



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 29, 2020 – 09:39 pm BST

PDB ID : 3OR1  
Title : Crystal structure of dissimilatory sulfite reductase I (DsrI)  
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Deposited on : 2010-09-06  
Resolution : 1.76 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

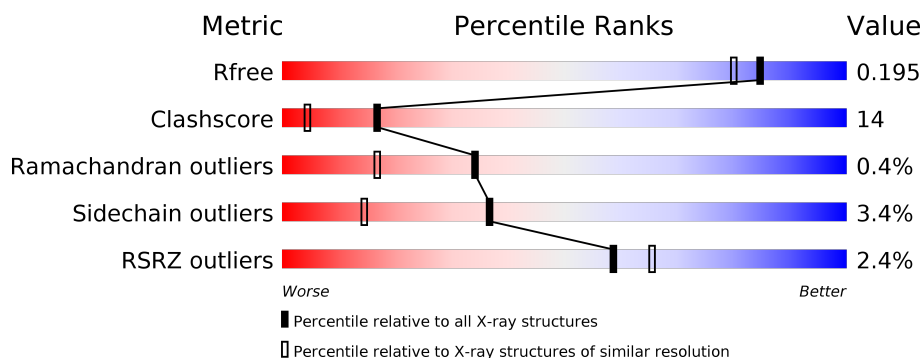
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.76 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))
$R_{free}$	130704	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	437	<div> <div>2%</div> <div> <div></div> <div>53%</div> <div>38%</div> <div>8%</div> </div> <div></div> </div>
1	D	437	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>25%</div> <div></div> </div> <div></div> </div>
2	B	386	<div> <div>2%</div> <div> <div></div> <div>69%</div> <div>25%</div> <div>5%</div> </div> <div></div> </div>
2	E	386	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>22%</div> <div></div> </div> <div></div> </div>
3	C	105	<div> <div>12%</div> <div> <div></div> <div>60%</div> <div>35%</div> <div></div> </div> <div></div> </div>
3	F	105	<div> <div>5%</div> <div> <div></div> <div>68%</div> <div>24%</div> <div>7%</div> </div> <div></div> </div>

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 15467 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 Sulfite reductase alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	S	0	0	0
			3422	2163	585	646	28			
1	D	435	Total	C	N	O	S	0	0	0
			3422	2163	585	646	28			

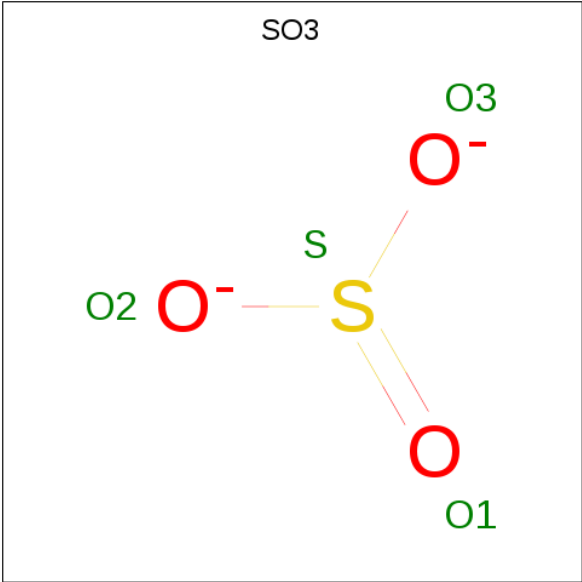
- Molecule 2 is a protein called Sulfite reductase beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	385	Total	C	N	O	S	0	0	0
			2999	1912	519	542	26			
2	E	385	Total	C	N	O	S	0	0	0
			2999	1912	519	542	26			

- Molecule 3 is a protein called Sulfite reductase gama.

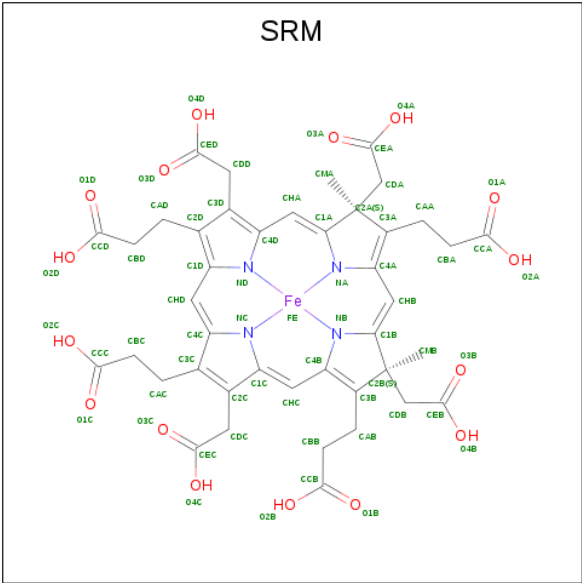
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	104	Total	C	N	O	S	0	0	0
			787	508	125	149	5			
3	F	104	Total	C	N	O	S	0	0	0
			787	508	125	149	5			

- Molecule 4 is SULFITE ION (three-letter code: SO3) (formula: O<sub>3</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			4	3	1		
4	D	1	Total	O	S	0	0
			4	3	1		

- Molecule 5 is SIROHEME (three-letter code: SRM) (formula: C<sub>42</sub>H<sub>42</sub>FeN<sub>4</sub>O<sub>16</sub>).



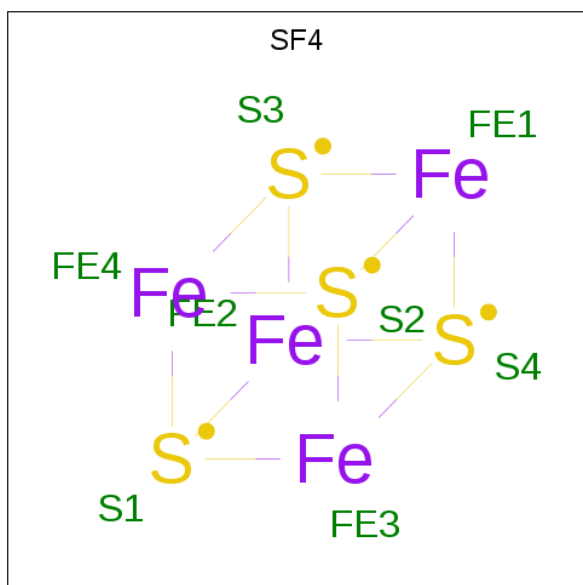
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	Fe	N	O	0
			63	42	1	4	16	
5	B	1	Total	C	N	O	0	0
			62	42	4	16		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	D	1	Total	C	Fe	N	O	0	0
			63	42	1	4	16		
5	E	1	Total	C	N	O	0	0	
			62	42	4	16			

- Molecule 6 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe<sub>4</sub>S<sub>4</sub>).



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

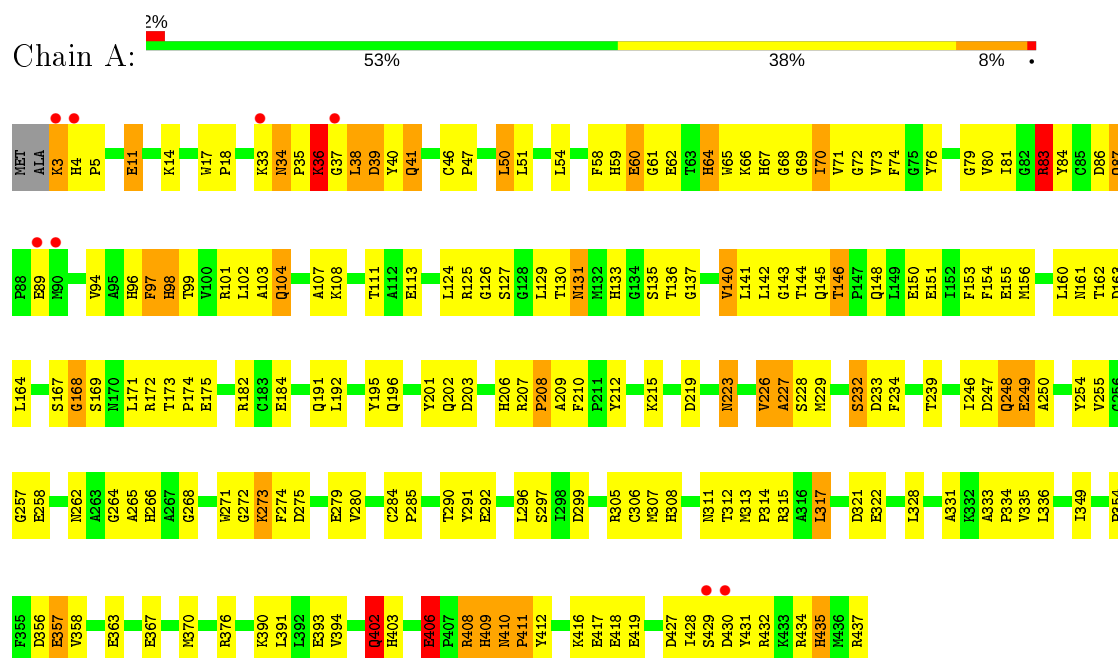
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	164	Total 164	O 164	0	0
7	B	160	Total 160	O 160	0	0
7	C	12	Total 12	O 12	0	0
7	D	220	Total 220	O 220	0	0
7	E	155	Total 155	O 155	0	0
7	F	18	Total 18	O 18	0	0

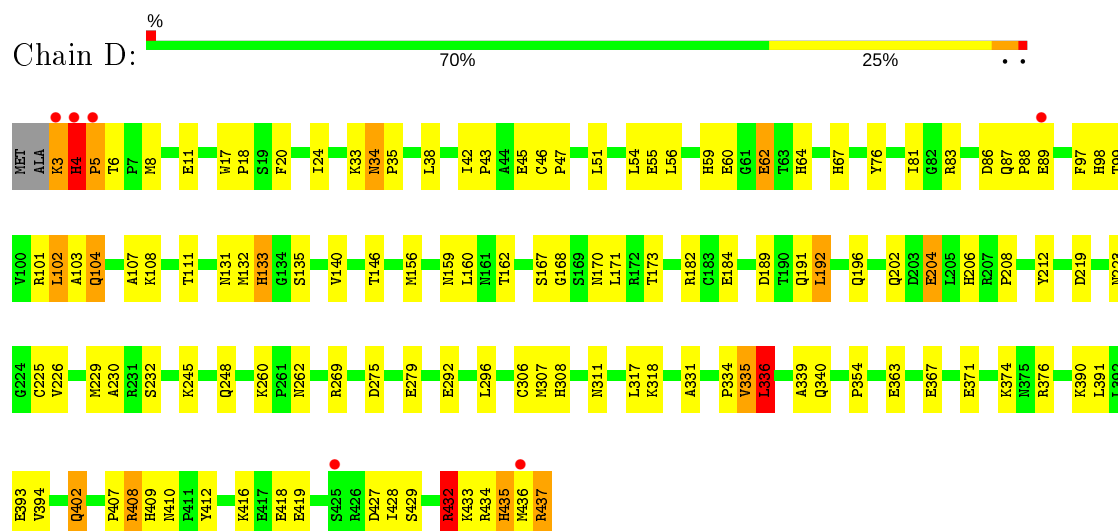
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

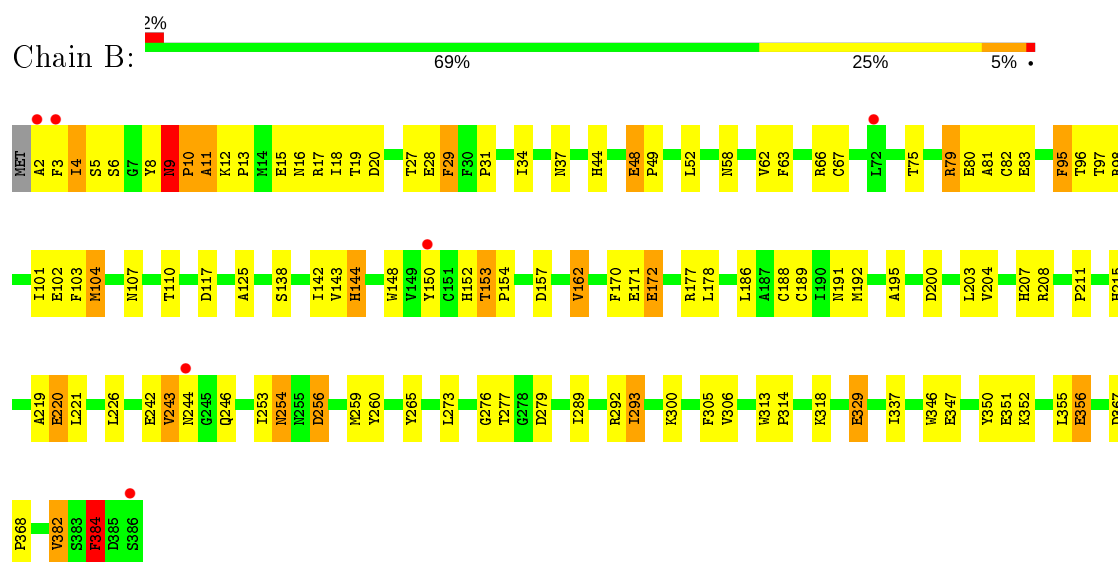
#### • Molecule 1: Sulfite reductase alpha



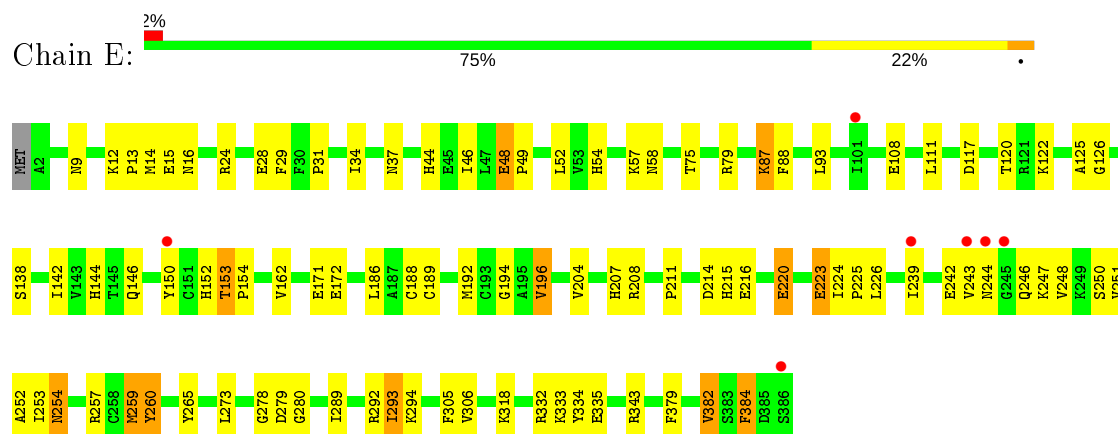
#### • Molecule 1: Sulfite reductase alpha



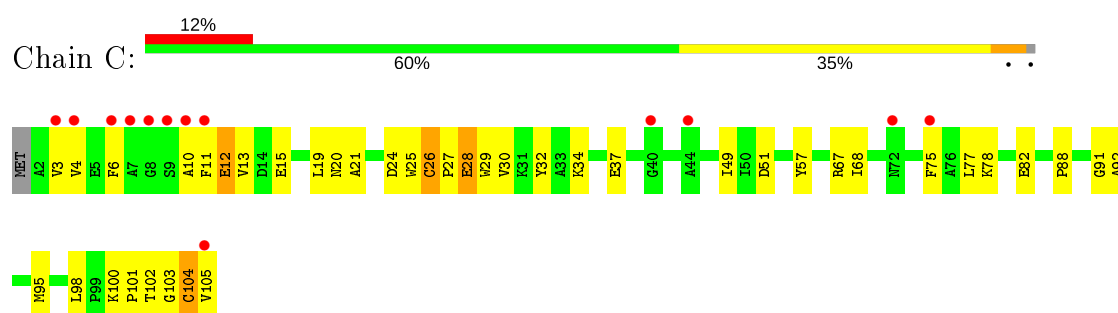
#### • Molecule 2: Sulfite reductase beta



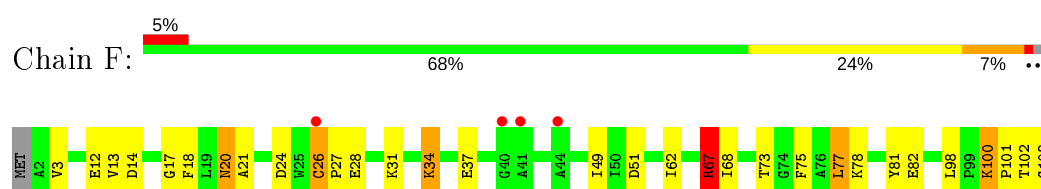
• Molecule 2: Sulfite reductase beta



• Molecule 3: Sulfite reductase gamma



• Molecule 3: Sulfite reductase gamma





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	114.72Å 60.48Å 132.91Å 90.00° 94.30° 90.00°	Depositor
Resolution (Å)	30.00 – 1.76 29.48 – 1.70	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-1.76) 99.0 (29.48-1.70)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.96 (at 1.70Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.193 , 0.212 0.192 , 0.195	Depositor DCC
$R_{free}$ test set	10018 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.0	Xtriage
Anisotropy	0.292	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 40.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	15467	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SF4, SO3, SRM

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	2.36	257/3512 (7.3%)	1.23	27/4751 (0.6%)
1	D	1.23	58/3512 (1.7%)	1.08	21/4751 (0.4%)
2	B	1.93	133/3076 (4.3%)	0.99	12/4170 (0.3%)
2	E	1.52	67/3076 (2.2%)	0.84	11/4170 (0.3%)
3	C	1.26	9/808 (1.1%)	0.77	1/1090 (0.1%)
3	F	1.60	17/808 (2.1%)	1.27	4/1090 (0.4%)
All	All	1.78	541/14792 (3.7%)	1.05	76/20022 (0.4%)

The worst 5 of 541 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	26	CYS	CB-SG	-20.23	1.47	1.82
2	B	384	PHE	CE1-CZ	-19.93	0.99	1.37
2	E	220	GLU	CB-CG	-19.51	1.15	1.52
1	A	406	GLU	CB-CG	-18.91	1.16	1.52
2	B	306	VAL	CB-CG2	-17.20	1.16	1.52

The worst 5 of 76 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	83	ARG	NE-CZ-NH2	-28.06	106.27	120.30
3	F	67	ARG	NE-CZ-NH1	25.12	132.86	120.30
1	D	432	ARG	NE-CZ-NH2	-22.41	109.09	120.30
1	D	437	ARG	NE-CZ-NH2	-18.76	110.92	120.30
3	F	67	ARG	NE-CZ-NH2	-17.91	111.34	120.30

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	3422	0	3275	99	0
1	D	3422	0	3275	113	0
2	B	2999	0	2963	79	0
2	E	2999	0	2963	78	0
3	C	787	0	757	49	0
3	F	787	0	758	46	0
4	A	4	0	0	0	0
4	D	4	0	0	0	0
5	A	63	0	33	12	0
5	B	62	0	35	8	0
5	D	63	0	35	10	0
5	E	62	0	34	5	0
6	A	16	0	0	0	0
6	B	16	0	0	0	0
6	D	16	0	0	0	0
6	E	16	0	0	0	0
7	A	164	0	0	10	0
7	B	160	0	0	4	0
7	C	12	0	0	1	0
7	D	220	0	0	8	0
7	E	155	0	0	3	0
7	F	18	0	0	0	0
All	All	15467	0	14128	408	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 14.

The worst 5 of 408 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)
5:A:580:SRM:HHA	3:C:104:CYS:SG	1.36	1.62
5:D:582:SRM:HHA	3:F:104:CYS:SG	1.47	1.51
5:A:580:SRM:CHA	3:C:104:CYS:SG	2.10	1.39
7:A:6631:HOH:O	3:C:105:VAL:CG2	1.75	1.34
5:D:582:SRM:CHA	3:F:104:CYS:SG	2.20	1.29

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	433/437 (99%)	418 (96%)	15 (4%)	0	100	100
1	D	433/437 (99%)	422 (98%)	10 (2%)	1 (0%)	47	29
2	B	383/386 (99%)	365 (95%)	16 (4%)	2 (0%)	29	12
2	E	383/386 (99%)	369 (96%)	12 (3%)	2 (0%)	29	12
3	C	102/105 (97%)	98 (96%)	3 (3%)	1 (1%)	15	4
3	F	102/105 (97%)	99 (97%)	2 (2%)	1 (1%)	15	4
All	All	1836/1856 (99%)	1771 (96%)	58 (3%)	7 (0%)	34	17

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	4	HIS
3	C	34	LYS
3	F	34	LYS
2	E	293	ILE
2	E	153	THR

### 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	364/365 (100%)	350 (96%)	14 (4%)	33	11

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	364/365 (100%)	352 (97%)	12 (3%)	38	15
2	B	324/325 (100%)	312 (96%)	12 (4%)	34	12
2	E	324/325 (100%)	319 (98%)	5 (2%)	65	49
3	C	79/80 (99%)	75 (95%)	4 (5%)	24	6
3	F	79/80 (99%)	74 (94%)	5 (6%)	18	3
All	All	1534/1540 (100%)	1482 (97%)	52 (3%)	37	14

5 of 52 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	329	GLU
3	C	77	LEU
3	F	34	LYS
2	B	356	GLU
3	C	12	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 54 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	254	ASN
1	D	104	GLN
2	E	207	HIS
2	B	263	ASN
1	D	64	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

14 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	SO3	A	6575	5	1,3,3	0.49	0	0,3,3	0.00	-
5	SRM	D	582	2,4	34,70,70	1.56	4 (11%)	38,112,112	2.49	10 (26%)
6	SF4	E	806	2	0,12,12	0.00	-	-	-	-
6	SF4	B	802	2	0,12,12	0.00	-	-	-	-
5	SRM	B	581	-	30,66,70	1.33	3 (10%)	37,98,112	6.22	11 (29%)
6	SF4	B	801	2	0,12,12	0.00	-	-	-	-
6	SF4	D	807	1	0,12,12	0.00	-	-	-	-
6	SF4	A	804	1	0,12,12	0.00	-	-	-	-
6	SF4	A	803	1	0,12,12	0.00	-	-	-	-
6	SF4	D	808	1	0,12,12	0.00	-	-	-	-
5	SRM	A	580	2,4	34,70,70	1.66	5 (14%)	38,112,112	3.23	9 (23%)
5	SRM	E	583	-	30,66,70	1.50	5 (16%)	37,98,112	5.30	11 (29%)
4	SO3	D	6576	5	1,3,3	1.28	0	0,3,3	0.00	-
6	SF4	E	805	2	0,12,12	0.00	-	-	-	-

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	SF4	D	808	1	-	-	0/6/5/5
6	SF4	E	806	2	-	-	0/6/5/5
6	SF4	B	802	2	-	-	0/6/5/5
5	SRM	B	581	-	-	3/34/94/126	0/4/4/8
6	SF4	B	801	2	-	-	0/6/5/5
6	SF4	A	804	1	-	-	0/6/5/5
6	SF4	A	803	1	-	-	0/6/5/5
5	SRM	D	582	2,4	-	6/22/126/126	-
6	SF4	D	807	1	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	SRM	E	583	-	-	5/34/94/126	0/4/4/8
5	SRM	A	580	2,4	-	6/22/126/126	-
6	SF4	E	805	2	-	-	0/6/5/5

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	581	SRM	CMB-C2B	4.05	1.61	1.54
5	A	580	SRM	CMB-C2B	4.00	1.61	1.54
5	A	580	SRM	FE-NA	3.82	2.10	1.95
5	D	582	SRM	CMB-C2B	3.77	1.61	1.54
5	E	583	SRM	CMB-C2B	3.72	1.61	1.54

The worst 5 of 41 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	581	SRM	CAA-C3A-C2A	-31.13	88.33	123.52
5	E	583	SRM	CAA-C3A-C2A	-20.37	100.49	123.52
5	E	583	SRM	CAB-C3B-C2B	-15.22	106.31	123.52
5	E	583	SRM	CAA-CBA-CCA	14.90	137.67	112.67
5	B	581	SRM	CAB-C3B-C2B	-14.50	107.13	123.52

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

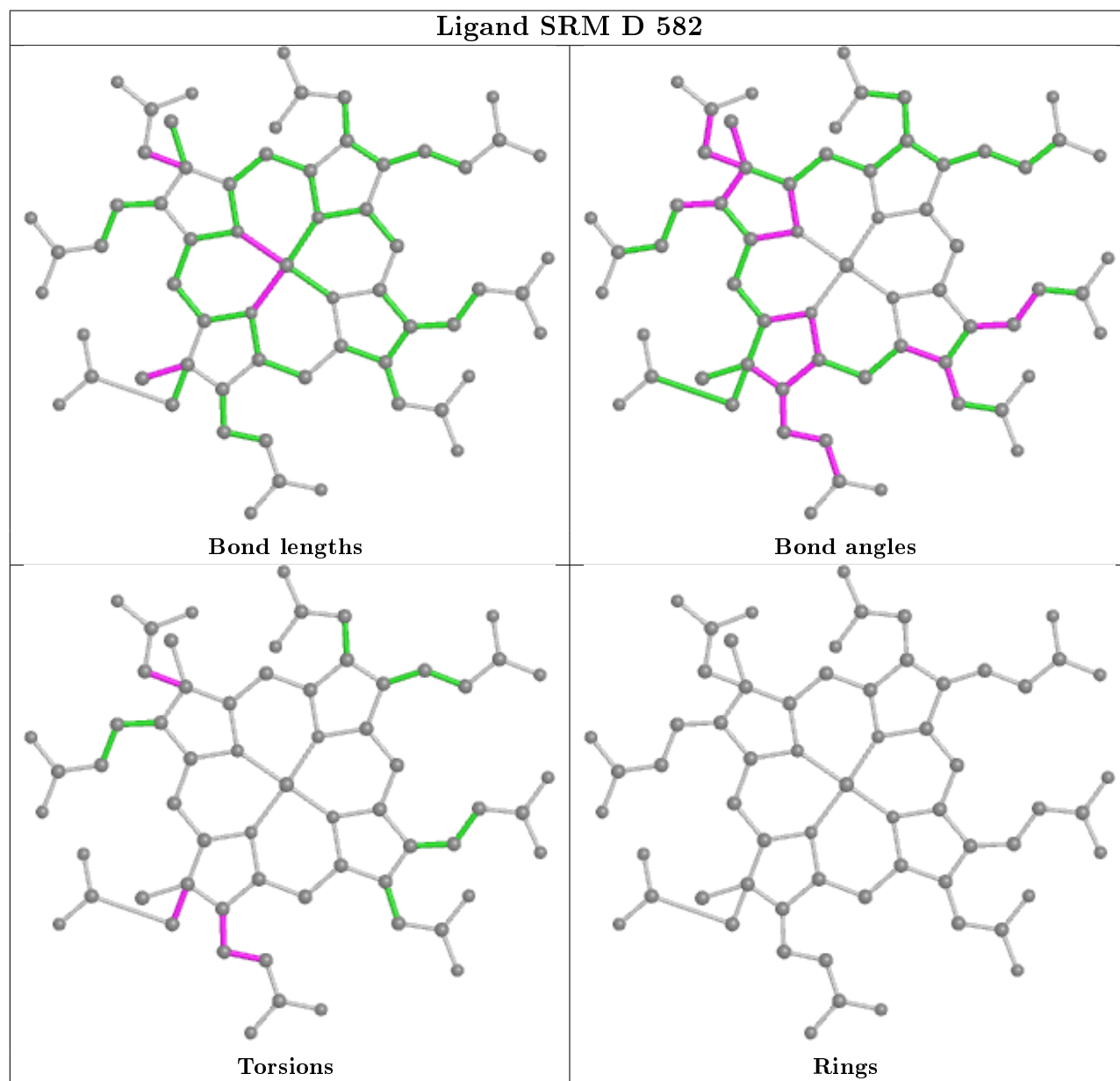
Mol	Chain	Res	Type	Atoms
5	D	582	SRM	C1A-C2A-CDA-CEA
5	D	582	SRM	CMA-C2A-CDA-CEA
5	D	582	SRM	C3A-C2A-CDA-CEA
5	B	581	SRM	C4A-C3A-CAA-CBA
5	E	583	SRM	C2A-C3A-CAA-CBA

There are no ring outliers.

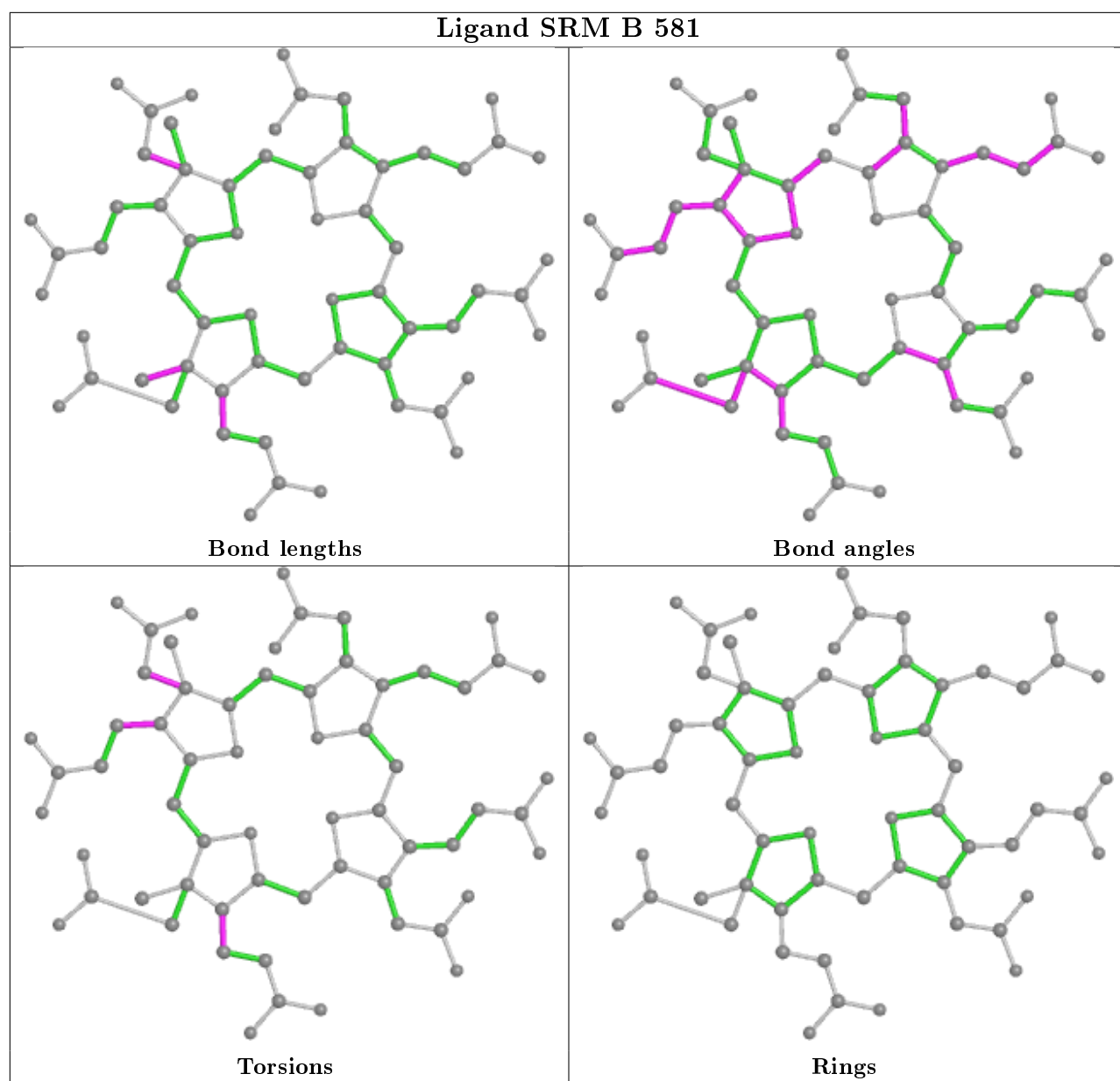
4 monomers are involved in 35 short contacts:

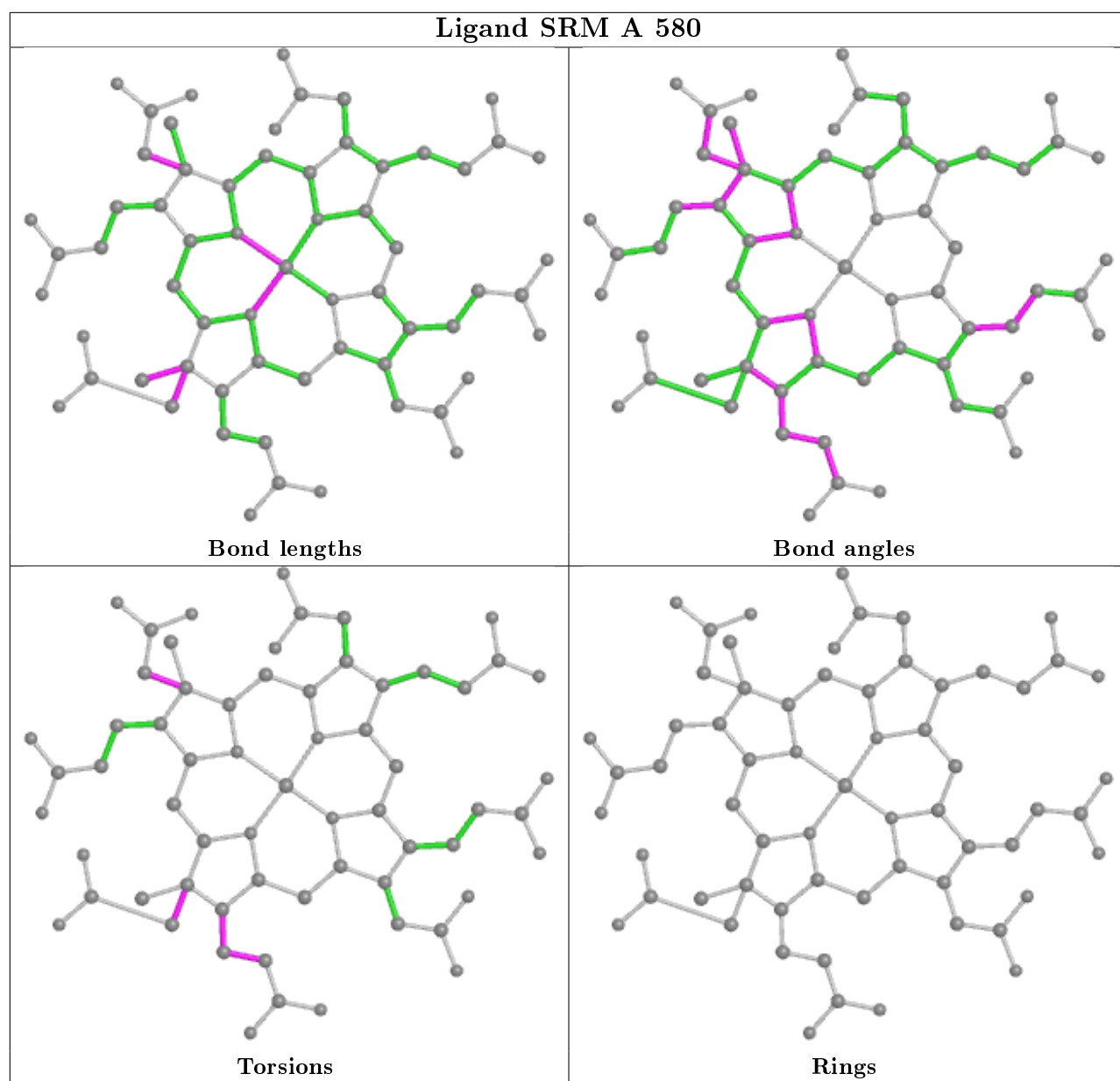
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	582	SRM	10	0
5	B	581	SRM	8	0
5	A	580	SRM	12	0
5	E	583	SRM	5	0

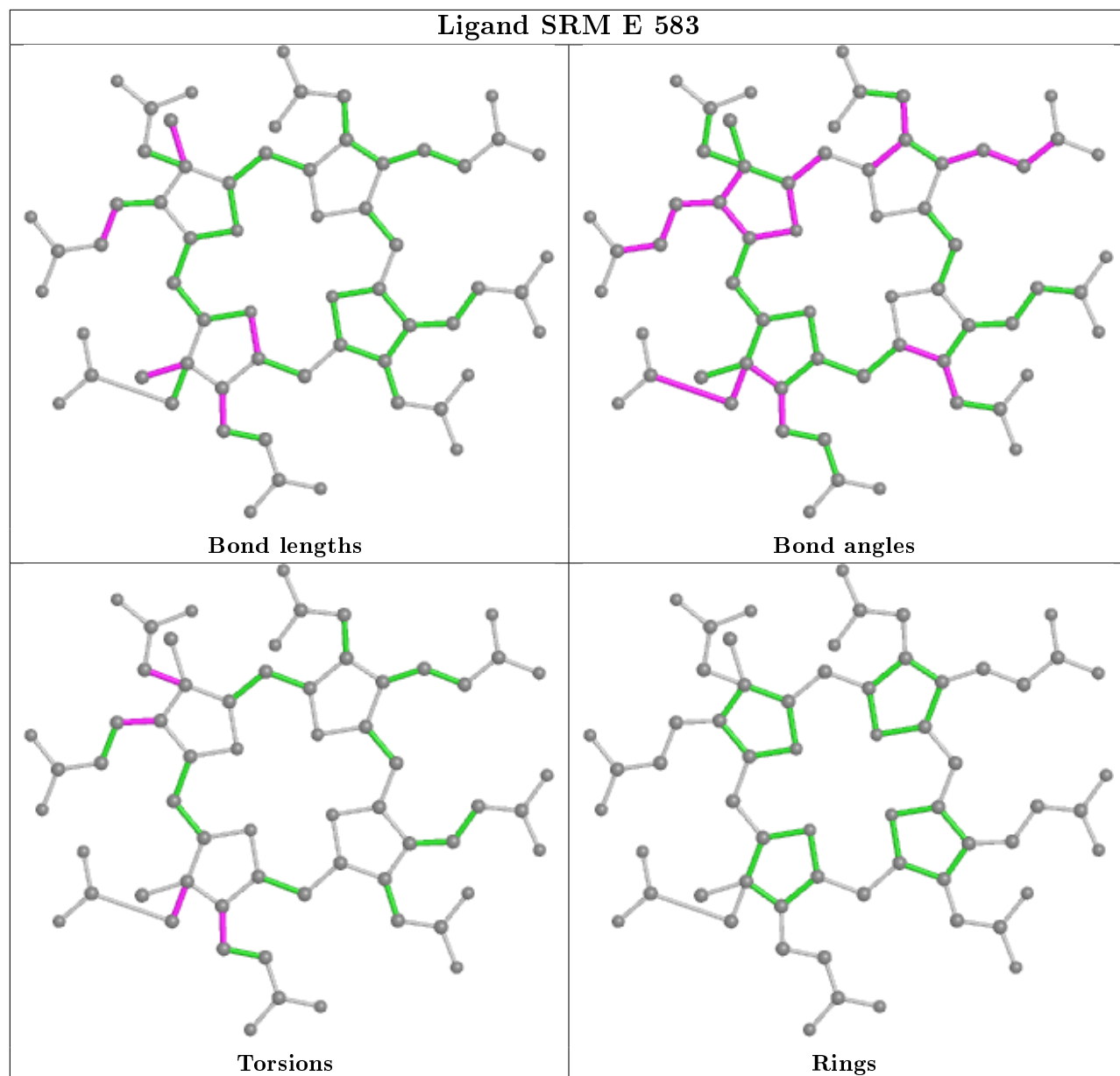
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.











## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	435/437 (99%)	0.19	8 (1%) 68 76	14, 23, 44, 74	0
1	D	435/437 (99%)	-0.10	6 (1%) 75 82	13, 19, 38, 73	0
2	B	385/386 (99%)	0.13	6 (1%) 72 79	15, 23, 40, 59	0
2	E	385/386 (99%)	0.02	7 (1%) 68 76	13, 21, 36, 78	0
3	C	104/105 (99%)	0.90	13 (12%) 3 5	22, 33, 64, 79	0
3	F	104/105 (99%)	0.33	5 (4%) 30 36	22, 29, 47, 66	0
All	All	1848/1856 (99%)	0.12	45 (2%) 59 65	13, 23, 44, 79	0

The worst 5 of 45 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	105	VAL	7.4
3	C	3	VAL	6.2
3	F	105	VAL	5.5
1	D	3	LYS	5.3
2	E	386	SER	5.1

### 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 carbohydrates in this entry.

## 6.4 Ligands

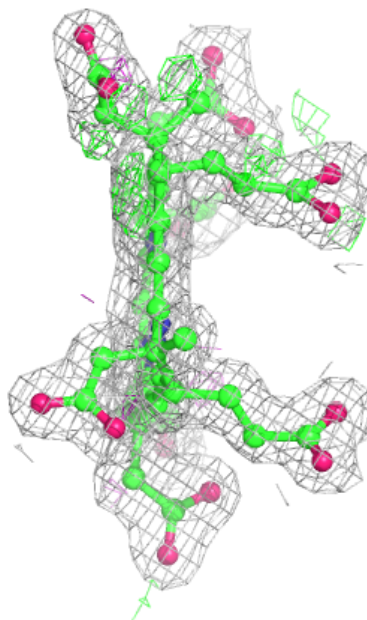
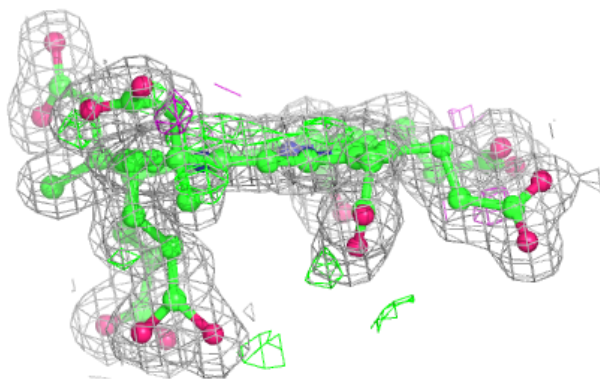
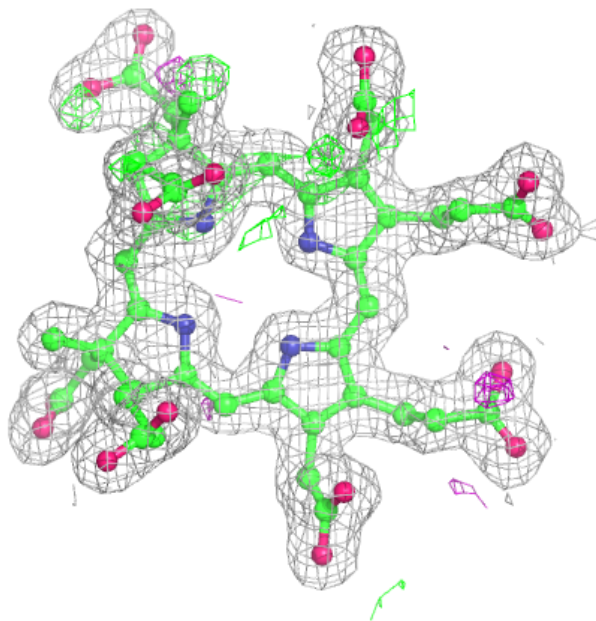
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	SRM	B	581	62/63	0.90	0.17	19,21,24,25	0
5	SRM	E	583	62/63	0.92	0.15	14,17,20,23	0
4	SO3	A	6575	4/4	0.93	0.14	39,40,40,40	0
5	SRM	D	582	63/63	0.93	0.13	18,21,28,31	0
5	SRM	A	580	63/63	0.93	0.13	20,24,31,33	0
6	SF4	A	803	8/8	0.95	0.08	22,23,25,26	0
6	SF4	E	806	8/8	0.95	0.08	22,23,25,25	0
6	SF4	B	801	8/8	0.95	0.08	17,23,24,24	0
6	SF4	E	805	8/8	0.95	0.08	18,21,24,24	0
6	SF4	A	804	8/8	0.96	0.11	17,20,23,23	0
6	SF4	D	808	8/8	0.96	0.10	13,18,21,22	0
4	SO3	D	6576	4/4	0.96	0.11	34,35,36,36	0
6	SF4	D	807	8/8	0.97	0.07	17,21,22,22	0
6	SF4	B	802	8/8	0.98	0.07	21,22,22,24	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.

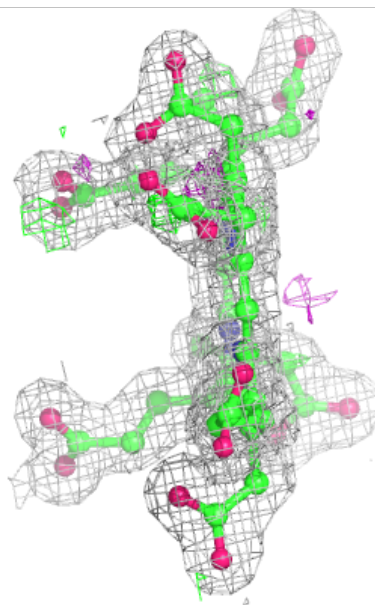
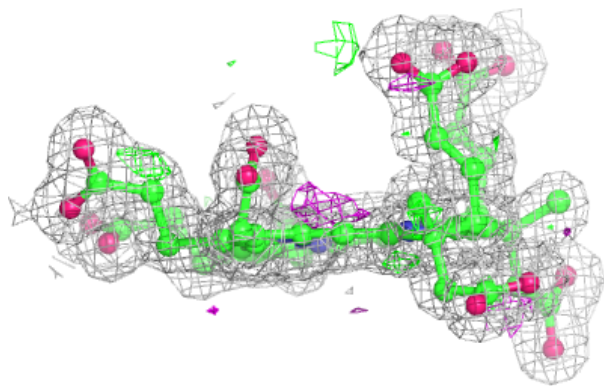
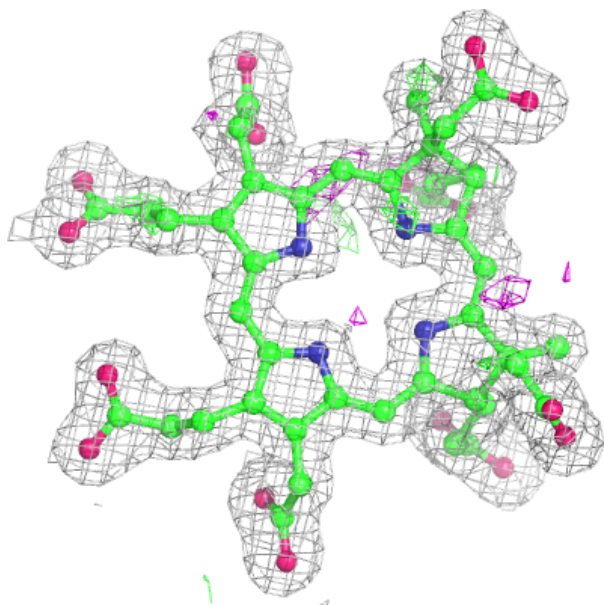
**Electron density around SRM B 581:**

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



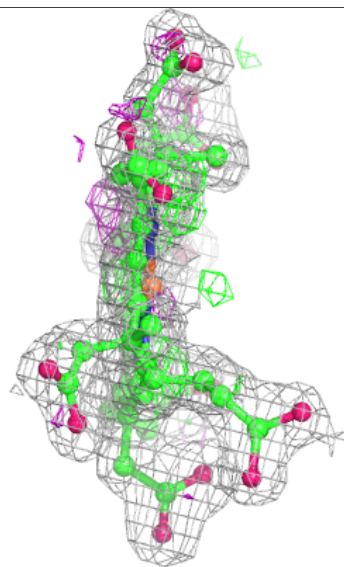
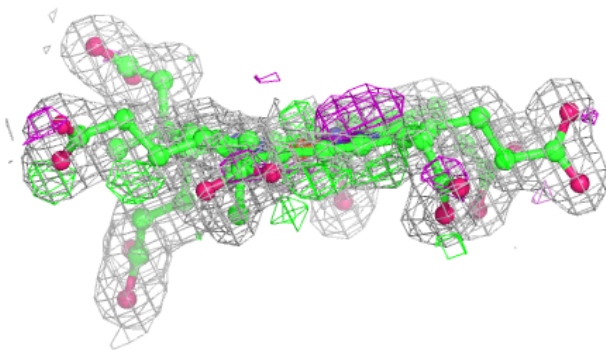
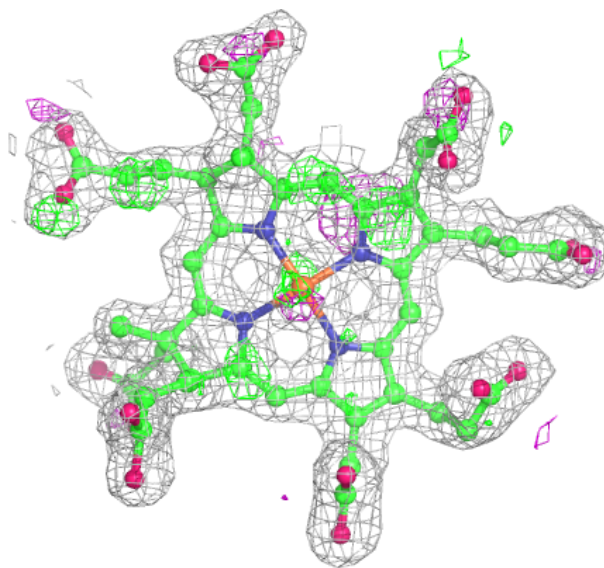
**Electron density around SRM E 583:**

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



**Electron density around SRM D 582:**

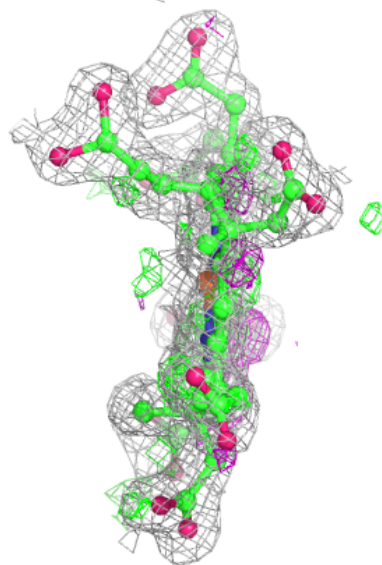
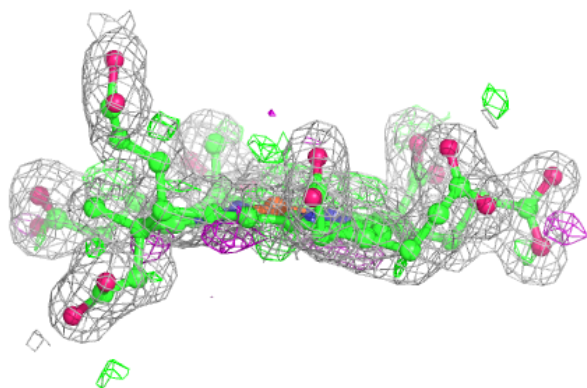
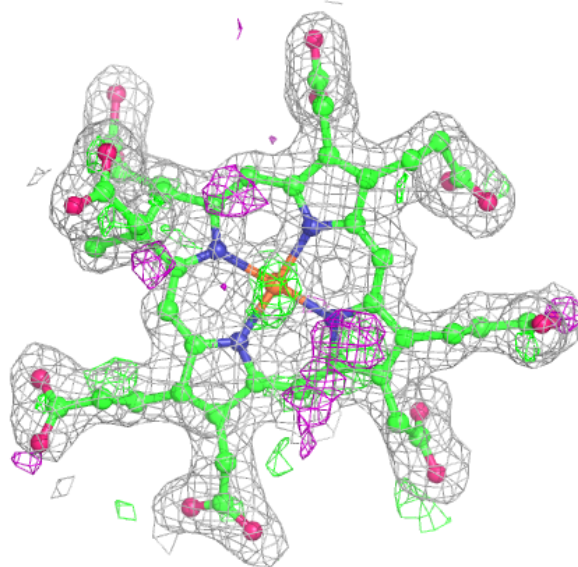
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around SRM A 580:**

$2mF_o-DF_c$  (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.