



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 9, 2023 – 11:25 PM EDT

PDB ID : 7MQX  
Title : P. putida mandelate racemase forms an oxobenzoxaborole adduct with 2-formylphenylboronic acid  
Authors : Grandinetti, L.; Bearne, S.L.; St.Maurice, M.  
Deposited on : 2021-05-06  
Resolution : 1.91 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](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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 2.35.1  
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) : 2.35.1

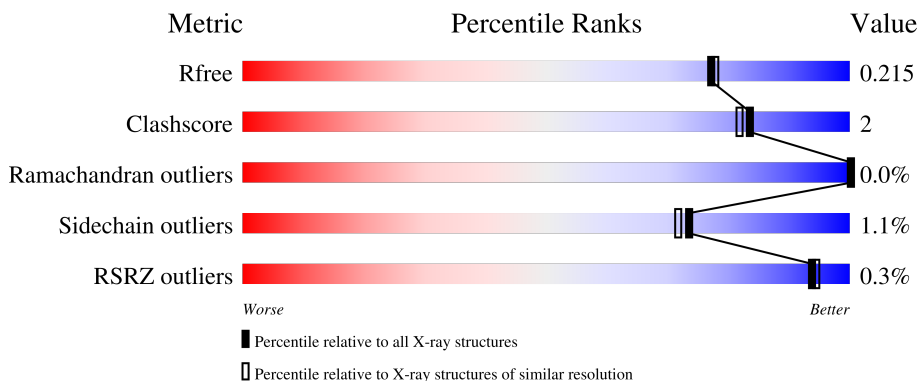
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.91 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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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  $\geq 3$ , 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	383	87% 5% • 7%
1	B	383	88% 5% 7%
1	C	383	87% 5% • 7%
1	D	383	87% 6% 7%
1	E	383	87% 6% 7%

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Mol	Chain	Length	Quality of chain
1	F	383	 <p>% 88% 5% 7%</p>
1	G	383	 <p>90% 7%</p>
1	H	383	 <p>% 87% 6% 7%</p>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 24058 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 Mandelate racemase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	357	2733	1754	467	499	13	0	7	0
1	B	357	2688	1724	458	493	13	0	3	0
1	C	357	2735	1749	467	506	13	0	7	0
1	D	357	2700	1730	462	494	14	0	4	0
1	E	358	2726	1746	461	506	13	2	7	0
1	F	357	2713	1739	461	500	13	1	6	0
1	G	357	2725	1744	465	503	13	1	6	0
1	H	357	2693	1725	461	494	13	0	1	0

There are 192 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	initiating methionine	UNP P11444
A	-22	ALA	-	expression tag	UNP P11444
A	-21	SER	-	expression tag	UNP P11444
A	-20	TRP	-	expression tag	UNP P11444
A	-19	SER	-	expression tag	UNP P11444
A	-18	HIS	-	expression tag	UNP P11444
A	-17	PRO	-	expression tag	UNP P11444
A	-16	GLN	-	expression tag	UNP P11444
A	-15	PHE	-	expression tag	UNP P11444
A	-14	GLU	-	expression tag	UNP P11444
A	-13	LYS	-	expression tag	UNP P11444
A	-12	GLY	-	expression tag	UNP P11444
A	-11	ALA	-	expression tag	UNP P11444

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	LEU	-	expression tag	UNP P11444
A	-9	GLU	-	expression tag	UNP P11444
A	-8	VAL	-	expression tag	UNP P11444
A	-7	LEU	-	expression tag	UNP P11444
A	-6	PHE	-	expression tag	UNP P11444
A	-5	GLN	-	expression tag	UNP P11444
A	-4	GLY	-	expression tag	UNP P11444
A	-3	PRO	-	expression tag	UNP P11444
A	-2	GLY	-	expression tag	UNP P11444
A	-1	TYR	-	expression tag	UNP P11444
A	0	HIS	-	expression tag	UNP P11444
B	-23	MET	-	initiating methionine	UNP P11444
B	-22	ALA	-	expression tag	UNP P11444
B	-21	SER	-	expression tag	UNP P11444
B	-20	TRP	-	expression tag	UNP P11444
B	-19	SER	-	expression tag	UNP P11444
B	-18	HIS	-	expression tag	UNP P11444
B	-17	PRO	-	expression tag	UNP P11444
B	-16	GLN	-	expression tag	UNP P11444
B	-15	PHE	-	expression tag	UNP P11444
B	-14	GLU	-	expression tag	UNP P11444
B	-13	LYS	-	expression tag	UNP P11444
B	-12	GLY	-	expression tag	UNP P11444
B	-11	ALA	-	expression tag	UNP P11444
B	-10	LEU	-	expression tag	UNP P11444
B	-9	GLU	-	expression tag	UNP P11444
B	-8	VAL	-	expression tag	UNP P11444
B	-7	LEU	-	expression tag	UNP P11444
B	-6	PHE	-	expression tag	UNP P11444
B	-5	GLN	-	expression tag	UNP P11444
B	-4	GLY	-	expression tag	UNP P11444
B	-3	PRO	-	expression tag	UNP P11444
B	-2	GLY	-	expression tag	UNP P11444
B	-1	TYR	-	expression tag	UNP P11444
B	0	HIS	-	expression tag	UNP P11444
C	-23	MET	-	initiating methionine	UNP P11444
C	-22	ALA	-	expression tag	UNP P11444
C	-21	SER	-	expression tag	UNP P11444
C	-20	TRP	-	expression tag	UNP P11444
C	-19	SER	-	expression tag	UNP P11444
C	-18	HIS	-	expression tag	UNP P11444
C	-17	PRO	-	expression tag	UNP P11444

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-16	GLN	-	expression tag	UNP P11444
C	-15	PHE	-	expression tag	UNP P11444
C	-14	GLU	-	expression tag	UNP P11444
C	-13	LYS	-	expression tag	UNP P11444
C	-12	GLY	-	expression tag	UNP P11444
C	-11	ALA	-	expression tag	UNP P11444
C	-10	LEU	-	expression tag	UNP P11444
C	-9	GLU	-	expression tag	UNP P11444
C	-8	VAL	-	expression tag	UNP P11444
C	-7	LEU	-	expression tag	UNP P11444
C	-6	PHE	-	expression tag	UNP P11444
C	-5	GLN	-	expression tag	UNP P11444
C	-4	GLY	-	expression tag	UNP P11444
C	-3	PRO	-	expression tag	UNP P11444
C	-2	GLY	-	expression tag	UNP P11444
C	-1	TYR	-	expression tag	UNP P11444
C	0	HIS	-	expression tag	UNP P11444
D	-23	MET	-	initiating methionine	UNP P11444
D	-22	ALA	-	expression tag	UNP P11444
D	-21	SER	-	expression tag	UNP P11444
D	-20	TRP	-	expression tag	UNP P11444
D	-19	SER	-	expression tag	UNP P11444
D	-18	HIS	-	expression tag	UNP P11444
D	-17	PRO	-	expression tag	UNP P11444
D	-16	GLN	-	expression tag	UNP P11444
D	-15	PHE	-	expression tag	UNP P11444
D	-14	GLU	-	expression tag	UNP P11444
D	-13	LYS	-	expression tag	UNP P11444
D	-12	GLY	-	expression tag	UNP P11444
D	-11	ALA	-	expression tag	UNP P11444
D	-10	LEU	-	expression tag	UNP P11444
D	-9	GLU	-	expression tag	UNP P11444
D	-8	VAL	-	expression tag	UNP P11444
D	-7	LEU	-	expression tag	UNP P11444
D	-6	PHE	-	expression tag	UNP P11444
D	-5	GLN	-	expression tag	UNP P11444
D	-4	GLY	-	expression tag	UNP P11444
D	-3	PRO	-	expression tag	UNP P11444
D	-2	GLY	-	expression tag	UNP P11444
D	-1	TYR	-	expression tag	UNP P11444
D	0	HIS	-	expression tag	UNP P11444
E	-23	MET	-	initiating methionine	UNP P11444

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-22	ALA	-	expression tag	UNP P11444
E	-21	SER	-	expression tag	UNP P11444
E	-20	TRP	-	expression tag	UNP P11444
E	-19	SER	-	expression tag	UNP P11444
E	-18	HIS	-	expression tag	UNP P11444
E	-17	PRO	-	expression tag	UNP P11444
E	-16	GLN	-	expression tag	UNP P11444
E	-15	PHE	-	expression tag	UNP P11444
E	-14	GLU	-	expression tag	UNP P11444
E	-13	LYS	-	expression tag	UNP P11444
E	-12	GLY	-	expression tag	UNP P11444
E	-11	ALA	-	expression tag	UNP P11444
E	-10	LEU	-	expression tag	UNP P11444
E	-9	GLU	-	expression tag	UNP P11444
E	-8	VAL	-	expression tag	UNP P11444
E	-7	LEU	-	expression tag	UNP P11444
E	-6	PHE	-	expression tag	UNP P11444
E	-5	GLN	-	expression tag	UNP P11444
E	-4	GLY	-	expression tag	UNP P11444
E	-3	PRO	-	expression tag	UNP P11444
E	-2	GLY	-	expression tag	UNP P11444
E	-1	TYR	-	expression tag	UNP P11444
E	0	HIS	-	expression tag	UNP P11444
F	-23	MET	-	initiating methionine	UNP P11444
F	-22	ALA	-	expression tag	UNP P11444
F	-21	SER	-	expression tag	UNP P11444
F	-20	TRP	-	expression tag	UNP P11444
F	-19	SER	-	expression tag	UNP P11444
F	-18	HIS	-	expression tag	UNP P11444
F	-17	PRO	-	expression tag	UNP P11444
F	-16	GLN	-	expression tag	UNP P11444
F	-15	PHE	-	expression tag	UNP P11444
F	-14	GLU	-	expression tag	UNP P11444
F	-13	LYS	-	expression tag	UNP P11444
F	-12	GLY	-	expression tag	UNP P11444
F	-11	ALA	-	expression tag	UNP P11444
F	-10	LEU	-	expression tag	UNP P11444
F	-9	GLU	-	expression tag	UNP P11444
F	-8	VAL	-	expression tag	UNP P11444
F	-7	LEU	-	expression tag	UNP P11444
F	-6	PHE	-	expression tag	UNP P11444
F	-5	GLN	-	expression tag	UNP P11444

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-4	GLY	-	expression tag	UNP P11444
F	-3	PRO	-	expression tag	UNP P11444
F	-2	GLY	-	expression tag	UNP P11444
F	-1	TYR	-	expression tag	UNP P11444
F	0	HIS	-	expression tag	UNP P11444
G	-23	MET	-	initiating methionine	UNP P11444
G	-22	ALA	-	expression tag	UNP P11444
G	-21	SER	-	expression tag	UNP P11444
G	-20	TRP	-	expression tag	UNP P11444
G	-19	SER	-	expression tag	UNP P11444
G	-18	HIS	-	expression tag	UNP P11444
G	-17	PRO	-	expression tag	UNP P11444
G	-16	GLN	-	expression tag	UNP P11444
G	-15	PHE	-	expression tag	UNP P11444
G	-14	GLU	-	expression tag	UNP P11444
G	-13	LYS	-	expression tag	UNP P11444
G	-12	GLY	-	expression tag	UNP P11444
G	-11	ALA	-	expression tag	UNP P11444
G	-10	LEU	-	expression tag	UNP P11444
G	-9	GLU	-	expression tag	UNP P11444
G	-8	VAL	-	expression tag	UNP P11444
G	-7	LEU	-	expression tag	UNP P11444
G	-6	PHE	-	expression tag	UNP P11444
G	-5	GLN	-	expression tag	UNP P11444
G	-4	GLY	-	expression tag	UNP P11444
G	-3	PRO	-	expression tag	UNP P11444
G	-2	GLY	-	expression tag	UNP P11444
G	-1	TYR	-	expression tag	UNP P11444
G	0	HIS	-	expression tag	UNP P11444
H	-23	MET	-	initiating methionine	UNP P11444
H	-22	ALA	-	expression tag	UNP P11444
H	-21	SER	-	expression tag	UNP P11444
H	-20	TRP	-	expression tag	UNP P11444
H	-19	SER	-	expression tag	UNP P11444
H	-18	HIS	-	expression tag	UNP P11444
H	-17	PRO	-	expression tag	UNP P11444
H	-16	GLN	-	expression tag	UNP P11444
H	-15	PHE	-	expression tag	UNP P11444
H	-14	GLU	-	expression tag	UNP P11444
H	-13	LYS	-	expression tag	UNP P11444
H	-12	GLY	-	expression tag	UNP P11444
H	-11	ALA	-	expression tag	UNP P11444

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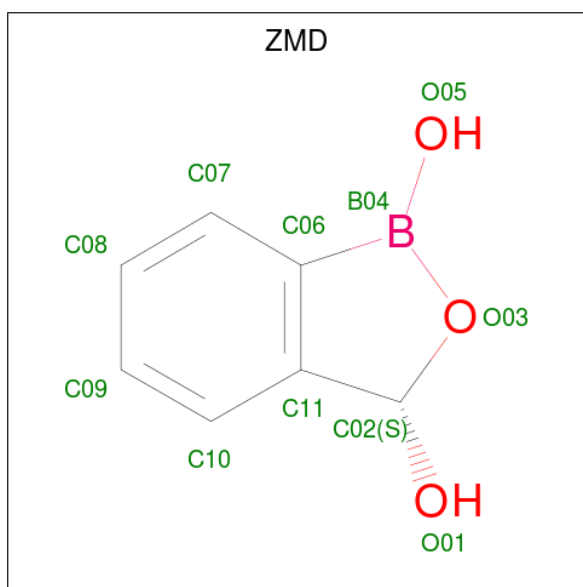
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Chain	Residue	Modelled	Actual	Comment	Reference
H	-10	LEU	-	expression tag	UNP P11444
H	-9	GLU	-	expression tag	UNP P11444
H	-8	VAL	-	expression tag	UNP P11444
H	-7	LEU	-	expression tag	UNP P11444
H	-6	PHE	-	expression tag	UNP P11444
H	-5	GLN	-	expression tag	UNP P11444
H	-4	GLY	-	expression tag	UNP P11444
H	-3	PRO	-	expression tag	UNP P11444
H	-2	GLY	-	expression tag	UNP P11444
H	-1	TYR	-	expression tag	UNP P11444
H	0	HIS	-	expression tag	UNP P11444

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

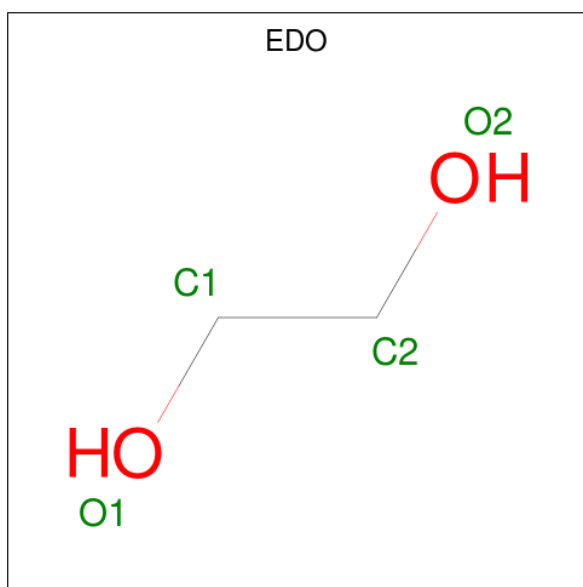
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	E	1	Total Mg 1 1	0	0
2	F	1	Total Mg 1 1	0	0
2	G	1	Total Mg 1 1	0	0
2	H	1	Total Mg 1 1	0	0

- Molecule 3 is (3S)-2,1-benzoxaborole-1,3(3H)-diol (three-letter code: ZMD) (formula: C<sub>7</sub>H<sub>7</sub>BO<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	B	C	O		
3	A	1	Total 11	B 1	C 7	O 3	0	0
3	B	1	Total 11	B 1	C 7	O 3	0	0
3	C	1	Total 11	B 1	C 7	O 3	0	0
3	D	1	Total 11	B 1	C 7	O 3	0	0
3	E	1	Total 11	B 1	C 7	O 3	0	0
3	F	1	Total 11	B 1	C 7	O 3	0	0
3	G	1	Total 11	B 1	C 7	O 3	0	0
3	H	1	Total 11	B 1	C 7	O 3	0	0

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	1	Total C O 4 2 2	0	0
4	F	1	Total C O 4 2 2	0	0
4	F	1	Total C O 4 2 2	0	0
4	F	1	Total C O 4 2 2	0	0
4	G	1	Total C O 4 2 2	0	0
4	G	1	Total C O 4 2 2	0	0
4	G	1	Total C O 4 2 2	0	0
4	H	1	Total C O 4 2 2	0	0
4	H	1	Total C O 4 2 2	0	0
4	H	1	Total C O 4 2 2	0	0


- Molecule 5 is water.

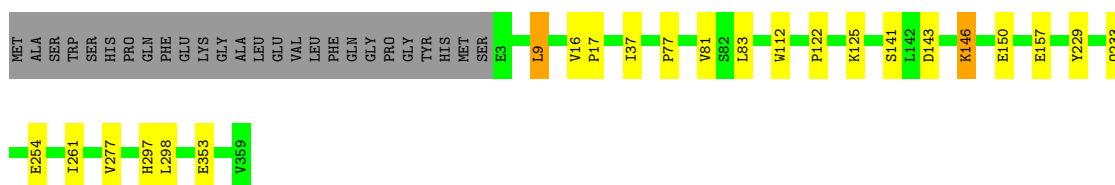
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	278	Total O 278 278	0	0
5	B	275	Total O 275 275	0	0
5	C	270	Total O 270 270	0	0
5	D	243	Total O 243 243	0	0
5	E	287	Total O 287 287	0	0
5	F	267	Total O 267 267	0	0
5	G	266	Total O 266 266	0	0
5	H	267	Total O 267 267	0	0

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


These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

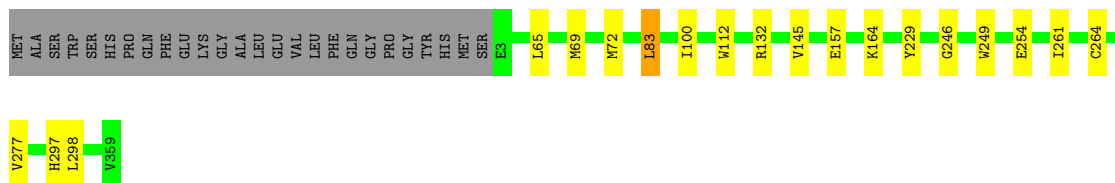
- Molecule 1: Mandelate racemase

Chain A: 



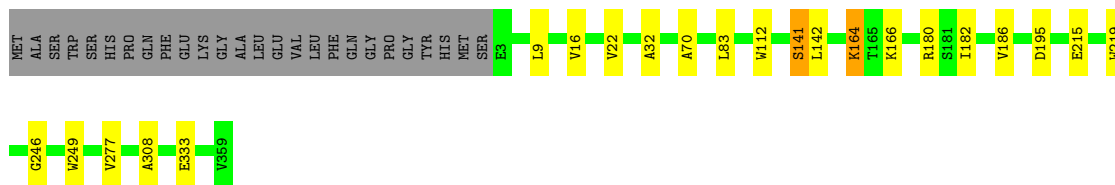
- Molecule 1: Mandelate racemase

Chain B: 



- Molecule 1: Mandelate racemase

Chain C: 



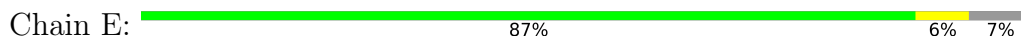
- Molecule 1: Mandelate racemase

Chain D: 

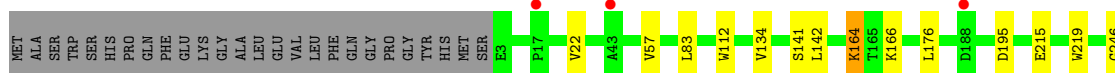
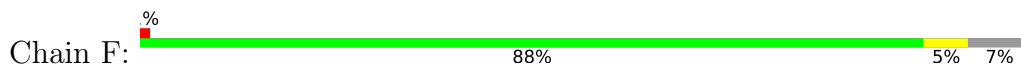




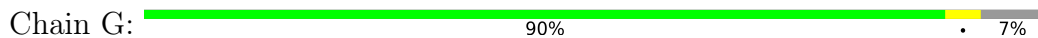
• Molecule 1: Mandelate racemase



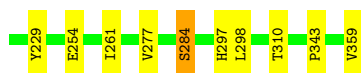
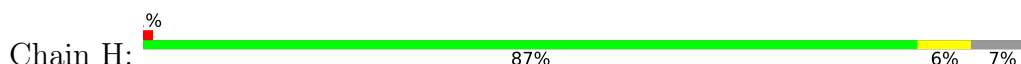
• Molecule 1: Mandelate racemase



• Molecule 1: Mandelate racemase



• Molecule 1: Mandelate racemase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	225.48Å 149.43Å 149.52Å 90.00° 129.34° 90.00°	Depositor
Resolution (Å)	40.61 – 1.91 40.61 – 1.91	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.61-1.91) 99.3 (40.61-1.91)	Depositor EDS
$R_{merge}$	0.19	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.08 (at 1.91Å)	Xtrriage
Refinement program	PHENIX 1.13_2998	Depositor
R, $R_{free}$	0.187 , 0.215 0.187 , 0.215	Depositor DCC
$R_{free}$ test set	14797 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.2	Xtrriage
Anisotropy	0.414	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 47.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.217 for -h+k-l,-l,-k 0.064 for -h-k-l,l,k 0.087 for -h-2*l,-k,l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	24058	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.87% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, ZMD, MG

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	0.43	0/2804	0.57	0/3818
1	B	0.39	0/2750	0.55	1/3747 (0.0%)
1	C	0.42	0/2795	0.57	0/3807
1	D	0.39	0/2768	0.55	0/3769
1	E	0.45	0/2803	0.57	0/3819
1	F	0.41	0/2784	0.55	0/3790
1	G	0.43	0/2790	0.57	0/3799
1	H	0.40	0/2752	0.55	0/3749
All	All	0.42	0/22246	0.56	1/30298 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	83	LEU	CA-CB-CG	5.21	127.28	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2733	0	2796	15	0
1	B	2688	0	2718	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2735	0	2771	15	0
1	D	2700	0	2738	13	0
1	E	2726	0	2770	16	0
1	F	2713	0	2757	14	0
1	G	2725	0	2767	9	0
1	H	2693	0	2737	14	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
3	A	11	0	0	0	0
3	B	11	0	0	0	0
3	C	11	0	0	0	0
3	D	11	0	0	1	0
3	E	11	0	0	0	0
3	F	11	0	0	1	0
3	G	11	0	0	0	0
3	H	11	0	0	0	0
4	A	12	0	18	1	0
4	B	16	0	24	2	0
4	C	8	0	12	0	0
4	D	12	0	18	1	0
4	E	12	0	18	3	0
4	F	12	0	18	1	0
4	G	12	0	18	0	0
4	H	12	0	18	1	0
5	A	278	0	0	4	0
5	B	275	0	0	5	0
5	C	270	0	0	1	0
5	D	243	0	0	2	0
5	E	287	0	0	5	0
5	F	267	0	0	3	0
5	G	266	0	0	0	0
5	H	267	0	0	5	0
All	All	24058	0	22198	107	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 2.

All (107) 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:230[A]:GLU:OE1	1:D:234[A]:ARG:NH1	1.97	0.98
1:H:254[B]:GLU:OE1	5:H:501:HOH:O	1.94	0.86
1:E:132:ARG:HH12	4:E:404:EDO:H12	1.42	0.84
1:H:92:CYS:HA	5:H:711:HOH:O	1.88	0.74
1:C:164:LYS:HE3	1:C:195:ASP:HB2	1.70	0.73
1:D:254[A]:GLU:OE1	5:D:501:HOH:O	2.05	0.73
1:A:254[A]:GLU:OE1	5:A:501:HOH:O	2.07	0.72
1:H:284:SER:HB2	1:H:310:THR:HG23	1.72	0.71
1:H:254[A]:GLU:OE1	5:H:502:HOH:O	2.11	0.68
1:B:254:GLU:OE1	5:B:501:HOH:O	2.13	0.67
1:A:233:GLN:OE1	5:A:502:HOH:O	2.13	0.66
1:B:132:ARG:HH12	4:B:405:EDO:H22	1.63	0.64
1:D:284:SER:HB2	1:D:310:THR:HG23	1.80	0.64
1:C:164:LYS:HE3	1:C:195:ASP:CB	2.30	0.62
1:E:132:ARG:NH1	4:E:404:EDO:H12	2.13	0.61
1:H:180:ARG:NH1	5:H:504:HOH:O	2.29	0.61
1:A:112:TRP:CD1	1:A:277:VAL:HB	2.36	0.60
1:C:112:TRP:CD1	1:C:277:VAL:HB	2.36	0.58
1:H:81:VAL:HG12	4:H:404:EDO:H11	1.85	0.58
1:E:72:MET:HE1	1:F:57:VAL:HG11	1.86	0.56
1:F:112:TRP:CD1	1:F:277:VAL:HB	2.41	0.56
1:E:311:PRO:HA	4:E:404:EDO:H11	1.88	0.56
1:H:112:TRP:CD1	1:H:277:VAL:HB	2.41	0.56
1:F:164:LYS:HE2	1:F:166:LYS:HG3	1.89	0.55
1:A:353:GLU:OE2	5:A:503:HOH:O	2.18	0.55
1:E:112:TRP:CD1	1:E:277:VAL:HB	2.42	0.55
1:B:112:TRP:CD1	1:B:277:VAL:HB	2.43	0.54
1:F:254[A]:GLU:OE1	5:F:501:HOH:O	2.18	0.54
1:D:112:TRP:CD1	1:D:277:VAL:HB	2.42	0.54
1:G:112:TRP:CD1	1:G:277:VAL:HB	2.43	0.53
1:G:137:TYR:HE2	1:G:164[B]:LYS:HE2	1.72	0.53
1:H:16:VAL:CG2	1:H:32:ALA:HB3	2.38	0.53
1:A:157:GLU:OE1	5:A:504:HOH:O	2.19	0.53
1:A:81:VAL:HG12	4:A:404:EDO:H11	1.92	0.52
1:A:16:VAL:HG13	1:A:17:PRO:HD2	1.92	0.51
1:B:264:CYS:O	5:B:502:HOH:O	2.19	0.51
1:F:164:LYS:HE3	1:F:195:ASP:CB	2.40	0.50
1:E:75[A]:ASN:ND2	5:E:501:HOH:O	2.17	0.49
1:B:65:LEU:HD21	1:B:100:ILE:HA	1.94	0.49
1:D:65:LEU:HD21	1:D:100:ILE:HA	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:23:HIS:O	1:E:141[B]:SER:OG	2.22	0.48
1:C:9:LEU:HD23	1:C:70:ALA:HB2	1.96	0.47
1:A:77:PRO:O	1:A:83:LEU:HD21	2.15	0.47
1:C:22:VAL:HG13	1:C:141[A]:SER:OG	2.15	0.47
1:C:182:ILE:O	1:C:186:VAL:HG22	2.15	0.47
1:E:76:GLU:OE2	5:E:502:HOH:O	2.21	0.47
1:D:202:VAL:HG21	1:D:234[A]:ARG:NH2	2.29	0.47
1:E:277:VAL:HG22	1:E:306:LEU:HD21	1.96	0.46
4:B:404:EDO:H11	1:C:308:ALA:O	2.16	0.46
1:E:230[B]:GLU:HG2	5:E:750:HOH:O	2.15	0.46
1:B:246:GLY:HA2	1:B:249:TRP:CD2	2.51	0.46
1:A:146[B]:LYS:HB2	1:A:146[B]:LYS:HE2	1.48	0.46
1:G:77:PRO:O	1:G:83:LEU:HD21	2.16	0.46
1:H:59:LEU:CD2	1:H:359:VAL:HG21	2.46	0.46
1:H:65:LEU:HD21	1:H:100:ILE:HA	1.97	0.46
1:A:122:PRO:HD2	1:A:125:LYS:HD2	1.97	0.45
1:C:333:GLU:OE2	5:C:501:HOH:O	2.20	0.45
1:F:284[A]:SER:HB2	1:F:310:THR:HG23	1.98	0.45
1:G:51:LEU:HD12	1:G:66:LEU:HD11	1.99	0.45
1:D:229:TYR:C	1:D:261:ILE:HD11	2.37	0.45
1:B:145:VAL:HG23	5:B:638:HOH:O	2.15	0.45
1:F:308:ALA:HB2	5:F:558:HOH:O	2.16	0.45
1:C:16:VAL:CG2	1:C:32:ALA:HB3	2.46	0.45
1:F:164:LYS:HE2	1:F:166:LYS:CG	2.47	0.45
1:C:16:VAL:HG22	1:C:32:ALA:HB3	1.99	0.45
1:H:297:HIS:CE1	1:H:298:LEU:HG	2.52	0.44
1:F:164:LYS:HE3	1:F:195:ASP:HB2	1.98	0.44
1:F:176:LEU:HD22	1:F:215:GLU:HG3	1.99	0.44
1:C:180:ARG:NH2	1:C:215:GLU:OE2	2.49	0.43
1:G:137:TYR:CE2	1:G:164[B]:LYS:HE2	2.52	0.43
1:E:65:LEU:HD21	1:E:100:ILE:HA	1.99	0.43
1:H:59:LEU:HD23	1:H:359:VAL:HG21	2.00	0.43
1:A:229:TYR:C	1:A:261:ILE:HD11	2.38	0.43
1:B:69:MET:HG2	1:B:72:MET:HE1	1.99	0.43
1:B:157:GLU:OE1	5:B:503:HOH:O	2.21	0.43
1:D:72:MET:HE1	1:D:91:PHE:CZ	2.54	0.43
1:G:16:VAL:CG2	1:G:32:ALA:HB3	2.49	0.43
1:A:141[A]:SER:OG	1:A:143:ASP:OD1	2.29	0.43
1:F:246:GLY:HA2	1:F:249:TRP:CD2	2.53	0.43
1:D:145:VAL:HG23	5:D:662:HOH:O	2.19	0.42
1:B:83:LEU:HD22	5:B:687:HOH:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:252:PRO:HD3	5:E:637:HOH:O	2.18	0.42
1:H:229:TYR:C	1:H:261:ILE:HD11	2.40	0.42
1:E:201:ASP:OD1	1:E:203:PRO:HD2	2.19	0.42
1:C:164:LYS:HE2	1:C:166:LYS:HG3	2.02	0.42
1:B:229:TYR:CB	1:B:261:ILE:HD11	2.50	0.42
1:G:167:ILE:HD12	1:G:196:TYR:CZ	2.55	0.42
4:F:404:EDO:H22	1:G:308:ALA:O	2.20	0.42
1:H:96:TYR:HB2	5:H:711:HOH:O	2.19	0.42
1:C:164:LYS:HE2	1:C:166:LYS:CG	2.49	0.42
1:C:246:GLY:HA2	1:C:249:TRP:CD2	2.55	0.42
1:F:22:VAL:HG13	1:F:141[B]:SER:OG	2.20	0.42
1:E:333:GLU:OE2	5:E:504:HOH:O	2.22	0.41
1:G:16:VAL:HG23	1:G:32:ALA:HB3	2.02	0.41
1:D:166:LYS:HE3	3:D:402:ZMD:C06	2.50	0.41
1:E:295:SER:OG	1:E:314:HIS:HB3	2.21	0.41
1:D:69:MET:HG2	1:D:72:MET:CE	2.50	0.41
1:C:83[B]:LEU:HD22	1:C:83[B]:LEU:HA	1.87	0.41
1:D:31:THR:O	1:D:33:PRO:HD3	2.20	0.41
1:E:195:ASP:HA	1:E:221:GLU:HB3	2.03	0.41
1:F:134:VAL:CG2	5:F:558:HOH:O	2.68	0.41
1:A:146[A]:LYS:HD3	1:A:150:GLU:OE2	2.20	0.40
1:B:297:HIS:CE1	1:B:298:LEU:HG	2.56	0.40
1:F:166:LYS:CE	3:F:402:ZMD:C06	2.99	0.40
1:D:308:ALA:O	4:D:404:EDO:H22	2.20	0.40
1:A:9[A]:LEU:HD21	1:A:37:ILE:HG23	2.03	0.40
1:A:297:HIS:CE1	1:A:298:LEU:HG	2.57	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	362/383 (94%)	353 (98%)	9 (2%)	0	100	100
1	B	358/383 (94%)	352 (98%)	6 (2%)	0	100	100
1	C	362/383 (94%)	353 (98%)	9 (2%)	0	100	100
1	D	359/383 (94%)	351 (98%)	7 (2%)	1 (0%)	41	31
1	E	363/383 (95%)	353 (97%)	10 (3%)	0	100	100
1	F	361/383 (94%)	354 (98%)	7 (2%)	0	100	100
1	G	361/383 (94%)	355 (98%)	6 (2%)	0	100	100
1	H	356/383 (93%)	349 (98%)	7 (2%)	0	100	100
All	All	2882/3064 (94%)	2820 (98%)	61 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	298	LEU

### 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	284/301 (94%)	280 (99%)	4 (1%)	67	63
1	B	274/301 (91%)	273 (100%)	1 (0%)	91	91
1	C	285/301 (95%)	280 (98%)	5 (2%)	59	53
1	D	277/301 (92%)	274 (99%)	3 (1%)	73	72
1	E	285/301 (95%)	283 (99%)	2 (1%)	84	83
1	F	282/301 (94%)	278 (99%)	4 (1%)	67	63
1	G	283/301 (94%)	282 (100%)	1 (0%)	91	91
1	H	278/301 (92%)	271 (98%)	7 (2%)	47	39
All	All	2248/2408 (93%)	2221 (99%)	27 (1%)	73	69

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

Mol	Chain	Res	Type
1	A	9[A]	LEU
1	A	9[B]	LEU
1	A	146[A]	LYS
1	A	146[B]	LYS
1	B	164	LYS
1	C	141[A]	SER
1	C	141[B]	SER
1	C	142	LEU
1	C	164	LYS
1	C	219	TRP
1	D	164	LYS
1	D	219	TRP
1	D	284	SER
1	E	83	LEU
1	E	142	LEU
1	F	83	LEU
1	F	142	LEU
1	F	164	LYS
1	F	219	TRP
1	G	214	GLN
1	H	10	ARG
1	H	15	ASN
1	H	142	LEU
1	H	164	LYS
1	H	219	TRP
1	H	284	SER
1	H	343	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	233	GLN

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 40 ligands modelled in this entry, 8 are monoatomic - leaving 32 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
4	EDO	H	405	-	3,3,3	0.44	0	2,2,2	0.38	0
4	EDO	G	404	-	3,3,3	0.53	0	2,2,2	0.30	0
4	EDO	G	403	-	3,3,3	0.41	0	2,2,2	0.24	0
4	EDO	F	405	-	3,3,3	0.46	0	2,2,2	0.28	0
4	EDO	C	403	-	3,3,3	0.43	0	2,2,2	0.62	0
4	EDO	D	403	-	3,3,3	0.47	0	2,2,2	0.31	0
3	ZMD	D	402	2,1	11,12,12	4.21	3 (27%)	15,17,17	2.17	3 (20%)
4	EDO	F	403	-	3,3,3	0.51	0	2,2,2	0.29	0
3	ZMD	E	402	2,1	11,12,12	4.20	4 (36%)	15,17,17	2.45	5 (33%)
4	EDO	D	404	-	3,3,3	0.51	0	2,2,2	0.16	0
4	EDO	A	404	-	3,3,3	0.48	0	2,2,2	0.14	0
4	EDO	E	404	-	3,3,3	0.45	0	2,2,2	0.47	0
3	ZMD	B	402	2,1	11,12,12	4.18	3 (27%)	15,17,17	2.39	3 (20%)
4	EDO	H	403	-	3,3,3	0.44	0	2,2,2	0.22	0
4	EDO	F	404	-	3,3,3	0.50	0	2,2,2	0.29	0
4	EDO	E	405	-	3,3,3	0.45	0	2,2,2	0.35	0
3	ZMD	G	402	2,1	11,12,12	3.93	4 (36%)	15,17,17	2.59	3 (20%)
4	EDO	A	405	-	3,3,3	0.53	0	2,2,2	0.36	0
4	EDO	B	406	-	3,3,3	0.46	0	2,2,2	0.39	0
4	EDO	H	404	-	3,3,3	0.48	0	2,2,2	0.57	0
4	EDO	B	403	-	3,3,3	0.45	0	2,2,2	0.46	0
4	EDO	G	405	-	3,3,3	0.46	0	2,2,2	0.42	0
3	ZMD	A	402	2,1	11,12,12	4.37	3 (27%)	15,17,17	2.63	5 (33%)
4	EDO	B	404	-	3,3,3	0.49	0	2,2,2	0.23	0
4	EDO	D	405	-	3,3,3	0.41	0	2,2,2	0.42	0
3	ZMD	H	402	2,1	11,12,12	4.10	3 (27%)	15,17,17	2.42	4 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	EDO	A	403	-	3,3,3	0.55	0	2,2,2	0.20	0
4	EDO	E	403	-	3,3,3	0.37	0	2,2,2	0.43	0
3	ZMD	F	402	2,1	11,12,12	4.51	3 (27%)	15,17,17	2.36	5 (33%)
3	ZMD	C	402	2,1	11,12,12	4.11	3 (27%)	15,17,17	2.28	3 (20%)
4	EDO	B	405	-	3,3,3	0.46	0	2,2,2	0.57	0
4	EDO	C	404	-	3,3,3	0.45	0	2,2,2	0.39	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	H	405	-	-	0/1/1/1	-
4	EDO	G	404	-	-	0/1/1/1	-
4	EDO	G	403	-	-	0/1/1/1	-
4	EDO	F	405	-	-	0/1/1/1	-
4	EDO	C	403	-	-	0/1/1/1	-
4	EDO	D	403	-	-	0/1/1/1	-
3	ZMD	D	402	2,1	-	-	0/2/2/2
4	EDO	F	403	-	-	0/1/1/1	-
4	EDO	D	404	-	-	0/1/1/1	-
3	ZMD	E	402	2,1	-	-	0/2/2/2
4	EDO	A	404	-	-	0/1/1/1	-
4	EDO	E	404	-	-	0/1/1/1	-
3	ZMD	B	402	2,1	-	-	0/2/2/2
4	EDO	H	403	-	-	0/1/1/1	-
4	EDO	F	404	-	-	0/1/1/1	-
4	EDO	E	405	-	-	0/1/1/1	-
4	EDO	A	405	-	-	0/1/1/1	-
3	ZMD	G	402	2,1	-	-	0/2/2/2
4	EDO	B	406	-	-	0/1/1/1	-
4	EDO	H	404	-	-	0/1/1/1	-
4	EDO	B	403	-	-	0/1/1/1	-
4	EDO	G	405	-	-	0/1/1/1	-
3	ZMD	A	402	2,1	-	-	0/2/2/2
4	EDO	B	404	-	-	1/1/1/1	-
4	EDO	D	405	-	-	0/1/1/1	-
4	EDO	A	403	-	-	0/1/1/1	-
3	ZMD	H	402	2,1	-	-	0/2/2/2
4	EDO	E	403	-	-	0/1/1/1	-
3	ZMD	F	402	2,1	-	-	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ZMD	C	402	2,1	-	-	0/2/2/2
4	EDO	B	405	-	-	0/1/1/1	-
4	EDO	C	404	-	-	0/1/1/1	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	402	ZMD	C11-C06	-14.07	1.30	1.40
3	A	402	ZMD	C11-C06	-13.75	1.31	1.40
3	B	402	ZMD	C11-C06	-13.17	1.31	1.40
3	D	402	ZMD	C11-C06	-13.13	1.31	1.40
3	E	402	ZMD	C11-C06	-12.88	1.31	1.40
3	C	402	ZMD	C11-C06	-12.80	1.31	1.40
3	H	402	ZMD	C11-C06	-12.62	1.31	1.40
3	G	402	ZMD	C11-C06	-12.01	1.32	1.40
3	H	402	ZMD	C11-C02	3.67	1.53	1.50
3	F	402	ZMD	C11-C02	2.97	1.53	1.50
3	E	402	ZMD	C11-C02	2.91	1.52	1.50
3	G	402	ZMD	C11-C02	2.90	1.52	1.50
3	D	402	ZMD	C11-C02	2.90	1.52	1.50
3	E	402	ZMD	C10-C11	2.87	1.43	1.39
3	C	402	ZMD	C10-C11	2.86	1.43	1.39
3	F	402	ZMD	C10-C11	2.70	1.43	1.39
3	G	402	ZMD	C10-C11	2.66	1.43	1.39
3	D	402	ZMD	C10-C11	2.58	1.43	1.39
3	A	402	ZMD	C10-C11	2.57	1.43	1.39
3	B	402	ZMD	C10-C11	2.56	1.43	1.39
3	E	402	ZMD	C07-C06	2.52	1.44	1.40
3	B	402	ZMD	C11-C02	2.51	1.52	1.50
3	A	402	ZMD	C11-C02	2.49	1.52	1.50
3	C	402	ZMD	C11-C02	2.29	1.52	1.50
3	H	402	ZMD	C10-C11	2.24	1.42	1.39
3	G	402	ZMD	C07-C06	2.17	1.44	1.40

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	402	ZMD	O05-B04-C06	-6.14	109.97	129.13
3	E	402	ZMD	O05-B04-C06	-6.07	110.17	129.13
3	G	402	ZMD	O03-B04-C06	-6.00	105.02	108.43
3	A	402	ZMD	O03-B04-C06	-5.73	105.17	108.43
3	B	402	ZMD	O05-B04-C06	-5.72	111.25	129.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	402	ZMD	O05-B04-C06	-5.54	111.83	129.13
3	A	402	ZMD	O05-B04-C06	-5.51	111.92	129.13
3	H	402	ZMD	O03-B04-C06	-5.47	105.32	108.43
3	C	402	ZMD	O03-B04-C06	-5.30	105.42	108.43
3	E	402	ZMD	O03-B04-C06	-5.27	105.43	108.43
3	A	402	ZMD	O05-B04-O03	-5.27	110.97	121.96
3	H	402	ZMD	O05-B04-C06	-5.22	112.83	129.13
3	G	402	ZMD	O05-B04-C06	-5.20	112.89	129.13
3	G	402	ZMD	O05-B04-O03	-5.13	111.26	121.96
3	C	402	ZMD	O05-B04-C06	-5.04	113.38	129.13
3	B	402	ZMD	O05-B04-O03	-4.53	112.51	121.96
3	B	402	ZMD	O03-B04-C06	-4.51	105.86	108.43
3	H	402	ZMD	O05-B04-O03	-4.37	112.84	121.96
3	D	402	ZMD	O05-B04-O03	-4.01	113.59	121.96
3	D	402	ZMD	O03-B04-C06	-3.87	106.23	108.43
3	F	402	ZMD	O03-B04-C06	-3.84	106.24	108.43
3	F	402	ZMD	O05-B04-O03	-3.79	114.06	121.96
3	C	402	ZMD	O05-B04-O03	-3.74	114.16	121.96
3	E	402	ZMD	O05-B04-O03	-3.51	114.64	121.96
3	F	402	ZMD	B04-C06-C07	-2.24	130.99	135.01
3	E	402	ZMD	B04-C06-C07	-2.20	131.07	135.01
3	F	402	ZMD	C10-C11-C06	-2.15	119.63	121.37
3	A	402	ZMD	O01-C02-O03	2.07	112.84	110.59
3	H	402	ZMD	B04-C06-C07	-2.01	131.41	135.01
3	E	402	ZMD	C10-C11-C06	-2.01	119.74	121.37
3	A	402	ZMD	B04-C06-C07	-2.00	131.43	135.01

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	404	EDO	O1-C1-C2-O2

There are no ring outliers.

9 monomers are involved in 11 short contacts:

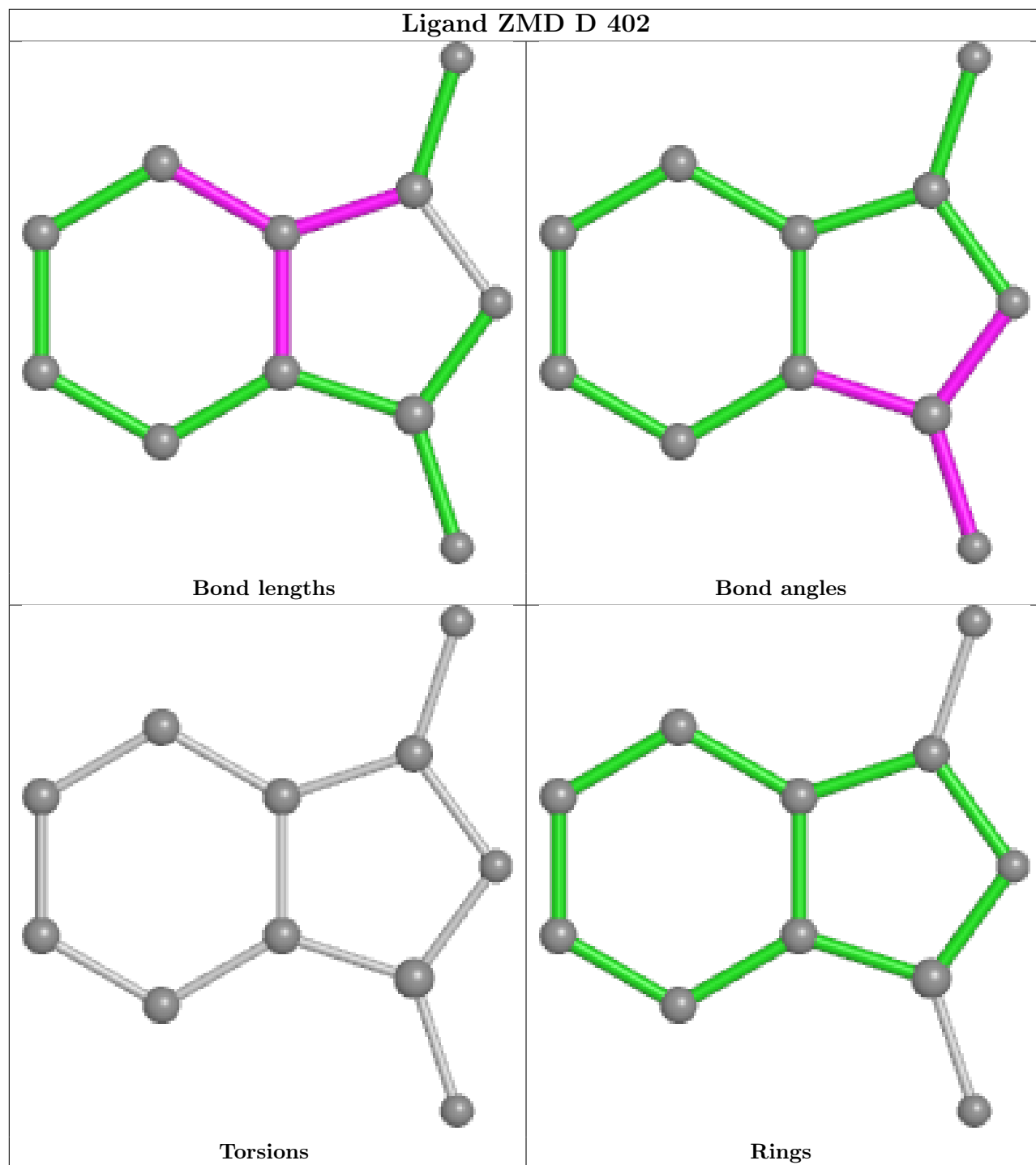
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	402	ZMD	1	0
4	D	404	EDO	1	0
4	A	404	EDO	1	0
4	E	404	EDO	3	0

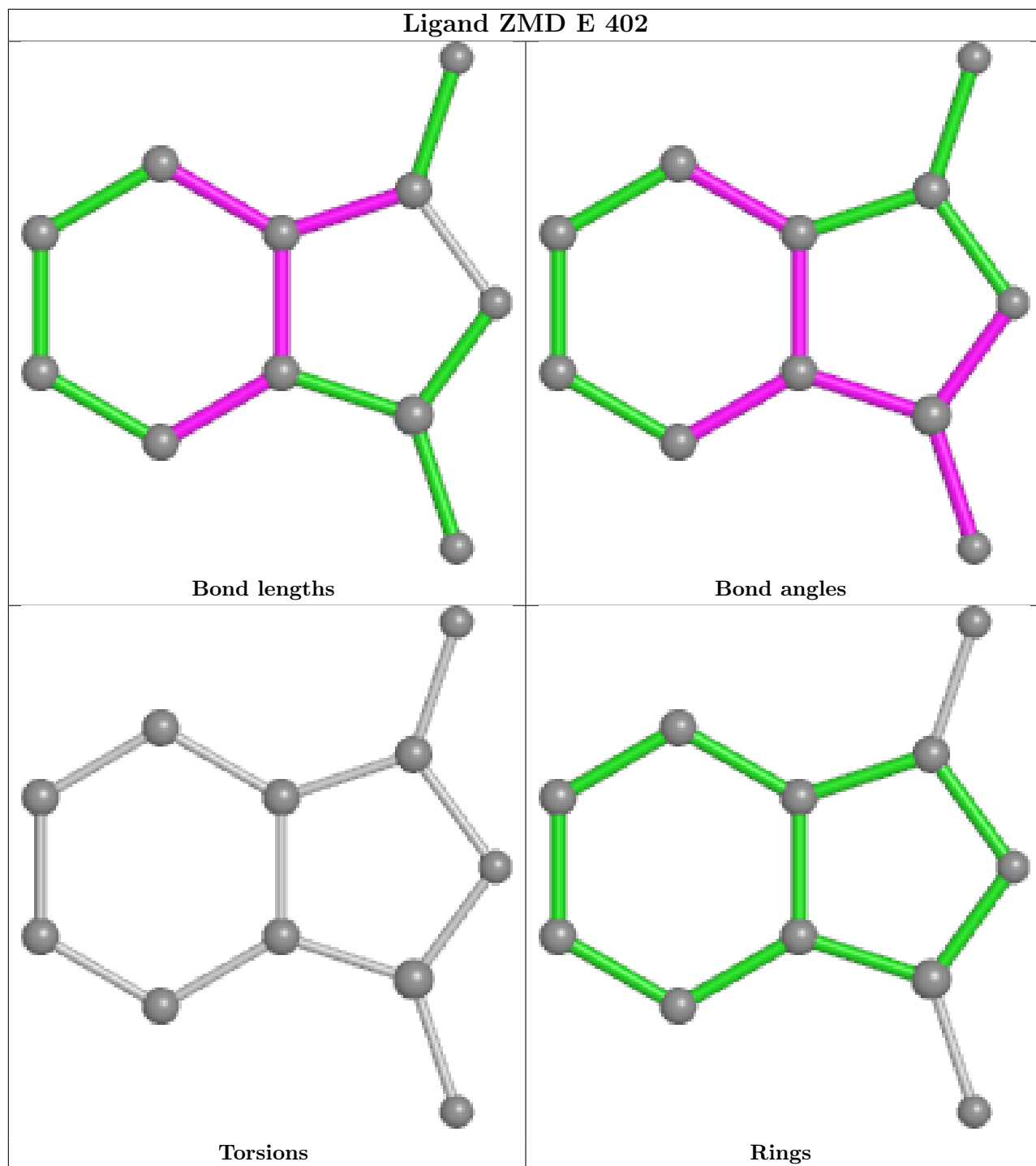
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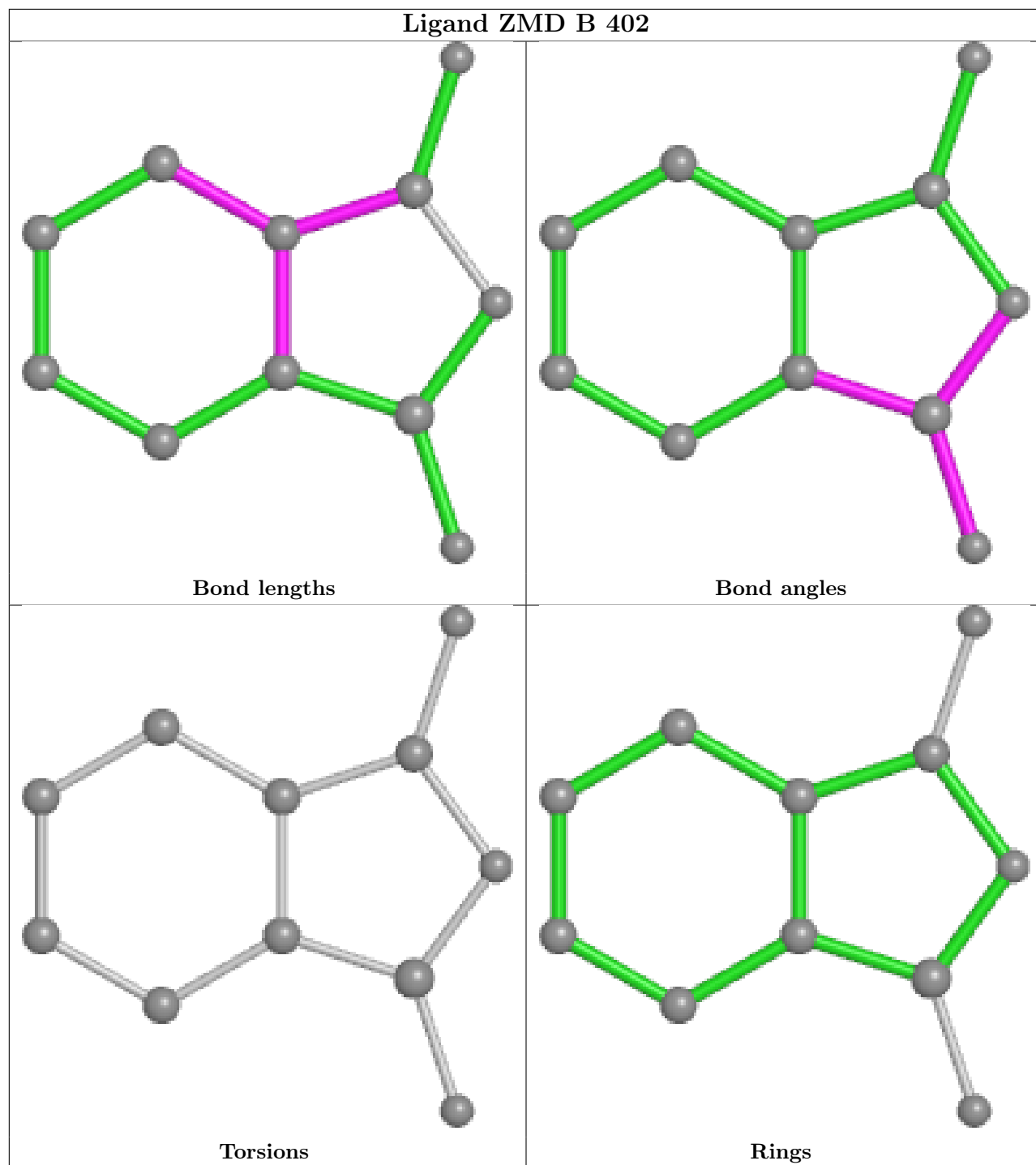
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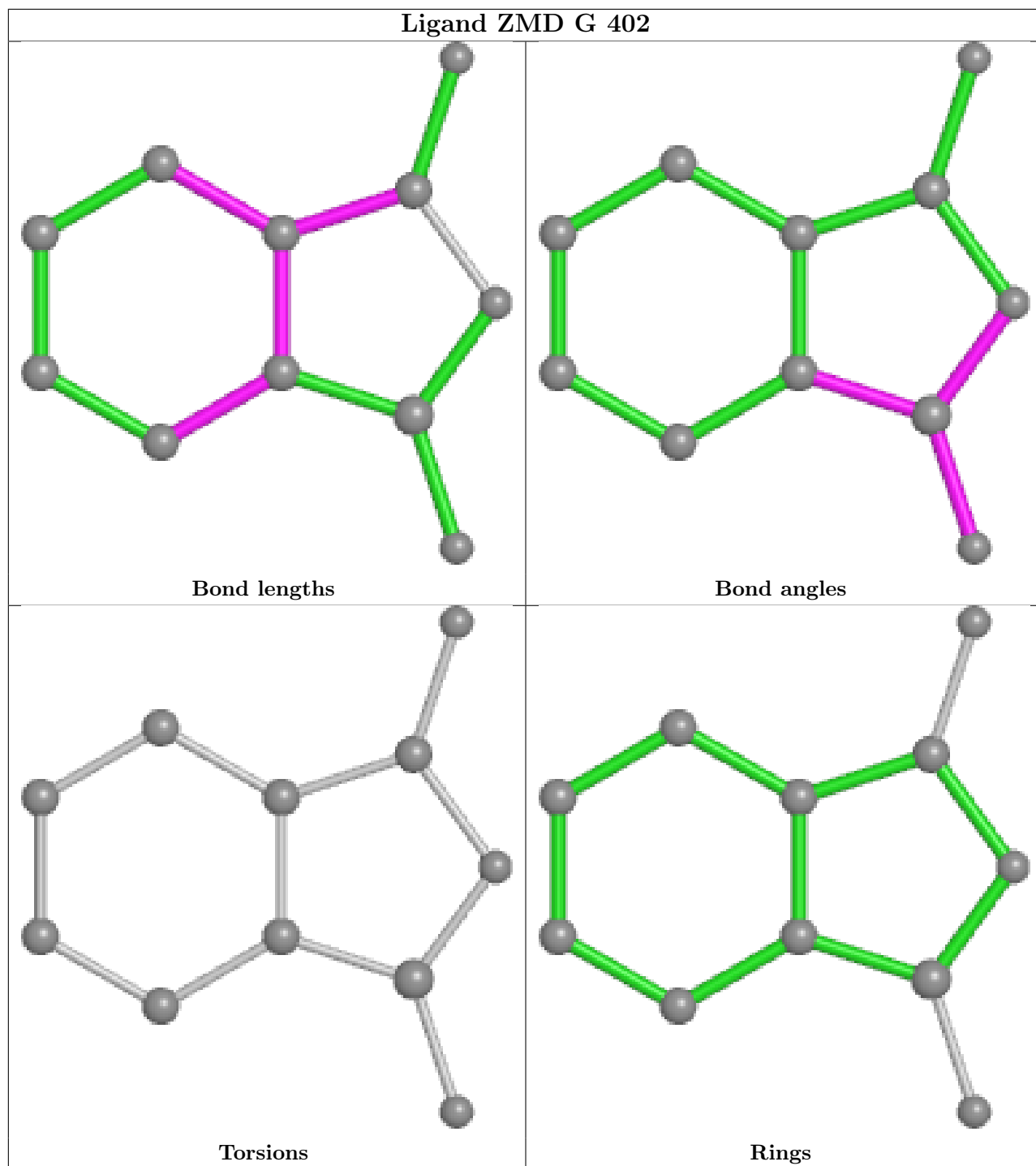
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	404	EDO	1	0
4	H	404	EDO	1	0
4	B	404	EDO	1	0
3	F	402	ZMD	1	0
4	B	405	EDO	1	0

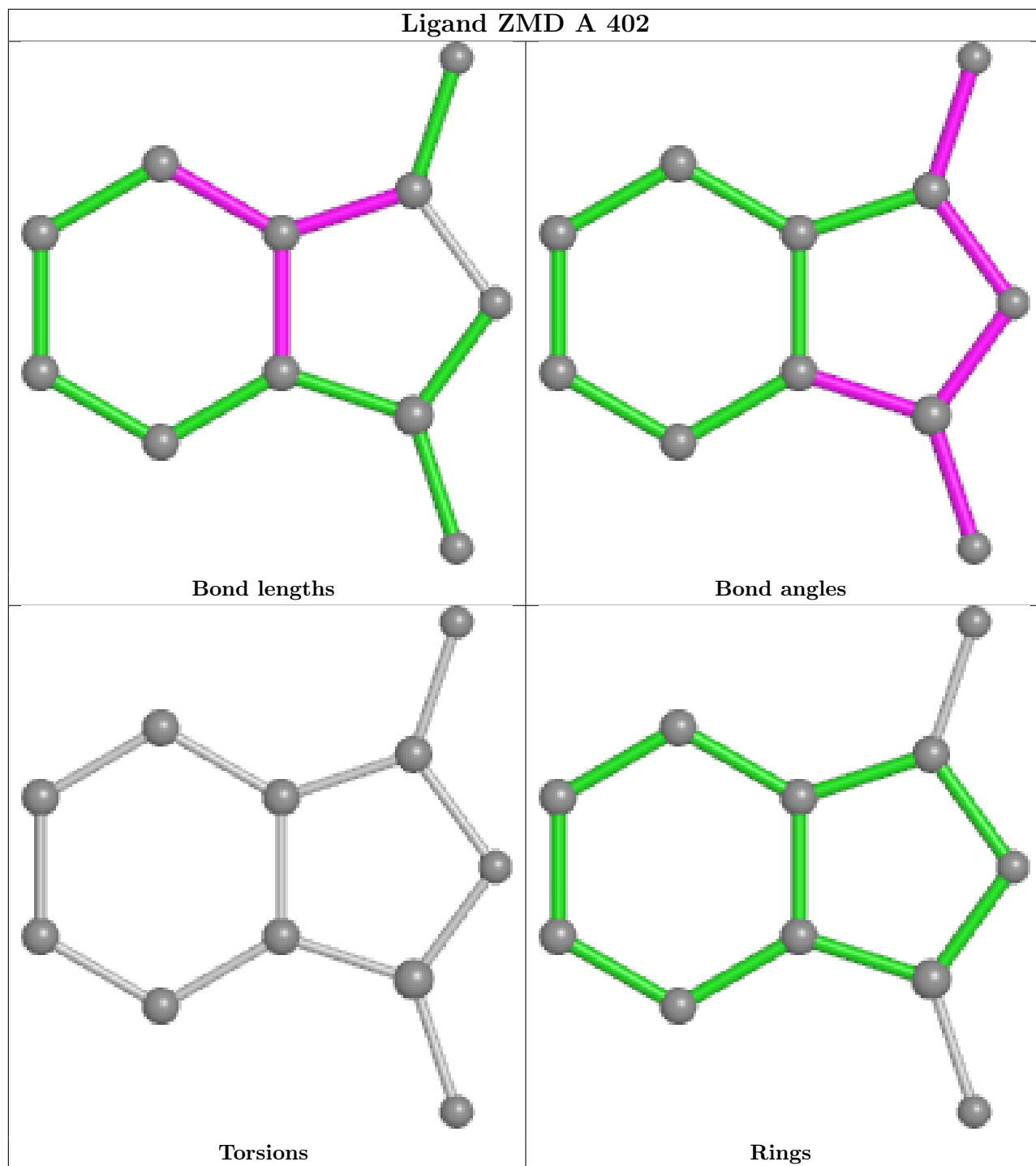
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



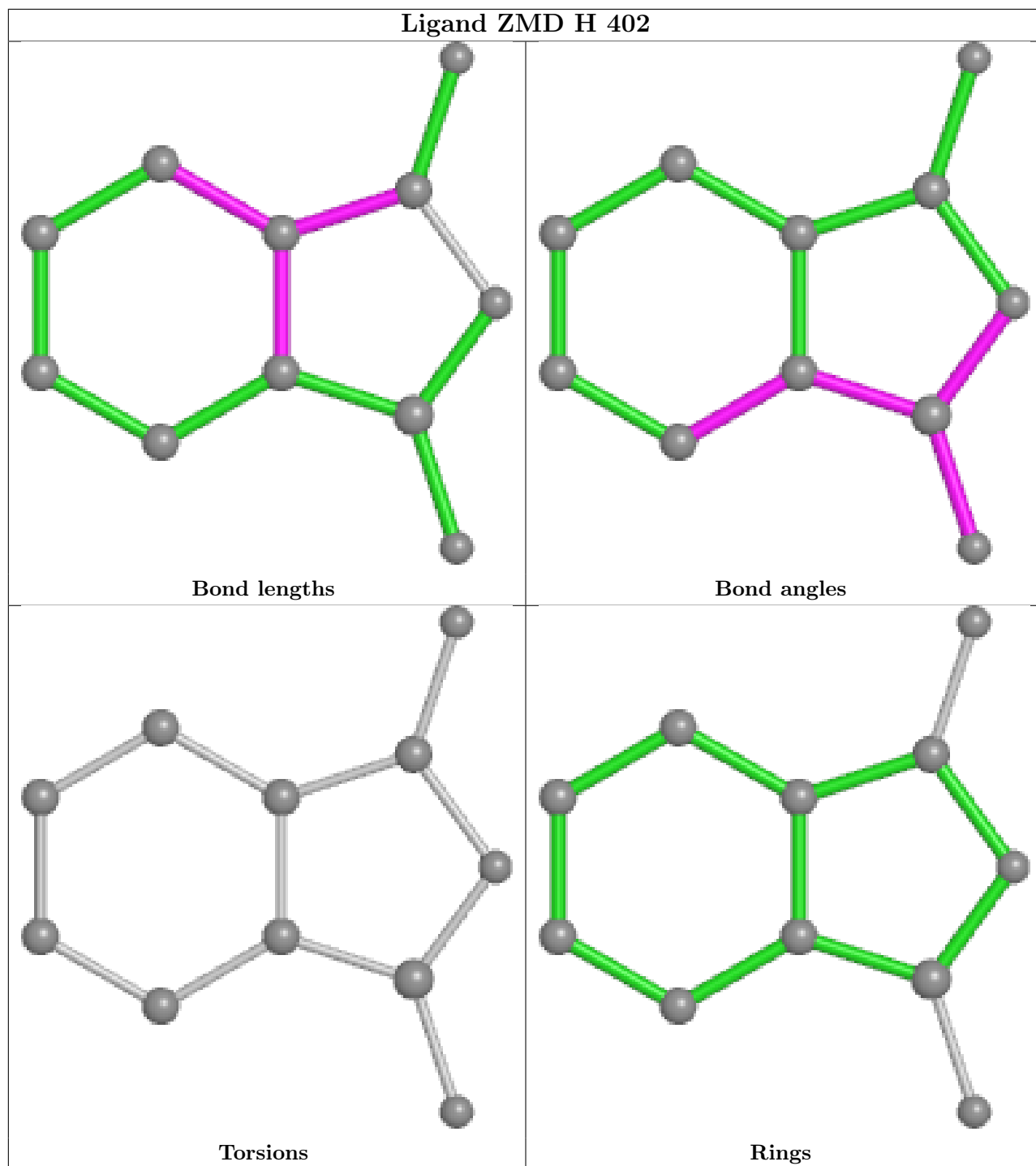


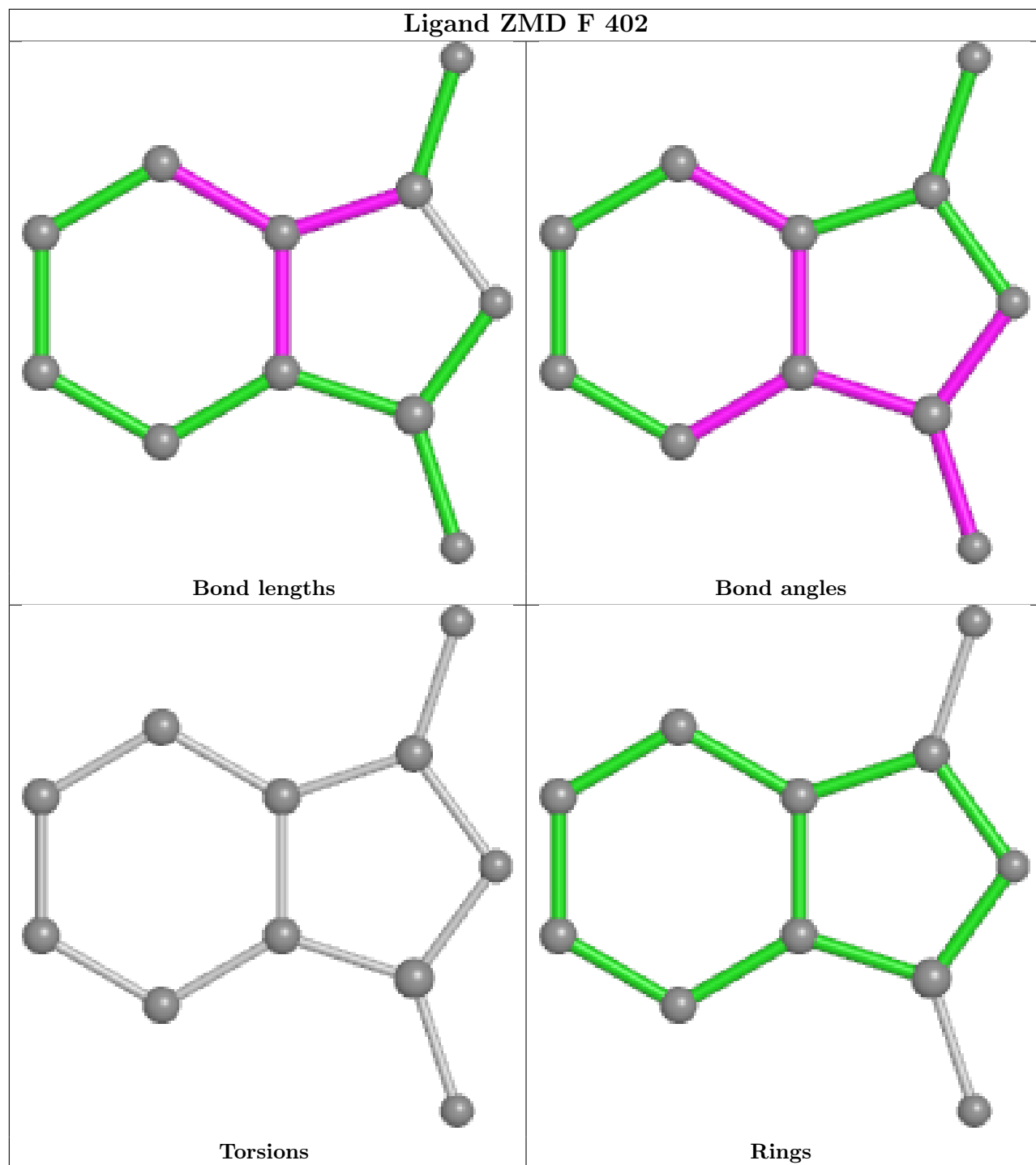


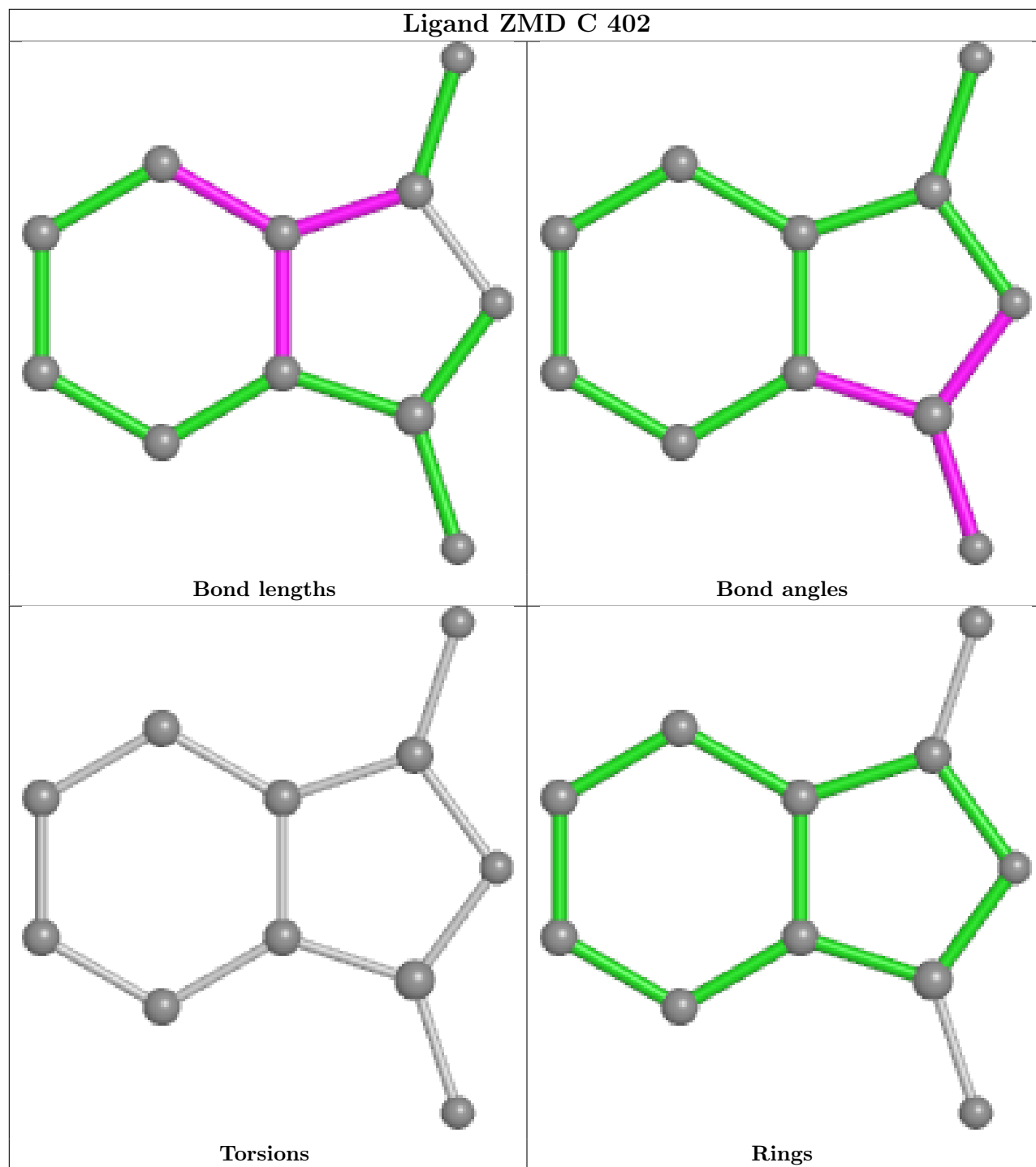












## 5.7 Other polymers [i](#)

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	357/383 (93%)	-0.34	0	100   100	12, 18, 29, 40	2 (0%)
1	B	357/383 (93%)	-0.31	0	100   100	15, 21, 32, 40	0
1	C	357/383 (93%)	-0.34	0	100   100	12, 18, 27, 37	2 (0%)
1	D	357/383 (93%)	-0.30	1 (0%)	94   94	14, 21, 32, 41	1 (0%)
1	E	358/383 (93%)	-0.39	1 (0%)	94   94	11, 18, 27, 48	3 (0%)
1	F	357/383 (93%)	-0.25	4 (1%)	80   82	14, 21, 32, 43	2 (0%)
1	G	357/383 (93%)	-0.41	1 (0%)	94   94	12, 18, 27, 42	3 (0%)
1	H	357/383 (93%)	-0.22	2 (0%)	89   90	14, 21, 33, 42	1 (0%)
All	All	2857/3064 (93%)	-0.32	9 (0%)	94   94	11, 20, 31, 48	14 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	2	SER	2.9
1	D	43	ALA	2.8
1	G	17	PRO	2.8
1	H	185	ALA	2.6
1	F	359	VAL	2.2
1	H	44	GLY	2.2
1	F	43	ALA	2.1
1	F	188	ASP	2.0
1	F	17	PRO	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 monosaccharides in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<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
4	EDO	E	404	4/4	0.81	0.16	21,22,32,34	0
4	EDO	H	404	4/4	0.82	0.17	24,27,28,32	0
4	EDO	F	404	4/4	0.83	0.16	25,27,29,33	0
4	EDO	B	405	4/4	0.87	0.21	27,30,31,31	0
3	ZMD	H	402	11/11	0.89	0.11	19,21,24,24	0
4	EDO	B	404	4/4	0.90	0.13	21,24,30,33	0
3	ZMD	B	402	11/11	0.90	0.10	19,22,23,24	0
4	EDO	D	405	4/4	0.90	0.11	24,26,32,33	0
3	ZMD	D	402	11/11	0.91	0.10	19,21,24,26	0
4	EDO	A	404	4/4	0.91	0.09	26,27,28,31	0
3	ZMD	C	402	11/11	0.92	0.10	15,19,21,22	0
3	ZMD	E	402	11/11	0.92	0.10	15,17,19,19	0
4	EDO	B	406	4/4	0.92	0.09	21,26,29,35	0
4	EDO	C	404	4/4	0.92	0.12	22,24,27,30	0
4	EDO	F	405	4/4	0.93	0.15	23,26,29,35	0
4	EDO	G	404	4/4	0.93	0.09	26,27,27,31	0
3	ZMD	A	402	11/11	0.93	0.09	17,19,22,22	0
4	EDO	H	405	4/4	0.93	0.11	25,28,31,33	0
4	EDO	A	405	4/4	0.94	0.07	22,23,25,34	0
3	ZMD	F	402	11/11	0.94	0.11	17,21,25,25	0
3	ZMD	G	402	11/11	0.94	0.08	15,17,20,21	0
4	EDO	H	403	4/4	0.95	0.09	26,28,29,30	0
4	EDO	D	404	4/4	0.95	0.10	27,28,33,34	0
4	EDO	D	403	4/4	0.95	0.08	22,26,28,30	0
4	EDO	B	403	4/4	0.96	0.07	25,27,28,30	0
4	EDO	G	403	4/4	0.96	0.08	20,24,25,25	0
4	EDO	E	403	4/4	0.96	0.08	20,21,22,22	0
4	EDO	A	403	4/4	0.96	0.06	24,25,25,25	0
4	EDO	E	405	4/4	0.96	0.08	22,23,23,25	0
4	EDO	C	403	4/4	0.96	0.08	19,21,22,24	0
2	MG	B	401	1/1	0.97	0.07	17,17,17,17	0
4	EDO	F	403	4/4	0.97	0.06	26,27,28,31	0

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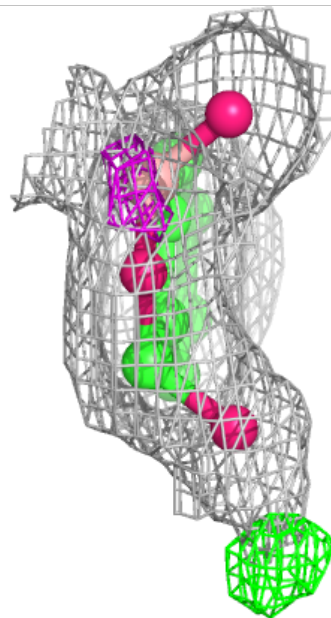
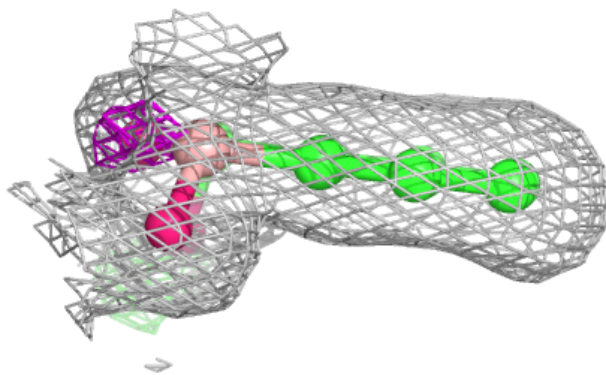
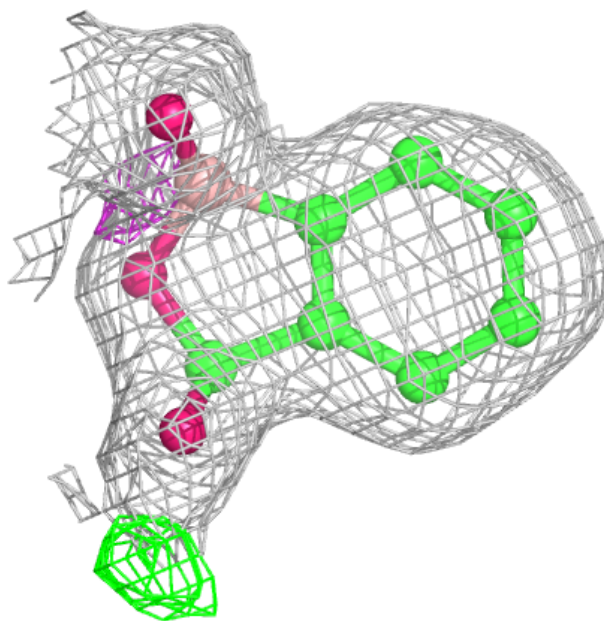
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	EDO	G	405	4/4	0.97	0.08	23,23,27,32	0
2	MG	F	401	1/1	0.98	0.07	18,18,18,18	0
2	MG	G	401	1/1	0.98	0.12	15,15,15,15	0
2	MG	A	401	1/1	0.99	0.12	13,13,13,13	0
2	MG	C	401	1/1	0.99	0.08	12,12,12,12	0
2	MG	H	401	1/1	0.99	0.08	17,17,17,17	0
2	MG	D	401	1/1	0.99	0.06	14,14,14,14	0
2	MG	E	401	1/1	0.99	0.06	13,13,13,13	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.

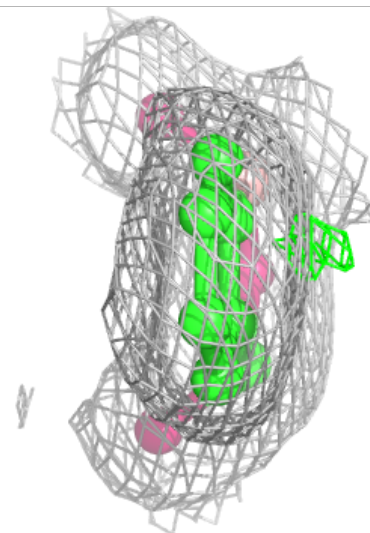
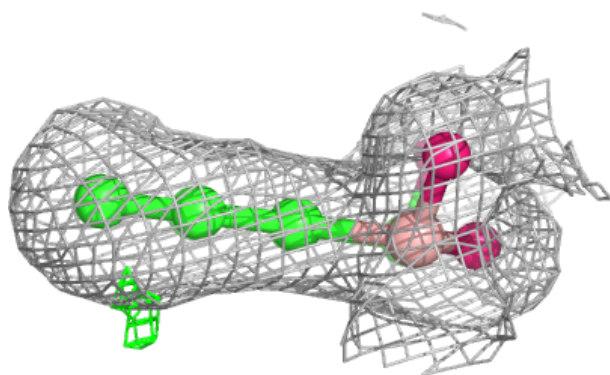
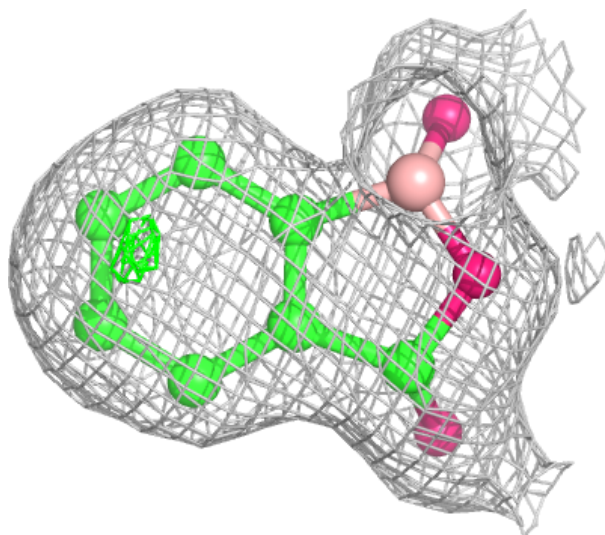
**Electron density around ZMD H 402:**

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



**Electron density around ZMD B 402:**

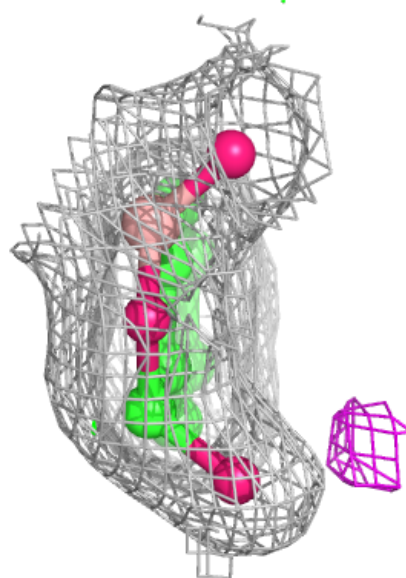
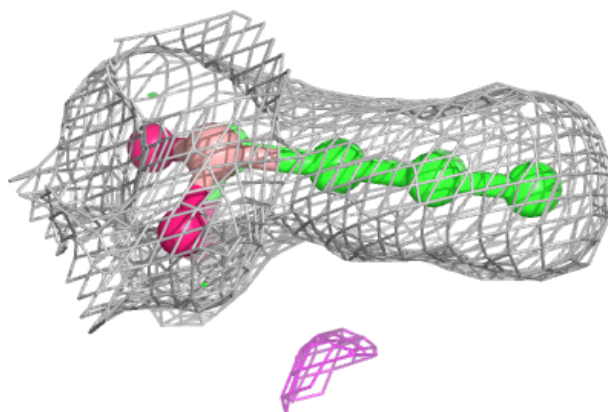
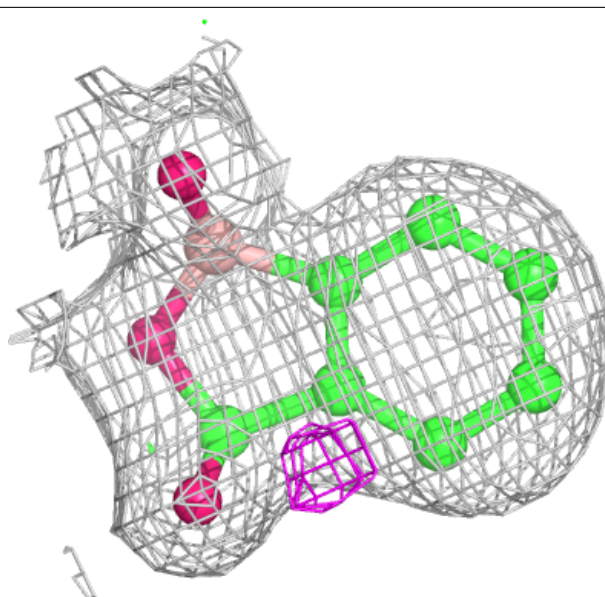
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





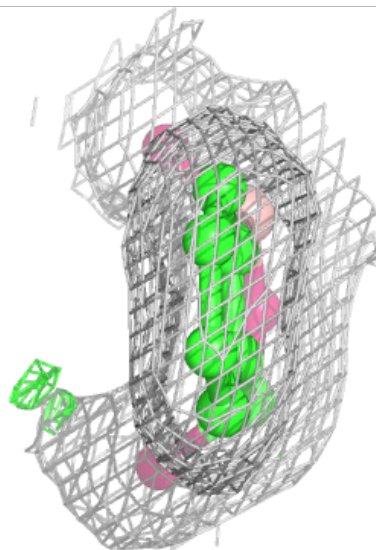
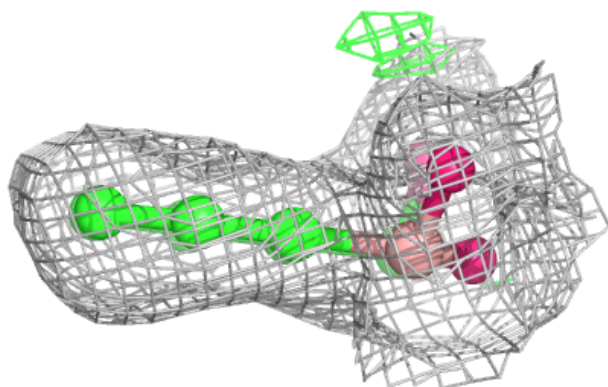
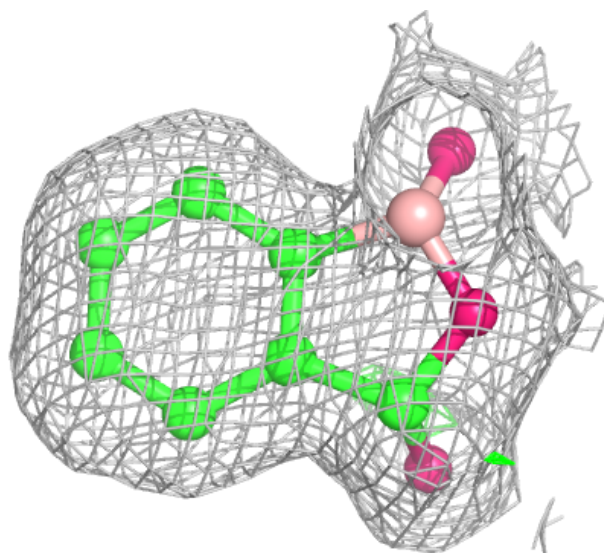
**Electron density around ZMD D 402:**

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



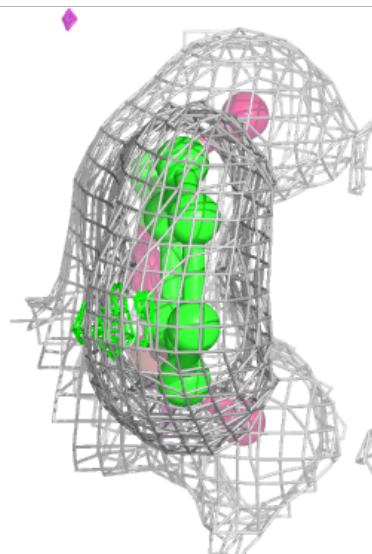
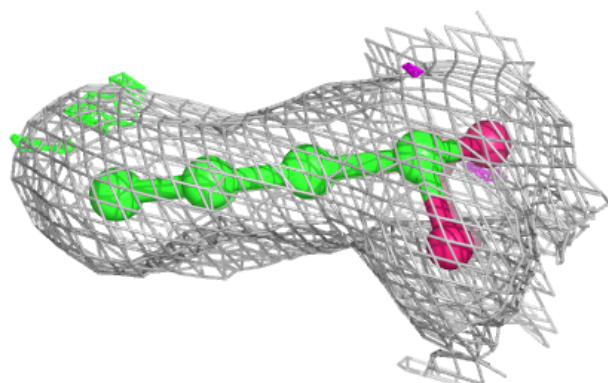
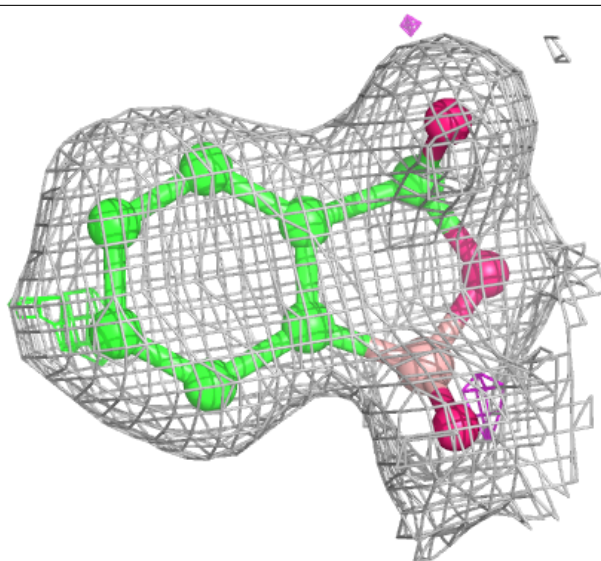
**Electron density around ZMD C 402:**

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



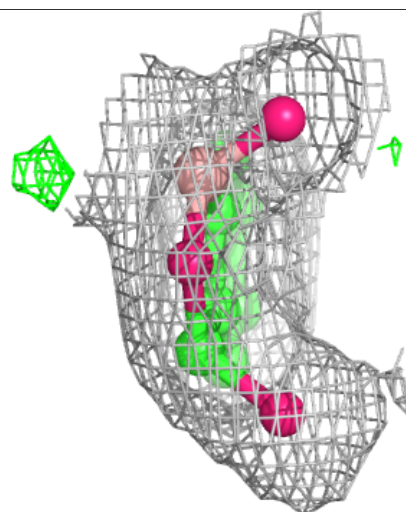
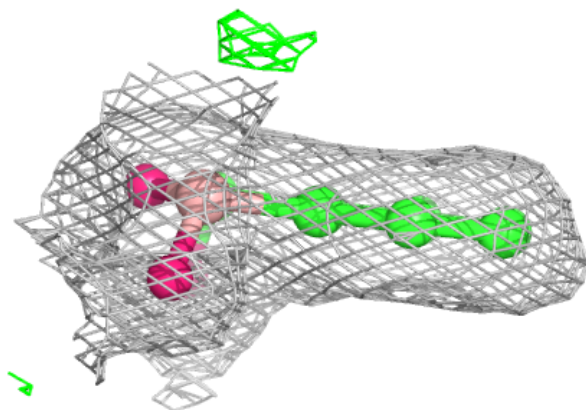
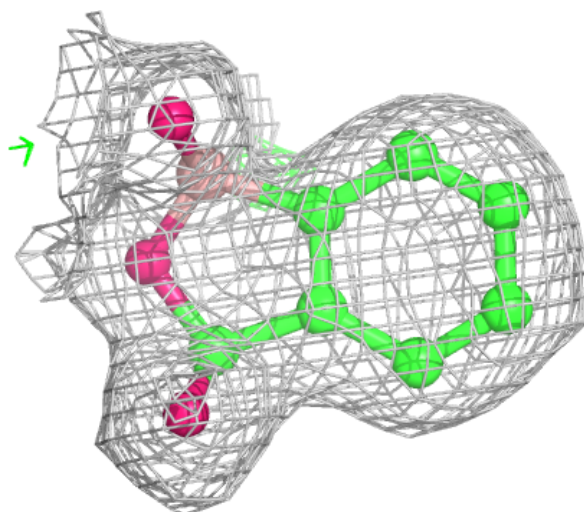
**Electron density around ZMD E 402:**

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



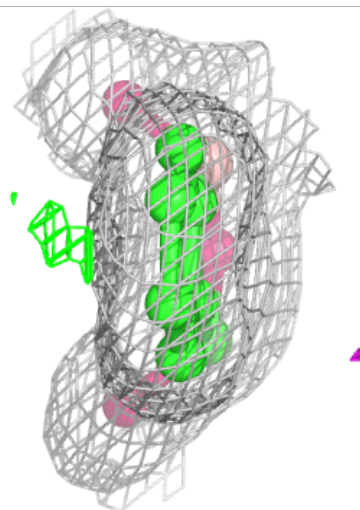
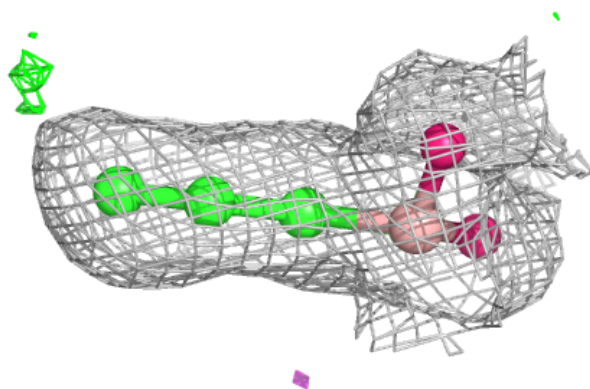
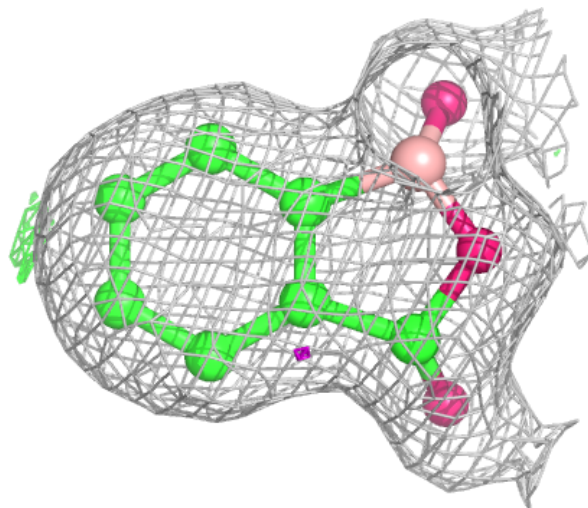
**Electron density around ZMD A 402:**

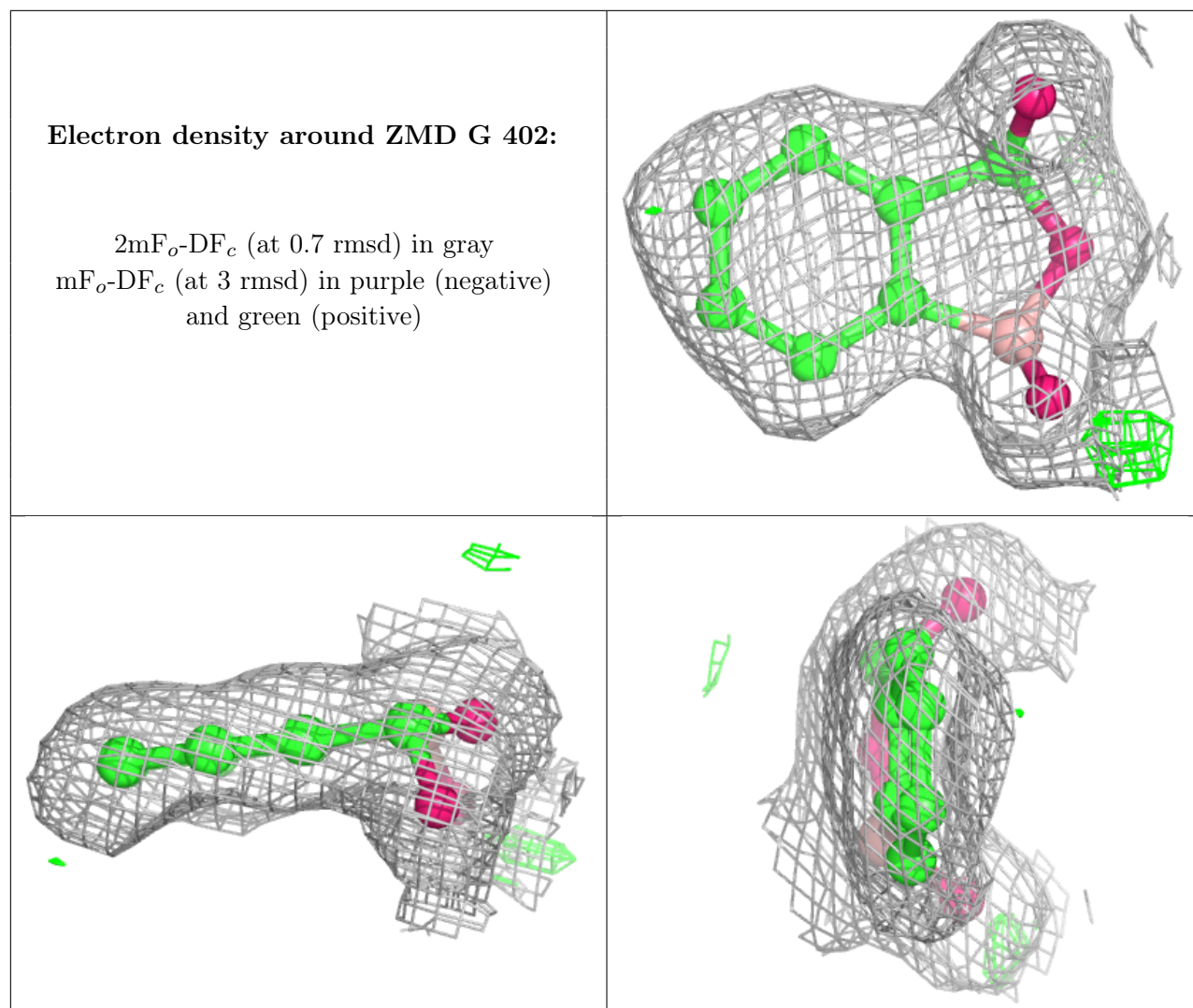
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around ZMD F 402:**

$2mF_o-DF_c$  (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.