4	B	508	5/5	0.87	0.16	62,67,74,100	0
3	GOL	A	504	6/6	0.87	0.26	54,64,70,78	0
5	SO4	C	507	5/5	0.87	0.39	73,83,99,102	0
3	GOL	A	505	6/6	0.88	0.24	40,55,65,70	0
3	GOL	B	503	6/6	0.89	0.28	46,56,61,74	0
3	GOL	A	502	6/6	0.90	0.26	53,62,76,77	0
5	SO4	A	507	5/5	0.90	0.32	87,87,109,110	0
3	GOL	B	504	6/6	0.91	0.30	37,60,66,67	0
3	GOL	B	502	6/6	0.92	0.12	32,42,52,63	0
5	SO4	B	506	5/5	0.96	0.17	49,49,81,82	0
2	FES	B	501	4/4	0.97	0.12	28,28,29,29	0

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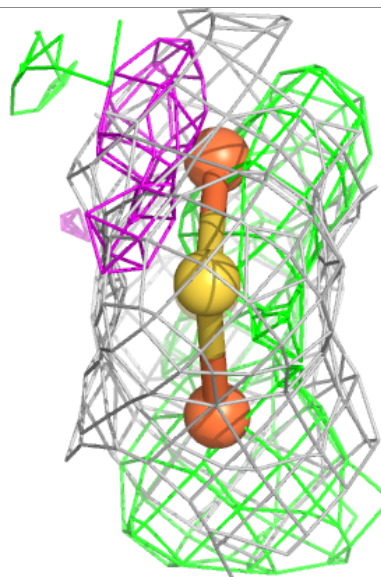
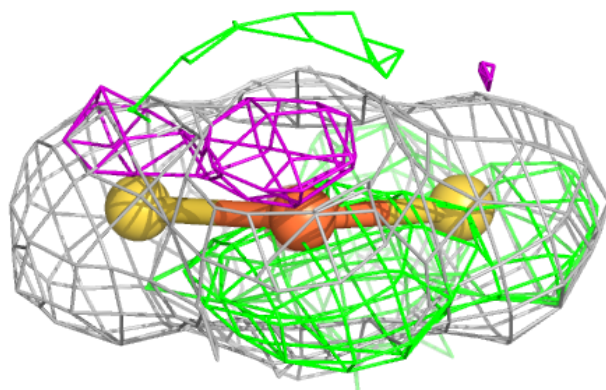
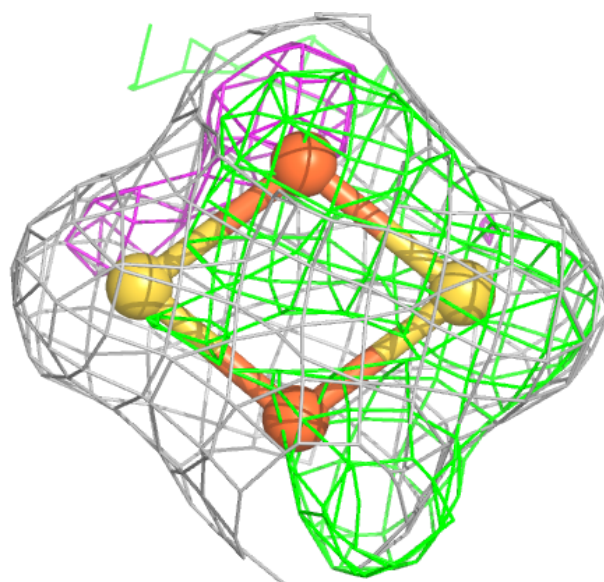
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	SO4	A	508	5/5	0.97	0.12	52,54,59,60	0
2	FES	C	501	4/4	0.98	0.08	30,31,32,33	0
2	FES	A	501	4/4	0.99	0.10	23,23,24,25	0
4	FE	C	504	1/1	1.00	0.09	30,30,30,30	0
4	FE	A	506	1/1	1.00	0.08	33,33,33,33	0
4	FE	B	505	1/1	1.00	0.12	26,26,26,26	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.

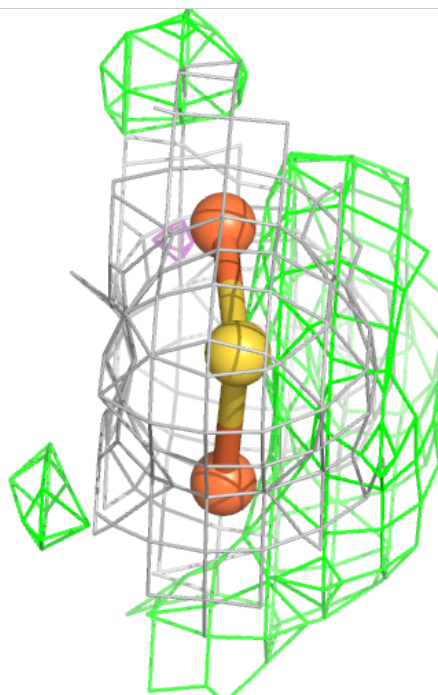
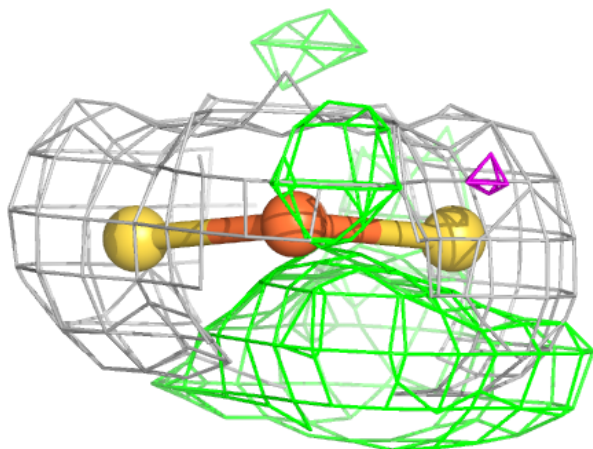
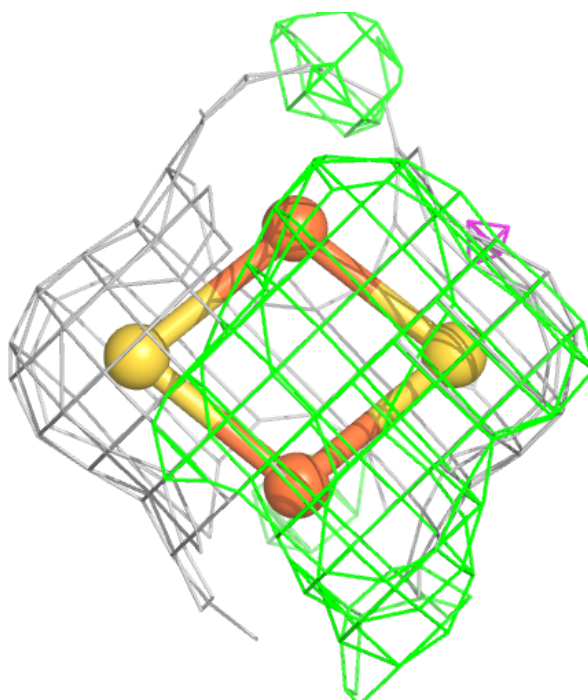
Electron density around FES B 501:

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



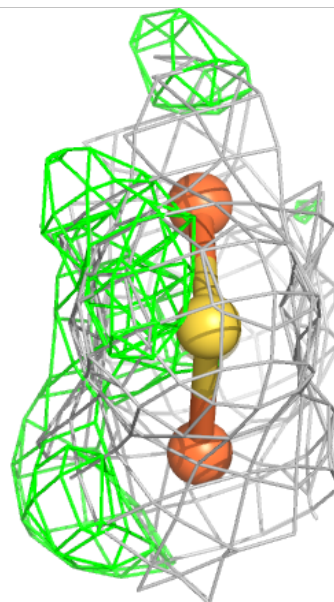
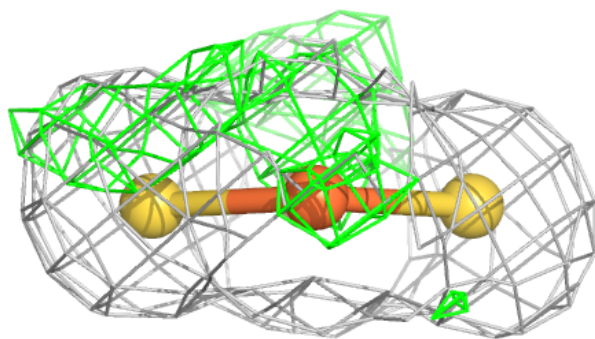
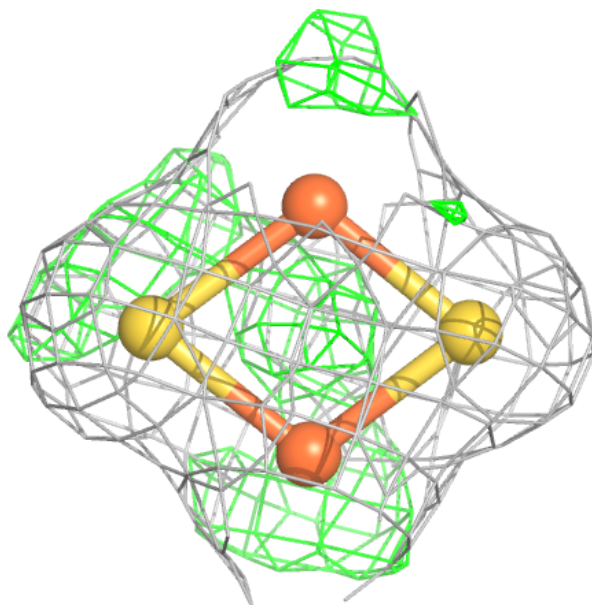
Electron density around FES C 501:

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



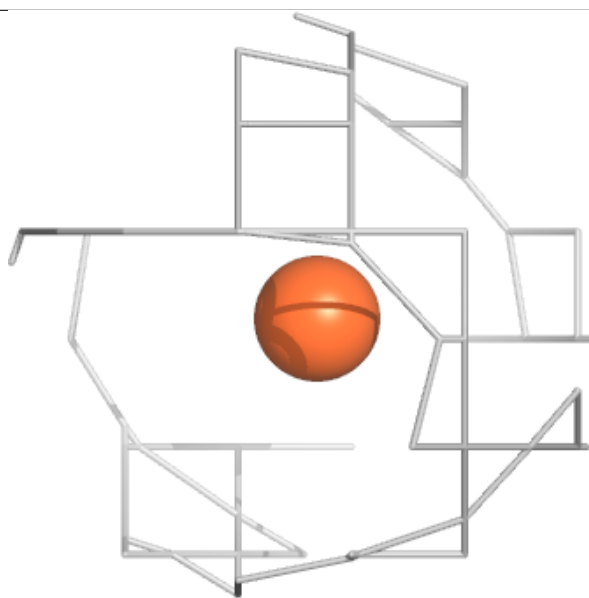
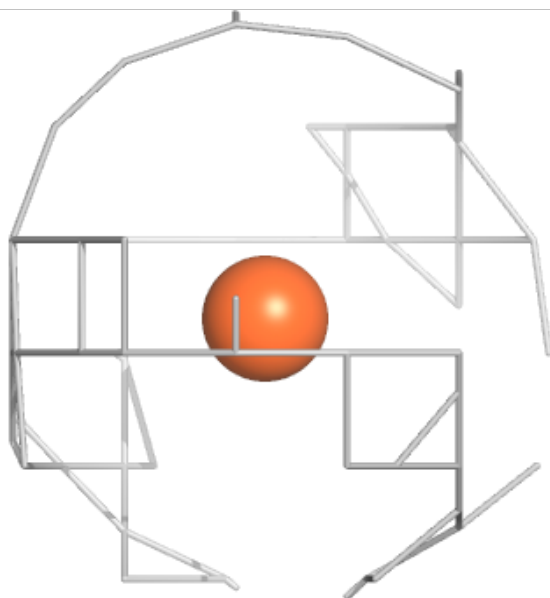
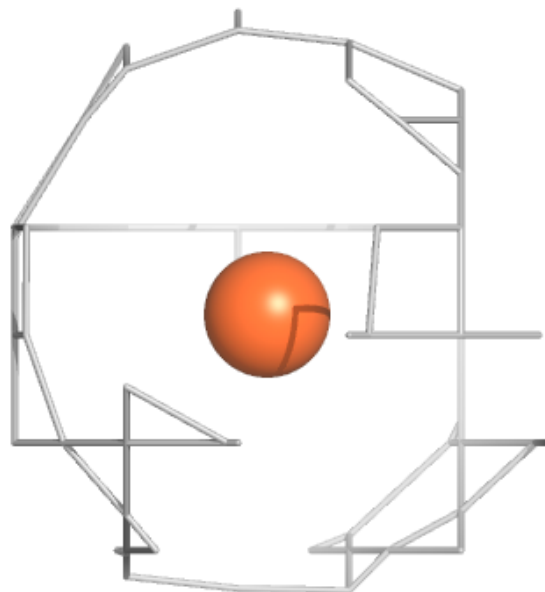
Electron density around FES A 501:

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



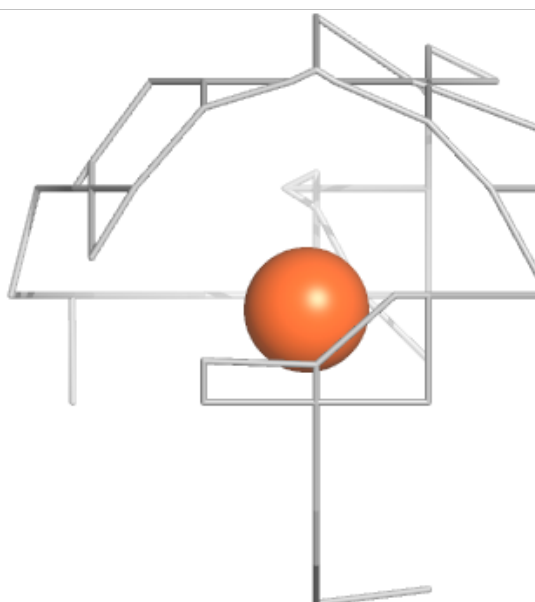
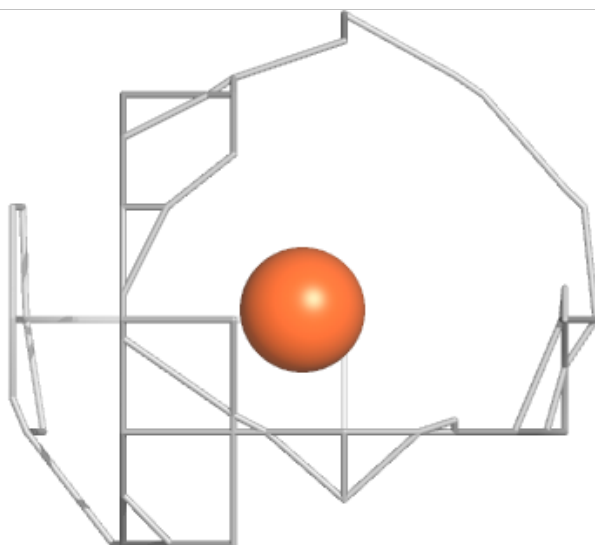
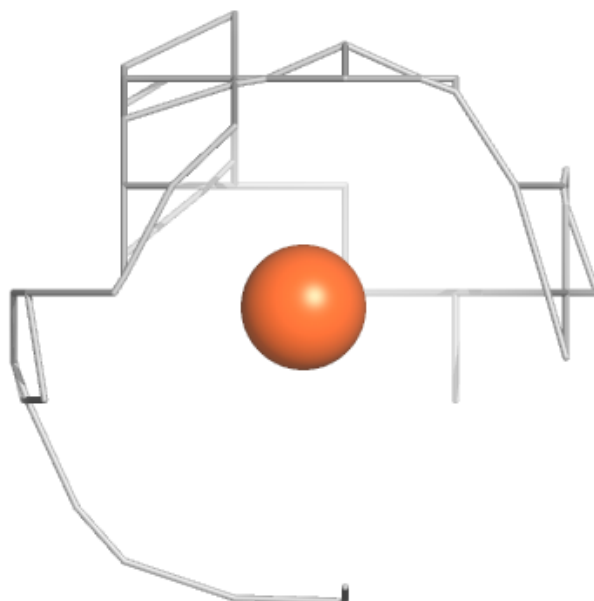
Electron density around FE C 504:

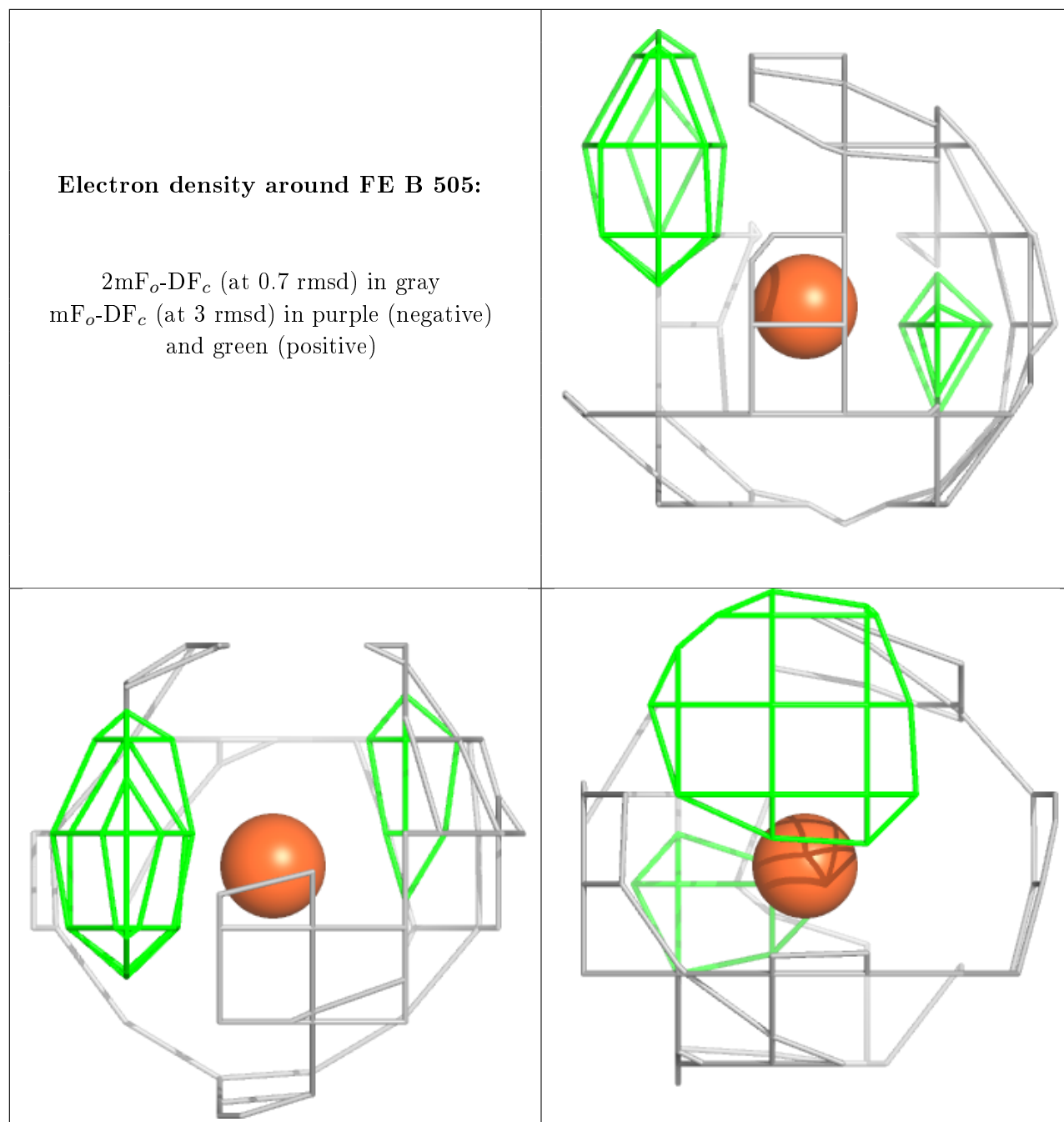
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FE A 506:

$2mF_o-DF_c$ (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.