wwPDB X-ray Structure Validation Summary Report (i)

Aug 7, 2026 – 07:59 AM JST

PDB ID	:	8JC1 / pdb_00008jc1
Title	:	Crystal structure of Pectocin M1 from Pectobacterium carotovorum
Authors	:	Jantarit, N.; Kurisu, G.; Tanaka, H.
Deposited on	:	2023-05-10
Resolution	:	2.04 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

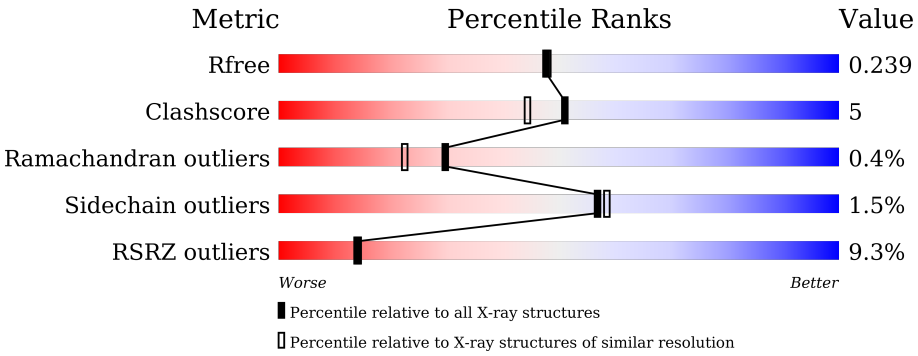
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.04 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	268	13% 81% 18%
1	B	268	13% 87% 12%
1	C	268	6% 85% 14%
1	D	268	4% 89% 10%

2 Entry composition (i)

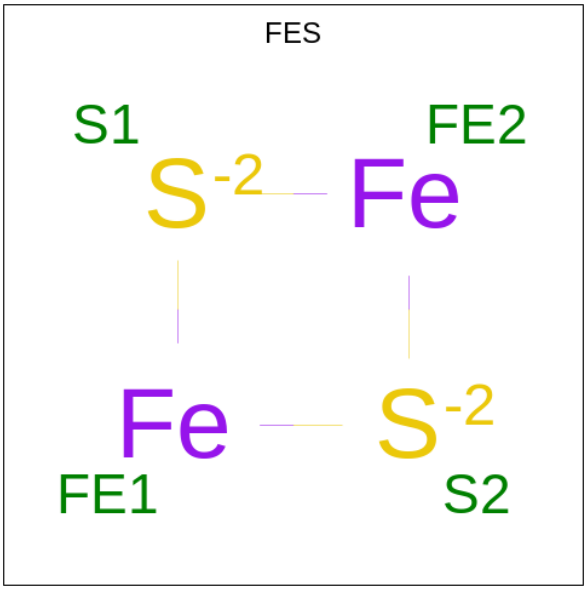
There are 9 unique types of molecules in this entry. The entry contains 8765 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 pectocin M1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			2037	1285	334	409	9			
1	B	267	Total	C	N	O	S	0	0	0
			2042	1287	336	410	9			
1	C	267	Total	C	N	O	S	0	0	0
			2044	1288	335	412	9			
1	D	267	Total	C	N	O	S	0	0	0
			2035	1284	333	409	9			

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



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			

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

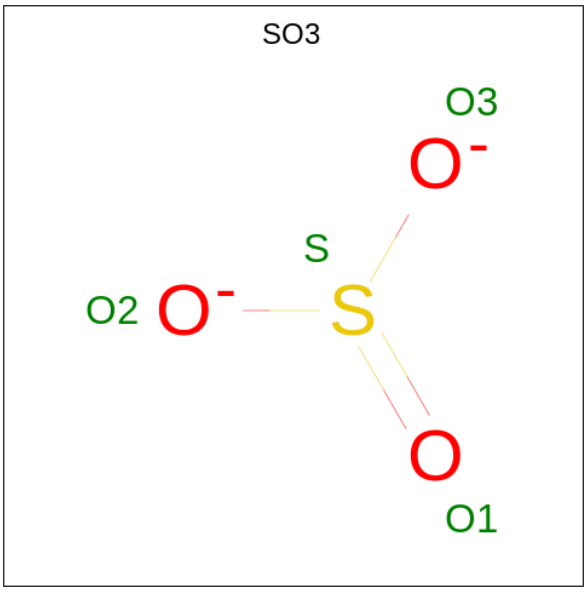
- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	4	Total Na 4 4	0	0
3	B	2	Total Na 2 2	0	0
3	C	1	Total Na 1 1	0	0
3	D	1	Total Na 1 1	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

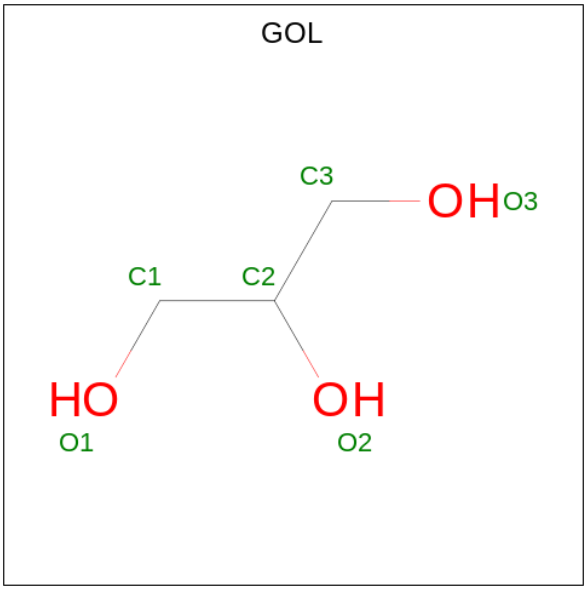
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mg 1 1	0	0
4	B	1	Total Mg 1 1	0	0
4	C	1	Total Mg 1 1	0	0
4	D	1	Total Mg 1 1	0	0

- Molecule 5 is SULFITE ION (CCD ID: SO3) (formula: O_3S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total O S 4 3 1	0	0
5	C	1	Total O S 4 3 1	0	0

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



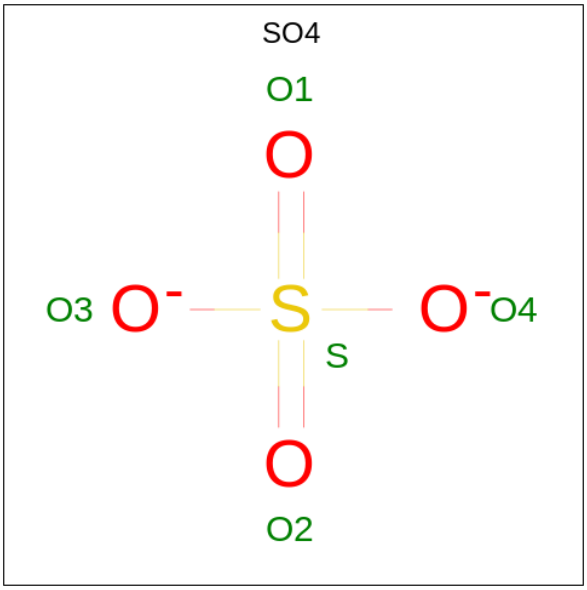
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total C O 6 3 3	0	0
6	C	1	Total C O 6 3 3	0	0

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

- Molecule 7 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

- Molecule 8 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	C	1	Total Cl 1 1	0	0
8	D	1	Total Cl 1 1	0	0

- Molecule 9 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	135	Total O 135 135	0	0
9	B	116	Total O 116 116	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	C	137	Total O 137 137	0	0
9	D	152	Total O 152 152	0	0

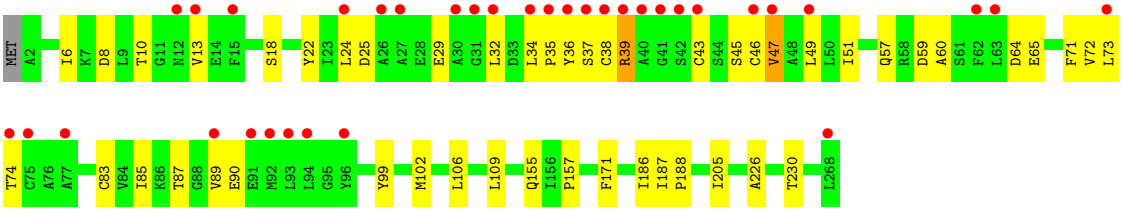
3 Residue-property plots (i)

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: pectocin M1

Chain A:

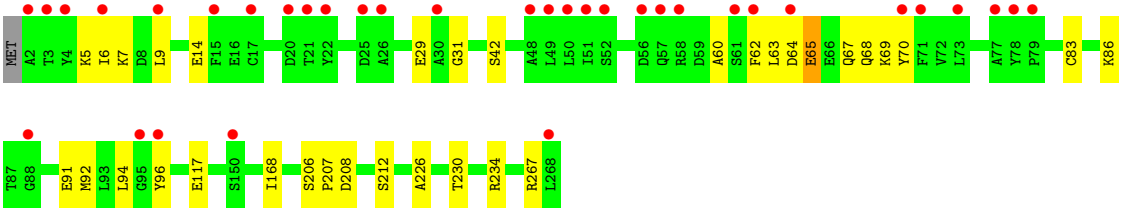




• Molecule 1: pectocin M1

Chain B:

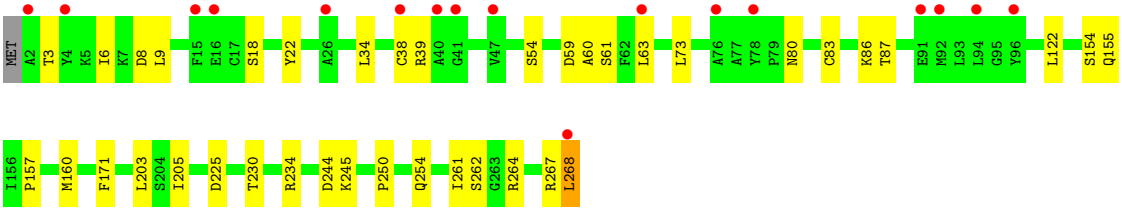




• Molecule 1: pectocin M1

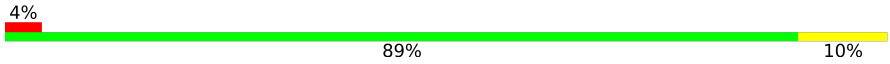
Chain C:

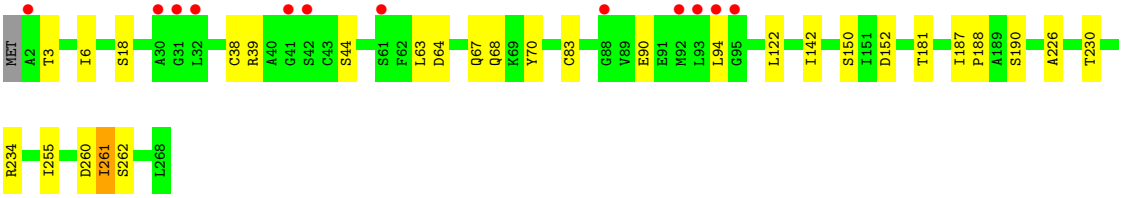




• Molecule 1: pectocin M1

Chain D:




MET
A2
T3
I6
S18
A30
G31
L32
C38
R39
A40
G41
S42
C43
S44
S61
F62
L63
D64
Q67
Q68
K69
Y70
C83
G88
V89
E90
E91
M92
L93
L94
G95
L122
I142
S150
I151
D152
T181
I187
P188
A189
S190
A226
T230
R234
I255
D260
I261
S262
L268

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	66.28Å 138.55Å 68.95Å 90.00° 94.18° 90.00°	Depositor
Resolution (Å)	49.49 – 2.04 49.49 – 2.04	Depositor EDS
% Data completeness (in resolution range)	98.6 (49.49-2.04) 98.6 (49.49-2.04)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.68 (at 2.05Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.195 , 0.238 0.195 , 0.239	Depositor DCC
R_{free} test set	1976 reflections (2.52%)	wwPDB-VP
Wilson B-factor (Å^2)	38.9	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 40.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.025 for l,-k,h	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8765	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% of the height of the origin peak. No significant pseudotranslation is detected.*

[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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, CL, SO3, NA, MG, 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	0.32	0/2078	0.47	0/2823
1	B	0.32	0/2083	0.50	0/2830
1	C	0.32	0/2085	0.48	0/2832
1	D	0.33	0/2076	0.50	0/2821
All	All	0.32	0/8322	0.49	0/11306

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	2037	0	1939	29	0
1	B	2042	0	1949	20	1
1	C	2044	0	1949	22	0
1	D	2035	0	1938	18	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
2	D	4	0	0	0	0
3	A	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	B	4	0	0	0	0
5	C	4	0	0	0	0
6	B	6	0	8	1	0
6	C	18	0	24	0	0
7	C	5	0	0	0	0
8	C	1	0	0	0	0
8	D	1	0	0	0	0
9	A	135	0	0	0	0
9	B	116	0	0	0	1
9	C	137	0	0	1	1
9	D	152	0	0	0	0
All	All	8765	0	7807	83	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

The worst 5 of 83 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:122:LEU:HD23	1:D:187:ILE:HD11	1.52	0.90
1:C:267:ARG:HG2	1:C:268:LEU:HD23	1.66	0.77
1:A:60:ALA:HB1	1:A:73:LEU:HD12	1.66	0.76
1:D:190:SER:HA	1:D:234:ARG:HG2	1.69	0.73
1:C:60:ALA:HB1	1:C:73:LEU:HD23	1.74	0.69

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:GLU:OE2	1:B:212:SER:OG[1_455]	2.03	0.17
9:B:452:HOH:O	9:C:527:HOH:O[1_654]	2.09	0.11

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	265/268 (99%)	257 (97%)	6 (2%)	2 (1%)	16 9
1	B	265/268 (99%)	248 (94%)	15 (6%)	2 (1%)	16 9
1	C	265/268 (99%)	253 (96%)	12 (4%)	0	100 100
1	D	265/268 (99%)	255 (96%)	10 (4%)	0	100 100
All	All	1060/1072 (99%)	1013 (96%)	43 (4%)	4 (0%)	30 22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	39	ARG
1	B	64	ASP
1	B	65	GLU
1	A	90	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	220/227 (97%)	216 (98%)	4 (2%)	51 51
1	B	222/227 (98%)	220 (99%)	2 (1%)	70 75
1	C	222/227 (98%)	218 (98%)	4 (2%)	51 51
1	D	221/227 (97%)	218 (99%)	3 (1%)	59 60
All	All	885/908 (98%)	872 (98%)	13 (2%)	57 58

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	261	ILE
1	C	262	SER
1	D	261	ILE
1	D	150	SER
1	D	152	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	67	GLN
1	D	214	ASN
1	B	161	ASN
1	B	254	GLN
1	C	80	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 oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 25 ligands modelled in this entry, 14 are monoatomic - leaving 11 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	SO3	B	302	-	1,3,3	0.67	0	0,3,3	-	-
5	SO3	C	306	-	1,3,3	0.69	0	0,3,3	-	-
6	GOL	C	301	3	5,5,5	0.86	0	5,5,5	1.23	0
7	SO4	C	305	-	4,4,4	0.35	0	6,6,6	0.41	0
2	FES	D	301	1	0,4,4	-	-	-		
2	FES	C	302	1	0,4,4	-	-	-		
2	FES	A	301	1	0,4,4	-	-	-		
6	GOL	C	303	-	5,5,5	0.89	0	5,5,5	0.96	0
2	FES	B	301	1	0,4,4	-	-	-		
6	GOL	B	303	-	5,5,5	0.79	0	5,5,5	1.03	0
6	GOL	C	304	-	5,5,5	1.14	0	5,5,5	0.90	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
6	GOL	C	301	3	-	2/4/4/4	-
2	FES	C	302	1	-	-	0/1/1/1
2	FES	A	301	1	-	-	0/1/1/1
2	FES	D	301	1	-	-	0/1/1/1
6	GOL	C	303	-	-	3/4/4/4	-
2	FES	B	301	1	-	-	0/1/1/1
6	GOL	B	303	-	-	4/4/4/4	-
6	GOL	C	304	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	303	GOL	O1-C1-C2-C3
6	C	303	GOL	C1-C2-C3-O3
6	C	303	GOL	O2-C2-C3-O3
6	C	304	GOL	C1-C2-C3-O3
6	C	301	GOL	O1-C1-C2-C3

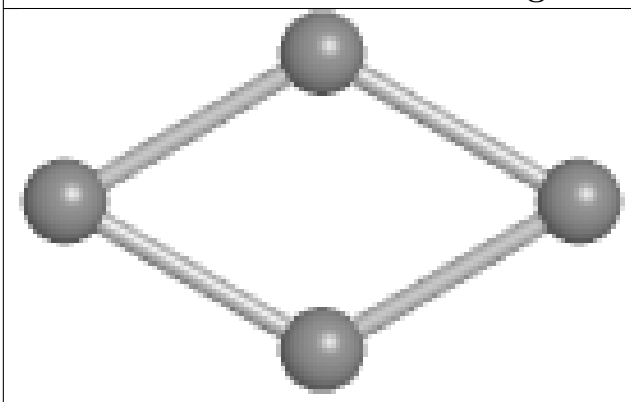
There are no ring outliers.

1 monomer is involved in 1 short contact:

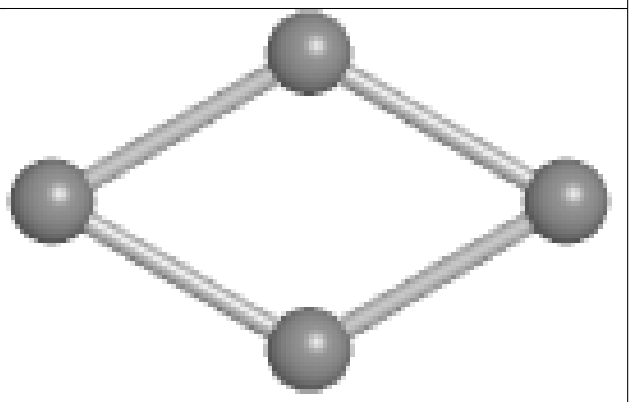
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	303	GOL	1	0

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.

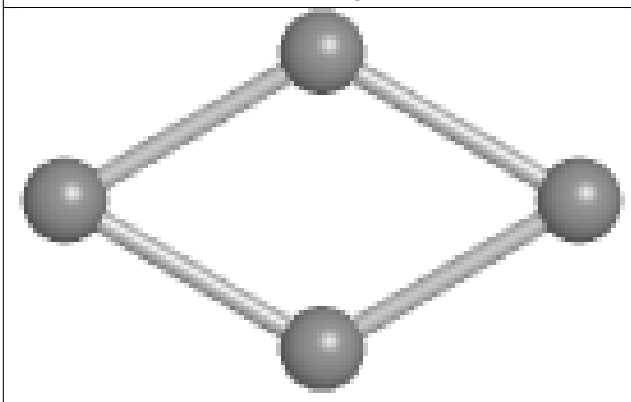
Ligand FES D 301



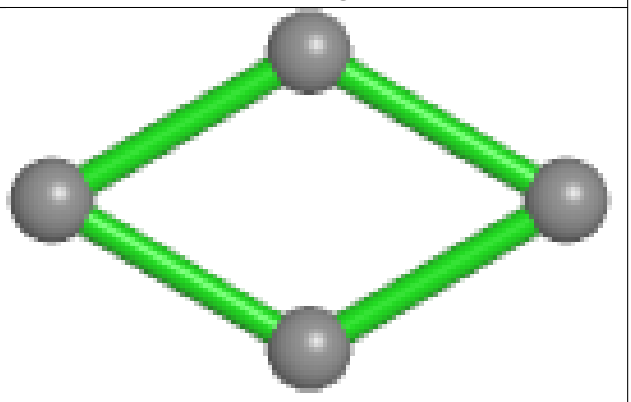
Bond lengths



Bond angles

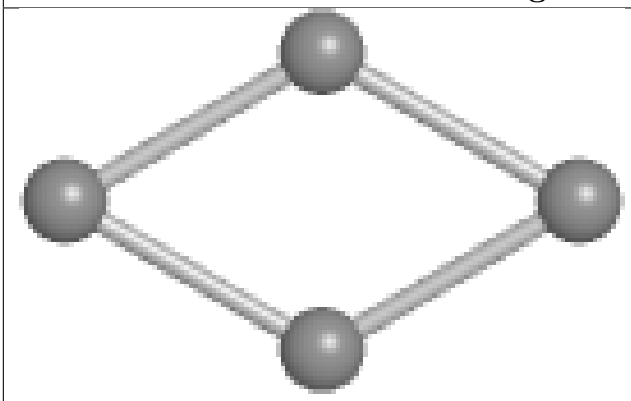


Torsions

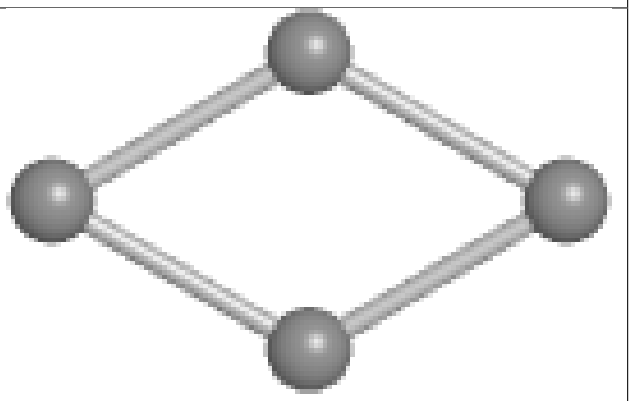


Rings

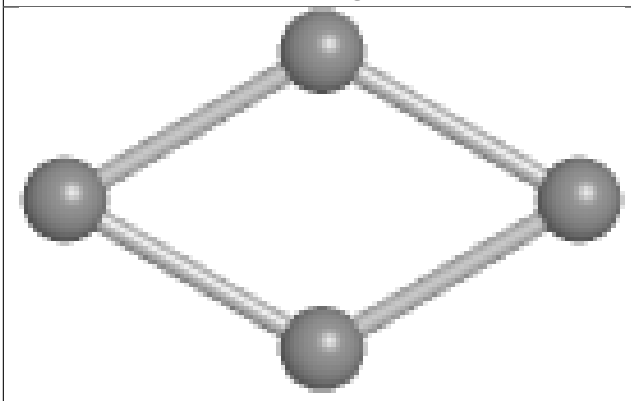
Ligand FES C 302



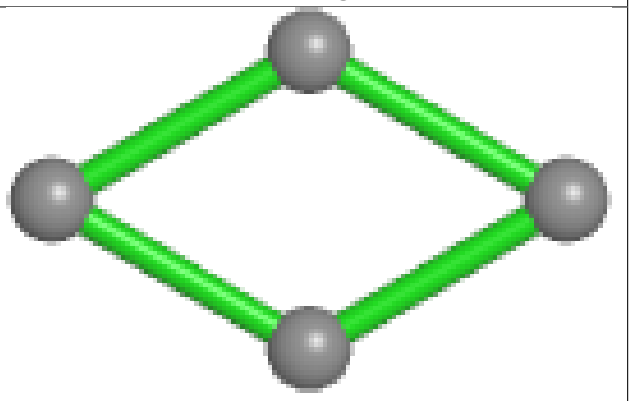
Bond lengths



Bond angles

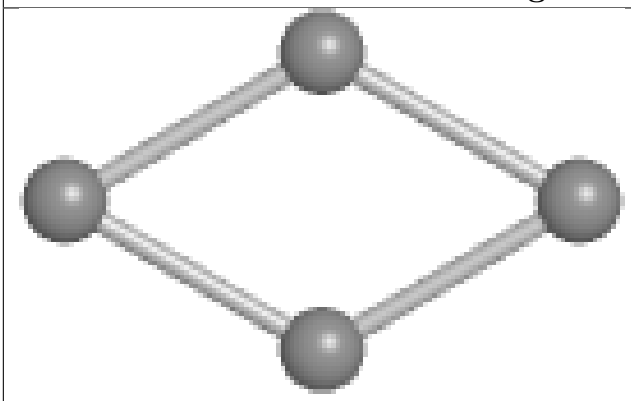


Torsions

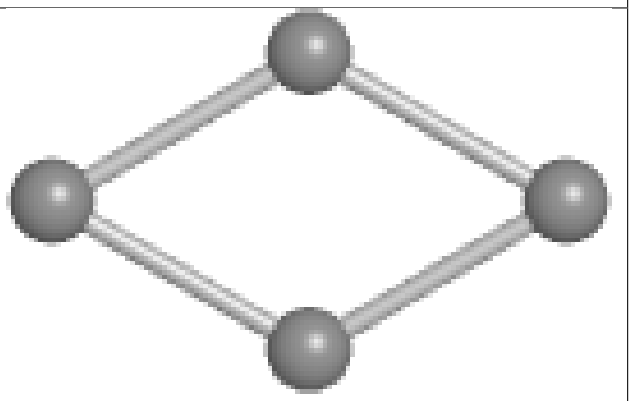


Rings

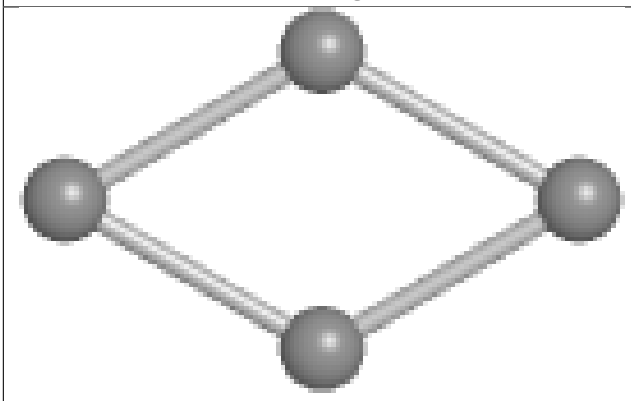
Ligand FES A 301



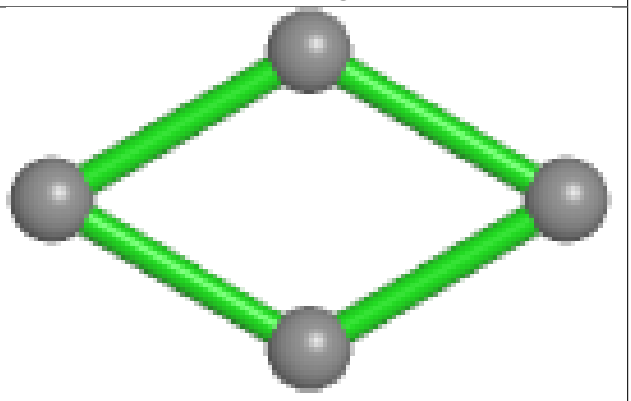
Bond lengths



Bond angles

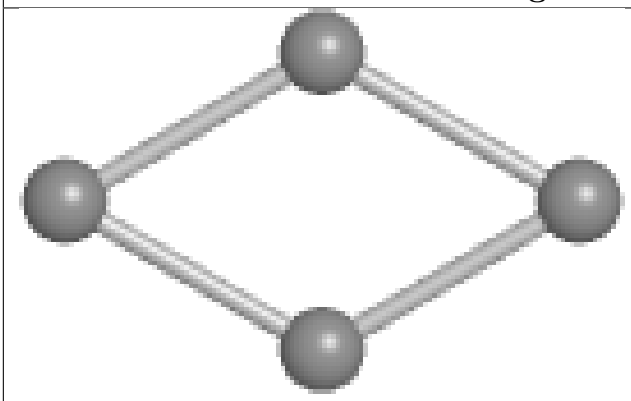


Torsions

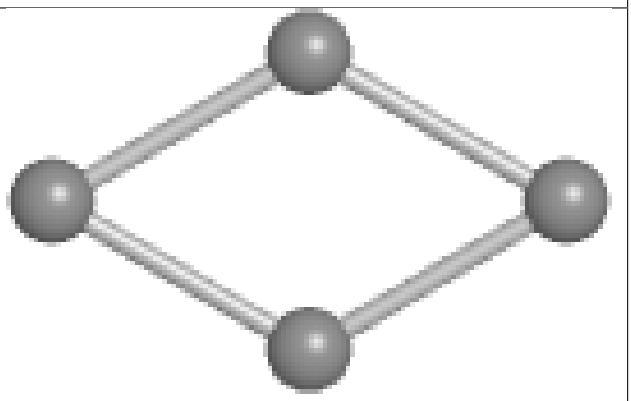


Rings

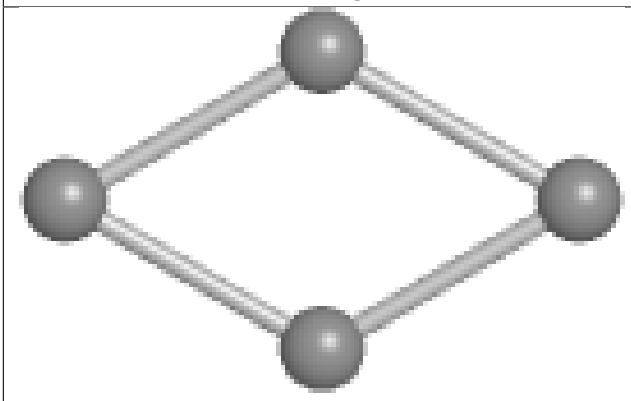
Ligand FES B 301



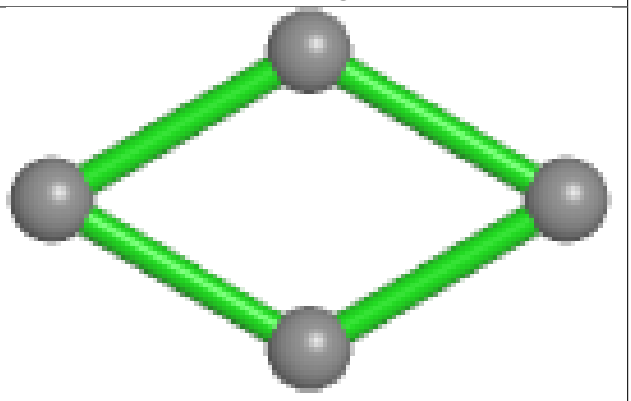
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{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	267/268 (99%)	0.56	35 (13%)	7	6	28, 43, 84, 92	0
1	B	267/268 (99%)	0.62	35 (13%)	7	6	30, 46, 77, 89	0
1	C	267/268 (99%)	0.37	17 (6%)	25	25	26, 40, 74, 79	0
1	D	267/268 (99%)	0.22	12 (4%)	38	38	26, 38, 68, 76	0
All	All	1068/1072 (99%)	0.44	99 (9%)	14	14	26, 42, 76, 92	0

The worst 5 of 99 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	89	VAL	4.6
1	A	74	THR	4.0
1	A	41	GLY	3.9
1	B	268	LEU	3.5
1	A	13	VAL	3.5

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

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

6.3 Carbohydrates

There are no oligosaccharides in this entry.

6.4 Ligands

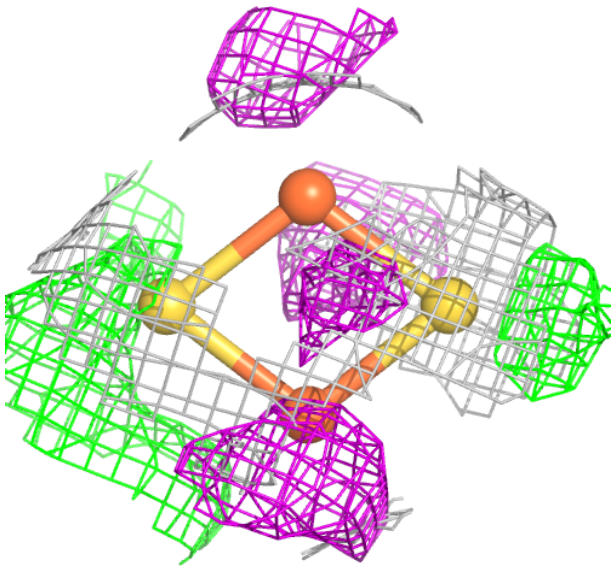
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

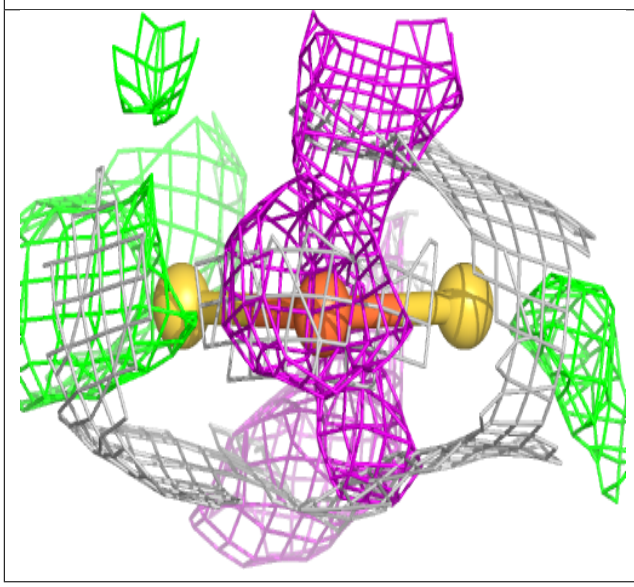
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
6	GOL	B	303	6/6	0.61	0.15	60,67,75,81	0
3	NA	B	304	1/1	0.77	0.18	74,74,74,74	0
6	GOL	C	303	6/6	0.79	0.15	73,75,76,78	0
3	NA	A	302	1/1	0.82	0.22	56,56,56,56	0
2	FES	A	301	4/4	0.82	0.12	78,90,96,97	0
6	GOL	C	304	6/6	0.82	0.15	50,53,57,57	0
3	NA	B	305	1/1	0.85	0.16	60,60,60,60	0
8	CL	D	303	1/1	0.85	0.12	70,70,70,70	0
3	NA	D	302	1/1	0.87	0.14	63,63,63,63	0
5	SO3	C	306	4/4	0.87	0.11	55,56,68,82	0
3	NA	A	305	1/1	0.89	0.11	45,45,45,45	0
6	GOL	C	301	6/6	0.90	0.15	50,54,56,57	0
7	SO4	C	305	5/5	0.90	0.17	41,43,54,73	0
3	NA	A	303	1/1	0.90	0.13	56,56,56,56	0
8	CL	C	308	1/1	0.93	0.10	71,71,71,71	0
4	MG	A	306	1/1	0.94	0.06	42,42,42,42	0
5	SO3	B	302	4/4	0.95	0.11	56,57,58,71	0
4	MG	D	304	1/1	0.96	0.06	39,39,39,39	0
3	NA	A	304	1/1	0.96	0.08	52,52,52,52	0
4	MG	B	306	1/1	0.97	0.04	49,49,49,49	0
4	MG	C	309	1/1	0.97	0.05	42,42,42,42	0
3	NA	C	307	1/1	0.97	0.05	48,48,48,48	0
2	FES	D	301	4/4	0.98	0.05	52,56,59,60	0
2	FES	C	302	4/4	0.98	0.07	50,52,58,60	0
2	FES	B	301	4/4	1.00	0.03	41,42,44,44	0

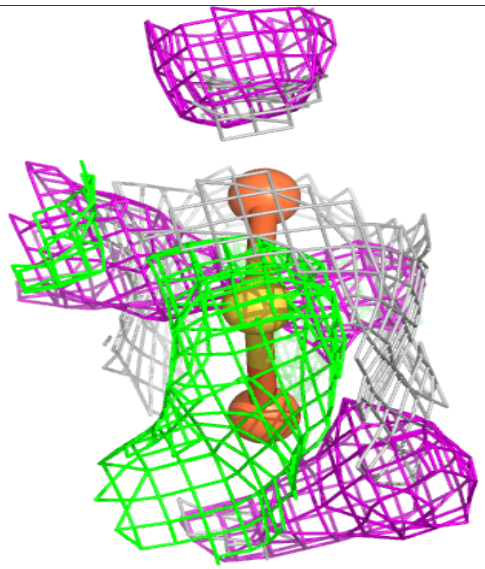
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FES A 301:

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

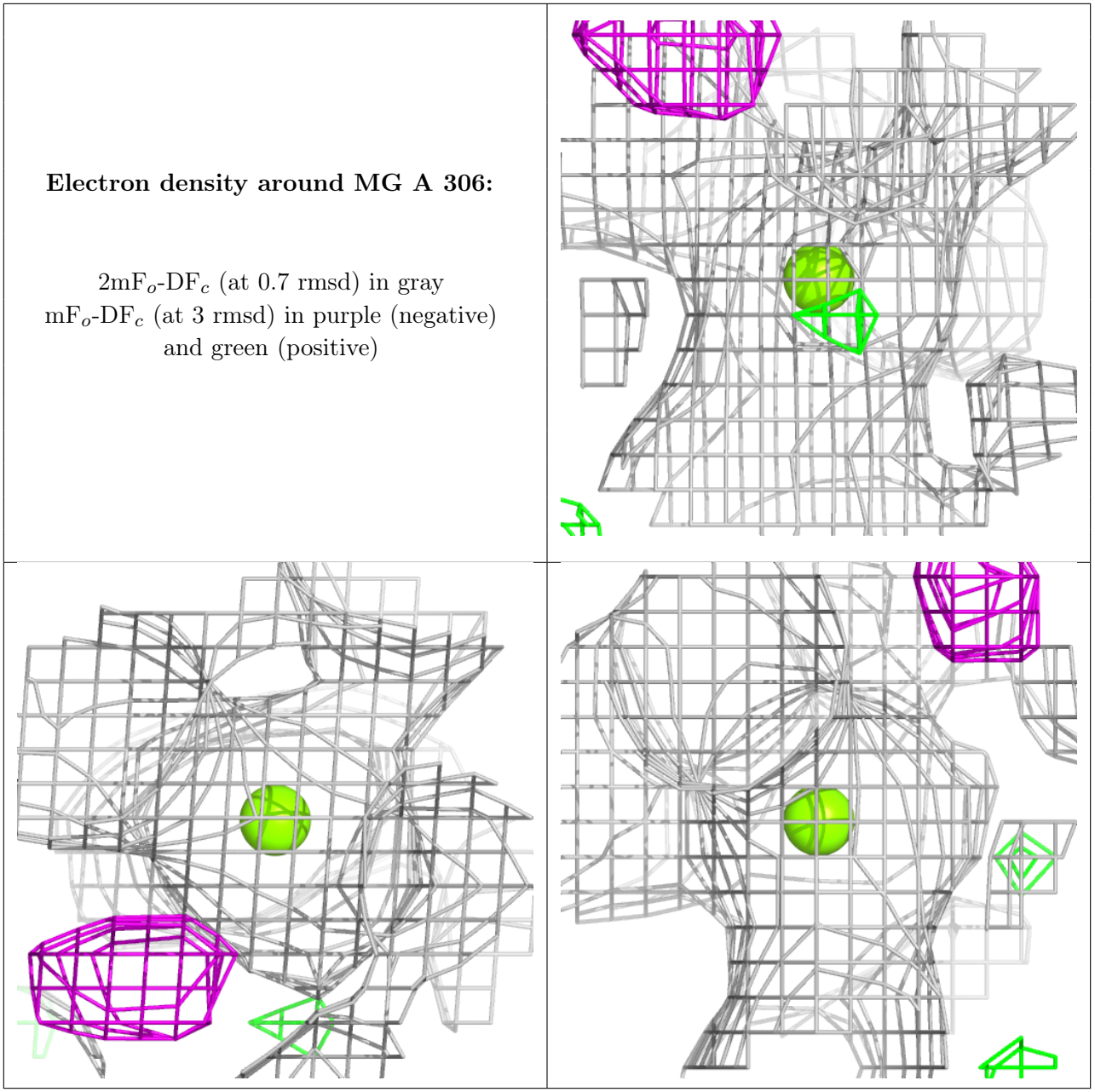






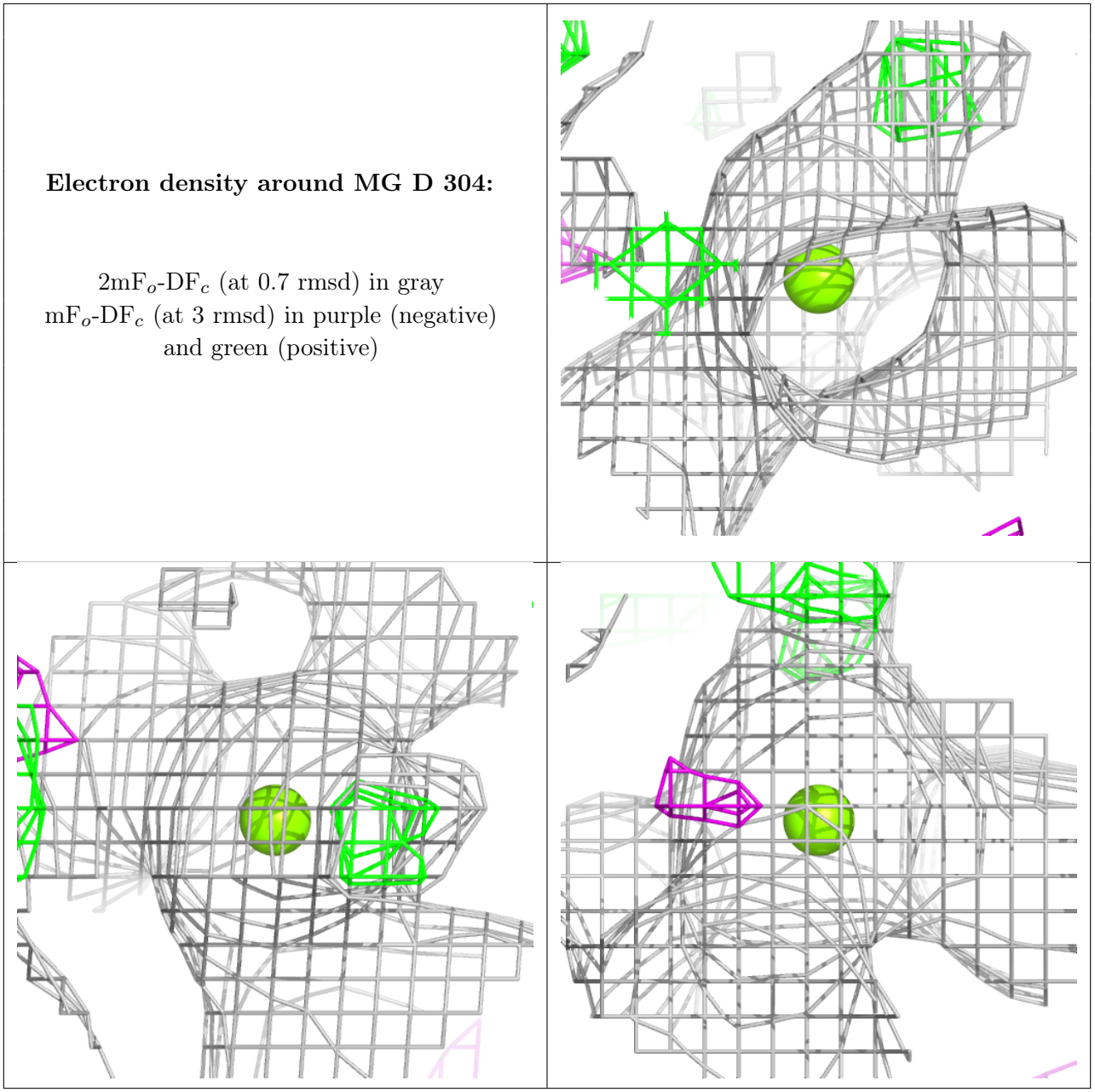
Electron density around MG A 306:

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



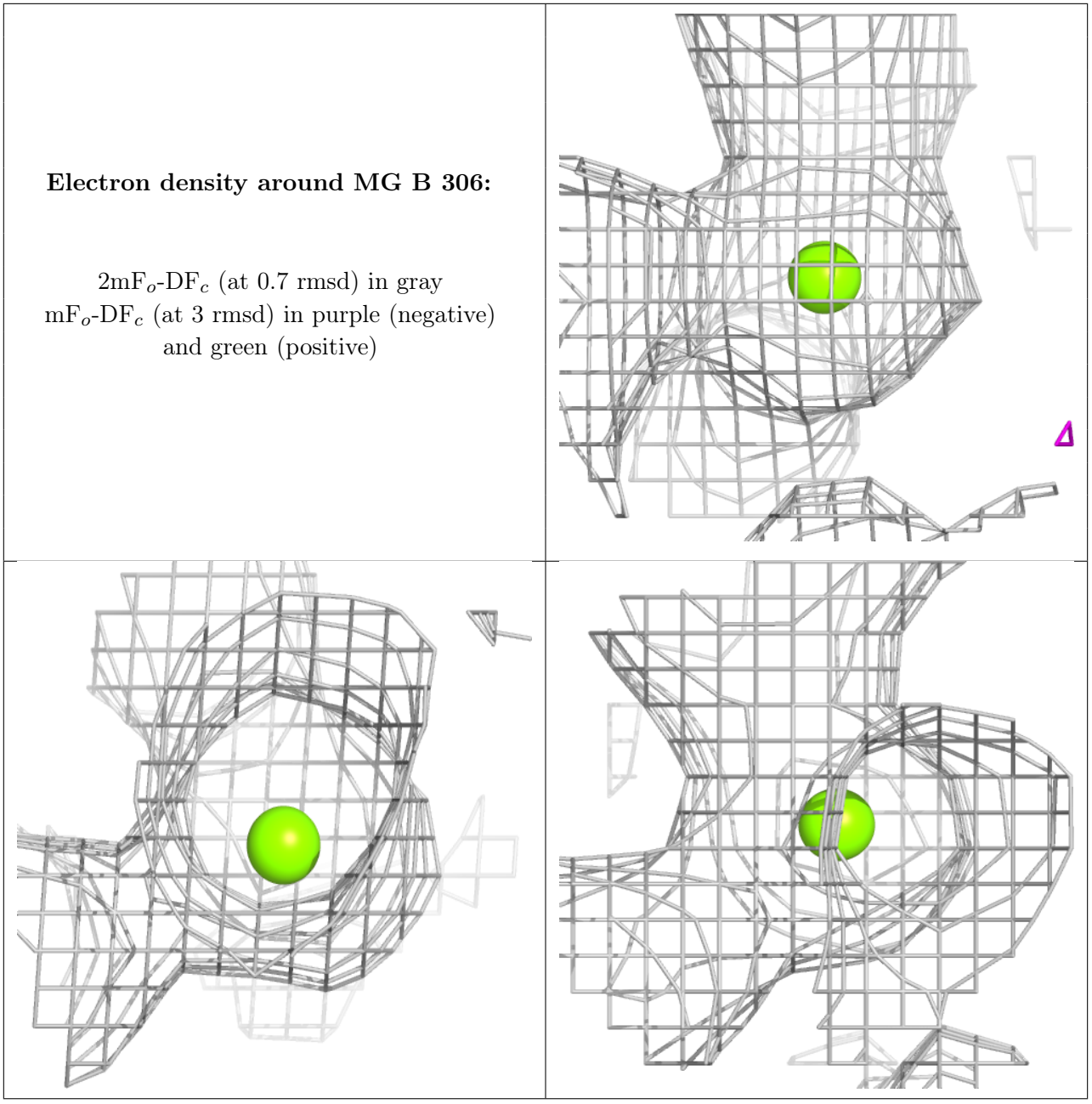
Electron density around MG D 304:

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



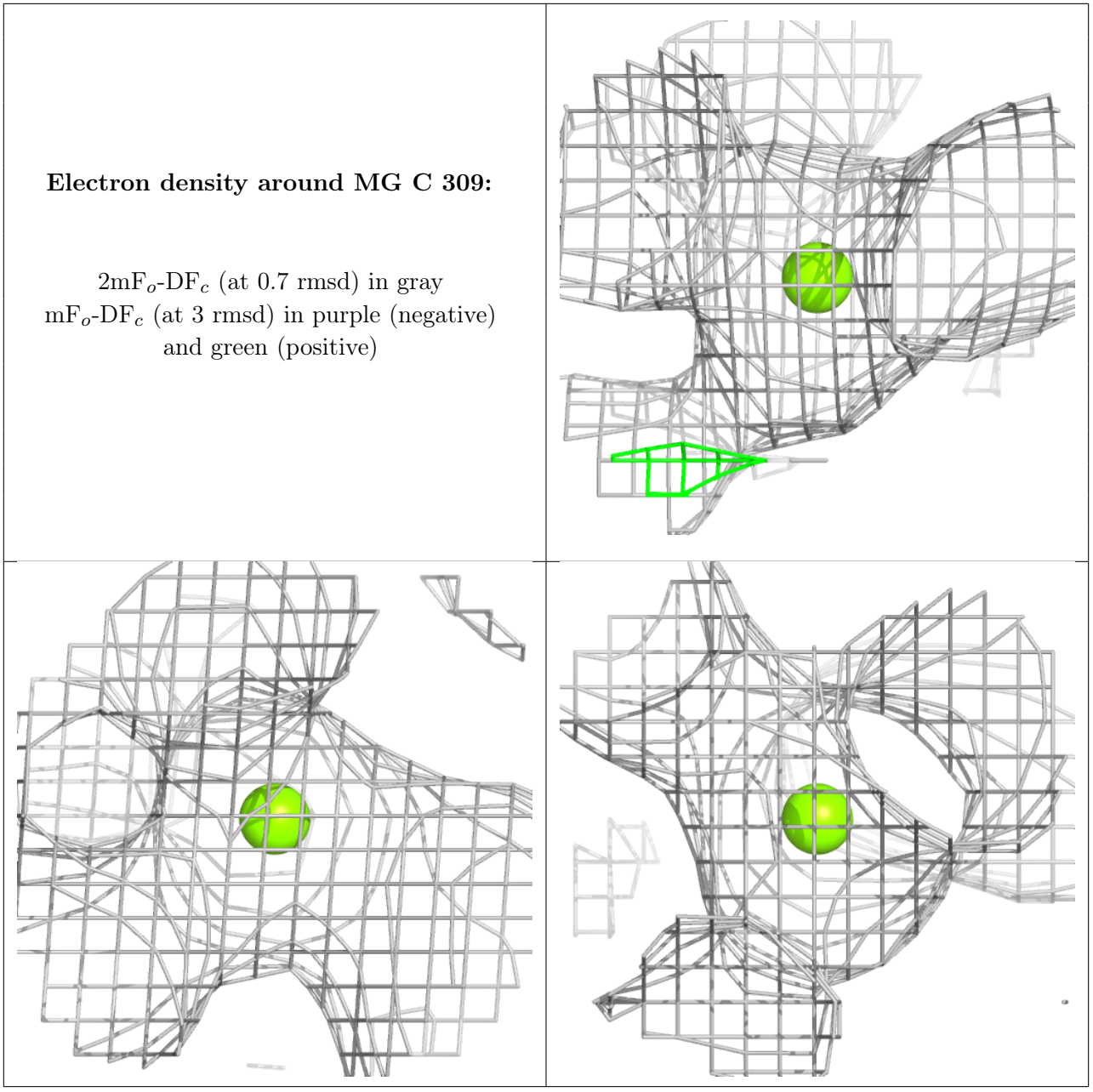
Electron density around MG B 306:

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



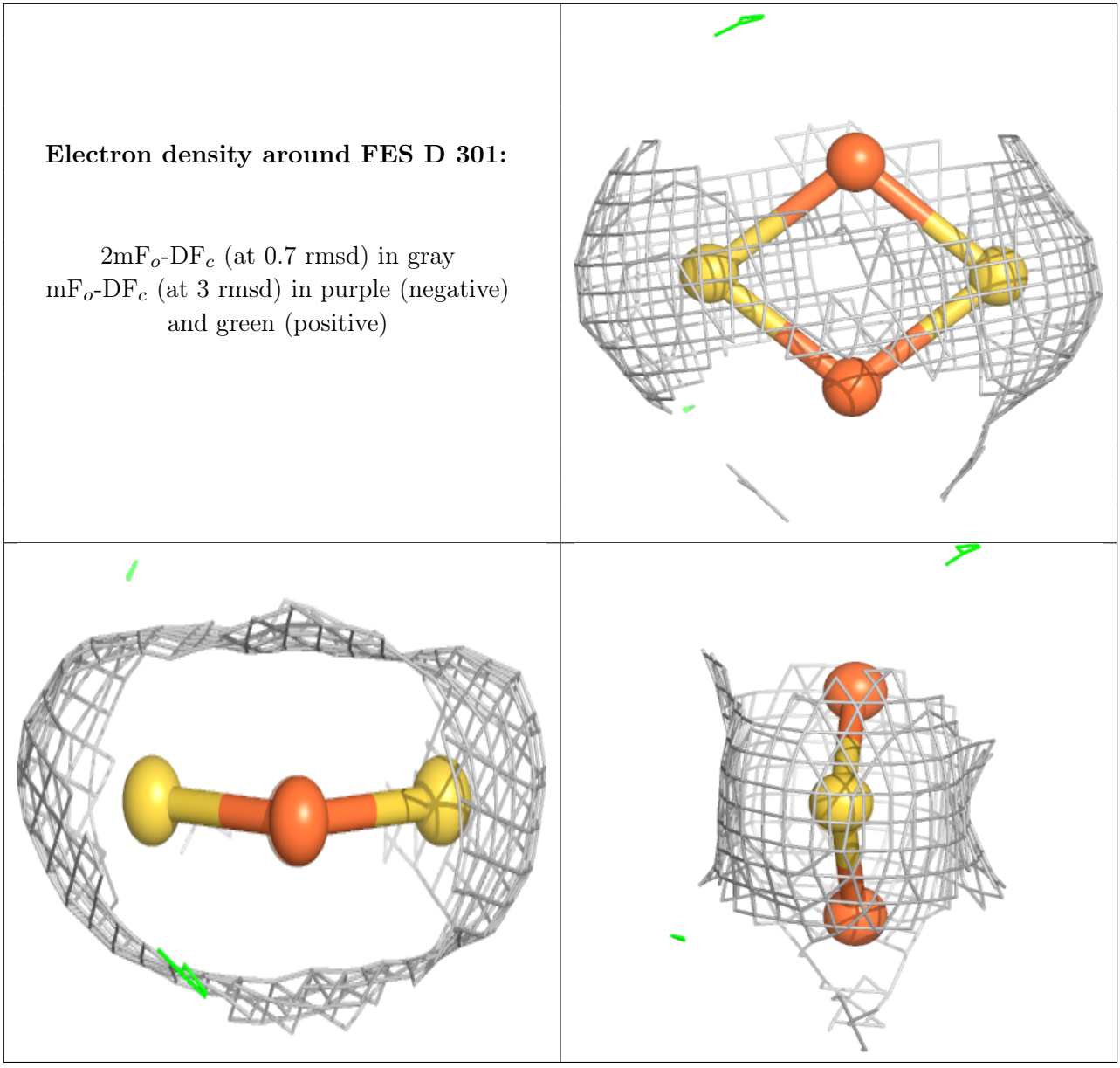
Electron density around MG C 309:

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



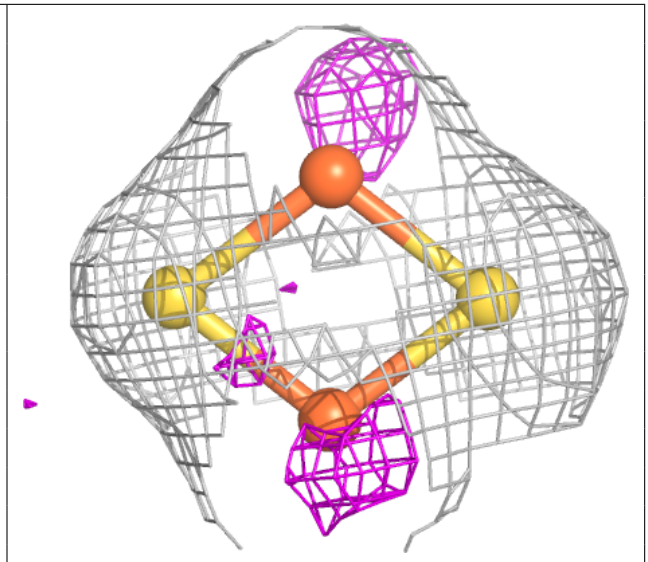
Electron density around FES D 301:

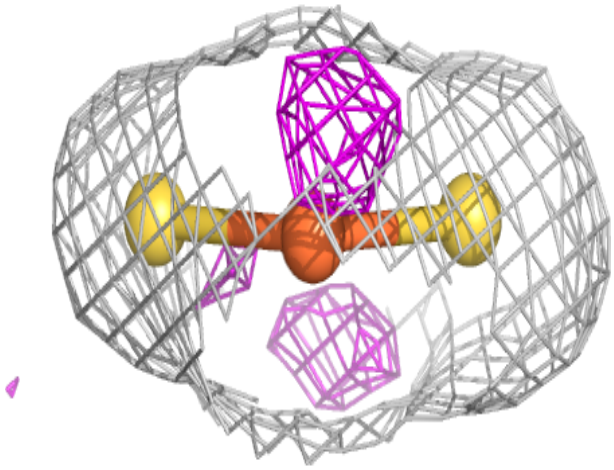
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

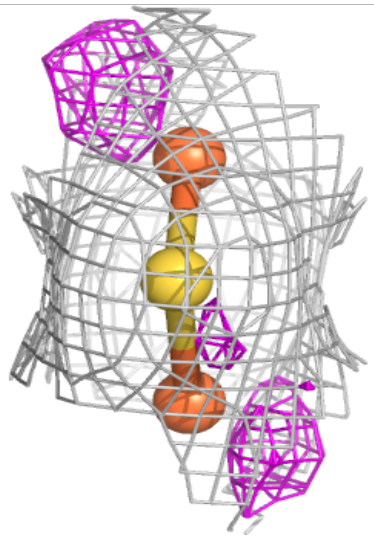


Electron density around FES C 302:

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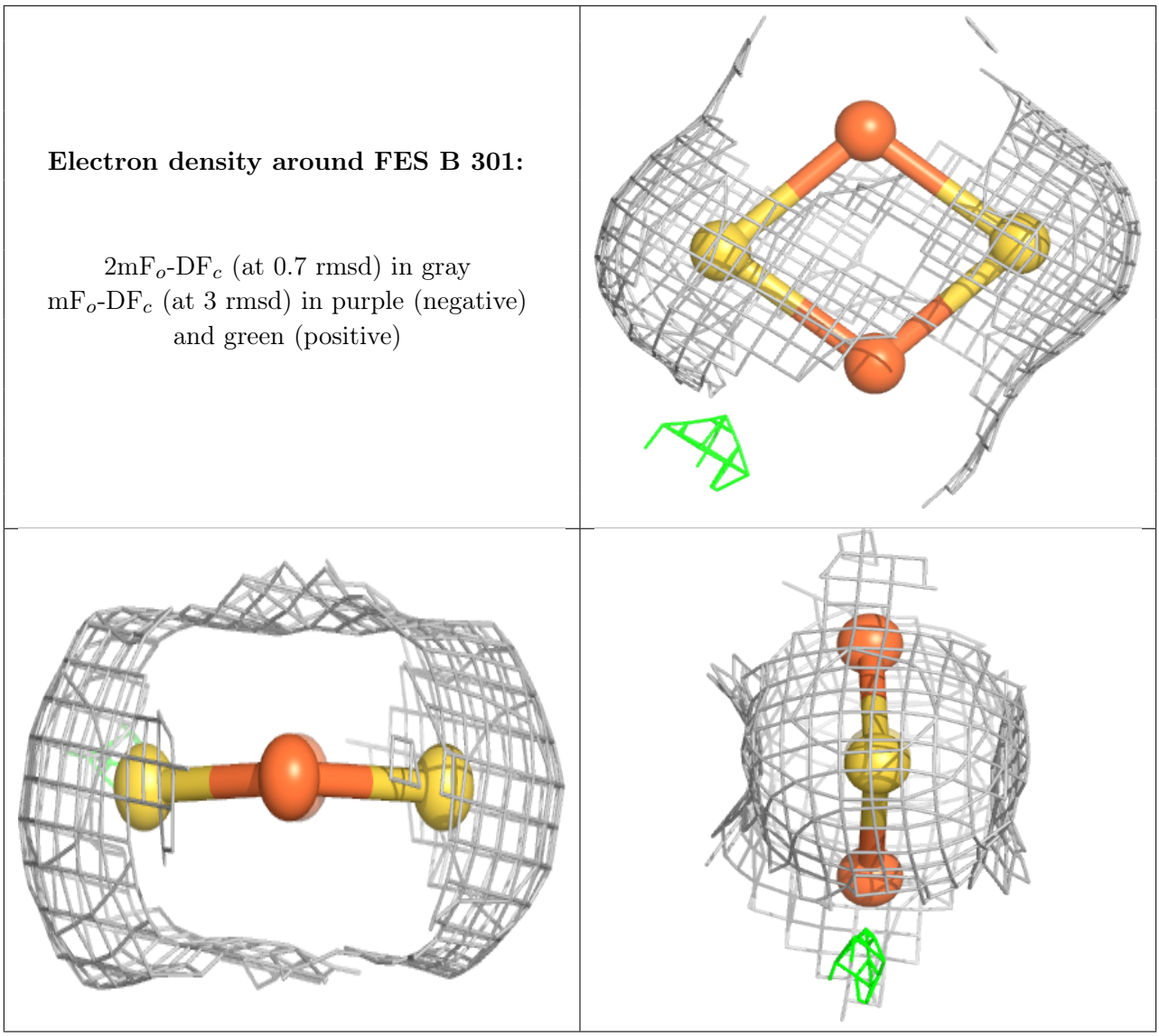






Electron density around FES B 301:

$2mF_o$-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.