



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 14, 2020 – 08:25 am BST

PDB ID : 1XDY  
Title : Structural and Biochemical Identification of a Novel Bacterial Oxidoreductase, W-containing cofactor  
Authors : Loschi, L.; Brokx, S.J.; Hills, T.L.; Zhang, G.; Bertero, M.G.; Lovering, A.L.; Weiner, J.H.; Strynadka, N.C.  
Deposited on : 2004-09-08  
Resolution : 2.20 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:

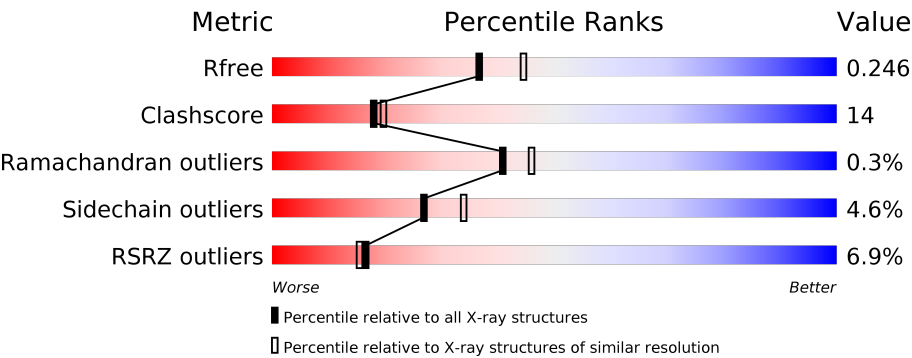
MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins)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

# 1 Overall quality at a glance i

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

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	298	<div><div>3%</div><div><div></div><div>67%</div><div>18%</div><div>•</div><div>12%</div></div></div>
1	B	298	<div><div>9%</div><div><div></div><div>60%</div><div>27%</div><div>•</div><div>12%</div></div></div>
1	C	298	<div><div>3%</div><div><div></div><div>67%</div><div>19%</div><div>•</div><div>11%</div></div></div>
1	D	298	<div><div>%</div><div><div></div><div>65%</div><div>22%</div><div>•</div><div>12%</div></div></div>
1	E	298	<div><div>3%</div><div><div></div><div>69%</div><div>18%</div><div>•</div><div>12%</div></div></div>
1	F	298	<div><div>16%</div><div><div></div><div>60%</div><div>26%</div><div>•</div><div>12%</div></div></div>

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Mol	Chain	Length	Quality of chain
1	G	298	
1	H	298	
1	I	298	
1	J	298	

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
3	MTE	A	301	X	-	-	-
3	MTE	B	302	X	-	-	-
3	MTE	C	303	X	-	-	-
3	MTE	D	304	X	-	-	-
3	MTE	E	305	X	-	-	-
3	MTE	F	306	X	-	-	-
3	MTE	G	307	X	-	-	-
3	MTE	H	308	X	-	-	-
3	MTE	I	309	X	-	-	-
3	MTE	J	310	X	-	-	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21369 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 Bacterial Sulfite Oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2093	1358	354	375	6			
1	B	262	Total	C	N	O	S	0	0	0
			2088	1354	353	375	6			
1	C	264	Total	C	N	O	S	0	0	0
			2102	1363	356	377	6			
1	D	263	Total	C	N	O	S	0	0	0
			2094	1357	355	376	6			
1	E	262	Total	C	N	O	S	0	0	0
			2085	1351	353	375	6			
1	F	262	Total	C	N	O	S	1	0	0
			2084	1352	352	374	6			
1	G	261	Total	C	N	O	S	0	0	0
			2079	1349	351	373	6			
1	H	264	Total	C	N	O	S	11	0	0
			2102	1363	356	377	6			
1	I	262	Total	C	N	O	S	0	0	0
			2085	1351	353	375	6			
1	J	262	Total	C	N	O	S	0	0	0
			2085	1351	353	375	6			

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	275	ASP	ALA	SEE REMARK 999	UNP P76342
A	291	LEU	-	EXPRESSION TAG	UNP P76342
A	292	GLU	-	EXPRESSION TAG	UNP P76342
A	293	HIS	-	EXPRESSION TAG	UNP P76342
A	294	HIS	-	EXPRESSION TAG	UNP P76342
A	295	HIS	-	EXPRESSION TAG	UNP P76342
A	296	HIS	-	EXPRESSION TAG	UNP P76342
A	297	HIS	-	EXPRESSION TAG	UNP P76342
A	298	HIS	-	EXPRESSION TAG	UNP P76342

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Chain	Residue	Modelled	Actual	Comment	Reference
B	275	ASP	ALA	SEE REMARK 999	UNP P76342
B	291	LEU	-	EXPRESSION TAG	UNP P76342
B	292	GLU	-	EXPRESSION TAG	UNP P76342
B	293	HIS	-	EXPRESSION TAG	UNP P76342
B	294	HIS	-	EXPRESSION TAG	UNP P76342
B	295	HIS	-	EXPRESSION TAG	UNP P76342
B	296	HIS	-	EXPRESSION TAG	UNP P76342
B	297	HIS	-	EXPRESSION TAG	UNP P76342
B	298	HIS	-	EXPRESSION TAG	UNP P76342
C	275	ASP	ALA	SEE REMARK 999	UNP P76342
C	291	LEU	-	EXPRESSION TAG	UNP P76342
C	292	GLU	-	EXPRESSION TAG	UNP P76342
C	293	HIS	-	EXPRESSION TAG	UNP P76342
C	294	HIS	-	EXPRESSION TAG	UNP P76342
C	295	HIS	-	EXPRESSION TAG	UNP P76342
C	296	HIS	-	EXPRESSION TAG	UNP P76342
C	297	HIS	-	EXPRESSION TAG	UNP P76342
C	298	HIS	-	EXPRESSION TAG	UNP P76342
D	275	ASP	ALA	SEE REMARK 999	UNP P76342
D	291	LEU	-	EXPRESSION TAG	UNP P76342
D	292	GLU	-	EXPRESSION TAG	UNP P76342
D	293	HIS	-	EXPRESSION TAG	UNP P76342
D	294	HIS	-	EXPRESSION TAG	UNP P76342
D	295	HIS	-	EXPRESSION TAG	UNP P76342
D	296	HIS	-	EXPRESSION TAG	UNP P76342
D	297	HIS	-	EXPRESSION TAG	UNP P76342
D	298	HIS	-	EXPRESSION TAG	UNP P76342
E	275	ASP	ALA	SEE REMARK 999	UNP P76342
E	291	LEU	-	EXPRESSION TAG	UNP P76342
E	292	GLU	-	EXPRESSION TAG	UNP P76342
E	293	HIS	-	EXPRESSION TAG	UNP P76342
E	294	HIS	-	EXPRESSION TAG	UNP P76342
E	295	HIS	-	EXPRESSION TAG	UNP P76342
E	296	HIS	-	EXPRESSION TAG	UNP P76342
E	297	HIS	-	EXPRESSION TAG	UNP P76342
E	298	HIS	-	EXPRESSION TAG	UNP P76342
F	275	ASP	ALA	SEE REMARK 999	UNP P76342
F	291	LEU	-	EXPRESSION TAG	UNP P76342
F	292	GLU	-	EXPRESSION TAG	UNP P76342
F	293	HIS	-	EXPRESSION TAG	UNP P76342
F	294	HIS	-	EXPRESSION TAG	UNP P76342
F	295	HIS	-	EXPRESSION TAG	UNP P76342

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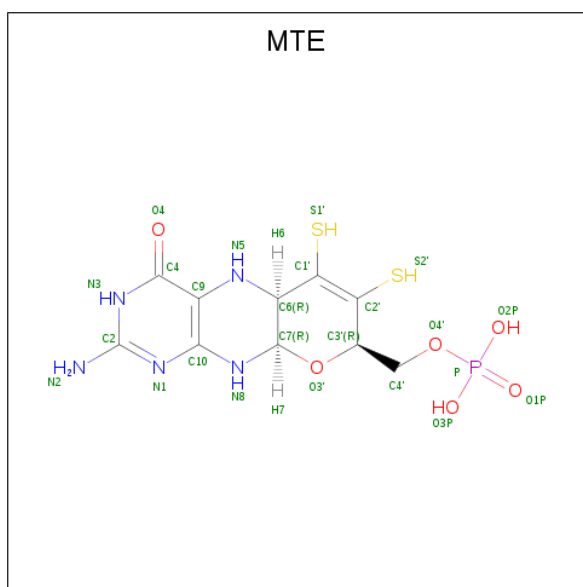
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Chain	Residue	Modelled	Actual	Comment	Reference
F	296	HIS	-	EXPRESSION TAG	UNP P76342
F	297	HIS	-	EXPRESSION TAG	UNP P76342
F	298	HIS	-	EXPRESSION TAG	UNP P76342
G	275	ASP	ALA	SEE REMARK 999	UNP P76342
G	291	LEU	-	EXPRESSION TAG	UNP P76342
G	292	GLU	-	EXPRESSION TAG	UNP P76342
G	293	HIS	-	EXPRESSION TAG	UNP P76342
G	294	HIS	-	EXPRESSION TAG	UNP P76342
G	295	HIS	-	EXPRESSION TAG	UNP P76342
G	296	HIS	-	EXPRESSION TAG	UNP P76342
G	297	HIS	-	EXPRESSION TAG	UNP P76342
G	298	HIS	-	EXPRESSION TAG	UNP P76342
H	275	ASP	ALA	SEE REMARK 999	UNP P76342
H	291	LEU	-	EXPRESSION TAG	UNP P76342
H	292	GLU	-	EXPRESSION TAG	UNP P76342
H	293	HIS	-	EXPRESSION TAG	UNP P76342
H	294	HIS	-	EXPRESSION TAG	UNP P76342
H	295	HIS	-	EXPRESSION TAG	UNP P76342
H	296	HIS	-	EXPRESSION TAG	UNP P76342
H	297	HIS	-	EXPRESSION TAG	UNP P76342
H	298	HIS	-	EXPRESSION TAG	UNP P76342
I	275	ASP	ALA	SEE REMARK 999	UNP P76342
I	291	LEU	-	EXPRESSION TAG	UNP P76342
I	292	GLU	-	EXPRESSION TAG	UNP P76342
I	293	HIS	-	EXPRESSION TAG	UNP P76342
I	294	HIS	-	EXPRESSION TAG	UNP P76342
I	295	HIS	-	EXPRESSION TAG	UNP P76342
I	296	HIS	-	EXPRESSION TAG	UNP P76342
I	297	HIS	-	EXPRESSION TAG	UNP P76342
I	298	HIS	-	EXPRESSION TAG	UNP P76342
J	275	ASP	ALA	SEE REMARK 999	UNP P76342
J	291	LEU	-	EXPRESSION TAG	UNP P76342
J	292	GLU	-	EXPRESSION TAG	UNP P76342
J	293	HIS	-	EXPRESSION TAG	UNP P76342
J	294	HIS	-	EXPRESSION TAG	UNP P76342
J	295	HIS	-	EXPRESSION TAG	UNP P76342
J	296	HIS	-	EXPRESSION TAG	UNP P76342
J	297	HIS	-	EXPRESSION TAG	UNP P76342
J	298	HIS	-	EXPRESSION TAG	UNP P76342

- Molecule 2 is TUNGSTEN ION (three-letter code: W) (formula: W).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	1	Total W 1 1	0	0
2	J	1	Total W 1 1	0	0
2	D	1	Total W 1 1	0	0
2	E	1	Total W 1 1	0	0
2	H	1	Total W 1 1	0	0
2	B	1	Total W 1 1	0	0
2	I	1	Total W 1 1	0	0
2	C	1	Total W 1 1	0	0
2	A	1	Total W 1 1	0	0
2	F	1	Total W 1 1	0	0

- Molecule 3 is PHOSPHONIC ACIDMONO-(2-AMINO-5,6-DIMERCAPTO-4-OXO-3,7,8A, 9,10,10A-HEXAHYDRO-4H-8-OXA-1,3,9,10-TETRAAZA-ANTHRACEN-7-YLMETHYL) ESTER (three-letter code: MTE) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>5</sub>O<sub>6</sub>P<sub>2</sub>).



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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	C	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	D	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	E	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	F	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	G	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	H	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	I	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	J	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		

- Molecule 4 is water.

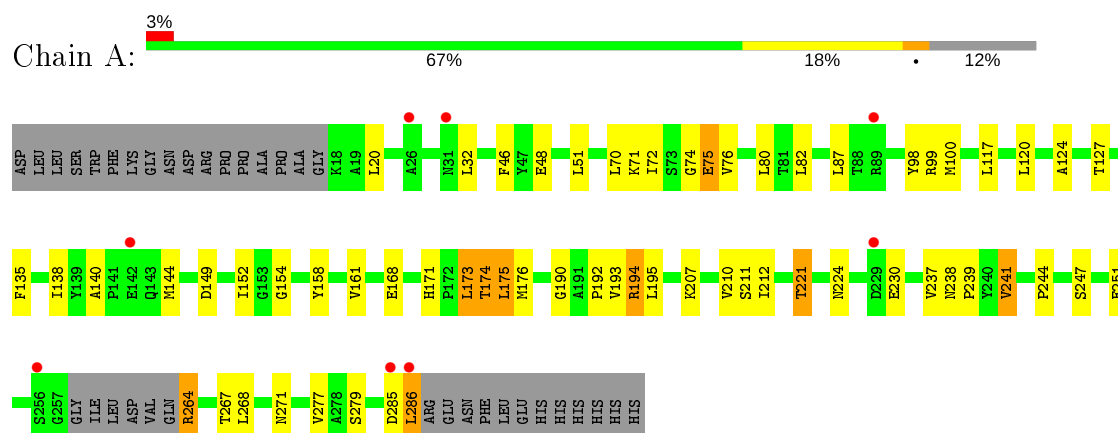
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	35	Total	O	0	0
			35	35		
4	B	13	Total	O	0	0
			13	13		
4	C	32	Total	O	0	0
			32	32		
4	D	30	Total	O	0	0
			30	30		
4	E	29	Total	O	0	0
			29	29		
4	F	10	Total	O	0	0
			10	10		
4	G	22	Total	O	0	0
			22	22		
4	H	4	Total	O	0	0
			4	4		
4	I	12	Total	O	0	0
			12	12		
4	J	35	Total	O	0	0
			35	35		



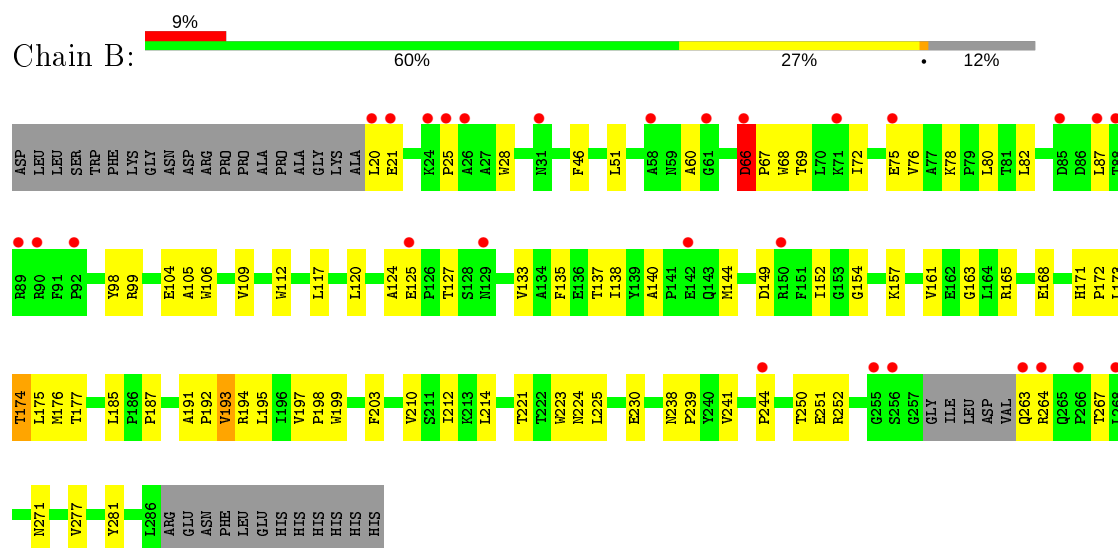
### 3 Residue-property plots [i](#)

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.

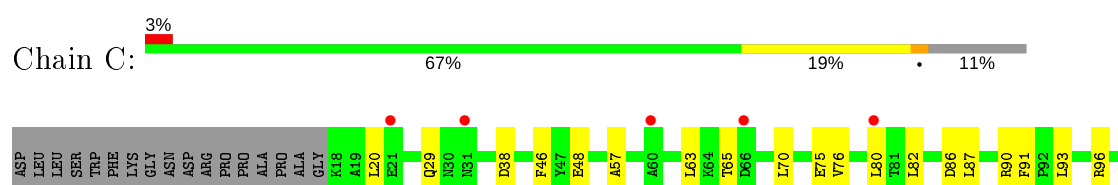
#### • Molecule 1: Bacterial Sulfite Oxidase

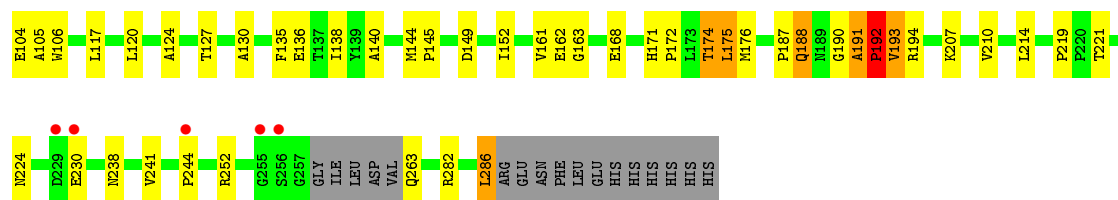


#### • Molecule 1: Bacterial Sulfite Oxidase

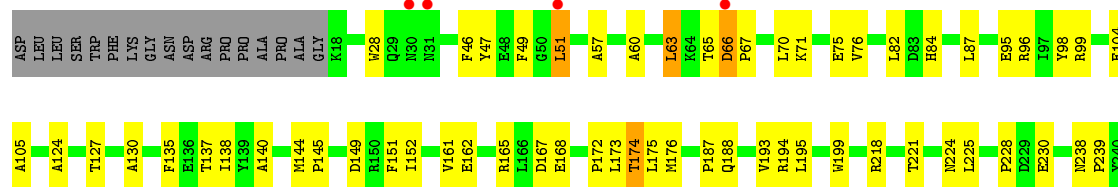


#### • Molecule 1: Bacterial Sulfite Oxidase

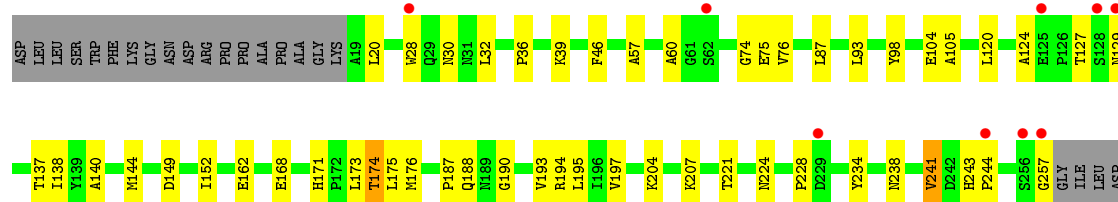




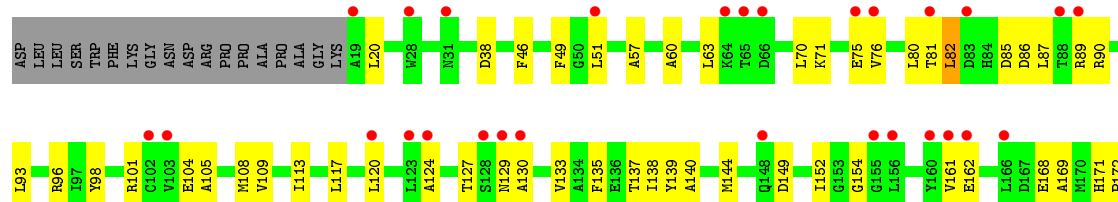
• Molecule 1: Bacterial Sulfite Oxidase

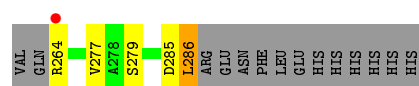


• Molecule 1: Bacterial Sulfite Oxidase

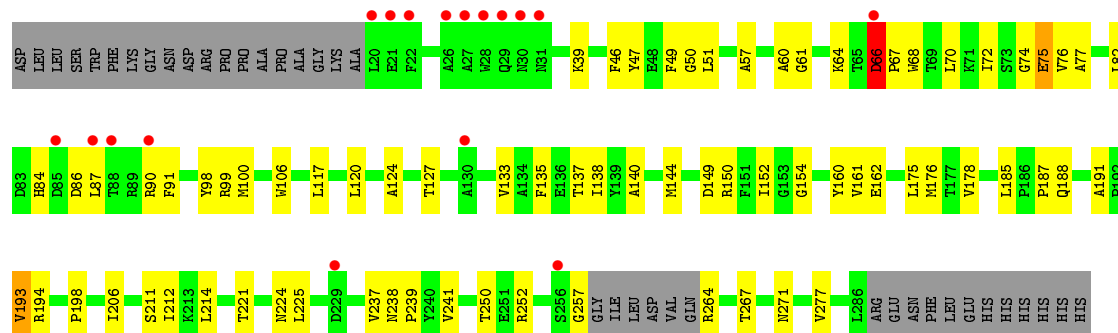


• Molecule 1: Bacterial Sulfite Oxidase

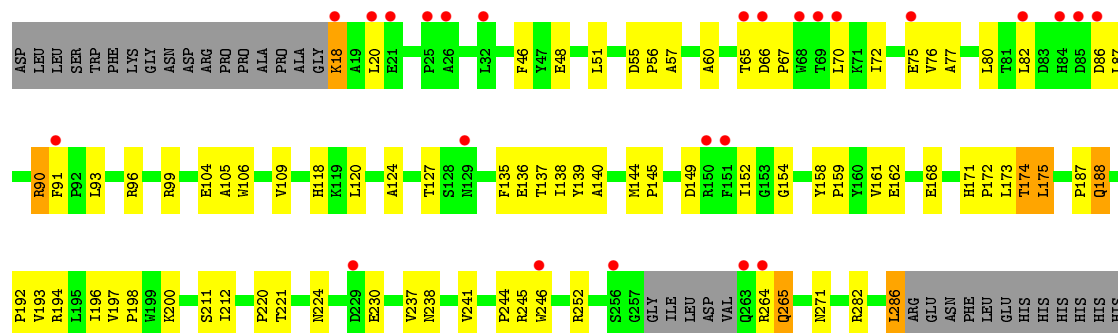




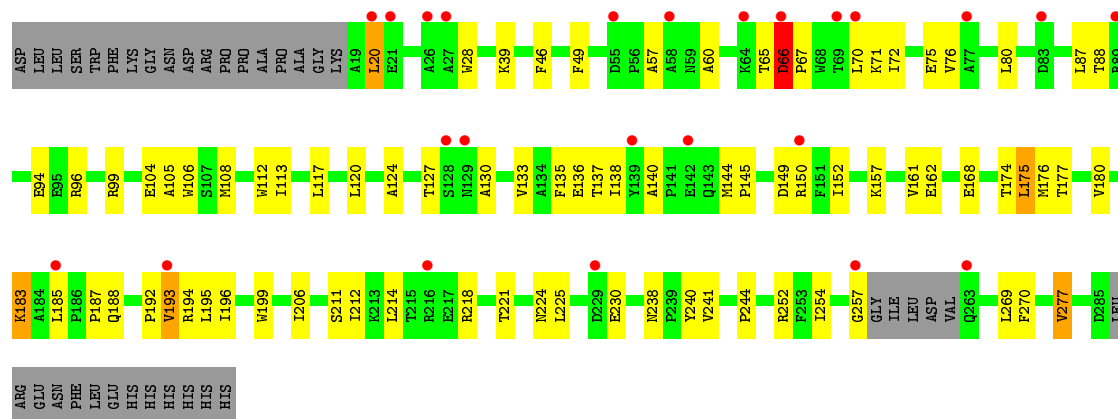
• Molecule 1: Bacterial Sulfite Oxidase



• Molecule 1: Bacterial Sulfite Oxidase

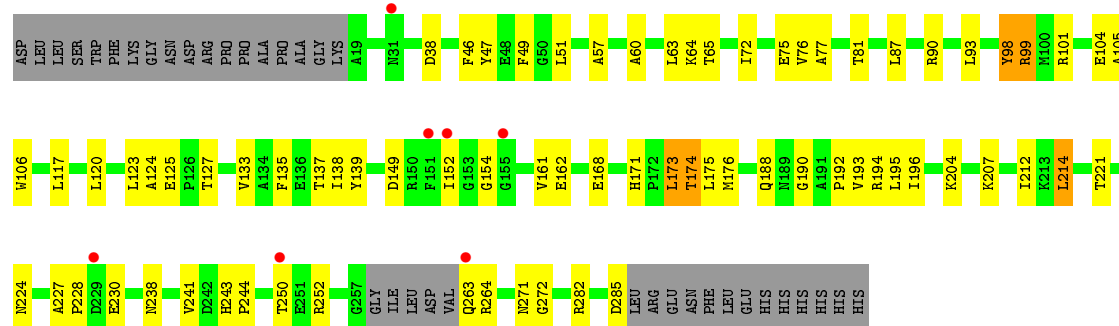


• Molecule 1: Bacterial Sulfite Oxidase



• Molecule 1: Bacterial Sulfite Oxidase

Chain J: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.84Å 177.51Å 104.21Å 90.00° 109.73° 90.00°	Depositor
Resolution (Å)	38.96 – 2.20 38.96 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.5 (38.96-2.20) 98.6 (38.96-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.69 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.226 , 0.258 0.213 , 0.246	Depositor DCC
$R_{free}$ test set	4389 reflections (3.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.8	Xtriage
Anisotropy	0.141	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 47.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	21369	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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: W, MTE

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	0.57	1/2158 (0.0%)	0.76	2/2942 (0.1%)
1	B	0.43	0/2153	0.69	2/2936 (0.1%)
1	C	0.51	0/2167	0.85	3/2954 (0.1%)
1	D	0.53	0/2159	0.77	1/2943 (0.0%)
1	E	0.55	0/2150	0.75	1/2932 (0.0%)
1	F	0.42	0/2149	0.69	0/2931
1	G	0.53	0/2144	0.75	1/2924 (0.0%)
1	H	0.40	0/2167	0.67	1/2954 (0.0%)
1	I	0.45	0/2150	0.70	1/2932 (0.0%)
1	J	0.58	0/2150	0.77	0/2932
All	All	0.50	1/21547 (0.0%)	0.74	12/29380 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	E	0	1
1	J	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	194	ARG	CB-CG	-5.40	1.38	1.52

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	191	ALA	C-N-CD	-20.12	76.34	120.60
1	C	191	ALA	C-N-CA	13.52	178.78	122.00
1	A	175	LEU	CA-CB-CG	7.14	131.72	115.30
1	G	66	ASP	C-N-CD	7.09	143.29	128.40
1	B	66	ASP	C-N-CD	6.70	142.46	128.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	98	TYR	Sidechain
1	E	234	TYR	Sidechain
1	J	98	TYR	Sidechain

## 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2093	0	2067	49	0
1	B	2088	0	2057	62	0
1	C	2102	0	2075	53	0
1	D	2094	0	2064	63	0
1	E	2085	0	2051	46	0
1	F	2084	0	2054	63	0
1	G	2079	0	2049	54	0
1	H	2102	0	2075	58	0
1	I	2085	0	2051	67	0
1	J	2085	0	2051	61	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	24	0	9	0	0
3	B	24	0	9	0	0
3	C	24	0	9	0	0
3	D	24	0	9	0	0
3	E	24	0	9	0	0
3	F	24	0	9	0	0
3	G	24	0	9	0	0
3	H	24	0	9	0	0
3	I	24	0	9	0	0
3	J	24	0	9	0	0
4	A	35	0	0	1	0
4	B	13	0	0	1	0
4	C	32	0	0	2	0
4	D	30	0	0	0	0
4	E	29	0	0	1	0
4	F	10	0	0	0	0
4	G	22	0	0	1	0
4	H	4	0	0	0	0
4	I	12	0	0	0	0
4	J	35	0	0	2	0
All	All	21369	0	20684	565	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 565 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:250:THR:HG23	1:D:264:ARG:HH21	1.33	0.92
1:H:57:ALA:HA	1:H:188:GLN:HG2	1.51	0.92
1:B:20:LEU:HG	1:B:21:GLU:H	1.35	0.90
1:A:140:ALA:HB3	1:A:144:MET:HE3	1.52	0.88
1:E:140:ALA:HB3	1:E:144:MET:HE3	1.56	0.86

There are no symmetry-related clashes.



## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	259/298 (87%)	249 (96%)	9 (4%)	1 (0%)	34	37
1	B	258/298 (87%)	242 (94%)	15 (6%)	1 (0%)	34	37
1	C	260/298 (87%)	246 (95%)	13 (5%)	1 (0%)	34	37
1	D	259/298 (87%)	248 (96%)	11 (4%)	0	100	100
1	E	258/298 (87%)	241 (93%)	17 (7%)	0	100	100
1	F	258/298 (87%)	243 (94%)	15 (6%)	0	100	100
1	G	257/298 (86%)	242 (94%)	13 (5%)	2 (1%)	19	19
1	H	260/298 (87%)	236 (91%)	23 (9%)	1 (0%)	34	37
1	I	258/298 (87%)	239 (93%)	18 (7%)	1 (0%)	34	37
1	J	258/298 (87%)	244 (95%)	14 (5%)	0	100	100
All	All	2585/2980 (87%)	2430 (94%)	148 (6%)	7 (0%)	41	46

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	192	PRO
1	G	75	GLU
1	A	75	GLU
1	B	66	ASP
1	G	66	ASP

### 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	219/249 (88%)	204 (93%)	15 (7%)	16	17
1	B	219/249 (88%)	213 (97%)	6 (3%)	44	57
1	C	220/249 (88%)	207 (94%)	13 (6%)	19	23
1	D	219/249 (88%)	211 (96%)	8 (4%)	34	43
1	E	218/249 (88%)	211 (97%)	7 (3%)	39	50
1	F	218/249 (88%)	206 (94%)	12 (6%)	21	26
1	G	218/249 (88%)	210 (96%)	8 (4%)	34	43
1	H	220/249 (88%)	207 (94%)	13 (6%)	19	23
1	I	218/249 (88%)	209 (96%)	9 (4%)	30	39
1	J	218/249 (88%)	208 (95%)	10 (5%)	27	34
All	All	2187/2490 (88%)	2086 (95%)	101 (5%)	27	34

5 of 101 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	174	THR
1	F	174	THR
1	J	81	THR
1	E	241	VAL
1	F	63	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	30	ASN
1	J	263	GLN
1	G	59	ASN
1	D	263	GLN
1	J	31	ASN

### 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 ⓘ

Of 20 ligands modelled in this entry, 10 are monoatomic - leaving 10 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
3	MTE	C	303	2	21,26,26	3.16	12 (57%)	21,40,40	6.51	13 (61%)
3	MTE	H	308	2	21,26,26	3.53	11 (52%)	21,40,40	5.96	15 (71%)
3	MTE	B	302	2	21,26,26	3.22	12 (57%)	21,40,40	5.99	15 (71%)
3	MTE	G	307	2	21,26,26	3.23	11 (52%)	21,40,40	6.40	12 (57%)
3	MTE	I	309	2	21,26,26	3.62	11 (52%)	21,40,40	6.14	13 (61%)
3	MTE	A	301	2	21,26,26	3.43	12 (57%)	21,40,40	6.27	12 (57%)
3	MTE	J	310	2	21,26,26	3.50	10 (47%)	21,40,40	6.55	13 (61%)
3	MTE	D	304	2	21,26,26	3.17	11 (52%)	21,40,40	6.04	13 (61%)
3	MTE	E	305	2	21,26,26	2.89	9 (42%)	21,40,40	6.53	13 (61%)
3	MTE	F	306	2	21,26,26	3.44	10 (47%)	21,40,40	6.29	16 (76%)

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
3	MTE	C	303	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	H	308	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	B	302	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	G	307	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	I	309	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	A	301	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	J	310	2	1/1/6/8	0/6/34/34	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MTE	D	304	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	E	305	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	F	306	2	1/1/6/8	0/6/34/34	0/3/3/3

The worst 5 of 109 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	309	MTE	C4'-C3'	-10.16	1.38	1.52
3	H	308	MTE	C4'-C3'	-9.50	1.39	1.52
3	J	310	MTE	C4'-C3'	-9.27	1.39	1.52
3	F	306	MTE	C4'-C3'	-9.02	1.39	1.52
3	G	307	MTE	C4'-C3'	-8.93	1.39	1.52

The worst 5 of 135 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	305	MTE	O3'-C7-C6	-24.51	92.61	108.96
3	J	310	MTE	O3'-C7-C6	-24.48	92.63	108.96
3	C	303	MTE	O3'-C7-C6	-24.35	92.72	108.96
3	G	307	MTE	O3'-C7-C6	-23.05	93.59	108.96
3	A	301	MTE	O3'-C7-C6	-22.85	93.73	108.96

5 of 10 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	I	309	MTE	C6
3	C	303	MTE	C6
3	G	307	MTE	C6
3	B	302	MTE	C6
3	E	305	MTE	C6

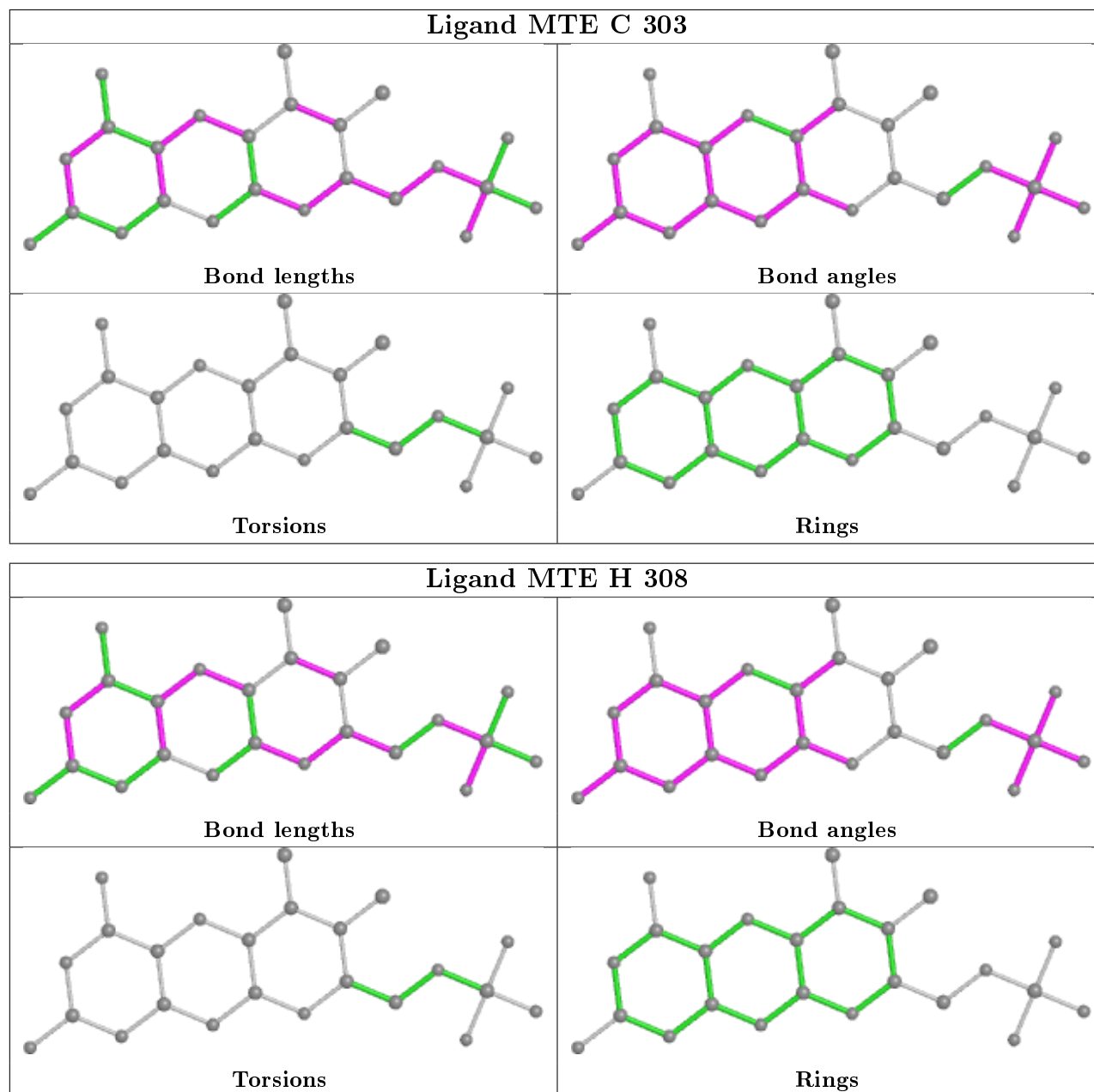
There are no torsion outliers.

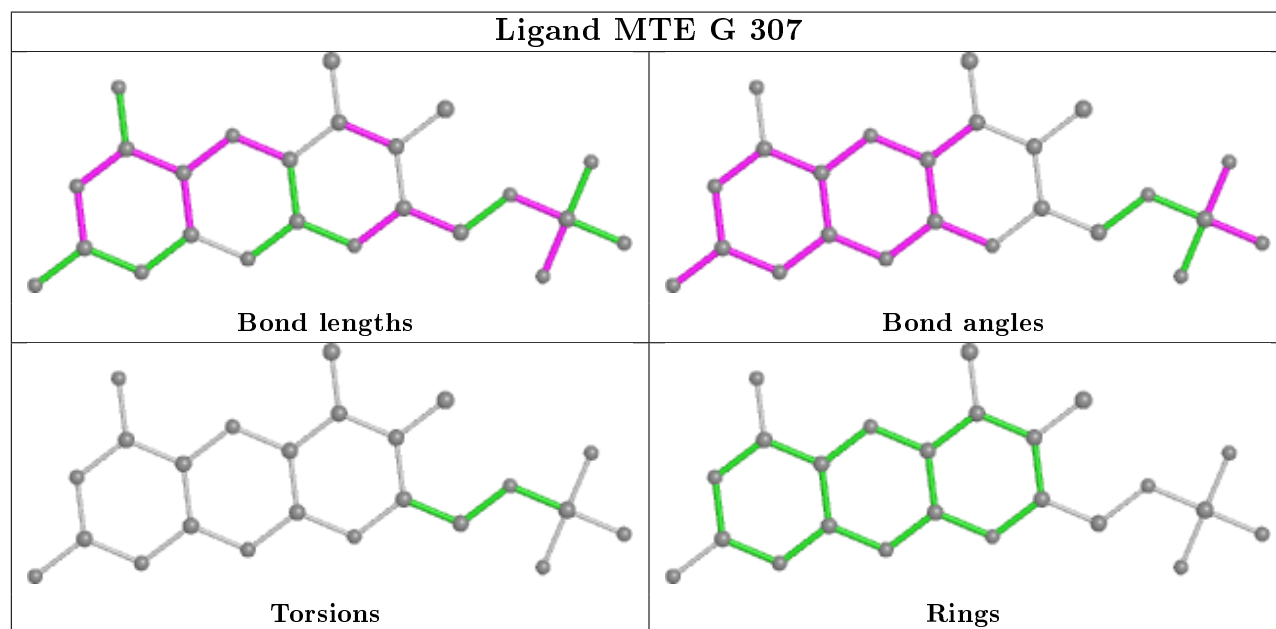
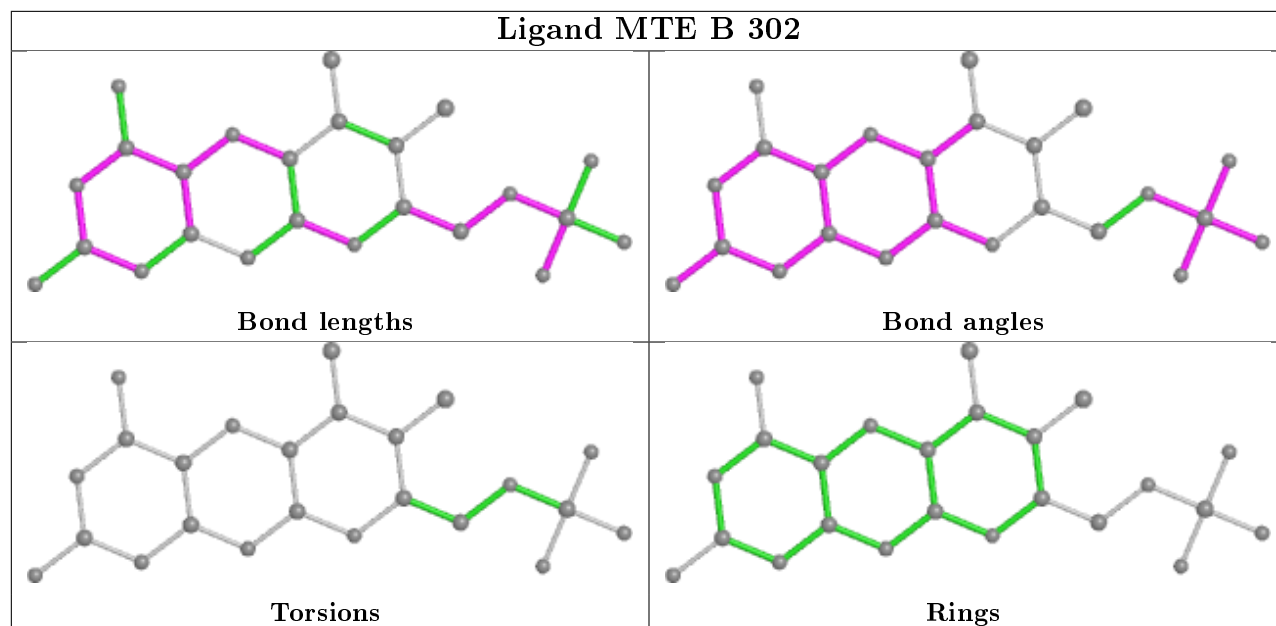
There are no ring outliers.

No monomer is involved in short contacts.

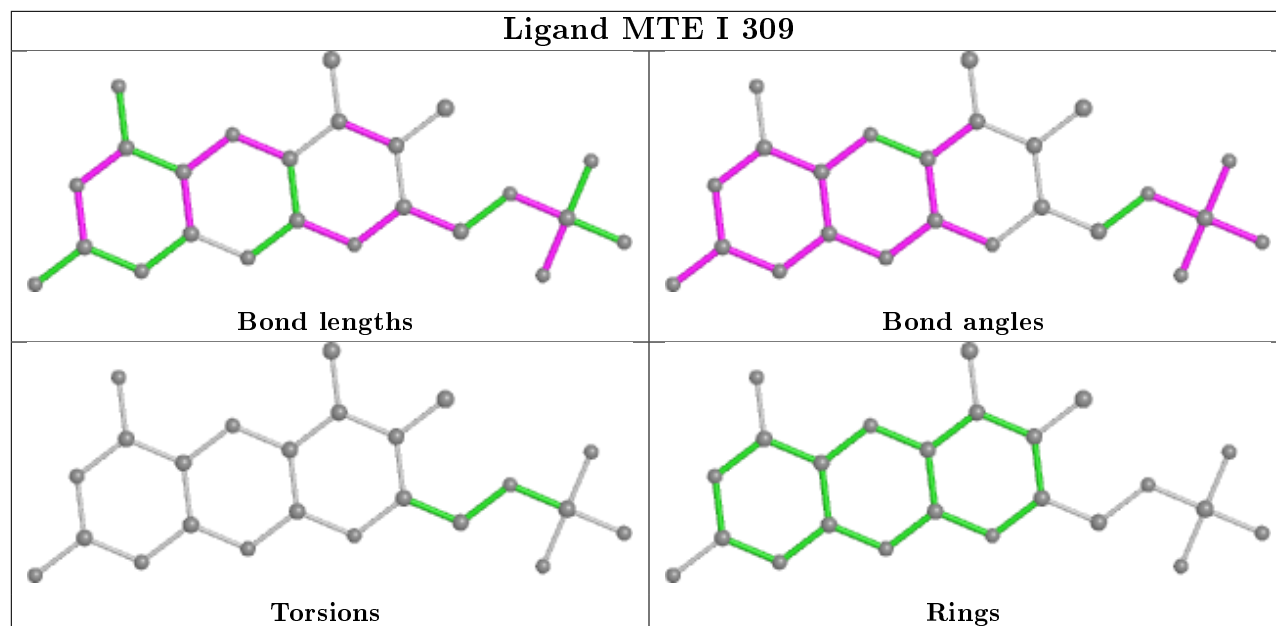
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

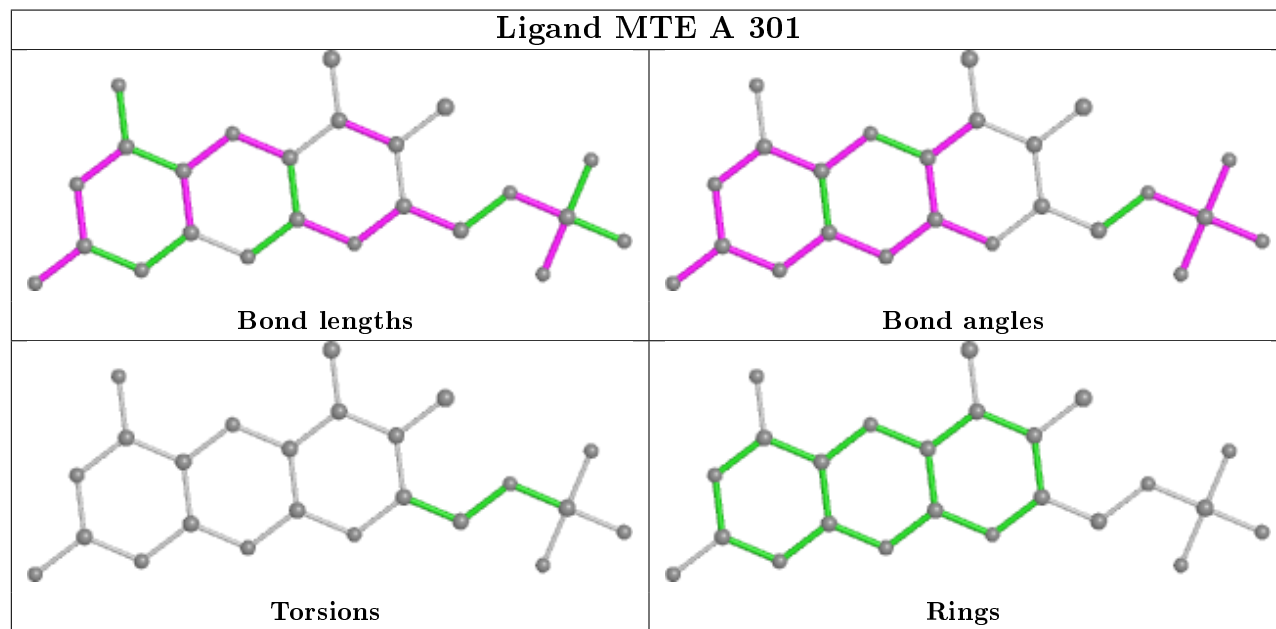




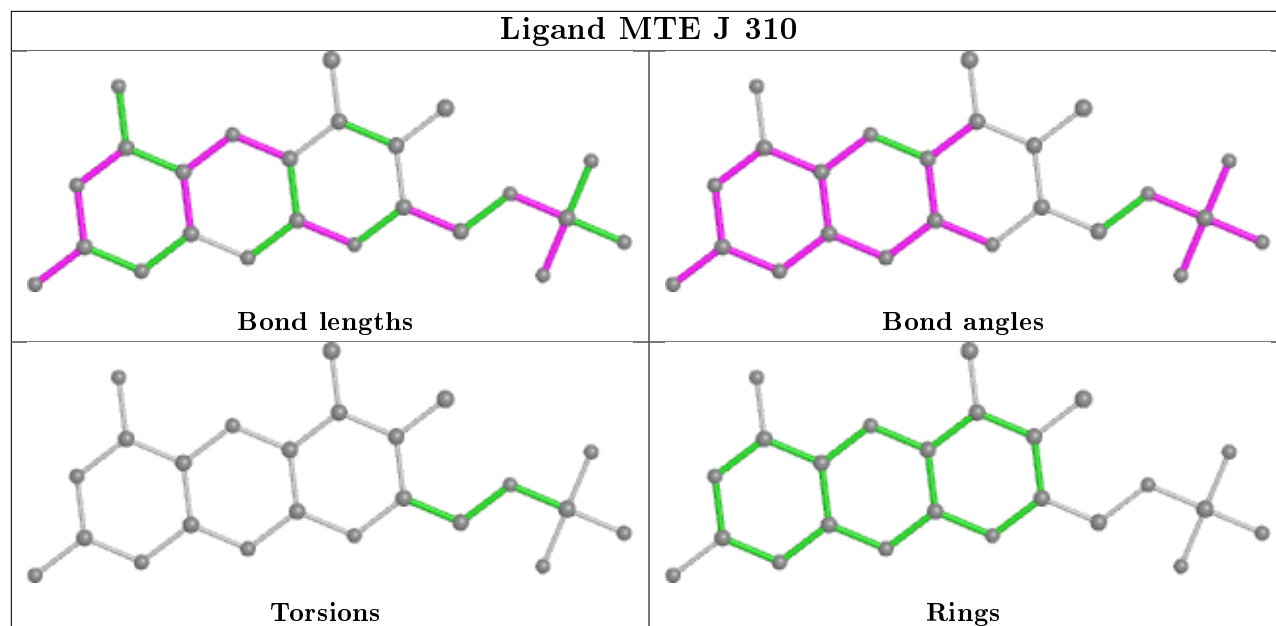
## Ligand MTE I 309



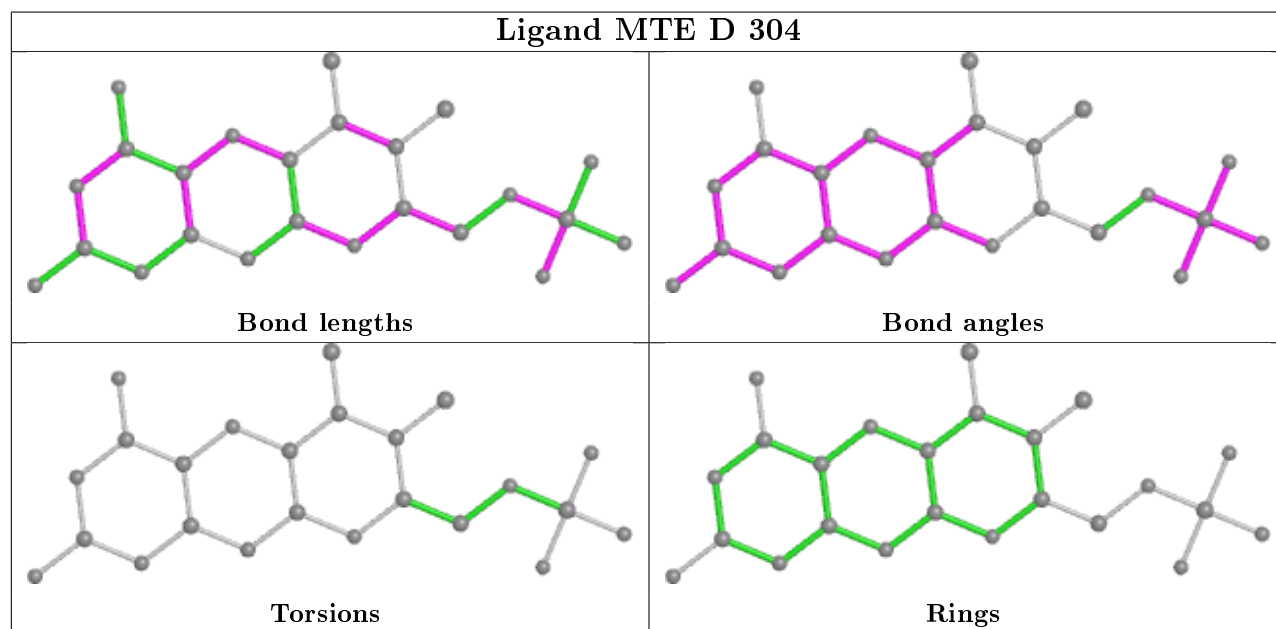
## Ligand MTE A 301



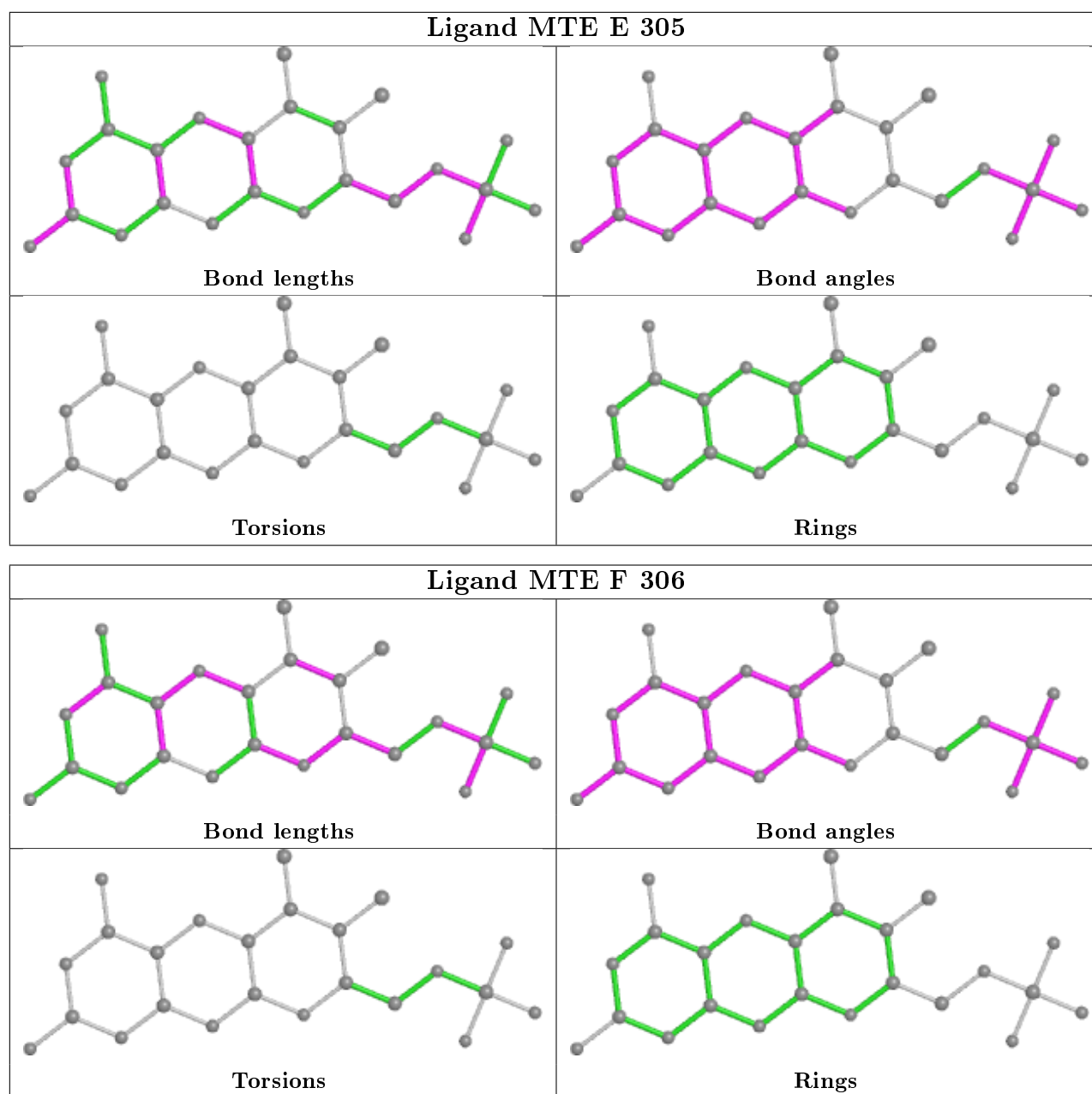
## Ligand MTE J 310



## Ligand MTE D 304







## 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<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	263/298 (88%)	0.17	8 (3%) 50 48	22, 46, 78, 95	0
1	B	262/298 (87%)	0.66	28 (10%) 6 5	33, 68, 94, 106	0
1	C	264/298 (88%)	0.19	10 (3%) 40 38	22, 49, 78, 90	0
1	D	263/298 (88%)	0.20	4 (1%) 73 72	24, 46, 76, 92	0
1	E	262/298 (87%)	0.19	10 (3%) 40 38	25, 46, 80, 97	0
1	F	262/298 (87%)	1.01	48 (18%) 1 1	36, 64, 88, 94	1 (0%)
1	G	261/298 (87%)	0.41	17 (6%) 18 17	21, 49, 84, 100	0
1	H	263/298 (88%)	0.66	25 (9%) 8 7	36, 68, 95, 102	0
1	I	262/298 (87%)	0.72	24 (9%) 9 7	34, 61, 90, 106	0
1	J	262/298 (87%)	0.16	7 (2%) 54 52	20, 44, 73, 91	0
All	All	2624/2980 (88%)	0.44	181 (6%) 16 15	20, 54, 88, 106	1 (0%)

The worst 5 of 181 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	31	ASN	7.6
1	F	130	ALA	7.1
1	G	66	ASP	5.5
1	F	120	LEU	5.3
1	B	256	SER	4.6

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

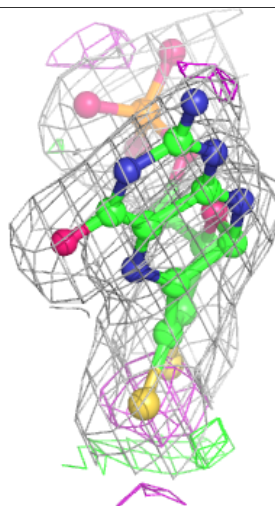
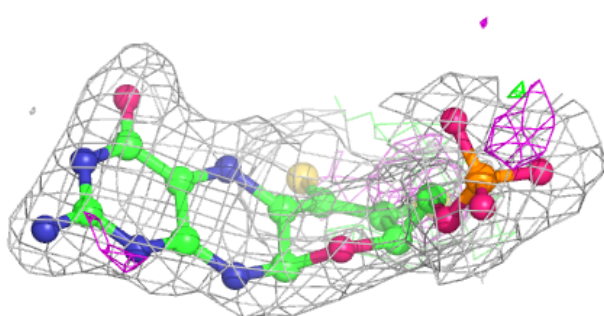
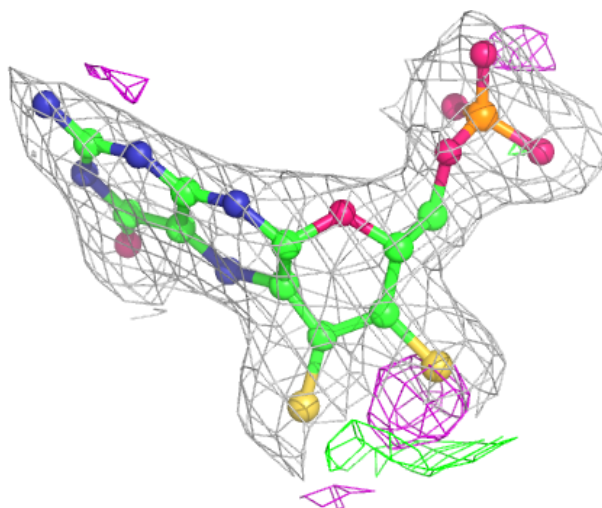
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
3	MTE	I	309	24/24	0.94	0.14	26,42,49,50	0
3	MTE	C	303	24/24	0.95	0.13	15,32,39,42	0
3	MTE	H	308	24/24	0.95	0.15	30,43,48,50	0
2	W	I	409	1/1	0.96	0.15	42,42,42,42	0
3	MTE	F	306	24/24	0.96	0.19	26,38,46,47	0
3	MTE	J	310	24/24	0.97	0.15	8,22,28,29	0
3	MTE	D	304	24/24	0.97	0.14	18,28,35,38	0
3	MTE	G	307	24/24	0.97	0.16	17,26,32,36	0
3	MTE	B	302	24/24	0.97	0.14	22,39,50,56	0
3	MTE	E	305	24/24	0.98	0.13	8,22,33,36	0
3	MTE	A	301	24/24	0.98	0.15	16,23,30,32	0
2	W	F	406	1/1	0.98	0.14	41,41,41,41	0
2	W	B	402	1/1	0.99	0.12	43,43,43,43	0
2	W	H	408	1/1	0.99	0.13	41,41,41,41	0
2	W	J	410	1/1	0.99	0.16	30,30,30,30	0
2	W	E	405	1/1	1.00	0.14	32,32,32,32	0
2	W	A	401	1/1	1.00	0.15	28,28,28,28	0
2	W	D	404	1/1	1.00	0.14	32,32,32,32	0
2	W	G	407	1/1	1.00	0.14	29,29,29,29	0
2	W	C	403	1/1	1.00	0.15	32,32,32,32	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.

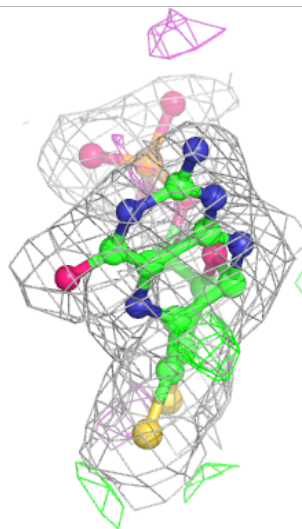
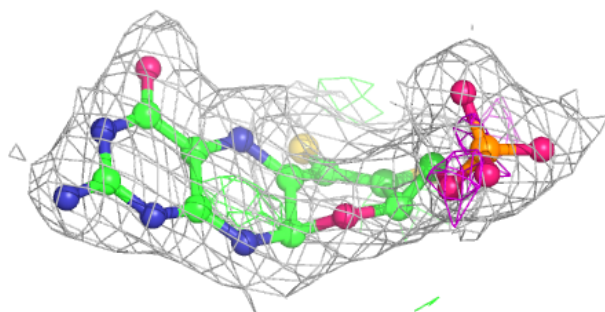
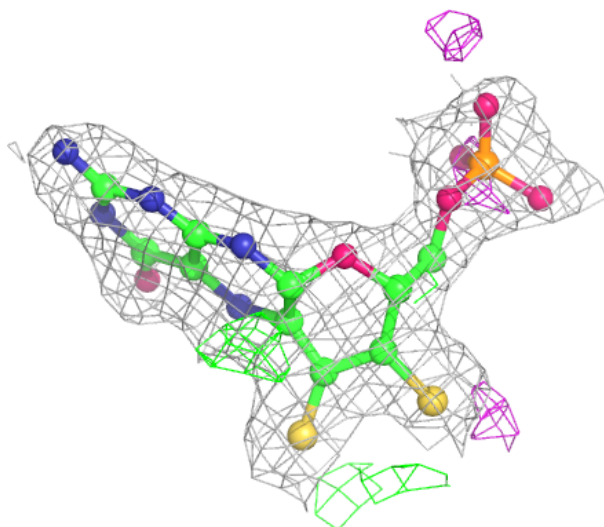
**Electron density around MTE I 309:**

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



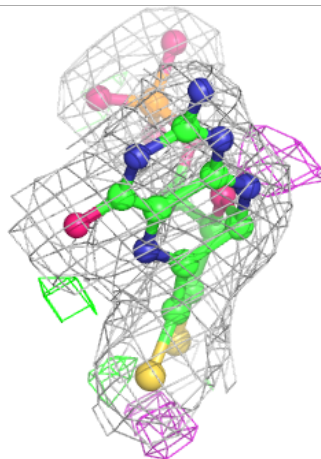
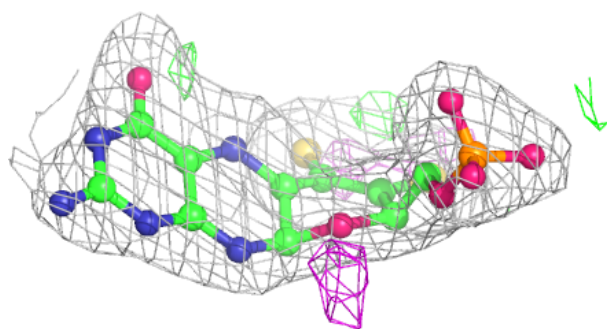
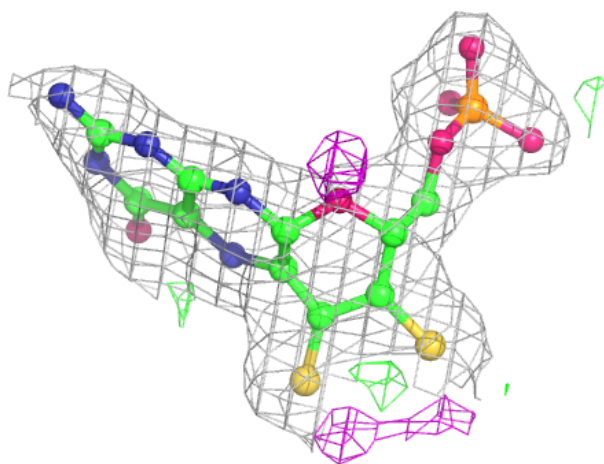
**Electron density around MTE C 303:**

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



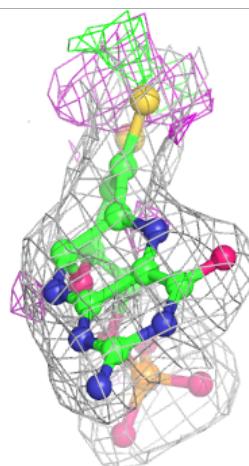
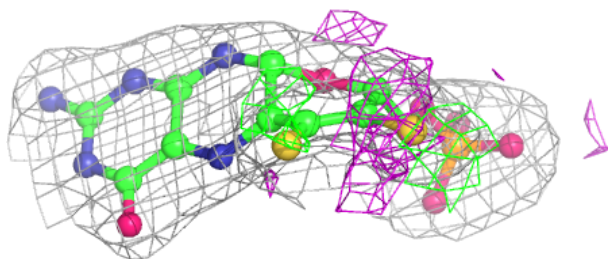
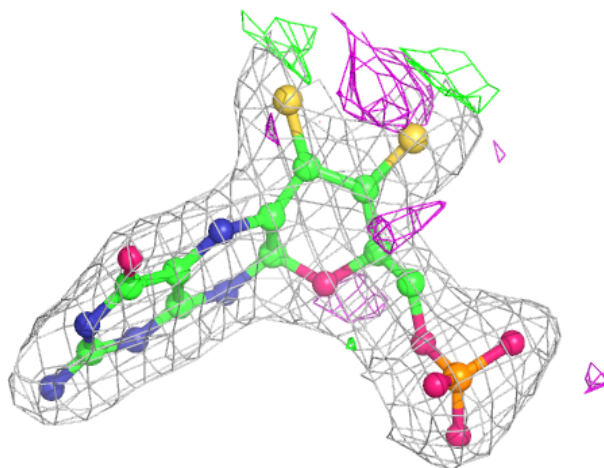
**Electron density around MTE H 308:**

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



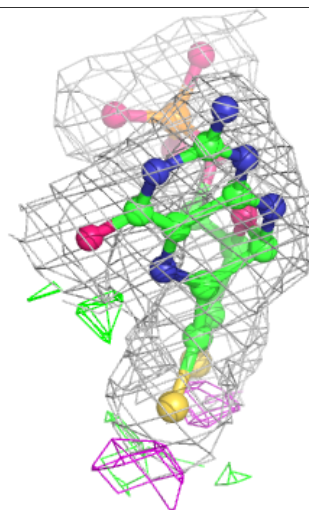
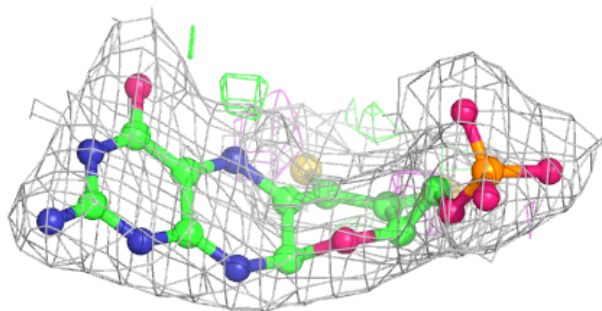
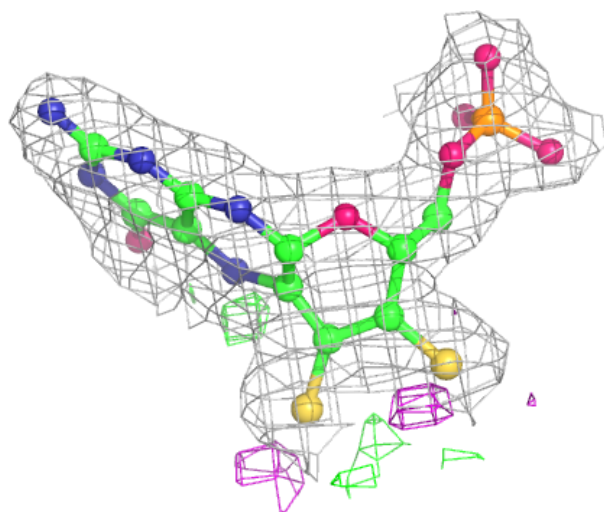
**Electron density around MTE F 306:**

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



**Electron density around MTE J 310:**

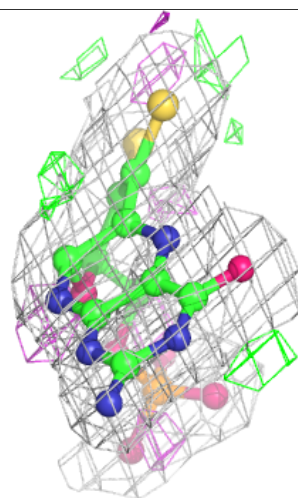
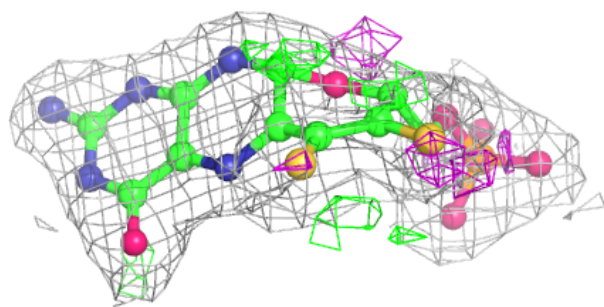
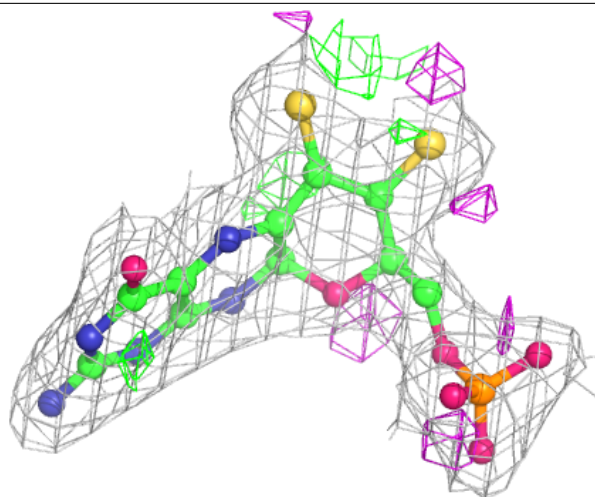
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





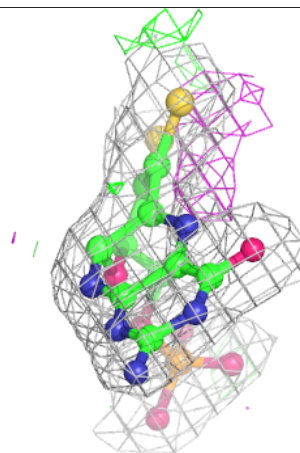
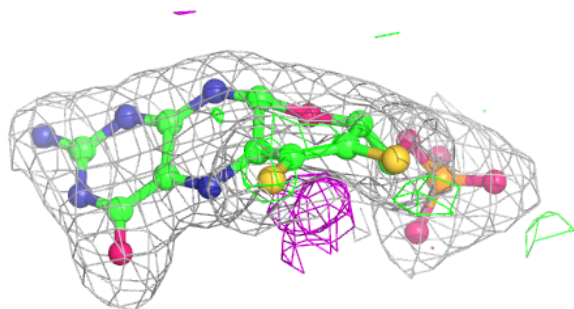
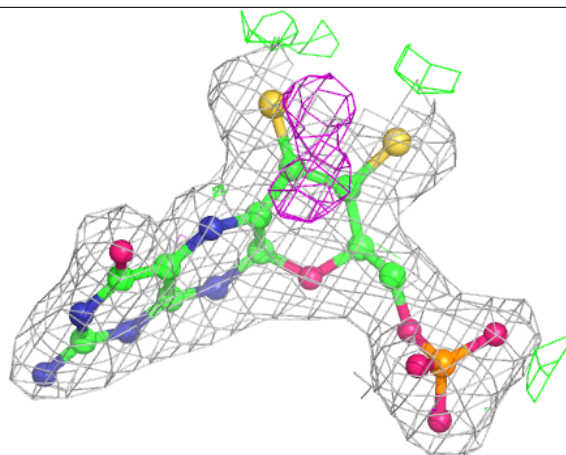
**Electron density around MTE D 304:**

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



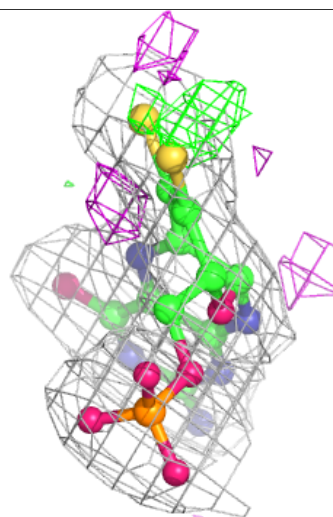
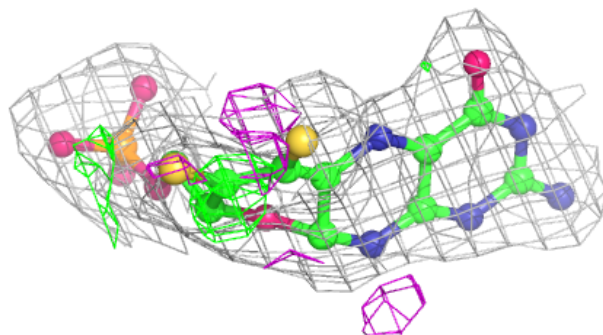
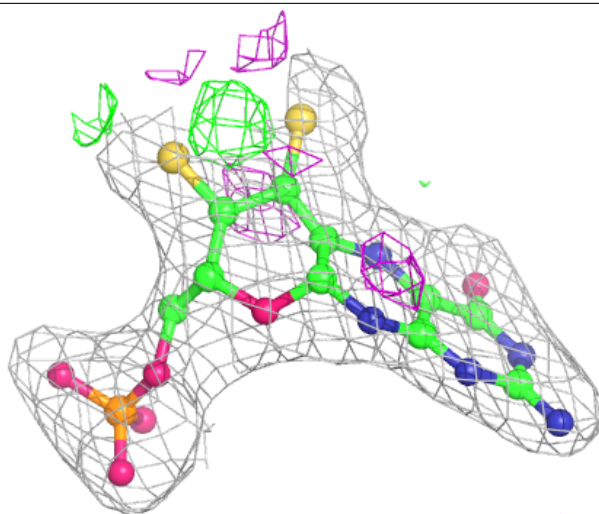
**Electron density around MTE G 307:**

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



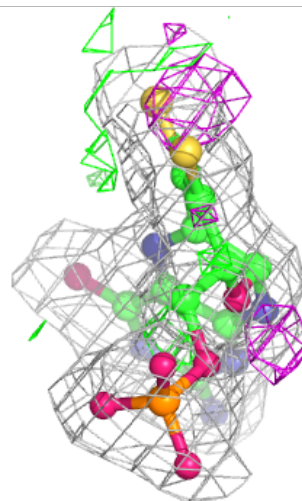
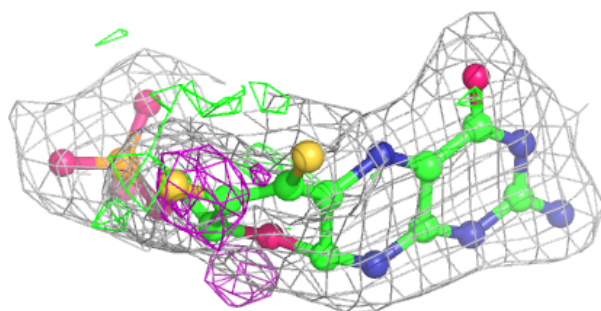
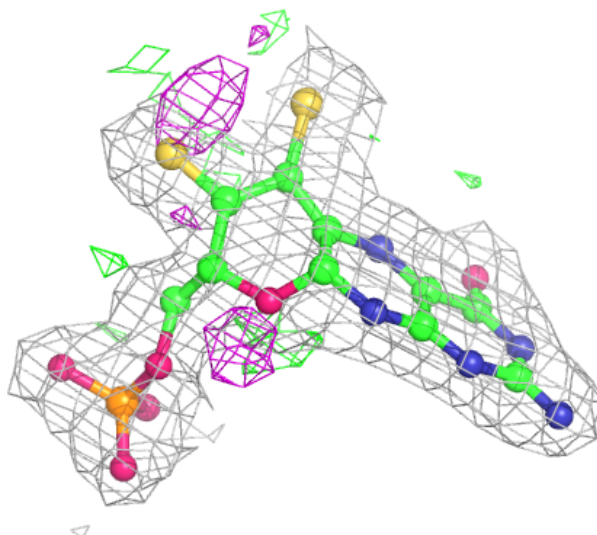
**Electron density around MTE B 302:**

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



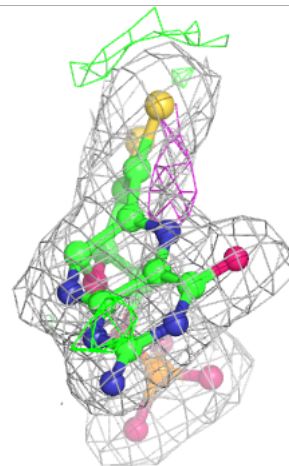
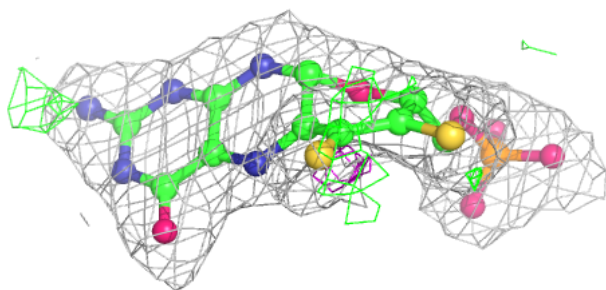
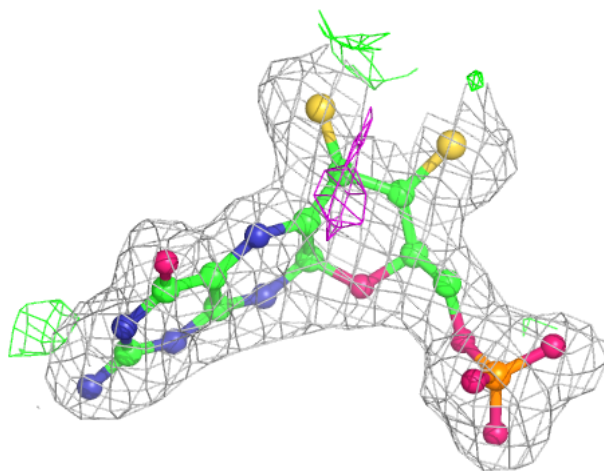
**Electron density around MTE E 305:**

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



**Electron density around MTE A 301:**

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



## 6.5 Other polymers [i](#)

There are no such residues in this entry.