



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:31 AM UTC

PDB ID : 8CM6 / pdb_00008cm6
Title : W-formate dehydrogenase C872A from *Desulfovibrio vulgaris* - with Formamide
Authors : Vilela-Alves, G.; Mota, C.; Oliveira, A.R.; Manuel, R.R.; Pereira, I.C.; Romao, M.J.
Deposited on : 2023-02-17
Resolution : 1.42 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 Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
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	:	9.0.010 (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

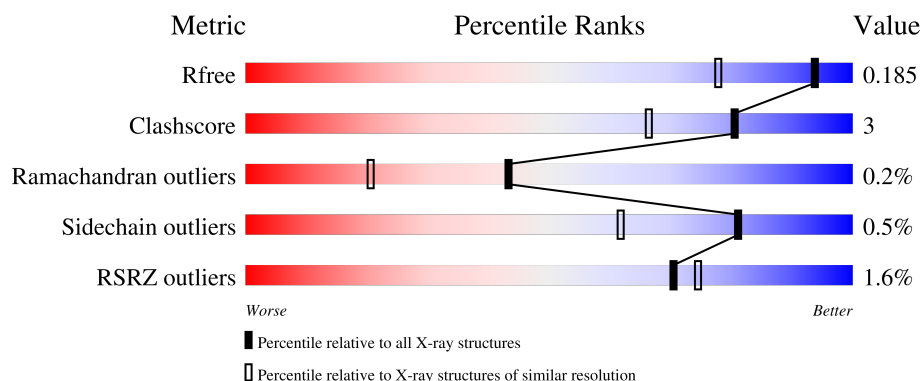
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.42 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	4041 (1.44-1.40)
Clashscore	190562	4154 (1.44-1.40)
Ramachandran outliers	187476	4083 (1.44-1.40)
Sidechain outliers	187428	4082 (1.44-1.40)
RSRZ outliers	180081	4039 (1.44-1.40)

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	1013	% 90% 6% .
1	C	1013	2% 89% 6% .
2	B	215	% 93% 6%
2	D	215	% 93% 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
10	PEG	A	1116	-	-	X	-
8	GOL	A	1115	-	-	X	-

2 Entry composition

There are 13 unique types of molecules in this entry. The entry contains 20783 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	972	Total	C	N	O	S	Se	0	8	0
			7658	4886	1332	1397	42	1			
1	C	971	Total	C	N	O	S	Se	0	3	0
			7623	4860	1326	1394	42	1			

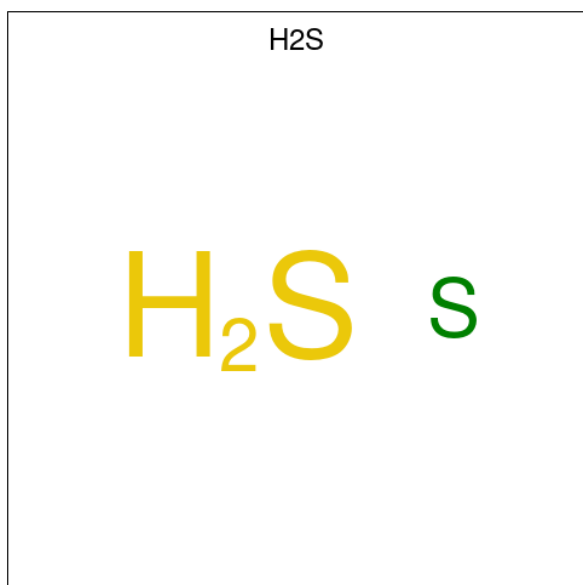
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	872	ALA	CYS	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
C	872	ALA	CYS	engineered mutation	UNP Q72EJ1
C	1006	TRP	-	expression tag	UNP Q72EJ1
C	1007	SER	-	expression tag	UNP Q72EJ1
C	1008	HIS	-	expression tag	UNP Q72EJ1
C	1009	PRO	-	expression tag	UNP Q72EJ1
C	1010	GLN	-	expression tag	UNP Q72EJ1
C	1011	PHE	-	expression tag	UNP Q72EJ1
C	1012	GLU	-	expression tag	UNP Q72EJ1
C	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	4	0
			1682	1054	292	320	16			
2	D	214	Total	C	N	O	S	0	0	0
			1664	1041	291	316	16			

- Molecule 3 is HYDROSULFURIC ACID (CCD ID: H2S) (formula: H₂S) (labeled as "Ligand of Interest" by depositor).

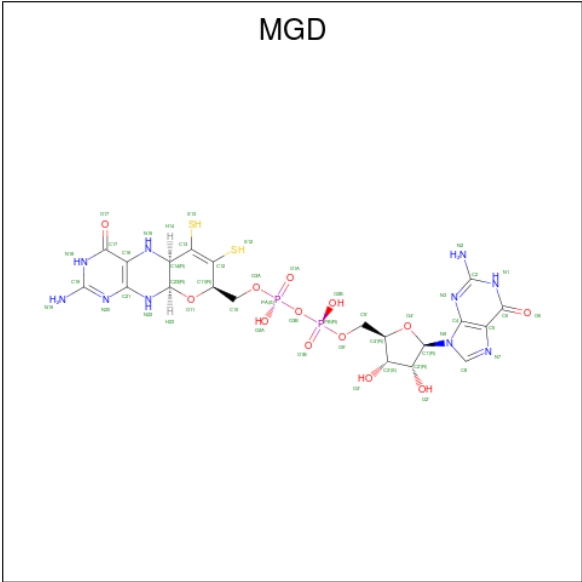


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	S	0	0
			1	1		
3	C	1	Total	S	0	0
			1	1		

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

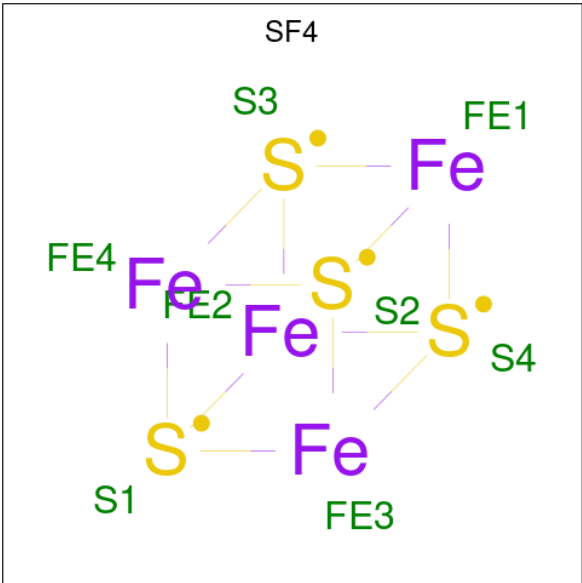
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	W	0	0
			1	1		
4	C	1	Total	W	0	0
			1	1		

- Molecule 5 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: C₂₀H₂₆N₁₀O₁₃P₂S₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 6 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄) (labeled as "Ligand of Interest" by depositor).

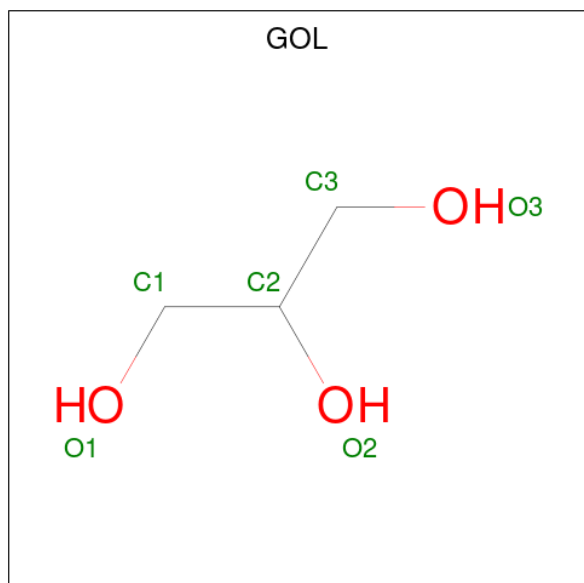


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	Fe	S	0	0
			8	4	4		
6	B	1	Total	Fe	S	0	0
			8	4	4		
6	B	1	Total	Fe	S	0	0
			8	4	4		
6	B	1	Total	Fe	S	0	0
			8	4	4		
6	C	1	Total	Fe	S	0	0
			8	4	4		
6	D	1	Total	Fe	S	0	0
			8	4	4		
6	D	1	Total	Fe	S	0	0
			8	4	4		
6	D	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

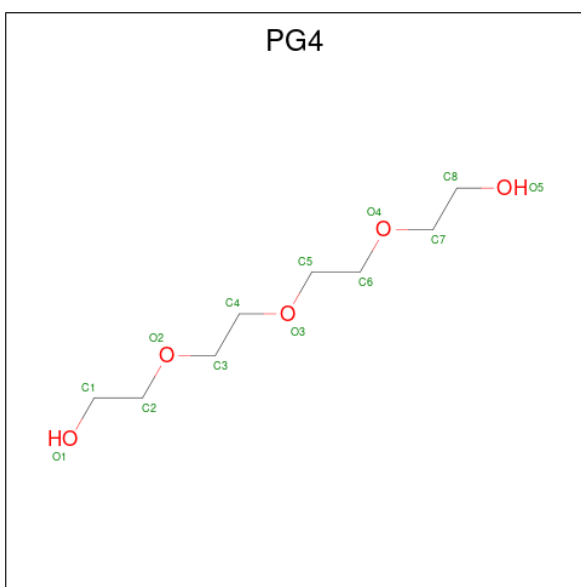
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Cl	0	0
			1	1		
7	B	1	Total	Cl	0	0
			1	1		
7	D	1	Total	Cl	0	0
			1	1		

- Molecule 8 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



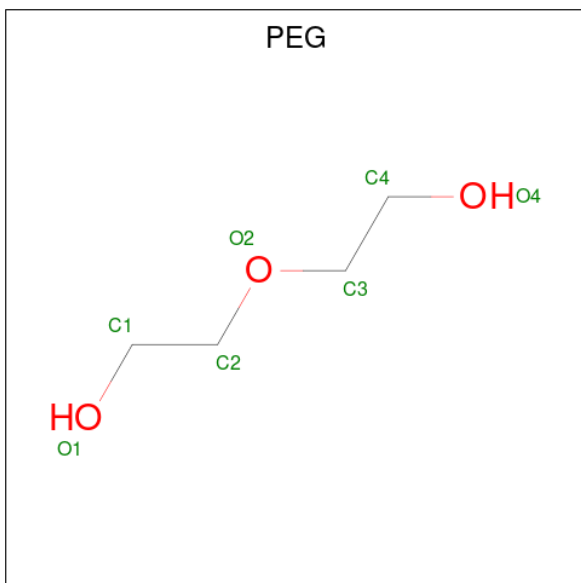
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 6 3 3	0	0
8	A	1	Total C O 6 3 3	0	0
8	A	1	Total C O 6 3 3	0	0
8	A	1	Total C O 6 3 3	0	0
8	A	1	Total C O 6 3 3	0	0
8	A	1	Total C O 6 3 3	0	0
8	A	1	Total C O 6 3 3	0	0
8	A	1	Total C O 6 3 3	0	0
8	A	1	Total C O 6 3 3	0	0
8	C	1	Total C O 6 3 3	0	0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

- Molecule 9 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



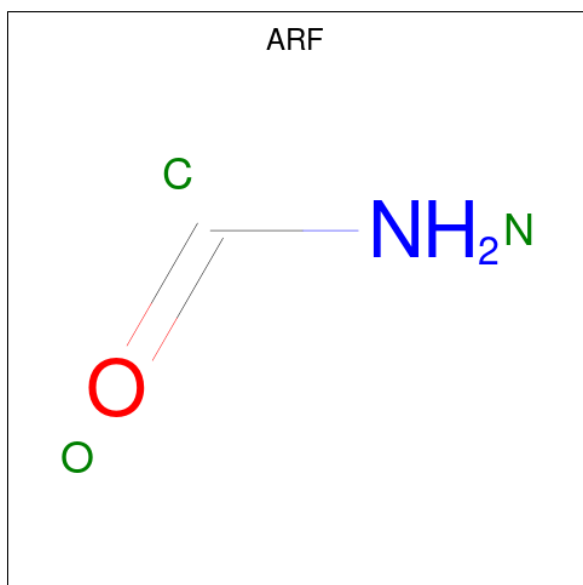
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			13	8	5		
9	A	1	Total	C	O	0	0
			13	8	5		
9	B	1	Total	C	O	0	0
			13	8	5		
9	C	1	Total	C	O	0	0
			13	8	5		

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



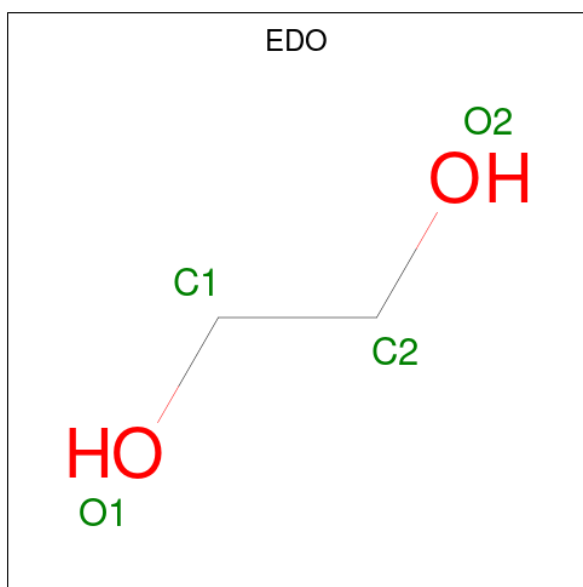
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			7	4	3		
10	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 11 is FORMAMIDE (CCD ID: ARF) (formula: CH_3NO) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 12 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	B	1	Total	C	O	0	0
			4	2	2		
12	D	1	Total	C	O	0	0
			4	2	2		

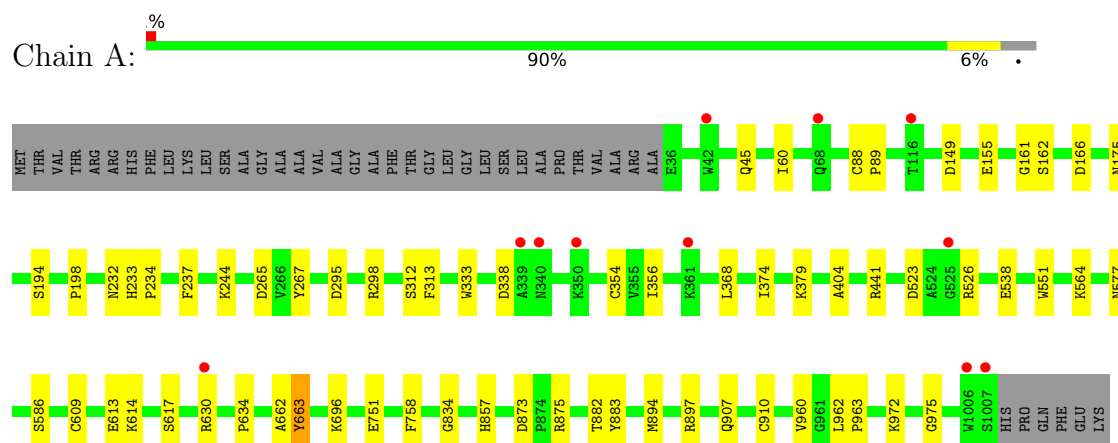
- Molecule 13 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	767	Total	O	0	0
			767	767		
13	B	193	Total	O	0	0
			193	193		
13	C	561	Total	O	0	0
			561	561		
13	D	197	Total	O	0	0
			197	197		

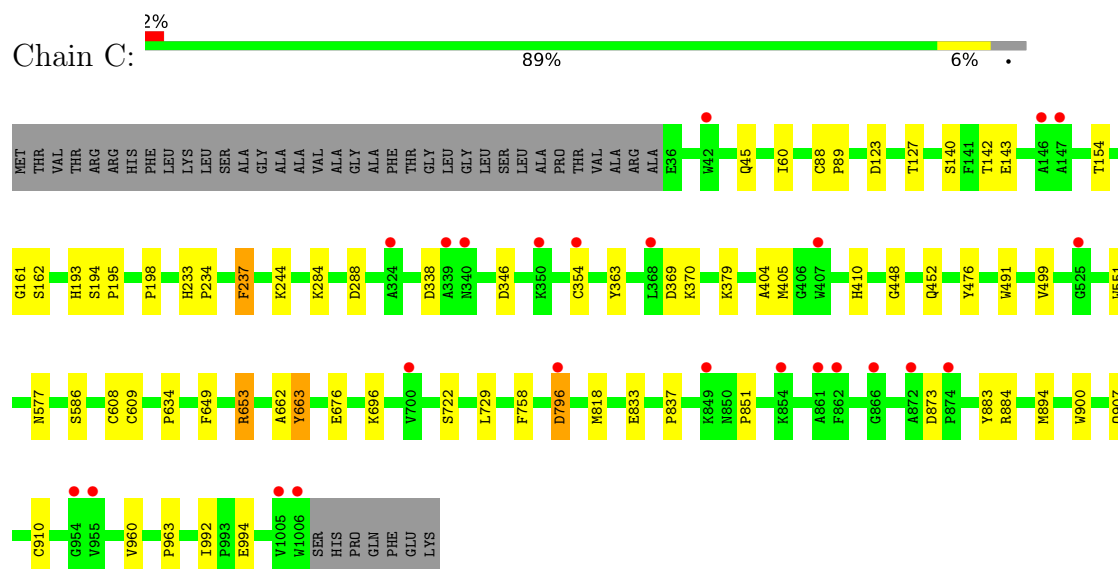
3 Residue-property plots [i](#)

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

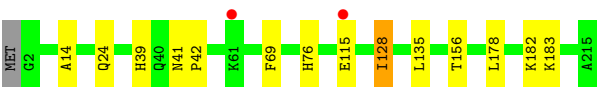


- Molecule 1: Formate dehydrogenase, alpha subunit, selenocysteine-containing

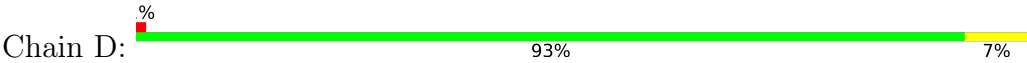


- Molecule 2: Formate dehydrogenase, beta subunit, putative





● Molecule 2: Formate dehydrogenase, beta subunit, putative



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	74.47Å 74.95Å 122.81Å 73.87° 89.14° 71.12°	Depositor
Resolution (Å)	117.57 – 1.42 117.57 – 1.42	Depositor EDS
% Data completeness (in resolution range)	75.8 (117.57-1.42) 75.8 (117.57-1.42)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 1.42Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.148 , 0.176 (Not available) , 0.185	Depositor DCC
R_{free} test set	17188 reflections (3.75%)	wwPDB-VP
Wilson B-factor (Å ²)	13.9	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	20783	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PG4, CL, MGD, H2S, PEG, GOL, ARF, W, SF4, EDO, 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	0.99	2/7880 (0.0%)	1.14	4/10687 (0.0%)
1	C	1.01	1/7836 (0.0%)	1.18	5/10631 (0.0%)
2	B	1.00	1/1729 (0.1%)	1.17	4/2341 (0.2%)
2	D	0.98	0/1699	1.17	1/2302 (0.0%)
All	All	1.00	4/19144 (0.0%)	1.16	14/25961 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	972	LYS	C-O	5.62	1.30	1.23
2	B	115	GLU	CD-OE2	5.35	1.35	1.25
1	C	758	PHE	N-CA	5.10	1.50	1.46
1	A	857	HIS	CE1-NE2	5.02	1.37	1.32

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	338	ASP	CA-CB-CG	7.27	119.87	112.60
2	B	115	GLU	CB-CG-CD	6.34	123.38	112.60
1	C	237	PHE	CA-CB-CG	-6.15	107.65	113.80
1	A	758	PHE	O-C-N	-6.10	118.72	121.53
1	A	175	ASN	CA-CB-CG	-6.10	106.50	112.60
1	A	338	ASP	CA-CB-CG	5.72	118.32	112.60
2	D	215	ALA	CA-C-O	-5.43	111.57	120.80
2	B	39	HIS	CA-CB-CG	-5.39	108.41	113.80
1	C	653	ARG	CG-CD-NE	5.30	123.67	112.00
1	C	142	THR	CA-CB-OG1	-5.26	101.70	109.60
1	C	143	GLU	CB-CG-CD	5.11	121.28	112.60
2	B	24	GLN	CB-CG-CD	5.08	121.24	112.60
2	B	76	HIS	CA-CB-CG	-5.08	108.72	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	175	ASN	CB-CA-C	5.08	119.48	110.85

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	7658	0	7513	46	0
1	C	7623	0	7455	38	0
2	B	1682	0	1662	5	0
2	D	1664	0	1633	8	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
5	A	94	0	44	2	0
5	C	94	0	44	3	0
6	A	8	0	0	0	0
6	B	24	0	0	0	0
6	C	8	0	0	0	0
6	D	24	0	0	1	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	D	1	0	0	0	0
8	A	54	0	72	4	0
8	C	30	0	40	0	0
9	A	26	0	36	3	0
9	B	13	0	18	0	0
9	C	13	0	18	2	0
10	A	7	0	10	4	0
10	C	7	0	10	1	0
11	A	6	0	6	0	0
11	B	3	0	3	1	0
11	C	12	0	12	0	0
12	B	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	D	4	0	6	0	0
13	A	767	0	0	8	0
13	B	193	0	0	1	0
13	C	561	0	0	2	0
13	D	197	0	0	1	0
All	All	20783	0	18588	97	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 3.

All (97) 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:614:LYS:HB2	13:A:1863:HOH:O	1.73	0.87
1:A:696:LYS:HE2	13:A:1432:HOH:O	1.88	0.73
1:A:313:PHE:H	8:A:1115:GOL:H12	1.59	0.67
1:A:374:ILE:HD11	13:A:1240:HOH:O	1.93	0.67
1:A:155:GLU:HG3	13:A:1702:HOH:O	1.95	0.66
1:A:295:ASP:OD1	1:A:298[B]:ARG:NH1	2.28	0.65
1:A:149:ASP:OD2	13:A:1201:HOH:O	2.15	0.64
1:C:608:CYS:C	1:C:609[B]:CYS:CA	2.70	0.64
2:D:211:GLU:HG2	13:D:440:HOH:O	1.98	0.62
1:C:609[C]:CYS:SG	1:C:634:PRO:HB3	2.40	0.62
1:A:897:ARG:HH22	10:A:1116:PEG:H22	1.67	0.59
1:C:284:LYS:NZ	1:C:288:ASP:OD2	2.38	0.56
2:D:140:MET:HG2	2:D:140:MET:O	2.05	0.56
1:A:354[B]:CYS:SG	1:A:356[B]:ILE:HG22	2.46	0.55
1:C:696:LYS:HE2	1:C:722:SER:OG	2.07	0.54
1:A:975:GLY:HA2	9:A:1110:PG4:H22	1.89	0.54
1:C:140:SER:OG	1:C:154:THR:HG23	2.08	0.53
1:A:313:PHE:H	8:A:1115:GOL:C1	2.20	0.53
1:C:193:HIS:HB2	13:C:1502:HOH:O	2.09	0.53
1:A:374:ILE:CD1	13:A:1240:HOH:O	2.55	0.52
1:C:894:MET:SD	5:C:1104:MGD:H2'	2.50	0.52
1:C:45:GLN:HG3	1:C:60:ILE:CG2	2.39	0.51
1:C:88:CYS:HB2	1:C:89:PRO:HD2	1.92	0.50
1:A:751:GLU:OE1	13:A:1202:HOH:O	2.19	0.50
1:C:370:LYS:NZ	1:C:851:PRO:O	2.45	0.50
1:A:894:MET:SD	5:A:1104:MGD:H2'	2.52	0.49
1:C:729:LEU:HD21	10:C:1107:PEG:H41	1.94	0.49
2:D:41:ASN:HA	2:D:42:PRO:C	2.35	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:41:ASN:HA	2:B:42:PRO:C	2.38	0.49
1:A:873:ASP:OD2	1:A:875:ARG:NH1	2.46	0.49
1:C:410:HIS:HA	1:C:994:GLU:HG3	1.93	0.49
2:D:128:ILE:HG12	2:D:135:LEU:CD2	2.43	0.48
1:C:910:CYS:SG	1:C:960:VAL:HG13	2.54	0.48
1:A:45:GLN:HG3	1:A:60:ILE:CG2	2.44	0.48
1:A:312:SER:HA	8:A:1115:GOL:H11	1.94	0.48
1:A:609[A]:CYS:SG	1:A:634:PRO:HB3	2.53	0.48
1:A:162:SER:HB2	1:A:551:TRP:O	2.14	0.47
1:C:162:SER:HB2	1:C:551:TRP:O	2.14	0.47
1:C:448:GLY:O	1:C:452:GLN:HG2	2.15	0.47
1:A:897:ARG:HH12	10:A:1116:PEG:H22	1.79	0.47
1:C:194:SER:O	1:C:198:PRO:HD2	2.15	0.47
1:A:897:ARG:HH22	10:A:1116:PEG:C2	2.26	0.46
2:B:14:ALA:HB2	2:B:69:PHE:CG	2.50	0.46
1:A:368:LEU:HD11	13:A:1942:HOH:O	2.16	0.45
1:C:476:TYR:HE1	1:C:499:VAL:HG21	1.81	0.45
1:C:195:PRO:HG3	1:C:992:ILE:HG21	1.99	0.45
1:C:662:ALA:O	1:C:663:TYR:C	2.60	0.45
1:C:346:ASP:HB3	1:C:354:CYS:SG	2.57	0.45
1:A:907:GLN:NE2	9:A:1110:PG4:H52	2.32	0.45
1:A:897:ARG:NH2	10:A:1116:PEG:H22	2.32	0.44
1:C:244:LYS:HE2	13:C:1309:HOH:O	2.16	0.44
2:D:81:PRO:HD2	6:D:302:SF4:S3	2.57	0.44
1:C:907:GLN:HE22	9:C:1108:PG4:H82	1.82	0.44
1:A:662:ALA:O	1:A:663:TYR:C	2.61	0.44
1:A:88:CYS:HB2	1:A:89:PRO:HD2	1.99	0.44
1:C:883:TYR:O	1:C:963:PRO:HA	2.16	0.44
1:C:369:ASP:OD1	1:C:379:LYS:HD2	2.18	0.44
1:A:577:ASN:OD1	1:A:586:SER:HB3	2.18	0.44
1:A:523:ASP:HB2	1:A:526:ARG:HG3	1.99	0.44
1:A:910:CYS:SG	1:A:960:VAL:HG13	2.57	0.44
2:B:178:LEU:HD11	2:B:182:LYS:HE3	1.99	0.44
1:A:161:GLY:HA3	1:A:551:TRP:O	2.18	0.43
1:A:166[B]:ASP:HB2	1:A:617:SER:O	2.19	0.43
1:A:194:SER:O	1:A:198:PRO:HD2	2.18	0.43
1:A:333:TRP:CH2	1:A:834:GLY:HA2	2.53	0.43
1:A:267:TYR:H	11:B:305:ARF:HN1	1.67	0.42
1:A:45:GLN:HG3	1:A:60:ILE:HG23	2.02	0.42
1:A:313:PHE:N	8:A:1115:GOL:H12	2.30	0.42
1:C:88:CYS:CB	1:C:89:PRO:HD2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:538:GLU:OE1	1:A:564:LYS:HE3	2.19	0.42
1:C:233:HIS:N	1:C:234:PRO:HD3	2.34	0.42
1:A:232:ASN:HA	5:A:1103:MGD:N20	2.34	0.42
1:A:538:GLU:OE1	1:A:564:LYS:CE	2.67	0.42
1:C:649:PHE:CE2	1:C:653:ARG:HD2	2.54	0.42
1:C:45:GLN:HG3	1:C:60:ILE:HG23	2.01	0.42
1:A:244:LYS:NZ	1:A:265:ASP:OD1	2.53	0.42
1:C:405:MET:HG2	5:C:1103:MGD:H101	2.01	0.42
1:C:577:ASN:OD1	1:C:586:SER:HB3	2.20	0.42
2:D:204:ARG:O	2:D:205:ASP:C	2.61	0.42
1:A:883:TYR:O	1:A:963:PRO:HA	2.20	0.41
1:C:476:TYR:CE1	1:C:499:VAL:HG21	2.55	0.41
2:D:14:ALA:HB2	2:D:69:PHE:CG	2.55	0.41
1:A:907:GLN:HE22	9:A:1110:PG4:H52	1.85	0.41
1:C:161:GLY:HA3	1:C:551:TRP:O	2.20	0.41
2:B:128:ILE:HB	2:B:135:LEU:HD23	2.02	0.41
1:C:884:ARG:HH22	5:C:1104:MGD:H15	1.68	0.41
1:C:900:TRP:CH2	2:D:24:GLN:HA	2.56	0.41
1:A:233:HIS:N	1:A:234:PRO:HD3	2.35	0.41
1:A:368:LEU:HB3	1:A:379:LYS:HG3	2.02	0.41
1:C:123:ASP:O	1:C:127[A]:THR:HG22	2.21	0.41
1:A:882:THR:HA	1:A:962:LEU:O	2.20	0.41
1:C:491:TRP:HB3	1:C:818:MET:SD	2.61	0.41
1:C:873:ASP:C	1:C:873:ASP:OD1	2.64	0.41
1:C:363:TYR:OH	1:C:837:PRO:HD2	2.21	0.41
1:A:613:GLU:O	1:A:630:ARG:NE	2.50	0.40
2:B:183:LYS:HE3	13:B:409:HOH:O	2.21	0.40
9:C:1108:PG4:H41	9:C:1108:PG4:H61	1.88	0.40

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	977/1013 (96%)	954 (98%)	21 (2%)	2 (0%)	43	20
1	C	972/1013 (96%)	946 (97%)	23 (2%)	3 (0%)	36	17
2	B	216/215 (100%)	207 (96%)	9 (4%)	0	100	100
2	D	212/215 (99%)	204 (96%)	8 (4%)	0	100	100
All	All	2377/2456 (97%)	2311 (97%)	61 (3%)	5 (0%)	43	20

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	796	ASP
1	A	404	ALA
1	C	404	ALA
1	A	663	TYR
1	C	663	TYR

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	797/818 (97%)	795 (100%)	2 (0%)	86	71
1	C	792/818 (97%)	788 (100%)	4 (0%)	81	62
2	B	189/186 (102%)	187 (99%)	2 (1%)	65	35
2	D	185/186 (100%)	184 (100%)	1 (0%)	81	62
All	All	1963/2008 (98%)	1954 (100%)	9 (0%)	81	62

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	237	PHE
1	A	441	ARG
2	B	128	ILE
2	B	156	THR
1	C	237	PHE
1	C	676	GLU

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Mol	Chain	Res	Type
1	C	796	ASP
1	C	833	GLU
2	D	156	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	GLN
1	A	383	GLN
1	A	863	GLN
1	C	45	GLN
1	C	383	GLN
1	C	671	ASN
1	C	710	GLN
1	C	857	HIS

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 48 ligands modelled in this entry, 2 are modelled with single atom and 5 are monoatomic - leaving 41 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
6	SF4	C	1105	1	0,12,12	-	-	-		
8	GOL	C	1112	-	5,5,5	0.19	0	5,5,5	0.69	0
6	SF4	D	303	2	0,12,12	-	-	-		
8	GOL	A	1114	-	5,5,5	0.20	0	5,5,5	0.55	0
5	MGD	C	1103	4	47,52,52	0.74	1 (2%)	58,81,81	1.14	7 (12%)
9	PG4	C	1108	-	12,12,12	0.32	0	11,11,11	0.30	0
11	ARF	C	1114	-	2,2,2	0.42	0	1,1,1	0.85	0
9	PG4	A	1110	-	12,12,12	0.22	0	11,11,11	0.28	0
6	SF4	D	301	2	0,12,12	-	-	-		
5	MGD	A	1104	4	47,52,52	0.73	1 (2%)	58,81,81	1.23	7 (12%)
11	ARF	C	1113	-	2,2,2	0.73	0	1,1,1	1.05	0
8	GOL	C	1111	-	5,5,5	0.17	0	5,5,5	0.49	0
6	SF4	B	303	2	0,12,12	-	-	-		
8	GOL	A	1117	-	5,5,5	0.42	0	5,5,5	0.98	0
8	GOL	A	1107	-	5,5,5	0.27	0	5,5,5	0.48	0
9	PG4	A	1112	-	12,12,12	0.19	0	11,11,11	0.20	0
8	GOL	A	1109	-	5,5,5	0.16	0	5,5,5	0.30	0
8	GOL	A	1115	-	5,5,5	0.20	0	5,5,5	0.20	0
11	ARF	B	305	-	2,2,2	2.12	1 (50%)	1,1,1	0.15	0
11	ARF	C	1115	-	2,2,2	0.79	0	1,1,1	0.93	0
8	GOL	A	1113	-	5,5,5	0.18	0	5,5,5	0.71	0
6	SF4	D	302	2	0,12,12	-	-	-		
11	ARF	A	1119	-	2,2,2	0.73	0	1,1,1	0.91	0
6	SF4	B	301	2	0,12,12	-	-	-		
10	PEG	A	1116	-	6,6,6	0.14	0	5,5,5	0.21	0
8	GOL	A	1108	-	5,5,5	0.13	0	5,5,5	0.46	0
8	GOL	C	1109	-	5,5,5	0.13	0	5,5,5	0.57	0
12	EDO	D	305	-	3,3,3	0.22	0	2,2,2	0.35	0
6	SF4	A	1105	1	0,12,12	-	-	-		
8	GOL	A	1118	-	5,5,5	0.16	0	5,5,5	0.34	0
12	EDO	B	307	-	3,3,3	0.14	0	2,2,2	0.10	0
5	MGD	C	1104	4	47,52,52	0.62	0	58,81,81	1.25	8 (13%)
8	GOL	A	1111	-	5,5,5	0.17	0	5,5,5	0.38	0
6	SF4	B	302	2	0,12,12	-	-	-		
10	PEG	C	1107	-	6,6,6	0.56	0	5,5,5	0.48	0
11	ARF	A	1120	-	2,2,2	0.12	0	1,1,1	0.84	0
11	ARF	C	1116	-	2,2,2	0.63	0	1,1,1	0.69	0
8	GOL	C	1110	-	5,5,5	0.33	0	5,5,5	0.68	0
9	PG4	B	306	-	12,12,12	0.24	0	11,11,11	0.19	0
8	GOL	C	1106	-	5,5,5	0.12	0	5,5,5	0.31	0
5	MGD	A	1103	4	47,52,52	0.80	1 (2%)	58,81,81	1.06	5 (8%)

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
8	GOL	C	1112	-	-	0/4/4/4	-
6	SF4	C	1105	1	-	-	0/6/5/5
6	SF4	D	303	2	-	-	0/6/5/5
8	GOL	A	1114	-	-	0/4/4/4	-
5	MGD	C	1103	4	-	3/22/66/66	0/6/6/6
9	PG4	C	1108	-	-	7/10/10/10	-
9	PG4	A	1110	-	-	4/10/10/10	-
6	SF4	D	301	2	-	-	0/6/5/5
5	MGD	A	1104	4	-	3/22/66/66	0/6/6/6
8	GOL	C	1111	-	-	0/4/4/4	-
6	SF4	B	303	2	-	-	0/6/5/5
8	GOL	A	1117	-	-	0/4/4/4	-
8	GOL	A	1107	-	-	0/4/4/4	-
9	PG4	A	1112	-	-	4/10/10/10	-
8	GOL	A	1109	-	-	0/4/4/4	-
8	GOL	A	1115	-	-	2/4/4/4	-
8	GOL	A	1113	-	-	2/4/4/4	-
6	SF4	D	302	2	-	-	0/6/5/5
6	SF4	B	301	2	-	-	0/6/5/5
10	PEG	A	1116	-	-	2/4/4/4	-
8	GOL	A	1108	-	-	0/4/4/4	-
8	GOL	C	1109	-	-	1/4/4/4	-
12	EDO	D	305	-	-	0/1/1/1	-
8	GOL	A	1118	-	-	3/4/4/4	-
12	EDO	B	307	-	-	0/1/1/1	-
6	SF4	A	1105	1	-	-	0/6/5/5
5	MGD	C	1104	4	-	3/22/66/66	0/6/6/6
8	GOL	A	1111	-	-	0/4/4/4	-
10	PEG	C	1107	-	-	2/4/4/4	-
6	SF4	B	302	2	-	-	0/6/5/5
8	GOL	C	1110	-	-	0/4/4/4	-
9	PG4	B	306	-	-	4/10/10/10	-
8	GOL	C	1106	-	-	0/4/4/4	-
5	MGD	A	1103	4	-	3/22/66/66	0/6/6/6

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1103	MGD	C23-C14	-2.70	1.51	1.53
11	B	305	ARF	O-C	2.63	1.31	1.22
5	C	1103	MGD	C23-C14	2.46	1.55	1.53
5	A	1104	MGD	C10-C11	2.42	1.55	1.51

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	1104	MGD	O11-C23-C14	3.96	111.60	108.96
5	C	1103	MGD	C23-C14-N15	3.95	111.75	107.87
5	A	1104	MGD	C17-C16-N15	3.72	126.42	116.27
5	A	1104	MGD	C19-N20-C21	3.42	119.40	113.36
5	A	1104	MGD	O11-C23-C14	3.40	111.23	108.96
5	A	1103	MGD	C19-N20-C21	3.14	118.90	113.36
5	C	1104	MGD	C23-C14-N15	3.04	110.85	107.87
5	C	1104	MGD	O4'-C1'-C2'	-3.02	100.16	106.62
5	C	1104	MGD	C19-N20-C21	2.97	118.61	113.36
5	A	1103	MGD	C23-C14-N15	2.96	110.78	107.87
5	C	1103	MGD	O11-C23-N22	2.95	111.28	108.61
5	C	1103	MGD	C19-N20-C21	2.59	117.93	113.36
5	A	1104	MGD	C23-C14-N15	2.59	110.41	107.87
5	C	1103	MGD	O11-C23-C14	-2.56	107.25	108.96
5	A	1103	MGD	C4'-O4'-C1'	-2.47	104.02	109.47
5	C	1104	MGD	C4'-O4'-C1'	-2.43	104.10	109.47
5	A	1104	MGD	C4'-O4'-C1'	-2.43	104.11	109.47
5	A	1104	MGD	O4'-C1'-C2'	-2.37	101.55	106.62
5	C	1104	MGD	O11-C23-N22	-2.36	106.46	108.61
5	C	1103	MGD	O4'-C1'-C2'	-2.17	101.97	106.62
5	A	1103	MGD	O2A-PA-O1A	2.15	122.43	112.44
5	A	1103	MGD	C17-C16-N15	2.13	122.07	116.27
5	C	1104	MGD	O2A-PA-O1A	2.09	122.18	112.44
5	A	1104	MGD	O11-C23-N22	-2.09	106.71	108.61
5	C	1104	MGD	C17-C16-N15	2.08	121.93	116.27
5	C	1103	MGD	O2A-PA-O1A	2.01	121.80	112.44
5	C	1103	MGD	C4'-O4'-C1'	-2.01	105.04	109.47

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1104	MGD	C4'-C5'-O5'-PB
5	C	1103	MGD	PA-O3B-PB-O5'

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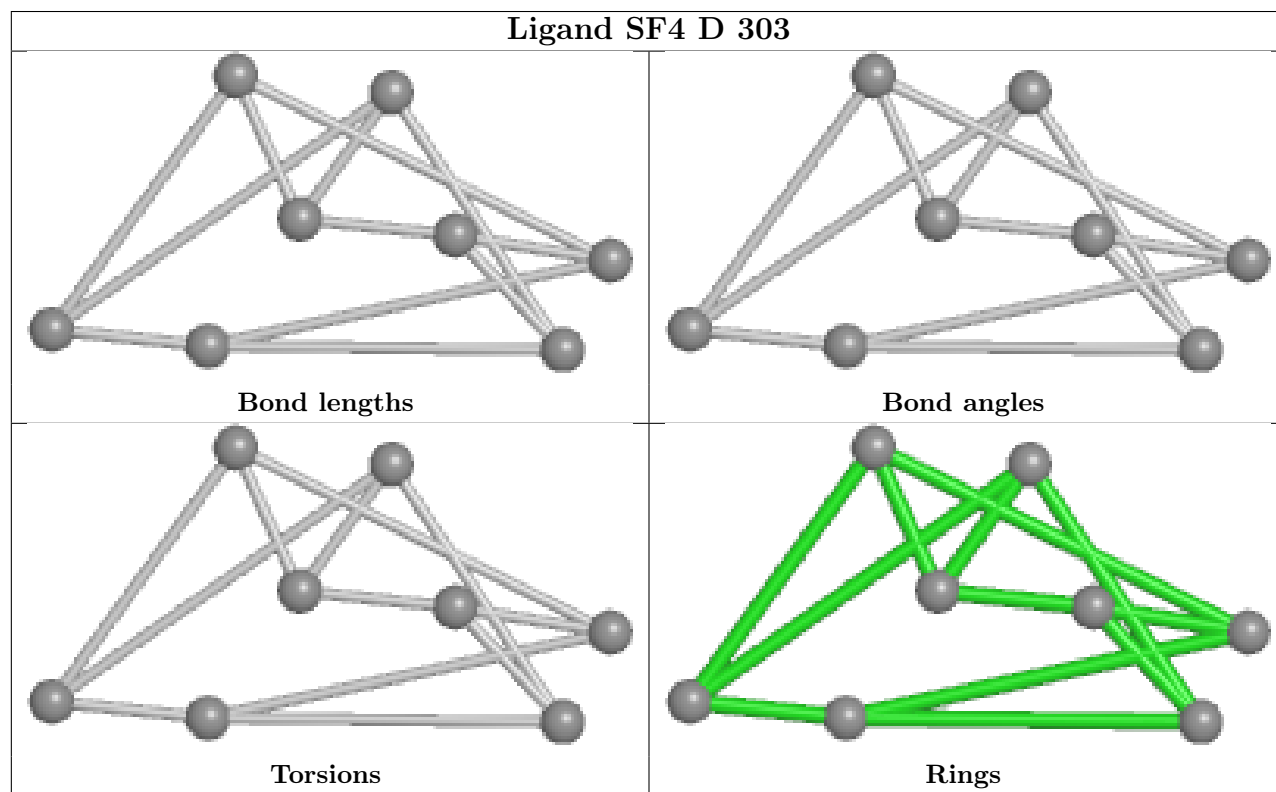
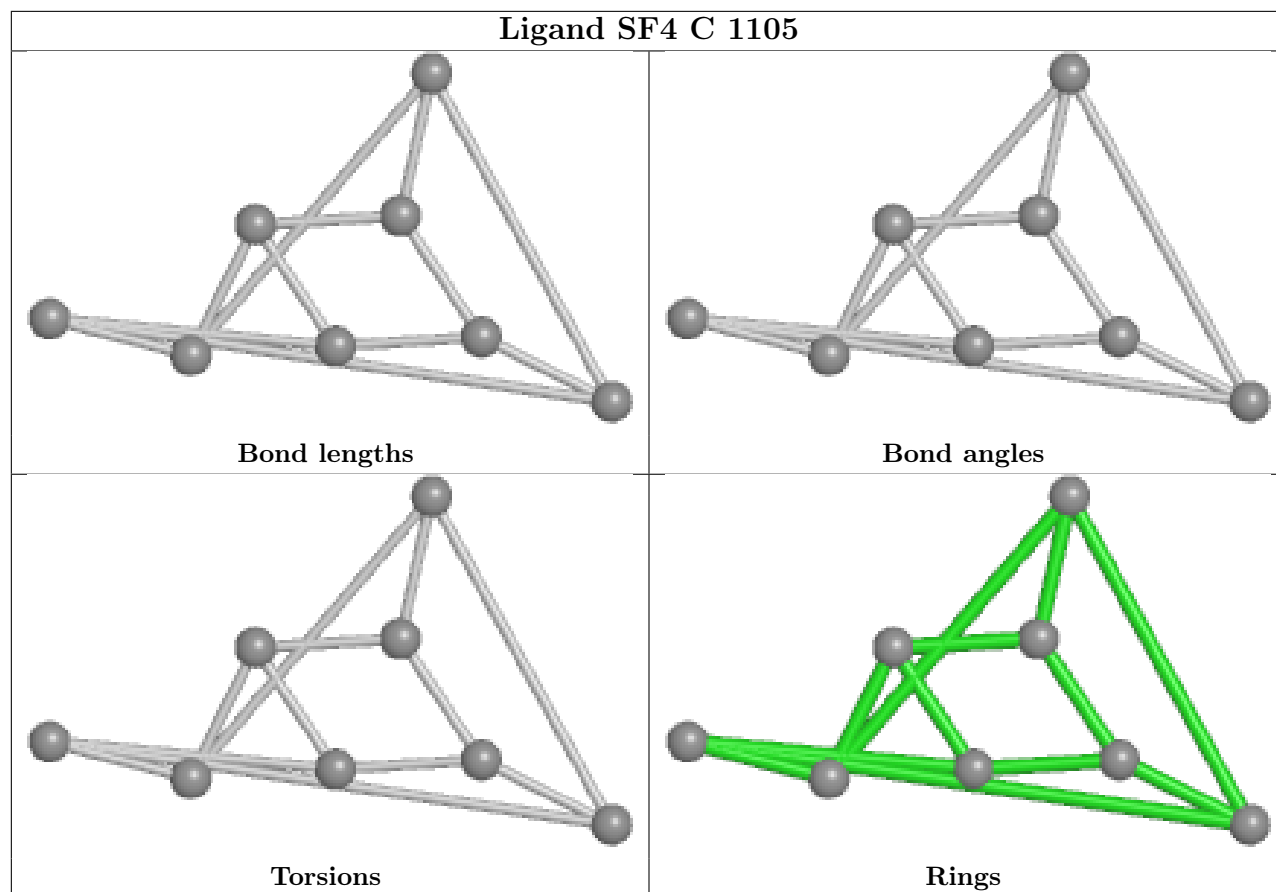
Mol	Chain	Res	Type	Atoms
5	C	1104	MGD	C4'-C5'-O5'-PB
8	A	1113	GOL	C1-C2-C3-O3
9	C	1108	PG4	C6-C5-O3-C4
9	C	1108	PG4	O3-C5-C6-O4
5	A	1104	MGD	O4'-C4'-C5'-O5'
5	C	1104	MGD	O4'-C4'-C5'-O5'
5	C	1104	MGD	C3'-C4'-C5'-O5'
9	C	1108	PG4	O2-C3-C4-O3
9	A	1110	PG4	O4-C7-C8-O5
9	B	306	PG4	O4-C7-C8-O5
10	A	1116	PEG	O1-C1-C2-O2
8	A	1118	GOL	O1-C1-C2-C3
10	A	1116	PEG	O2-C3-C4-O4
8	A	1113	GOL	O2-C2-C3-O3
10	C	1107	PEG	O1-C1-C2-O2
9	C	1108	PG4	O4-C7-C8-O5
5	A	1104	MGD	C3'-C4'-C5'-O5'
8	A	1115	GOL	O2-C2-C3-O3
5	A	1103	MGD	PA-O3B-PB-O5'
9	C	1108	PG4	C4-C3-O2-C2
8	C	1109	GOL	O2-C2-C3-O3
8	A	1118	GOL	C1-C2-C3-O3
9	C	1108	PG4	C8-C7-O4-C6
9	A	1112	PG4	O4-C7-C8-O5
9	B	306	PG4	C6-C5-O3-C4
9	B	306	PG4	C5-C6-O4-C7
9	B	306	PG4	C1-C2-O2-C3
9	A	1112	PG4	C1-C2-O2-C3
10	C	1107	PEG	C1-C2-O2-C3
8	A	1118	GOL	O1-C1-C2-O2
9	A	1110	PG4	O1-C1-C2-O2
9	C	1108	PG4	C1-C2-O2-C3
9	A	1110	PG4	O2-C3-C4-O3
8	A	1115	GOL	C1-C2-C3-O3
5	A	1103	MGD	PA-O3B-PB-O1B
5	C	1103	MGD	PA-O3B-PB-O1B
9	A	1112	PG4	O3-C5-C6-O4
9	A	1112	PG4	C3-C4-O3-C5
9	A	1110	PG4	C5-C6-O4-C7
5	A	1103	MGD	PA-O3B-PB-O2B
5	C	1103	MGD	PA-O3B-PB-O2B

There are no ring outliers.

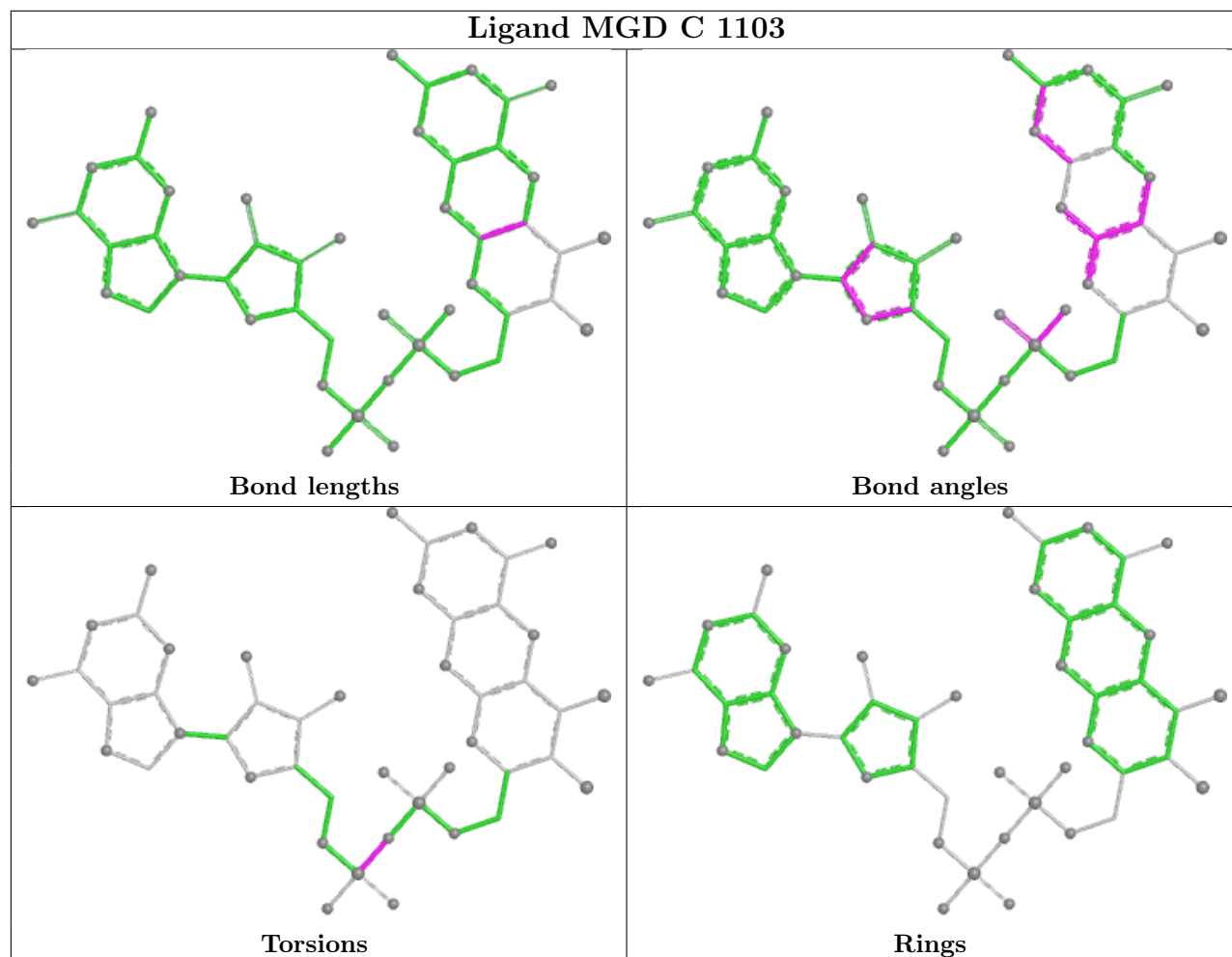
11 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	1103	MGD	1	0
9	C	1108	PG4	2	0
9	A	1110	PG4	3	0
5	A	1104	MGD	1	0
8	A	1115	GOL	4	0
11	B	305	ARF	1	0
6	D	302	SF4	1	0
10	A	1116	PEG	4	0
5	C	1104	MGD	2	0
10	C	1107	PEG	1	0
5	A	1103	MGD	1	0

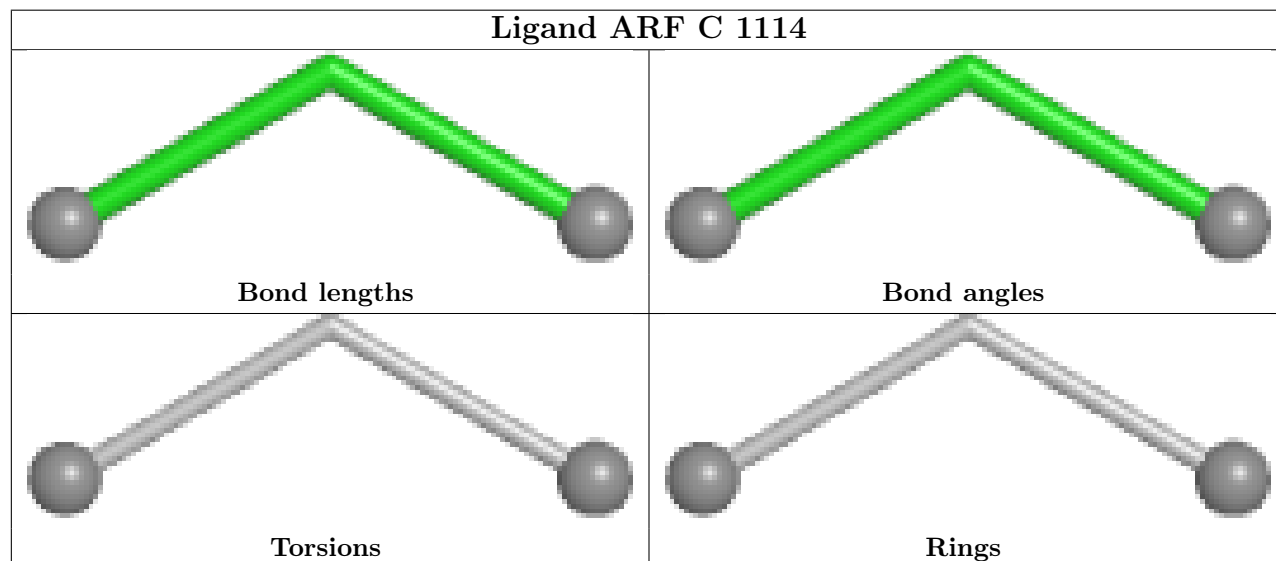
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

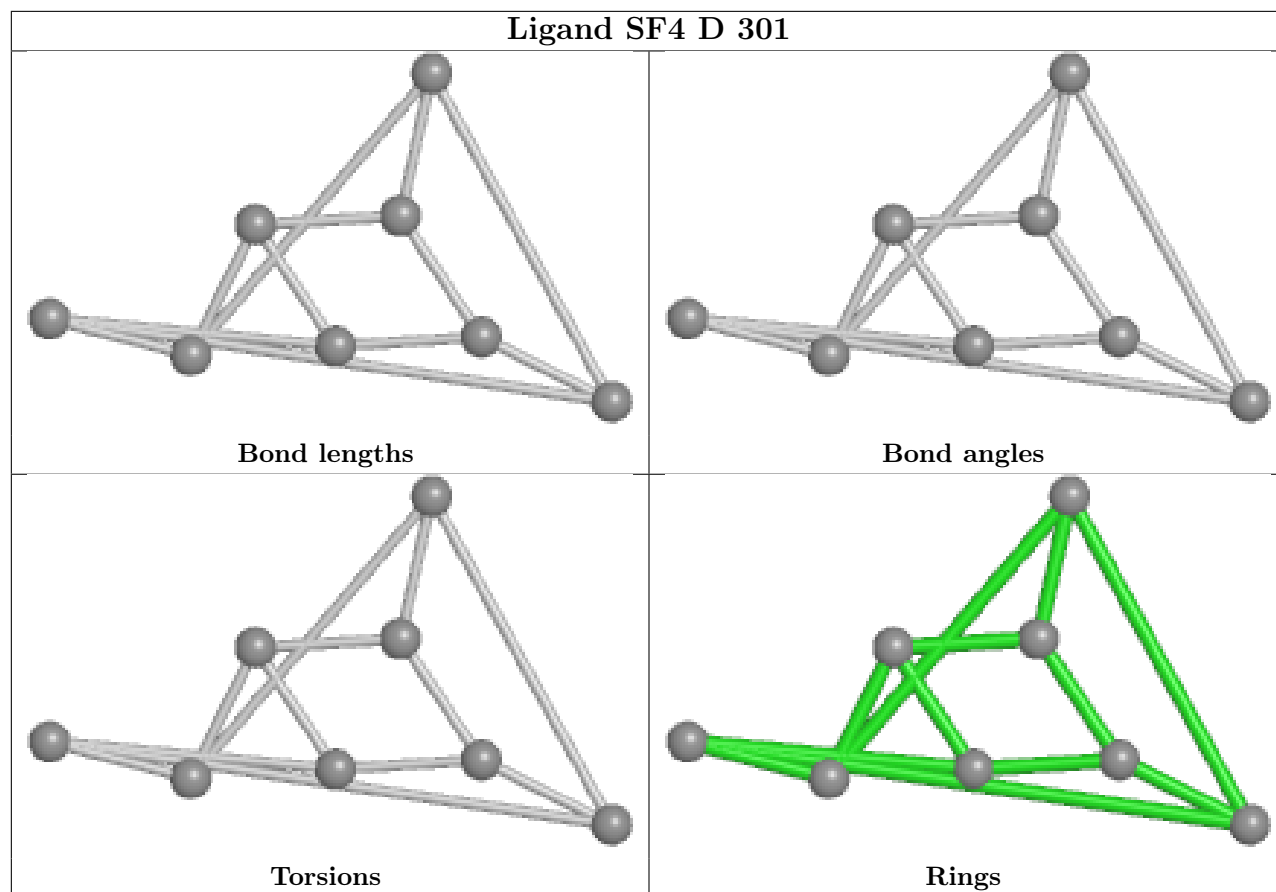


Ligand MGD C 1103

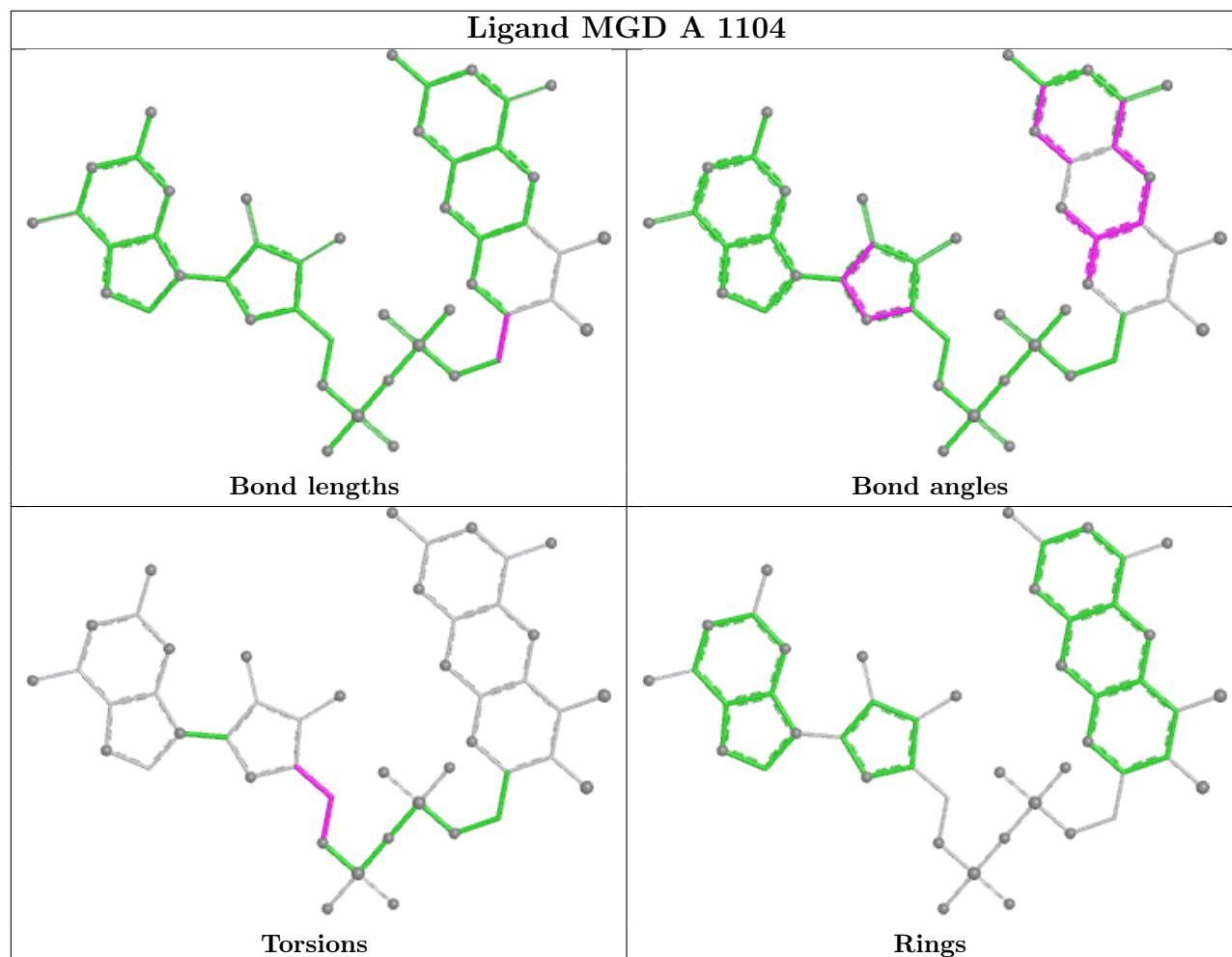


Ligand ARF C 1114

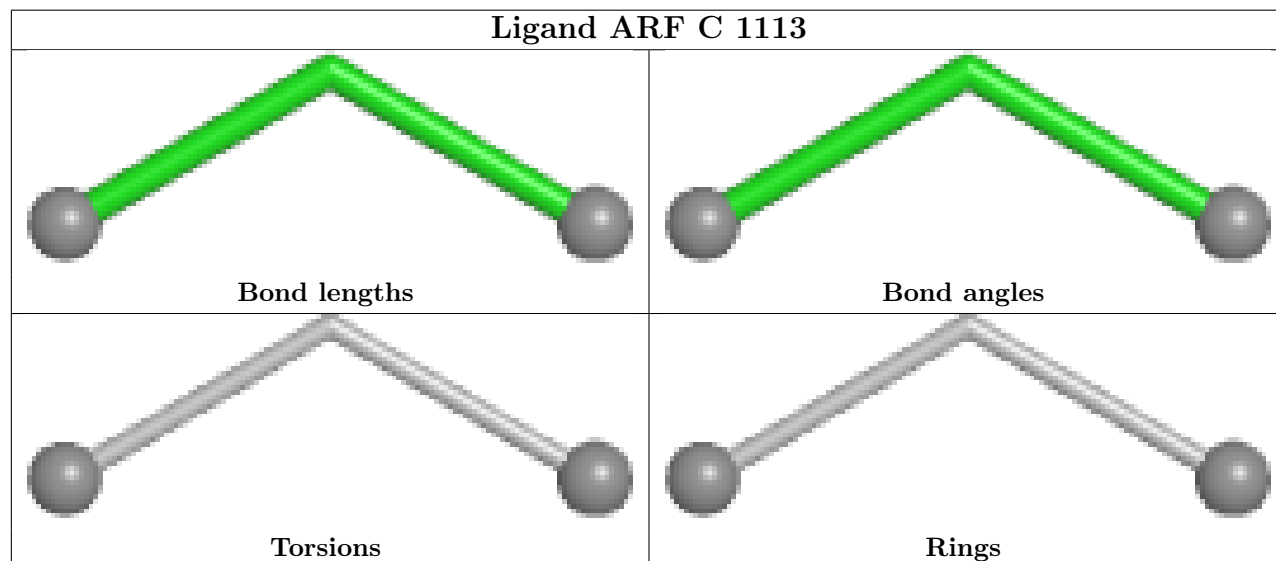


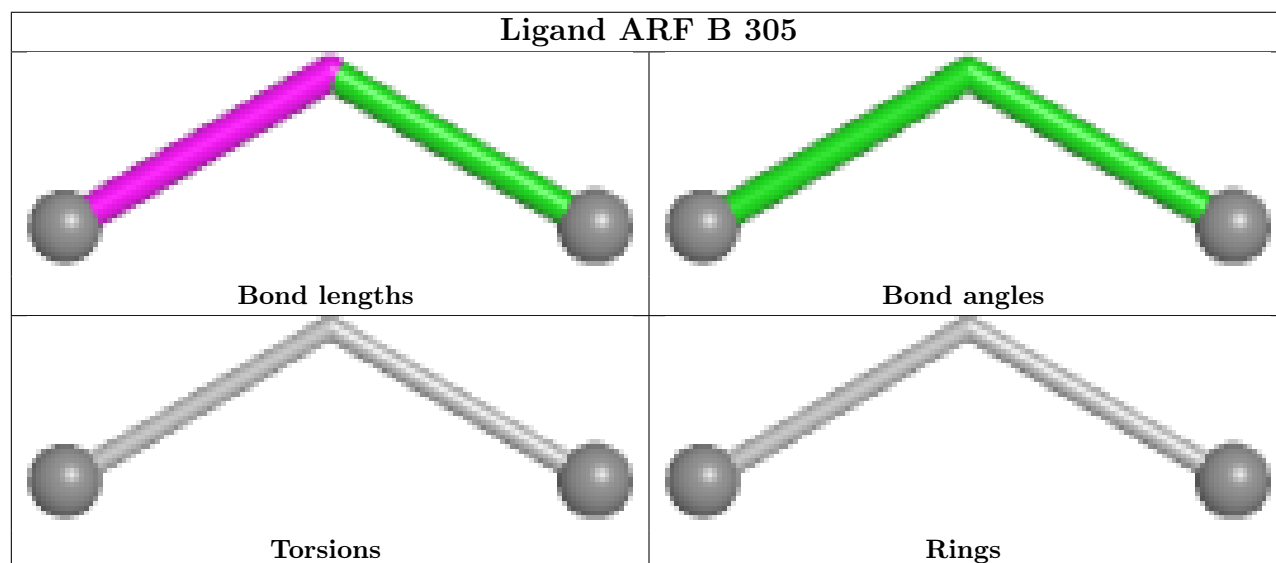
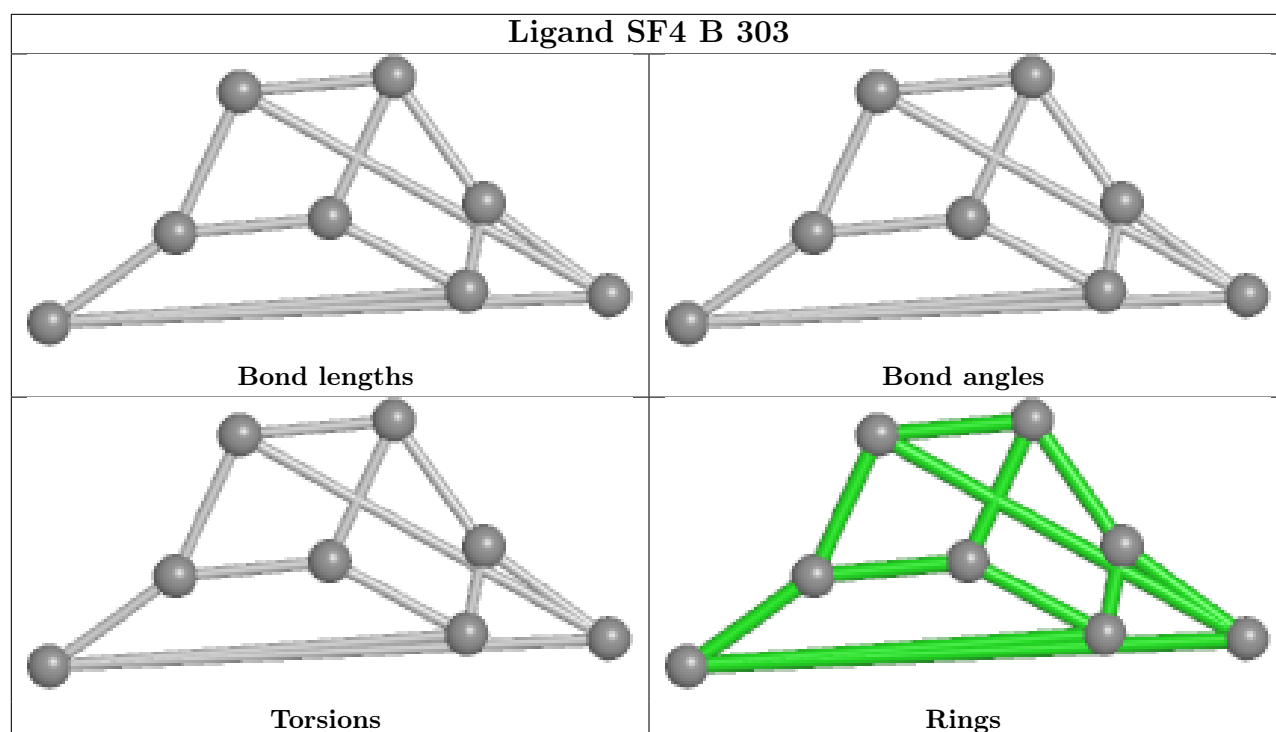


Ligand MGD A 1104

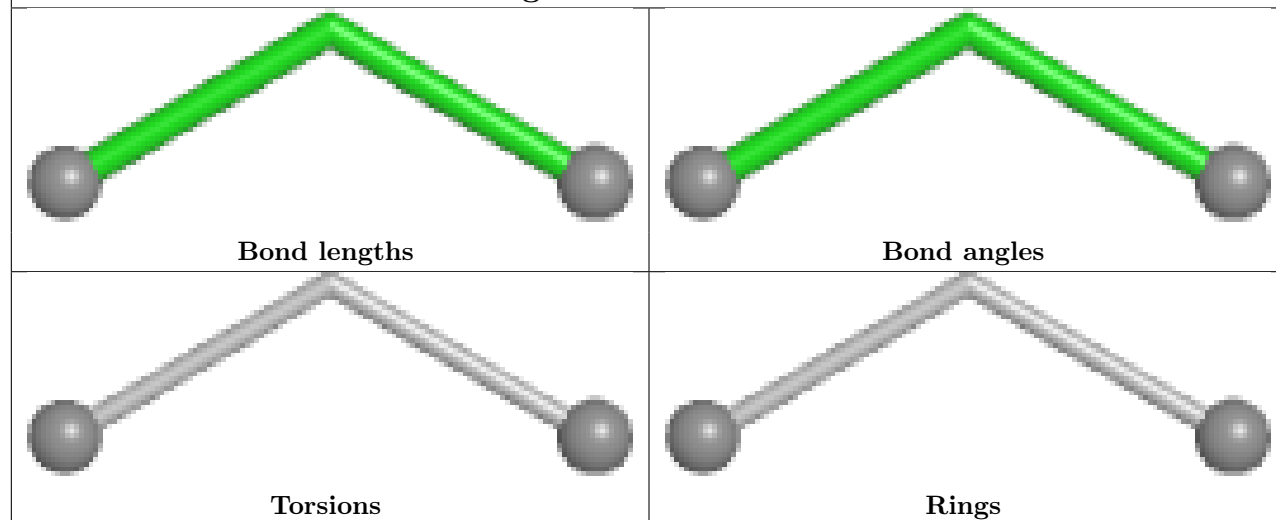


Ligand ARF C 1113

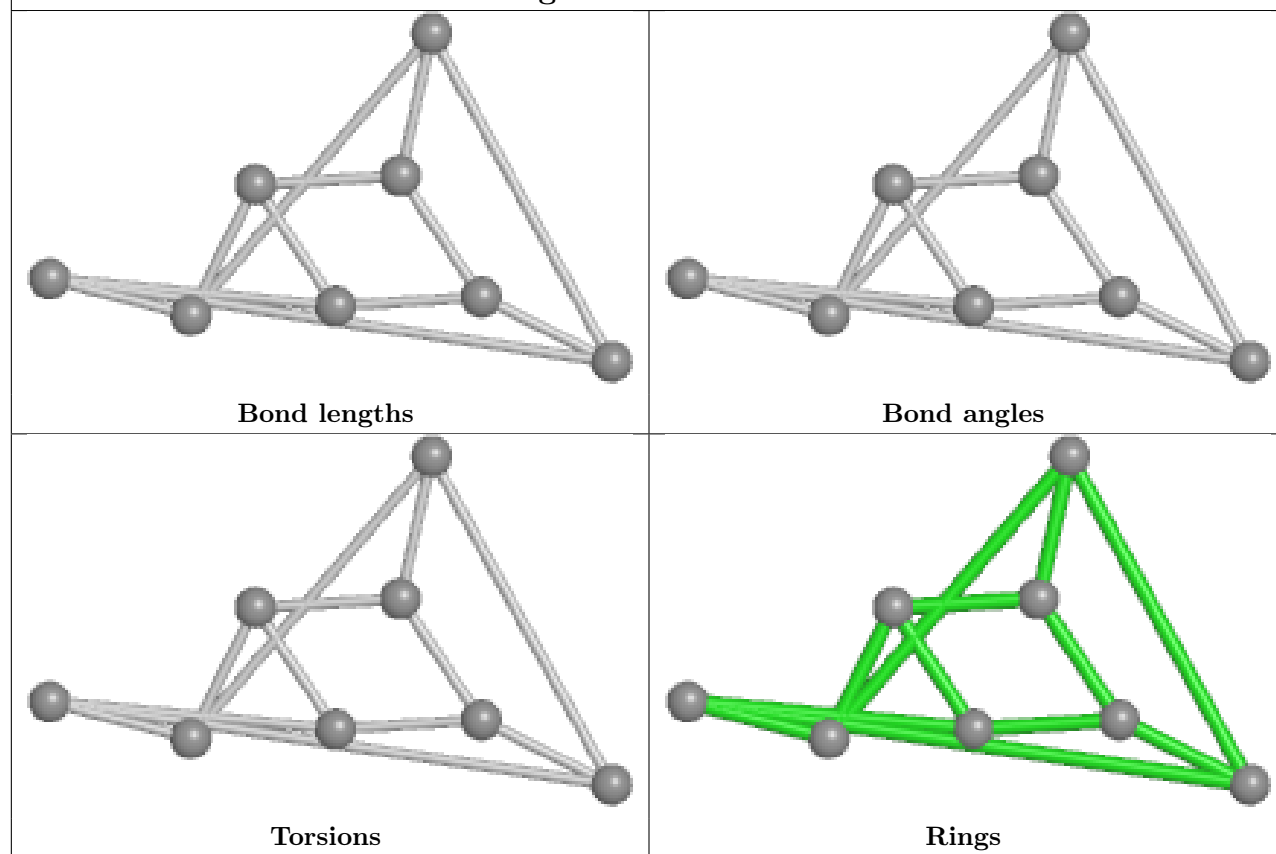


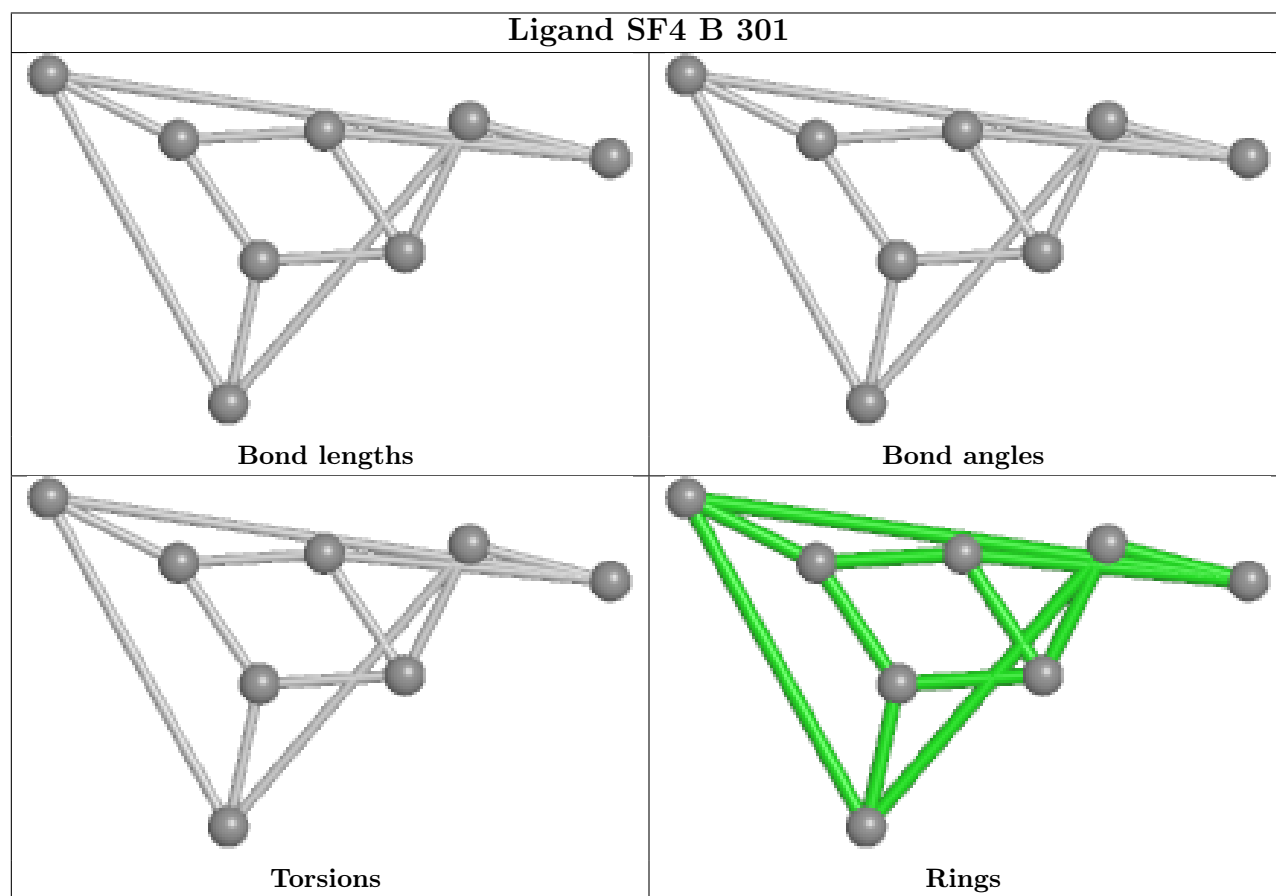
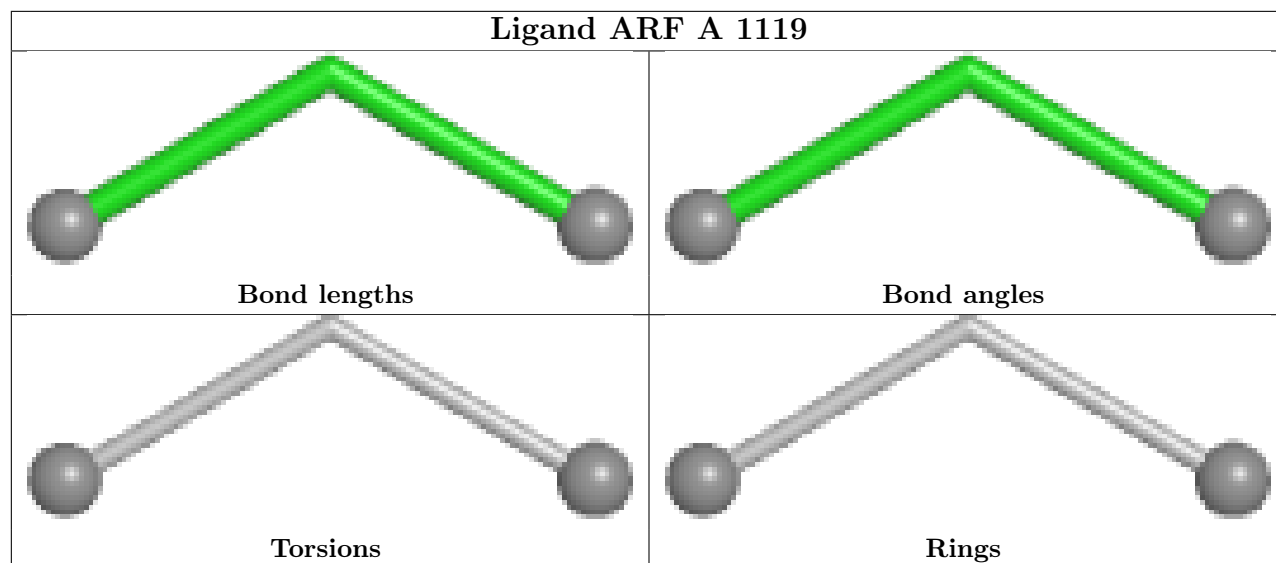


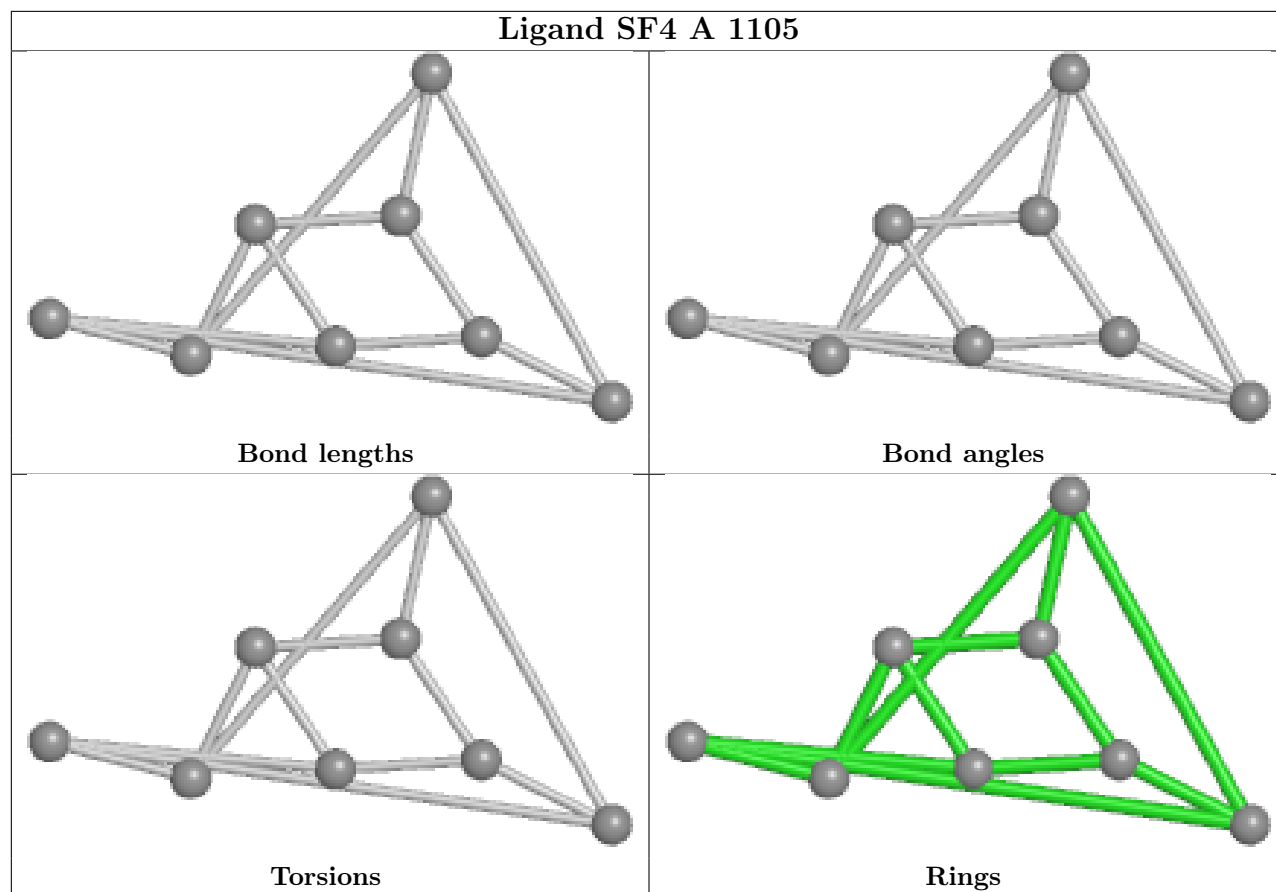
Ligand ARF C 1115



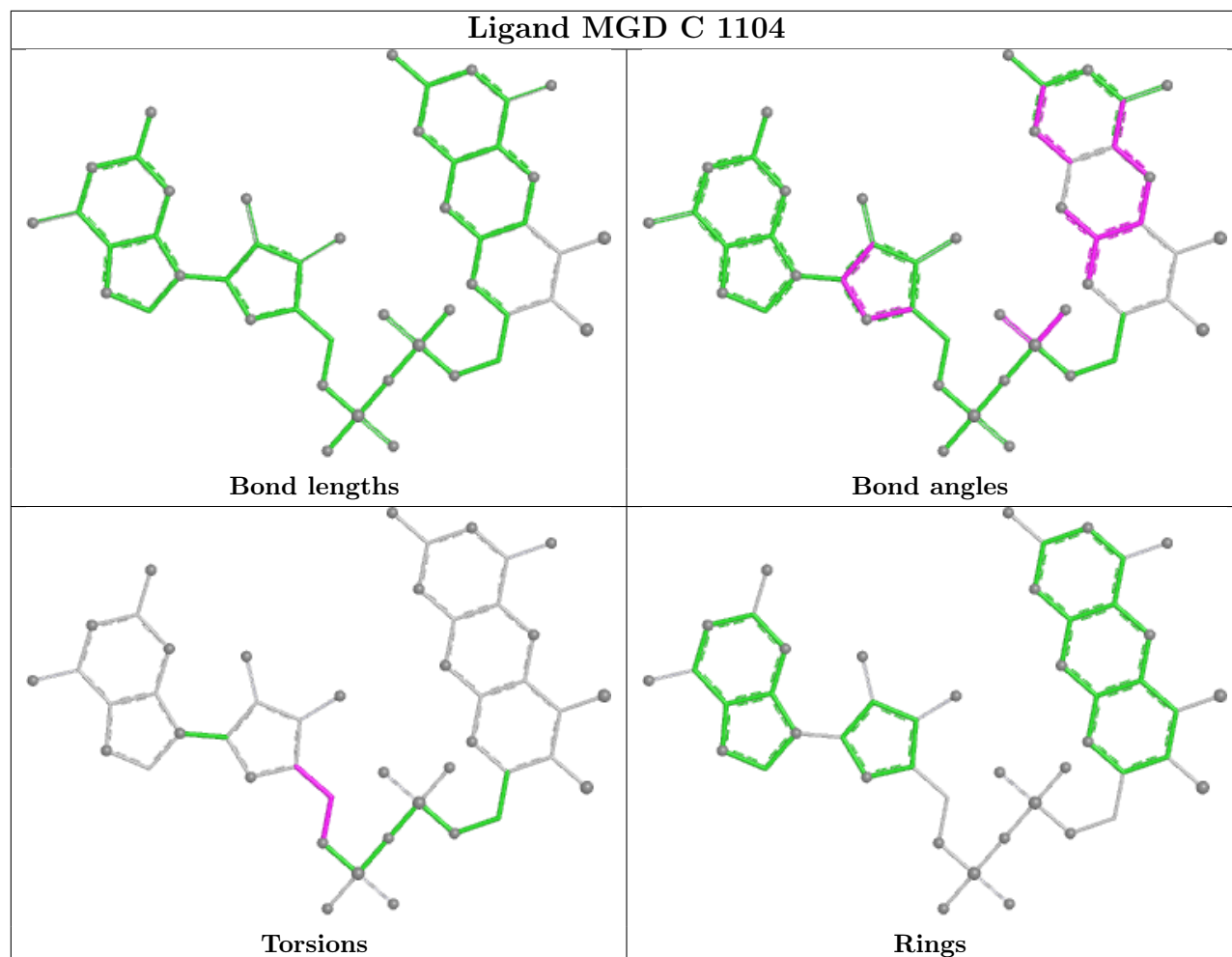
Ligand SF4 D 302

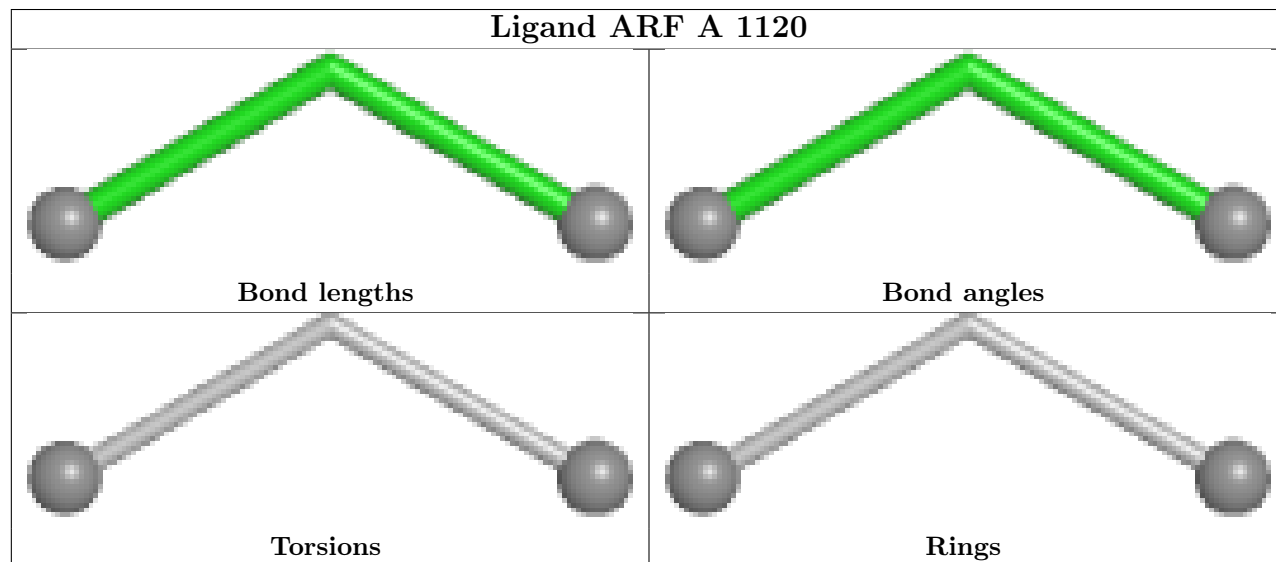
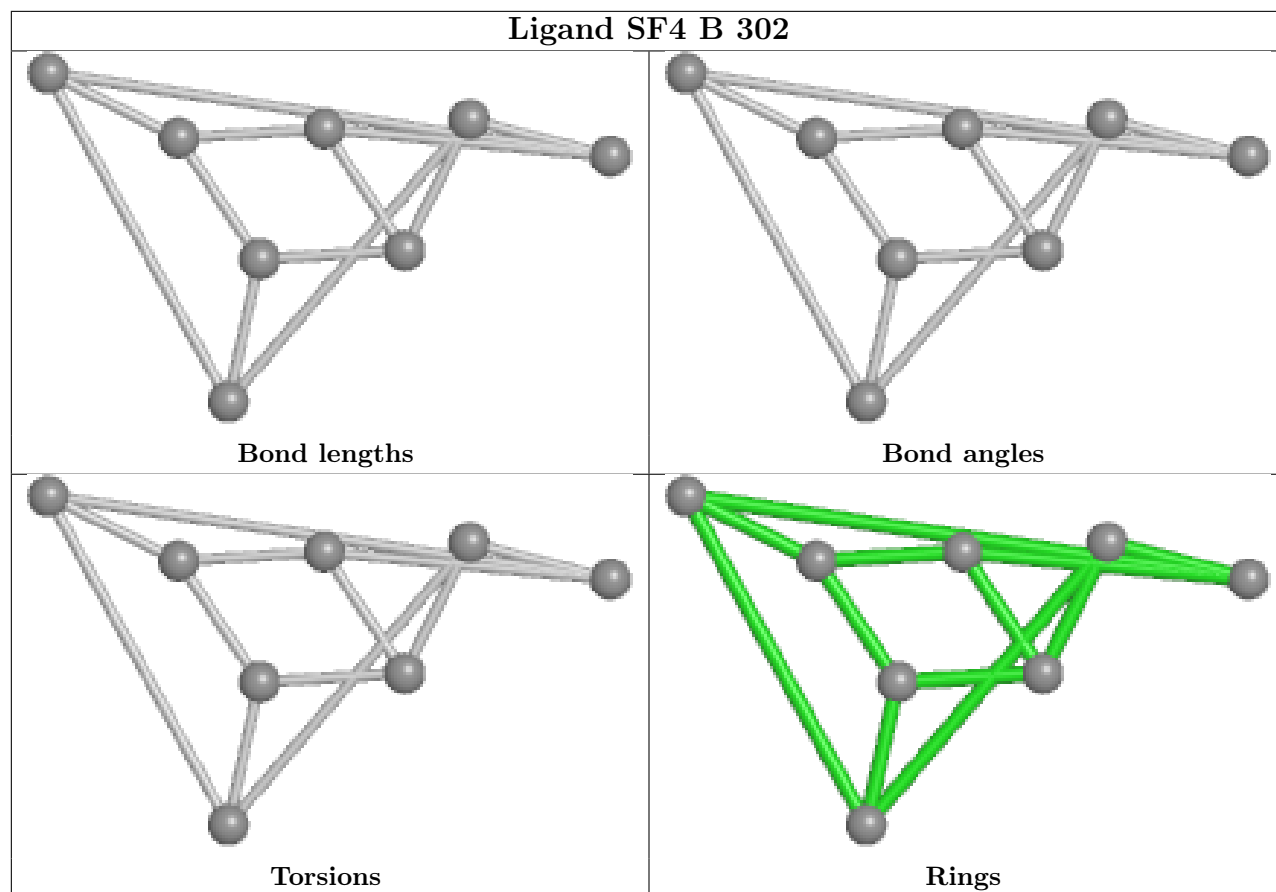


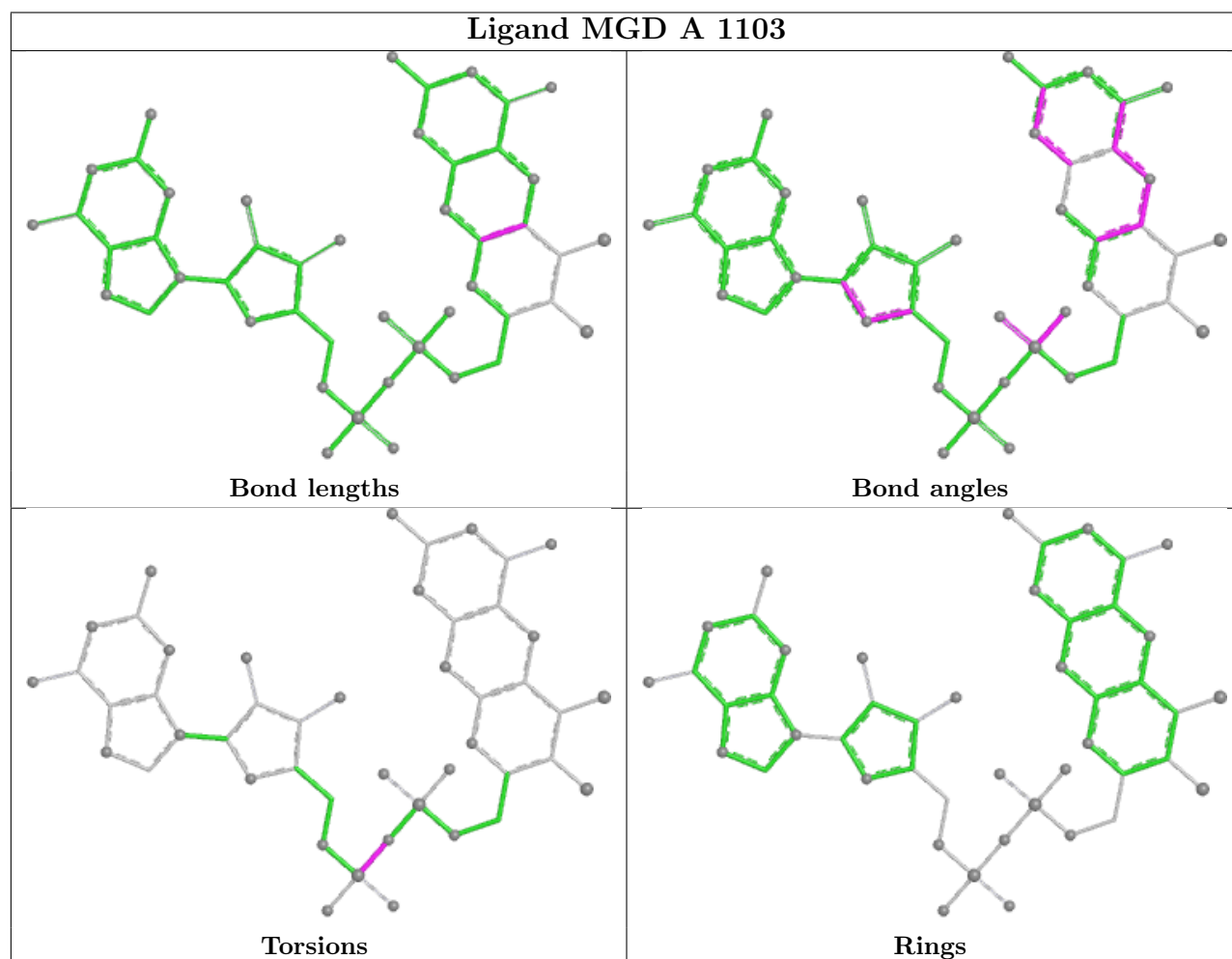
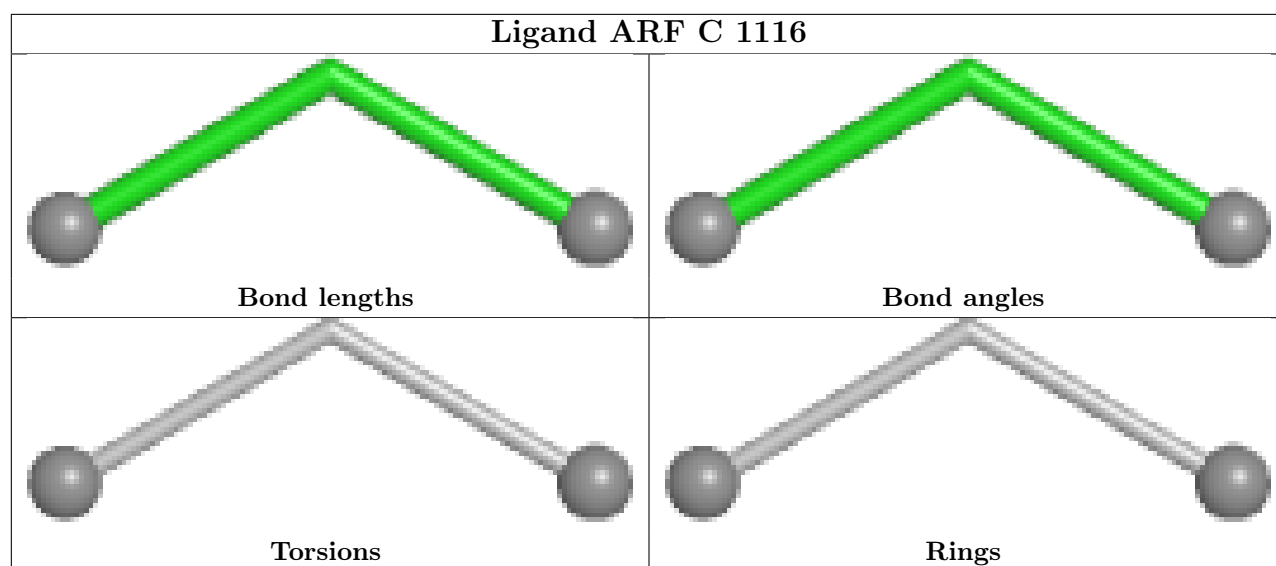




Ligand MGD C 1104







5.7 Other polymers [\(i\)](#)

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, 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	971/1013 (95%)	-0.29	11 (1%) 78 81	6, 14, 27, 67	12 (1%)
1	C	970/1013 (95%)	0.21	24 (2%) 58 62	6, 20, 38, 61	3 (0%)
2	B	214/215 (99%)	-0.22	2 (0%) 81 84	10, 15, 27, 47	4 (1%)
2	D	214/215 (99%)	-0.17	2 (0%) 81 84	11, 16, 30, 56	0
All	All	2369/2456 (96%)	-0.07	39 (1%) 70 74	6, 16, 33, 67	19 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	42	TRP	6.0
1	C	1006	TRP	5.4
1	C	42	TRP	4.5
1	A	339	ALA	4.1
1	A	525	GLY	3.8
1	A	1006	TRP	3.5
1	A	630	ARG	3.0
1	C	339	ALA	2.9
1	C	872	ALA	2.9
1	A	1007	SER	2.9
1	A	116	THR	2.9
1	C	866	GLY	2.9
1	A	340	ASN	2.8
1	C	146	ALA	2.8
1	C	955	VAL	2.7
1	C	796	ASP	2.7
1	A	68	GLN	2.7
1	C	862	PHE	2.7
1	C	854	LYS	2.6
2	D	61	LYS	2.6
1	C	340	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	525	GLY	2.6
1	C	147	ALA	2.5
1	C	874	PRO	2.5
1	C	350	LYS	2.4
1	C	861	ALA	2.4
1	C	954	GLY	2.3
1	C	324	ALA	2.3
1	C	354	CYS	2.3
1	C	1005	VAL	2.2
1	A	350	LYS	2.2
1	C	849	LYS	2.2
2	B	61	LYS	2.2
2	B	115	GLU	2.2
1	C	407	TRP	2.1
1	A	361	LYS	2.0
2	D	63	PRO	2.0
1	C	700	VAL	2.0
1	C	368	LEU	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, 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($\text{\AA}^2$)	Q<0.9
11	ARF	C	1116	3/3	0.75	0.19	40,40,43,46	0
8	GOL	A	1113	6/6	0.80	0.18	29,38,40,44	0
11	ARF	C	1115	3/3	0.81	0.16	36,36,37,39	0
10	PEG	A	1116	7/7	0.81	0.17	39,46,53,56	0
11	ARF	B	305	3/3	0.82	0.18	20,20,28,29	0

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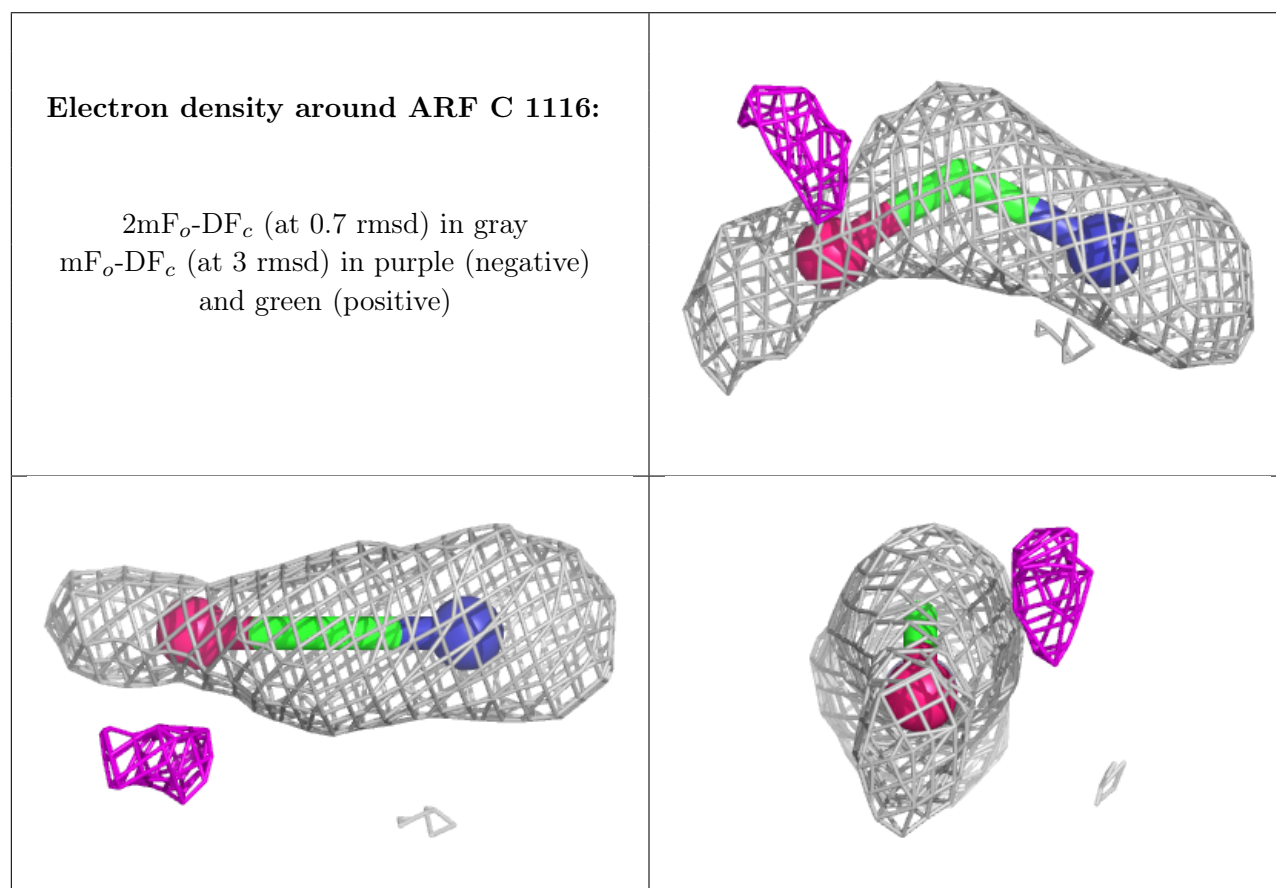
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	PG4	C	1108	13/13	0.82	0.16	33,38,51,52	0
11	ARF	A	1119	3/3	0.82	0.13	40,40,41,41	0
9	PG4	A	1110	13/13	0.84	0.16	32,36,45,48	0
10	PEG	C	1107	7/7	0.84	0.13	22,25,31,34	0
8	GOL	A	1115	6/6	0.85	0.15	25,31,41,43	0
8	GOL	A	1118	6/6	0.86	0.14	26,37,42,47	0
9	PG4	A	1112	13/13	0.86	0.14	33,40,50,53	0
11	ARF	A	1120	3/3	0.86	0.14	31,31,32,35	0
9	PG4	B	306	13/13	0.87	0.15	33,38,46,53	0
11	ARF	C	1114	3/3	0.87	0.14	50,50,52,54	0
8	GOL	A	1114	6/6	0.91	0.09	26,29,31,33	0
11	ARF	C	1113	3/3	0.92	0.11	28,28,31,33	0
8	GOL	C	1109	6/6	0.92	0.10	31,34,35,39	0
8	GOL	C	1110	6/6	0.94	0.07	17,21,21,22	0
8	GOL	C	1106	6/6	0.95	0.08	18,22,23,24	0
8	GOL	C	1111	6/6	0.95	0.08	25,27,29,30	0
8	GOL	C	1112	6/6	0.95	0.08	23,27,29,29	0
8	GOL	A	1117	6/6	0.95	0.07	14,18,19,20	0
8	GOL	A	1111	6/6	0.96	0.07	22,24,27,30	0
8	GOL	A	1108	6/6	0.96	0.07	19,21,24,26	0
8	GOL	A	1109	6/6	0.96	0.06	15,19,20,22	0
12	EDO	B	307	4/4	0.96	0.06	13,15,16,16	0
12	EDO	D	305	4/4	0.96	0.06	14,16,16,16	0
8	GOL	A	1107	6/6	0.97	0.06	18,20,20,22	0
5	MGD	C	1104	47/47	0.98	0.05	12,15,17,18	0
3	H2S	A	1101	1/1	0.99	0.06	15,15,15,15	0
6	SF4	B	302	8/8	0.99	0.03	12,13,14,14	0
7	CL	A	1106	1/1	0.99	0.04	23,23,23,23	0
3	H2S	C	1101	1/1	0.99	0.08	21,21,21,21	0
4	W	C	1102	1/1	0.99	0.03	15,15,15,15	0
5	MGD	A	1103	47/47	0.99	0.03	8,9,10,12	0
5	MGD	A	1104	47/47	0.99	0.03	8,10,11,13	0
5	MGD	C	1103	47/47	0.99	0.04	11,14,17,19	0
7	CL	D	304	1/1	1.00	0.02	14,14,14,14	0
6	SF4	B	301	8/8	1.00	0.02	10,11,11,11	0
4	W	A	1102	1/1	1.00	0.01	10,10,10,10	0
6	SF4	B	303	8/8	1.00	0.02	11,11,11,12	0
6	SF4	C	1105	8/8	1.00	0.02	11,12,12,12	0
6	SF4	D	301	8/8	1.00	0.02	11,12,12,12	0
6	SF4	D	302	8/8	1.00	0.02	12,13,14,14	0
6	SF4	D	303	8/8	1.00	0.02	12,12,12,12	0
6	SF4	A	1105	8/8	1.00	0.01	9,9,10,10	0

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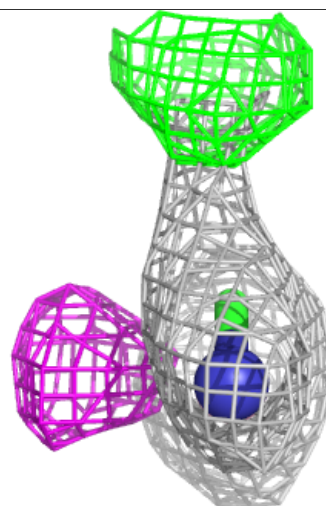
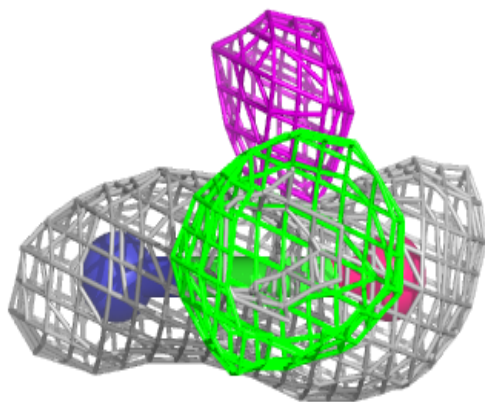
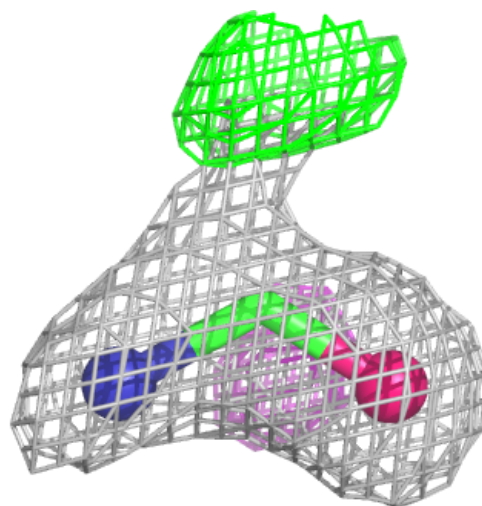
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	CL	B	304	1/1	1.00	0.02	12,12,12,12	0

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.



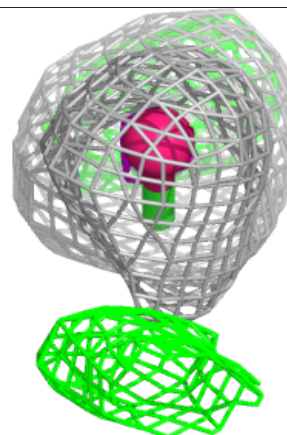
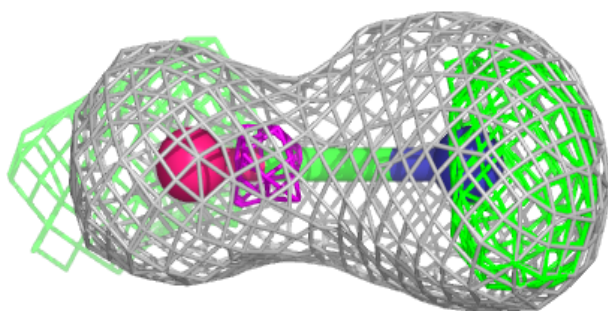
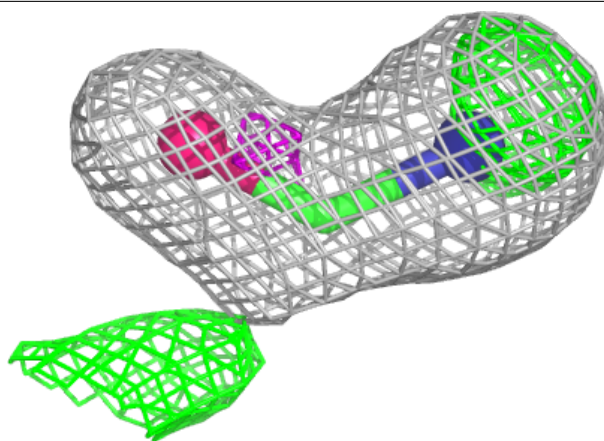
Electron density around ARF C 1115:

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



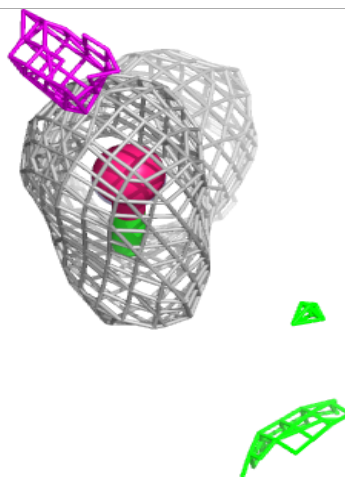
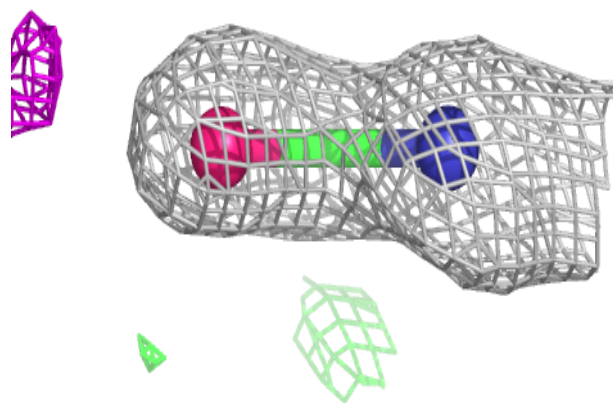
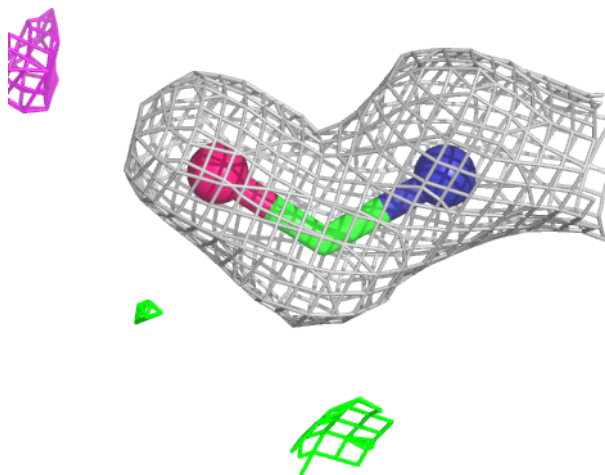
Electron density around ARF B 305:

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



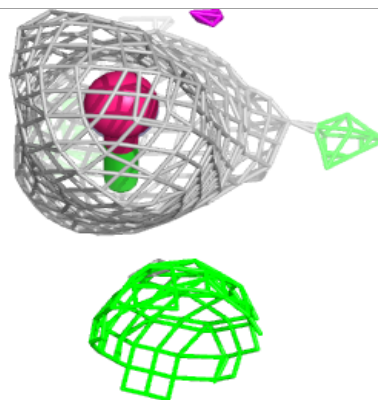
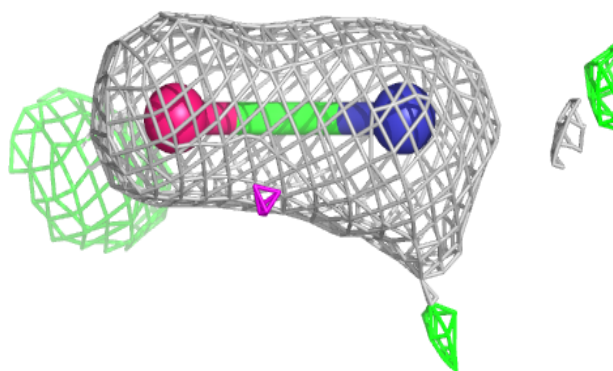
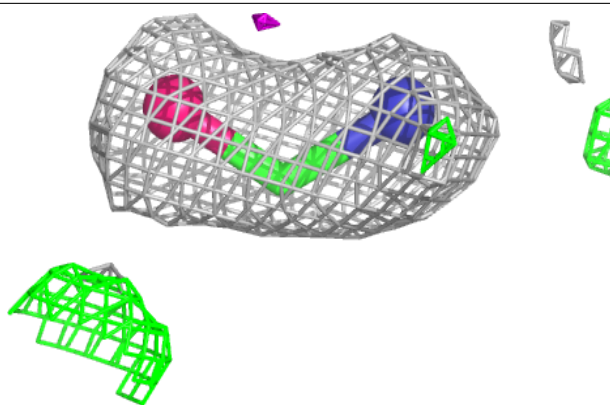
Electron density around ARF A 1119:

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

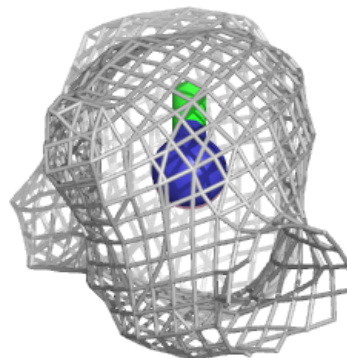
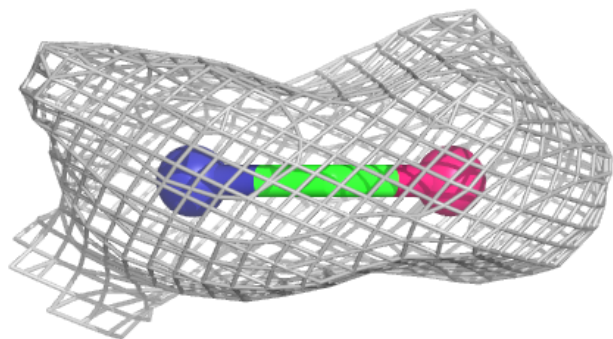
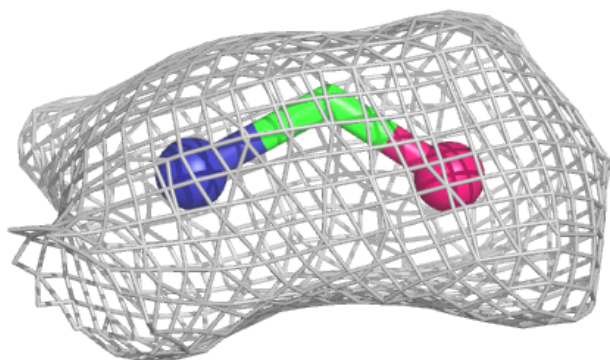


Electron density around ARF A 1120:

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

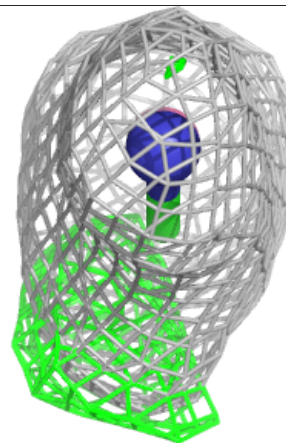
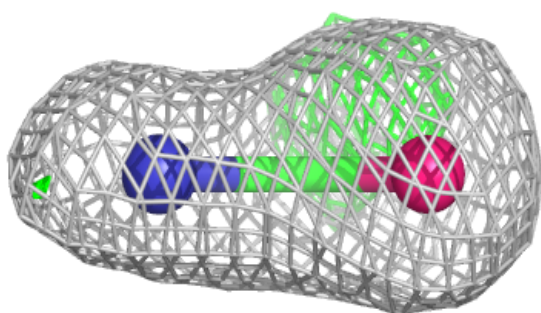
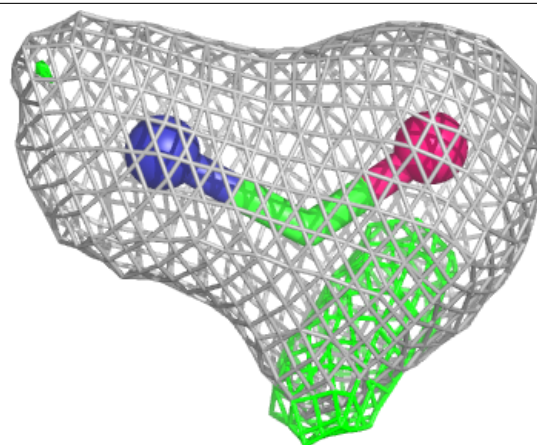
**Electron density around ARF C 1114:**

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

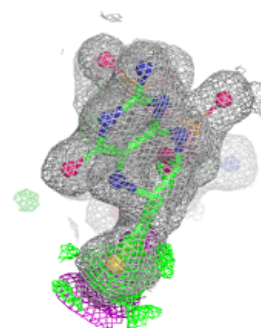
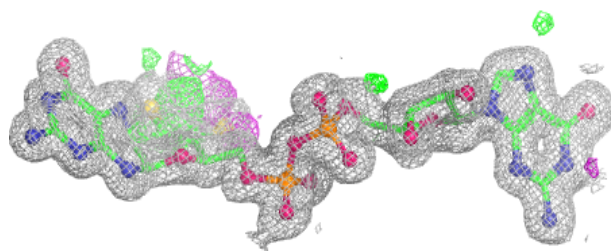
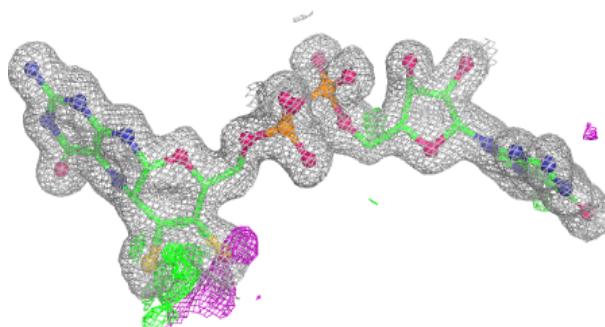


Electron density around ARF C 1113:

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

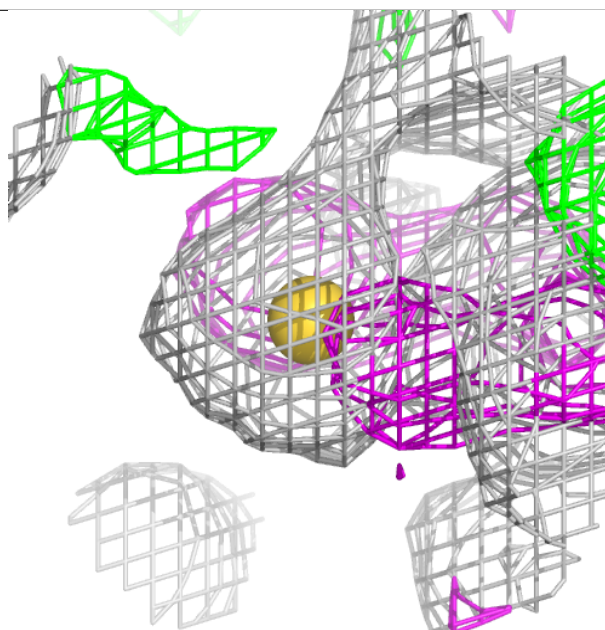
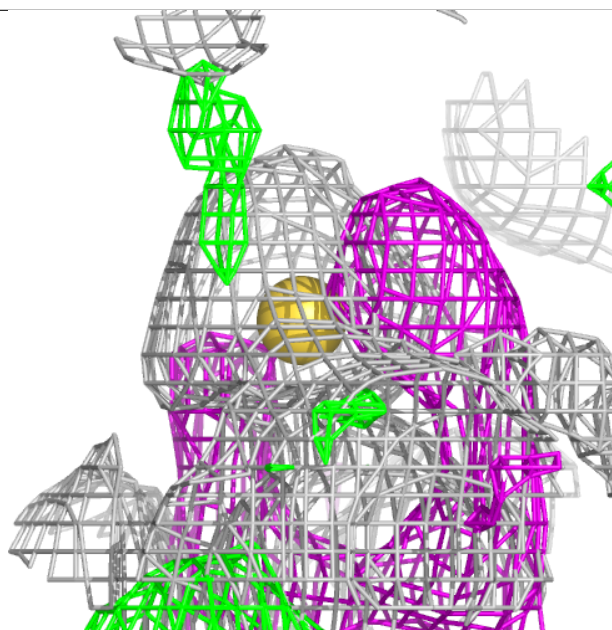
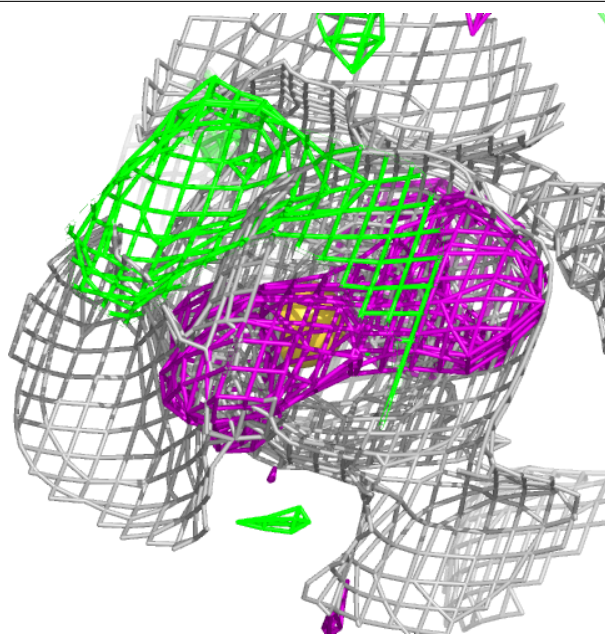
**Electron density around MGD C 1104:**

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



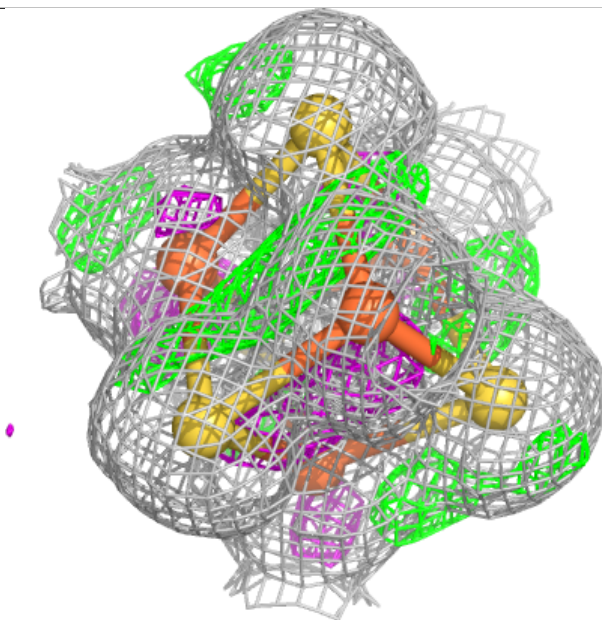
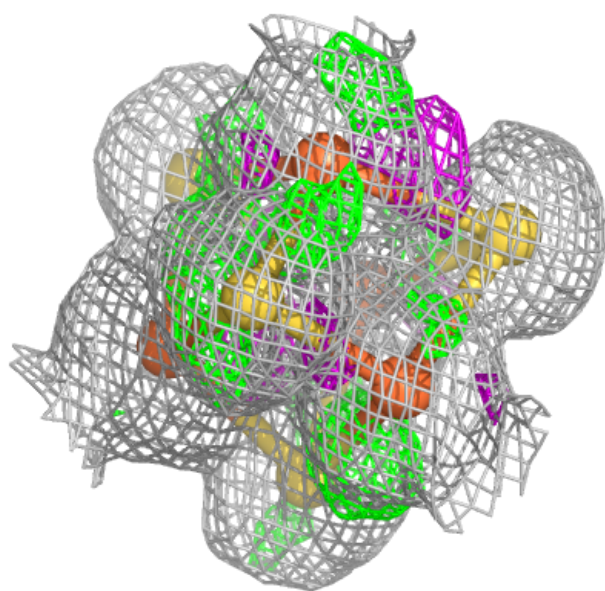
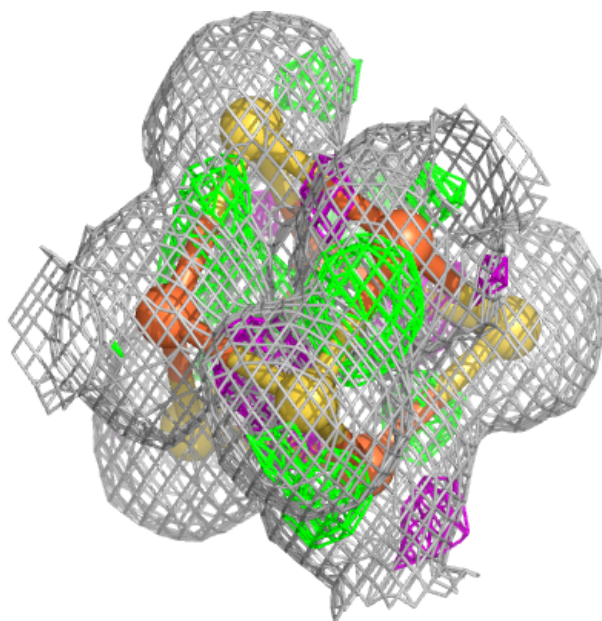
Electron density around H2S A 1101:

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



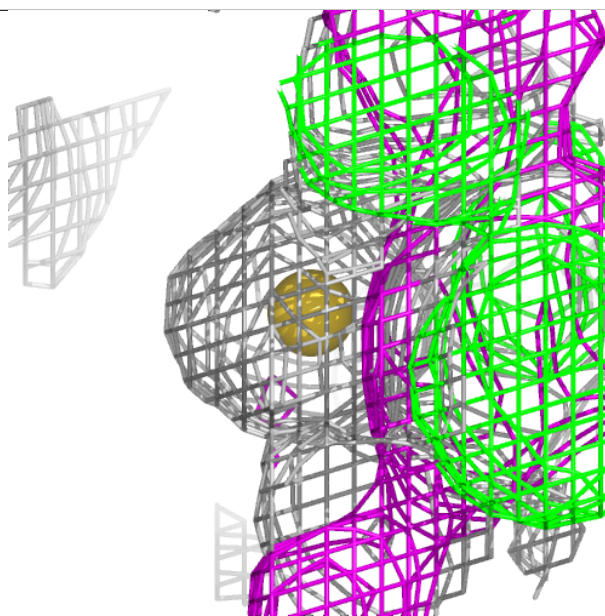
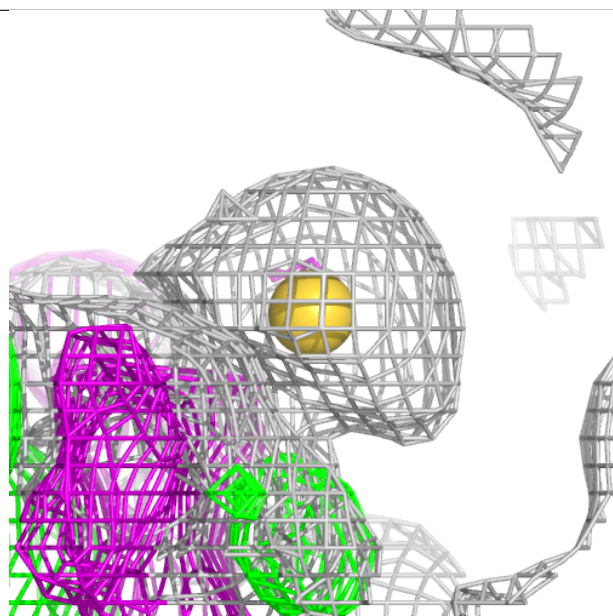
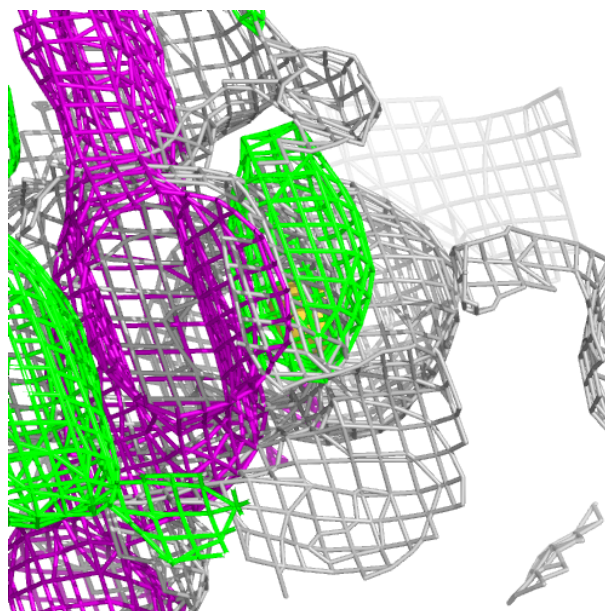
Electron density around SF4 B 302:

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



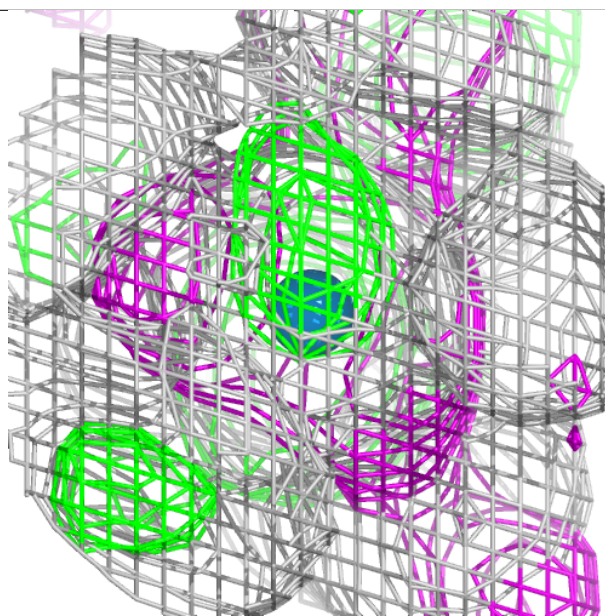
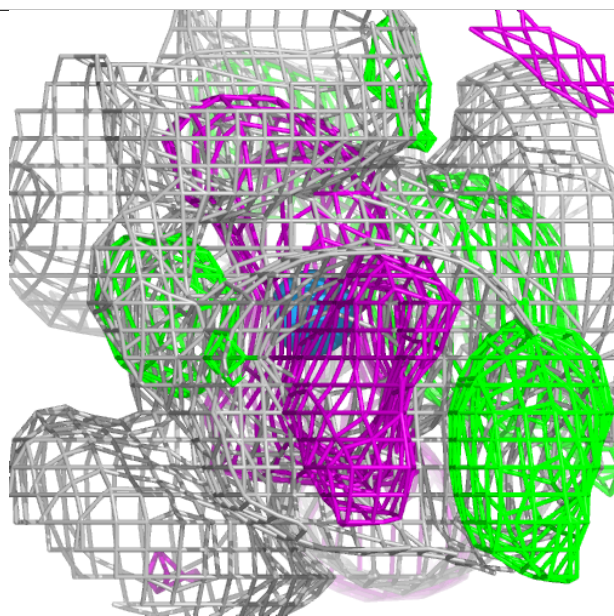
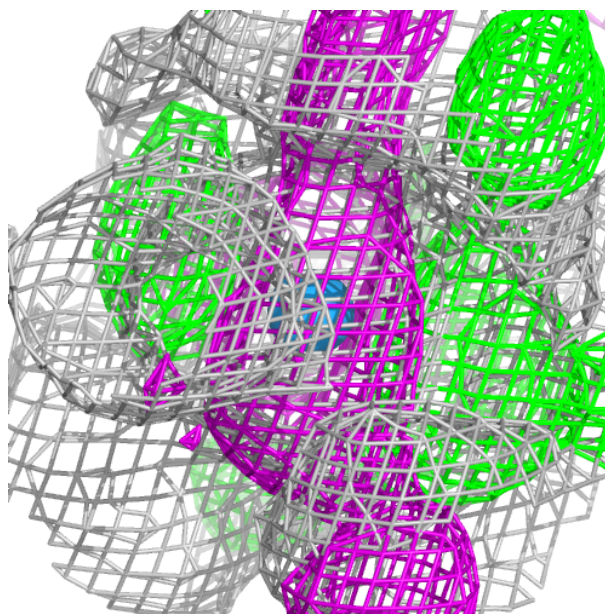
Electron density around H2S C 1101:

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



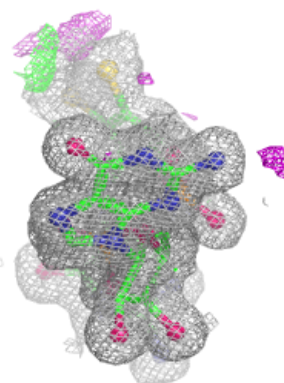
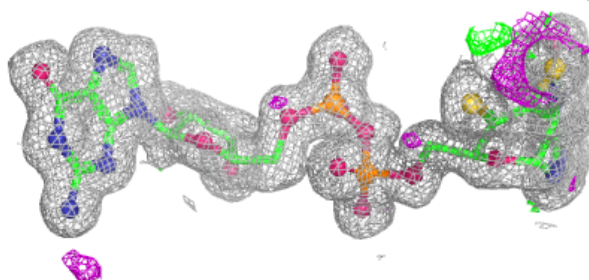
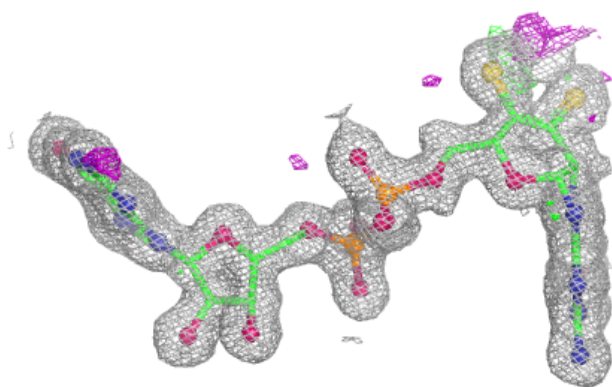
Electron density around W C 1102:

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

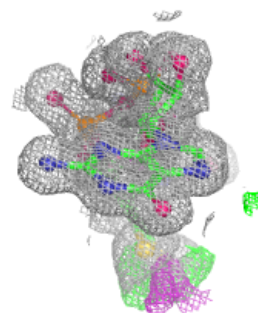
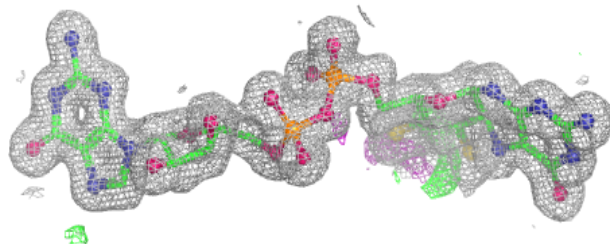
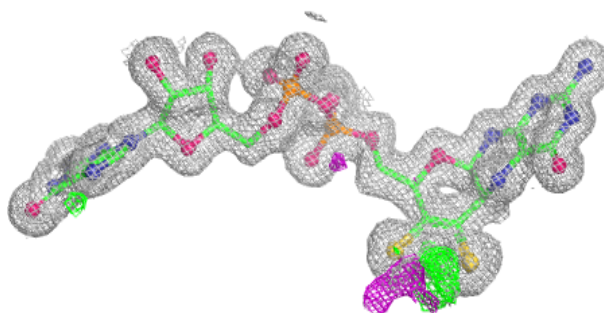


Electron density around MGD A 1103:

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

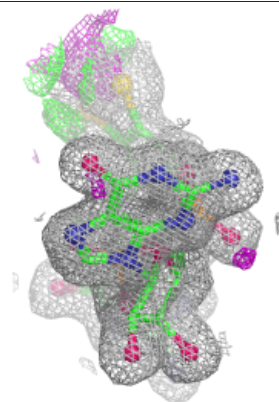
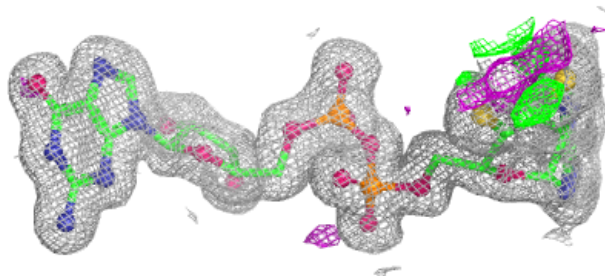
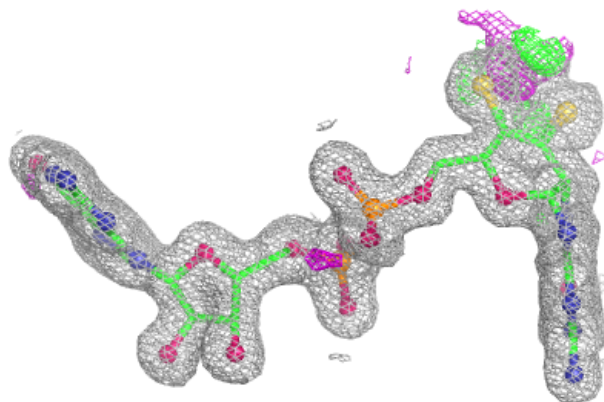
**Electron density around MGD A 1104:**

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



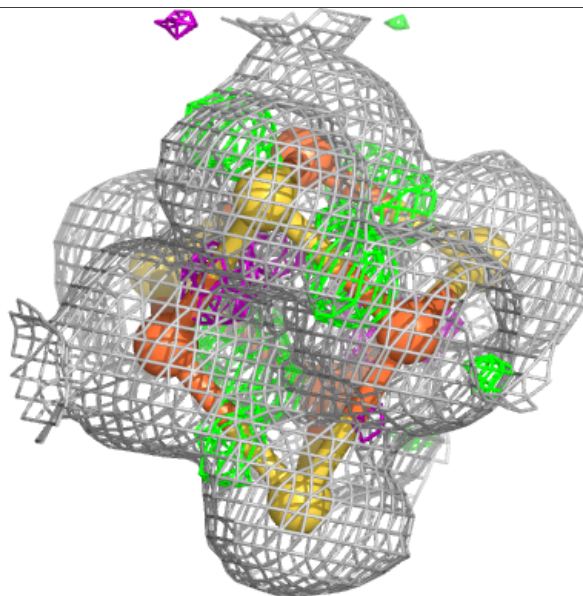
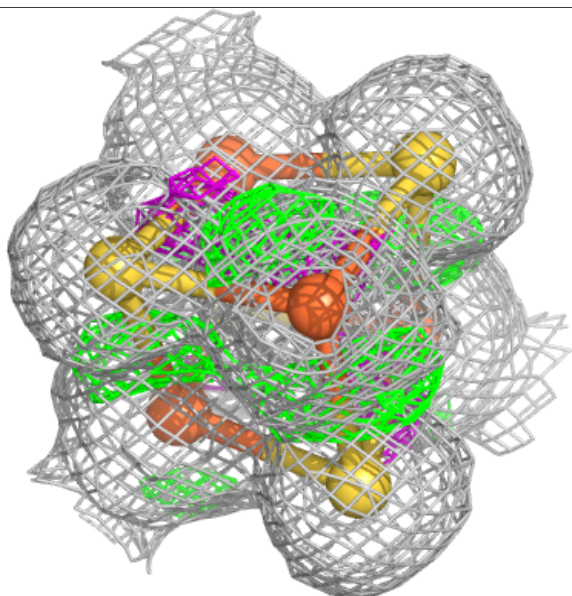
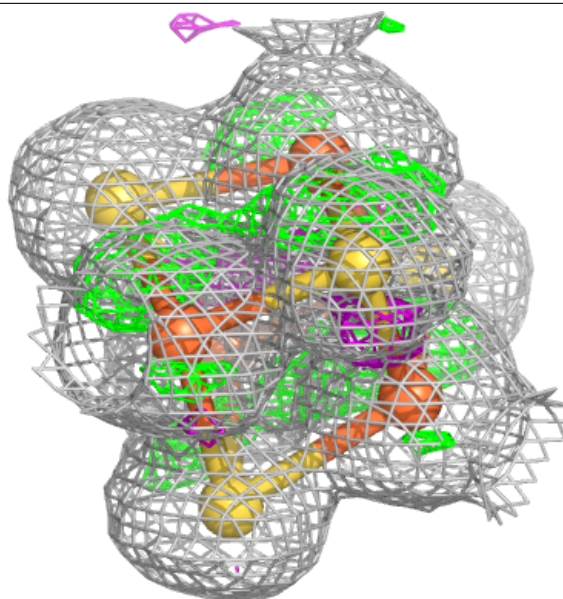
Electron density around MGD C 1103:

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



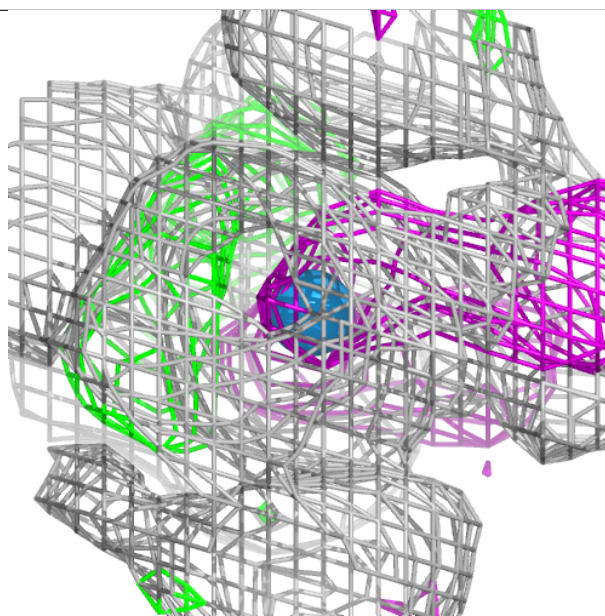
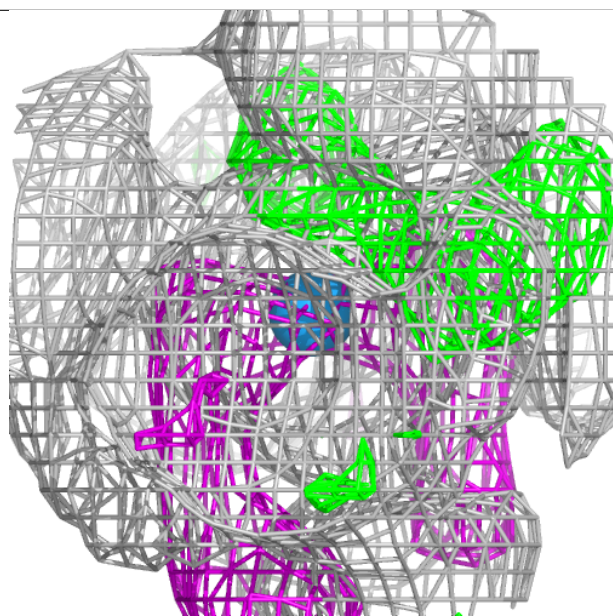
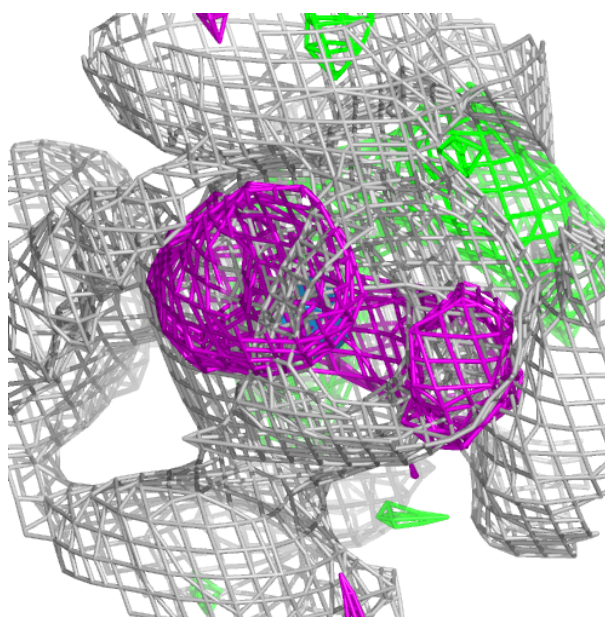
Electron density around SF4 B 301:

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



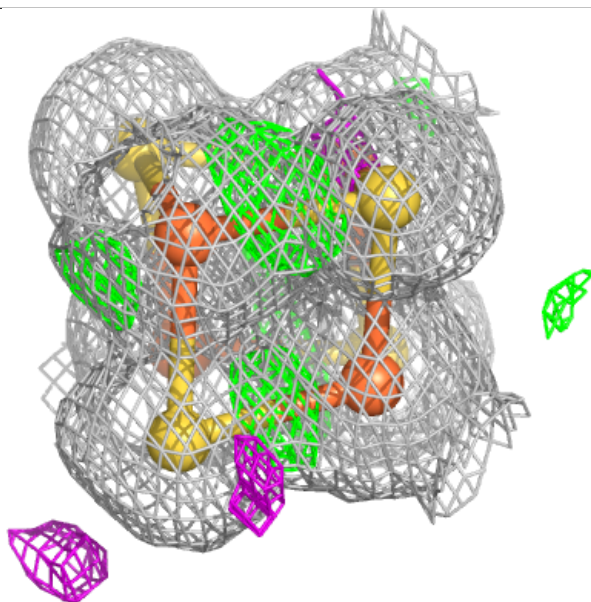
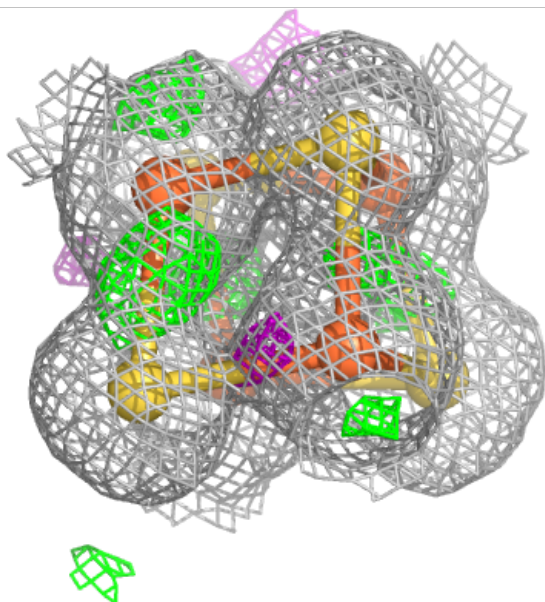
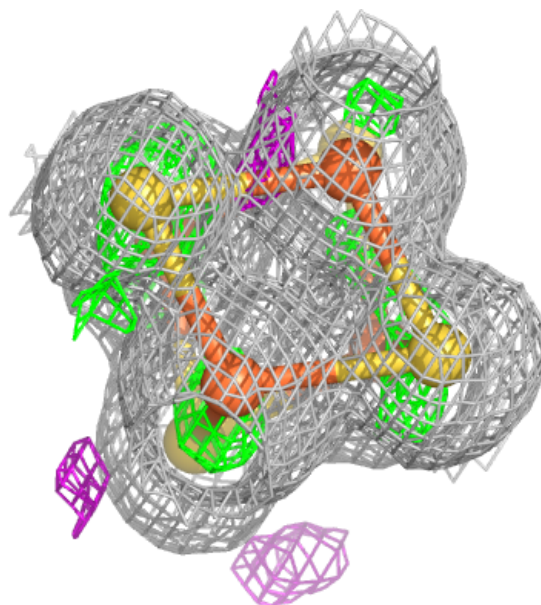
Electron density around W A 1102:

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



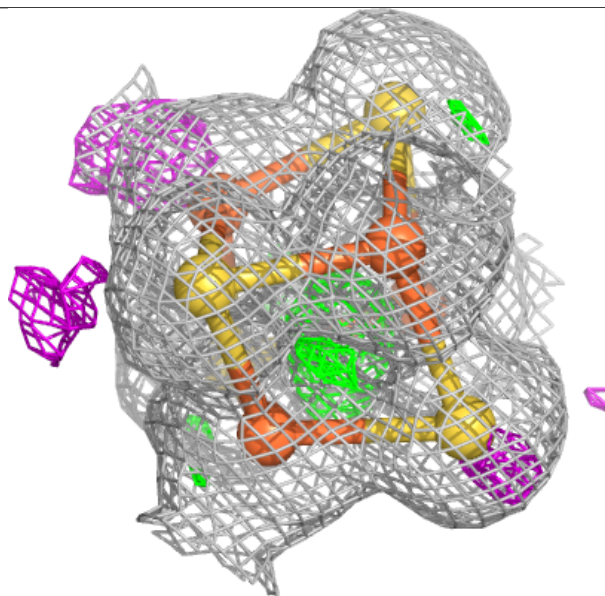
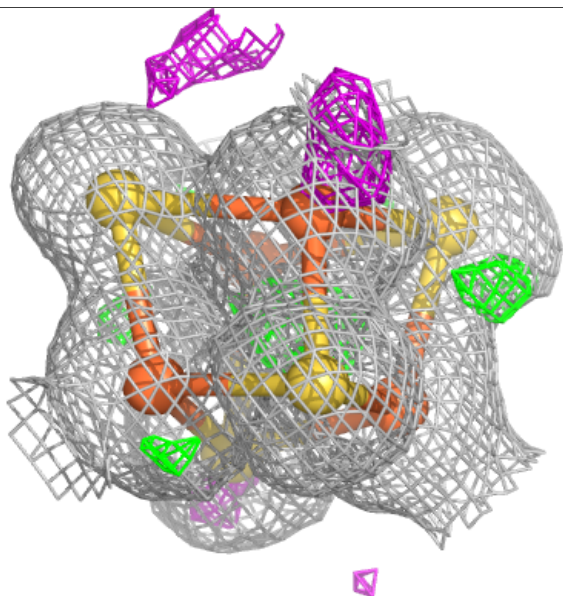
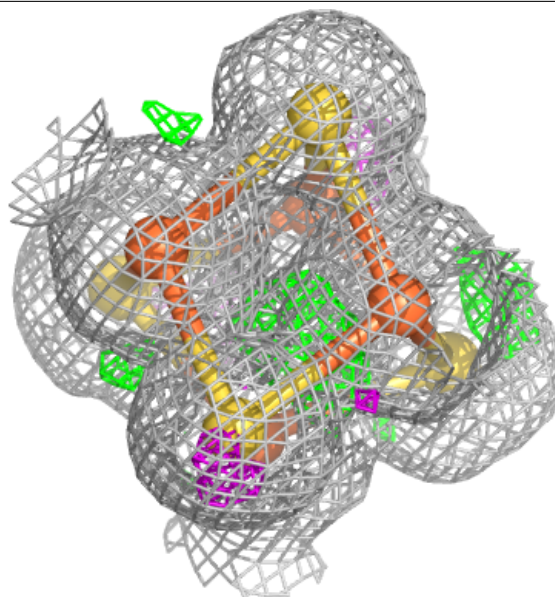
Electron density around SF4 B 303:

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



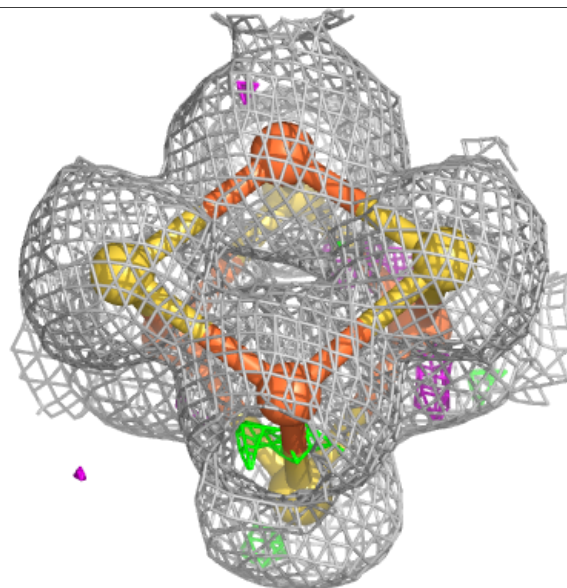
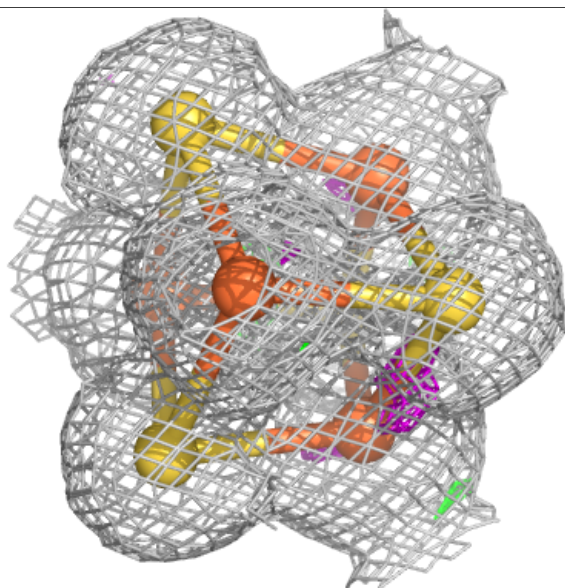
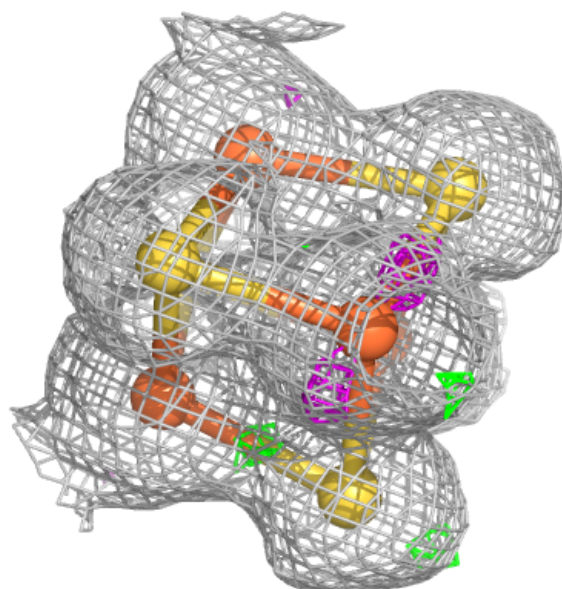
Electron density around SF4 C 1105:

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



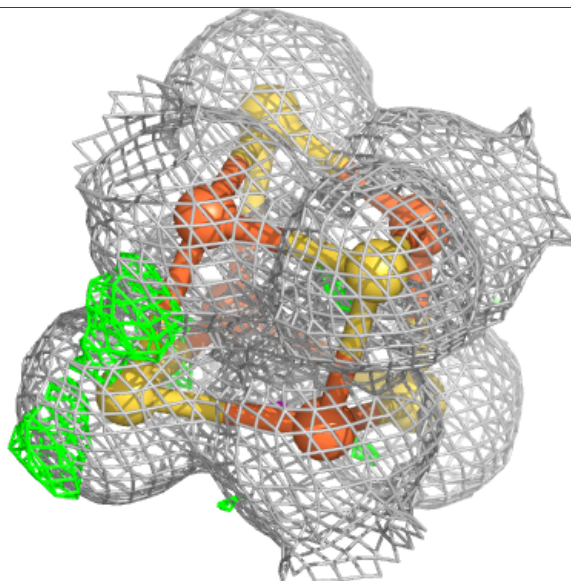
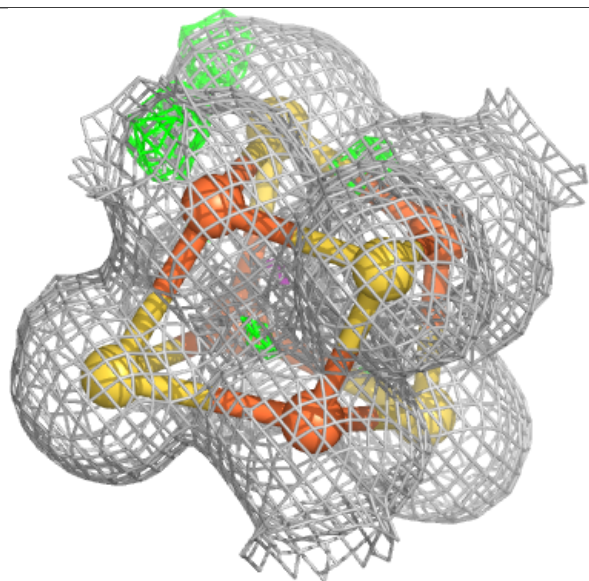
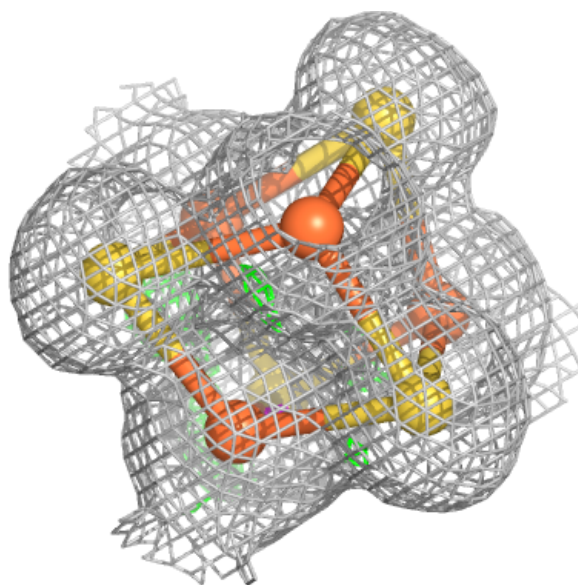
Electron density around SF4 D 301:

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



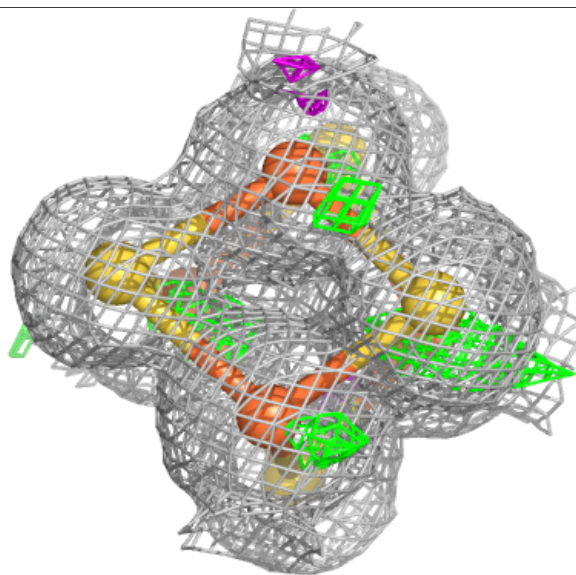
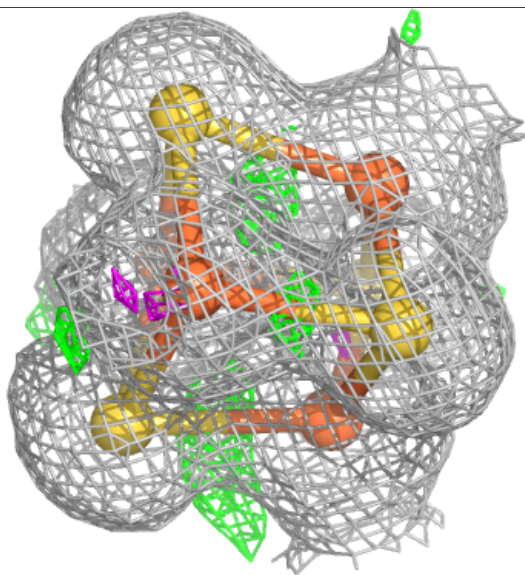
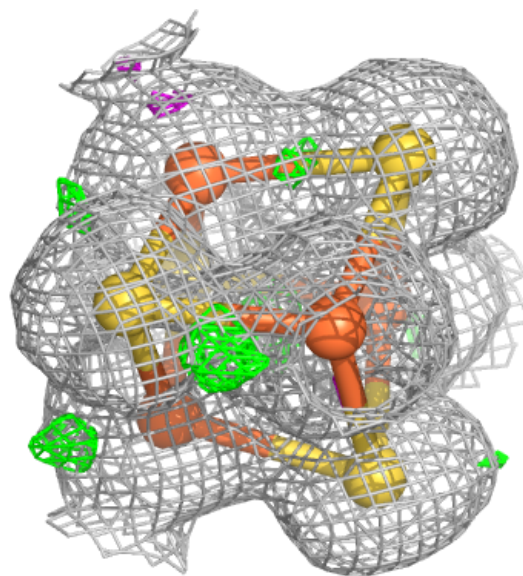
Electron density around SF4 D 302:

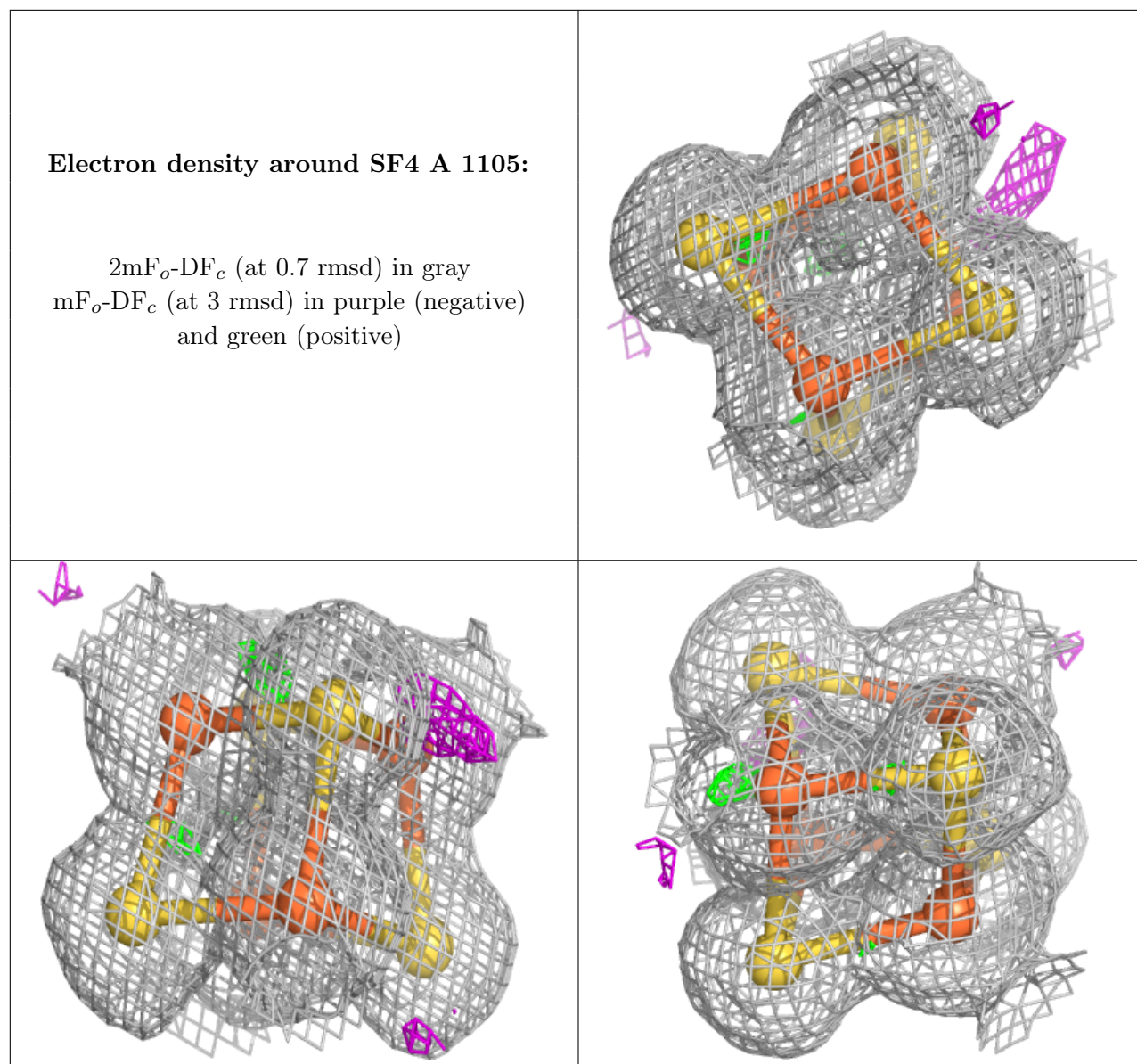
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around SF4 D 303:

$2mF_o-DF_c$ (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.