



wwPDB X-ray Structure Validation Summary Report

Mar 9, 2026 – 09:32 PM UTC

PDB ID : 9FDJ / pdb_00009fdj
Title : Crystal structure of the NuoEF variant R66G (NuoF) from Aquifex aeolicus bound to NADH under anoxic conditions (short soak)
Authors : Wohlwend, D.; Friedrich, T.; Goeppert-Asadollahpour, S.
Deposited on : 2024-05-17
Resolution : 1.70 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  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
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

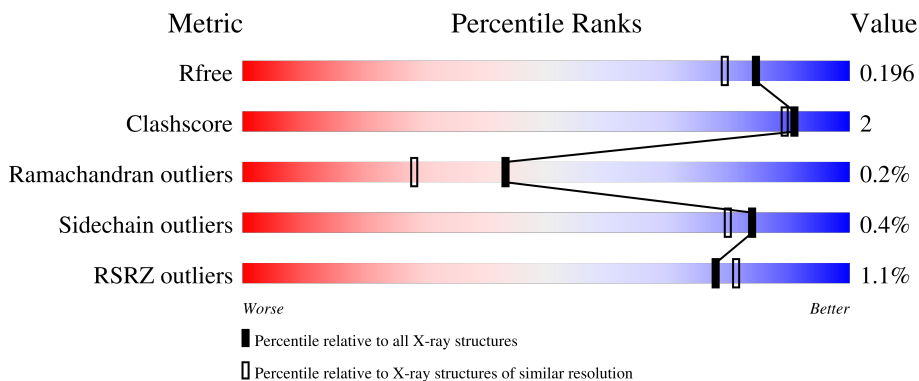
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	160	94%
1	C	160	93%
2	B	434	91% 5%
2	D	434	90% 6%

2 Entry composition i

There are 11 unique types of molecules in this entry. The entry contains 10206 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 NADH-quinone oxidoreductase subunit E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	156	1268	821	204	234	9	0	0	0
1	C	156	1268	821	204	234	9	0	0	0

- Molecule 2 is a protein called NADH-quinone oxidoreductase subunit F.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	416	3274	2107	541	613	13	0	1	0
2	D	418	3306	2125	547	620	14	0	3	0

There are 18 discrepancies between the modelled and reference sequences:

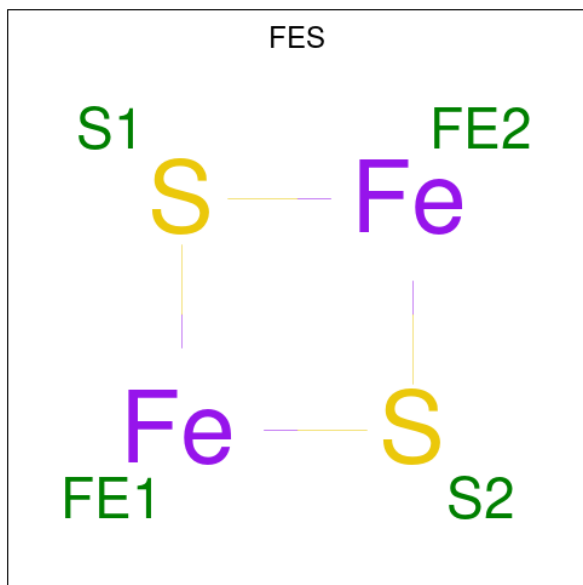
Chain	Residue	Modelled	Actual	Comment	Reference
B	66	GLY	ARG	engineered mutation	UNP O66841
B	427	ALA	-	expression tag	UNP O66841
B	428	GLY	-	expression tag	UNP O66841
B	429	HIS	-	expression tag	UNP O66841
B	430	HIS	-	expression tag	UNP O66841
B	431	HIS	-	expression tag	UNP O66841
B	432	HIS	-	expression tag	UNP O66841
B	433	HIS	-	expression tag	UNP O66841
B	434	HIS	-	expression tag	UNP O66841
D	66	GLY	ARG	engineered mutation	UNP O66841
D	427	ALA	-	expression tag	UNP O66841
D	428	GLY	-	expression tag	UNP O66841
D	429	HIS	-	expression tag	UNP O66841
D	430	HIS	-	expression tag	UNP O66841
D	431	HIS	-	expression tag	UNP O66841
D	432	HIS	-	expression tag	UNP O66841

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Chain	Residue	Modelled	Actual	Comment	Reference
D	433	HIS	-	expression tag	UNP O66841
D	434	HIS	-	expression tag	UNP O66841

- Molecule 3 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



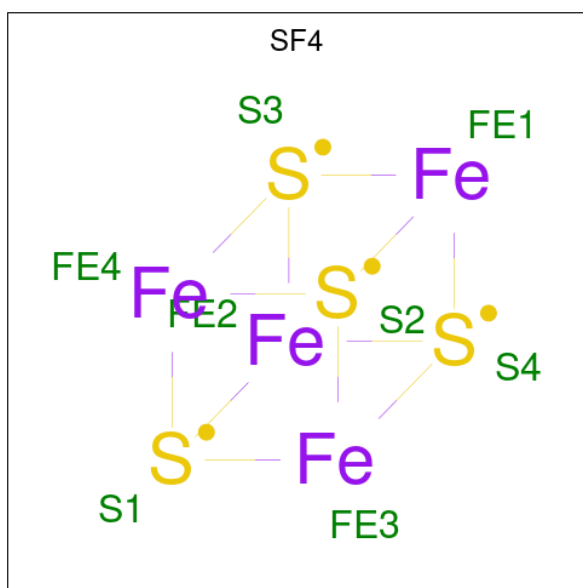
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	D	1	Total O S 5 4 1	0	0
4	D	1	Total O S 5 4 1	0	0

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Cl 1 1	0	0
5	B	1	Total Cl 1 1	0	0
5	C	2	Total Cl 2 2	0	0

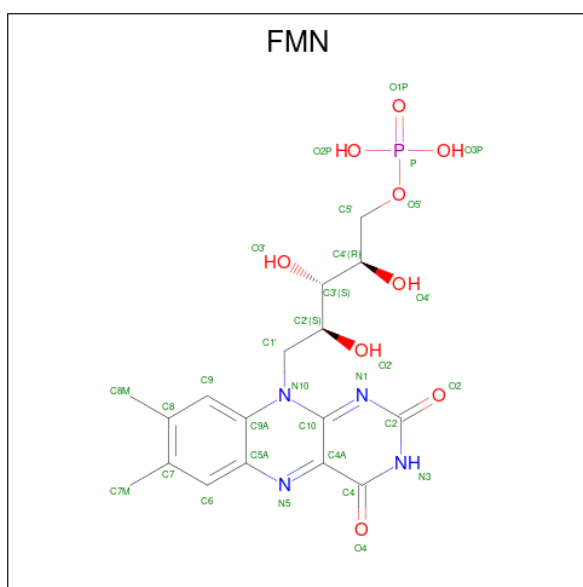
- Molecule 6 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄) (labeled as "Lig-

and of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total Fe S 8 4 4	0	0
6	D	1	Total Fe S 8 4 4	0	0

- Molecule 7 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$) (labeled as "Ligand of Interest" by depositor).





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

- Molecule 10 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	B	3	Total Na 3 3	0	0
10	D	5	Total Na 5 5	0	0

- Molecule 11 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	112	Total O 114 114	0	2
11	B	319	Total O 319 319	0	0
11	C	116	Total O 116 116	0	0
11	D	335	Total O 336 336	0	1

4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.24Å 116.16Å 189.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.93 – 1.70 47.93 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.93-1.70) 99.9 (47.93-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 1.70Å)	Xtrriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.158 , 0.184 0.170 , 0.196	Depositor DCC
R_{free} test set	7788 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtrriage
Anisotropy	0.219	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 32.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10206	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.28 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.0872e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: NAI, NA, SO4, SF4, FMN, CL, GOL, FES

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	0.51	0/1297	0.90	1/1752 (0.1%)
1	C	0.51	0/1297	0.90	1/1752 (0.1%)
2	B	0.52	0/3355	0.92	1/4542 (0.0%)
2	D	0.52	0/3387	0.90	0/4583
All	All	0.52	0/9336	0.91	3/12629 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	C	0	1
All	All	0	3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	74	ASP	CA-CB-CG	5.64	118.24	112.60
1	C	74	ASP	CA-CB-CG	5.46	118.06	112.60
2	B	298	ASP	CA-CB-CG	5.07	117.67	112.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	27	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	A	82	ARG	Sidechain
1	C	82	ARG	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1268	0	1269	2	0
1	C	1268	0	1269	4	0
2	B	3274	0	3245	14	0
2	D	3306	0	3273	16	0
3	A	4	0	0	0	0
3	C	4	0	0	0	0
4	A	15	0	0	0	0
4	B	5	0	0	0	0
4	C	10	0	0	0	0
4	D	10	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	2	0	0	0	0
6	B	8	0	0	0	0
6	D	8	0	0	0	0
7	B	31	0	19	0	0
7	D	31	0	19	0	0
8	B	28	0	12	1	0
8	D	27	0	11	0	0
9	B	6	0	8	0	0
9	D	6	0	8	0	0
10	B	3	0	0	0	0
10	D	5	0	0	0	0
11	A	114	0	0	0	0
11	B	319	0	0	5	0
11	C	116	0	0	1	0
11	D	336	0	0	2	0
All	All	10206	0	9133	32	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

The worst 5 of 32 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:190:GLU:OE1	2:D:195:LYS:HE2	1.93	0.68
2:D:233:MET:HE1	2:D:241:ILE:HD11	1.79	0.65
2:D:384:ASN:HD22	2:D:404:ARG:HH21	1.48	0.60
1:C:62:HIS:HE1	2:D:194:GLY:O	1.86	0.59
2:B:195:LYS:NZ	11:B:601:HOH:O	2.35	0.59

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	154/160 (96%)	149 (97%)	5 (3%)	0	100	100
1	C	154/160 (96%)	149 (97%)	5 (3%)	0	100	100
2	B	415/434 (96%)	405 (98%)	9 (2%)	1 (0%)	43	28
2	D	419/434 (96%)	410 (98%)	8 (2%)	1 (0%)	43	28
All	All	1142/1188 (96%)	1113 (98%)	27 (2%)	2 (0%)	43	28

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	95	GLU
2	D	95	GLU

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	142/146 (97%)	142 (100%)	0	100	100
1	C	142/146 (97%)	140 (99%)	2 (1%)	59	46
2	B	342/356 (96%)	341 (100%)	1 (0%)	86	83
2	D	345/356 (97%)	344 (100%)	1 (0%)	86	83
All	All	971/1004 (97%)	967 (100%)	4 (0%)	84	80

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

Mol	Chain	Res	Type
2	B	412	GLU
1	C	5	GLU
1	C	76	GLU
2	D	388	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	62	HIS
1	C	153	ASN
2	D	384	ASN
2	D	286	ASN
2	D	349	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 30 ligands modelled in this entry, 12 are monoatomic - leaving 18 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
4	SO4	A	203	-	4,4,4	0.31	0	6,6,6	0.11	0
4	SO4	D	506	-	4,4,4	0.29	0	6,6,6	0.07	0
4	SO4	D	505	-	4,4,4	0.33	0	6,6,6	0.25	0
6	SF4	B	501	2	0,12,12	-	-	-		
7	FMN	D	502	10	33,33,33	0.48	0	48,50,50	0.68	1 (2%)
9	GOL	B	504	-	5,5,5	0.11	0	5,5,5	0.25	0
3	FES	A	201	1	0,4,4	-	-	-		
7	FMN	B	502	-	33,33,33	0.60	0	48,50,50	0.69	1 (2%)
8	NAI	D	503	10	28,29,48	0.74	1 (3%)	43,45,73	0.74	1 (2%)
4	SO4	A	202	-	4,4,4	0.27	0	6,6,6	0.12	0
8	NAI	B	503	-	30,30,48	0.56	1 (3%)	43,46,73	0.59	1 (2%)
3	FES	C	201	1	0,4,4	-	-	-		
4	SO4	C	203	-	4,4,4	0.29	0	6,6,6	0.08	0
4	SO4	A	204	-	4,4,4	0.29	0	6,6,6	0.07	0
4	SO4	C	202	-	4,4,4	0.30	0	6,6,6	0.19	0
4	SO4	B	505	-	4,4,4	0.27	0	6,6,6	0.22	0
9	GOL	D	504	-	5,5,5	0.12	0	5,5,5	0.11	0
6	SF4	D	501	2	0,12,12	-	-	-		

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
7	FMN	D	502	10	-	3/18/18/18	0/3/3/3
6	SF4	B	501	2	-	-	0/6/5/5
9	GOL	B	504	-	-	0/4/4/4	-
8	NAI	D	503	10	-	1/16/32/72	0/3/3/5
7	FMN	B	502	-	-	2/18/18/18	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	FES	A	201	1	-	-	0/1/1/1
8	NAI	B	503	-	-	1/19/35/72	0/3/3/5
3	FES	C	201	1	-	-	0/1/1/1
9	GOL	D	504	-	-	0/4/4/4	-
6	SF4	D	501	2	-	-	0/6/5/5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	D	503	NAI	PN-O5D	3.06	1.66	1.54
8	B	503	NAI	PN-O5D	2.32	1.66	1.58

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	D	503	NAI	O5D-PN-O3	-3.18	93.96	104.64
7	D	502	FMN	O2P-P-O5'	-2.31	100.64	106.67
8	B	503	NAI	PN-O5D-C5D	2.26	134.03	121.91
7	B	502	FMN	C5'-C4'-C3'	-2.02	108.41	112.22

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	B	503	NAI	C5B-O5B-PA-O1A
8	D	503	NAI	C5B-O5B-PA-O1A
7	B	502	FMN	C4'-C5'-O5'-P
7	D	502	FMN	C4'-C5'-O5'-P
7	D	502	FMN	O2'-C2'-C3'-C4'

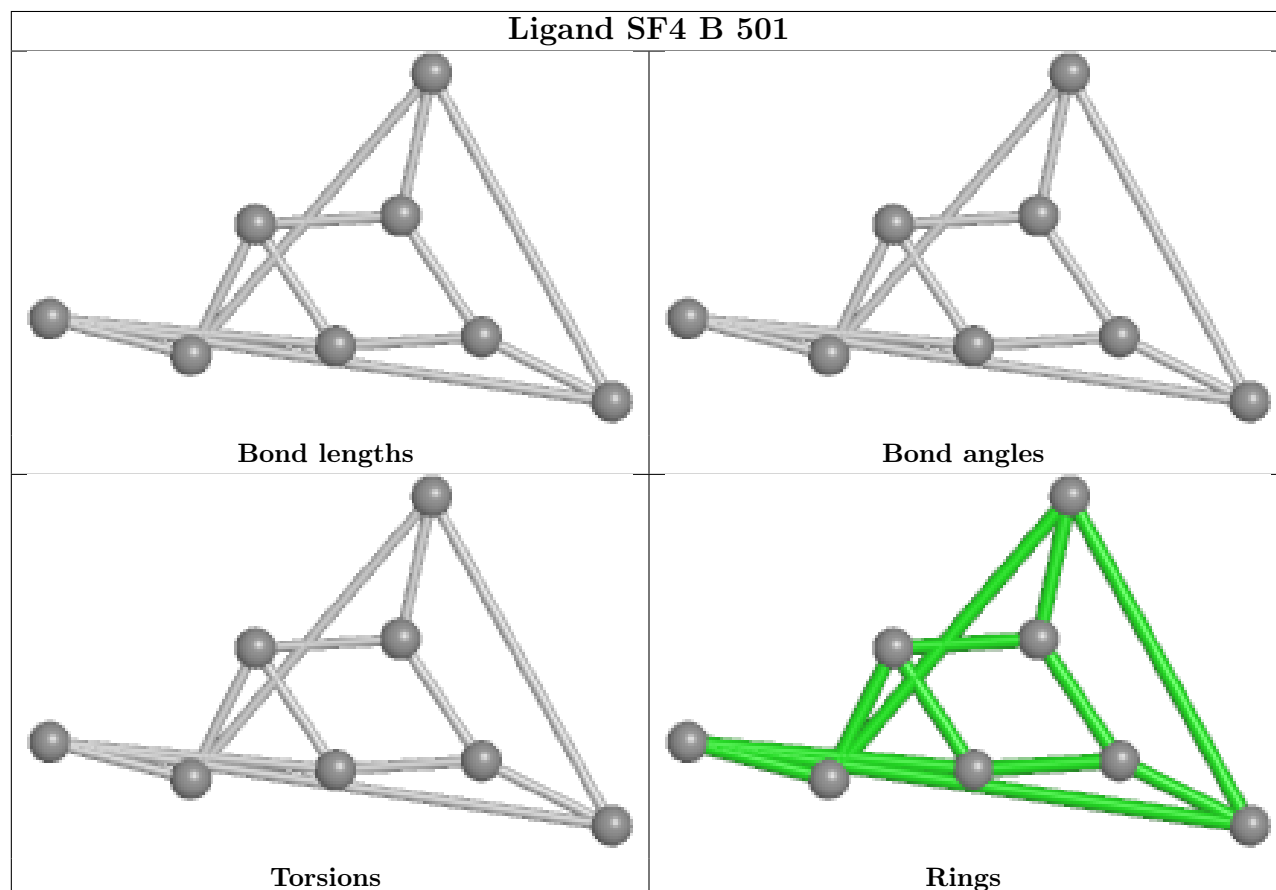
There are no ring outliers.

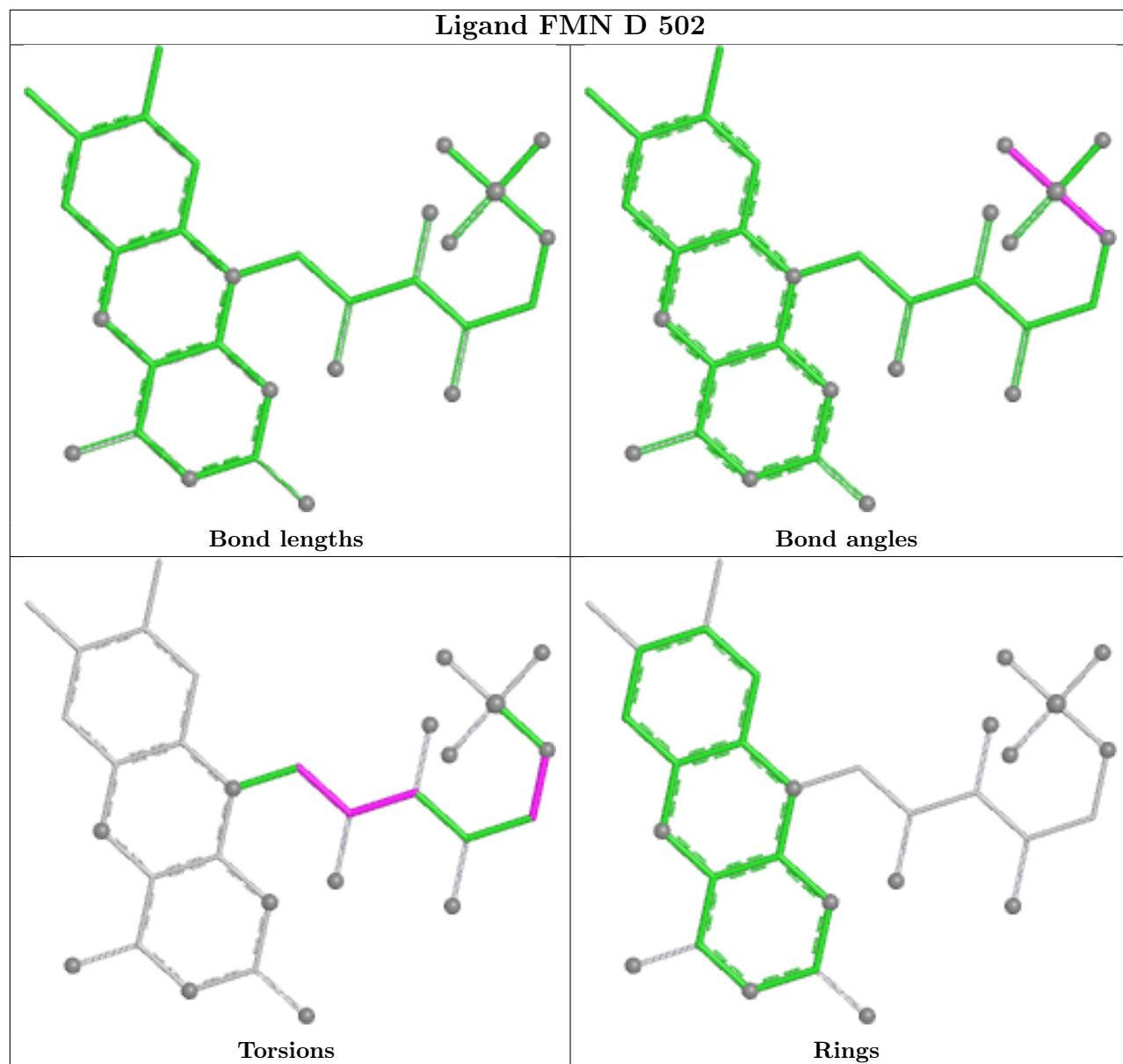
1 monomer is involved in 1 short contact:

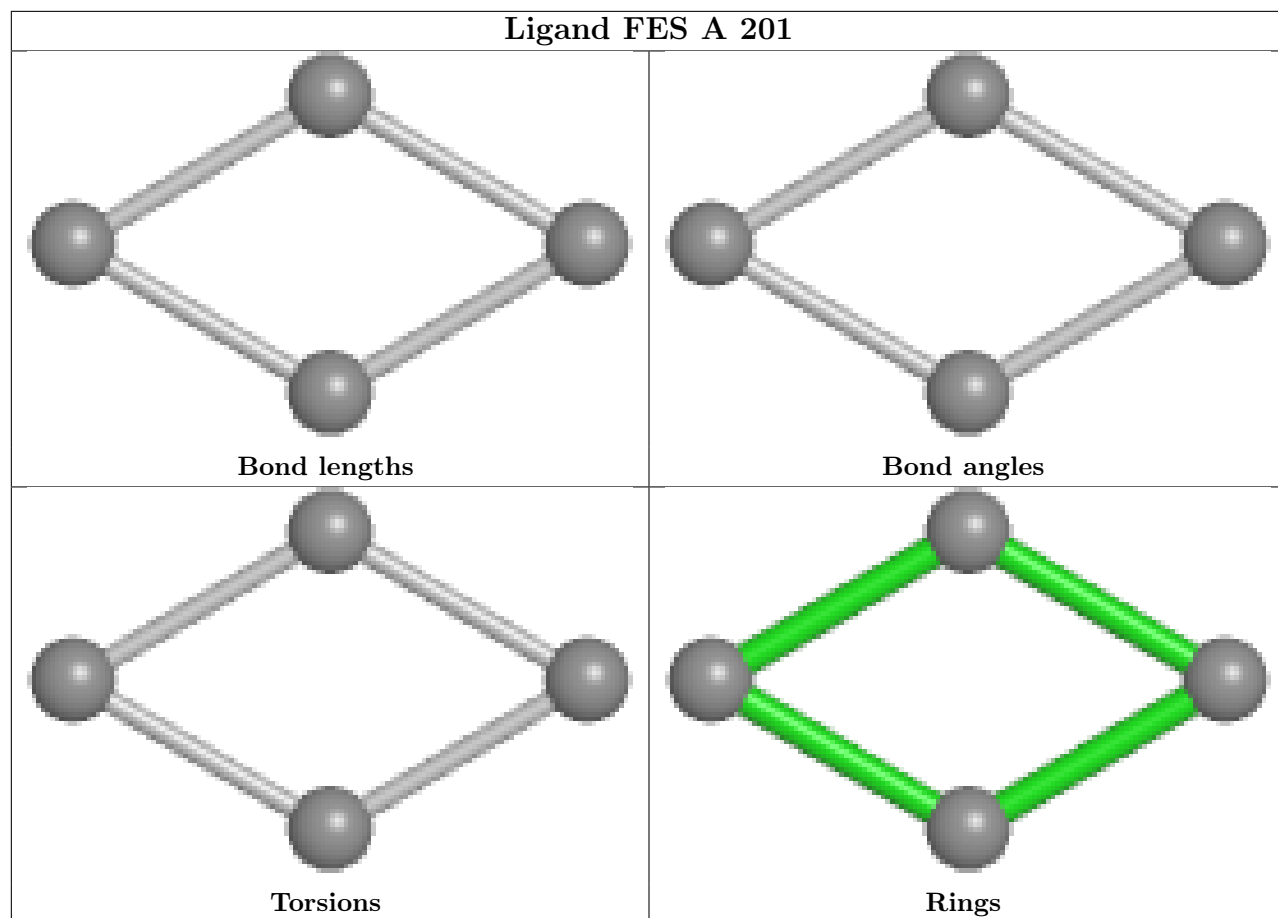
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	503	NAI	1	0

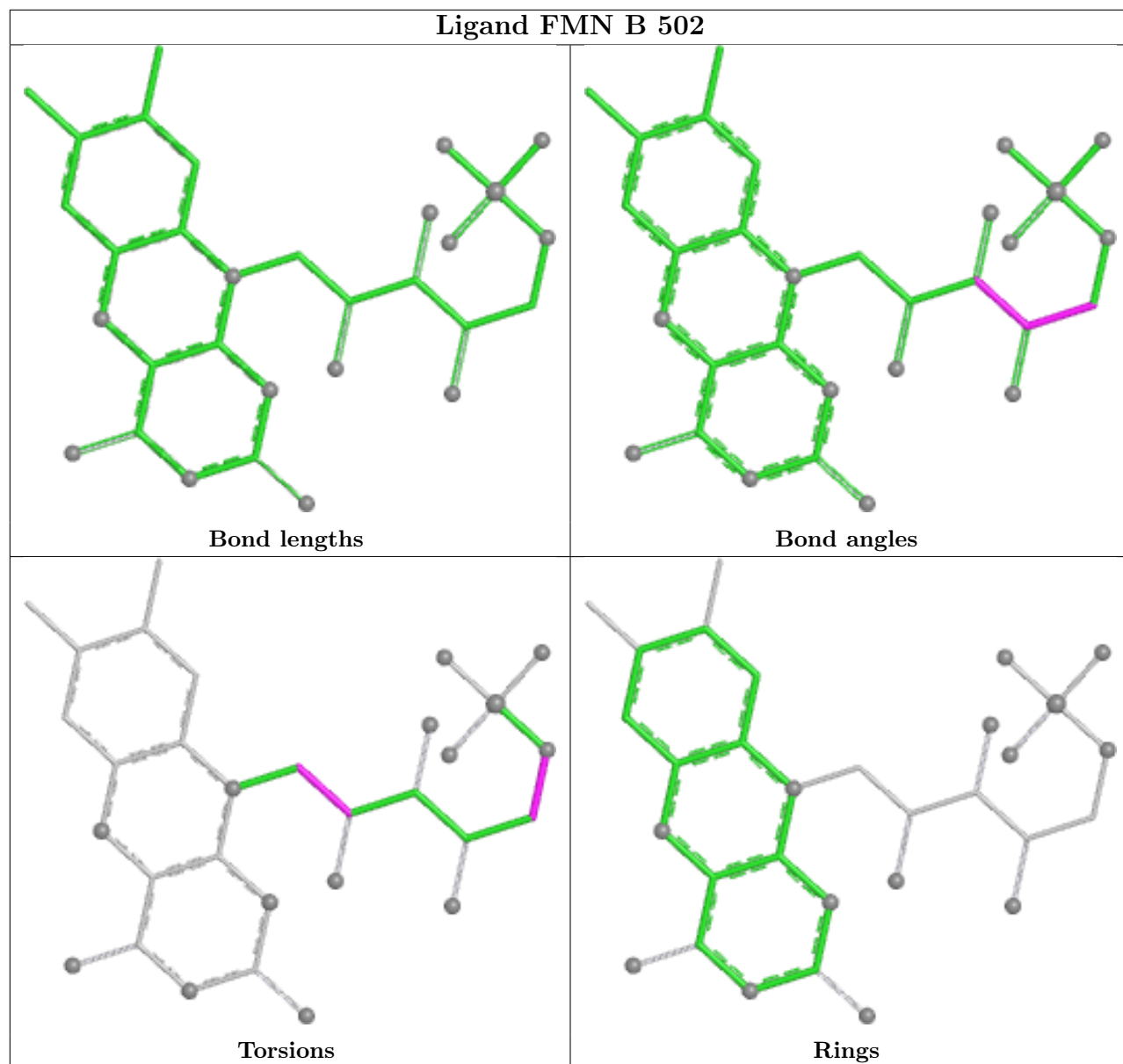
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.

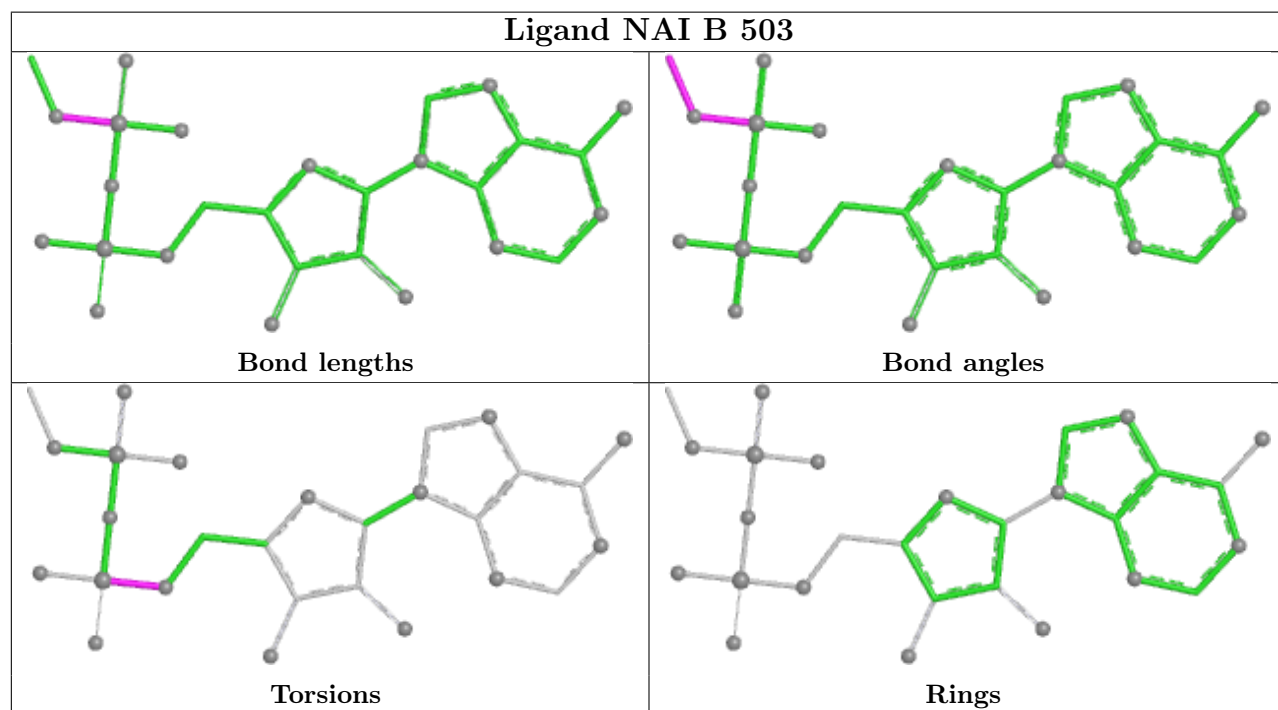
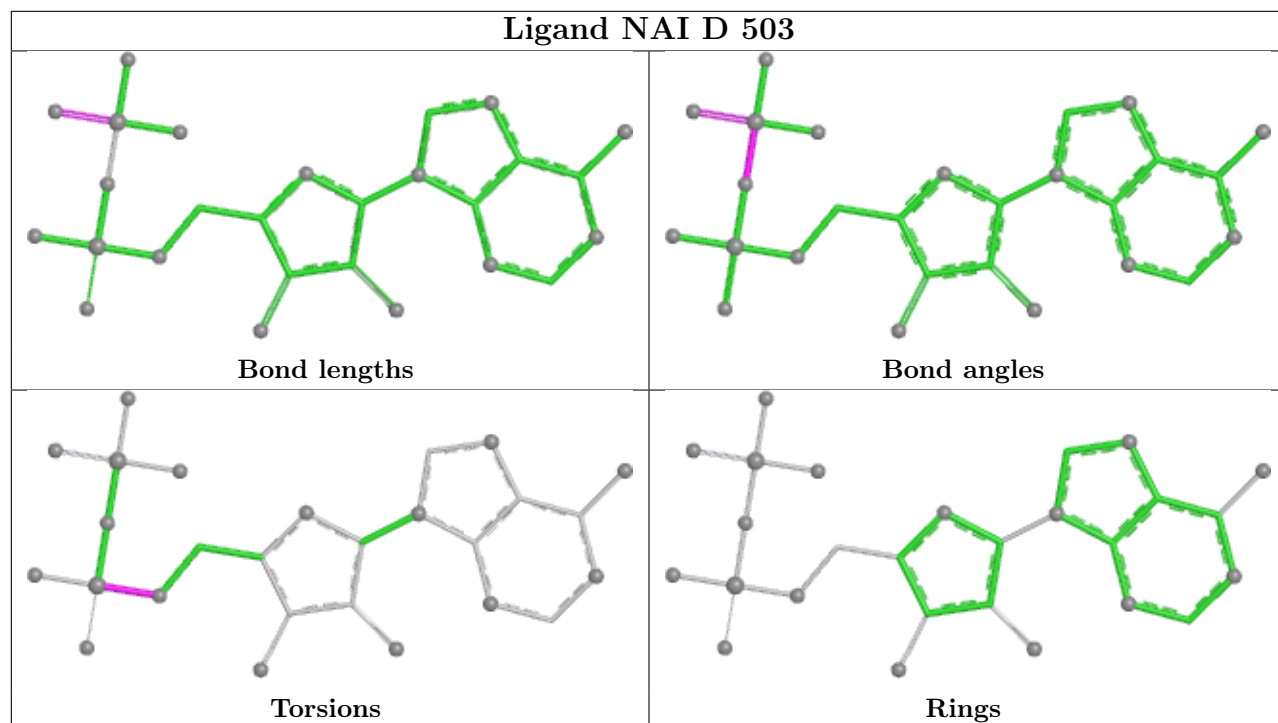
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD 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.

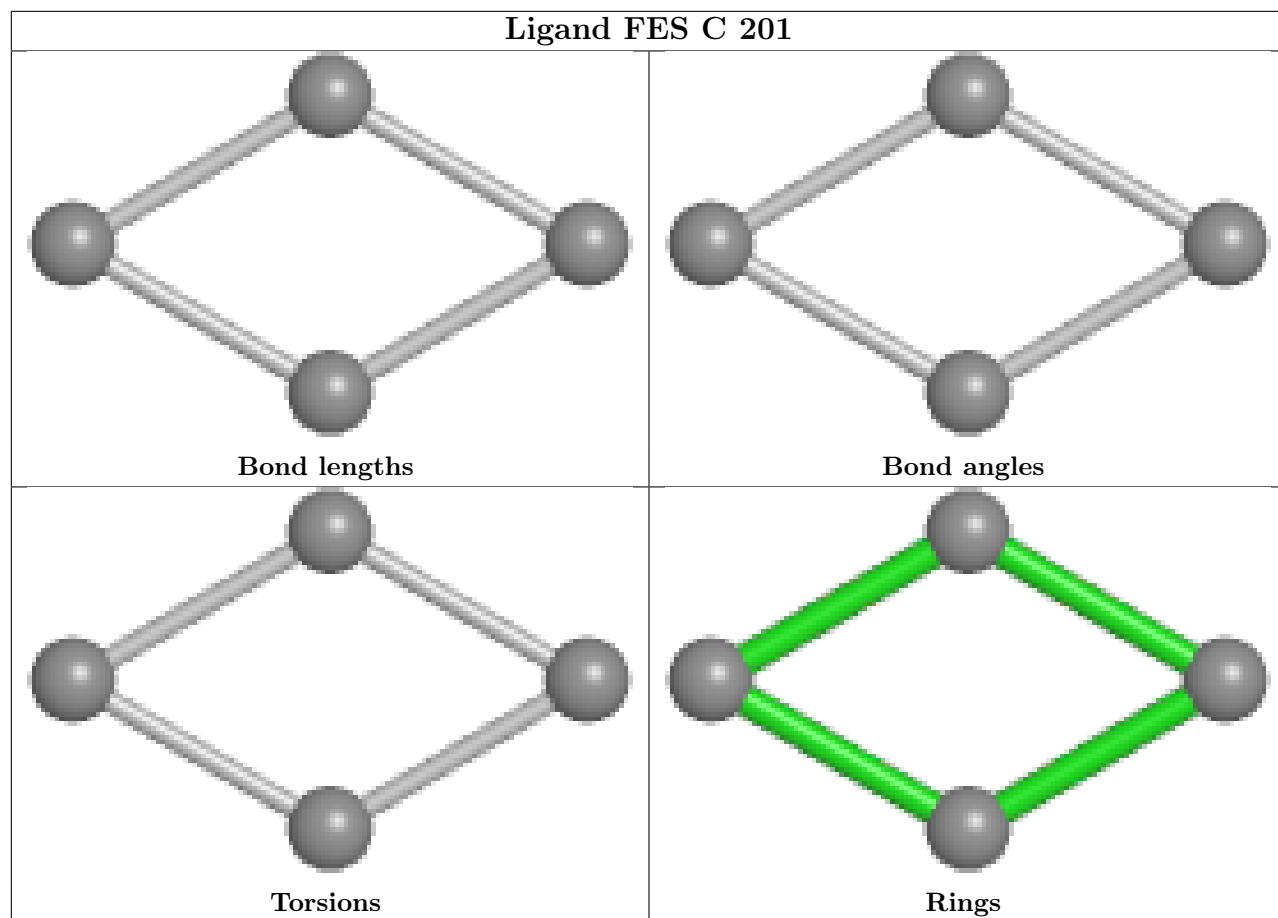


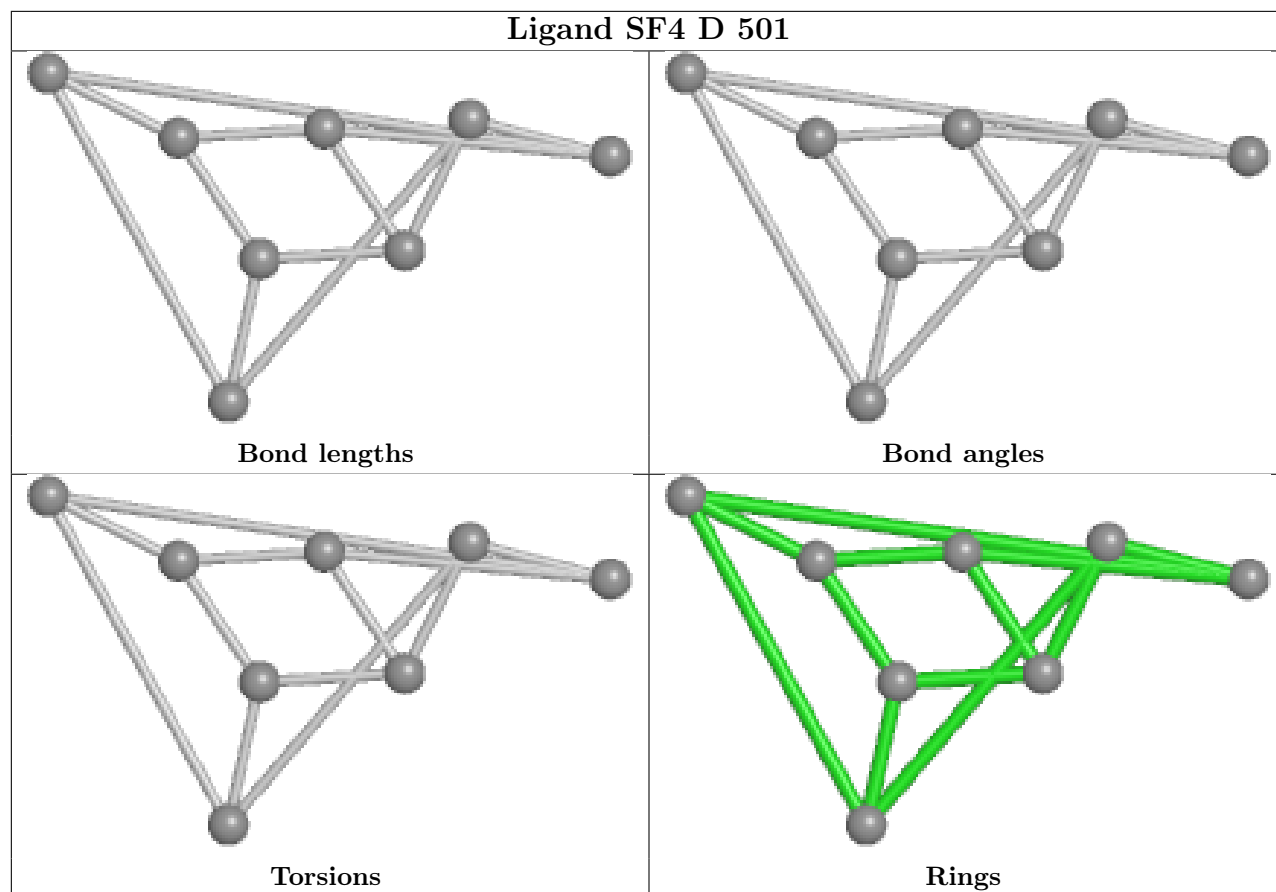












5.7 Other polymers [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	156/160 (97%)	0.00	0 100 100	21, 34, 53, 94	0
1	C	156/160 (97%)	0.01	1 (0%) 85 88	20, 33, 54, 84	0
2	B	416/434 (95%)	-0.16	7 (1%) 69 73	14, 27, 40, 67	1 (0%)
2	D	418/434 (96%)	-0.23	5 (1%) 76 80	11, 26, 40, 83	3 (0%)
All	All	1146/1188 (96%)	-0.14	13 (1%) 78 81	11, 28, 47, 94	4 (0%)

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	1	MET	6.3
2	B	377	TRP	4.0
2	B	418	ARG	3.7
2	D	377	TRP	2.7
2	D	2	ARG	2.6

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](#)

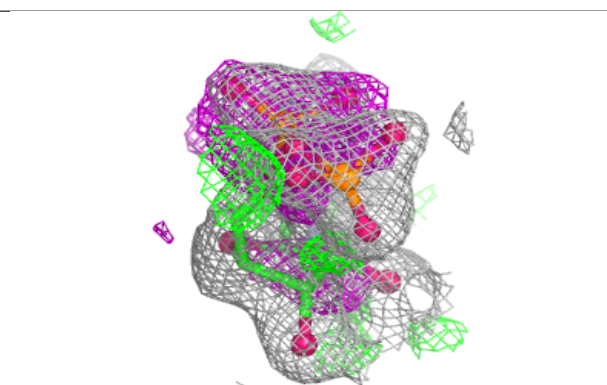
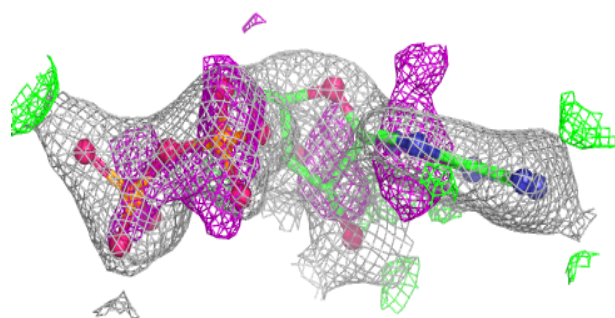
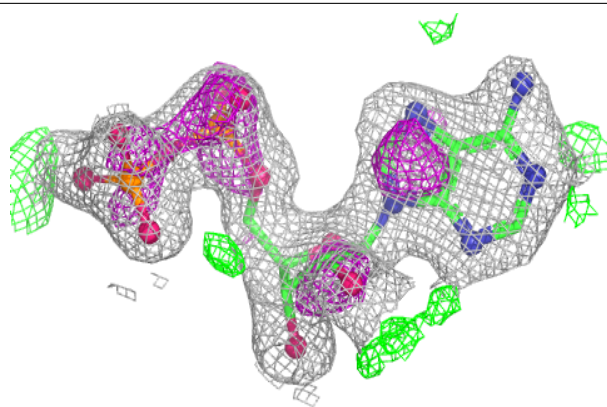
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	SO4	D	506	5/5	0.66	0.12	78,79,81,87	0
4	SO4	A	202	5/5	0.75	0.12	73,80,82,83	0
10	NA	D	510	1/1	0.84	0.17	47,47,47,47	0
4	SO4	B	505	5/5	0.85	0.12	30,47,51,55	0
5	CL	A	205	1/1	0.86	0.13	68,68,68,68	0
4	SO4	C	202	5/5	0.86	0.09	55,59,60,62	0
9	GOL	D	504	6/6	0.88	0.11	31,36,37,38	0
10	NA	B	508	1/1	0.88	0.27	53,53,53,53	0
5	CL	C	205	1/1	0.88	0.11	68,68,68,68	0
4	SO4	C	203	5/5	0.89	0.10	47,50,56,61	0
4	SO4	A	203	5/5	0.89	0.10	46,52,61,62	0
9	GOL	B	504	6/6	0.89	0.11	30,35,37,38	0
10	NA	D	511	1/1	0.89	0.16	56,56,56,56	0
5	CL	C	204	1/1	0.90	0.18	66,66,66,66	0
4	SO4	D	505	5/5	0.91	0.10	33,43,46,47	0
10	NA	D	509	1/1	0.91	0.24	50,50,50,50	0
4	SO4	A	204	5/5	0.92	0.07	53,54,63,66	0
8	NAI	D	503	27/44	0.92	0.10	28,40,46,47	0
5	CL	B	506	1/1	0.93	0.26	59,59,59,59	0
8	NAI	B	503	28/44	0.95	0.07	27,35,40,45	0
10	NA	B	509	1/1	0.95	0.10	56,56,56,56	0
10	NA	D	508	1/1	0.95	0.15	36,36,36,36	0
10	NA	B	507	1/1	0.98	0.06	32,32,32,32	0
10	NA	D	507	1/1	0.98	0.08	32,32,32,32	0
7	FMN	B	502	31/31	0.98	0.04	19,21,23,23	0
7	FMN	D	502	31/31	0.99	0.04	18,20,23,23	0
6	SF4	B	501	8/8	1.00	0.02	21,22,22,22	0
6	SF4	D	501	8/8	1.00	0.02	20,20,21,21	0
3	FES	A	201	4/4	1.00	0.02	21,21,21,21	0
3	FES	C	201	4/4	1.00	0.01	20,20,20,20	0

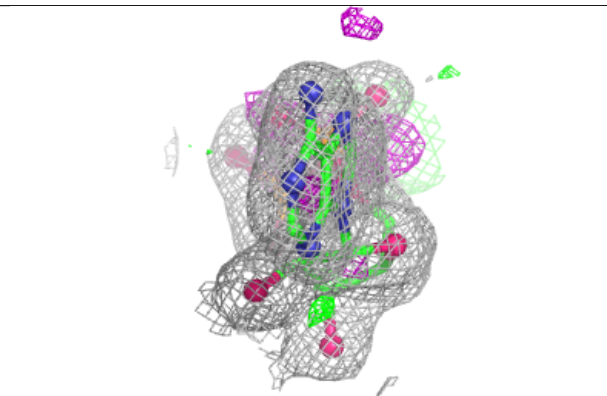
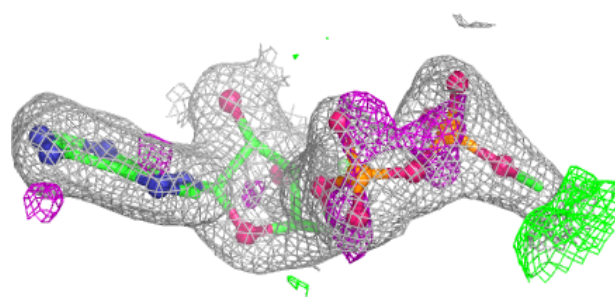
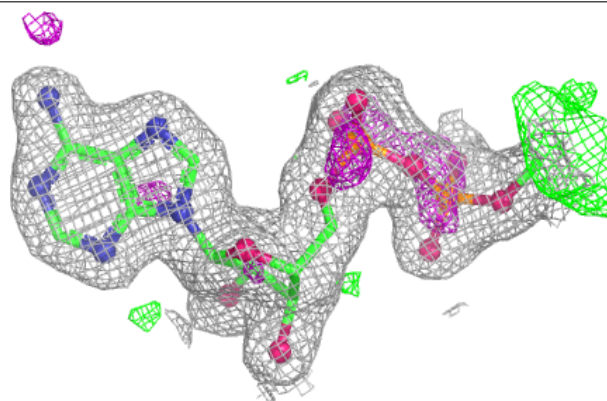
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAI D 503:

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

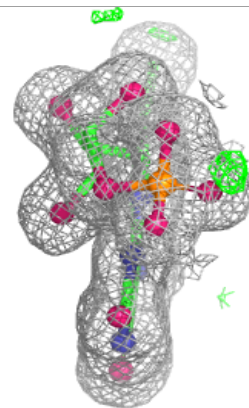
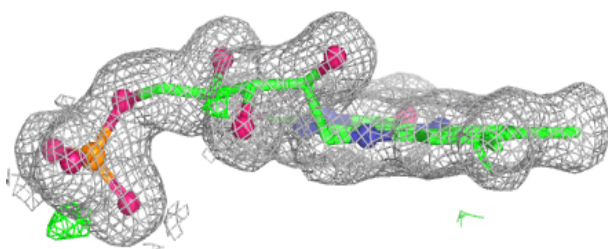
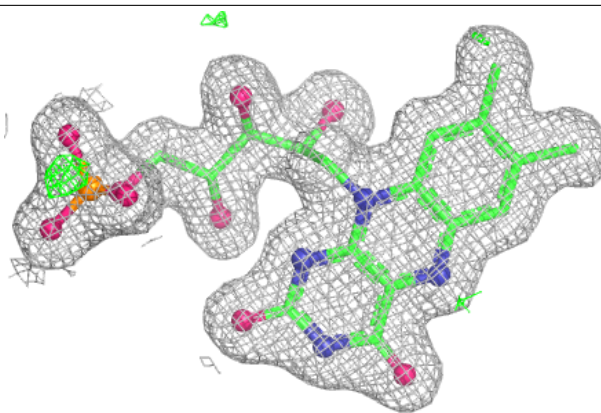
**Electron density around NAI B 503:**

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

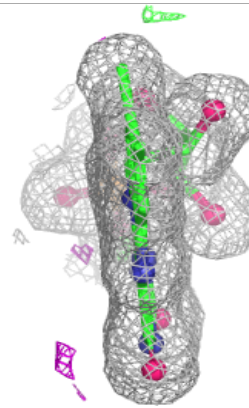
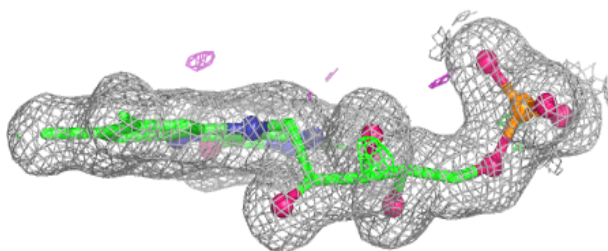
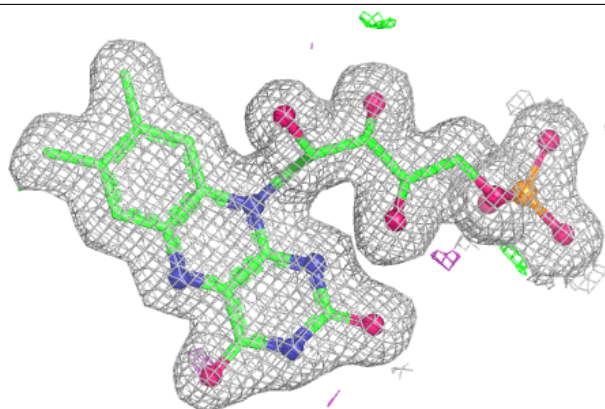


Electron density around FMN B 502:

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

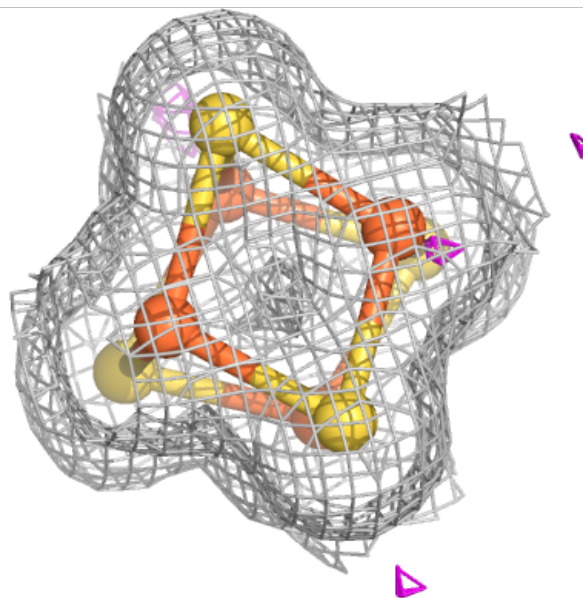
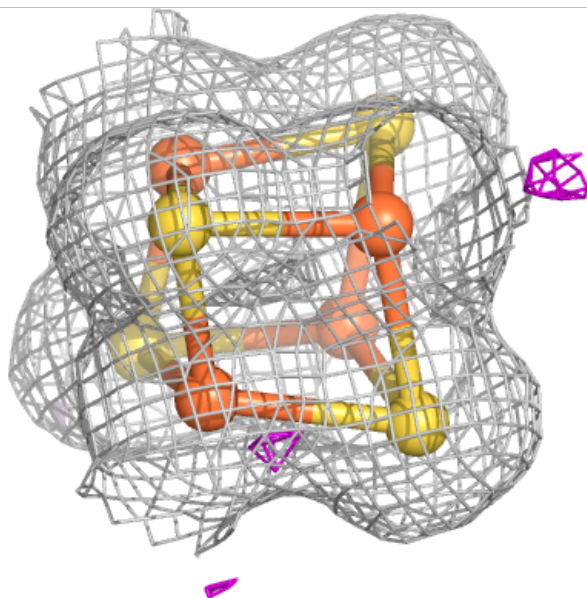
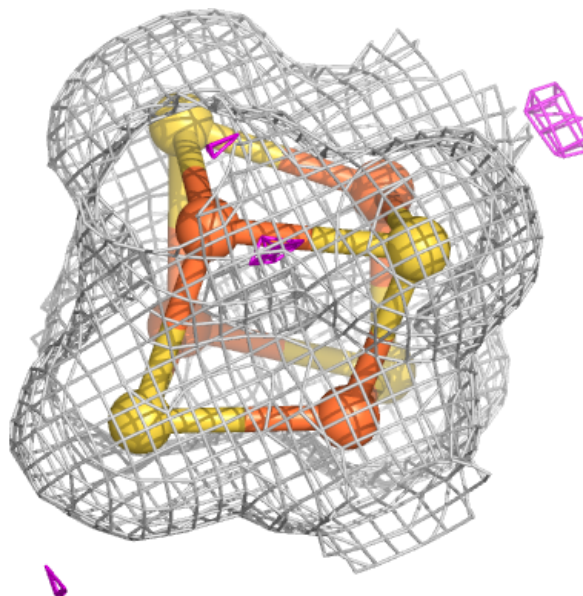
**Electron density around FMN D 502:**

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



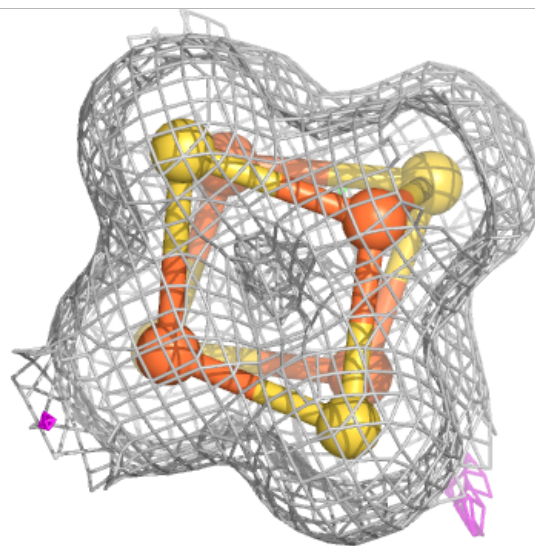
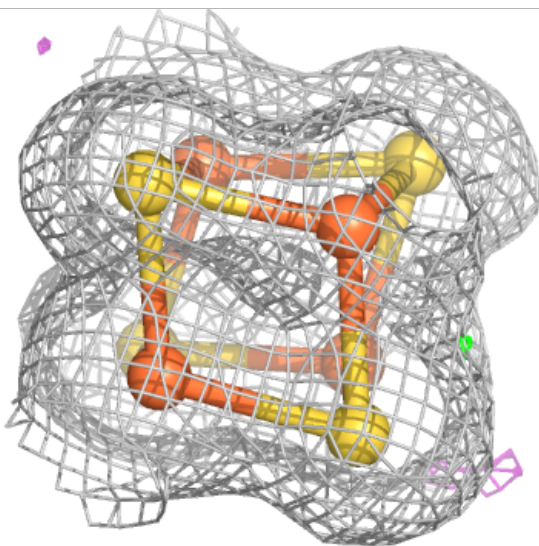
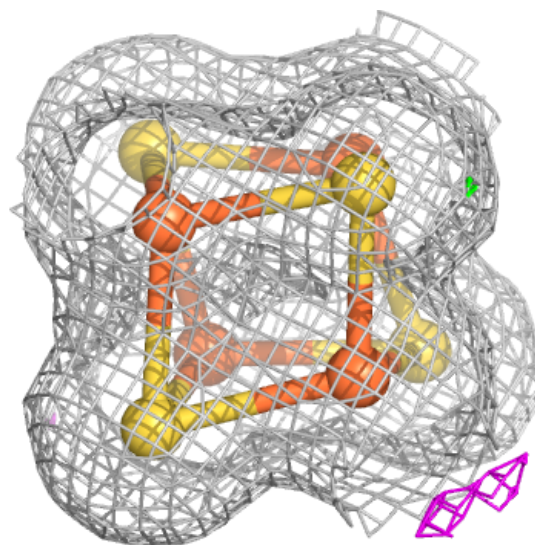
Electron density around SF4 B 501:

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



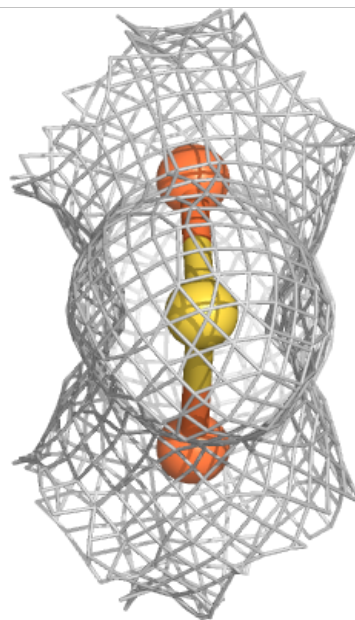
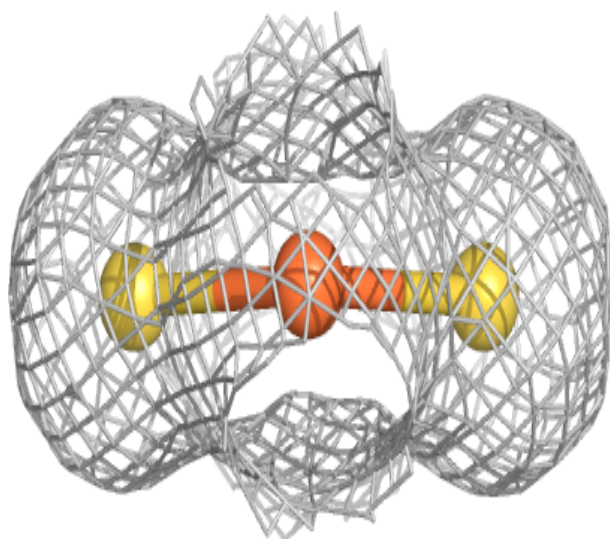
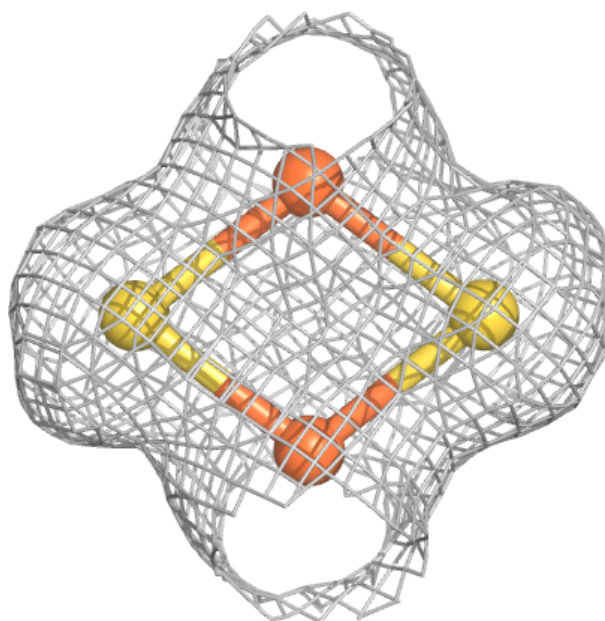
Electron density around SF4 D 501:

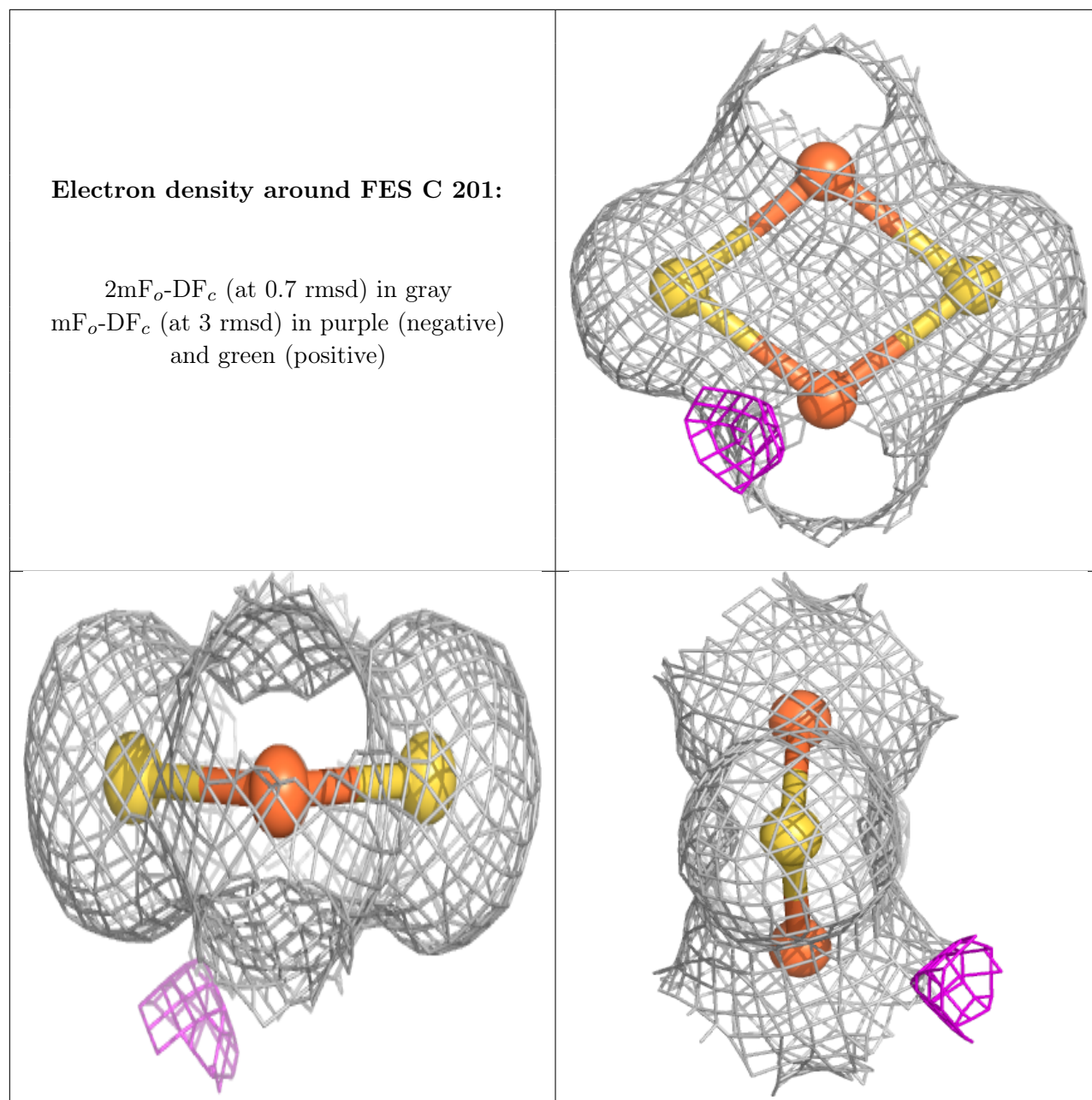
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FES A 201:

$2mF_o-DF_c$ (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.