Full wwPDB X-ray Structure Validation Report ⓘ

Nov 11, 2026 – 03:41 pm GMT

PDB ID	:	30OV / pdb_000030ov
Title	:	W-formate dehydrogenase H193A from Nitratidesulfovibrio vulgaris (Desulfovibrio vulgaris)
Authors	:	Vilela-Alves, G.; Manuel, R.R.; Oliveira, A.R.; Martins, G.; Pereira, I.C.; Romao, M.J.; Mota, C.
Deposited on	:	2026-05-06
Resolution	:	1.83 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

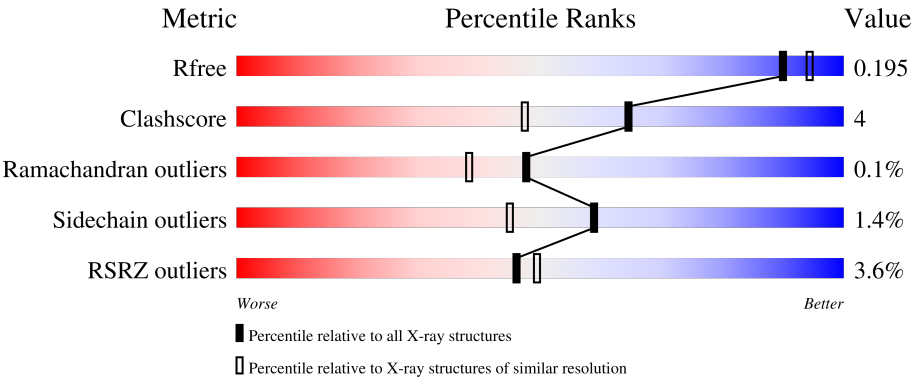
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

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 1.83 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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	1013	2% 87% 8% • 5%
1	E	1013	2% 89% 6% 5%
2	B	214	4% 88% 11% •
2	F	214	17% 91% 7% •

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
6	PG4	A	1105	-	-	X	-
6	PG4	E	1104	-	-	X	-
7	PEG	B	304	-	-	X	-

2 Entry composition (i)

There are 10 unique types of molecules in this entry. The entry contains 19680 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 Formate dehydrogenase, alpha subunit, selenocysteine-containing.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	963	Total	C	N	O	S	Se	0	2	0
			7541	4806	1314	1378	42	1			
1	E	963	Total	C	N	O	S	Se	0	0	0
			7528	4797	1311	1378	41	1			

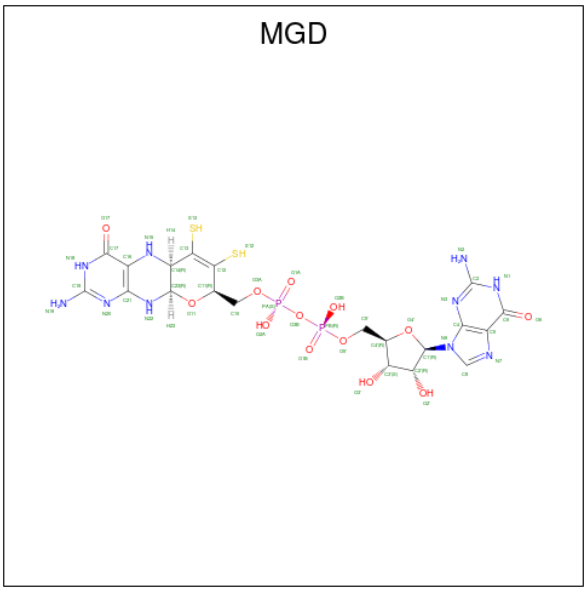
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	193	ALA	HIS	engineered mutation	UNP Q72EJ1
A	1006	TRP	-	expression tag	UNP Q72EJ1
A	1007	SER	-	expression tag	UNP Q72EJ1
A	1008	HIS	-	expression tag	UNP Q72EJ1
A	1009	PRO	-	expression tag	UNP Q72EJ1
A	1010	GLN	-	expression tag	UNP Q72EJ1
A	1011	PHE	-	expression tag	UNP Q72EJ1
A	1012	GLU	-	expression tag	UNP Q72EJ1
A	1013	LYS	-	expression tag	UNP Q72EJ1
E	193	ALA	HIS	engineered mutation	UNP Q72EJ1
E	1006	TRP	-	expression tag	UNP Q72EJ1
E	1007	SER	-	expression tag	UNP Q72EJ1
E	1008	HIS	-	expression tag	UNP Q72EJ1
E	1009	PRO	-	expression tag	UNP Q72EJ1
E	1010	GLN	-	expression tag	UNP Q72EJ1
E	1011	PHE	-	expression tag	UNP Q72EJ1
E	1012	GLU	-	expression tag	UNP Q72EJ1
E	1013	LYS	-	expression tag	UNP Q72EJ1

- Molecule 2 is a protein called Formate dehydrogenase, beta subunit, putative.

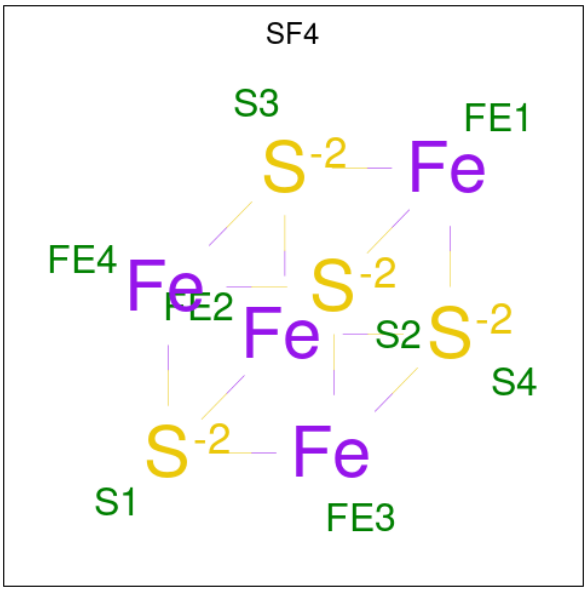
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	214	Total	C	N	O	S	0	0	0
			1664	1041	291	316	16			
2	F	214	Total	C	N	O	S	0	0	0
			1664	1041	291	316	16			

- Molecule 3 is 2-AMINO-5,6-DIMERCAPTO-7-METHYL-3,7,8A,9-TETRAHYDRO-8-OXA -1,3,9,10-TETRAAZA-ANTHRACEN-4-ONE GUANOSINE DINUCLEOTIDE (CCD ID: MGD) (formula: $C_{20}H_{26}N_{10}O_{13}P_2S_2$) (labeled as "Ligand of Interest" by depositor).



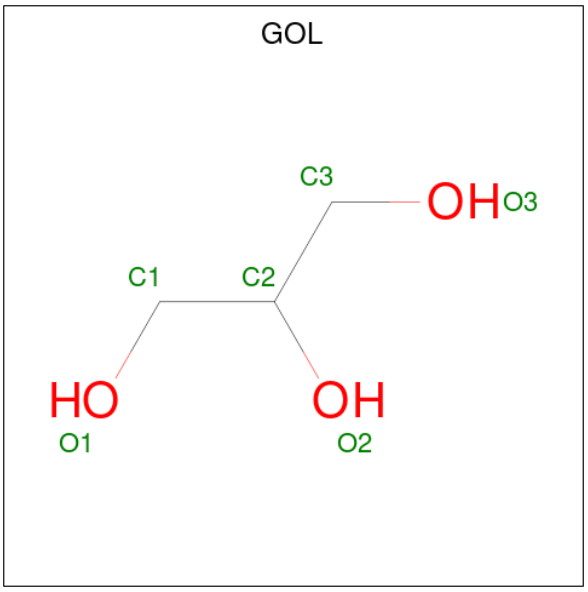
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			47	20	10	13	2	2		
3	A	1	Total	C	N	O	P	S	0	0
			47	20	10	13	2	2		
3	E	1	Total	C	N	O	P	S	0	0
			47	20	10	13	2	2		
3	E	1	Total	C	N	O	P	S	0	0
			47	20	10	13	2	2		

- Molecule 4 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4) (labeled as "Ligand of Interest" by depositor).



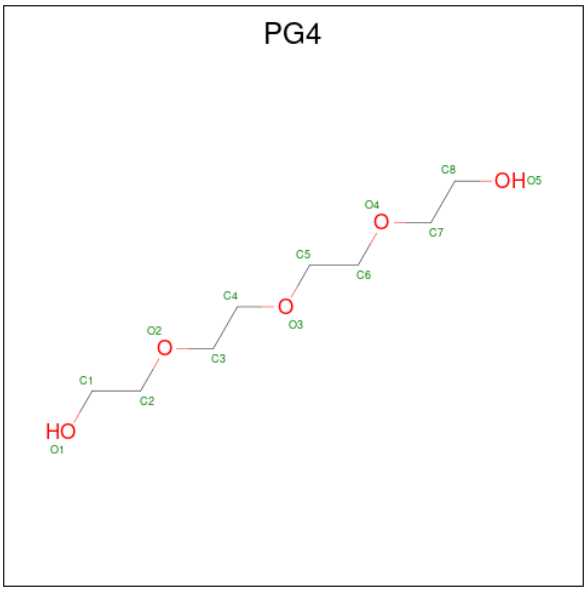
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Fe S 8 4 4	0	0
4	B	1	Total Fe S 8 4 4	0	0
4	B	1	Total Fe S 8 4 4	0	0
4	B	1	Total Fe S 8 4 4	0	0
4	E	1	Total Fe S 8 4 4	0	0
4	F	1	Total Fe S 8 4 4	0	0
4	F	1	Total Fe S 8 4 4	0	0
4	F	1	Total Fe S 8 4 4	0	0

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



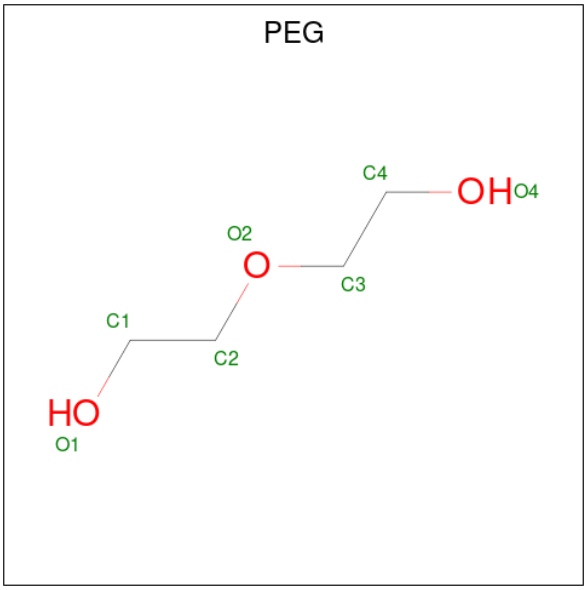
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	E	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 13 8 5	0	0
6	E	1	Total C O 13 8 5	0	0

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 7 4 3	0	0
7	A	1	Total C O 7 4 3	0	0
7	B	1	Total C O 7 4 3	0	0
7	E	1	Total C O 7 4 3	0	0
7	E	1	Total C O 7 4 3	0	0
7	E	1	Total C O 7 4 3	0	0

- Molecule 8 is TUNGSTEN ION (CCD ID: W) (formula: W) (labeled as "Ligand of Interest" by depositor).

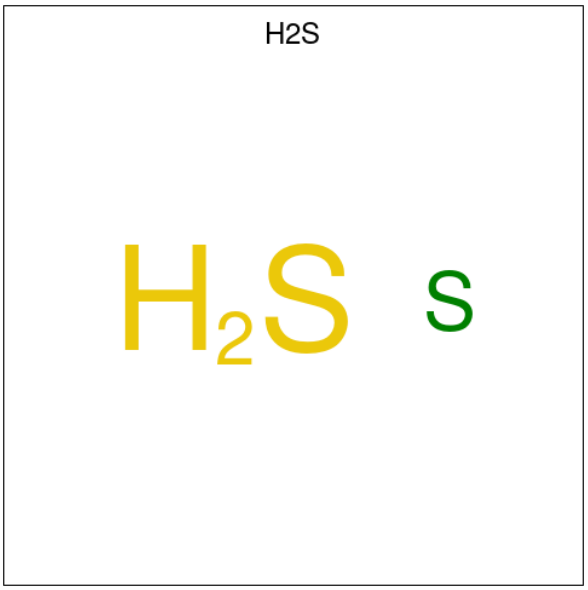
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total W 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	E	1	Total W 1 1	0	0

- Molecule 9 is HYDROSULFURIC ACID (CCD ID: H2S) (formula: H_2S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total S 1 1	0	0
9	E	1	Total S 1 1	0	0

- Molecule 10 is water.

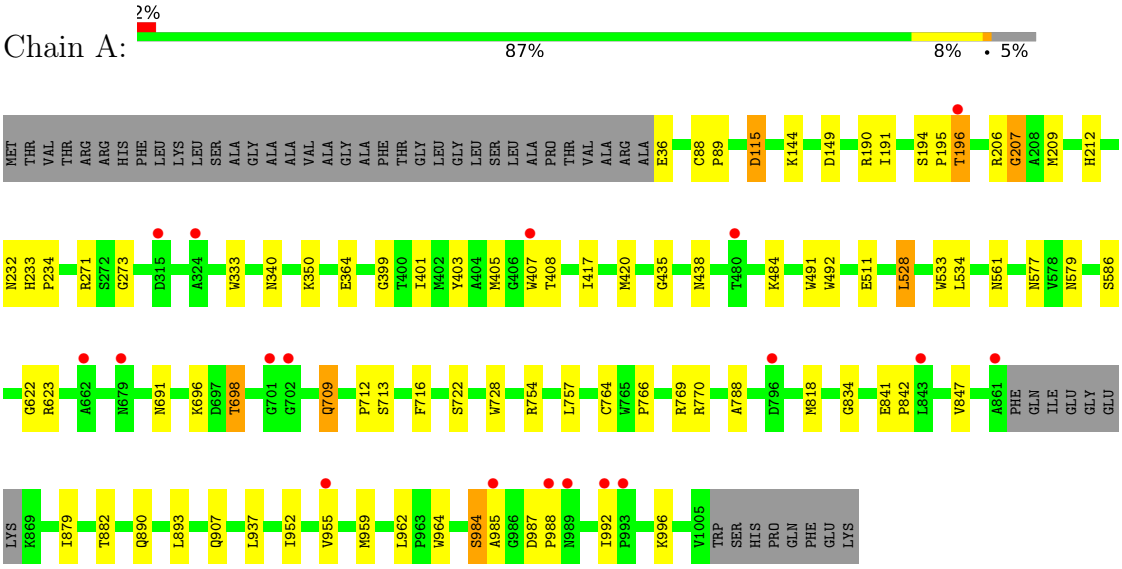
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	416	Total O 416 416	0	0
10	B	74	Total O 74 74	0	0
10	E	395	Total O 395 395	0	0
10	F	56	Total O 56 56	0	0

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.

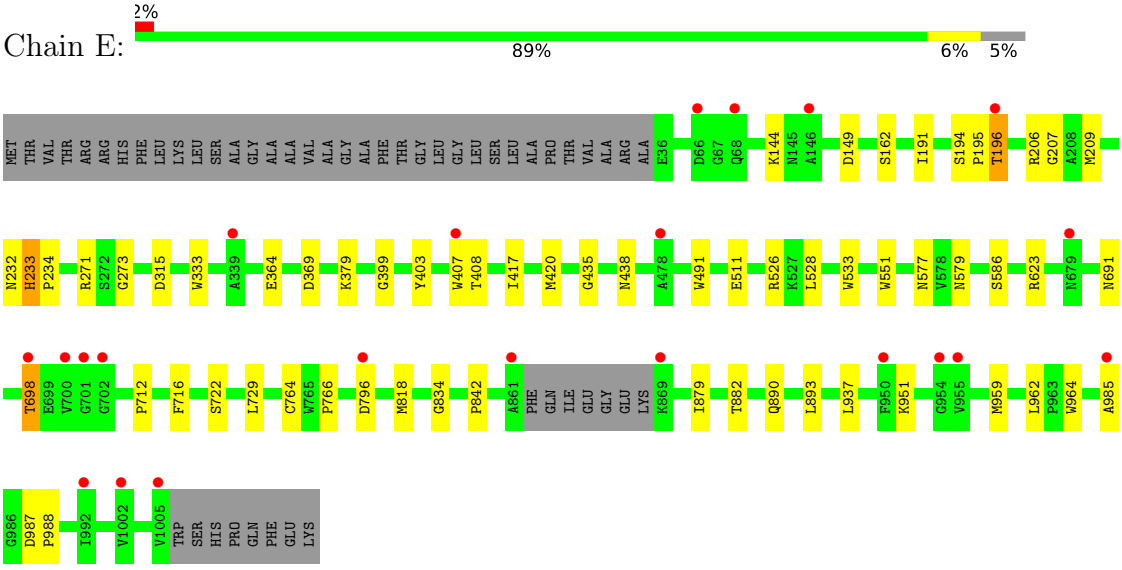
• Molecule 1: Formate dehydrogenase, alpha subunit, selenocysteine-containing

Chain A:



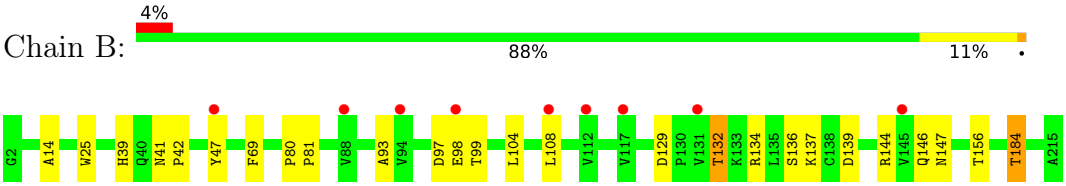
• Molecule 1: Formate dehydrogenase, alpha subunit, selenocysteine-containing

Chain E:



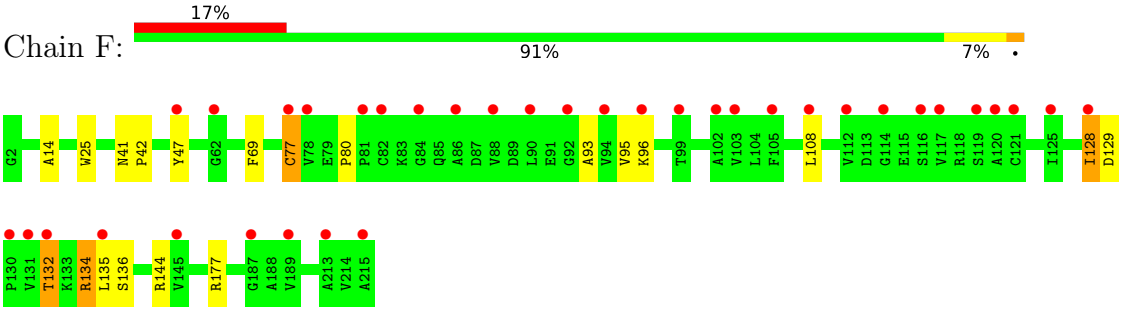
• Molecule 2: Formate dehydrogenase, beta subunit, putative

Chain B:



• Molecule 2: Formate dehydrogenase, beta subunit, putative

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	64.53Å 149.25Å 123.22Å 90.00° 90.71° 90.00°	Depositor
Resolution (Å)	59.23 – 1.83 59.23 – 1.83	Depositor EDS
% Data completeness (in resolution range)	97.2 (59.23-1.83) 71.7 (59.23-1.83)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.02 (at 1.83Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.167 , 0.186 0.178 , 0.195	Depositor DCC
R_{free} test set	6437 reflections (3.13%)	wwPDB-VP
Wilson B-factor (Å^2)	20.2	Xtriage
Anisotropy	0.162	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 43.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.149 for h,-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	19680	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PG4, PEG, W, MGD, H2S, SF4, SEC

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	1.05	2/7746 (0.0%)	1.25	4/10506 (0.0%)
1	E	1.05	3/7727 (0.0%)	1.25	2/10482 (0.0%)
2	B	1.09	0/1699	1.33	4/2302 (0.2%)
2	F	1.09	0/1699	1.33	1/2302 (0.0%)
All	All	1.06	5/18871 (0.0%)	1.27	11/25592 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	985	ALA	C-O	13.98	1.41	1.24
1	E	985	ALA	C-O	13.34	1.40	1.24
1	E	207	GLY	C-O	5.65	1.30	1.23
1	E	233	HIS	CE1-NE2	5.21	1.37	1.32
1	A	207	GLY	C-O	5.18	1.30	1.23

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	709	GLN	CB-CA-C	7.68	123.59	109.46
1	E	511	GLU	CB-CA-C	7.51	120.30	109.26
1	A	511	GLU	CB-CA-C	7.01	119.57	109.26
1	A	985	ALA	O-C-N	6.74	130.88	123.13
1	E	985	ALA	O-C-N	5.89	129.91	123.13
2	B	39	HIS	CA-CB-CG	-5.67	108.13	113.80
2	F	80	PRO	N-CA-C	5.49	117.40	110.70
1	A	115	ASP	CB-CA-C	5.32	118.90	109.65
2	B	97	ASP	CA-C-N	5.24	127.82	120.28
2	B	97	ASP	C-N-CA	5.24	127.82	120.28
2	B	80	PRO	N-CA-C	5.03	116.83	110.70

There are no chirality outliers.

There are no planarity outliers.

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	7541	0	7386	68	0
1	E	7528	0	7364	40	0
2	B	1664	0	1633	20	0
2	F	1664	0	1634	15	0
3	A	94	0	44	3	0
3	E	94	0	44	3	0
4	A	8	0	0	0	0
4	B	24	0	0	1	0
4	E	8	0	0	0	0
4	F	24	0	0	1	0
5	A	12	0	16	0	0
5	E	6	0	8	0	0
6	A	13	0	18	13	0
6	E	13	0	18	7	0
7	A	14	0	20	3	0
7	B	7	0	10	6	0
7	E	21	0	30	0	0
8	A	1	0	0	0	0
8	E	1	0	0	0	0
9	A	1	0	0	0	0
9	E	1	0	0	0	0
10	A	416	0	0	2	1
10	B	74	0	0	2	0
10	E	395	0	0	1	1
10	F	56	0	0	0	0
All	All	19680	0	18225	143	1

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 4.

All (143) 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:401:ILE:HD13	1:A:420[B]:MET:SD	2.02	0.99
1:A:579:ASN:OD1	3:A:1102:MGD:O2'	1.83	0.96
1:A:209:MET:HE2	1:A:622:GLY:HA2	1.48	0.93
1:E:579:ASN:OD1	3:E:1102:MGD:O2'	1.89	0.91
2:B:104:LEU:HA	7:B:304:PEG:H12	1.57	0.87
1:E:698:THR:HG21	1:E:722:SER:HB3	1.63	0.77
1:A:209:MET:HE3	1:A:212:HIS:HA	1.66	0.77
1:A:698:THR:HG21	1:A:722:SER:HB3	1.66	0.76
1:E:766:PRO:HB3	6:E:1104:PG4:H61	1.68	0.75
1:A:209:MET:CE	1:A:212:HIS:HA	2.20	0.71
1:A:209:MET:HE2	1:A:622:GLY:CA	2.21	0.70
1:E:766:PRO:CB	6:E:1104:PG4:H61	2.23	0.68
1:A:196:THR:HB	1:A:403:TYR:OH	1.93	0.68
1:E:407:TRP:CD1	1:E:408:THR:HG23	2.29	0.68
1:A:407:TRP:CD1	1:A:408:THR:HG23	2.29	0.68
2:F:77:CYS:SG	4:F:302:SF4:FE1	1.85	0.67
1:E:196:THR:HB	1:E:403:TYR:OH	1.95	0.67
2:B:104:LEU:CA	7:B:304:PEG:H12	2.24	0.67
1:A:407:TRP:CH2	1:A:420[A]:MET:HE3	2.30	0.66
1:E:206:ARG:HH21	6:E:1104:PG4:H72	1.61	0.66
2:F:132:THR:HG23	2:F:134:ARG:HB2	1.78	0.66
1:E:206:ARG:HE	6:E:1104:PG4:H62	1.60	0.65
2:F:132:THR:CG2	2:F:134:ARG:HB2	2.26	0.65
2:F:128:ILE:HG13	2:F:135:LEU:CD2	2.28	0.62
1:A:407:TRP:CH2	1:A:420[A]:MET:CE	2.83	0.62
1:A:401:ILE:HG21	1:A:420[B]:MET:SD	2.41	0.60
1:A:696:LYS:NZ	10:A:1204:HOH:O	2.34	0.60
1:E:407:TRP:CH2	1:E:420:MET:SD	2.95	0.60
1:A:190:ARG:O	1:A:992:ILE:HD11	2.03	0.59
1:E:407:TRP:CH2	1:E:420:MET:HE3	2.37	0.59
2:B:184:THR:CG2	10:B:434:HOH:O	2.51	0.58
1:E:407:TRP:CH2	1:E:420:MET:CE	2.86	0.58
1:A:407:TRP:CH2	1:A:420[A]:MET:SD	2.97	0.58
2:B:139:ASP:HB2	7:B:304:PEG:H11	1.86	0.57
2:F:128:ILE:HD12	2:F:128:ILE:N	2.19	0.57
7:A:1107:PEG:O4	10:A:1201:HOH:O	2.18	0.56
1:E:315:ASP:HA	10:E:1484:HOH:O	2.05	0.56
1:A:209:MET:CE	1:A:622:GLY:HA2	2.28	0.55
2:B:139:ASP:CB	7:B:304:PEG:H11	2.35	0.55
1:A:401:ILE:CD1	1:A:420[B]:MET:SD	2.88	0.54
2:B:41:ASN:HA	2:B:42:PRO:C	2.31	0.54
1:A:728:TRP:CH2	6:A:1105:PG4:H42	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:769:ARG:HH12	6:A:1105:PG4:C2	2.21	0.54
1:A:340:ASN:O	1:A:484:LYS:HG3	2.08	0.53
1:E:937:LEU:C	1:E:937:LEU:HD12	2.34	0.53
1:A:206:ARG:HE	6:A:1105:PG4:H31	1.73	0.53
2:F:41:ASN:HA	2:F:42:PRO:C	2.32	0.53
2:B:129:ASP:HB3	2:B:132:THR:HG22	1.90	0.53
1:A:818:MET:HE2	6:A:1105:PG4:H51	1.92	0.52
1:A:937:LEU:C	1:A:937:LEU:HD12	2.35	0.51
2:F:129:ASP:HB3	2:F:132:THR:HG22	1.92	0.51
1:A:206:ARG:HH21	6:A:1105:PG4:C3	2.24	0.51
2:B:132:THR:HG23	2:B:134:ARG:HG3	1.93	0.51
1:A:407:TRP:CZ2	1:A:417:ILE:CD1	2.94	0.50
1:A:273:GLY:HA3	1:A:842:PRO:HG2	1.93	0.50
1:A:209:MET:HG3	1:A:438:ASN:HA	1.94	0.49
1:E:206:ARG:NE	6:E:1104:PG4:H62	2.26	0.49
1:A:209:MET:CE	1:A:622:GLY:CA	2.88	0.49
1:A:407:TRP:CZ2	1:A:417:ILE:HD13	2.48	0.49
1:A:191:ILE:HG21	3:A:1102:MGD:H14	1.94	0.49
1:A:207:GLY:H	6:A:1105:PG4:H81	1.78	0.49
2:B:81:PRO:HD2	4:B:302:SF4:S3	2.53	0.49
1:E:407:TRP:CZ2	1:E:417:ILE:CD1	2.96	0.48
2:B:132:THR:HG21	2:B:134:ARG:HD2	1.95	0.48
1:E:232:ASN:HA	3:E:1101:MGD:N20	2.29	0.48
1:A:232:ASN:HA	3:A:1101:MGD:N20	2.28	0.48
1:E:407:TRP:CZ2	1:E:420:MET:SD	3.07	0.48
1:A:407:TRP:CZ2	1:A:420[A]:MET:SD	3.07	0.47
1:A:561:ASN:HD22	1:A:984:SER:HG	1.58	0.47
2:B:104:LEU:HA	7:B:304:PEG:C1	2.37	0.47
2:B:137:LYS:O	7:B:304:PEG:H31	2.14	0.47
1:A:907:GLN:HE22	7:A:1107:PEG:H32	1.79	0.47
2:F:128:ILE:HG13	2:F:135:LEU:HD23	1.96	0.47
2:B:184:THR:HG23	10:B:434:HOH:O	2.13	0.46
1:E:273:GLY:HA3	1:E:842:PRO:HG2	1.98	0.46
1:A:766:PRO:HB3	6:A:1105:PG4:H41	1.96	0.46
1:E:407:TRP:CZ2	1:E:417:ILE:HD13	2.50	0.46
2:F:47:TYR:CE1	2:F:136:SER:HB3	2.51	0.46
1:E:369:ASP:OD1	1:E:379:LYS:HD2	2.16	0.46
2:F:93:ALA:HA	2:F:108:LEU:HB2	1.98	0.46
1:A:194:SER:N	1:A:195:PRO:CD	2.78	0.46
1:A:233:HIS:N	1:A:234:PRO:HD3	2.31	0.46
1:A:754:ARG:HD2	1:A:788:ALA:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:528:LEU:HD12	1:E:528:LEU:HA	1.84	0.46
1:E:191:ILE:HG21	3:E:1102:MGD:H14	1.98	0.45
1:E:729:LEU:HD21	6:E:1104:PG4:H41	1.98	0.45
1:E:209:MET:HG3	1:E:438:ASN:HA	1.98	0.45
1:E:194:SER:N	1:E:195:PRO:CD	2.79	0.45
1:E:577:ASN:OD1	1:E:586:SER:HB3	2.17	0.45
1:A:36:GLU:N	2:B:147:ASN:HD21	2.14	0.45
1:A:577:ASN:OD1	1:A:586:SER:HB3	2.16	0.45
2:F:25:TRP:CE2	2:F:144:ARG:HD2	2.52	0.45
1:A:333:TRP:CH2	1:A:834:GLY:HA2	2.52	0.45
2:F:128:ILE:HG13	2:F:135:LEU:HD21	1.99	0.44
1:A:959:MET:HE2	1:A:959:MET:HB3	1.86	0.44
1:E:712:PRO:HG2	1:E:716:PHE:CD2	2.53	0.44
1:E:623:ARG:O	1:E:764:CYS:HA	2.18	0.44
1:A:491:TRP:CE2	6:A:1105:PG4:H11	2.53	0.44
1:A:757:LEU:HD22	1:A:770:ARG:HB2	1.99	0.44
1:E:491:TRP:HB3	1:E:818:MET:SD	2.58	0.44
1:A:890:GLN:HA	1:A:964:TRP:CH2	2.53	0.43
2:F:177:ARG:HG2	2:F:177:ARG:HH11	1.84	0.43
2:F:95:VAL:HG22	2:F:96:LYS:N	2.34	0.43
1:A:769:ARG:NH1	6:A:1105:PG4:H22	2.34	0.43
1:A:623:ARG:O	1:A:764:CYS:HA	2.19	0.43
2:B:99:THR:OG1	2:B:146:GLN:NE2	2.35	0.43
1:E:882:THR:HA	1:E:962:LEU:O	2.19	0.43
1:E:987:ASP:HB2	1:E:988:PRO:HD2	2.01	0.43
1:A:88:CYS:HB2	1:A:89:PRO:HD2	2.00	0.43
2:B:25:TRP:CE2	2:B:144:ARG:HD2	2.53	0.43
2:B:132:THR:HG23	2:B:134:ARG:CG	2.49	0.43
1:A:952:ILE:O	1:A:955:VAL:HG22	2.19	0.42
1:A:206:ARG:HH21	6:A:1105:PG4:H32	1.82	0.42
1:E:162:SER:HB2	1:E:551:TRP:O	2.19	0.42
1:E:333:TRP:CH2	1:E:834:GLY:HA2	2.54	0.42
1:A:987:ASP:HB2	1:A:988:PRO:HD2	2.00	0.42
2:B:93:ALA:HA	2:B:108:LEU:HB2	1.99	0.42
1:E:491:TRP:CD1	6:E:1104:PG4:H82	2.54	0.42
1:E:879:ILE:O	1:E:959:MET:HA	2.20	0.42
1:A:769:ARG:NH1	6:A:1105:PG4:C2	2.83	0.42
1:E:233:HIS:N	1:E:234:PRO:HD3	2.34	0.42
1:A:206:ARG:HH21	6:A:1105:PG4:H21	1.85	0.42
2:B:47:TYR:CE1	2:B:136:SER:HB3	2.55	0.42
1:A:712:PRO:HG2	1:A:716:PHE:CD2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:MET:SD	1:A:996:LYS:HE3	2.60	0.41
1:A:144:LYS:HA	1:A:149:ASP:O	2.20	0.41
2:F:14:ALA:HB2	2:F:69:PHE:CG	2.54	0.41
1:A:534:LEU:HD23	1:A:534:LEU:HA	1.96	0.41
1:A:882:THR:HA	1:A:962:LEU:O	2.21	0.41
1:A:841:GLU:HG2	1:A:847:VAL:HG11	2.03	0.41
1:A:907:GLN:NE2	7:A:1107:PEG:H32	2.36	0.41
1:E:144:LYS:HA	1:E:149:ASP:O	2.21	0.41
1:E:890:GLN:HA	1:E:964:TRP:CH2	2.55	0.41
1:A:206:ARG:HE	6:A:1105:PG4:C3	2.34	0.41
1:A:399:GLY:O	1:A:435:GLY:HA2	2.21	0.41
1:A:492:TRP:CD1	1:A:818:MET:HE3	2.56	0.41
2:B:14:ALA:HB2	2:B:69:PHE:CG	2.56	0.41
1:E:399:GLY:O	1:E:435:GLY:HA2	2.21	0.41
1:A:879:ILE:O	1:A:959:MET:HA	2.21	0.40
1:A:893:LEU:C	1:A:893:LEU:HD23	2.46	0.40
1:E:893:LEU:C	1:E:893:LEU:HD23	2.46	0.40
1:A:209:MET:HE1	1:A:212:HIS:HA	2.00	0.40
1:A:528:LEU:HA	1:A:528:LEU:HD12	1.83	0.40

All (1) 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)
10:A:1556:HOH:O	10:E:1428:HOH:O[2_546]	2.18	0.02

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	960/1013 (95%)	930 (97%)	29 (3%)	1 (0%)	48 38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	E	958/1013 (95%)	930 (97%)	27 (3%)	1 (0%)	48 38
2	B	212/214 (99%)	202 (95%)	10 (5%)	0	100 100
2	F	212/214 (99%)	202 (95%)	10 (5%)	0	100 100
All	All	2342/2454 (95%)	2264 (97%)	76 (3%)	2 (0%)	48 38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	533	TRP
1	E	533	TRP

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	783/818 (96%)	772 (99%)	11 (1%)	59 45
1	E	781/818 (96%)	773 (99%)	8 (1%)	68 58
2	B	185/185 (100%)	181 (98%)	4 (2%)	45 29
2	F	185/185 (100%)	181 (98%)	4 (2%)	45 29
All	All	1934/2006 (96%)	1907 (99%)	27 (1%)	59 45

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	ASP
1	A	196	THR
1	A	271	ARG
1	A	350	LYS
1	A	364	GLU
1	A	528	LEU
1	A	691	ASN
1	A	698	THR
1	A	709	GLN

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Mol	Chain	Res	Type
1	A	713	SER
1	A	984	SER
2	B	98	GLU
2	B	132	THR
2	B	156	THR
2	B	184	THR
1	E	196	THR
1	E	271	ARG
1	E	364	GLU
1	E	526	ARG
1	E	691	ASN
1	E	698	THR
1	E	796	ASP
1	E	951	LYS
2	F	77	CYS
2	F	128	ILE
2	F	132	THR
2	F	134	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	GLN
1	A	989	ASN
1	E	106	GLN
1	E	289	ASN
1	E	383	GLN
1	E	463	ASN
2	F	85	GLN

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 (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 27 ligands modelled in this entry, 2 are monoatomic and 2 are modelled with single atom - leaving 23 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	$\#\|Z\| > 2$
7	PEG	E	1108	-	6,6,6	0.31	0	5,5,5	0.13	0
4	SF4	B	301	2	0,12,12	-	-	-		
6	PG4	A	1105	-	12,12,12	0.48	0	11,11,11	0.57	0
4	SF4	B	302	2	0,12,12	-	-	-		
5	GOL	A	1106	-	5,5,5	0.18	0	5,5,5	0.44	0
4	SF4	E	1103	1	0,12,12	-	-	-		
3	MGD	A	1101	8	45,52,52	0.65	0	54,81,81	0.88	2 (3%)
3	MGD	E	1102	8	45,52,52	0.78	1 (2%)	54,81,81	0.92	3 (5%)
5	GOL	E	1106	-	5,5,5	0.11	0	5,5,5	0.33	0
5	GOL	A	1104	-	5,5,5	0.15	0	5,5,5	0.47	0
7	PEG	A	1108	-	6,6,6	0.30	0	5,5,5	0.28	0
4	SF4	A	1103	1	0,12,12	-	-	-		
4	SF4	B	303	2	0,12,12	-	-	-		
7	PEG	A	1107	-	6,6,6	0.16	0	5,5,5	0.24	0
4	SF4	F	302	2	0,12,12	-	-	-		
3	MGD	E	1101	8	45,52,52	0.53	0	54,81,81	0.97	3 (5%)
7	PEG	E	1107	-	6,6,6	0.38	0	5,5,5	0.30	0
7	PEG	E	1105	-	6,6,6	0.38	0	5,5,5	0.28	0
6	PG4	E	1104	-	12,12,12	0.41	0	11,11,11	0.67	0
4	SF4	F	301	2	0,12,12	-	-	-		
7	PEG	B	304	-	6,6,6	0.80	0	5,5,5	0.67	0
3	MGD	A	1102	8	45,52,52	0.55	0	54,81,81	0.98	4 (7%)
4	SF4	F	303	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	PEG	E	1108	-	-	2/4/4/4	-
4	SF4	B	301	2	-	-	0/6/5/5
6	PG4	A	1105	-	-	6/10/10/10	-
4	SF4	B	302	2	-	-	0/6/5/5
5	GOL	A	1106	-	-	2/4/4/4	-
4	SF4	E	1103	1	-	-	0/6/5/5
3	MGD	A	1101	8	-	2/22/66/66	0/6/6/6
3	MGD	E	1102	8	-	4/22/66/66	0/6/6/6
5	GOL	E	1106	-	-	0/4/4/4	-
5	GOL	A	1104	-	-	0/4/4/4	-
7	PEG	A	1108	-	-	3/4/4/4	-
4	SF4	A	1103	1	-	-	0/6/5/5
4	SF4	B	303	2	-	-	0/6/5/5
7	PEG	A	1107	-	-	2/4/4/4	-
4	SF4	F	302	2	-	-	0/6/5/5
3	MGD	E	1101	8	-	1/22/66/66	0/6/6/6
7	PEG	E	1107	-	-	2/4/4/4	-
7	PEG	E	1105	-	-	1/4/4/4	-
6	PG4	E	1104	-	-	3/10/10/10	-
4	SF4	F	301	2	-	-	0/6/5/5
7	PEG	B	304	-	-	4/4/4/4	-
3	MGD	A	1102	8	-	4/22/66/66	0/6/6/6
4	SF4	F	303	2	-	-	0/6/5/5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	1102	MGD	C23-C14	3.65	1.56	1.53

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	E	1102	MGD	O11-C23-C14	3.25	111.13	108.96
3	A	1102	MGD	O11-C23-N22	-3.00	105.48	108.57
3	E	1101	MGD	O11-C23-C14	2.96	110.94	108.96
3	E	1101	MGD	C19-N20-C21	2.82	118.53	113.43
3	A	1101	MGD	C19-N20-C21	2.66	118.23	113.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	E	1102	MGD	C19-N20-C21	2.55	118.03	113.43
3	A	1102	MGD	O11-C23-C14	2.48	110.62	108.96
3	A	1102	MGD	C17-C16-N15	2.45	123.32	116.76
3	A	1102	MGD	C19-N20-C21	2.40	117.76	113.43
3	A	1101	MGD	C4'-O4'-C1'	-2.28	104.44	109.47
3	E	1102	MGD	C17-C16-N15	2.11	122.44	116.76
3	E	1101	MGD	C4'-O4'-C1'	-2.00	105.06	109.47

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1102	MGD	C4'-C5'-O5'-PB
3	E	1102	MGD	C4'-C5'-O5'-PB
3	A	1102	MGD	O4'-C4'-C5'-O5'
3	A	1102	MGD	C3'-C4'-C5'-O5'
3	E	1102	MGD	O4'-C4'-C5'-O5'
3	E	1102	MGD	C3'-C4'-C5'-O5'
6	E	1104	PG4	O2-C3-C4-O3
7	E	1108	PEG	O1-C1-C2-O2
6	A	1105	PG4	O2-C3-C4-O3
7	E	1105	PEG	O2-C3-C4-O4
5	A	1106	GOL	C1-C2-C3-O3
6	E	1104	PG4	O1-C1-C2-O2
7	B	304	PEG	O1-C1-C2-O2
5	A	1106	GOL	O2-C2-C3-O3
7	E	1107	PEG	O2-C3-C4-O4
7	A	1107	PEG	O1-C1-C2-O2
7	A	1108	PEG	O1-C1-C2-O2
7	A	1108	PEG	O2-C3-C4-O4
7	B	304	PEG	O2-C3-C4-O4
6	A	1105	PG4	O3-C5-C6-O4
3	A	1101	MGD	PA-O3B-PB-O5'
3	E	1101	MGD	PA-O3B-PB-O5'
6	A	1105	PG4	C6-C5-O3-C4
7	B	304	PEG	C4-C3-O2-C2
7	A	1107	PEG	O2-C3-C4-O4
6	E	1104	PG4	C1-C2-O2-C3
6	A	1105	PG4	O1-C1-C2-O2
7	E	1107	PEG	O1-C1-C2-O2
7	E	1108	PEG	C4-C3-O2-C2
7	A	1108	PEG	C1-C2-O2-C3

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Mol	Chain	Res	Type	Atoms
6	A	1105	PG4	C5-C6-O4-C7
3	A	1101	MGD	PA-O3B-PB-O1B
3	A	1102	MGD	C5'-O5'-PB-O1B
3	E	1102	MGD	C5'-O5'-PB-O1B
7	B	304	PEG	C1-C2-O2-C3
6	A	1105	PG4	C4-C3-O2-C2

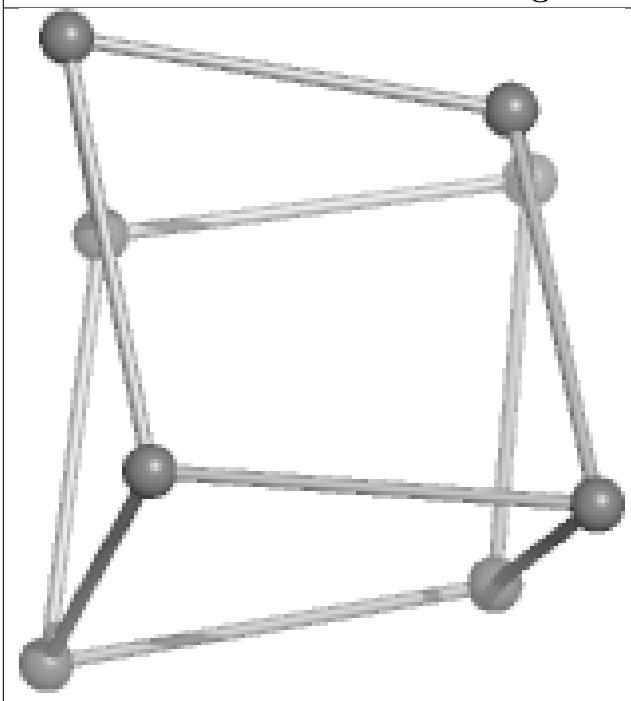
There are no ring outliers.

10 monomers are involved in 37 short contacts:

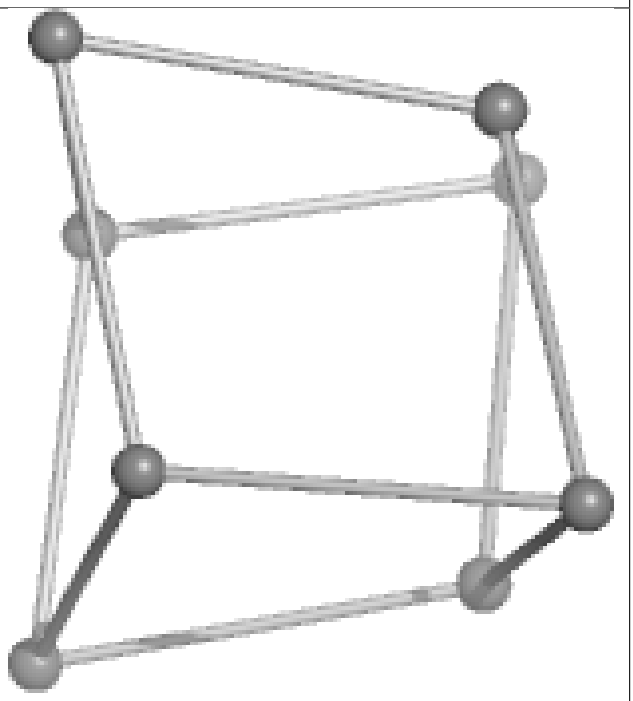
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1105	PG4	13	0
4	B	302	SF4	1	0
3	A	1101	MGD	1	0
3	E	1102	MGD	2	0
7	A	1107	PEG	3	0
4	F	302	SF4	1	0
3	E	1101	MGD	1	0
6	E	1104	PG4	7	0
7	B	304	PEG	6	0
3	A	1102	MGD	2	0

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 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.

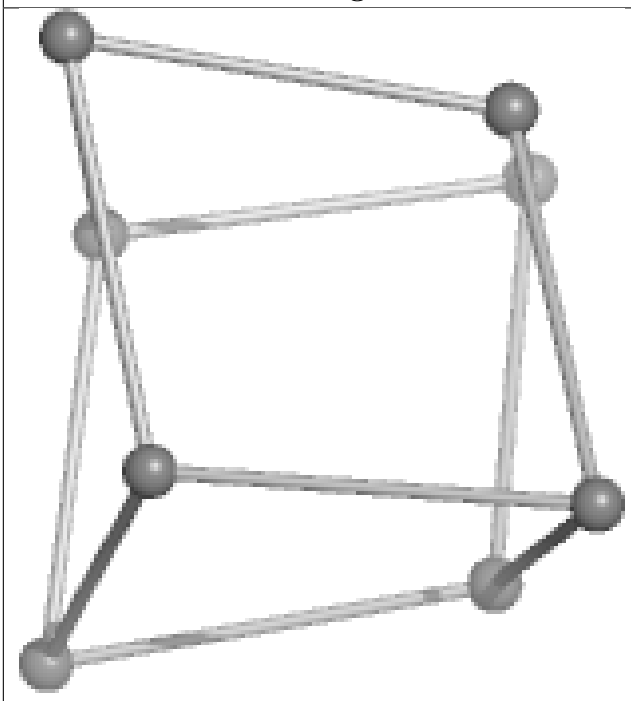
Ligand SF4 B 301



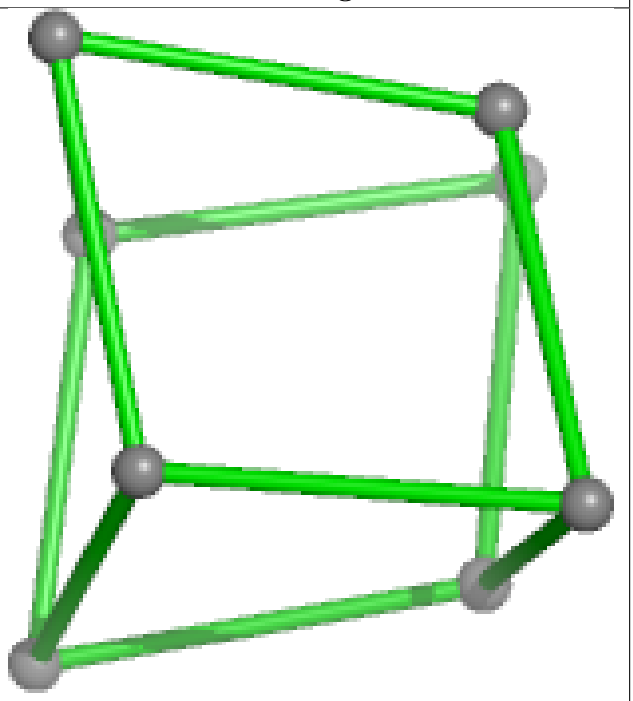
Bond lengths



Bond angles

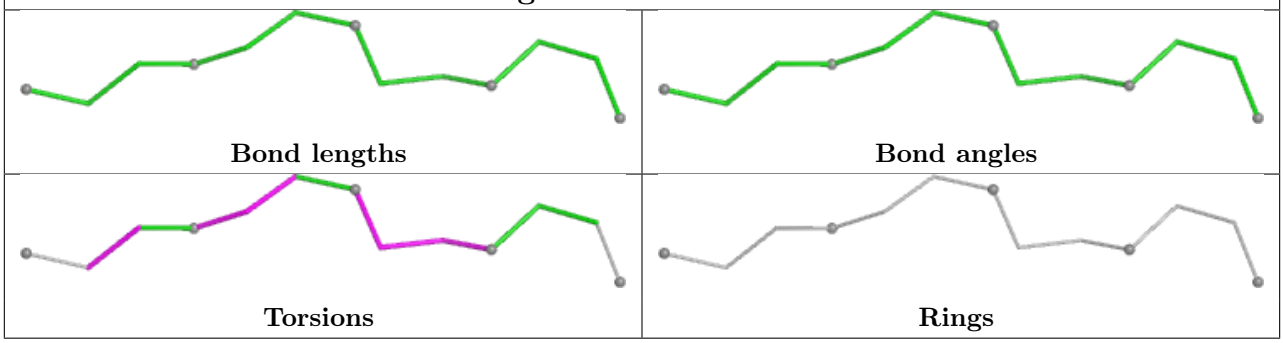


Torsions

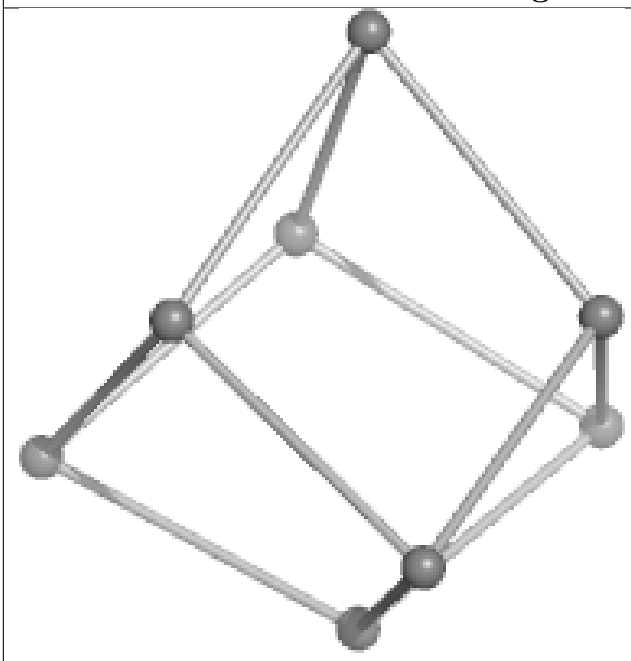


Rings

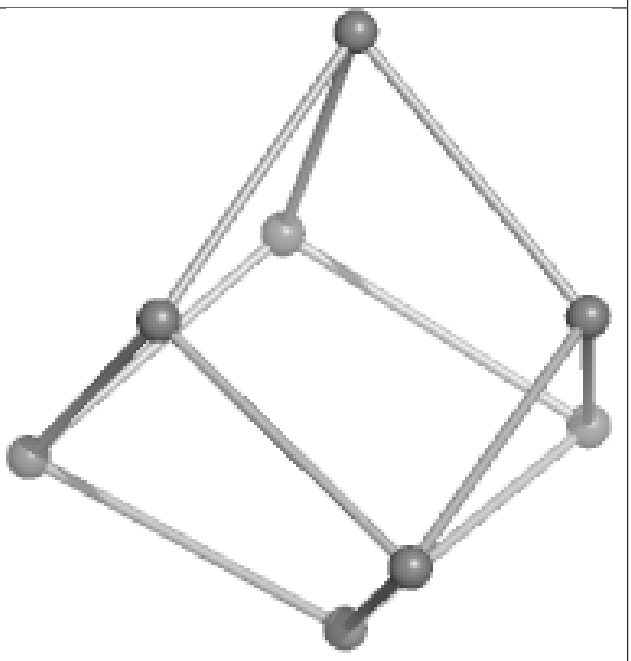
Ligand PG4 A 1105



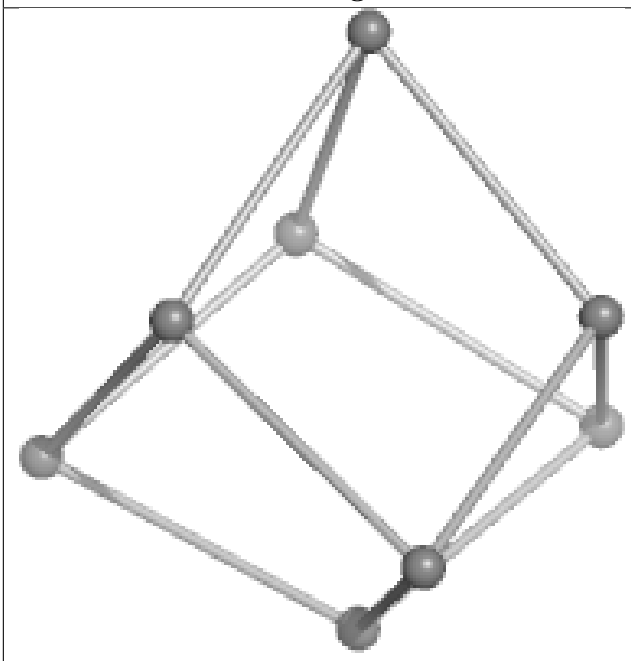
Ligand SF4 B 302



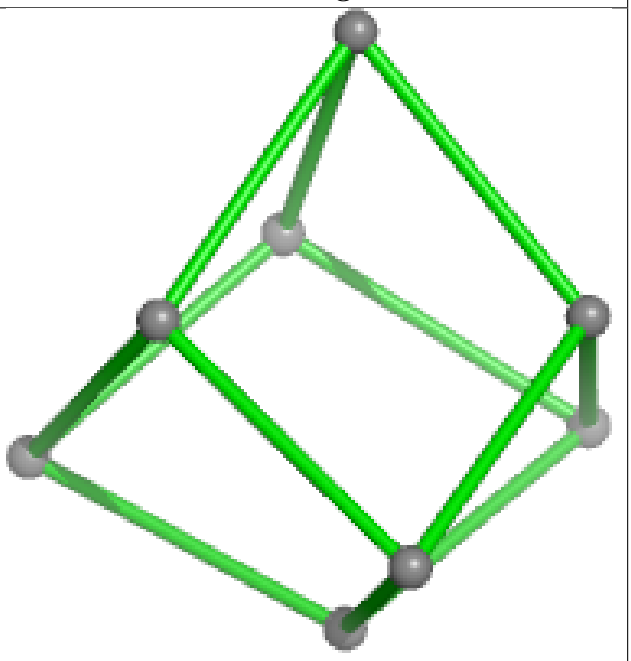
Bond lengths



Bond angles

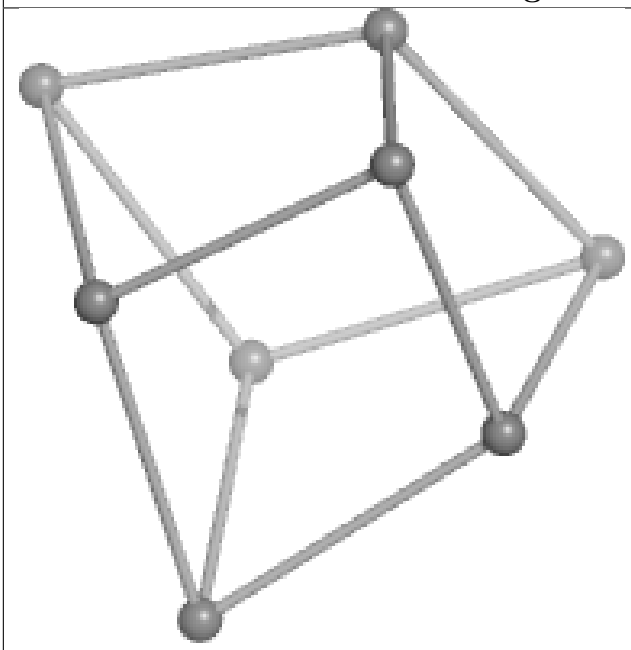


Torsions

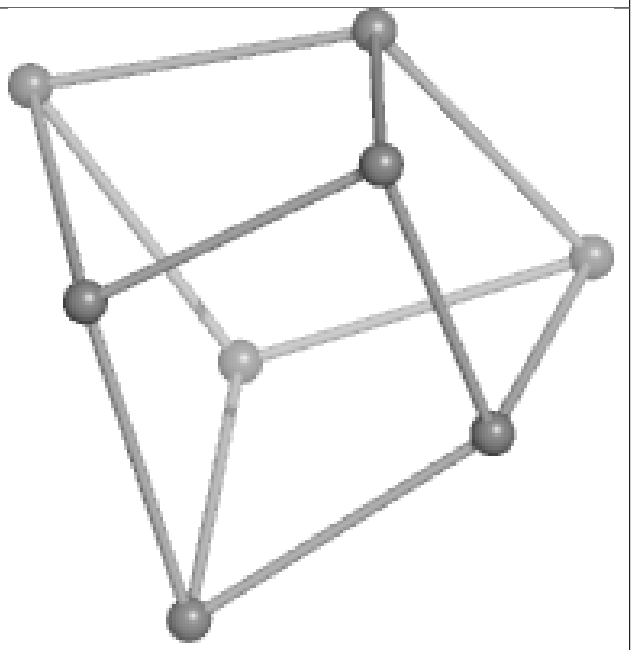


Rings

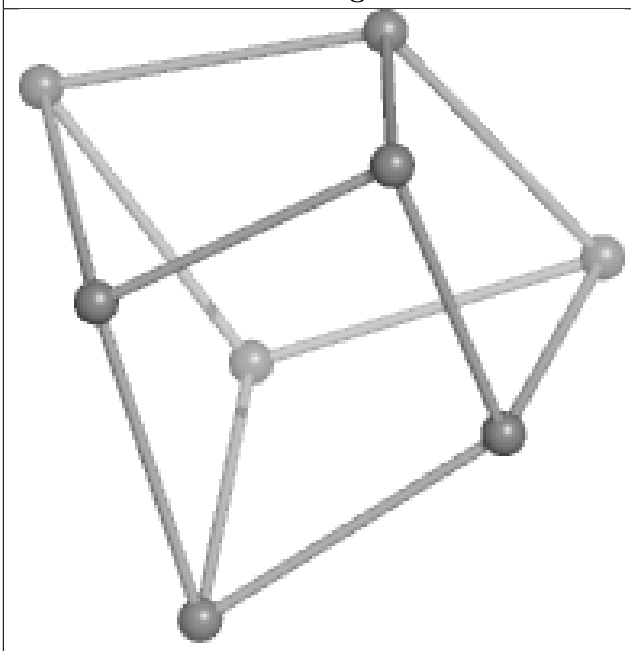
Ligand SF4 E 1103



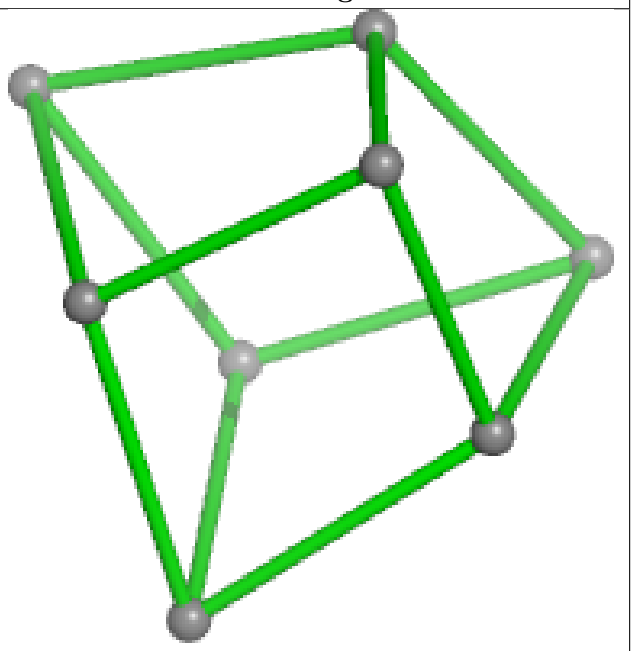
Bond lengths



Bond angles

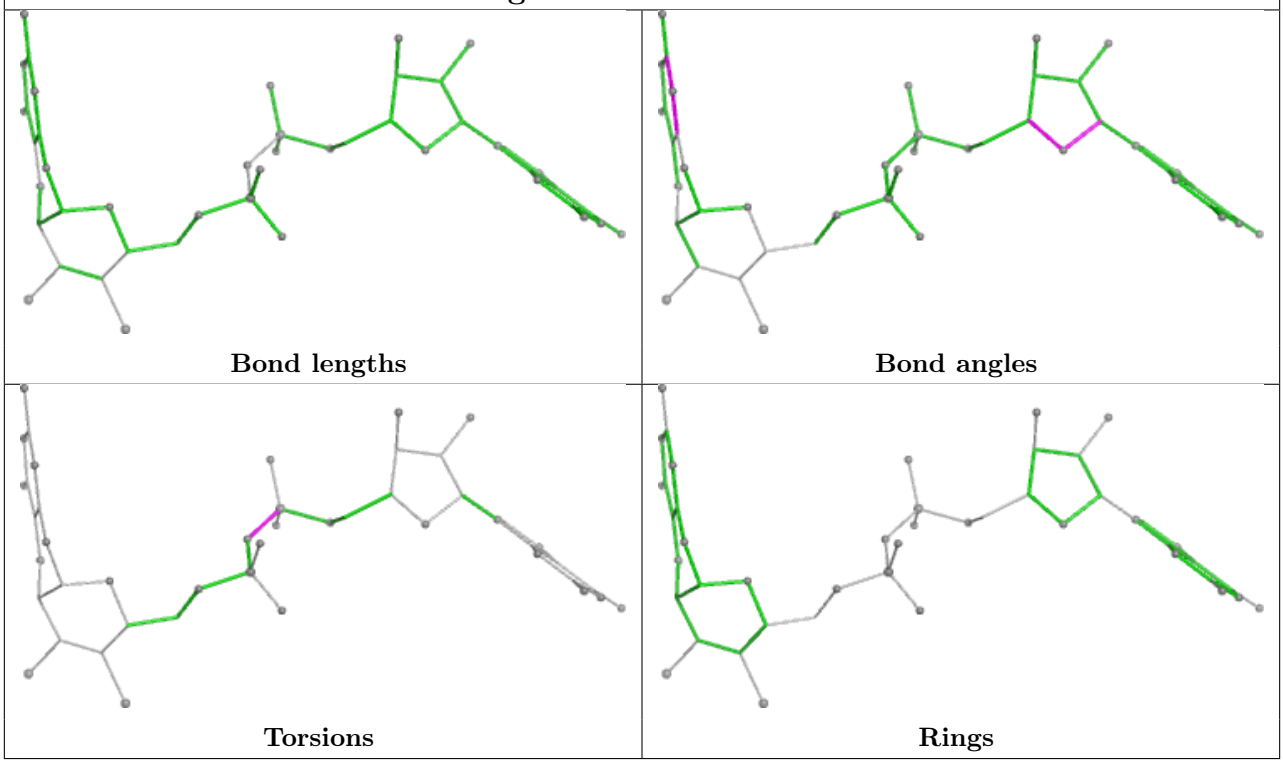


Torsions

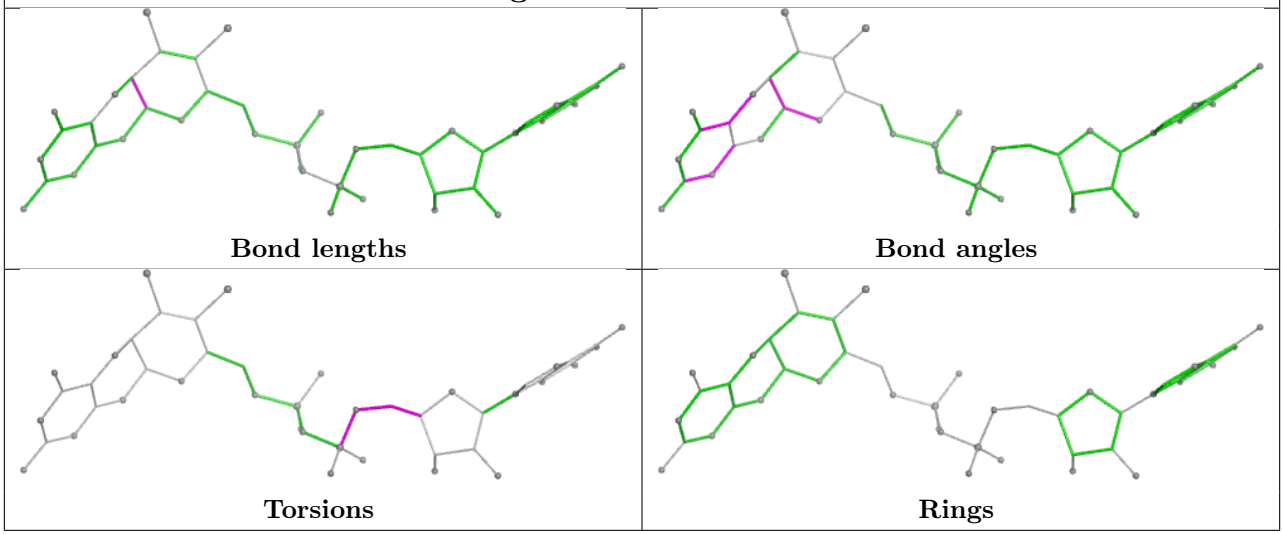


Rings

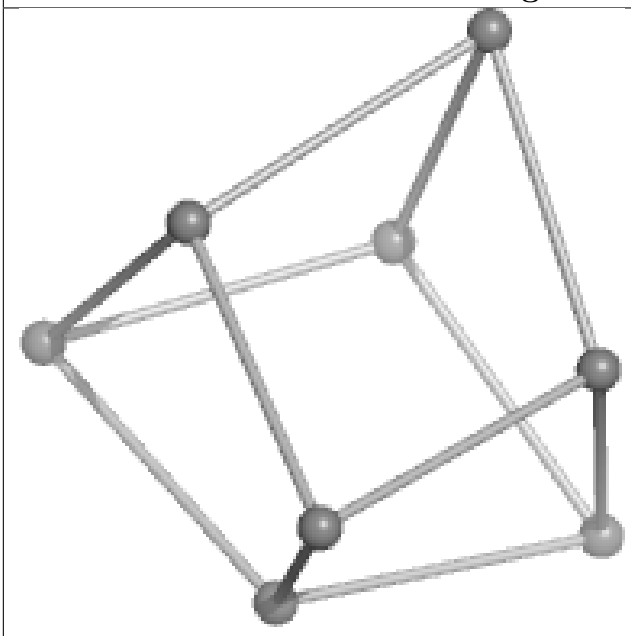
Ligand MGD A 1101



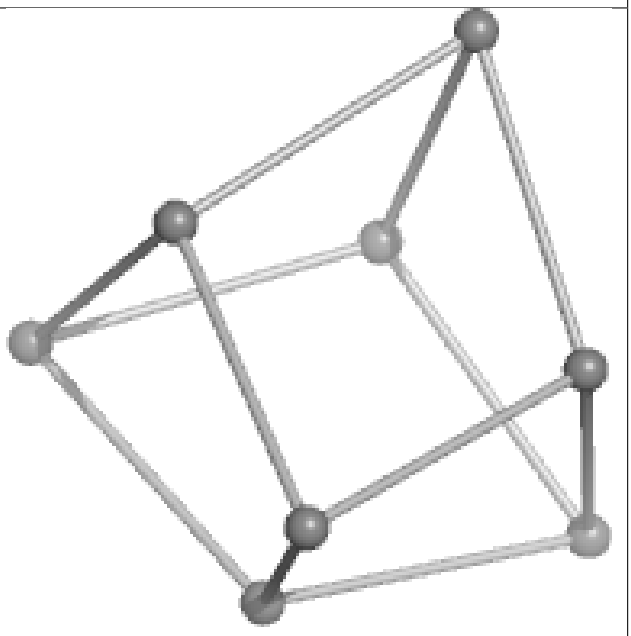
Ligand MGD E 1102



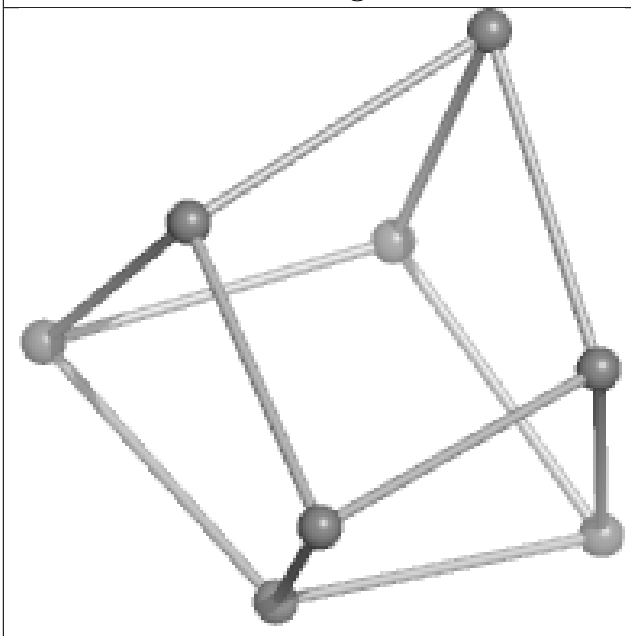
Ligand SF4 A 1103



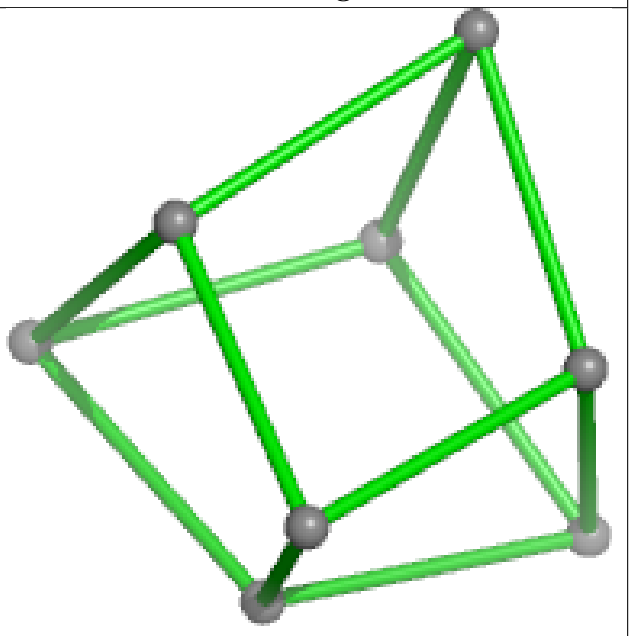
Bond lengths



Bond angles

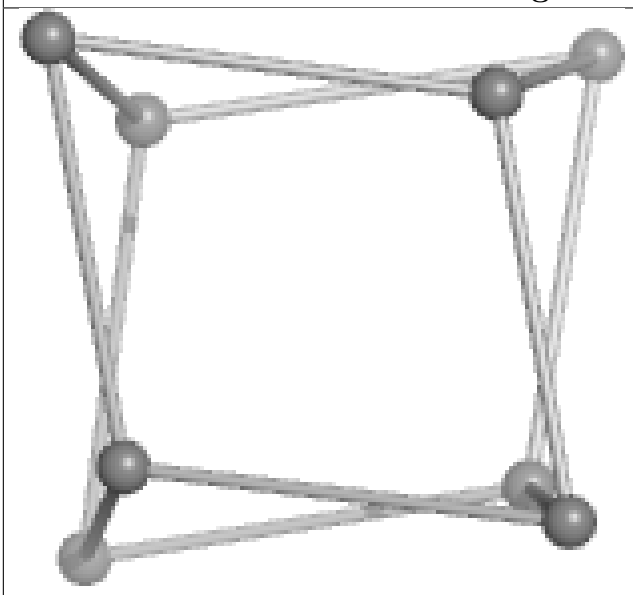


Torsions

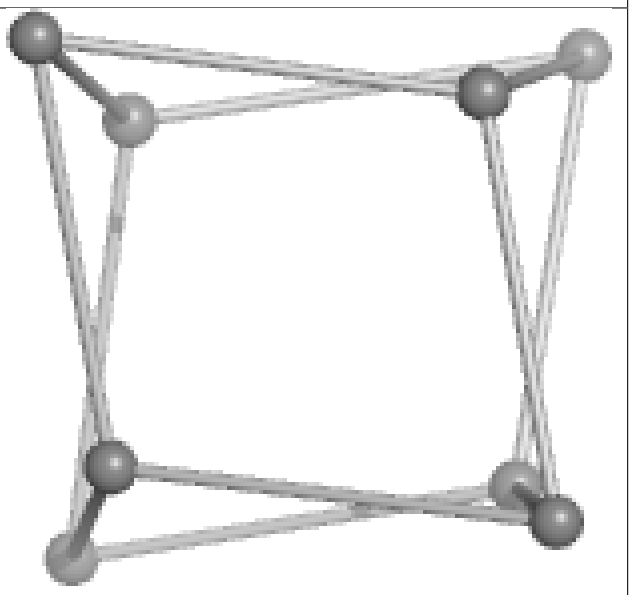


Rings

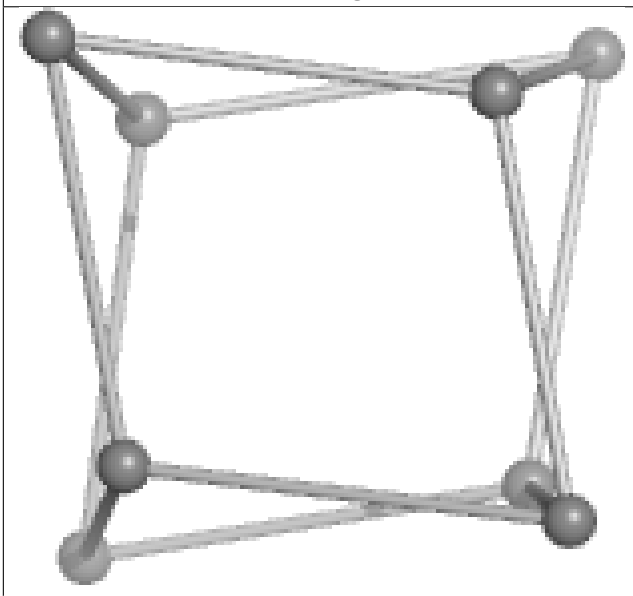
Ligand SF4 B 303



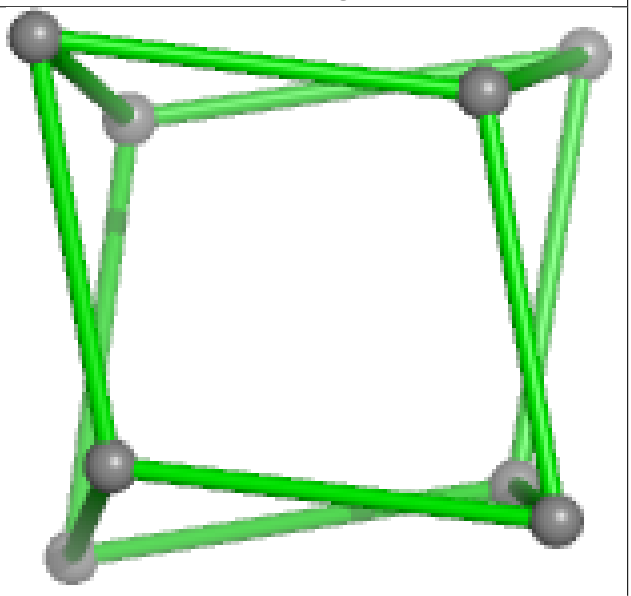
Bond lengths



Bond angles

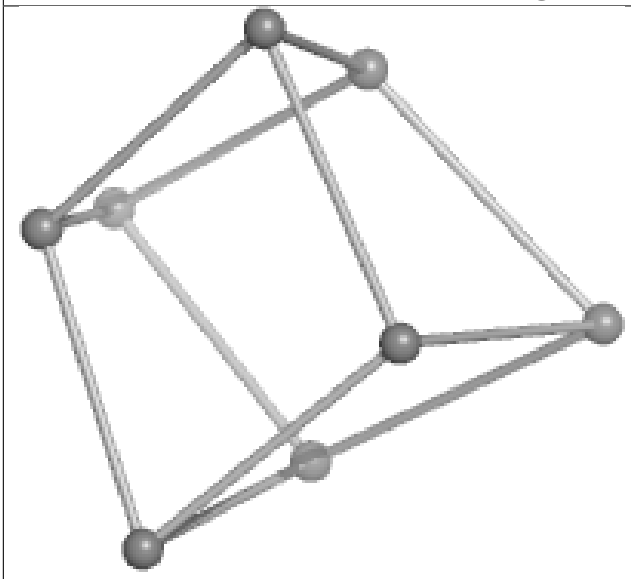


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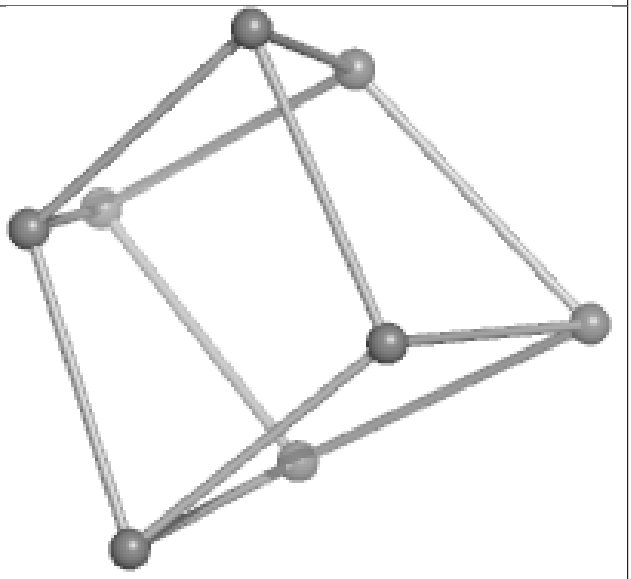


Rings

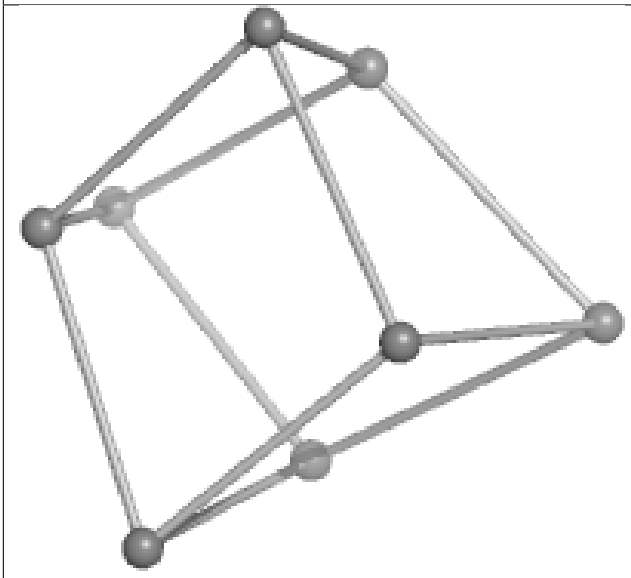
Ligand SF4 F 302



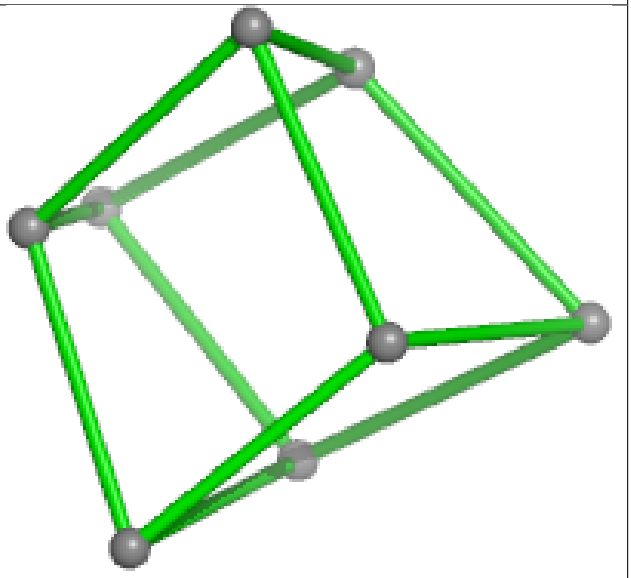
Bond lengths



Bond angles

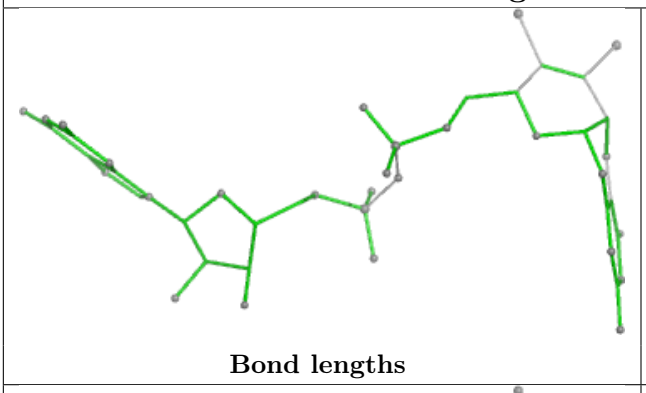


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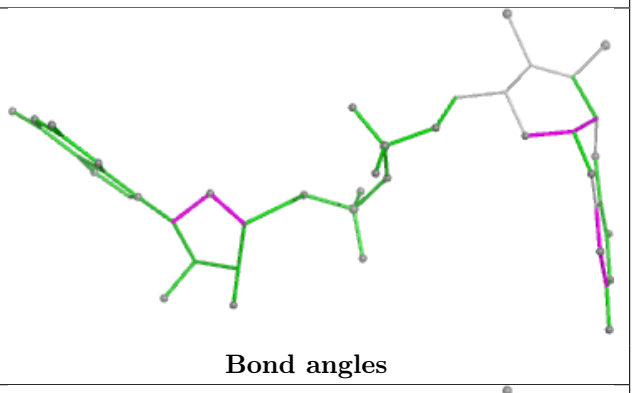


Rings

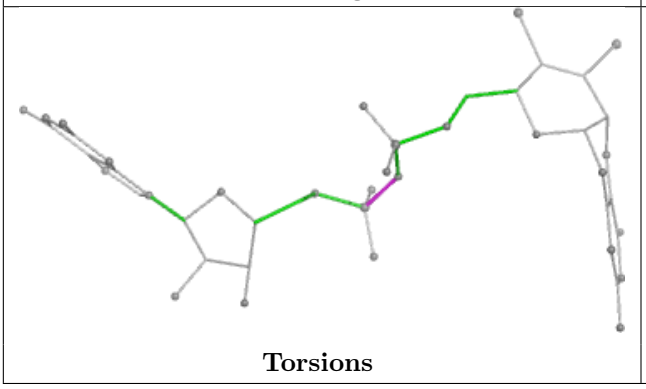
Ligand MGD E 1101



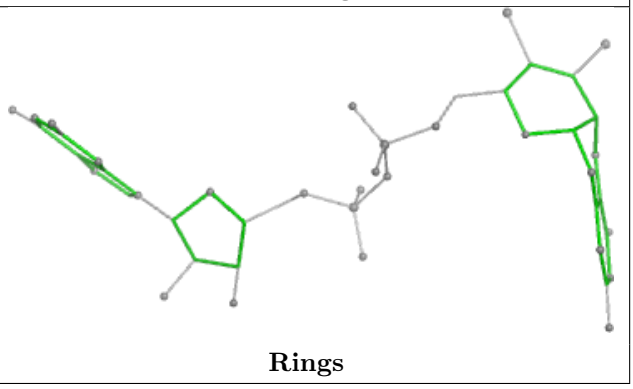
Bond lengths



Bond angles

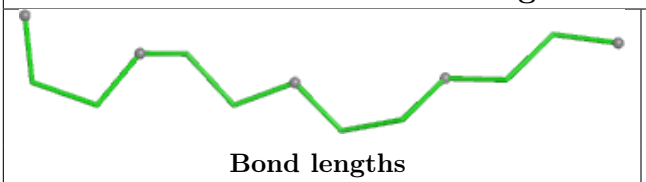


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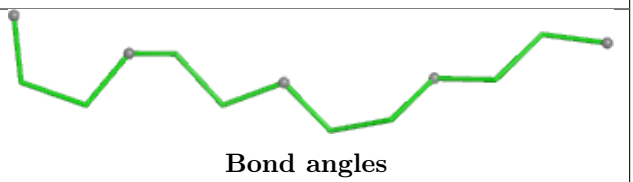


Rings

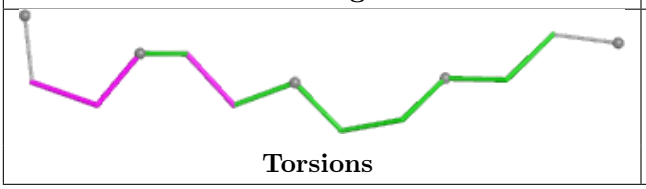
Ligand PG4 E 1104



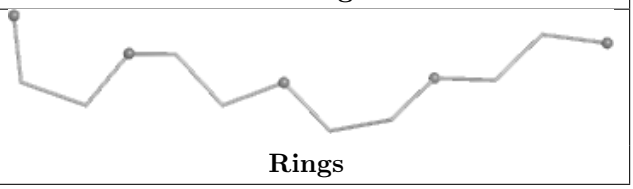
Bond lengths



Bond angles

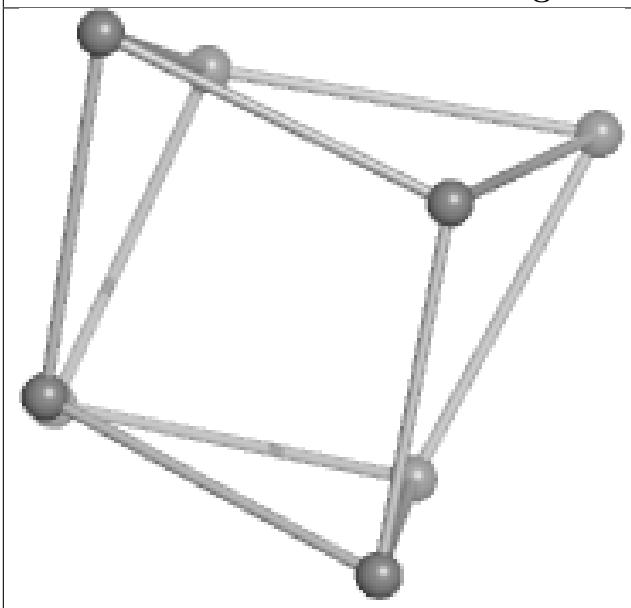


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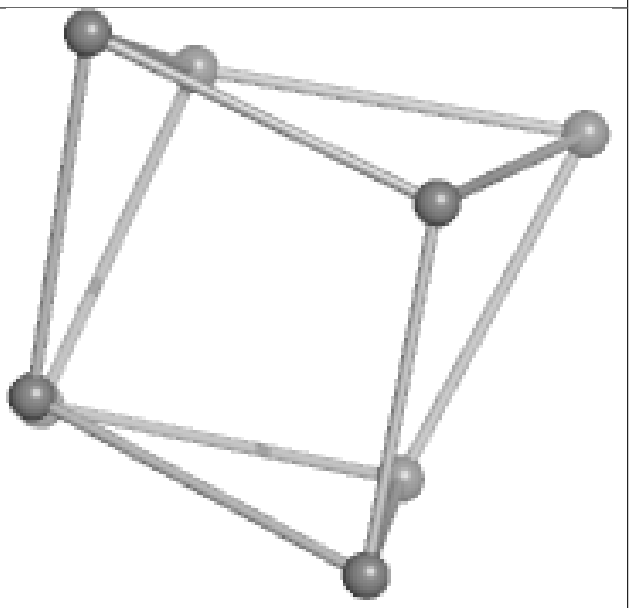


Rings

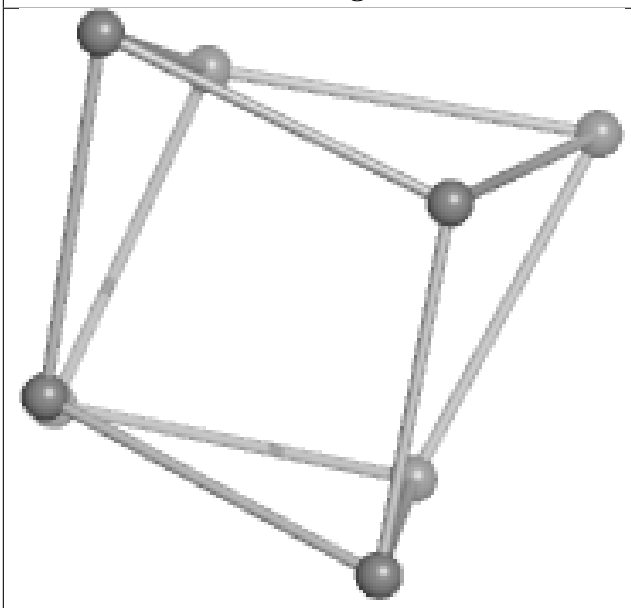
Ligand SF4 F 301



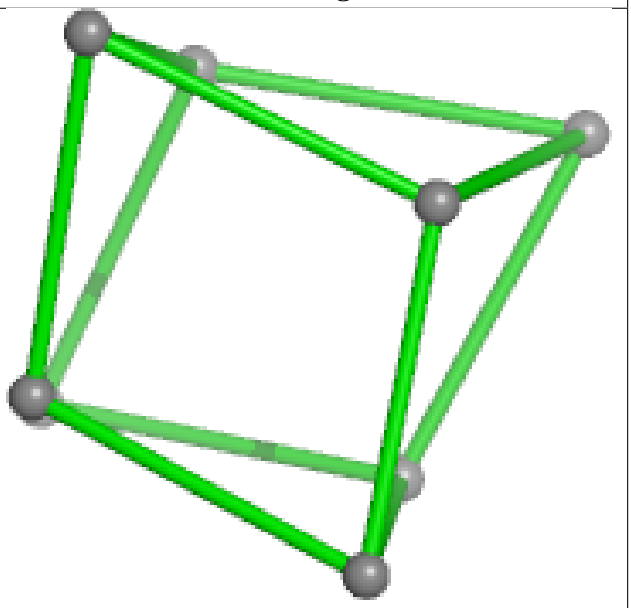
Bond lengths



Bond angles

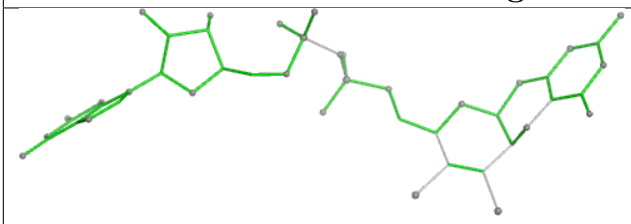


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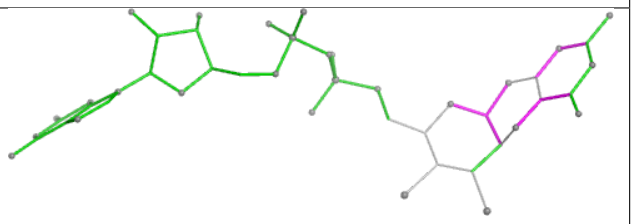


Rings

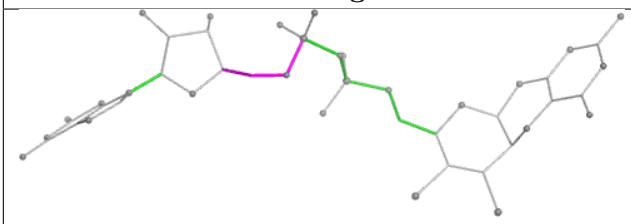
Ligand MGD A 1102



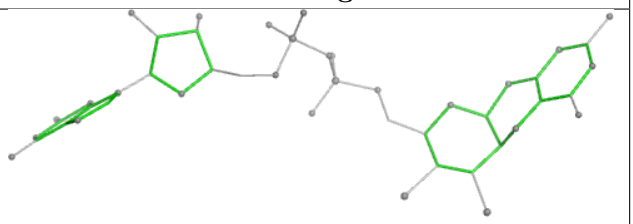
Bond lengths



Bond angles

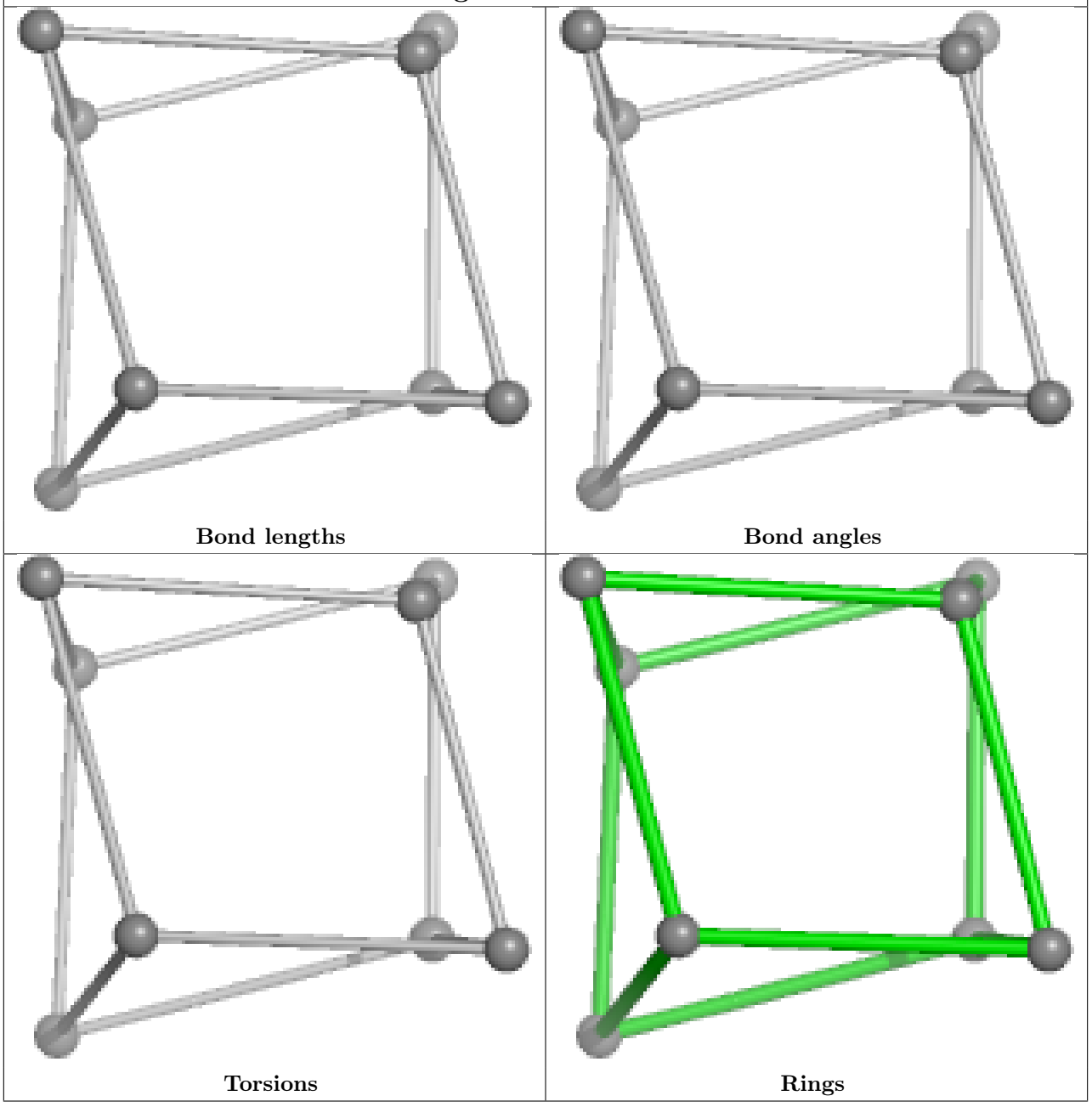


Torsions



Rings

Ligand SF4 F 303



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	962/1013 (94%)	0.05	18 (1%) 66 73	11, 22, 47, 88	2 (0%)
1	E	962/1013 (94%)	0.10	22 (2%) 61 67	13, 23, 48, 98	0
2	B	214/214 (100%)	0.47	9 (4%) 40 43	14, 27, 58, 72	0
2	F	214/214 (100%)	1.10	36 (16%) 4 4	16, 34, 64, 85	0
All	All	2352/2454 (95%)	0.20	85 (3%) 46 49	11, 24, 52, 98	2 (0%)

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	985	ALA	4.7
2	F	145	VAL	4.3
2	F	117	VAL	4.2
2	B	117	VAL	4.0
2	F	96	LYS	3.8
1	E	985	ALA	3.7
2	F	125	ILE	3.7
2	F	108	LEU	3.5
2	F	86	ALA	3.5
2	F	120	ALA	3.4
2	F	82	CYS	3.3
2	F	47	TYR	3.3
1	A	992	ILE	3.1
1	E	146	ALA	3.0
2	F	131	VAL	3.0
1	A	407	TRP	3.0
1	A	861	ALA	3.0
2	F	187	GLY	2.8
2	F	189	VAL	2.8
1	A	315	ASP	2.8
2	F	135	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
2	F	213	ALA	2.8
2	F	128	ILE	2.8
2	B	145	VAL	2.8
2	F	77	CYS	2.7
2	F	215	ALA	2.7
2	F	132	THR	2.7
2	B	112	VAL	2.7
2	F	116	SER	2.6
1	E	796	ASP	2.6
1	E	954	GLY	2.6
2	F	81	PRO	2.6
1	A	702	GLY	2.5
1	E	702	GLY	2.5
2	F	119	SER	2.5
2	B	98	GLU	2.5
1	E	407	TRP	2.5
1	E	339	ALA	2.5
1	E	1005	VAL	2.5
2	F	88	VAL	2.5
2	F	112	VAL	2.5
2	F	121	CYS	2.4
2	B	108	LEU	2.4
1	E	861	ALA	2.4
1	A	988	PRO	2.4
2	F	114	GLY	2.4
2	F	102	ALA	2.4
1	A	993	PRO	2.4
1	A	955	VAL	2.4
1	A	843	LEU	2.4
2	F	90	LEU	2.4
1	E	66	ASP	2.4
1	A	196	THR	2.3
1	E	679	ASN	2.3
1	E	701	GLY	2.3
1	E	1002	VAL	2.3
1	A	679	ASN	2.3
2	F	92	GLY	2.3
1	E	950	PHE	2.3
2	F	94	VAL	2.3
2	F	99	THR	2.3
2	F	105	PHE	2.2
2	B	88	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	131	VAL	2.2
2	F	103	VAL	2.2
1	E	196	THR	2.2
1	E	700	VAL	2.2
1	E	955	VAL	2.2
1	A	662	ALA	2.2
1	A	989	ASN	2.1
1	E	992	ILE	2.1
1	E	68	GLN	2.1
1	A	480	THR	2.1
2	B	94	VAL	2.1
1	E	869	LYS	2.1
1	A	796	ASP	2.1
2	F	130	PRO	2.1
1	A	324	ALA	2.1
1	A	701	GLY	2.1
2	F	78	VAL	2.1
2	B	47	TYR	2.0
2	F	84	GLY	2.0
1	E	478	ALA	2.0
1	E	698	THR	2.0
2	F	62	GLY	2.0

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

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

6.3 Carbohydrates (i)

There are no oligosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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
7	PEG	E	1108	7/7	0.71	0.24	47,54,61,74	0

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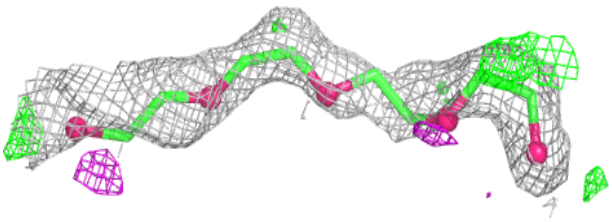
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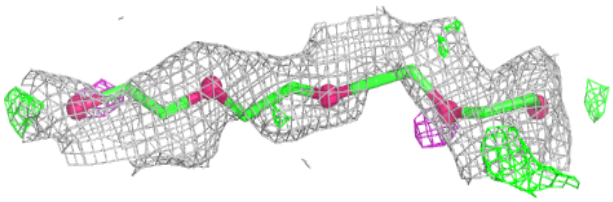
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
6	PG4	E	1104	13/13	0.81	0.16	28,32,44,53	0
6	PG4	A	1105	13/13	0.82	0.15	34,38,45,46	0
7	PEG	B	304	7/7	0.84	0.20	12,30,39,58	0
7	PEG	E	1107	7/7	0.87	0.14	34,39,45,48	0
7	PEG	A	1108	7/7	0.88	0.13	36,38,50,54	0
4	SF4	F	302	8/8	0.92	0.08	29,32,34,36	0
7	PEG	A	1107	7/7	0.93	0.12	37,39,51,59	0
7	PEG	E	1105	7/7	0.93	0.15	37,50,55,66	0
4	SF4	B	302	8/8	0.94	0.09	25,27,29,30	0
4	SF4	F	303	8/8	0.95	0.08	22,23,24,25	0
8	W	E	1109	1/1	0.95	0.05	17,17,17,17	0
5	GOL	A	1106	6/6	0.96	0.07	23,29,30,31	0
9	H2S	E	1110	1/1	0.96	0.06	18,18,18,18	0
3	MGD	E	1101	47/47	0.97	0.05	12,16,21,26	0
3	MGD	A	1102	47/47	0.97	0.07	15,18,23,34	0
5	GOL	A	1104	6/6	0.97	0.06	15,17,20,22	0
8	W	A	1109	1/1	0.97	0.04	16,16,16,16	0
4	SF4	F	301	8/8	0.97	0.06	18,18,19,20	0
5	GOL	E	1106	6/6	0.97	0.06	18,21,26,26	0
4	SF4	B	303	8/8	0.98	0.07	19,21,21,23	0
3	MGD	E	1102	47/47	0.98	0.06	14,18,22,27	0
3	MGD	A	1101	47/47	0.98	0.04	9,14,17,18	0
4	SF4	E	1103	8/8	0.99	0.04	13,14,15,15	0
4	SF4	A	1103	8/8	0.99	0.04	13,13,14,14	0
9	H2S	A	1110	1/1	0.99	0.03	15,15,15,15	0
4	SF4	B	301	8/8	0.99	0.04	15,15,17,17	0

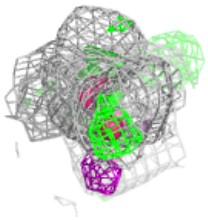
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PG4 E 1104:

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

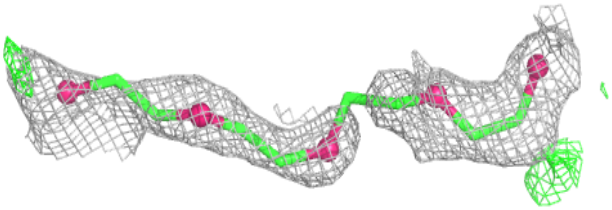


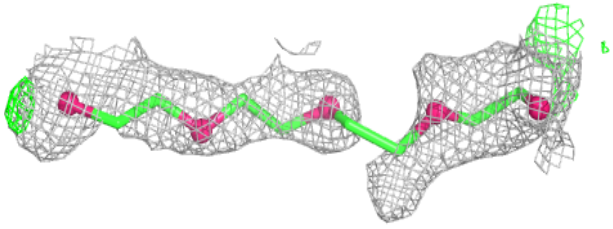


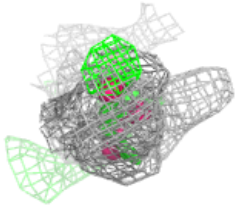


Electron density around PG4 A 1105:

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

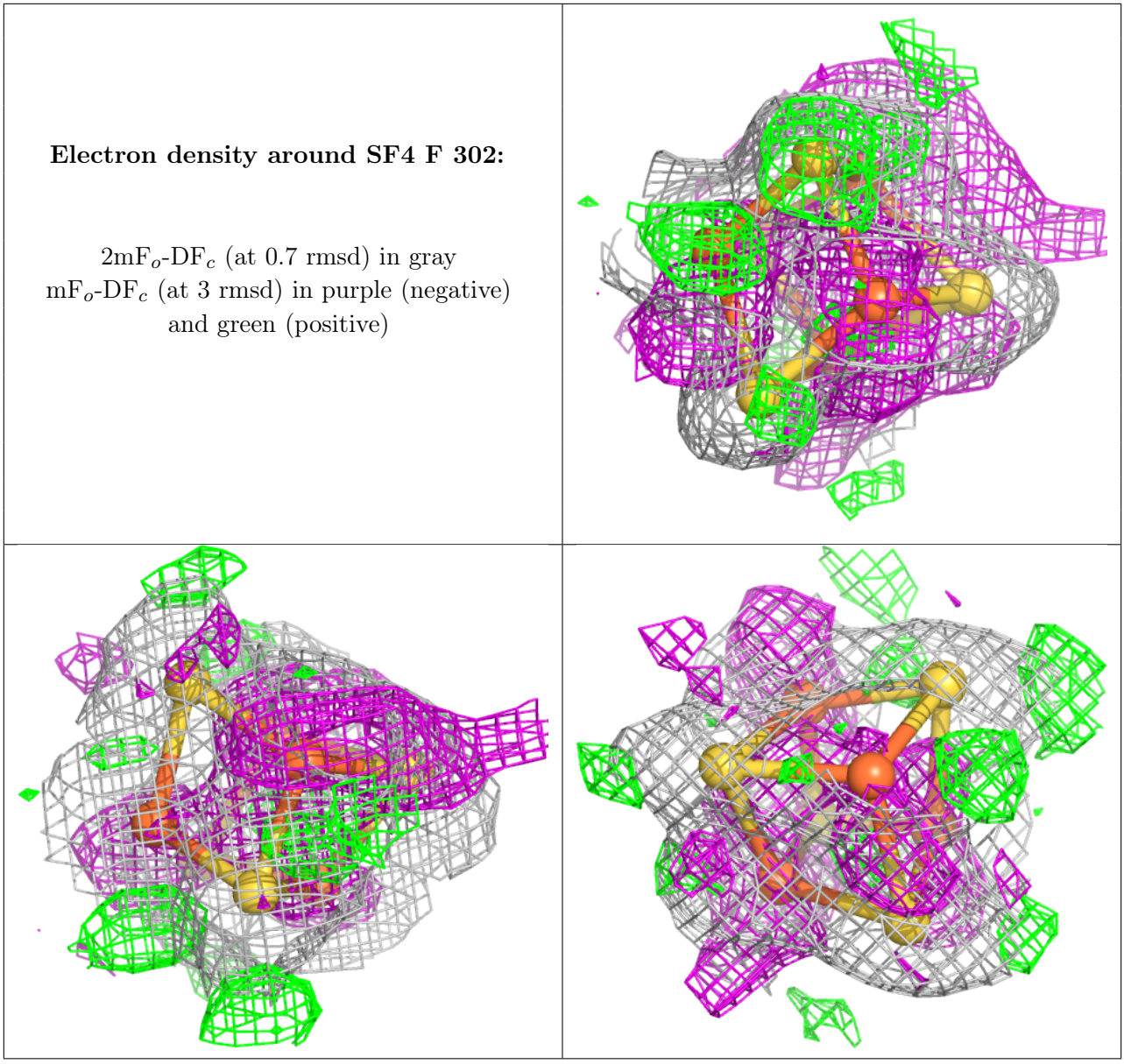






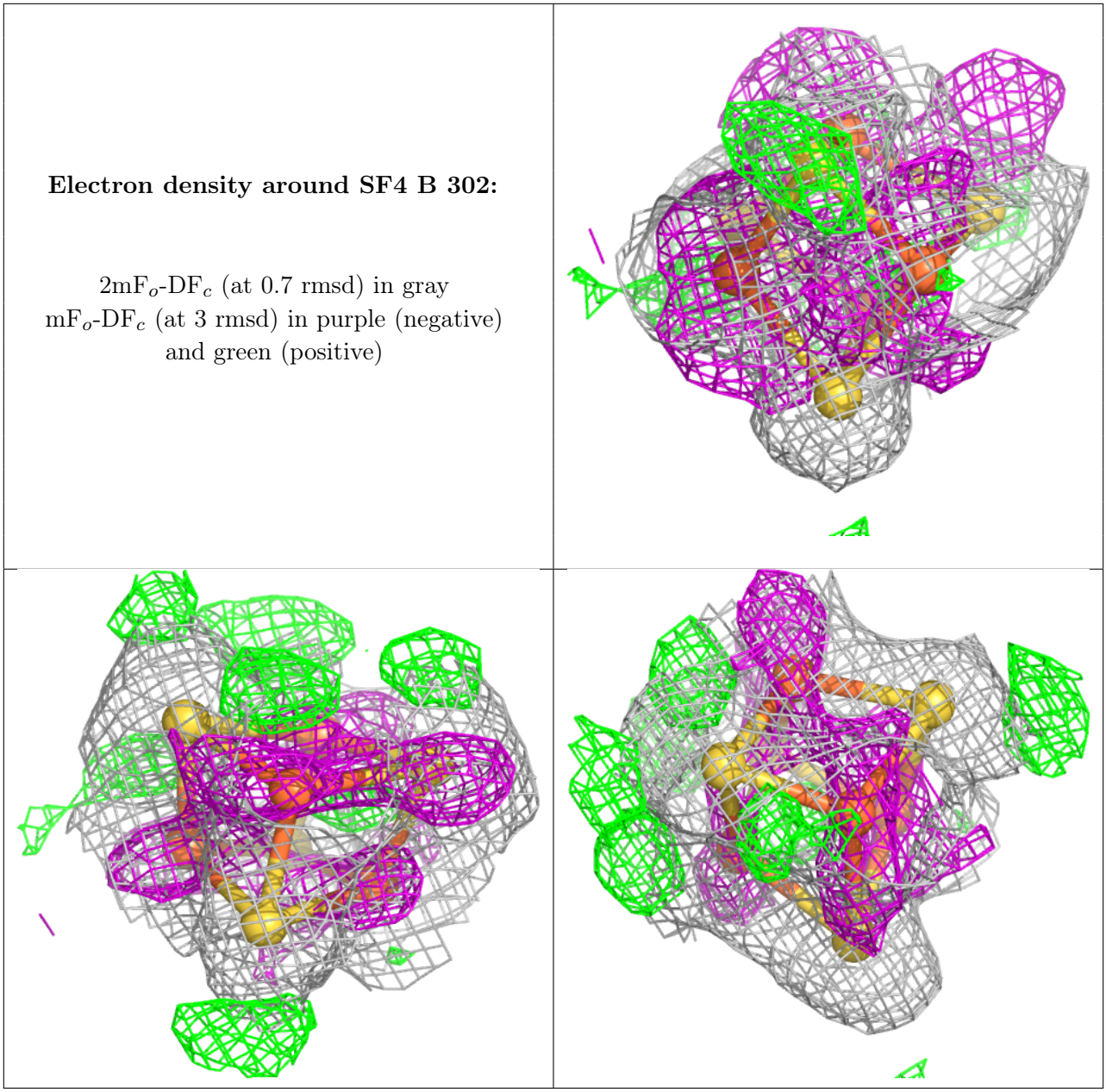
Electron density around SF4 F 302:

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



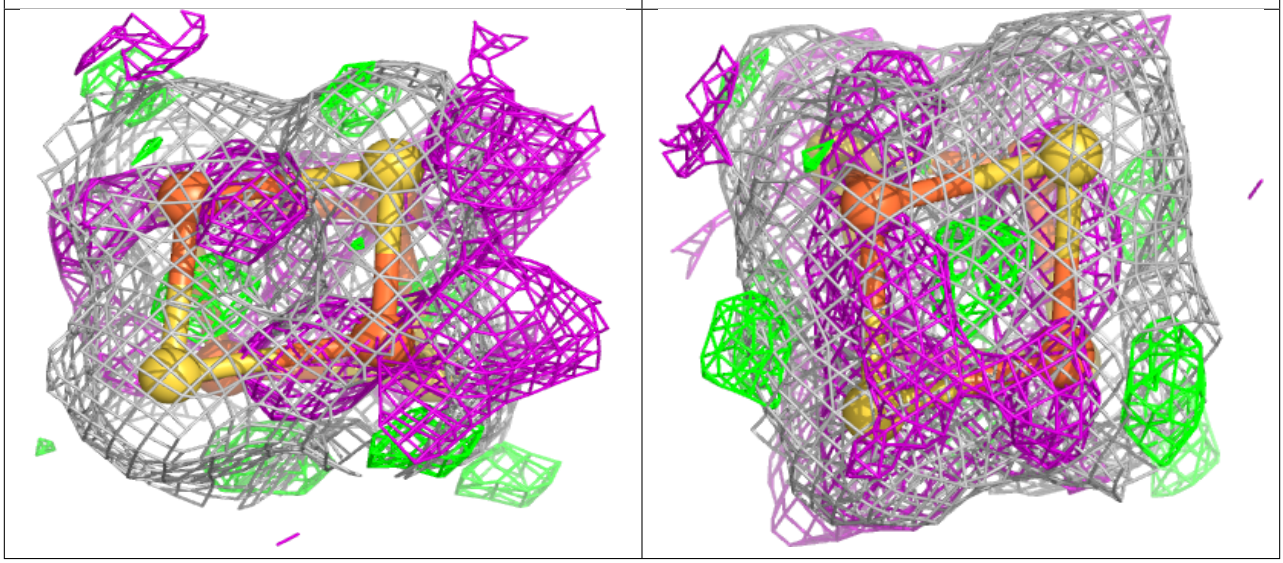
Electron density around SF4 B 302:

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



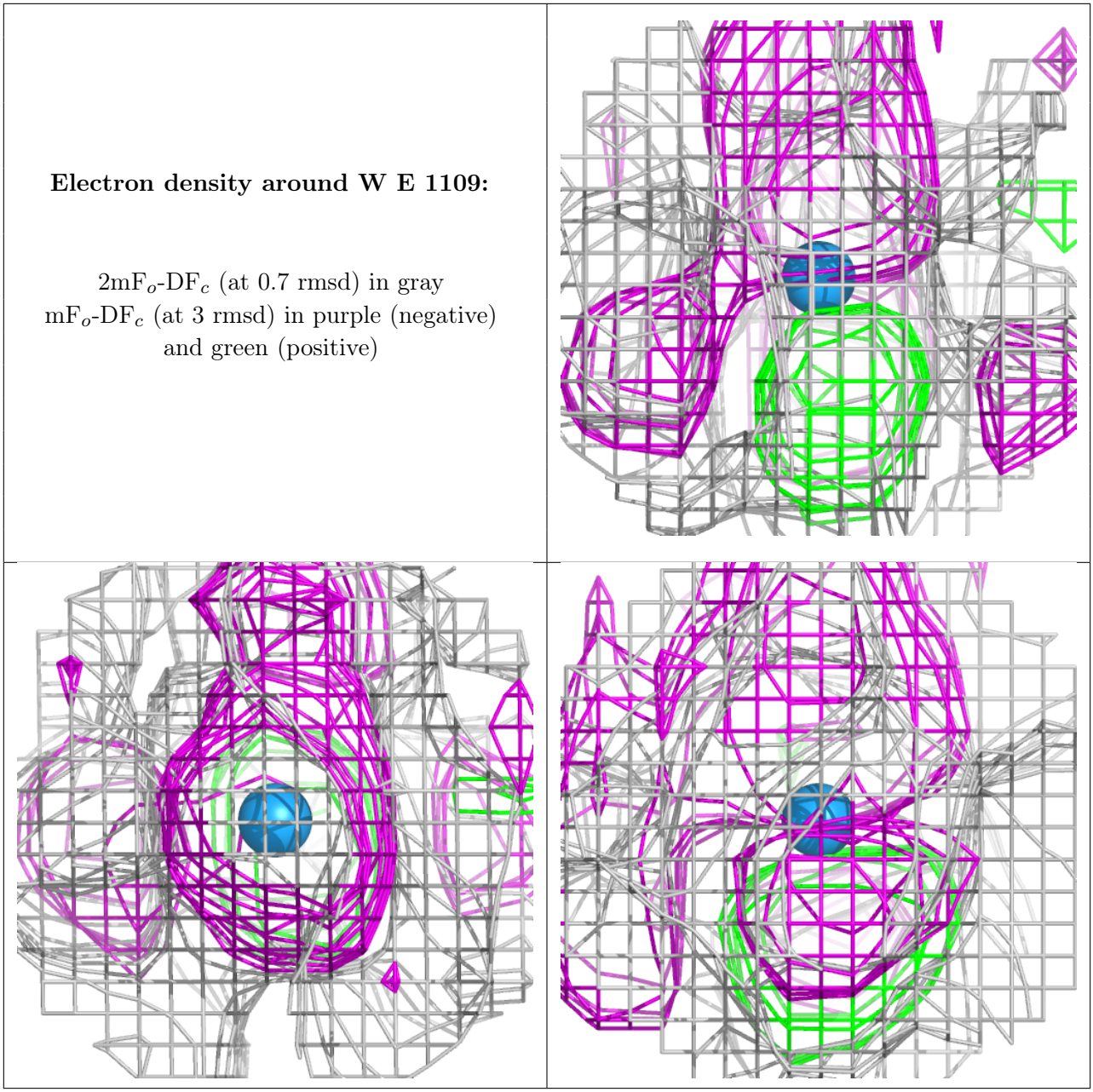
Electron density around SF4 F 303:

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



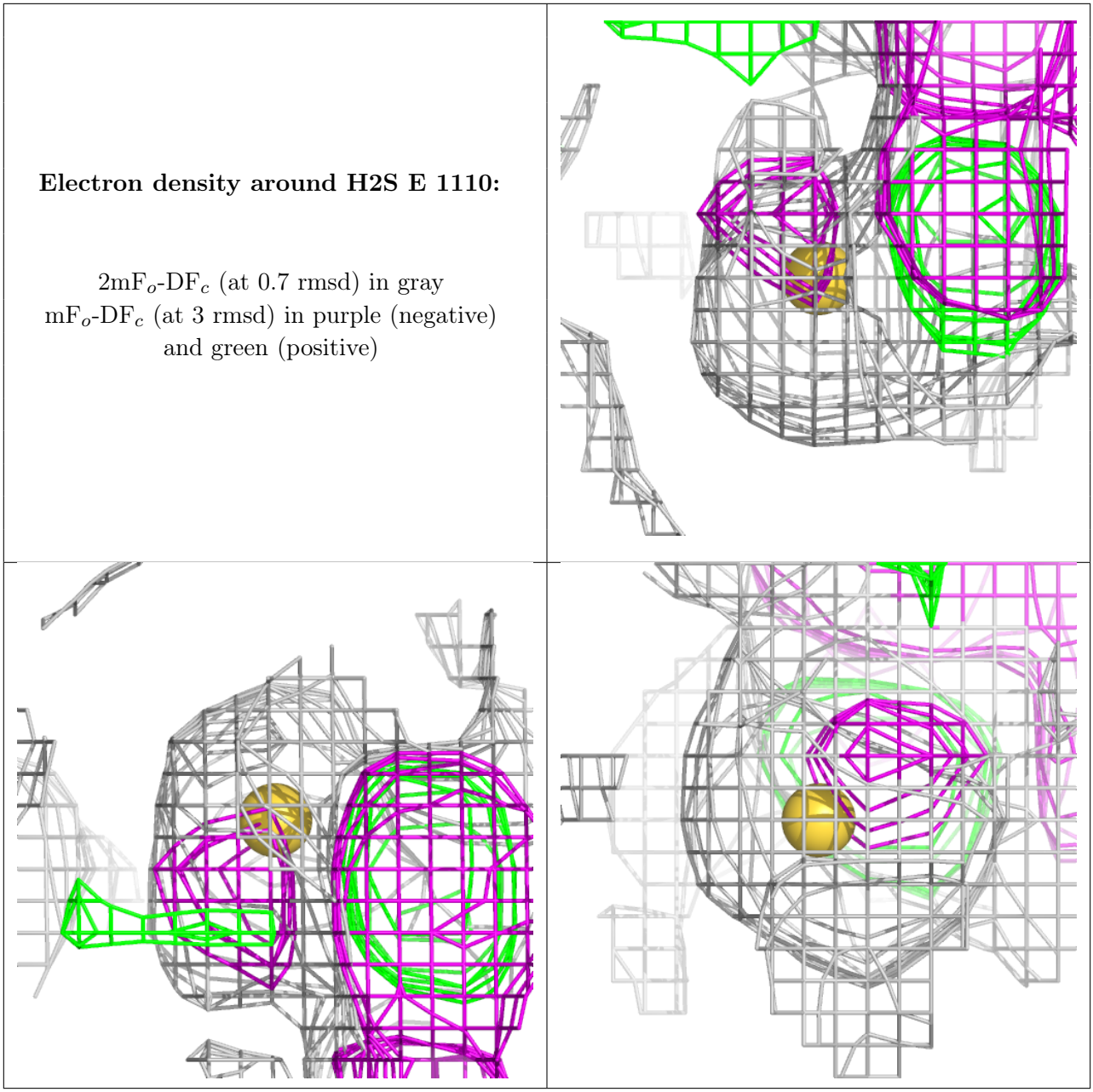
Electron density around W E 1109:

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



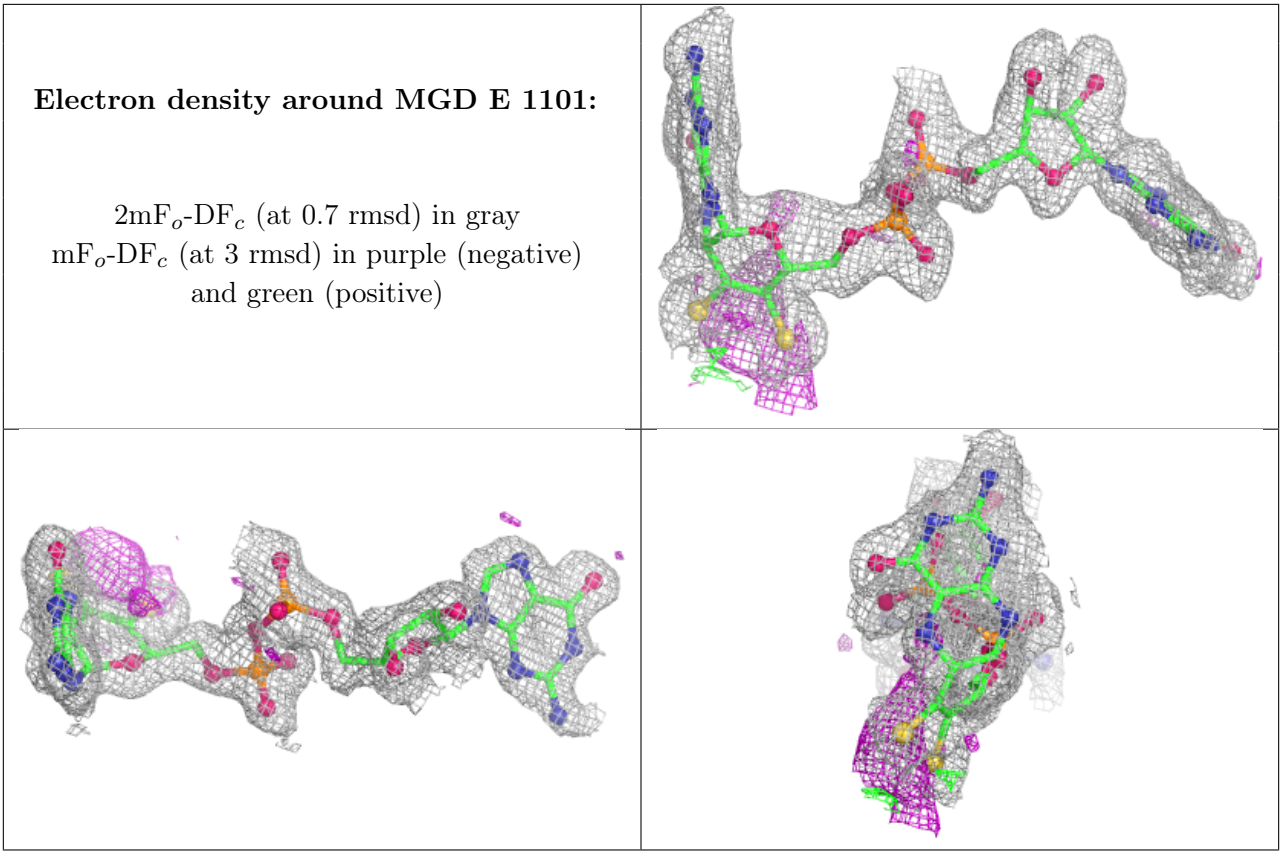
Electron density around H2S E 1110:

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



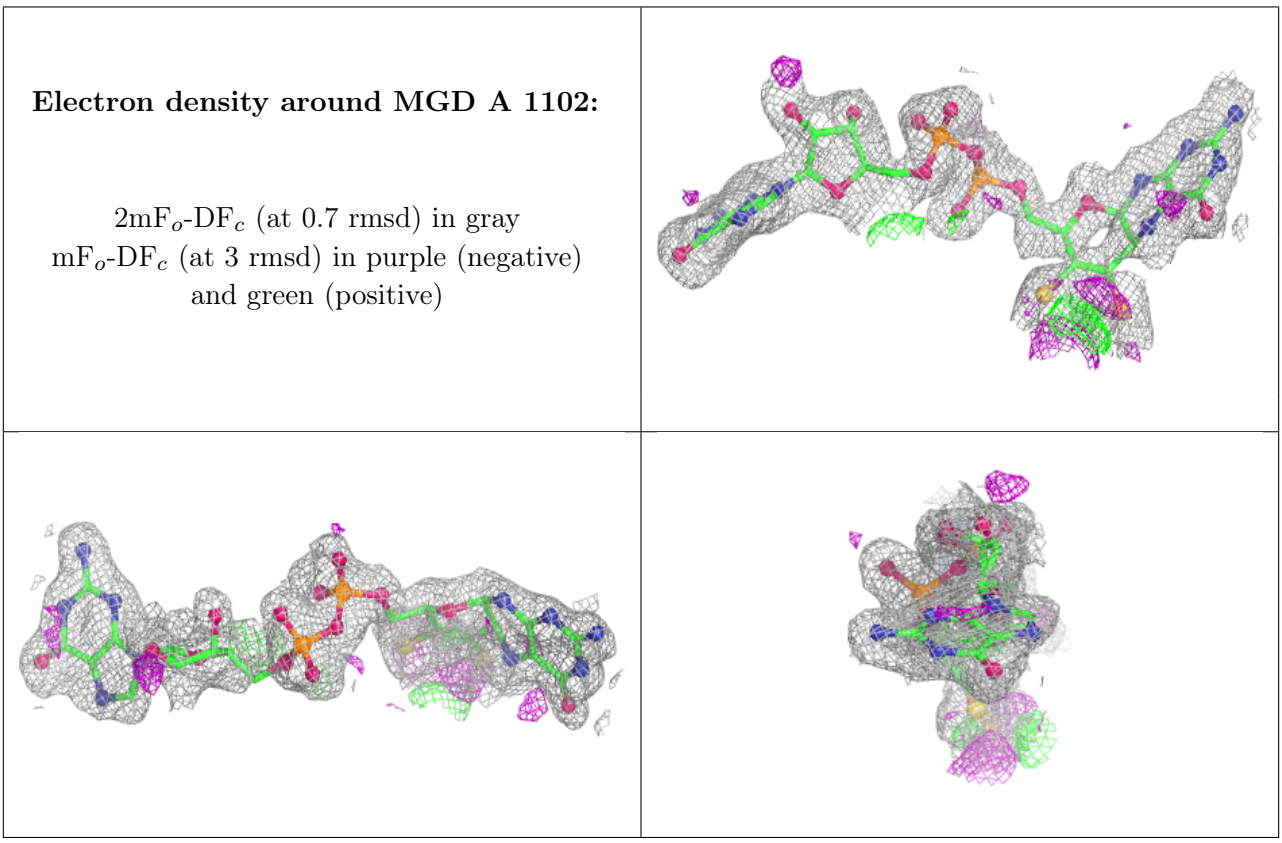
Electron density around MGD E 1101:

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



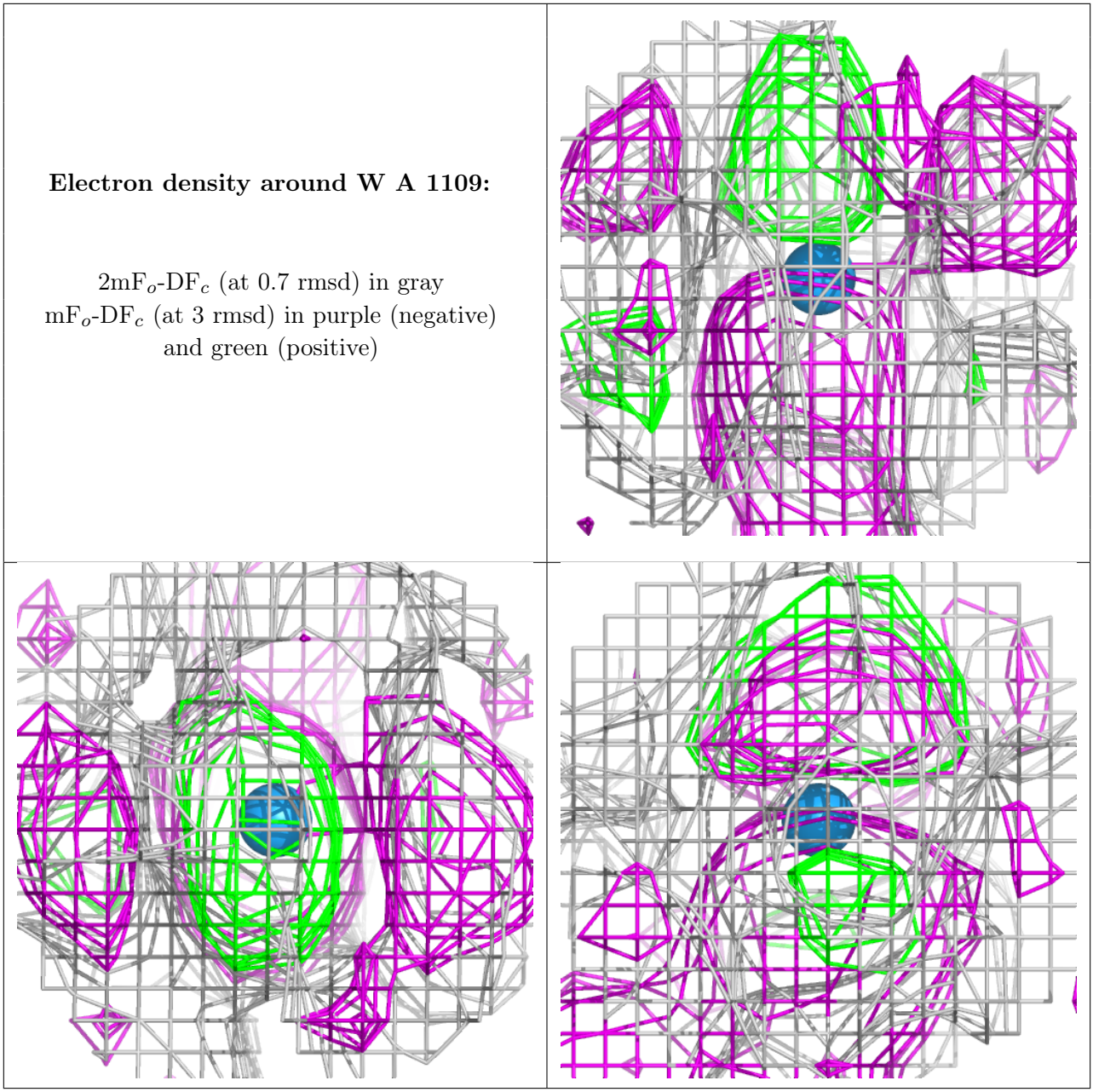
Electron density around MGD A 1102:

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



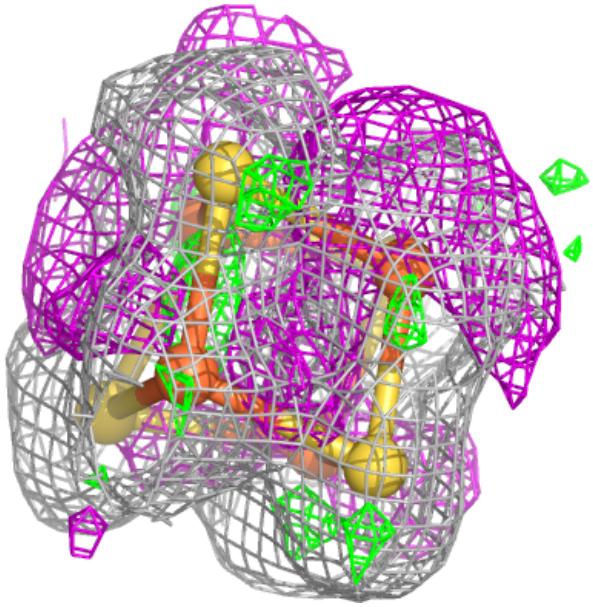
Electron density around W A 1109:

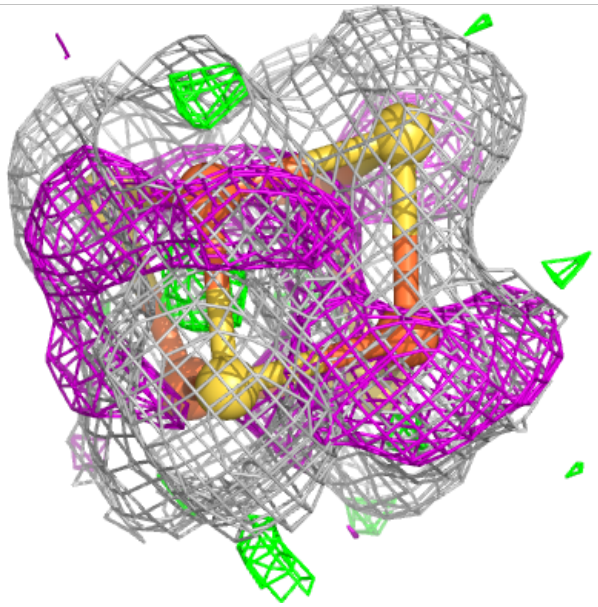
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

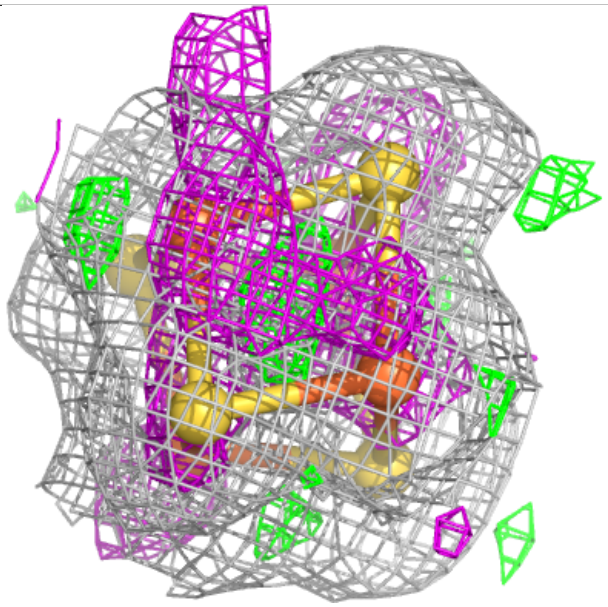


Electron density around SF4 F 301:

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

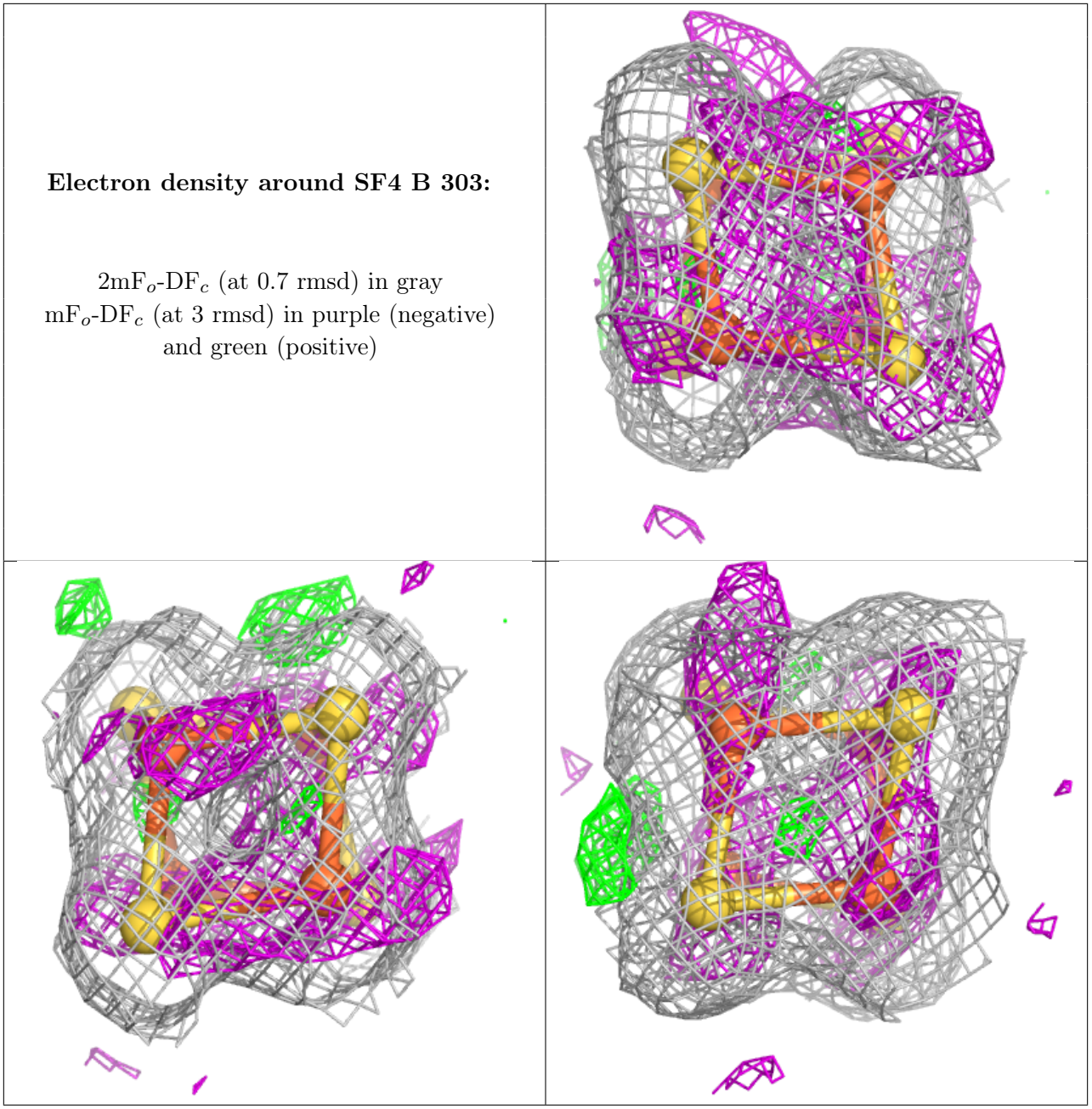






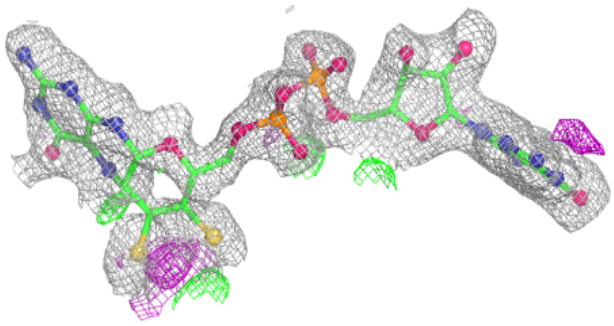
Electron density around SF4 B 303:

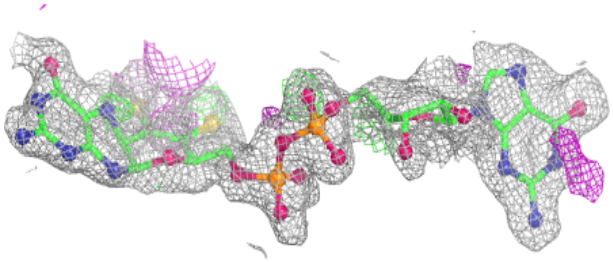
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

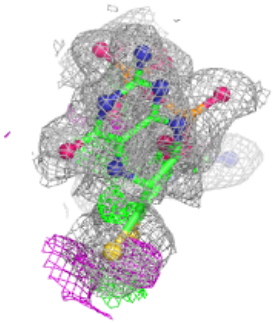


Electron density around MGD E 1102:

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

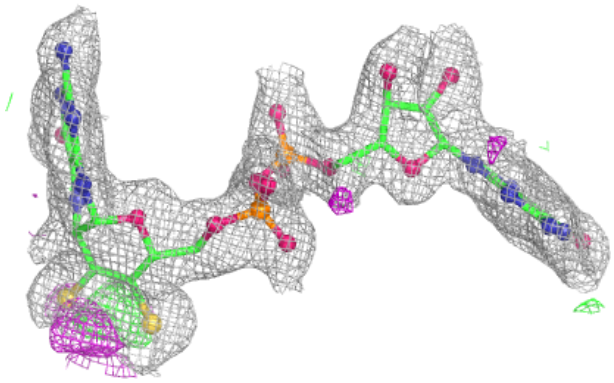


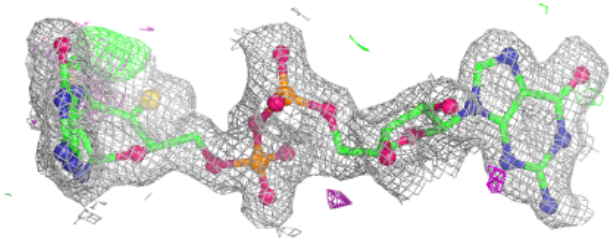


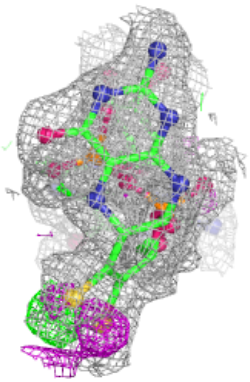


Electron density around MGD A 1101:

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

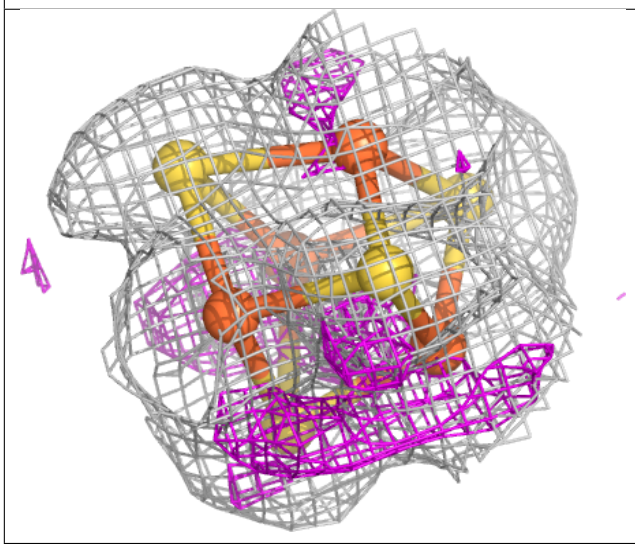


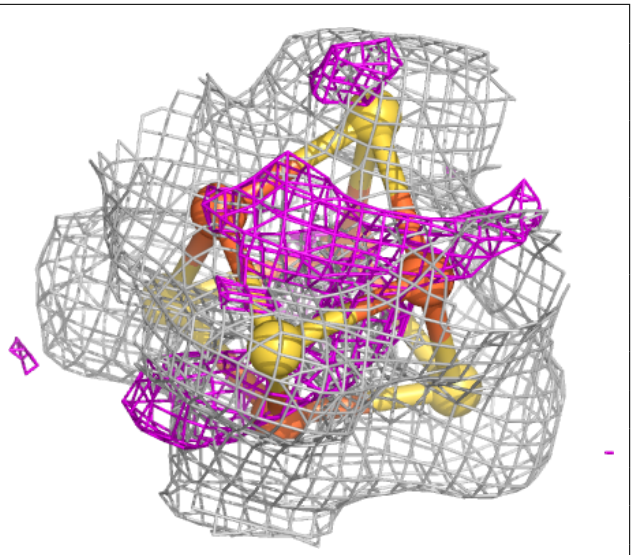


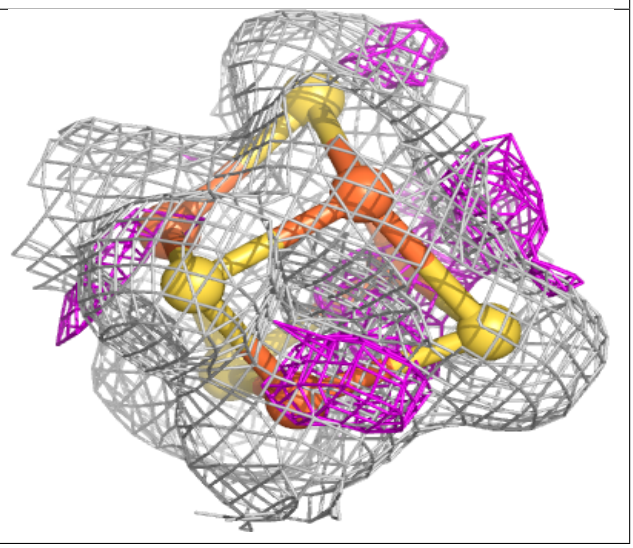


Electron density around SF4 E 1103:

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

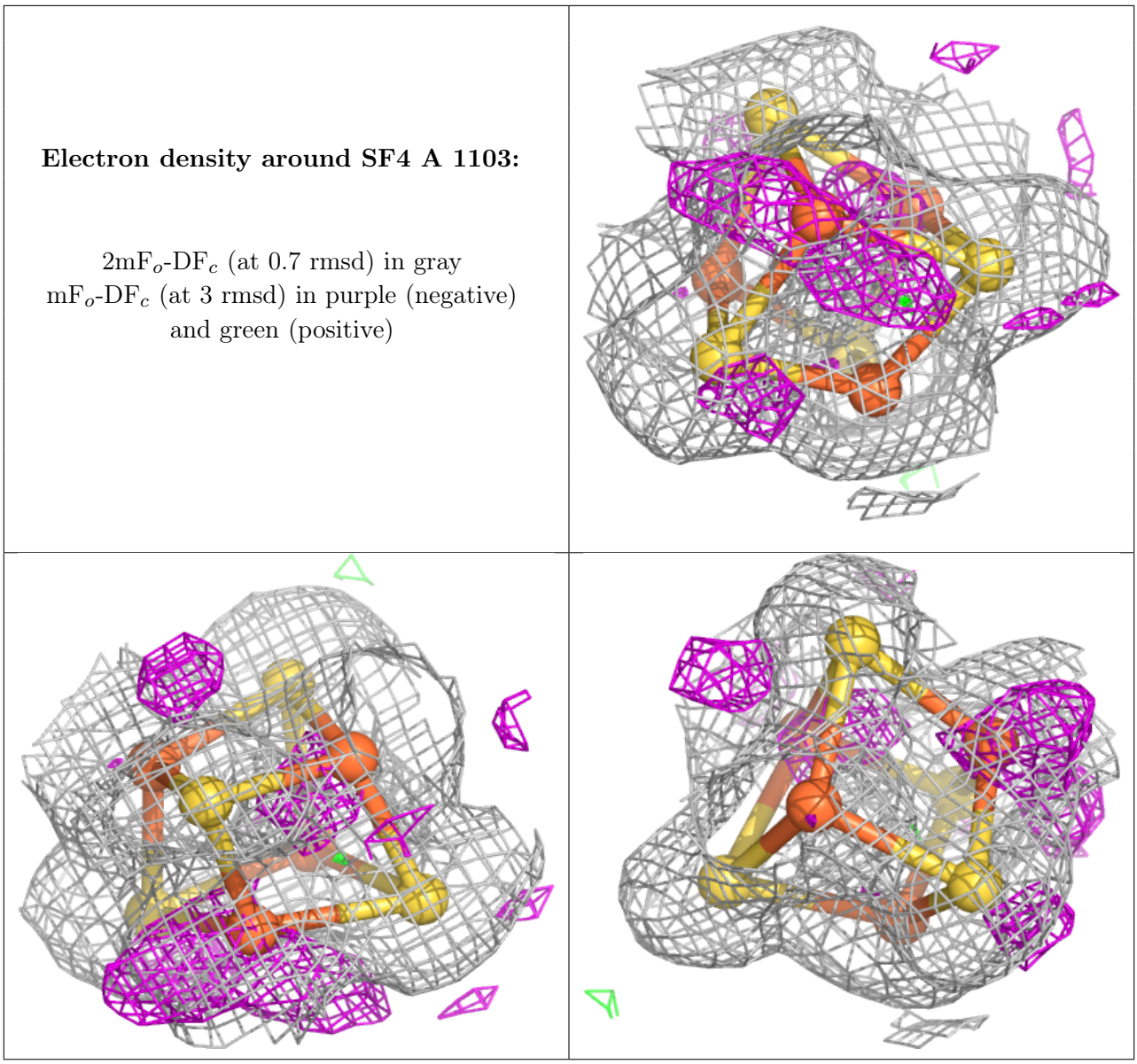






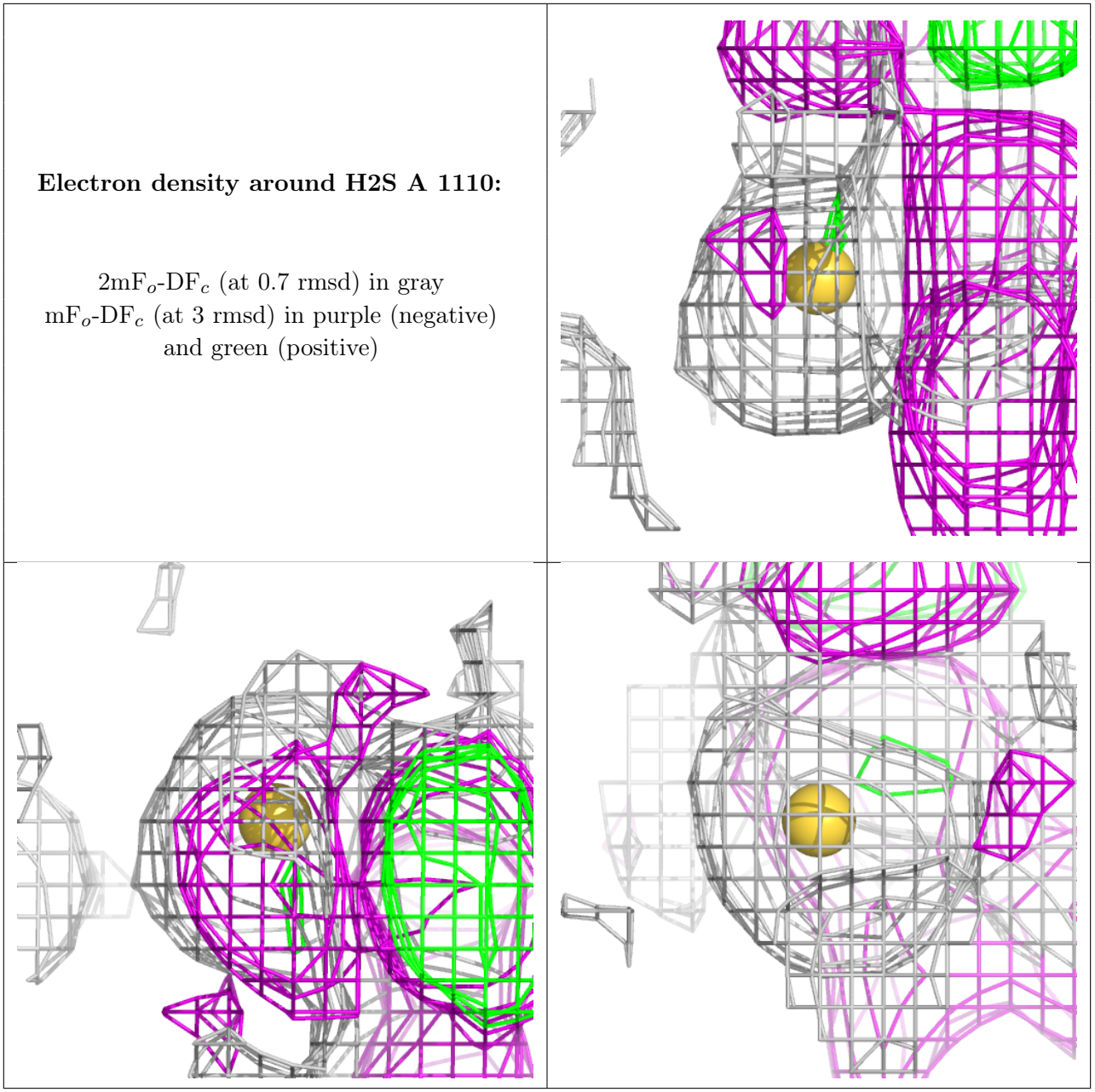
Electron density around SF4 A 1103:

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



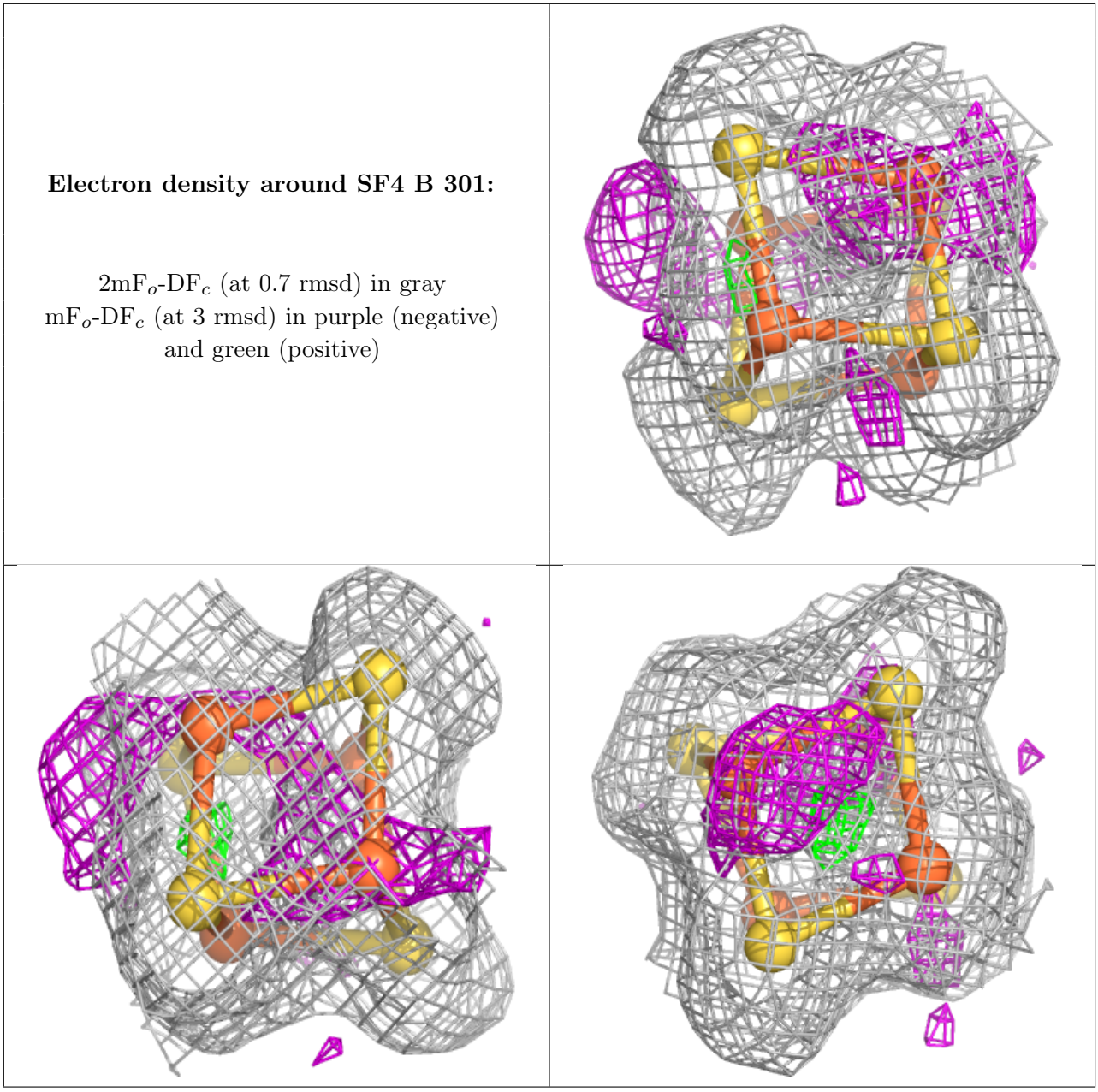
Electron density around H2S A 1110:

$2mF_o$-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 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 (i)

There are no such residues in this entry.