



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 24, 2024 – 06:44 PM EDT

PDB ID : 6YRC
Title : Spectroscopically-validated structure of DtpB from *Streptomyces lividans* in the ferric state
Authors : Lucic, M.; Dworkowski, F.S.N.; Worrall, J.A.R.; Hough, M.A.
Deposited on : 2020-04-20
Resolution : 1.99 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

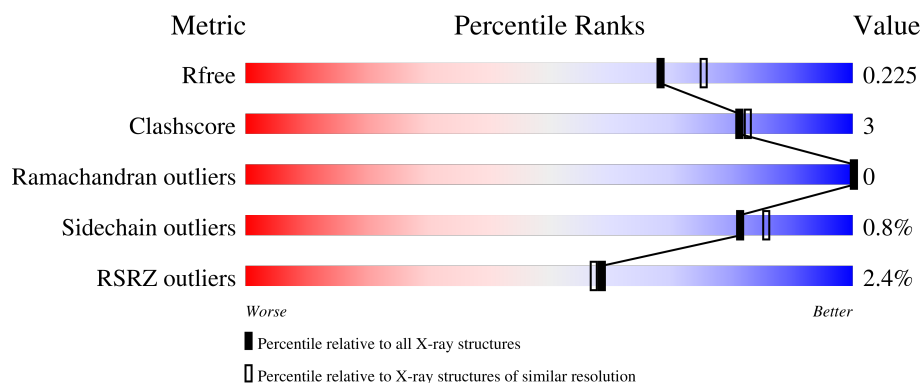
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	316	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>7%</div> <div>.</div> </div> </div>
1	B	316	<div> <div>3%</div> <div> <div></div> <div>93%</div> <div>.</div> <div>.</div> </div> </div>
1	C	316	<div> <div>2%</div> <div> <div></div> <div>92%</div> <div>5%</div> <div>.</div> </div> </div>
1	D	316	<div> <div>2%</div> <div> <div></div> <div>92%</div> <div>5%</div> <div>.</div> </div> </div>
1	E	316	<div> <div>3%</div> <div> <div></div> <div>95%</div> <div>.</div> <div>.</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	316	

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	7PE	A	402	-	-	-	X

2 Entry composition [i](#)

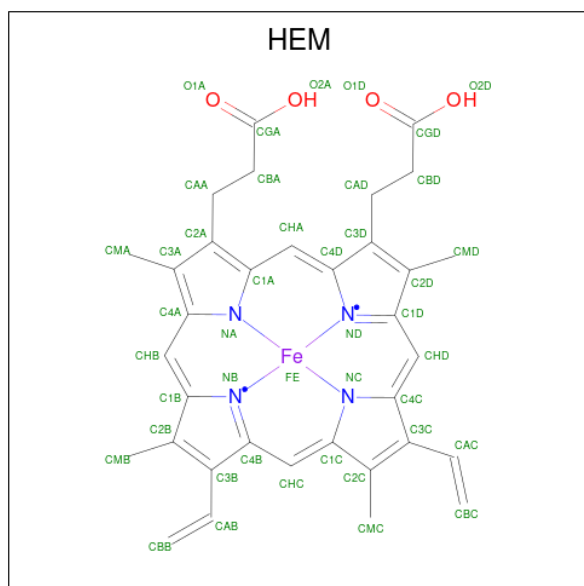
There are 6 unique types of molecules in this entry. The entry contains 15579 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Putative iron-dependent peroxidase.

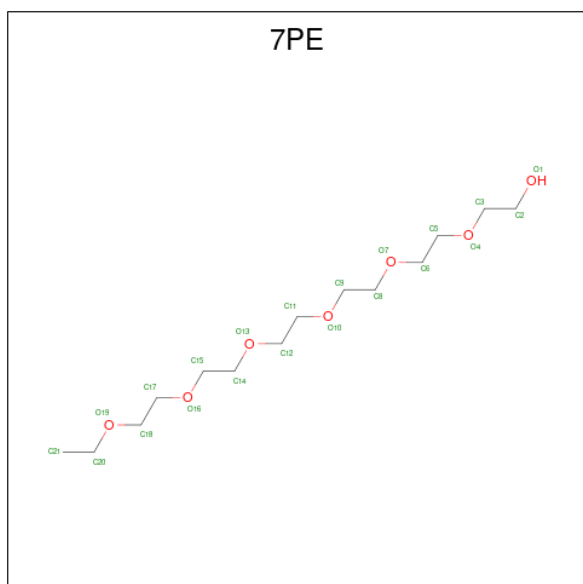
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	5	0
			2372	1488	414	461	9			
1	B	306	Total	C	N	O	S	0	1	0
			2328	1466	404	449	9			
1	C	306	Total	C	N	O	S	0	1	0
			2330	1468	398	454	10			
1	D	306	Total	C	N	O	S	0	2	0
			2348	1474	407	458	9			
1	E	306	Total	C	N	O	S	0	2	0
			2335	1471	401	454	9			
1	F	308	Total	C	N	O	S	0	5	0
			2349	1476	401	463	9			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
2	B	1	Total	C	Fe	N	O	
			43	34	1	4	4	
2	C	1	Total	C	Fe	N	O	
			43	34	1	4	4	
2	D	1	Total	C	Fe	N	O	
			43	34	1	4	4	
2	E	1	Total	C	Fe	N	O	
			43	34	1	4	4	
2	F	1	Total	C	Fe	N	O	
			43	34	1	4	4	

- Molecule 3 is 2-(2-(2-(2-(2-ETHOXYETHOXY)ETHOXY)ETHOXY)ETHOXY)ETHOXY)ETHANOL (three-letter code: 7PE) (formula: C₁₄H₃₀O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O		
			6	4	2		
3	A	1	Total	C	O		
			4	2	2		
3	B	1	Total	C	O		
			5	4	1		
3	B	1	Total	C	O		
			5	3	2		
3	D	1	Total	C	O		
			6	4	2		

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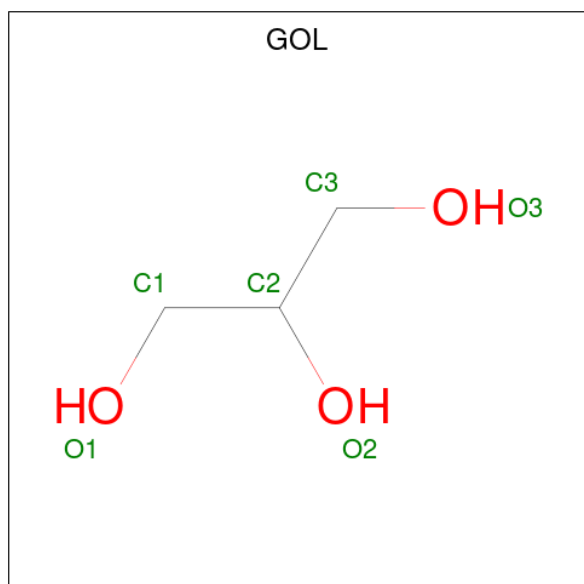
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	C	O	0	0
			9	6	3		
3	F	1	Total	C	O	0	1
			6	4	2		
3	F	1	Total	C	O	0	0
			5	3	2		
3	F	1	Total	C	O	0	0
			7	5	2		
3	F	1	Total	C	O	0	0
			5	3	2		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Mg	0	0
			3	3		
4	C	1	Total	Mg	0	0
			1	1		
4	E	1	Total	Mg	0	0
			1	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	C	1	Total	C	O	0	0
			6	3	3		
5	F	1	Total	C	O	0	0
			6	3	3		
5	F	1	Total	C	O	0	0
			6	3	3		
5	F	1	Total	C	O	0	0
			6	3	3		
5	F	1	Total	C	O	0	0
			6	3	3		

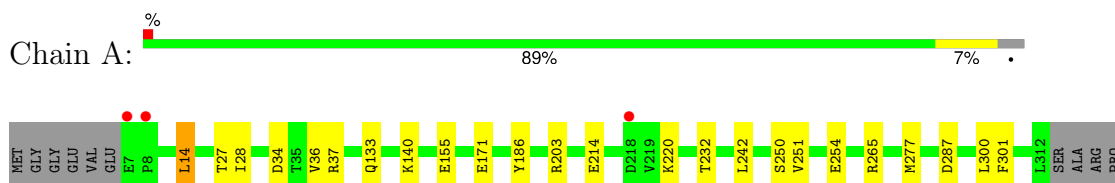
- Molecule 6 is water.

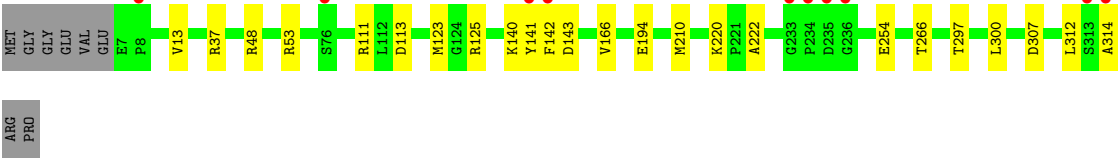
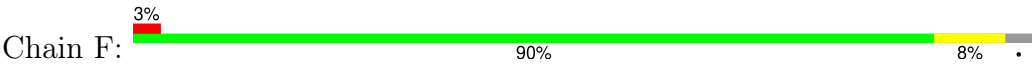
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	216	Total	O	0	0
			216	216		
6	B	147	Total	O	0	0
			147	147		
6	C	205	Total	O	0	0
			205	205		
6	D	198	Total	O	0	0
			198	198		
6	E	189	Total	O	0	0
			189	189		
6	F	187	Total	O	0	0
			187	187		

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: Putative iron-dependent peroxidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.79Å 120.27Å 196.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.05 – 1.99 49.00 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.05-1.99) 99.1 (49.00-1.99)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.176 , 0.218 0.185 , 0.225	Depositor DCC
R_{free} test set	6824 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	30.7	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15579	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, 7PE, HEM, 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	0.76	0/2422	0.84	0/3289
1	B	0.71	0/2378	0.81	0/3231
1	C	0.72	0/2380	0.83	0/3233
1	D	0.75	1/2398 (0.0%)	0.85	0/3256
1	E	0.73	0/2385	0.82	0/3241
1	F	0.74	1/2398 (0.0%)	0.81	0/3259
All	All	0.74	2/14361 (0.0%)	0.83	0/19509

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	F	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	199	GLU	CD-OE2	6.57	1.32	1.25
1	F	254	GLU	CD-OE1	-5.49	1.19	1.25

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	312	LEU	Peptide

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	2372	0	2288	16	0
1	B	2328	0	2251	8	0
1	C	2330	0	2251	11	0
1	D	2348	0	2272	13	0
1	E	2335	0	2249	5	0
1	F	2349	0	2233	20	0
2	A	43	0	30	3	0
2	B	43	0	30	2	0
2	C	43	0	30	0	0
2	D	43	0	30	0	0
2	E	43	0	30	0	0
2	F	43	0	30	0	0
3	A	10	0	10	6	0
3	B	10	0	8	1	0
3	D	6	0	6	1	0
3	F	32	0	28	5	0
4	A	3	0	0	0	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
5	B	12	0	16	0	0
5	C	18	0	24	2	0
5	F	24	0	32	1	0
6	A	216	0	0	4	0
6	B	147	0	0	1	0
6	C	205	0	0	4	0
6	D	198	0	0	5	0
6	E	189	0	0	0	0
6	F	187	0	0	3	0
All	All	15579	0	13848	73	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (73) 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:C:214:GLU:OE1	6:C:501:HOH:O	2.05	0.74
1:E:48:ARG:CZ	3:F:410:7PE:H61	2.19	0.73
1:D:214:GLU:OE1	6:D:501:HOH:O	2.13	0.66
1:D:111[B]:ARG:NH1	1:F:142:PHE:HB3	2.10	0.65
1:F:37:ARG:HD2	1:F:314:ALA:HA	1.78	0.65
1:E:143:ASP:N	1:E:143:ASP:OD1	2.30	0.64
1:F:297:THR:HG22	3:F:405[A]:7PE:H62	1.80	0.62
1:D:214:GLU:OE2	1:D:220:LYS:NZ	2.21	0.62
1:D:90:ASP:OD1	1:D:95:ARG:NH1	2.33	0.61
1:F:111[B]:ARG:HD2	1:F:113:ASP:OD1	2.01	0.61
2:B:401:HEM:HBC2	2:B:401:HEM:HMC2	1.83	0.59
1:A:214:GLU:OE2	1:A:220:LYS:NZ	2.20	0.59
1:A:232:THR:HG23	6:A:691:HOH:O	2.03	0.59
1:C:37:ARG:HD3	6:C:681:HOH:O	2.03	0.58
1:A:34:ASP:OD1	1:A:37:ARG:NH2	2.35	0.58
1:D:111[B]:ARG:NH1	1:F:142:PHE:CB	2.66	0.58
1:C:218:ASP:OD1	6:C:502:HOH:O	2.17	0.58
1:F:194:GLU:OE1	6:F:501:HOH:O	2.16	0.58
2:A:401:HEM:HMC2	2:A:401:HEM:HBC2	1.87	0.55
6:B:609:HOH:O	1:F:210:MET:HE2	2.07	0.55
1:B:53:ARG:HD2	1:C:143:ASP:OD2	2.07	0.54
1:F:140:LYS:HG3	1:F:141:TYR:O	2.07	0.53
1:D:171:GLU:OE2	3:D:402:7PE:C8	2.57	0.52
1:C:210[A]:MET:SD	5:C:405:GOL:H31	2.51	0.51
3:A:402:7PE:H91	1:F:48:ARG:CD	2.41	0.51
1:A:155:GLU:OE1	3:A:402:7PE:O13	2.29	0.50
1:F:140:LYS:CG	1:F:141:TYR:O	2.60	0.49
3:A:402:7PE:H91	1:F:48:ARG:HD3	1.93	0.49
3:A:402:7PE:H111	1:F:48:ARG:CZ	2.42	0.49
1:F:13:VAL:HG11	1:F:166:VAL:CG1	2.42	0.48
2:A:401:HEM:HBC2	2:A:401:HEM:CMC	2.44	0.48
1:A:203[B]:ARG:NH1	6:A:501:HOH:O	2.20	0.47
1:D:185:LYS:NZ	6:D:504:HOH:O	2.41	0.47
1:C:13:VAL:HG11	1:C:166:VAL:CG1	2.45	0.47
1:F:13:VAL:HG11	1:F:166:VAL:HG11	1.96	0.47
1:A:171:GLU:HG2	3:A:403:7PE:H171	1.96	0.47
1:D:113:ASP:HB3	6:F:611:HOH:O	2.13	0.47
1:D:196:LEU:HB3	1:D:200:ALA:HB3	1.97	0.46
2:B:401:HEM:HBC2	2:B:401:HEM:CMC	2.46	0.46
1:C:228:LEU:HD12	1:C:287:ASP:HA	1.99	0.45
1:E:48:ARG:NE	3:F:410:7PE:H61	2.31	0.45
1:A:14:LEU:HD23	1:A:242:LEU:HD22	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:34:ASP:OD1	1:C:37:ARG:NH1	2.50	0.44
5:C:403:GOL:H11	6:C:649:HOH:O	2.16	0.44
1:F:220:LYS:O	3:F:402:7PE:H61	2.16	0.44
1:B:300:LEU:C	1:B:300:LEU:HD23	2.38	0.44
1:A:254:GLU:OE1	6:A:502:HOH:O	2.21	0.44
3:A:403:7PE:C18	6:A:521:HOH:O	2.66	0.44
1:F:300:LEU:C	1:F:300:LEU:HD23	2.39	0.44
1:E:300:LEU:HD23	1:E:301:PHE:N	2.32	0.43
1:D:276:ARG:CZ	6:D:562:HOH:O	2.66	0.43
1:B:55:GLN:HA	1:B:55:GLN:OE1	2.17	0.43
1:B:90:ASP:OD1	1:B:95:ARG:NE	2.51	0.43
1:E:34:ASP:OD1	1:E:37[A]:ARG:NH2	2.40	0.43
1:B:7:GLU:N	1:B:8:PRO:CD	2.82	0.43
1:D:251:VAL:HG12	1:F:123:MET:HG3	2.01	0.43
1:C:90:ASP:OD1	1:C:95:ARG:NH1	2.52	0.42
1:A:27:THR:HG23	1:A:133[A]:GLN:HG3	2.01	0.42
1:A:186:TYR:CG	2:A:401:HEM:HBB1	2.55	0.42
1:A:251:VAL:HG12	1:C:123:MET:HG3	2.01	0.42
1:A:300:LEU:HD23	1:A:301:PHE:N	2.34	0.42
1:B:300:LEU:HD23	1:B:301:PHE:N	2.35	0.42
1:B:202:GLU:OE1	3:B:402:7PE:H141	2.20	0.41
1:F:222:ALA:O	3:F:402:7PE:C11	2.68	0.41
1:D:276:ARG:HD2	6:D:562:HOH:O	2.20	0.41
1:D:203:ARG:NE	6:D:502:HOH:O	2.26	0.41
1:F:125:ARG:HD3	6:F:538:HOH:O	2.21	0.41
1:A:140:LYS:O	1:C:113:ASP:HB3	2.21	0.41
1:F:143[A]:ASP:OD1	5:F:407:GOL:H11	2.20	0.41
1:A:277:MET:HA	1:A:287:ASP:HB2	2.03	0.40
1:B:12:MET:O	1:B:12:MET:HG3	2.21	0.40
1:A:27:THR:HG23	1:A:133[A]:GLN:CG	2.52	0.40
1:A:28:ILE:HG12	1:A:36:VAL:HG21	2.03	0.40

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	309/316 (98%)	300 (97%)	9 (3%)	0	100	100
1	B	305/316 (96%)	299 (98%)	6 (2%)	0	100	100
1	C	305/316 (96%)	298 (98%)	7 (2%)	0	100	100
1	D	306/316 (97%)	300 (98%)	6 (2%)	0	100	100
1	E	306/316 (97%)	299 (98%)	7 (2%)	0	100	100
1	F	311/316 (98%)	307 (99%)	4 (1%)	0	100	100
All	All	1842/1896 (97%)	1803 (98%)	39 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	248/252 (98%)	244 (98%)	4 (2%)	62	67
1	B	241/252 (96%)	238 (99%)	3 (1%)	71	76
1	C	243/252 (96%)	242 (100%)	1 (0%)	91	93
1	D	247/252 (98%)	245 (99%)	2 (1%)	81	86
1	E	242/252 (96%)	241 (100%)	1 (0%)	91	93
1	F	241/252 (96%)	238 (99%)	3 (1%)	71	76
All	All	1462/1512 (97%)	1448 (99%)	14 (1%)	81	81

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	A	250	SER
1	A	265[A]	ARG
1	A	265[B]	ARG

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Mol	Chain	Res	Type
1	B	143	ASP
1	B	210	MET
1	B	283	SER
1	C	127	ARG
1	D	307[A]	ASP
1	D	307[B]	ASP
1	E	143	ASP
1	F	53	ARG
1	F	266	THR
1	F	307	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 30 ligands modelled in this entry, 5 are monoatomic - leaving 25 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	7PE	B	404	-	4,4,20	0.33	0	3,3,19	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GOL	F	406	-	5,5,5	0.10	0	5,5,5	0.24	0
5	GOL	C	402	-	5,5,5	0.13	0	5,5,5	0.20	0
5	GOL	B	403	-	5,5,5	0.11	0	5,5,5	0.26	0
2	HEM	F	401	1	42,50,50	1.32	6 (14%)	46,82,82	1.82	13 (28%)
5	GOL	B	405	-	5,5,5	0.14	0	5,5,5	0.34	0
5	GOL	F	404	-	5,5,5	0.09	0	5,5,5	0.38	0
3	7PE	F	405[A]	-	5,5,20	0.08	0	4,4,19	0.38	0
3	7PE	A	402	-	5,5,20	0.35	0	4,4,19	0.21	0
3	7PE	F	402	-	8,8,20	0.34	0	7,7,19	0.16	0
5	GOL	F	407	-	5,5,5	0.15	0	5,5,5	0.37	0
3	7PE	B	402	-	4,4,20	0.31	0	3,3,19	0.20	0
3	7PE	D	402	-	5,5,20	0.29	0	4,4,19	0.28	0
3	7PE	A	403	-	3,3,20	0.34	0	2,2,19	0.11	0
3	7PE	F	408	-	4,4,20	0.36	0	3,3,19	0.24	0
2	HEM	E	401	1	42,50,50	1.36	7 (16%)	46,82,82	2.03	17 (36%)
5	GOL	C	403	-	5,5,5	0.12	0	5,5,5	0.11	0
2	HEM	D	401	1	42,50,50	1.27	7 (16%)	46,82,82	1.77	10 (21%)
5	GOL	C	405	-	5,5,5	0.16	0	5,5,5	0.39	0
3	7PE	F	409	-	6,6,20	0.41	0	5,5,19	0.17	0
3	7PE	F	410	-	4,4,20	0.26	0	3,3,19	0.17	0
2	HEM	A	401	1	42,50,50	1.37	6 (14%)	46,82,82	1.94	15 (32%)
2	HEM	C	401	1	42,50,50	1.18	5 (11%)	46,82,82	1.72	13 (28%)
5	GOL	F	403	-	5,5,5	0.15	0	5,5,5	0.39	0
2	HEM	B	401	1	42,50,50	1.28	6 (14%)	46,82,82	1.82	11 (23%)

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	7PE	B	404	-	-	2/2/2/18	-
5	GOL	F	406	-	-	0/4/4/4	-
5	GOL	C	402	-	-	2/4/4/4	-
5	GOL	B	403	-	-	2/4/4/4	-
2	HEM	F	401	1	-	4/12/54/54	-
5	GOL	B	405	-	-	3/4/4/4	-
5	GOL	F	404	-	-	2/4/4/4	-
3	7PE	F	405[A]	-	-	1/3/3/18	-
3	7PE	A	402	-	-	1/3/3/18	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	7PE	F	402	-	-	3/6/6/18	-
5	GOL	F	407	-	-	0/4/4/4	-
3	7PE	B	402	-	-	1/2/2/18	-
3	7PE	D	402	-	-	3/3/3/18	-
3	7PE	A	403	-	-	0/1/1/18	-
3	7PE	F	408	-	-	2/2/2/18	-
2	HEM	E	401	1	-	4/12/54/54	-
5	GOL	C	403	-	-	2/4/4/4	-
2	HEM	D	401	1	-	4/12/54/54	-
5	GOL	C	405	-	-	2/4/4/4	-
3	7PE	F	409	-	-	1/4/4/18	-
3	7PE	F	410	-	-	0/2/2/18	-
2	HEM	A	401	1	-	4/12/54/54	-
2	HEM	C	401	1	-	5/12/54/54	-
5	GOL	F	403	-	-	0/4/4/4	-
2	HEM	B	401	1	-	4/12/54/54	-

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	HEM	C1B-NB	-3.90	1.33	1.40
2	F	401	HEM	C1B-NB	-3.65	1.33	1.40
2	E	401	HEM	C4B-NB	-3.37	1.32	1.38
2	E	401	HEM	C4D-ND	-3.18	1.34	1.40
2	D	401	HEM	C4B-NB	-3.13	1.32	1.38
2	A	401	HEM	C4D-C3D	2.98	1.50	1.45
2	A	401	HEM	C3C-C4C	2.91	1.45	1.41
2	A	401	HEM	C4D-ND	-2.89	1.35	1.40
2	A	401	HEM	C1B-NB	-2.88	1.35	1.40
2	C	401	HEM	C1B-NB	-2.77	1.35	1.40
2	B	401	HEM	CHB-C1B	2.62	1.41	1.34
2	A	401	HEM	C3B-C4B	2.61	1.50	1.44
2	F	401	HEM	CHB-C1B	2.59	1.40	1.34
2	E	401	HEM	O1A-CGA	2.38	1.29	1.22
2	D	401	HEM	C3B-C4B	2.33	1.49	1.44
2	E	401	HEM	C3C-C4C	2.32	1.44	1.41
2	F	401	HEM	CHA-C4D	2.32	1.40	1.34
2	D	401	HEM	C3D-C2D	-2.32	1.31	1.36
2	A	401	HEM	C4B-NB	-2.30	1.34	1.38
2	D	401	HEM	C4D-ND	-2.27	1.36	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	401	HEM	CHB-C1B	2.26	1.40	1.34
2	B	401	HEM	C4D-ND	-2.20	1.36	1.40
2	E	401	HEM	C3C-C2C	-2.19	1.37	1.40
2	D	401	HEM	C1B-NB	-2.18	1.36	1.40
2	C	401	HEM	C4B-NB	-2.17	1.34	1.38
2	C	401	HEM	CHA-C4D	2.17	1.39	1.34
2	E	401	HEM	C1B-NB	-2.16	1.36	1.40
2	F	401	HEM	C1D-ND	-2.15	1.34	1.38
2	D	401	HEM	O2A-CGA	-2.15	1.23	1.30
2	C	401	HEM	O2A-CGA	-2.13	1.23	1.30
2	D	401	HEM	C3C-C4C	2.11	1.44	1.41
2	C	401	HEM	C3B-C4B	2.07	1.48	1.44
2	F	401	HEM	C4D-C3D	2.06	1.48	1.45
2	B	401	HEM	C4B-NB	-2.06	1.34	1.38
2	B	401	HEM	C3C-C2C	-2.04	1.37	1.40
2	B	401	HEM	C4D-C3D	2.03	1.48	1.45
2	F	401	HEM	C4B-NB	-2.00	1.34	1.38

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	HEM	CHC-C4B-NB	5.21	130.04	124.44
2	B	401	HEM	CHC-C4B-NB	4.85	129.66	124.44
2	A	401	HEM	C3B-C4B-NB	-4.78	106.04	109.47
2	A	401	HEM	C1B-NB-C4B	4.78	110.86	105.21
2	E	401	HEM	CHC-C4B-NB	4.60	129.39	124.44
2	E	401	HEM	C1B-NB-C4B	4.58	110.63	105.21
2	C	401	HEM	CHC-C4B-NB	4.52	129.30	124.44
2	D	401	HEM	CHC-C4B-NB	4.33	129.10	124.44
2	B	401	HEM	C1B-NB-C4B	4.22	110.20	105.21
2	B	401	HEM	CHD-C1D-ND	3.85	128.57	124.44
2	E	401	HEM	C3B-C4B-NB	-3.71	106.80	109.47
2	F	401	HEM	CHD-C1D-ND	3.47	128.17	124.44
2	F	401	HEM	C1B-NB-C4B	3.46	109.31	105.21
2	D	401	HEM	CHD-C1D-ND	3.43	128.12	124.44
2	A	401	HEM	CHD-C1D-C2D	-3.40	119.67	125.03
2	A	401	HEM	CHD-C1D-ND	3.32	128.00	124.44
2	D	401	HEM	C1B-NB-C4B	3.31	109.13	105.21
2	E	401	HEM	CHA-C4D-ND	3.27	128.42	124.37
2	E	401	HEM	CMD-C2D-C1D	3.26	130.13	125.03
2	E	401	HEM	CMA-C3A-C4A	-3.23	123.72	128.46
2	C	401	HEM	CHD-C1D-ND	3.19	127.86	124.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	HEM	CHC-C4B-NB	3.13	127.80	124.44
2	F	401	HEM	CHD-C1D-C2D	-3.11	120.12	125.03
2	B	401	HEM	CHD-C1D-C2D	-3.10	120.14	125.03
2	C	401	HEM	CMA-C3A-C4A	-3.03	124.02	128.46
2	D	401	HEM	CMA-C3A-C4A	-3.02	124.03	128.46
2	E	401	HEM	CMC-C2C-C3C	3.00	130.67	124.68
2	E	401	HEM	CHB-C1B-NB	2.94	128.02	124.37
2	D	401	HEM	CHD-C1D-C2D	-2.91	120.43	125.03
2	C	401	HEM	C1B-NB-C4B	2.91	108.65	105.21
2	D	401	HEM	CHA-C4D-C3D	-2.90	119.89	125.23
2	E	401	HEM	CHD-C1D-C2D	-2.86	120.51	125.03
2	B	401	HEM	C4B-CHC-C1C	2.81	126.27	122.56
2	D	401	HEM	CHA-C4D-ND	2.77	127.81	124.37
2	C	401	HEM	CHD-C1D-C2D	-2.75	120.69	125.03
2	C	401	HEM	CMC-C2C-C3C	2.71	130.10	124.68
2	B	401	HEM	CAD-C3D-C4D	2.69	129.38	124.70
2	B	401	HEM	CHA-C4D-ND	2.61	127.61	124.37
2	F	401	HEM	O2D-CGD-CBD	2.59	122.18	114.00
2	E	401	HEM	CBA-CAA-C2A	2.57	116.85	112.54
2	C	401	HEM	CHA-C4D-ND	2.56	127.55	124.37
2	F	401	HEM	CHA-C4D-C3D	-2.56	120.50	125.23
2	B	401	HEM	CBA-CAA-C2A	2.55	116.83	112.54
2	A	401	HEM	CMB-C2B-C1B	2.55	129.02	125.03
2	A	401	HEM	CHB-C1B-NB	2.49	127.46	124.37
2	E	401	HEM	CHA-C4D-C3D	-2.49	120.64	125.23
2	C	401	HEM	CMA-C3A-C2A	2.49	129.63	124.94
2	A	401	HEM	O2D-CGD-CBD	2.48	121.84	114.00
2	D	401	HEM	CAD-C3D-C4D	2.42	128.91	124.70
2	E	401	HEM	CHD-C1D-ND	2.40	127.02	124.44
2	A	401	HEM	CAD-C3D-C4D	2.40	128.88	124.70
2	E	401	HEM	C2C-C3C-C4C	-2.39	105.23	106.90
2	F	401	HEM	CHB-C1B-NB	2.37	127.31	124.37
2	E	401	HEM	O2D-CGD-CBD	2.35	121.42	114.00
2	A	401	HEM	CHA-C4D-ND	2.32	127.25	124.37
2	C	401	HEM	CHA-C4D-C3D	-2.32	120.95	125.23
2	D	401	HEM	CMD-C2D-C1D	2.31	128.65	125.03
2	A	401	HEM	CHA-C4D-C3D	-2.28	121.02	125.23
2	D	401	HEM	CMC-C2C-C3C	2.27	129.22	124.68
2	C	401	HEM	C4B-CHC-C1C	2.27	125.55	122.56
2	B	401	HEM	CHA-C4D-C3D	-2.27	121.05	125.23
2	A	401	HEM	C4B-CHC-C1C	2.25	125.53	122.56
2	F	401	HEM	CAD-C3D-C4D	2.23	128.59	124.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	HEM	CHA-C4D-ND	2.22	127.12	124.37
2	C	401	HEM	C3B-C4B-NB	-2.22	107.88	109.47
2	E	401	HEM	C2D-C1D-ND	2.22	112.46	109.90
2	F	401	HEM	O1D-CGD-CBD	-2.21	116.09	123.09
2	C	401	HEM	CMD-C2D-C1D	2.19	128.46	125.03
2	A	401	HEM	CAB-C3B-C2B	-2.19	121.30	128.43
2	E	401	HEM	CAD-C3D-C4D	2.19	128.51	124.70
2	B	401	HEM	CHB-C1B-NB	2.15	127.04	124.37
2	F	401	HEM	CBA-CAA-C2A	2.13	116.12	112.54
2	A	401	HEM	C2D-C1D-ND	2.10	112.33	109.90
2	E	401	HEM	CMA-C3A-C2A	2.08	128.87	124.94
2	A	401	HEM	CMD-C2D-C1D	2.08	128.28	125.03
2	B	401	HEM	CMD-C2D-C1D	2.06	128.25	125.03
2	C	401	HEM	CAD-C3D-C4D	2.06	128.28	124.70
2	F	401	HEM	O2A-CGA-CBA	2.05	120.49	114.00
2	F	401	HEM	CMA-C3A-C4A	-2.04	125.46	128.46

There are no chirality outliers.

All (52) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	402	GOL	O1-C1-C2-O2
5	C	402	GOL	O1-C1-C2-C3
5	C	403	GOL	C1-C2-C3-O3
5	C	403	GOL	O2-C2-C3-O3
5	C	405	GOL	O1-C1-C2-C3
3	F	402	7PE	O7-C8-C9-O10
5	B	405	GOL	O1-C1-C2-O2
3	F	409	7PE	O13-C14-C15-O16
3	F	405[A]	7PE	O4-C5-C6-O7
5	B	403	GOL	C1-C2-C3-O3
5	B	405	GOL	O1-C1-C2-C3
5	F	404	GOL	C1-C2-C3-O3
3	F	402	7PE	O4-C5-C6-O7
5	F	404	GOL	O2-C2-C3-O3
3	B	404	7PE	O10-C11-C12-O13
3	F	408	7PE	O13-C14-C15-O16
3	A	402	7PE	C8-C9-O10-C11
3	B	402	7PE	C15-C14-O13-C12
5	B	403	GOL	O2-C2-C3-O3
5	B	405	GOL	O2-C2-C3-O3
3	B	404	7PE	C12-C11-O10-C9

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Mol	Chain	Res	Type	Atoms
3	F	402	7PE	C8-C9-O10-C11
3	D	402	7PE	O10-C11-C12-O13
3	F	408	7PE	C14-C15-O16-C17
2	F	401	HEM	CAD-CBD-CGD-O2D
5	C	405	GOL	O1-C1-C2-O2
2	C	401	HEM	CAD-CBD-CGD-O2D
2	A	401	HEM	CAD-CBD-CGD-O1D
2	A	401	HEM	CAD-CBD-CGD-O2D
2	C	401	HEM	CAD-CBD-CGD-O1D
2	C	401	HEM	CAA-CBA-CGA-O2A
2	F	401	HEM	CAD-CBD-CGD-O1D
2	A	401	HEM	CAA-CBA-CGA-O1A
2	B	401	HEM	CAA-CBA-CGA-O2A
2	F	401	HEM	CAA-CBA-CGA-O1A
2	A	401	HEM	CAA-CBA-CGA-O2A
2	D	401	HEM	CAA-CBA-CGA-O1A
2	F	401	HEM	CAA-CBA-CGA-O2A
2	E	401	HEM	CAA-CBA-CGA-O1A
2	C	401	HEM	CAA-CBA-CGA-O1A
2	E	401	HEM	CAA-CBA-CGA-O2A
3	D	402	7PE	C12-C11-O10-C9
2	B	401	HEM	CAD-CBD-CGD-O2D
2	D	401	HEM	CAA-CBA-CGA-O2A
2	B	401	HEM	CAD-CBD-CGD-O1D
2	D	401	HEM	CAD-CBD-CGD-O2D
2	E	401	HEM	CAD-CBD-CGD-O2D
2	C	401	HEM	C4B-C3B-CAB-CBB
2	D	401	HEM	CAD-CBD-CGD-O1D
2	E	401	HEM	CAD-CBD-CGD-O1D
2	B	401	HEM	CAA-CBA-CGA-O1A
3	D	402	7PE	C8-C9-O10-C11

There are no ring outliers.

12 monomers are involved in 21 short contacts:

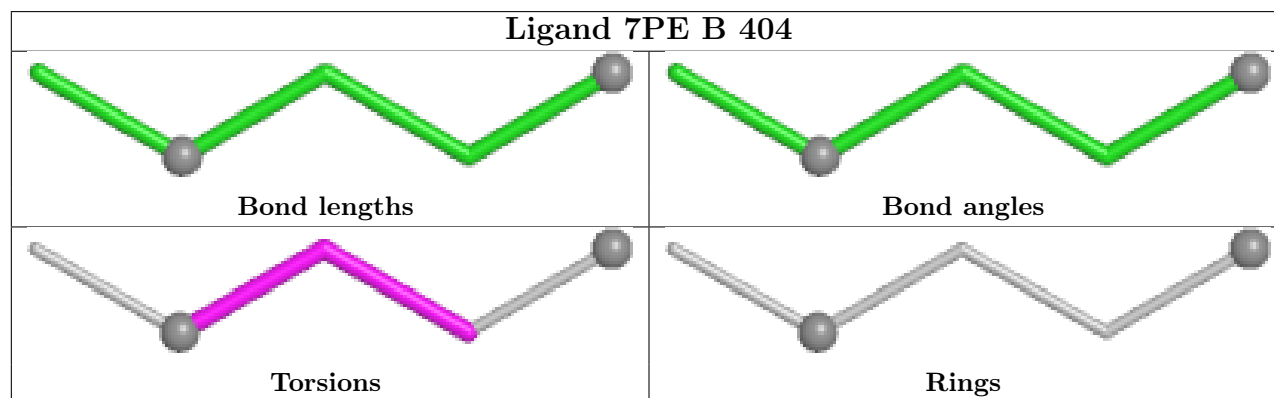
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	405[A]	7PE	1	0
3	A	402	7PE	4	0
3	F	402	7PE	2	0
5	F	407	GOL	1	0
3	B	402	7PE	1	0
3	D	402	7PE	1	0

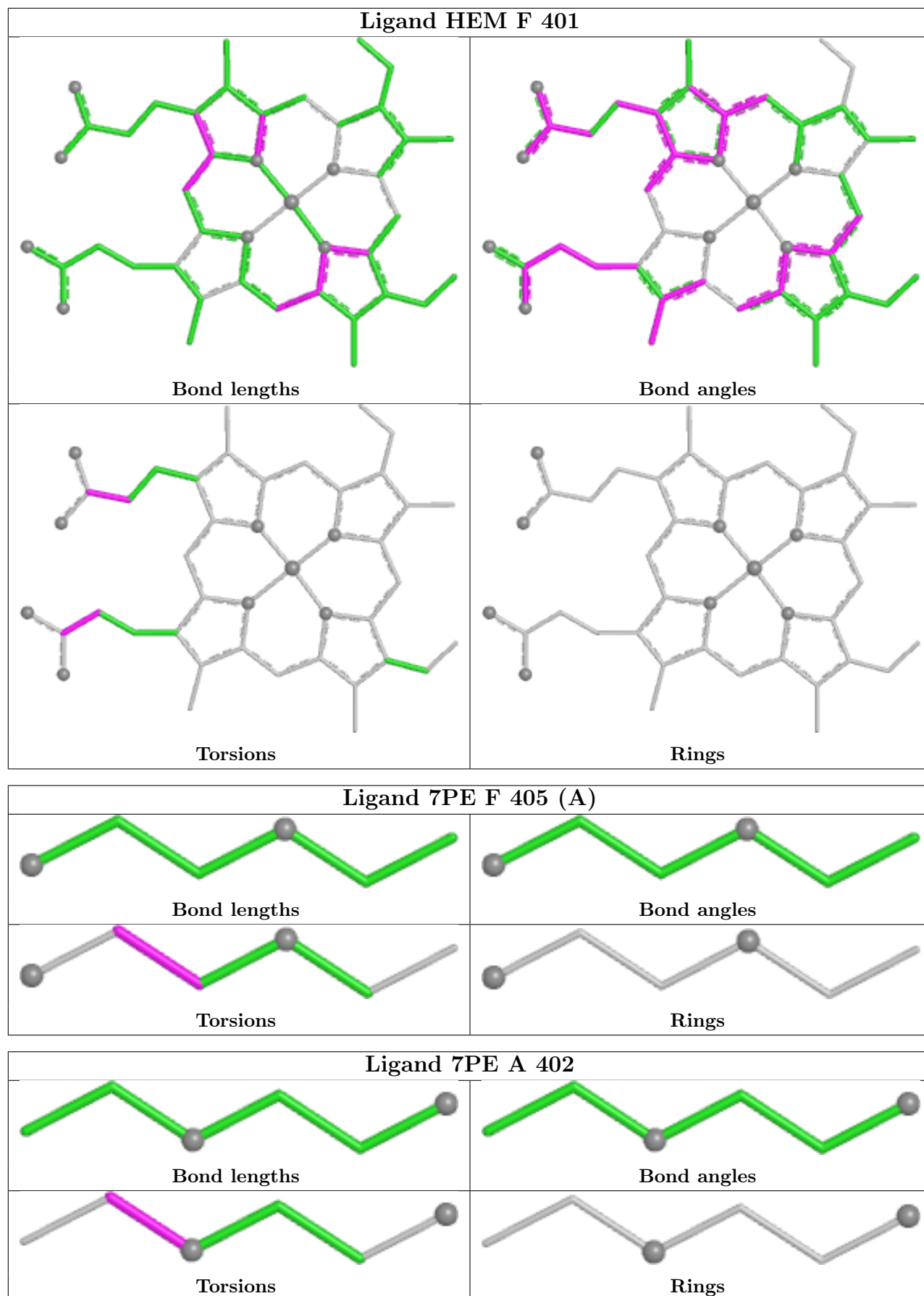
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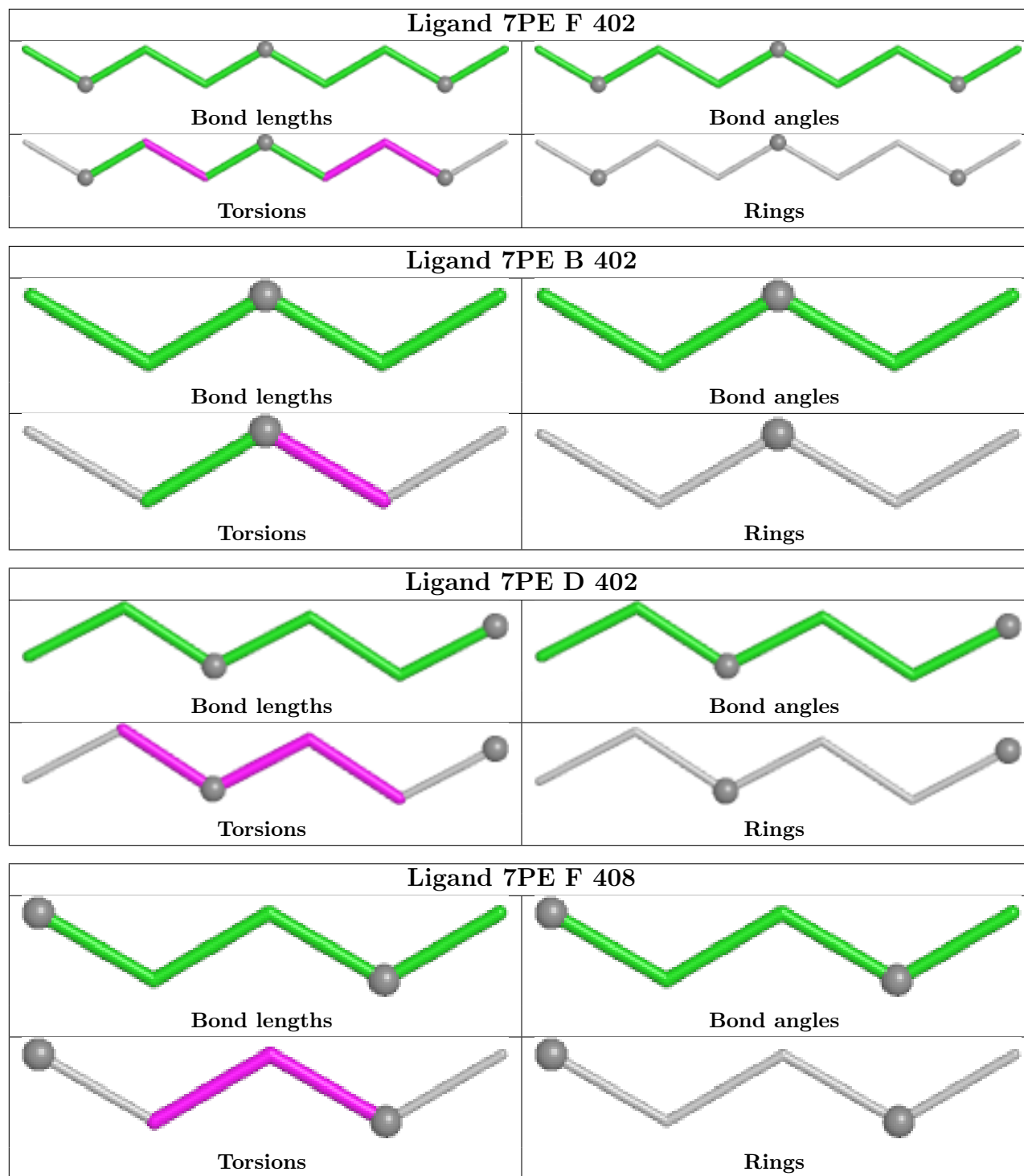
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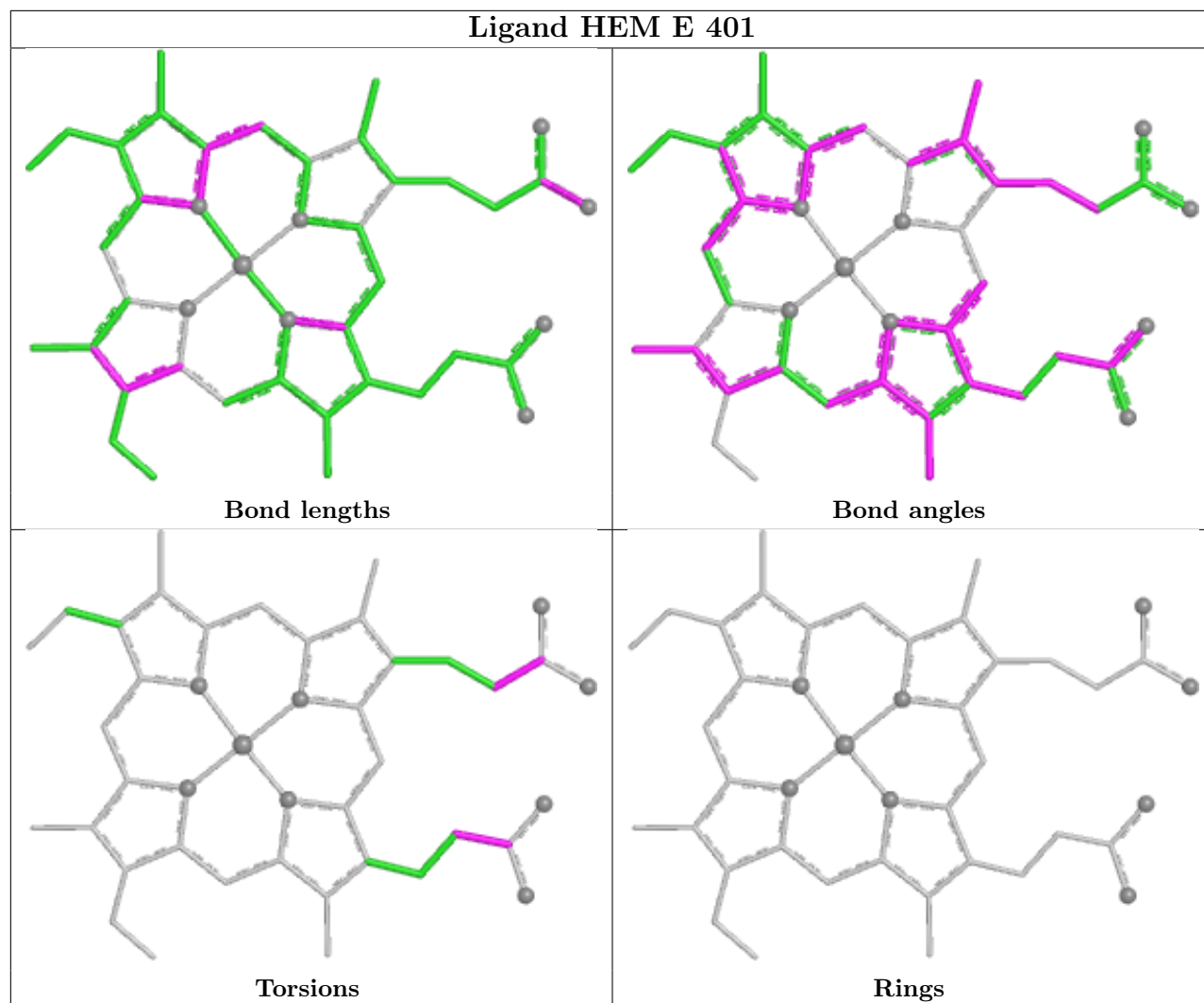
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	7PE	2	0
5	C	403	GOL	1	0
5	C	405	GOL	1	0
3	F	410	7PE	2	0
2	A	401	HEM	3	0
2	B	401	HEM	2	0

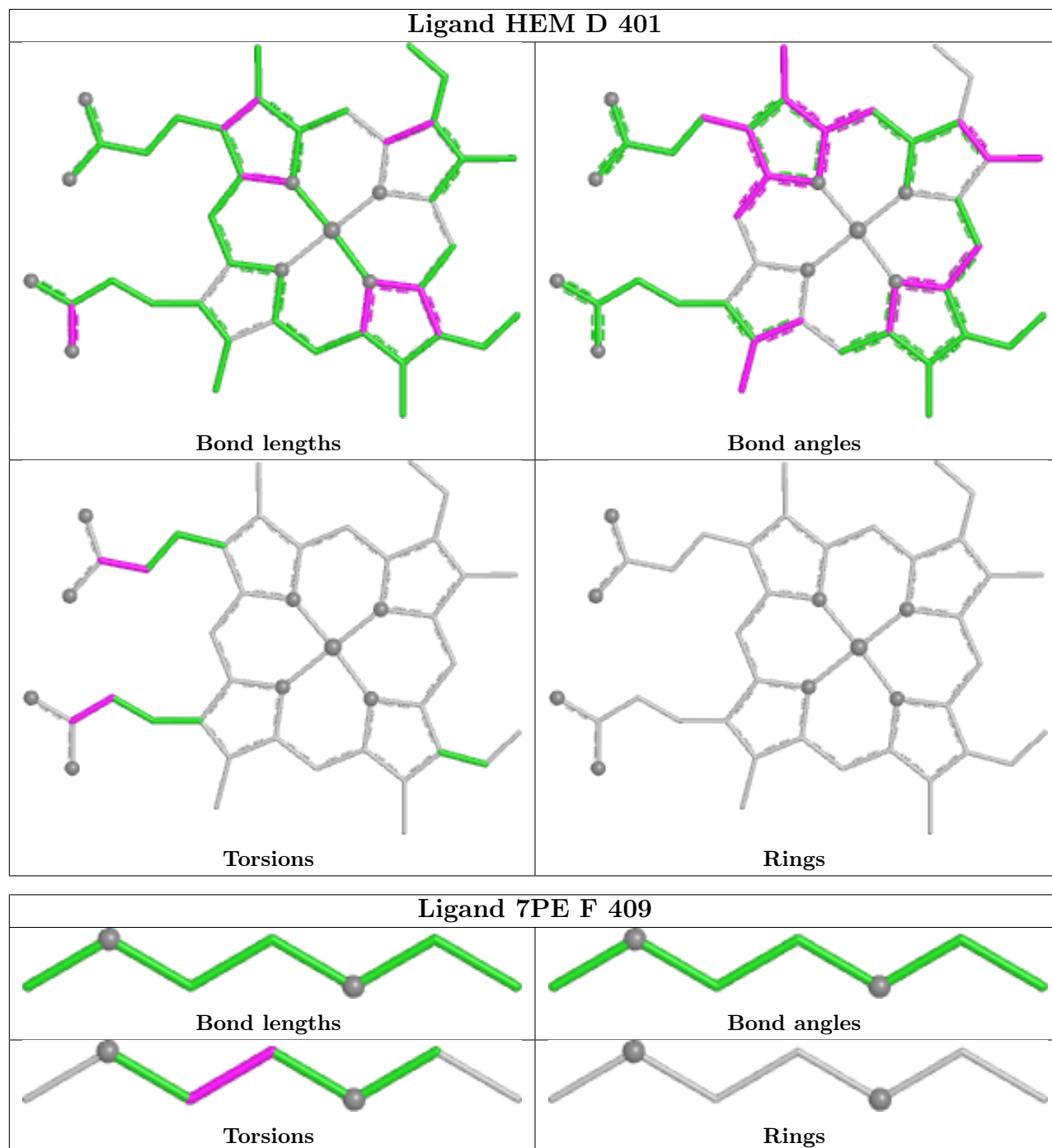
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

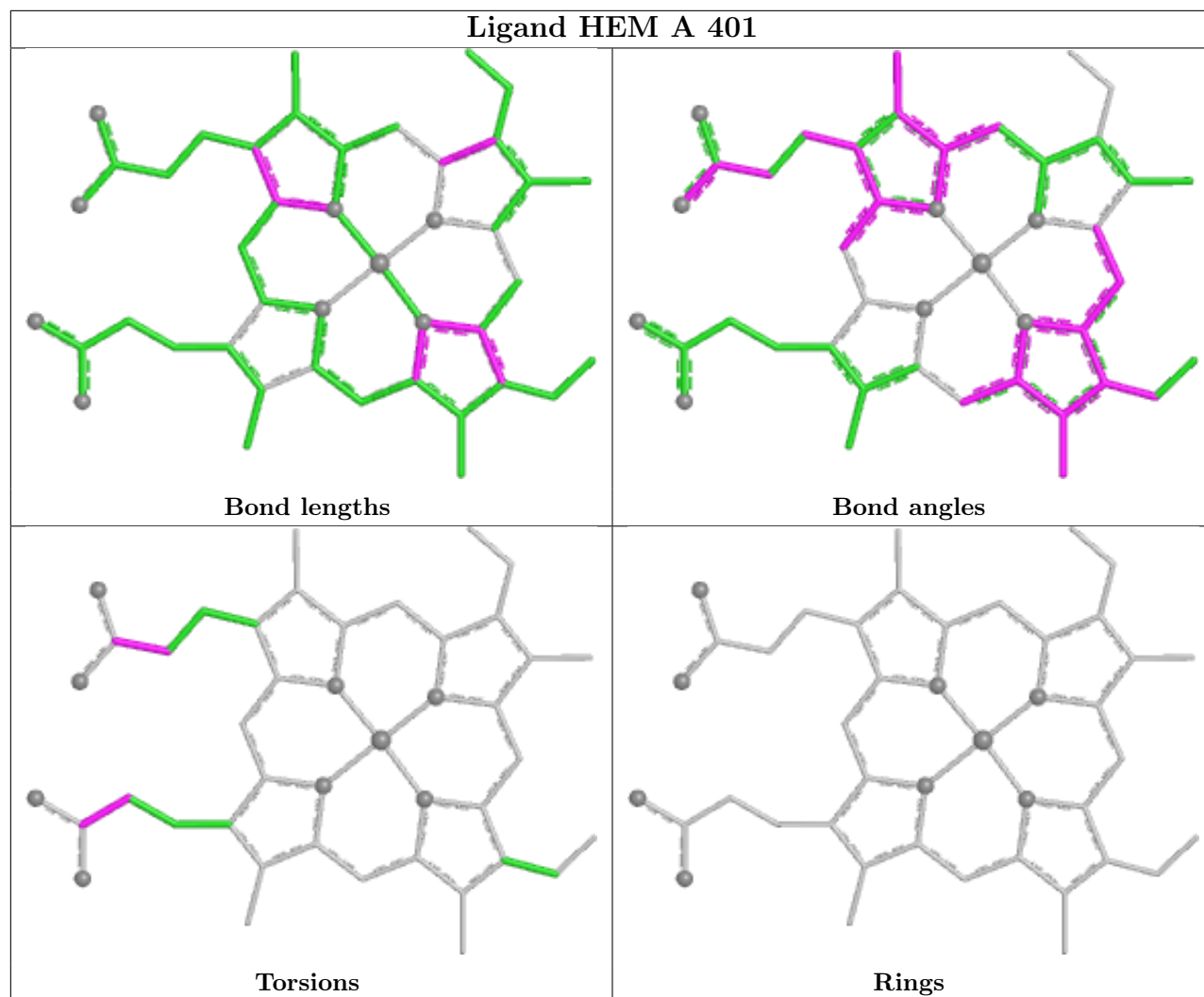


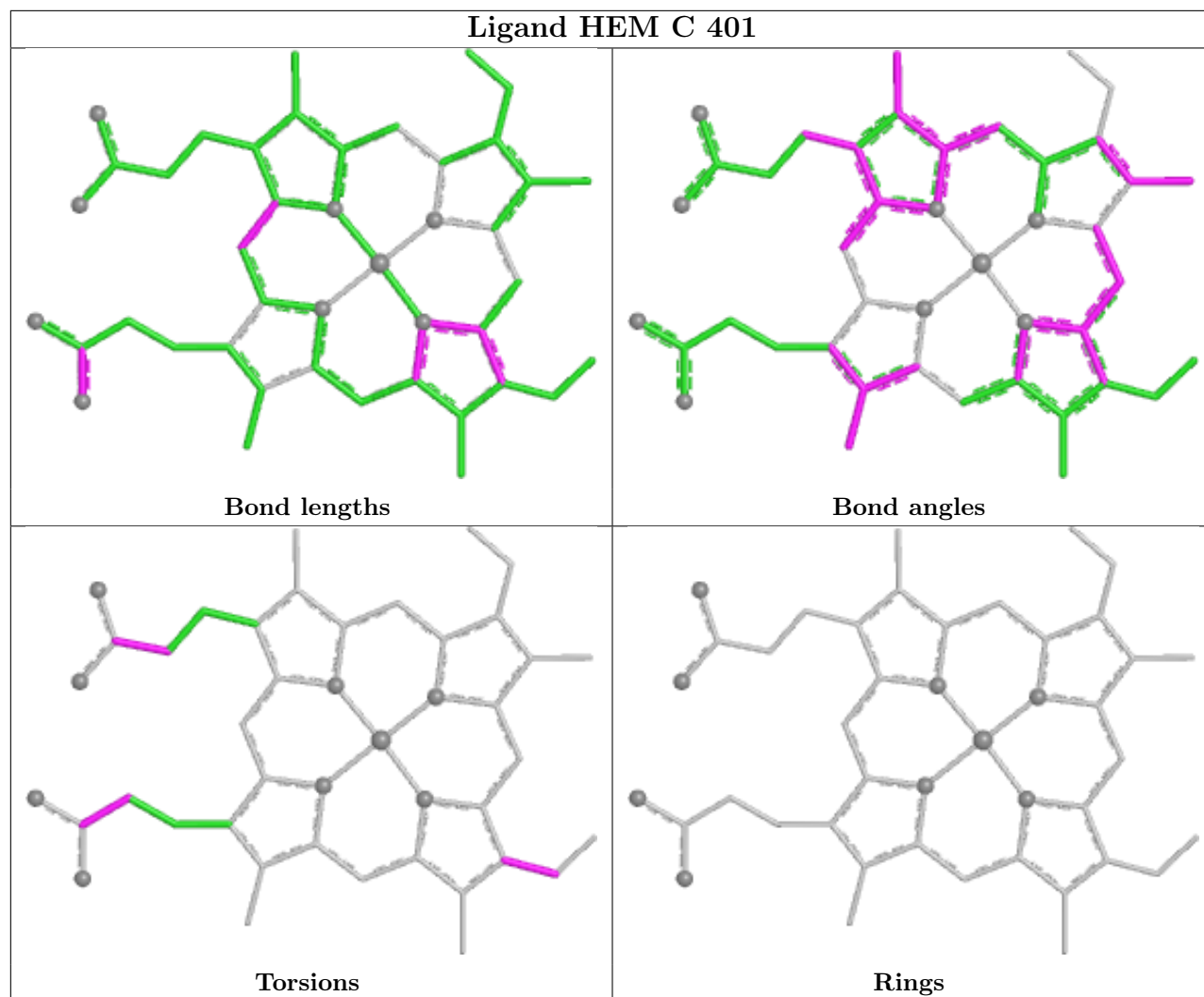


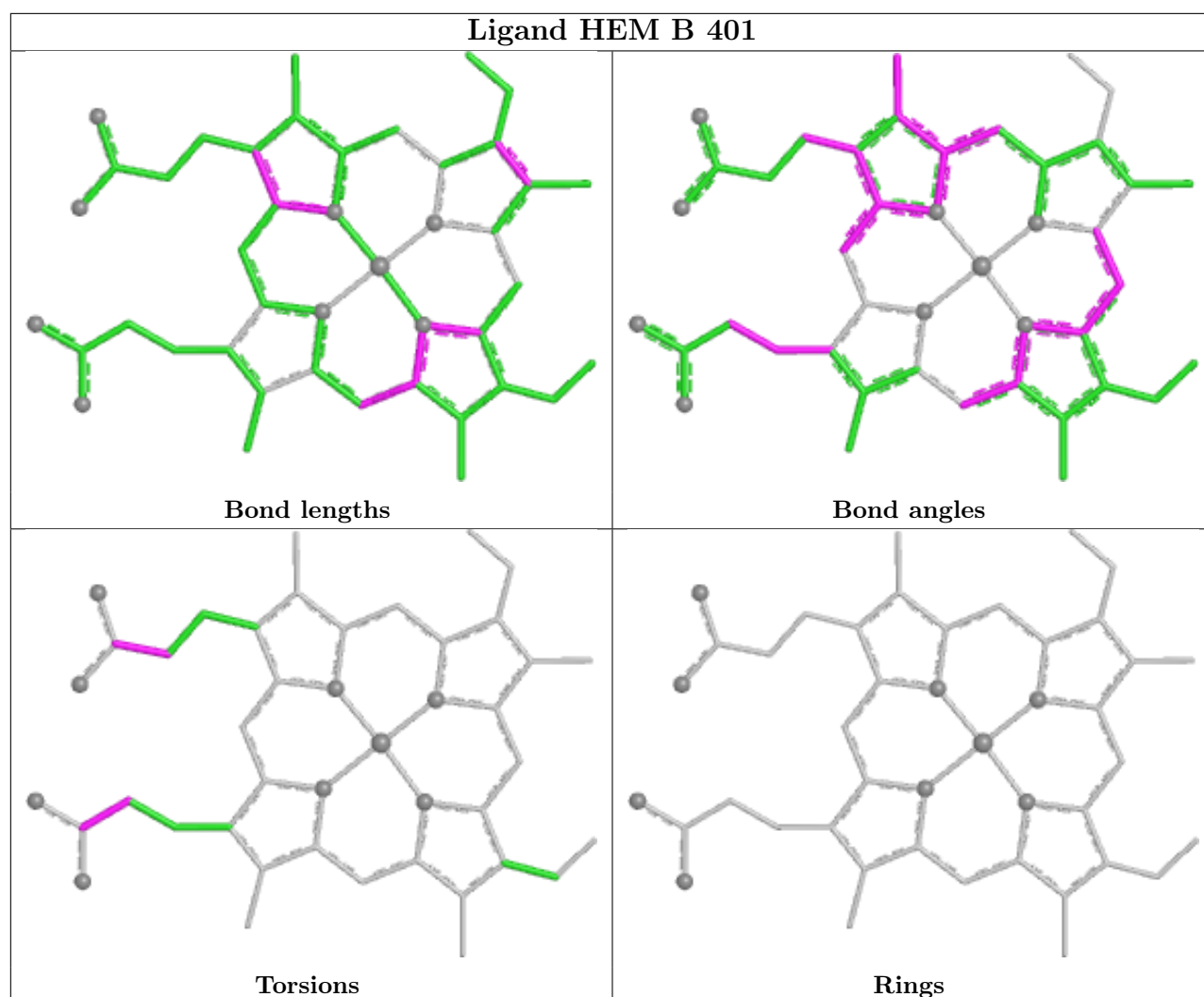












5.7 Other polymers [i](#)

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	306/316 (96%)	-0.10	3 (0%) 82 81	22, 28, 45, 69	0
1	B	306/316 (96%)	0.20	11 (3%) 42 42	24, 33, 53, 80	0
1	C	306/316 (96%)	0.00	6 (1%) 65 63	22, 29, 46, 70	0
1	D	306/316 (96%)	0.02	5 (1%) 72 70	23, 29, 48, 67	0
1	E	306/316 (96%)	0.16	9 (2%) 51 50	24, 30, 49, 91	0
1	F	308/316 (97%)	0.02	10 (3%) 47 46	23, 29, 48, 73	0
All	All	1838/1896 (96%)	0.05	44 (2%) 59 57	22, 30, 49, 91	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	234	PRO	6.6
1	C	78	ALA	5.0
1	E	235	ASP	3.8
1	F	142	PHE	3.5
1	F	314	ALA	3.5
1	D	312	LEU	3.4
1	F	76[A]	SER	3.3
1	C	77	GLY	3.1
1	F	313	SER	3.1
1	F	235	ASP	3.1
1	D	141	TYR	3.0
1	E	142	PHE	3.0
1	B	31	GLY	3.0
1	E	161	ALA	3.0
1	A	8	PRO	3.0
1	C	8	PRO	3.0
1	D	170	ALA	3.0
1	E	160	ALA	3.0
1	D	313	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	142	PHE	2.9
1	F	234	PRO	2.9
1	B	30	SER	2.8
1	A	218	ASP	2.8
1	B	75	PHE	2.8
1	B	8	PRO	2.7
1	B	78	ALA	2.7
1	E	236	GLY	2.6
1	B	127	ARG	2.5
1	C	170	ALA	2.5
1	A	7	GLU	2.5
1	B	76	SER	2.5
1	B	70	ALA	2.4
1	E	233	GLY	2.3
1	B	142	PHE	2.3
1	F	233	GLY	2.2
1	C	234	PRO	2.2
1	F	8	PRO	2.2
1	F	141	TYR	2.2
1	F	236	GLY	2.1
1	E	217	ASP	2.1
1	B	170	ALA	2.1
1	E	114	LEU	2.1
1	C	210[A]	MET	2.0
1	B	233	GLY	2.0

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 monosaccharides in this entry.

6.4 Ligands ⓘ

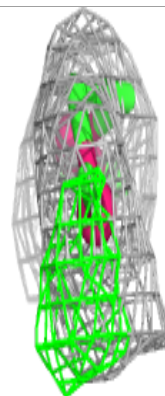
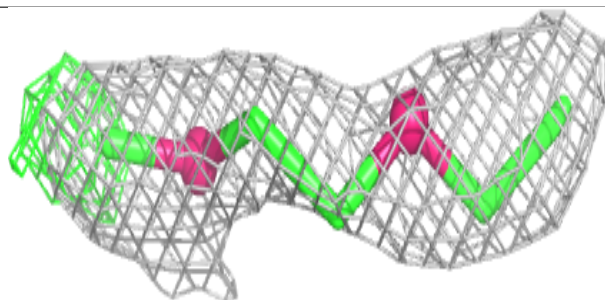
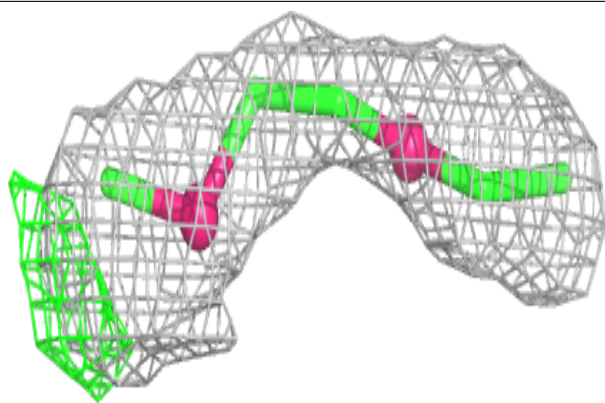
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	7PE	F	409	7/21	0.69	0.22	62,69,74,77	0
4	MG	A	406	1/1	0.72	0.14	56,56,56,56	0
5	GOL	F	407	6/6	0.72	0.21	59,63,67,69	0
3	7PE	A	402	6/21	0.73	0.43	48,52,55,57	0
5	GOL	F	403	6/6	0.74	0.17	58,66,68,69	0
5	GOL	C	405	6/6	0.74	0.32	48,58,60,62	0
5	GOL	B	403	6/6	0.75	0.19	49,59,64,67	0
5	GOL	B	405	6/6	0.79	0.14	54,58,59,60	0
3	7PE	F	402	9/21	0.80	0.37	39,57,65,72	0
5	GOL	F	404	6/6	0.80	0.30	59,62,64,65	0
3	7PE	D	402	6/21	0.80	0.33	49,56,56,59	0
5	GOL	C	403	6/6	0.83	0.18	47,61,62,68	0
5	GOL	C	402	6/6	0.84	0.13	44,52,54,63	0
3	7PE	B	404	5/21	0.84	0.19	49,56,59,60	0
5	GOL	F	406	6/6	0.85	0.20	38,60,65,67	0
3	7PE	F	408	5/21	0.86	0.28	51,54,56,57	0
3	7PE	B	402	5/21	0.88	0.17	50,53,56,58	0
3	7PE	F	410	5/21	0.90	0.31	38,46,49,50	0
3	7PE	A	403	4/21	0.91	0.26	42,42,43,45	0
3	7PE	F	405[A]	6/21	0.92	0.28	14,21,25,25	6
4	MG	E	402	1/1	0.94	0.03	54,54,54,54	0
4	MG	A	405	1/1	0.96	0.04	41,41,41,41	0
2	HEM	B	401	43/43	0.97	0.10	23,25,28,30	0
4	MG	C	404	1/1	0.97	0.05	30,30,30,30	0
2	HEM	C	401	43/43	0.97	0.11	21,25,28,29	0
2	HEM	E	401	43/43	0.97	0.09	22,25,28,29	0
2	HEM	F	401	43/43	0.97	0.10	21,24,26,27	0
2	HEM	A	401	43/43	0.97	0.10	21,24,26,32	0
4	MG	A	404	1/1	0.98	0.04	27,27,27,27	0
2	HEM	D	401	43/43	0.98	0.10	21,23,26,27	0

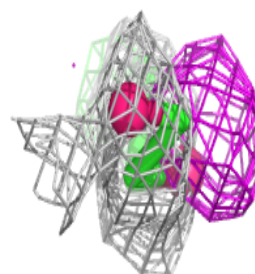
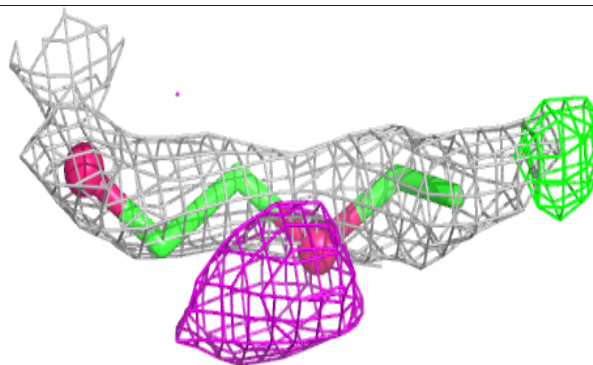
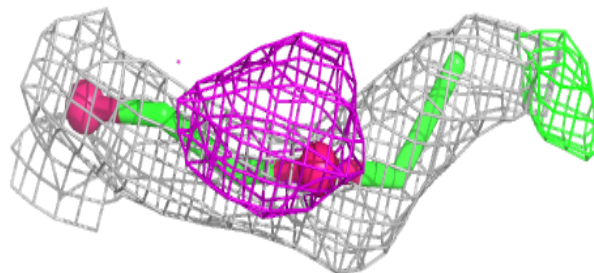
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 7PE F 409:

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

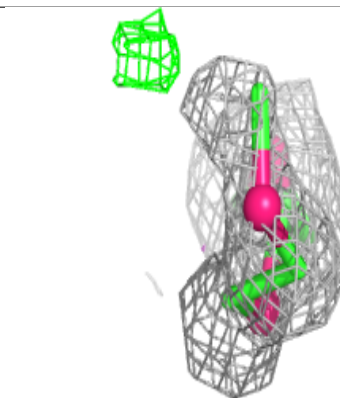
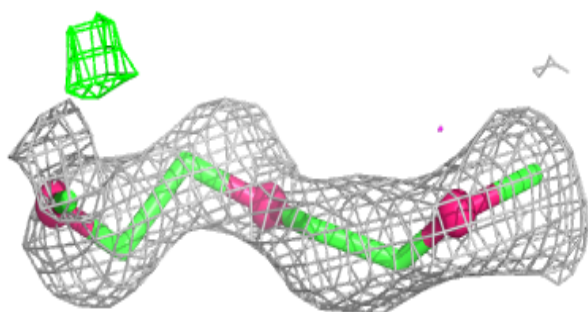
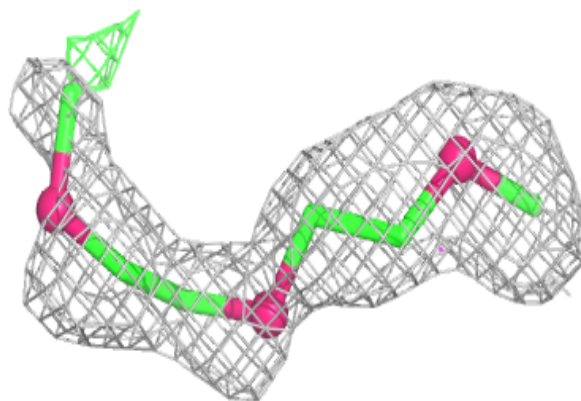
**Electron density around 7PE A 402:**

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

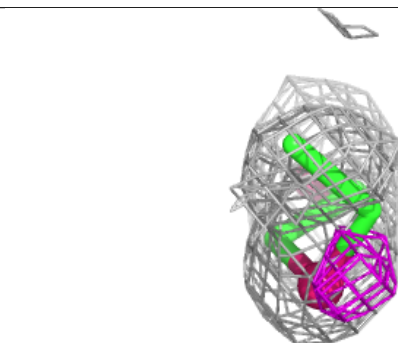
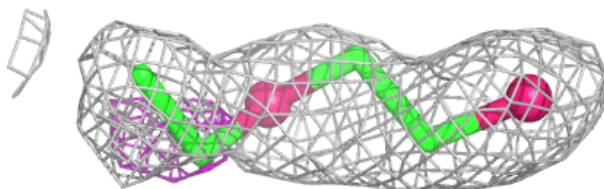
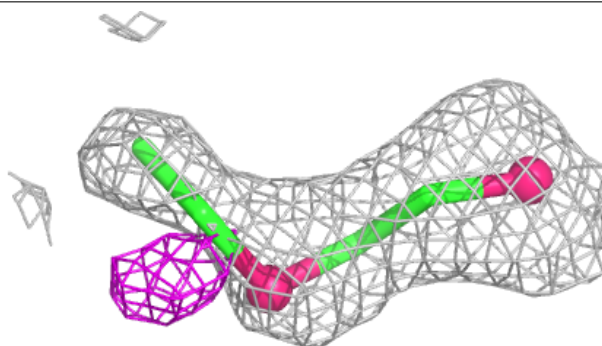


Electron density around 7PE F 402:

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

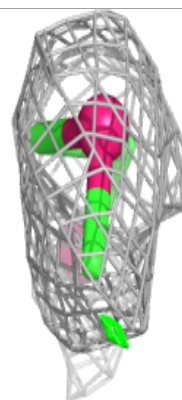
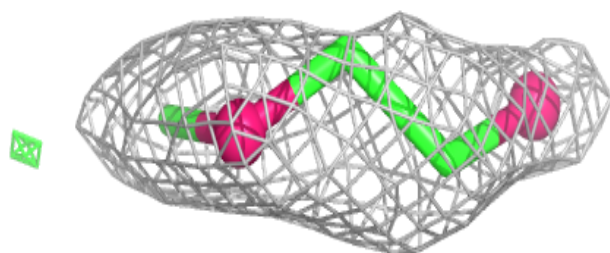
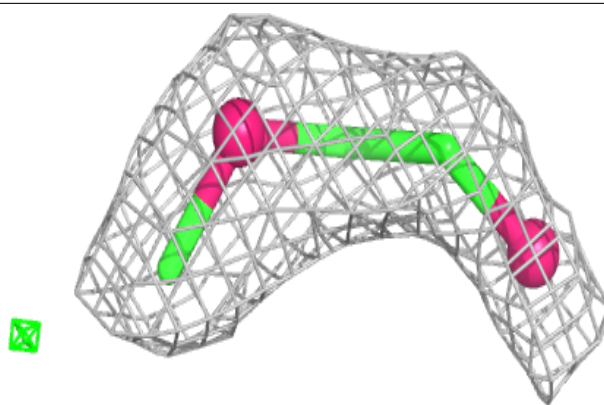
**Electron density around 7PE D 402:**

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

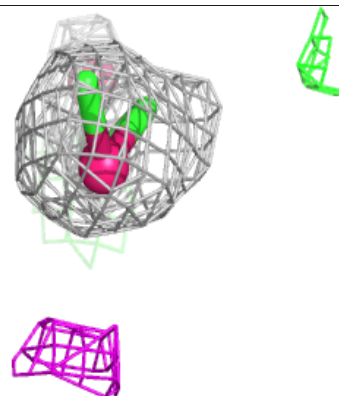
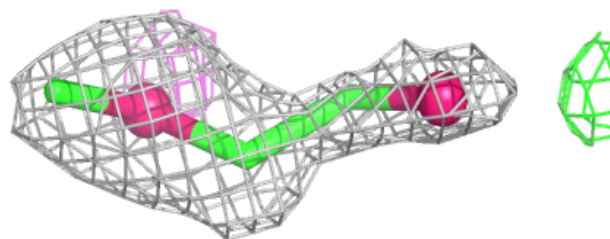
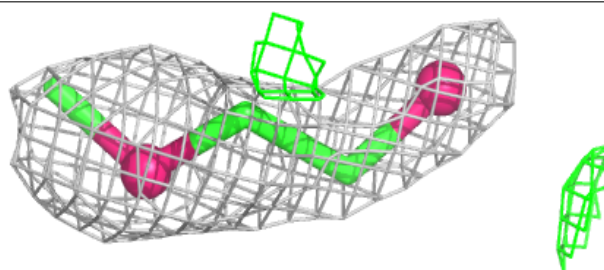


Electron density around 7PE B 404:

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

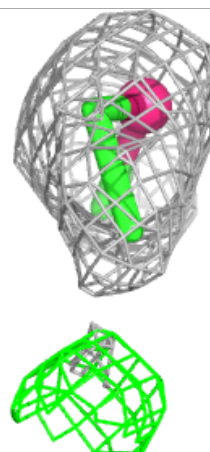
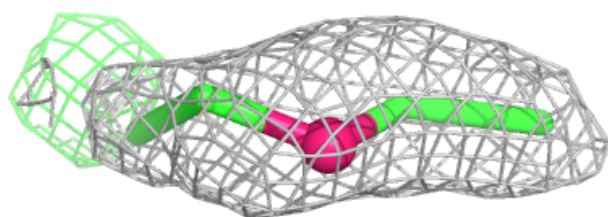
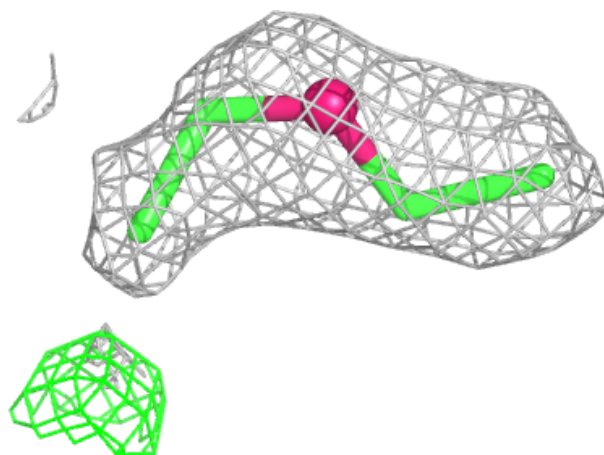
**Electron density around 7PE F 408:**

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



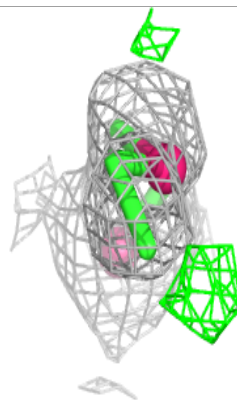
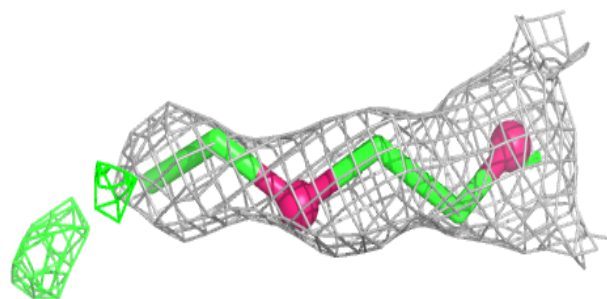
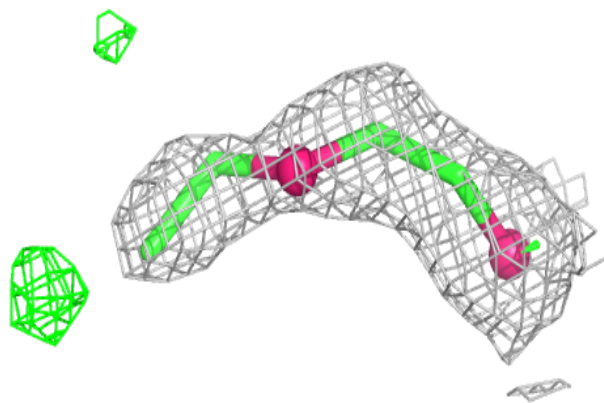
Electron density around 7PE B 402:

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



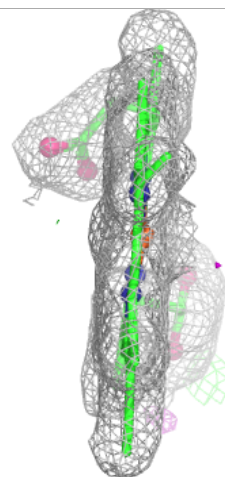
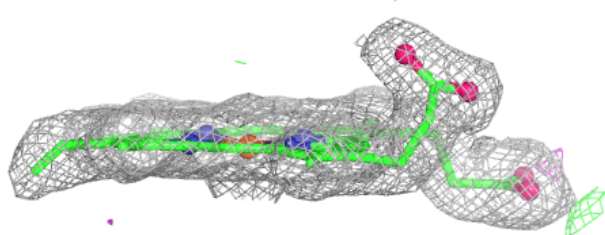
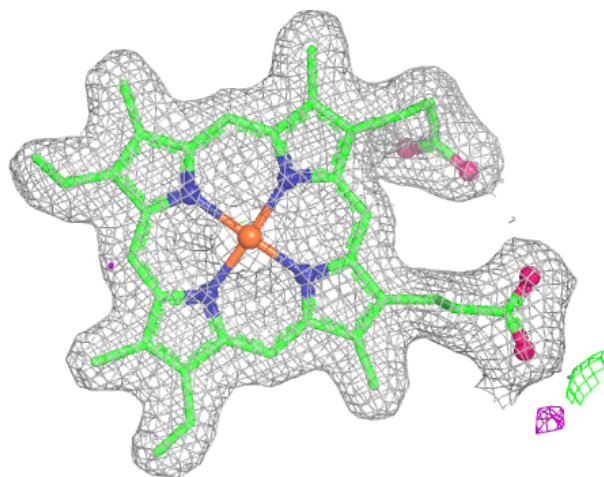
Electron density around 7PE F 405 (A):

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



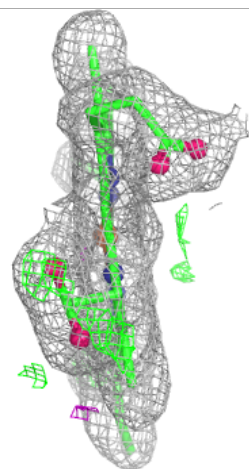
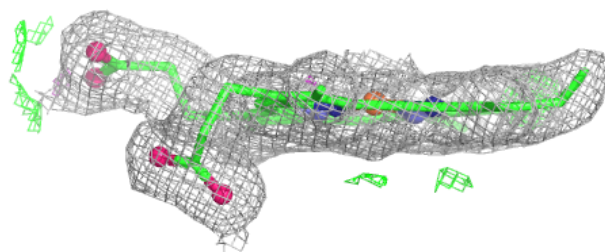
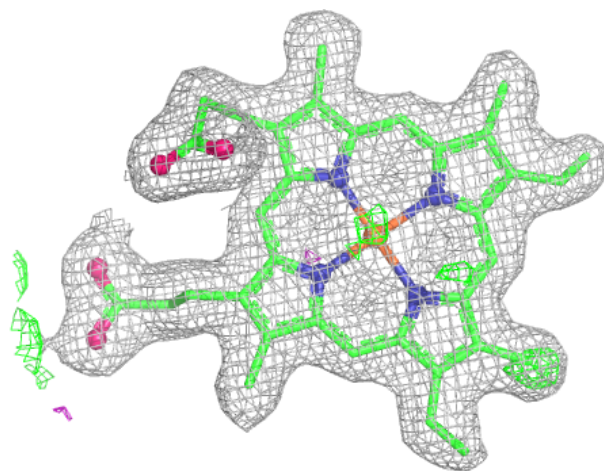
Electron density around HEM B 401:

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



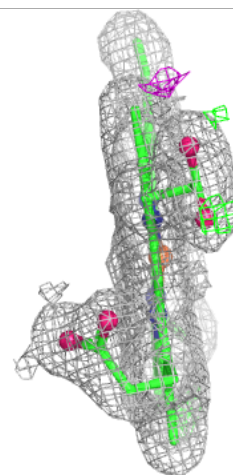
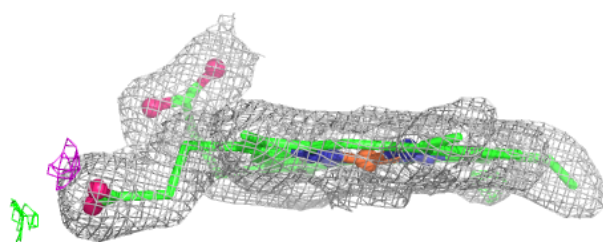
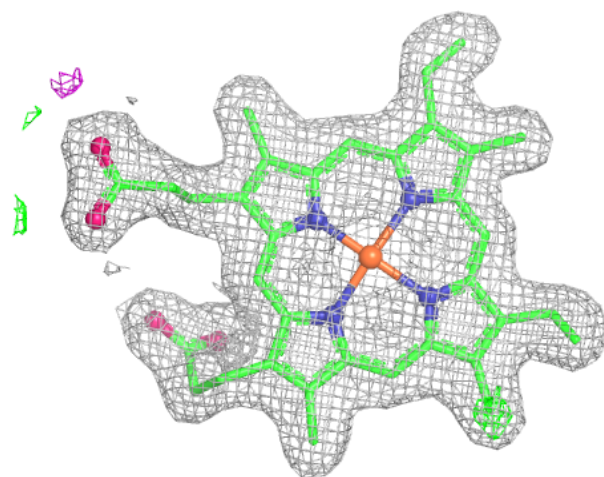
Electron density around HEM C 401:

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



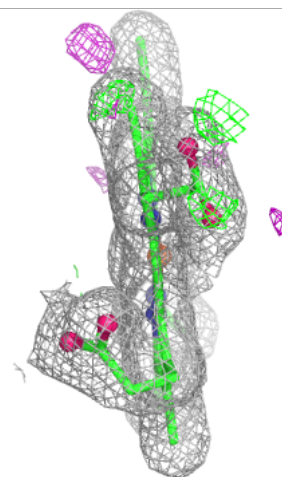
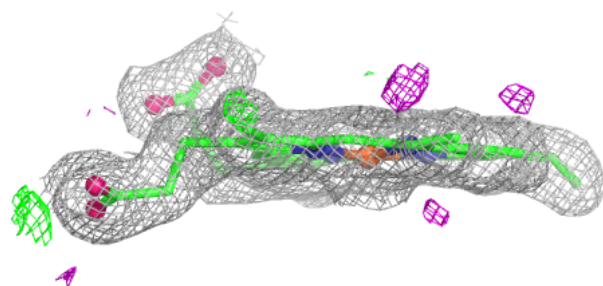
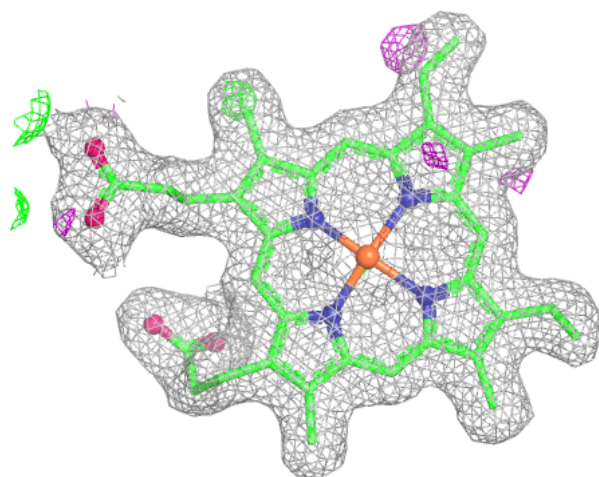
Electron density around HEM E 401:

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



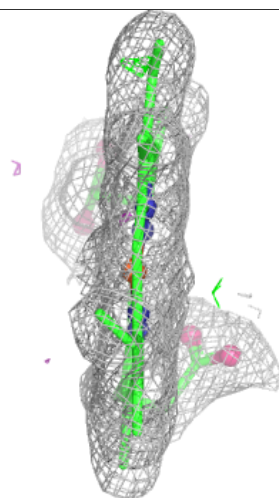
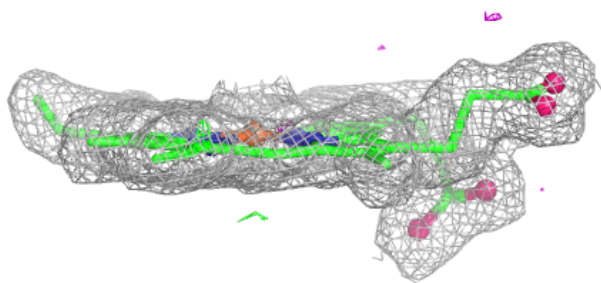
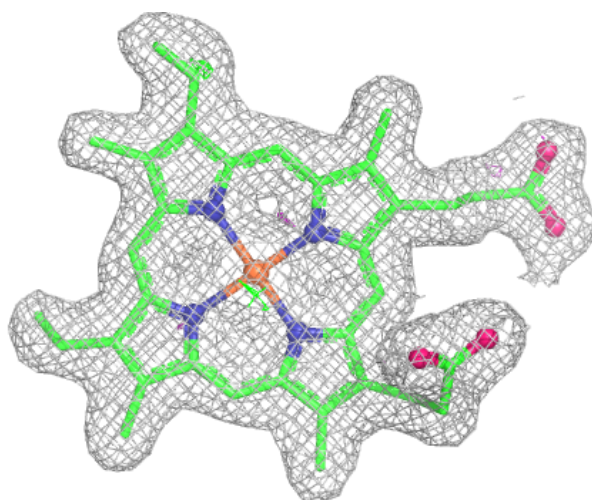
Electron density around HEM F 401:

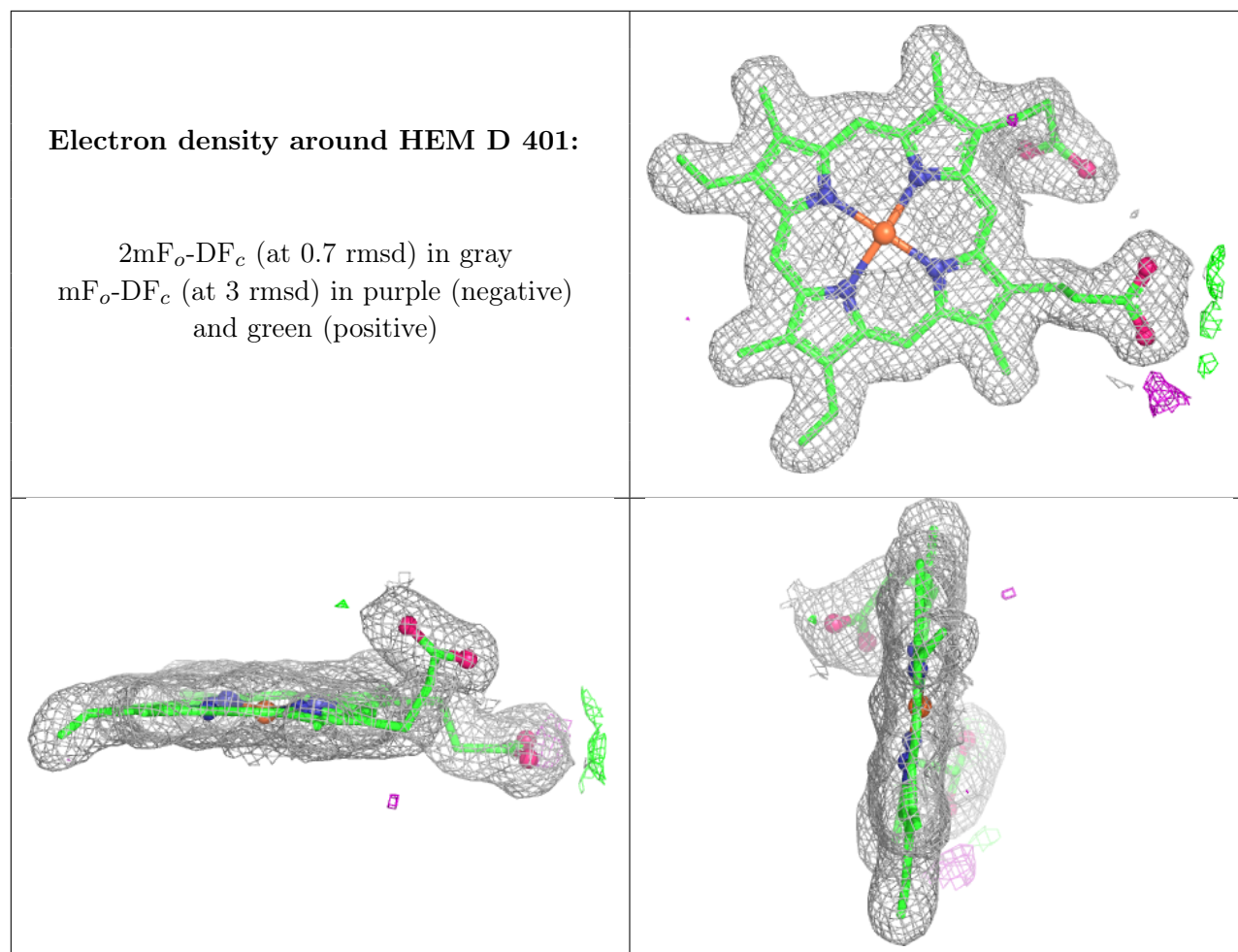
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 401:

$2mF_o-DF_c$ (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.