



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 25, 2025 – 02:02 am BST

PDB ID : 9HO7 / pdb_00009ho7
Title : DtpB in complex with photocaged nitric oxide, 10 ms, 30 microjoule, SFX
Authors : Smyth, P.; Williams, L.J.; Hough, M.A.; Worrall, J.A.R.; Owen, R.L.
Deposited on : 2024-12-11
Resolution : 1.69 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

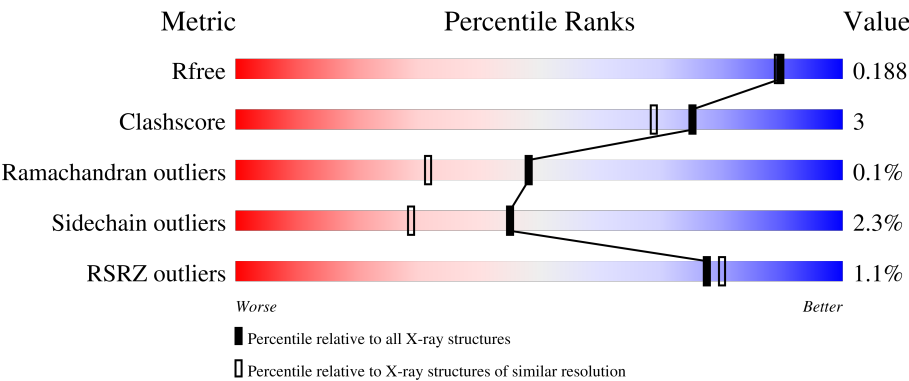
MolProbity : 4-5-2 with Phenix2.0rc1
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

1 Overall quality at a glance i

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

The reported resolution of this entry is 1.69 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.



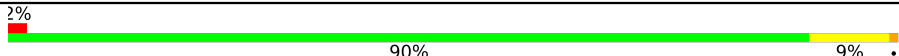
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	164625	5161 (1.70-1.70)
Clashscore	180529	5671 (1.70-1.70)
Ramachandran outliers	177936	5594 (1.70-1.70)
Sidechain outliers	177891	5594 (1.70-1.70)
RSRZ outliers	164620	5159 (1.70-1.70)

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	307	<div><div></div><div>92%6% .</div></div>
1	B	307	<div><div>2%</div><div>92%7% ..</div></div>
1	C	307	<div><div>%</div><div>88%11%</div></div>
1	D	307	<div><div>2%</div><div>90%8% .</div></div>
1	E	307	<div><div>%</div><div>88%11% .</div></div>

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Mol	Chain	Length	Quality of chain
1	F	307	

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	NO	B	402	-	-	X	-
3	NO	C	402	-	-	X	-
3	NO	E	402	-	-	X	-

2 Entry composition [i](#)

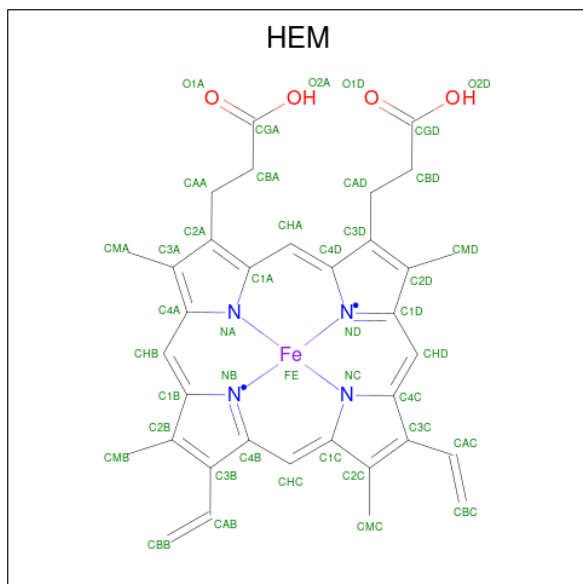
There are 4 unique types of molecules in this entry. The entry contains 15447 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dyp-type peroxidase family.

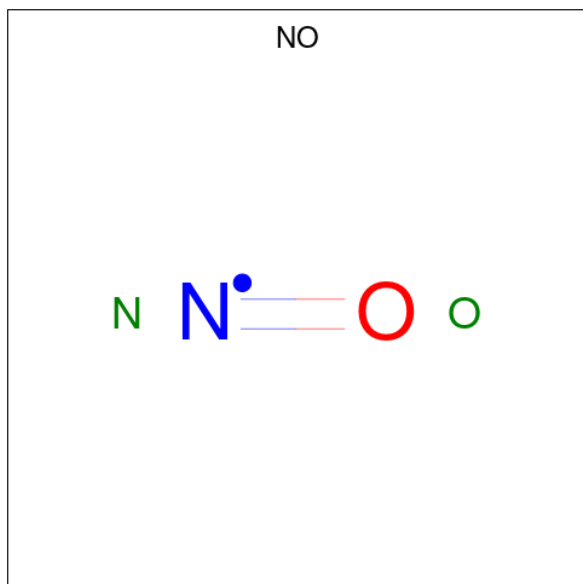
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	4	0
			2368	1487	412	459	10			
1	B	306	Total	C	N	O	S	0	3	0
			2343	1475	400	459	9			
1	C	306	Total	C	N	O	S	0	6	0
			2383	1498	411	465	9			
1	D	306	Total	C	N	O	S	0	3	0
			2353	1479	409	456	9			
1	E	306	Total	C	N	O	S	0	9	0
			2403	1509	418	467	9			
1	F	306	Total	C	N	O	S	0	3	0
			2355	1479	407	460	9			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: 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 NITRIC OXIDE (CCD ID: NO) (formula: NO) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	N	O		
			2	1	1		
3	B	1	Total	N	O		
			2	1	1		
3	C	1	Total	N	O		
			2	1	1		
3	D	1	Total	N	O		
			2	1	1		
3	E	1	Total	N	O		
			2	1	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	N	O	0	0
			2	1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	180	Total	O	0	0
			180	180		
4	B	148	Total	O	0	0
			148	148		
4	C	150	Total	O	0	0
			150	150		
4	D	179	Total	O	0	0
			179	179		
4	E	167	Total	O	0	0
			167	167		
4	F	148	Total	O	0	0
			148	148		

3 Residue-property plots [i](#)

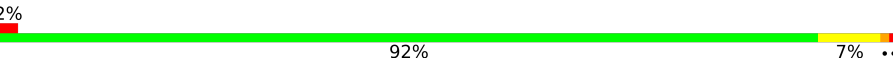
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Dyp-type peroxidase family

Chain A: 




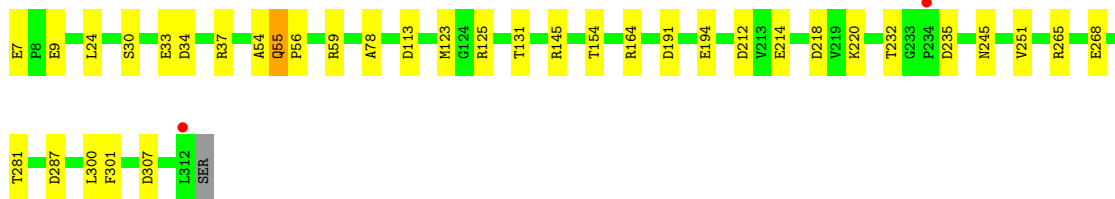
- Molecule 1: Dyp-type peroxidase family

Chain B: 



- Molecule 1: Dyp-type peroxidase family

Chain C: 




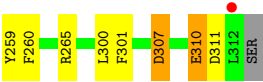
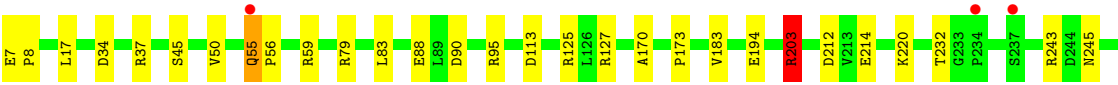
- Molecule 1: Dyp-type peroxidase family

Chain D: 

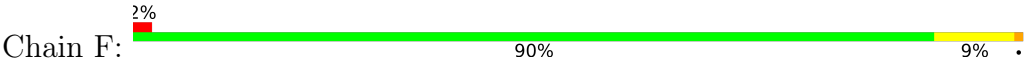


- Molecule 1: Dyp-type peroxidase family

Chain E: 



● Molecule 1: Dyp-type peroxidase family



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.17Å 123.06Å 195.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	104.11 – 1.69 104.11 – 1.69	Depositor EDS
% Data completeness (in resolution range)	100.0 (104.11-1.69) 99.9 (104.11-1.69)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 1.69Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.190 , 0.218 0.184 , 0.188	Depositor DCC
R_{free} test set	1958 reflections (0.84%)	wwPDB-VP
Wilson B-factor (Å ²)	20.8	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15447	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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: HEM, NO

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	0.77	0/2421	1.18	11/3287 (0.3%)
1	B	0.74	0/2399	1.20	11/3261 (0.3%)
1	C	0.71	0/2433	1.16	13/3305 (0.4%)
1	D	0.79	0/2409	1.21	12/3270 (0.4%)
1	E	0.75	0/2456	1.18	15/3336 (0.4%)
1	F	0.74	0/2411	1.16	11/3273 (0.3%)
All	All	0.75	0/14529	1.18	73/19732 (0.4%)

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	B	0	1
1	C	0	2
1	D	0	4
1	E	0	3
1	F	0	3
All	All	0	13

There are no bond length outliers.

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	55	GLN	CB-CA-C	12.60	127.79	109.26
1	D	210	MET	CG-SD-CE	-11.68	75.20	100.90
1	C	55	GLN	CB-CA-C	11.02	126.93	109.52
1	B	12	MET	CG-SD-CE	9.98	122.85	100.90
1	B	210	MET	CG-SD-CE	-9.39	80.24	100.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	232	THR	CA-CB-OG1	-8.27	97.19	109.60
1	C	55	GLN	N-CA-CB	-8.26	97.00	109.98
1	A	232	THR	CA-CB-OG1	-7.82	97.88	109.60
1	E	232	THR	CA-CB-OG1	-7.51	98.33	109.60
1	A	9	GLU	CB-CA-C	7.12	117.97	109.85
1	D	113	ASP	CA-CB-CG	7.01	119.61	112.60
1	F	55	GLN	CB-CA-C	6.91	120.44	109.52
1	B	113	ASP	CA-CB-CG	6.82	119.42	112.60
1	E	113	ASP	CA-CB-CG	6.64	119.24	112.60
1	D	8	PRO	N-CA-CB	-6.64	96.28	103.25
1	E	203[A]	ARG	NE-CZ-NH1	-6.61	114.89	121.50
1	E	203[B]	ARG	NE-CZ-NH1	-6.61	114.89	121.50
1	E	212	ASP	CA-CB-CG	6.33	118.93	112.60
1	C	281	THR	CA-CB-OG1	-6.30	100.15	109.60
1	A	47	GLU	CG-CD-OE2	-6.29	103.93	118.40
1	E	265	ARG	N-CA-CB	-6.25	100.02	110.39
1	C	145	ARG	CA-CB-CG	-6.19	101.72	114.10
1	D	145	ARG	CA-CB-CG	-6.04	102.02	114.10
1	C	287	ASP	CA-CB-CG	6.02	118.62	112.60
1	C	251	VAL	N-CA-CB	5.89	119.42	110.58
1	D	265	ARG	N-CA-CB	-5.79	100.78	110.39
1	C	232	THR	CA-CB-OG1	-5.76	100.96	109.60
1	D	199	GLU	CB-CG-CD	5.76	122.39	112.60
1	A	9	GLU	N-CA-CB	-5.74	101.63	110.01
1	B	79	ARG	CB-CA-C	-5.73	100.91	108.76
1	F	55	GLN	N-CA-CB	-5.73	100.98	109.98
1	E	88	GLU	N-CA-CB	5.68	118.31	109.97
1	C	212	ASP	CA-CB-CG	5.66	118.25	112.60
1	D	212	ASP	CA-CB-CG	5.60	118.20	112.60
1	E	88	GLU	CB-CA-C	-5.58	100.11	109.65
1	C	154	THR	CA-CB-OG1	-5.57	101.24	109.60
1	F	266	THR	CA-CB-OG1	-5.54	101.28	109.60
1	D	244	ASP	CA-CB-CG	5.54	118.14	112.60
1	F	163	ARG	N-CA-CB	5.51	118.01	110.01
1	C	113[A]	ASP	CA-CB-CG	5.51	118.11	112.60
1	C	113[B]	ASP	CA-CB-CG	5.51	118.11	112.60
1	A	145	ARG	CA-CB-CG	-5.50	103.11	114.10
1	B	199	GLU	CB-CG-CD	5.44	121.85	112.60
1	E	127	ARG	N-CA-CB	5.40	117.95	109.69
1	F	194	GLU	CB-CG-CD	5.39	121.77	112.60
1	E	307	ASP	CA-CB-CG	5.38	117.98	112.60
1	F	72	ASP	CA-CB-CG	5.37	117.97	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	143	ASP	CA-CB-CG	5.37	117.97	112.60
1	A	212	ASP	CA-CB-CG	5.36	117.96	112.60
1	B	307	ASP	CB-CA-C	-5.34	100.74	110.63
1	A	287	ASP	CA-CB-CG	5.32	117.92	112.60
1	B	194	GLU	CB-CG-CD	5.31	121.63	112.60
1	C	194	GLU	CB-CG-CD	5.29	121.59	112.60
1	E	203[A]	ARG	NE-CZ-NH2	5.28	123.95	119.20
1	E	203[B]	ARG	NE-CZ-NH2	5.28	123.95	119.20
1	F	113	ASP	CA-CB-CG	5.25	117.85	112.60
1	E	50	VAL	CA-C-N	5.22	125.78	119.98
1	E	50	VAL	C-N-CA	5.22	125.78	119.98
1	B	212	ASP	CA-CB-CG	5.20	117.80	112.60
1	F	175	PHE	CA-CB-CG	-5.18	108.62	113.80
1	F	287	ASP	CA-CB-CG	5.17	117.77	112.60
1	E	311	ASP	CA-CB-CG	5.17	117.77	112.60
1	A	56	PRO	N-CA-CB	-5.13	96.96	102.60
1	A	171	GLU	CB-CG-CD	5.13	121.32	112.60
1	B	234	PRO	N-CA-CB	-5.12	97.87	103.25
1	B	158	THR	CA-CB-OG1	-5.12	101.93	109.60
1	D	219	VAL	N-CA-CB	-5.11	105.30	112.35
1	D	255	GLU	CB-CA-C	5.11	119.11	110.79
1	C	131	THR	CA-CB-OG1	-5.08	101.97	109.60
1	F	172	ASP	CA-CB-CG	5.08	117.68	112.60
1	D	287	ASP	CA-CB-CG	5.07	117.67	112.60
1	A	113	ASP	CA-CB-CG	5.04	117.64	112.60
1	A	307	ASP	CB-CA-C	-5.01	101.06	110.67

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	203	ARG	Sidechain
1	C	125	ARG	Sidechain
1	C	59	ARG	Sidechain
1	D	125	ARG	Sidechain
1	D	253	ARG	Sidechain
1	D	59[A]	ARG	Sidechain
1	D	59[B]	ARG	Sidechain
1	E	125[A]	ARG	Sidechain
1	E	203[A]	ARG	Sidechain
1	E	59	ARG	Sidechain
1	F	111	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	F	125	ARG	Sidechain
1	F	59	ARG	Sidechain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2368	0	2298	12	0
1	B	2343	0	2252	12	0
1	C	2383	0	2305	10	0
1	D	2353	0	2287	12	0
1	E	2403	0	2322	21	0
1	F	2355	0	2285	8	0
2	A	43	0	30	1	0
2	B	43	0	30	2	0
2	C	43	0	30	3	0
2	D	43	0	30	1	0
2	E	43	0	30	0	0
2	F	43	0	30	1	0
3	A	2	0	0	1	0
3	B	2	0	0	3	0
3	C	2	0	0	3	0
3	D	2	0	0	1	0
3	E	2	0	0	2	0
3	F	2	0	0	1	0
4	A	180	0	0	5	0
4	B	148	0	0	4	0
4	C	150	0	0	2	0
4	D	179	0	0	2	0
4	E	167	0	0	5	0
4	F	148	0	0	0	0
All	All	15447	0	13929	81	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 (81) 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:59[A]:ARG:NH1	4:D:501:HOH:O	1.72	1.18
1:D:245:ASN:ND2	3:D:402:NO:O	2.05	0.89
1:E:203[A]:ARG:HD2	4:E:654:HOH:O	1.73	0.86
1:E:55[B]:GLN:OE1	1:E:55[B]:GLN:HA	1.79	0.83
1:A:55[A]:GLN:OE1	4:A:501:HOH:O	1.98	0.82
1:A:191:ASP:OD2	4:A:502:HOH:O	2.00	0.79
1:B:55:GLN:HG3	4:B:501:HOH:O	1.83	0.78
1:A:55[B]:GLN:HA	4:A:531:HOH:O	1.88	0.74
1:A:55[A]:GLN:HA	4:A:531:HOH:O	1.88	0.73
1:E:90:ASP:OD2	1:E:95:ARG:NH1	2.23	0.70
1:D:34:ASP:OD1	1:D:37[B]:ARG:NH1	2.25	0.69
1:A:34:ASP:OD1	1:A:37[A]:ARG:NH2	2.26	0.68
1:B:55:GLN:NE2	4:B:501:HOH:O	2.27	0.68
1:E:245[B]:ASN:ND2	3:E:402:NO:O	2.27	0.68
1:B:55:GLN:CG	4:B:501:HOH:O	2.41	0.68
1:D:127:ARG:NH2	1:F:254:GLU:OE2	2.29	0.65
1:A:245:ASN:ND2	3:A:402:NO:O	2.29	0.65
2:C:401:HEM:HMC2	2:C:401:HEM:HBC2	1.78	0.65
1:E:55[A]:GLN:HG2	1:E:56:PRO:HA	1.79	0.64
1:F:245:ASN:ND2	3:F:402:NO:O	2.33	0.62
1:B:245:ASN:ND2	3:B:402:NO:O	2.33	0.61
1:A:147:MET:HG3	1:A:256:PHE:HB3	1.81	0.61
1:D:147:MET:HE2	1:D:256:PHE:HB3	1.83	0.61
1:C:33:GLU:O	1:C:37[B]:ARG:HG3	2.02	0.59
1:C:78:ALA:HA	4:C:554:HOH:O	2.03	0.58
1:E:95:ARG:HG3	4:E:510:HOH:O	2.03	0.58
1:E:194:GLU:HG2	4:E:650:HOH:O	2.05	0.56
1:B:243:ARG:HD2	3:B:402:NO:N	2.20	0.55
1:E:170:ALA:O	1:E:173:PRO:HD3	2.07	0.55
1:A:216:SER:HB3	1:A:218:ASP:OD1	2.06	0.55
1:D:7:GLU:HB2	1:D:8:PRO:HD3	1.88	0.55
1:E:214:GLU:OE2	1:E:220:LYS:NZ	2.31	0.55
1:A:56:PRO:HA	4:A:531:HOH:O	2.08	0.54
1:B:55:GLN:CD	4:B:501:HOH:O	2.51	0.54
2:B:401:HEM:NC	3:B:402:NO:N	2.56	0.53
2:B:401:HEM:HBC2	2:B:401:HEM:HMC2	1.90	0.53
2:F:401:HEM:HBC2	2:F:401:HEM:HMC2	1.91	0.53
1:A:57:ASP:O	1:A:59:ARG:NH1	2.41	0.53
2:A:401:HEM:HMC2	2:A:401:HEM:HBC2	1.90	0.53
1:C:164:ARG:O	1:C:265:ARG:HD2	2.10	0.52
1:B:52:PHE:CD1	1:B:55:GLN:OE1	2.62	0.52
1:E:7:GLU:N	1:E:8:PRO:CD	2.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:34:ASP:OD1	1:F:37:ARG:NH1	2.44	0.51
1:C:34:ASP:HA	1:C:37[B]:ARG:HG3	1.95	0.48
1:C:245:ASN:ND2	3:C:402:NO:O	2.46	0.47
1:E:34:ASP:OD1	1:E:37:ARG:NH2	2.44	0.47
1:C:24[A]:LEU:HD21	1:C:123:MET:SD	2.55	0.47
1:C:214:GLU:OE2	1:C:220:LYS:NZ	2.44	0.47
1:B:57:ASP:O	1:B:59:ARG:NH1	2.48	0.46
1:D:127:ARG:CG	4:D:669:HOH:O	2.64	0.46
1:E:183[B]:VAL:CG1	1:E:259:TYR:OH	2.64	0.46
1:C:55:GLN:HA	1:C:56:PRO:HA	1.78	0.45
1:E:300:LEU:HD23	1:E:301:PHE:N	2.32	0.45
1:E:90:ASP:HA	1:E:95:ARG:HG2	2.00	0.44
1:D:170:ALA:O	1:D:173:PRO:HD3	2.18	0.44
1:D:231:VAL:HG23	1:D:239:LEU:HB2	1.99	0.43
1:B:241:ILE:HG22	1:B:264:ALA:HB2	1.99	0.43
2:C:401:HEM:NB	3:C:402:NO:N	2.67	0.43
2:C:401:HEM:NC	3:C:402:NO:N	2.66	0.43
1:F:271:GLU:O	1:F:275:GLU:HG3	2.19	0.43
1:E:300:LEU:HD23	1:E:300:LEU:C	2.43	0.43
1:E:90:ASP:OD2	1:E:95:ARG:HG2	2.19	0.42
1:E:245[A]:ASN:HB3	1:E:260:PHE:HA	2.00	0.42
1:A:162:ALA:O	1:A:166:VAL:HG13	2.19	0.42
1:C:300:LEU:HD23	1:C:301:PHE:N	2.34	0.42
1:D:231:VAL:CG2	1:D:239:LEU:HB2	2.49	0.42
1:D:251:VAL:HG12	1:F:123:MET:HG3	2.00	0.42
1:F:90:ASP:OD2	1:F:95:ARG:HD2	2.20	0.42
1:B:309:LEU:HA	1:B:312:LEU:HD23	2.01	0.42
1:A:55[A]:GLN:HG2	1:A:56:PRO:HA	2.02	0.41
1:C:54:ALA:HB1	4:C:501:HOH:O	2.20	0.41
1:E:243:ARG:HD2	3:E:402:NO:N	2.36	0.41
1:B:12:MET:HG3	1:B:15:SER:HB3	2.01	0.41
1:F:211:THR:O	1:F:212:ASP:HB3	2.20	0.41
1:D:59[A]:ARG:HG2	1:D:110:SER:HB3	2.03	0.41
1:F:127:ARG:HH11	1:F:127:ARG:CG	2.34	0.40
1:B:39:LEU:C	1:B:39:LEU:HD13	2.46	0.40
1:E:79:ARG:NH2	1:E:83:LEU:O	2.55	0.40
2:D:401:HEM:HMC2	2:D:401:HEM:HBC2	2.02	0.40
1:E:310:GLU:OE2	4:E:501:HOH:O	2.22	0.40
1:E:203[B]:ARG:HD3	4:E:654:HOH:O	2.21	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	308/307 (100%)	303 (98%)	5 (2%)	0	100	100
1	B	307/307 (100%)	300 (98%)	6 (2%)	1 (0%)	37	23
1	C	310/307 (101%)	305 (98%)	5 (2%)	0	100	100
1	D	307/307 (100%)	300 (98%)	6 (2%)	1 (0%)	37	23
1	E	313/307 (102%)	305 (97%)	8 (3%)	0	100	100
1	F	308/307 (100%)	301 (98%)	7 (2%)	0	100	100
All	All	1853/1842 (101%)	1814 (98%)	37 (2%)	2 (0%)	48	32

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	8	PRO
1	B	234	PRO

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	249/246 (101%)	243 (98%)	6 (2%)	44	27
1	B	244/246 (99%)	239 (98%)	5 (2%)	50	34
1	C	251/246 (102%)	243 (97%)	8 (3%)	34	17
1	D	248/246 (101%)	246 (99%)	2 (1%)	79	71
1	E	252/246 (102%)	246 (98%)	6 (2%)	44	27

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	249/246 (101%)	240 (96%)	9 (4%)	30	14
All	All	1493/1476 (101%)	1457 (98%)	36 (2%)	45	27

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55[A]	GLN
1	A	55[B]	GLN
1	A	56	PRO
1	A	147	MET
1	A	166	VAL
1	A	307	ASP
1	B	12	MET
1	B	24	LEU
1	B	210	MET
1	B	234	PRO
1	B	307	ASP
1	C	7	GLU
1	C	9	GLU
1	C	30	SER
1	C	191	ASP
1	C	218	ASP
1	C	235	ASP
1	C	268	GLU
1	C	307	ASP
1	D	127	ARG
1	D	307	ASP
1	E	17	LEU
1	E	45	SER
1	E	55[A]	GLN
1	E	55[B]	GLN
1	E	307	ASP
1	E	310	GLU
1	F	9	GLU
1	F	38	ASP
1	F	55	GLN
1	F	87[A]	ARG
1	F	87[B]	ARG
1	F	127	ARG
1	F	235	ASP
1	F	283	SER
1	F	307	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	55	GLN
1	C	55	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

12 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	HEM	E	401	3,1	41,50,50	1.35	3 (7%)	45,82,82	1.99	13 (28%)
3	NO	D	402	2	0,1,1	-	-	-		
3	NO	F	402	2	0,1,1	-	-	-		
2	HEM	D	401	3,1	41,50,50	1.38	6 (14%)	45,82,82	1.81	12 (26%)
2	HEM	B	401	3,1	41,50,50	1.34	3 (7%)	45,82,82	1.95	13 (28%)
3	NO	C	402	2	0,1,1	-	-	-		
2	HEM	F	401	3,1	41,50,50	1.25	6 (14%)	45,82,82	1.97	11 (24%)
3	NO	E	402	2	0,1,1	-	-	-		
3	NO	A	402	2	0,1,1	-	-	-		

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NO	B	402	2	0,1,1	-	-	-		
2	HEM	C	401	3,1	41,50,50	1.51	7 (17%)	45,82,82	2.09	15 (33%)
2	HEM	A	401	3,1	41,50,50	1.44	9 (21%)	45,82,82	2.00	16 (35%)

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
2	HEM	E	401	3,1	-	6/12/54/54	-
2	HEM	D	401	3,1	-	6/12/54/54	-
2	HEM	B	401	3,1	-	5/12/54/54	-
2	HEM	F	401	3,1	-	4/12/54/54	-
2	HEM	C	401	3,1	-	5/12/54/54	-
2	HEM	A	401	3,1	-	6/12/54/54	-

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	HEM	C4D-ND	-3.92	1.33	1.40
2	B	401	HEM	C1B-NB	-3.80	1.33	1.40
2	D	401	HEM	C1B-NB	-3.80	1.33	1.40
2	F	401	HEM	C1B-NB	-3.53	1.34	1.40
2	E	401	HEM	C1B-NB	-3.47	1.34	1.40
2	C	401	HEM	C1B-NB	-3.26	1.34	1.40
2	E	401	HEM	C4D-ND	-3.19	1.34	1.40
2	B	401	HEM	CAA-C2A	3.14	1.56	1.52
2	A	401	HEM	CHA-C4D	3.07	1.42	1.35
2	C	401	HEM	C4D-C3D	3.02	1.50	1.45
2	A	401	HEM	C3C-C2C	-3.00	1.36	1.40
2	D	401	HEM	C3C-C2C	-2.86	1.36	1.40
2	D	401	HEM	C4D-ND	-2.65	1.35	1.40
2	C	401	HEM	C4B-NB	-2.62	1.33	1.38
2	B	401	HEM	C1D-ND	-2.49	1.33	1.38
2	F	401	HEM	C4A-NA	2.38	1.41	1.36
2	A	401	HEM	C3C-CAC	2.37	1.52	1.47
2	A	401	HEM	C3B-C4B	2.35	1.49	1.44
2	D	401	HEM	C1D-ND	-2.31	1.34	1.38
2	D	401	HEM	CHA-C4D	2.28	1.40	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	HEM	C1A-NA	2.26	1.40	1.36
2	A	401	HEM	CAB-C3B	2.26	1.53	1.47
2	A	401	HEM	C1B-NB	-2.24	1.36	1.40
2	C	401	HEM	FE-NB	2.24	2.07	1.96
2	E	401	HEM	CMB-C2B	2.24	1.55	1.50
2	D	401	HEM	CMB-C2B	2.23	1.55	1.50
2	F	401	HEM	C3C-C2C	-2.22	1.37	1.40
2	A	401	HEM	O1D-CGD	2.19	1.29	1.22
2	C	401	HEM	C1D-C2D	2.18	1.48	1.44
2	C	401	HEM	CMD-C2D	-2.15	1.46	1.50
2	F	401	HEM	C1D-ND	-2.14	1.34	1.38
2	F	401	HEM	C4D-C3D	2.13	1.48	1.45
2	F	401	HEM	CHB-C1B	2.06	1.40	1.35
2	A	401	HEM	C1D-ND	-2.04	1.34	1.38

All (80) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	HEM	CHC-C4B-NB	6.57	131.57	124.43
2	C	401	HEM	CHC-C4B-NB	6.41	131.39	124.43
2	C	401	HEM	C1B-NB-C4B	5.46	110.71	105.07
2	D	401	HEM	CHC-C4B-NB	5.19	130.07	124.43
2	B	401	HEM	CHC-C4B-NB	5.16	130.04	124.43
2	B	401	HEM	C1B-NB-C4B	5.08	110.32	105.07
2	E	401	HEM	CHC-C4B-NB	5.07	129.94	124.43
2	E	401	HEM	C1B-NB-C4B	5.00	110.23	105.07
2	F	401	HEM	C1B-NB-C4B	4.78	110.01	105.07
2	A	401	HEM	C1B-NB-C4B	4.52	109.74	105.07
2	B	401	HEM	CHB-C1B-NB	4.12	129.47	124.38
2	A	401	HEM	CHC-C4B-NB	4.01	128.79	124.43
2	C	401	HEM	C4A-C3A-C2A	3.96	109.75	107.00
2	E	401	HEM	CHB-C1B-NB	3.93	129.23	124.38
2	F	401	HEM	CHB-C1B-NB	3.86	129.15	124.38
2	A	401	HEM	C3B-C2B-C1B	3.82	109.32	106.49
2	D	401	HEM	CHA-C4D-ND	3.52	128.73	124.38
2	D	401	HEM	C1B-NB-C4B	3.45	108.63	105.07
2	C	401	HEM	CMA-C3A-C4A	-3.42	123.20	128.46
2	A	401	HEM	CHB-C1B-NB	3.25	128.40	124.38
2	C	401	HEM	CMD-C2D-C1D	3.19	129.90	125.04
2	A	401	HEM	CHA-C4D-ND	3.18	128.31	124.38
2	E	401	HEM	CHA-C4D-ND	3.14	128.25	124.38
2	E	401	HEM	CAD-C3D-C4D	3.11	130.09	124.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	401	HEM	CHA-C4D-ND	3.10	128.22	124.38
2	A	401	HEM	CHD-C1D-ND	2.94	127.62	124.43
2	D	401	HEM	CHA-C4D-C3D	-2.93	119.82	125.33
2	F	401	HEM	C4A-C3A-C2A	-2.93	104.96	107.00
2	B	401	HEM	CMA-C3A-C4A	-2.91	123.99	128.46
2	B	401	HEM	CHA-C4D-C3D	-2.91	119.87	125.33
2	D	401	HEM	CHD-C1D-C2D	-2.89	120.46	124.98
2	A	401	HEM	CHD-C1D-C2D	-2.88	120.48	124.98
2	A	401	HEM	CHA-C4D-C3D	-2.84	119.99	125.33
2	A	401	HEM	CMA-C3A-C4A	-2.84	124.10	128.46
2	D	401	HEM	CMD-C2D-C1D	2.83	129.34	125.04
2	F	401	HEM	CHD-C1D-C2D	-2.79	120.62	124.98
2	E	401	HEM	C3B-C2B-C1B	2.78	108.55	106.49
2	E	401	HEM	CHD-C1D-C2D	-2.75	120.69	124.98
2	A	401	HEM	CAD-CBD-CGD	-2.68	107.83	113.60
2	B	401	HEM	CHA-C4D-ND	2.66	127.67	124.38
2	A	401	HEM	C4B-C3B-C2B	-2.61	105.04	107.11
2	E	401	HEM	C4B-C3B-C2B	-2.58	105.06	107.11
2	E	401	HEM	CHA-C4D-C3D	-2.57	120.50	125.33
2	D	401	HEM	C2C-C3C-C4C	-2.52	105.14	106.90
2	B	401	HEM	CMA-C3A-C2A	2.51	129.68	124.94
2	B	401	HEM	CAB-C3B-C2B	-2.50	120.35	128.60
2	C	401	HEM	CMB-C2B-C1B	2.50	128.85	125.04
2	C	401	HEM	O2A-CGA-CBA	2.50	122.06	114.03
2	C	401	HEM	CBA-CAA-C2A	2.46	116.81	112.62
2	D	401	HEM	C2D-C1D-ND	2.40	112.76	109.88
2	A	401	HEM	CAB-C3B-C2B	-2.39	120.74	128.60
2	F	401	HEM	C2D-C1D-ND	2.38	112.73	109.88
2	C	401	HEM	CHA-C4D-C3D	-2.34	120.94	125.33
2	D	401	HEM	CBA-CAA-C2A	2.32	116.58	112.62
2	B	401	HEM	O2A-CGA-CBA	2.31	121.44	114.03
2	C	401	HEM	CAA-CBA-CGA	-2.24	107.48	113.76
2	E	401	HEM	CHD-C1D-ND	2.24	126.86	124.43
2	B	401	HEM	CAD-C3D-C4D	2.22	128.54	124.66
2	A	401	HEM	CMB-C2B-C3B	-2.20	122.91	128.30
2	E	401	HEM	CAD-C3D-C2D	-2.19	123.80	127.88
2	F	401	HEM	CMA-C3A-C2A	2.17	129.03	124.94
2	C	401	HEM	C3D-C4D-ND	2.16	112.57	110.17
2	E	401	HEM	C2D-C1D-ND	2.14	112.45	109.88
2	D	401	HEM	CHD-C1D-ND	2.13	126.75	124.43
2	A	401	HEM	CBA-CAA-C2A	2.13	116.26	112.62
2	C	401	HEM	O1A-CGA-CBA	-2.12	116.28	123.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	HEM	C3C-C4C-NC	-2.11	106.96	110.94
2	B	401	HEM	CMC-C2C-C3C	2.11	128.62	124.68
2	F	401	HEM	CHA-C4D-C3D	-2.09	121.39	125.33
2	C	401	HEM	CHB-C1B-NB	2.07	126.94	124.38
2	C	401	HEM	CHD-C1D-ND	2.07	126.68	124.43
2	D	401	HEM	CAB-C3B-C2B	2.06	135.40	128.60
2	B	401	HEM	C3D-C4D-ND	2.06	112.46	110.17
2	C	401	HEM	C3C-C4C-NC	-2.05	107.07	110.94
2	E	401	HEM	CMA-C3A-C4A	-2.04	125.32	128.46
2	B	401	HEM	CBA-CAA-C2A	2.04	116.11	112.62
2	F	401	HEM	CHC-C4B-C3B	-2.03	121.46	124.57
2	D	401	HEM	CAD-C3D-C4D	2.02	128.18	124.66
2	F	401	HEM	CHD-C1D-ND	2.01	126.61	124.43
2	A	401	HEM	CAB-C3B-C4B	2.01	133.82	124.47

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	401	HEM	C2B-C3B-CAB-CBB
2	A	401	HEM	C2B-C3B-CAB-CBB
2	B	401	HEM	C2B-C3B-CAB-CBB
2	A	401	HEM	C4B-C3B-CAB-CBB
2	C	401	HEM	C4B-C3B-CAB-CBB
2	D	401	HEM	C4B-C3B-CAB-CBB
2	E	401	HEM	C4B-C3B-CAB-CBB
2	D	401	HEM	C2B-C3B-CAB-CBB
2	C	401	HEM	CAD-CBD-CGD-O2D
2	B	401	HEM	CAD-CBD-CGD-O2D
2	D	401	HEM	CAD-CBD-CGD-O2D
2	E	401	HEM	CAD-CBD-CGD-O2D
2	F	401	HEM	CAD-CBD-CGD-O2D
2	D	401	HEM	CAD-CBD-CGD-O1D
2	F	401	HEM	CAD-CBD-CGD-O1D
2	E	401	HEM	CAD-CBD-CGD-O1D
2	C	401	HEM	CAD-CBD-CGD-O1D
2	F	401	HEM	CAA-CBA-CGA-O2A
2	B	401	HEM	CAD-CBD-CGD-O1D
2	A	401	HEM	CAD-CBD-CGD-O2D
2	B	401	HEM	CAA-CBA-CGA-O2A
2	C	401	HEM	CAA-CBA-CGA-O1A
2	E	401	HEM	CAA-CBA-CGA-O1A

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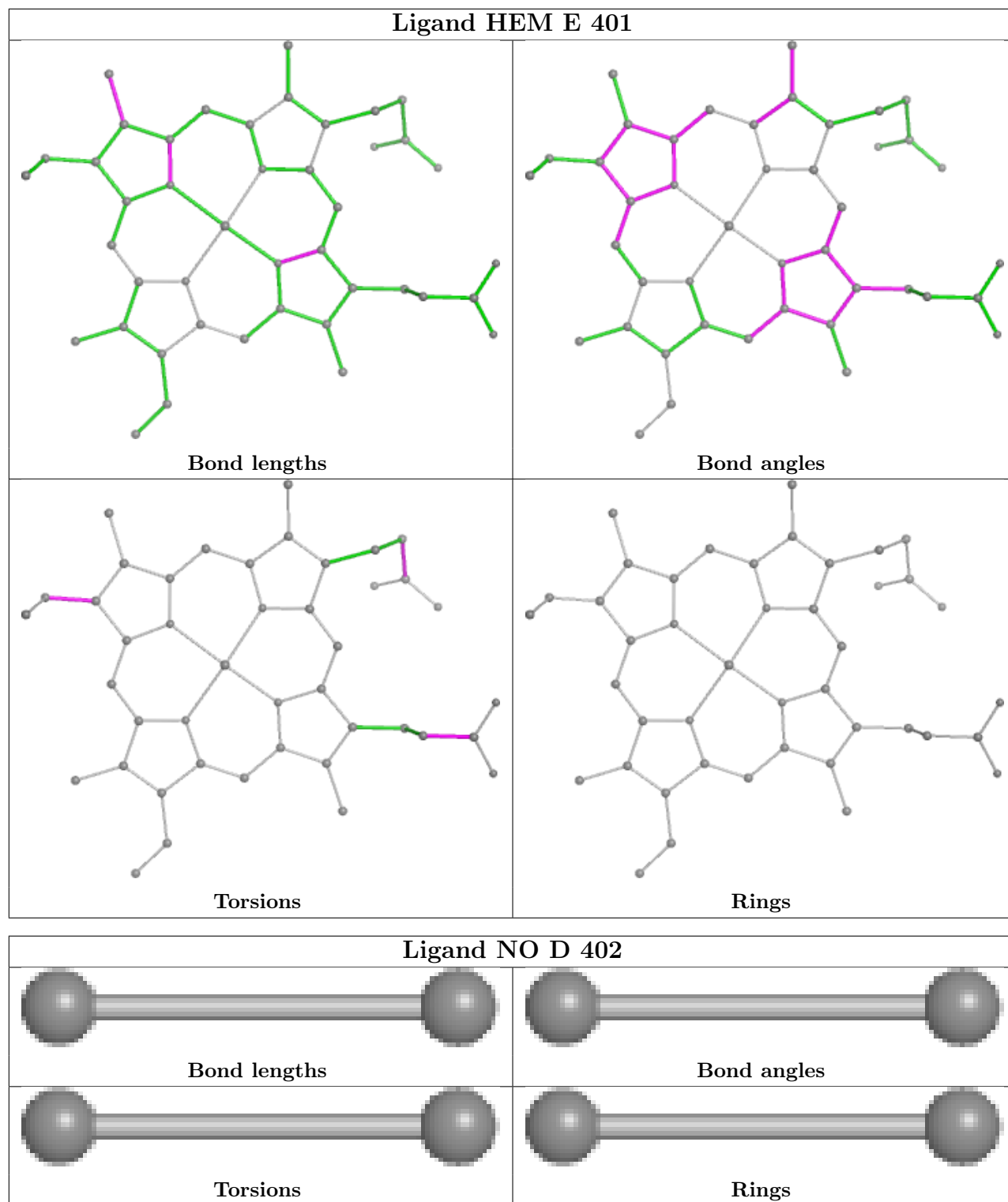
Mol	Chain	Res	Type	Atoms
2	A	401	HEM	CAA-CBA-CGA-O2A
2	A	401	HEM	CAD-CBD-CGD-O1D
2	E	401	HEM	CAA-CBA-CGA-O2A
2	B	401	HEM	CAA-CBA-CGA-O1A
2	D	401	HEM	CAA-CBA-CGA-O2A
2	C	401	HEM	CAA-CBA-CGA-O2A
2	F	401	HEM	CAA-CBA-CGA-O1A
2	D	401	HEM	CAA-CBA-CGA-O1A
2	A	401	HEM	CAA-CBA-CGA-O1A

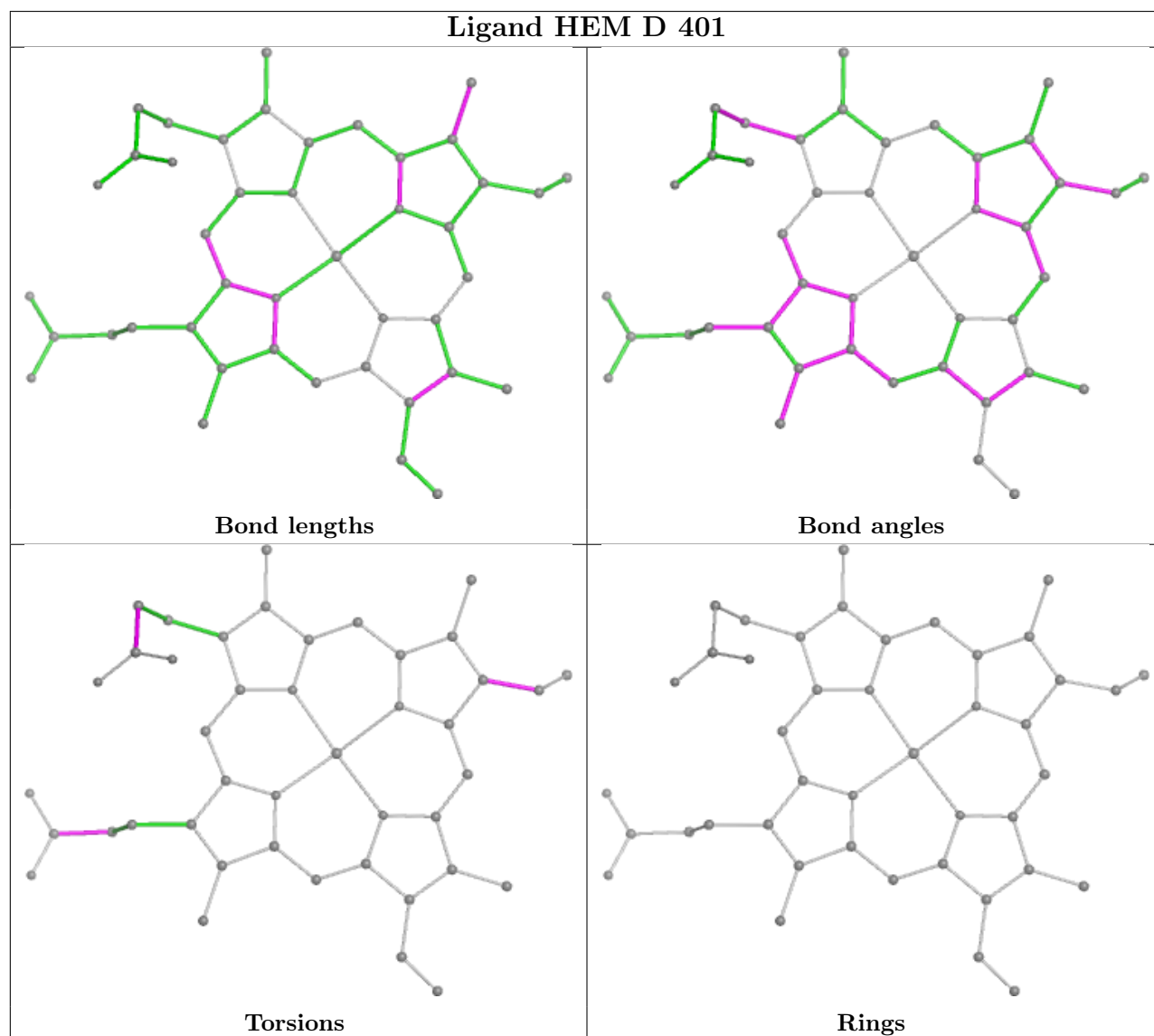
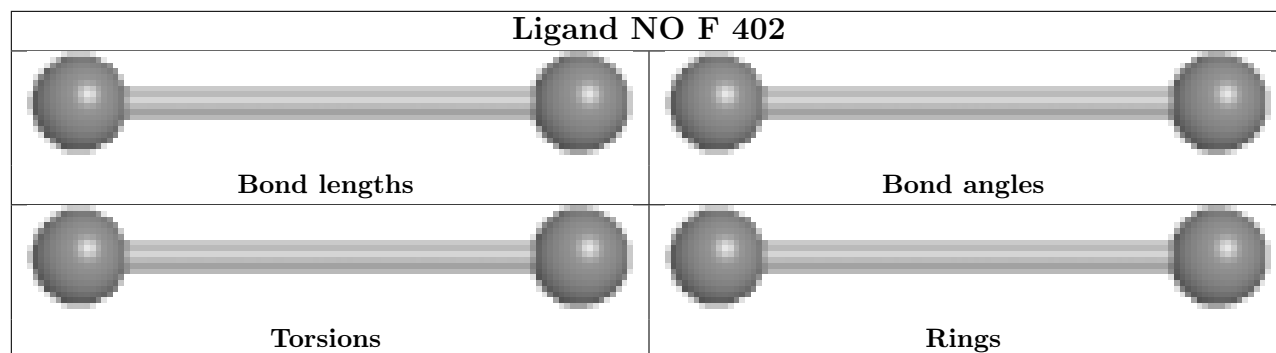
There are no ring outliers.

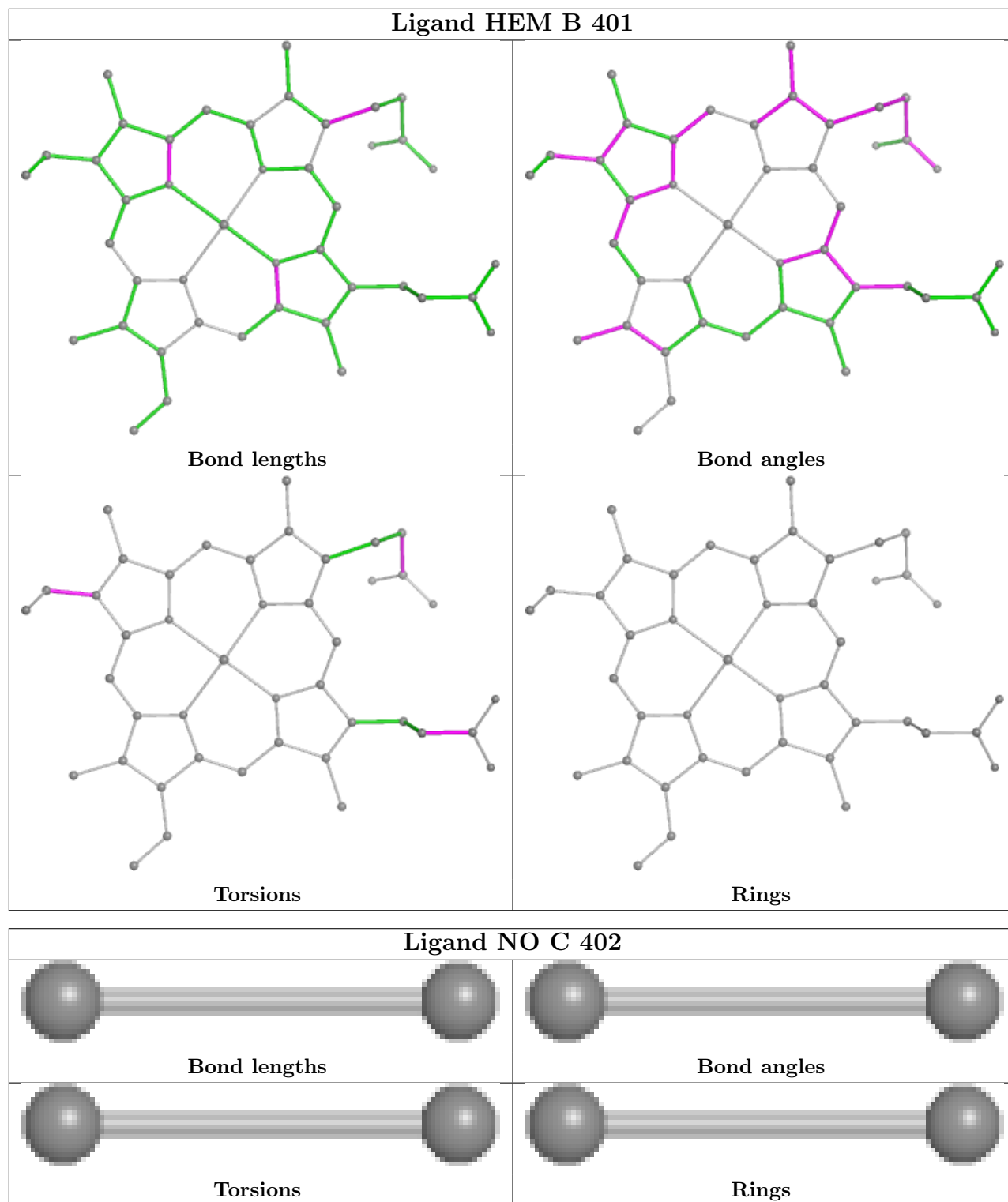
11 monomers are involved in 16 short contacts:

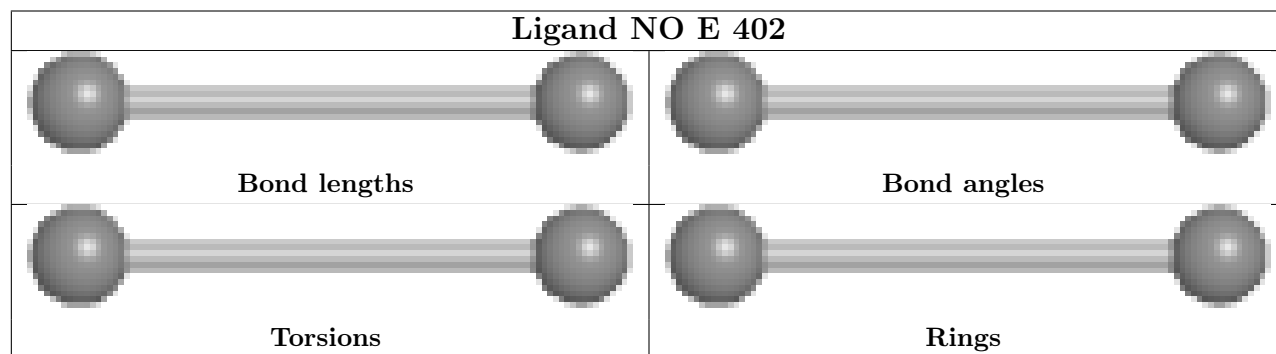
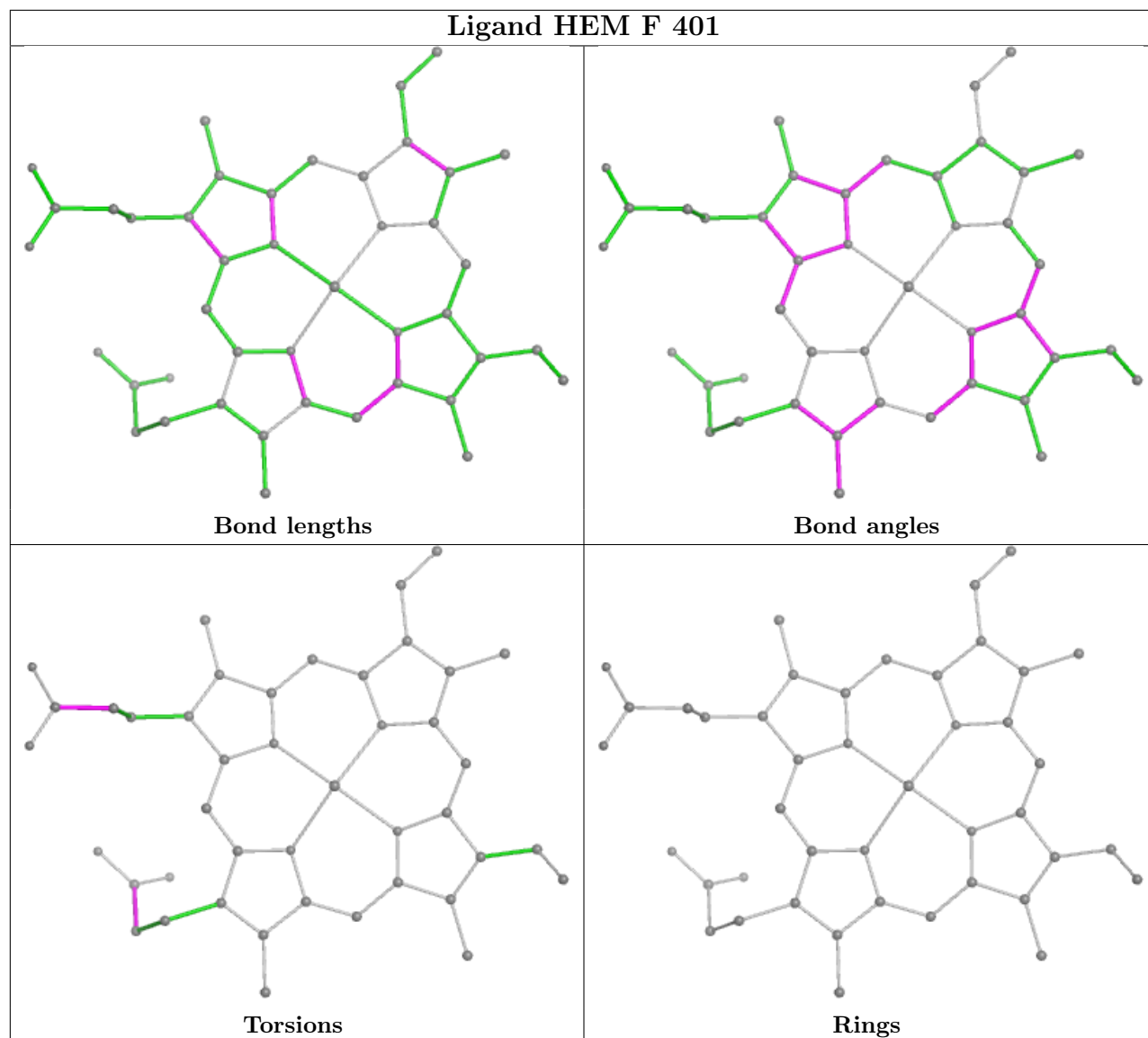
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	402	NO	1	0
3	F	402	NO	1	0
2	D	401	HEM	1	0
2	B	401	HEM	2	0
3	C	402	NO	3	0
2	F	401	HEM	1	0
3	E	402	NO	2	0
3	A	402	NO	1	0
3	B	402	NO	3	0
2	C	401	HEM	3	0
2	A	401	HEM	1	0





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







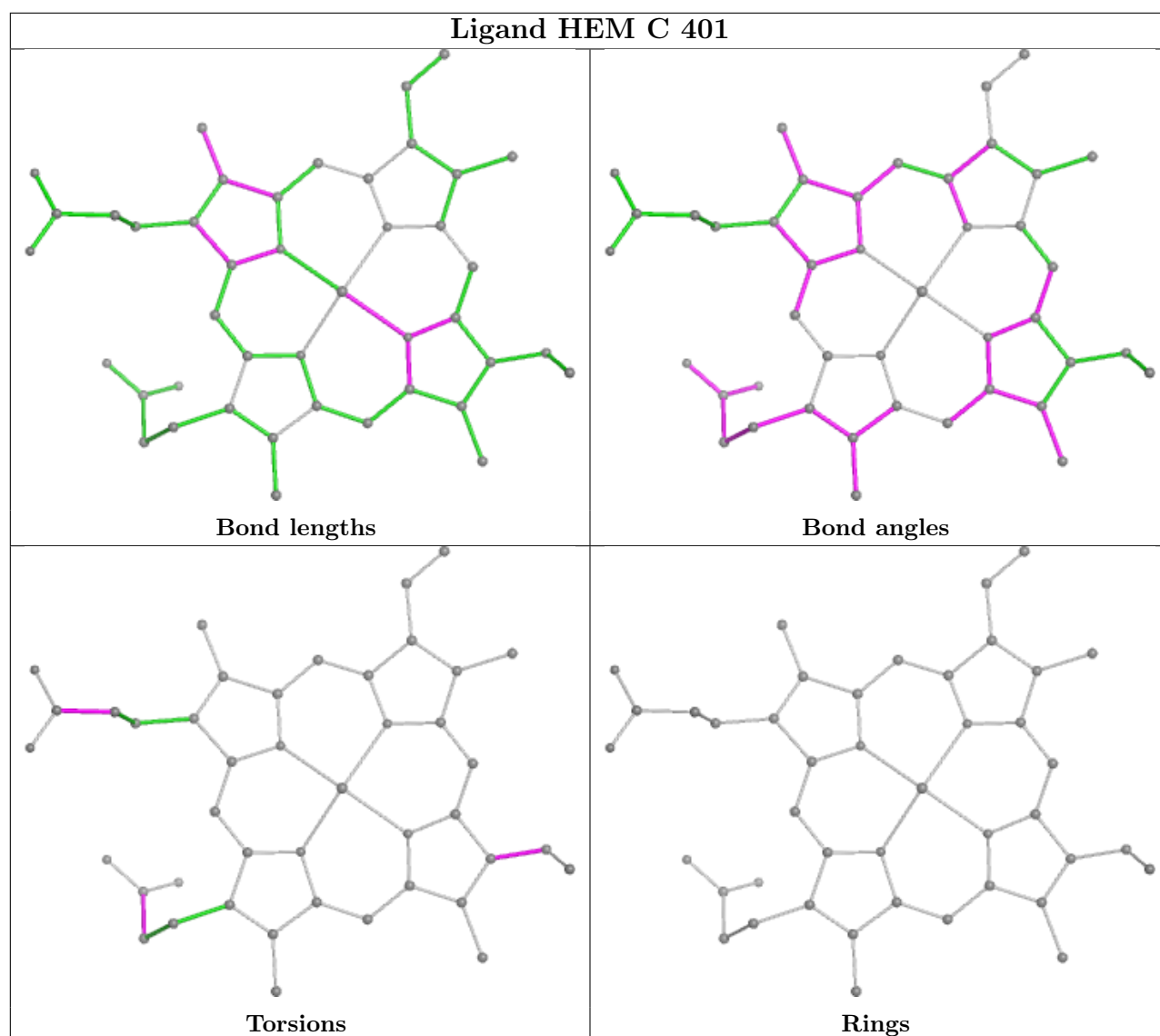


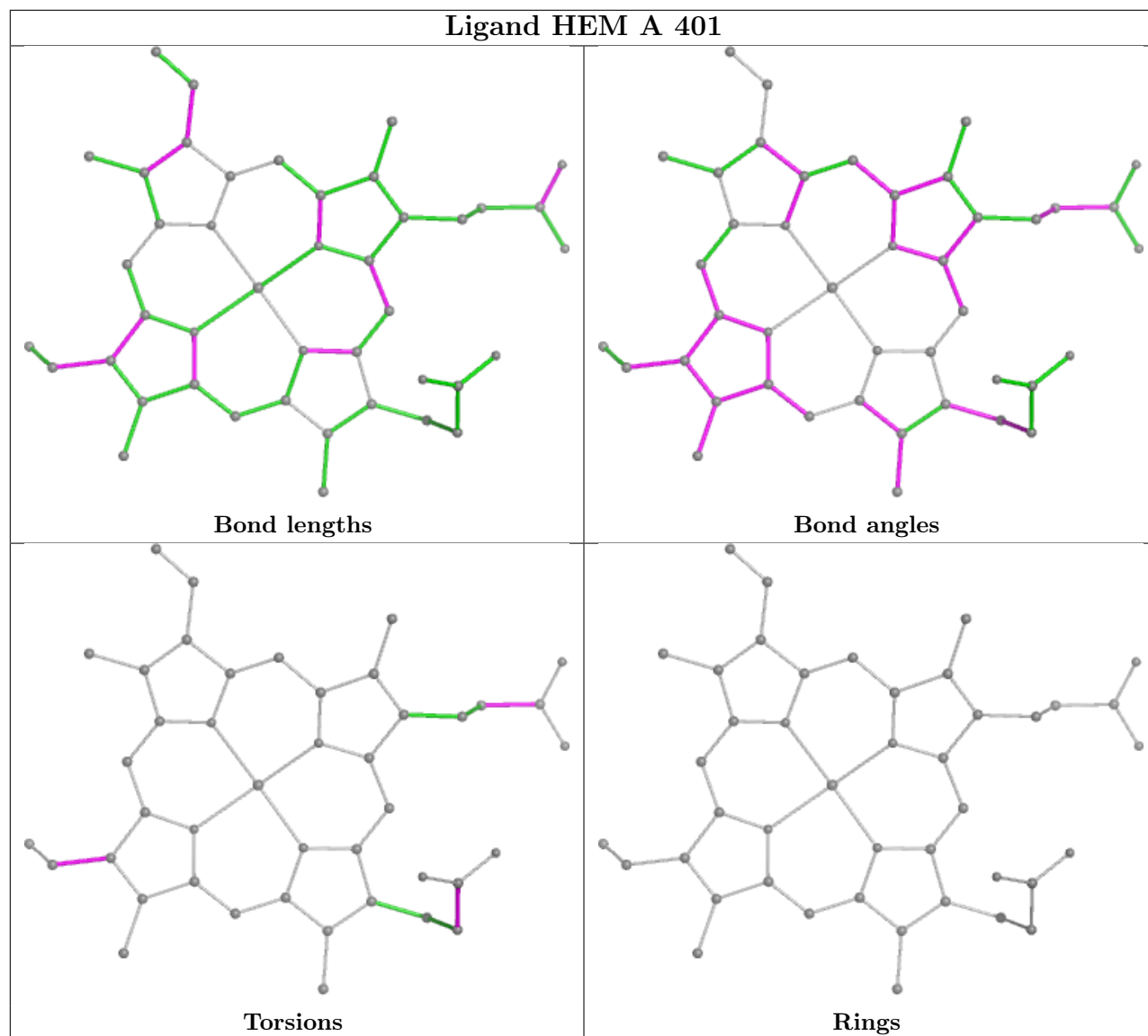




Ligand NO A 402			
 Bond lengths		 Bond angles	
 Torsions		 Rings	

Ligand NO B 402			
 Bond lengths		 Bond angles	
 Torsions		 Rings	





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	306/307 (99%)	-0.39	0 100 100	10, 19, 40, 74	4 (1%)
1	B	306/307 (99%)	-0.18	5 (1%) 70 73	13, 25, 52, 81	3 (0%)
1	C	306/307 (99%)	-0.19	2 (0%) 84 86	7, 24, 50, 102	6 (1%)
1	D	306/307 (99%)	-0.32	5 (1%) 70 73	11, 20, 46, 86	3 (0%)
1	E	306/307 (99%)	-0.31	4 (1%) 74 77	8, 21, 47, 74	9 (2%)
1	F	306/307 (99%)	-0.19	5 (1%) 70 73	6, 25, 50, 86	3 (0%)
All	All	1836/1842 (99%)	-0.26	21 (1%) 77 80	6, 22, 48, 102	28 (1%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	234	PRO	4.3
1	B	312	LEU	3.7
1	F	8	PRO	3.3
1	D	234	PRO	3.3
1	F	234	PRO	3.1
1	D	312	LEU	2.8
1	E	237	SER	2.8
1	D	8	PRO	2.7
1	F	282	ALA	2.6
1	C	312	LEU	2.5
1	B	234	PRO	2.5
1	E	55[A]	GLN	2.4
1	E	312	LEU	2.4
1	E	234	PRO	2.3
1	B	219	VAL	2.3
1	F	313	SER	2.2
1	B	55	GLN	2.1
1	D	311	ASP	2.1
1	F	76	SER	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	8	PRO	2.0
1	D	7	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

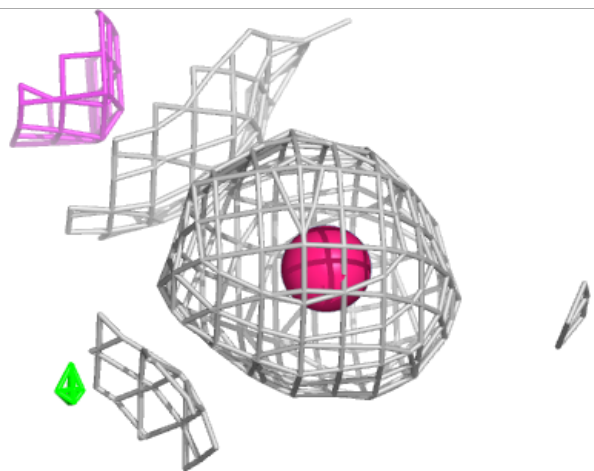
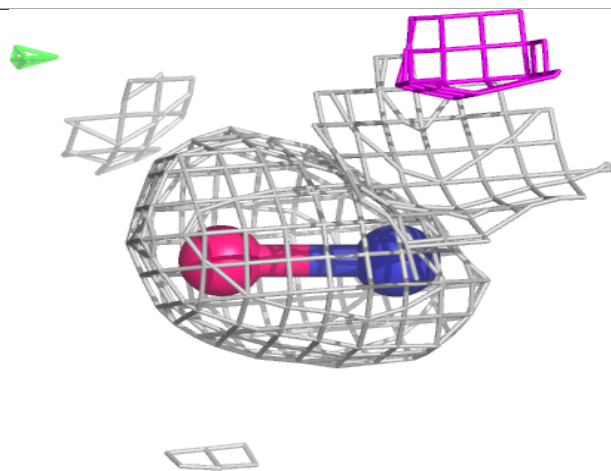
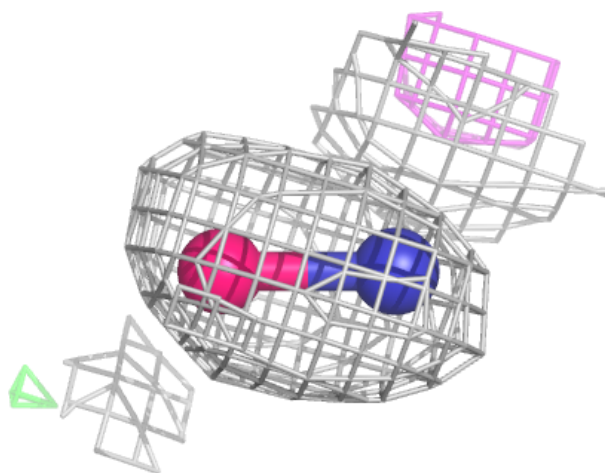
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NO	C	402	2/2	0.95	0.20	42,42,42,43	0
3	NO	F	402	2/2	0.95	0.26	43,43,43,51	0
3	NO	A	402	2/2	0.96	0.20	37,37,37,39	0
3	NO	E	402	2/2	0.97	0.13	34,34,34,44	0
3	NO	D	402	2/2	0.98	0.12	37,37,37,40	0
3	NO	B	402	2/2	0.98	0.20	38,38,38,45	0
2	HEM	F	401	43/43	0.98	0.06	17,20,25,31	0
2	HEM	C	401	43/43	0.99	0.05	16,20,25,29	0
2	HEM	D	401	43/43	0.99	0.05	13,15,21,24	0
2	HEM	E	401	43/43	0.99	0.04	14,16,20,23	0
2	HEM	A	401	43/43	0.99	0.05	15,17,21,23	0
2	HEM	B	401	43/43	0.99	0.04	14,18,23,25	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.

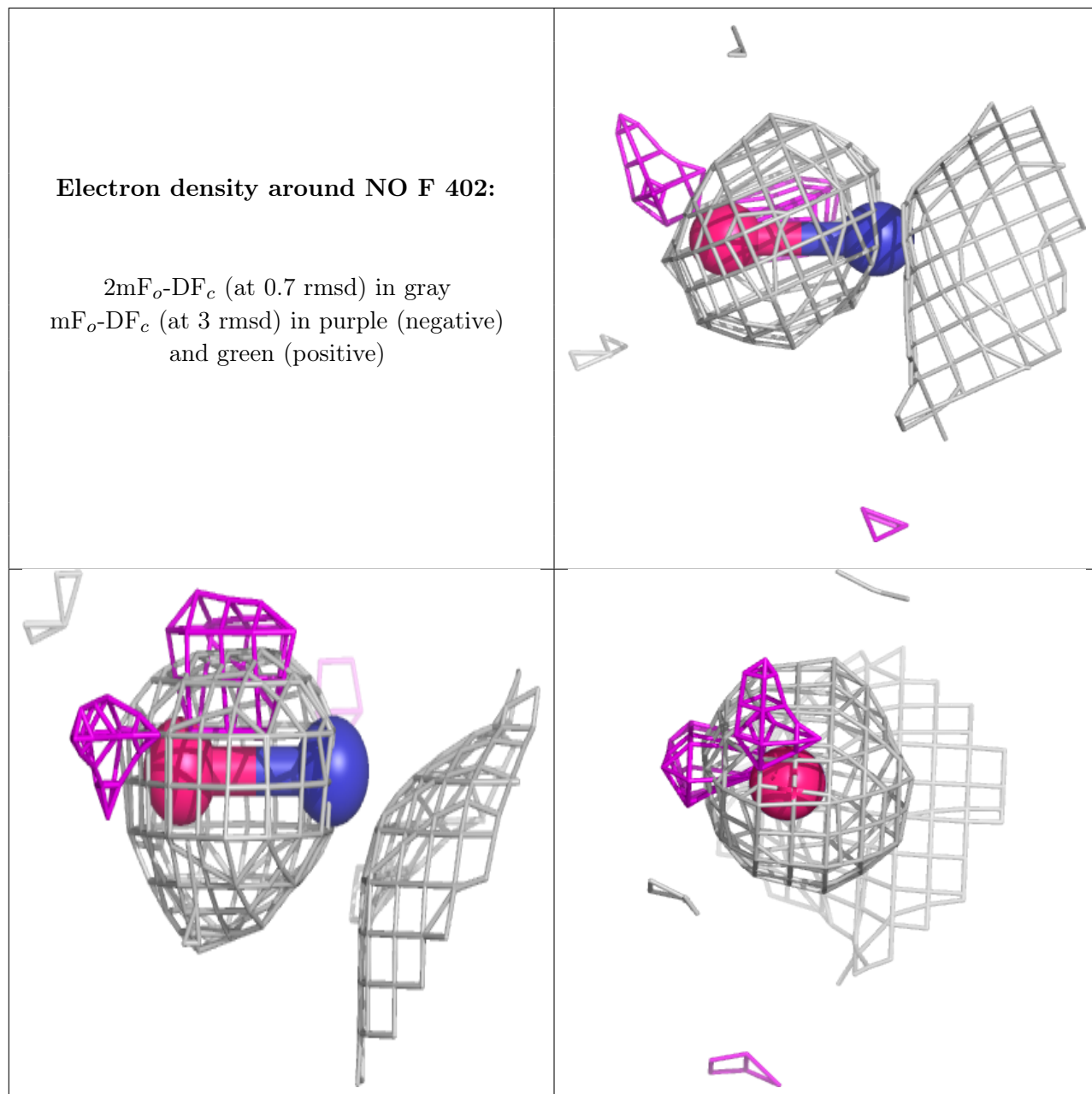
Electron density around NO C 402:

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



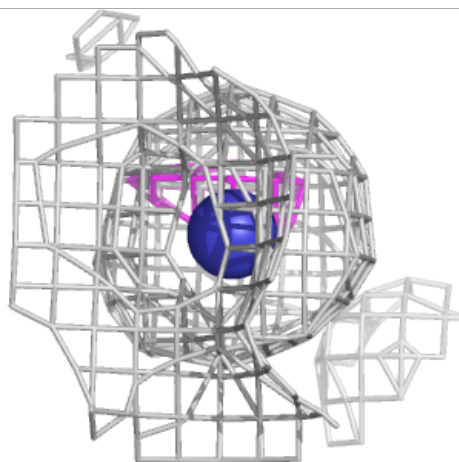
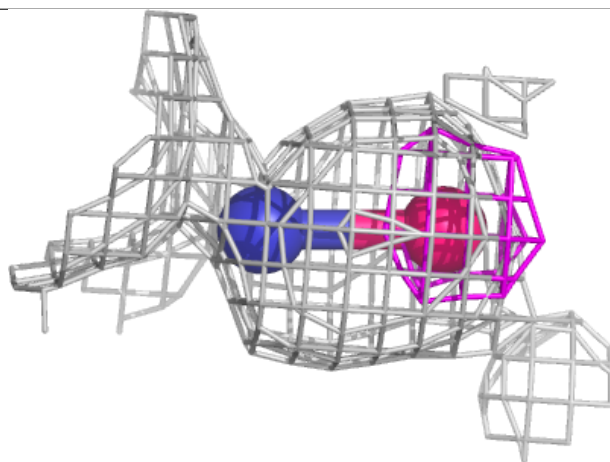
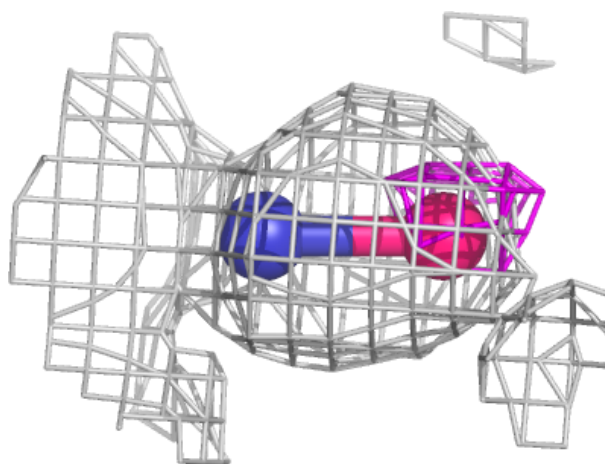
Electron density around NO F 402:

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



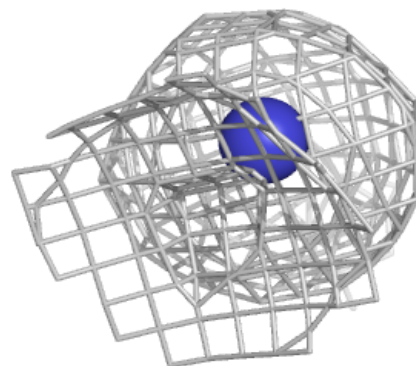
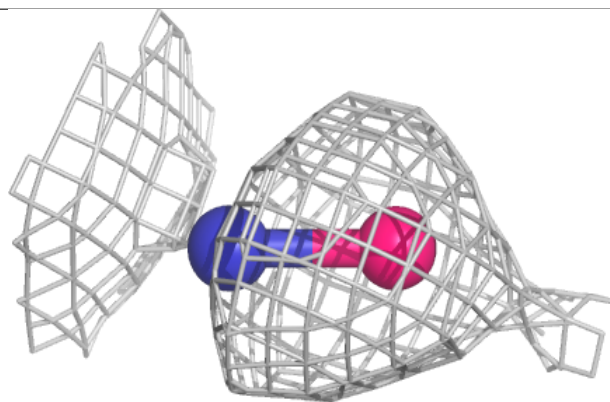
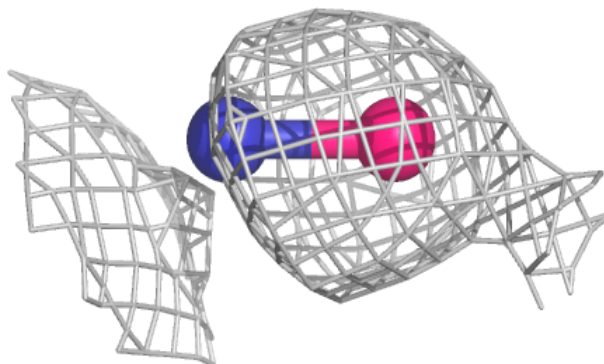
Electron density around NO A 402:

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

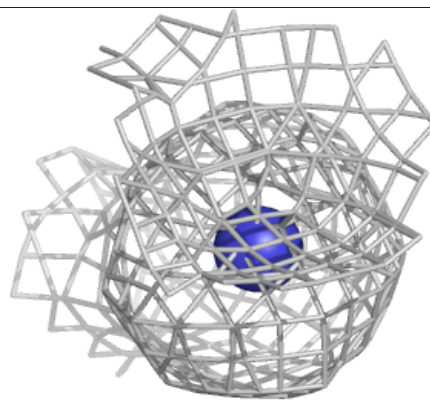
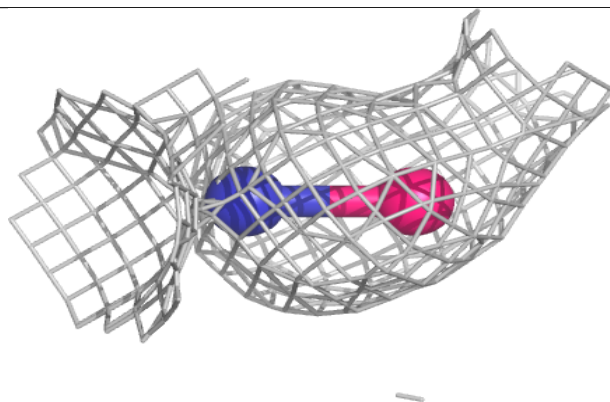
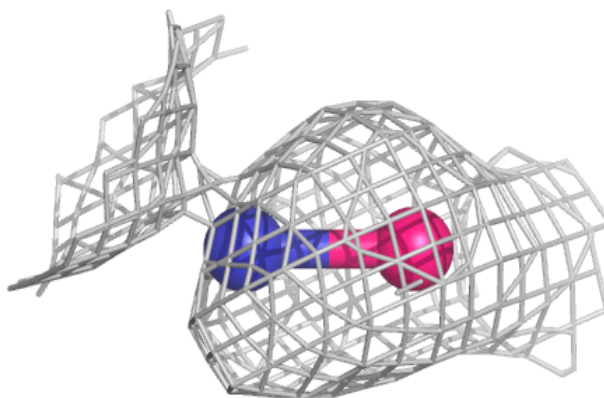


Electron density around NO E 402:

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

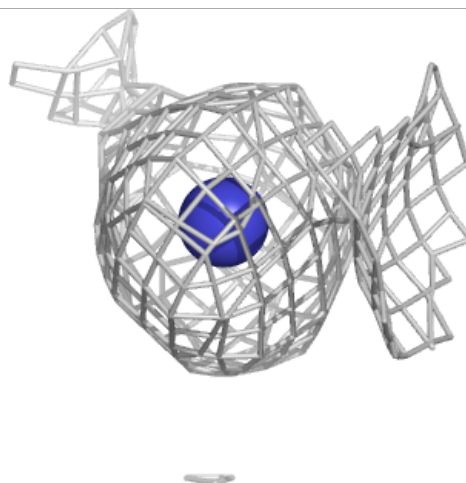
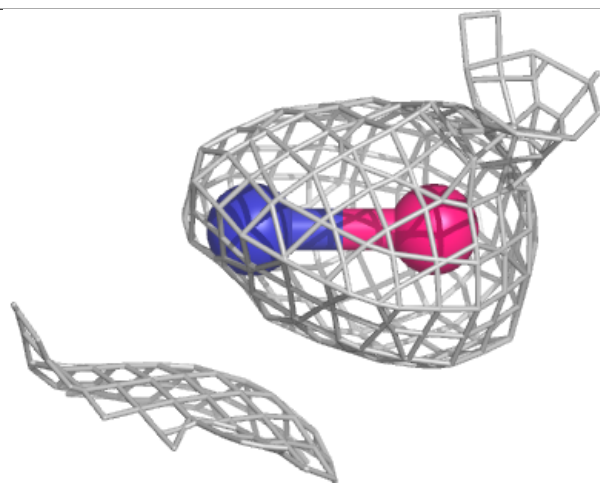
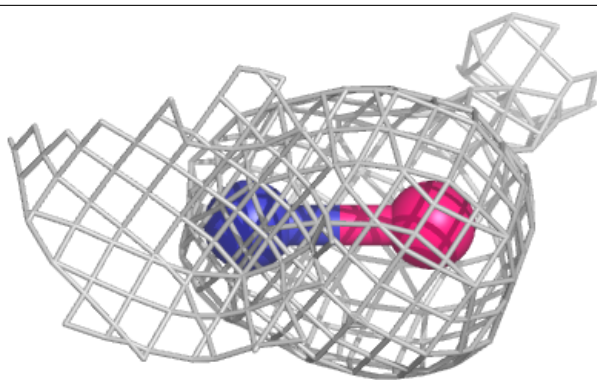
**Electron density around NO D 402:**

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



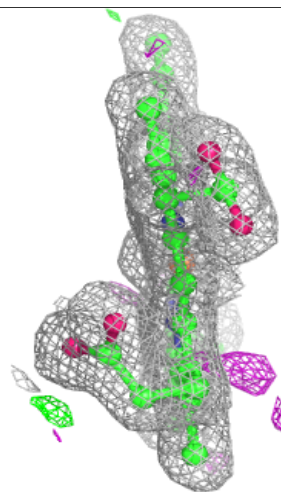
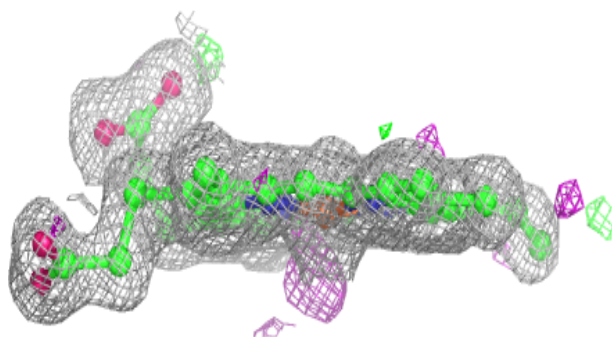
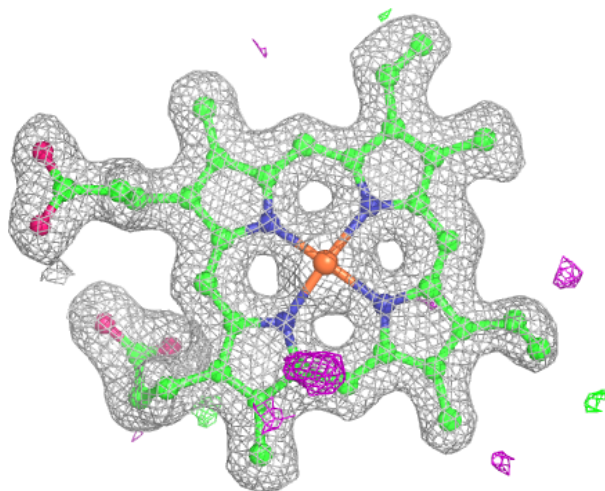
Electron density around NO B 402:

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and green (positive)



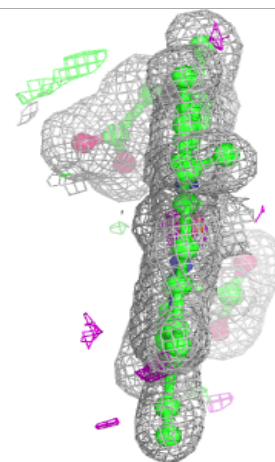
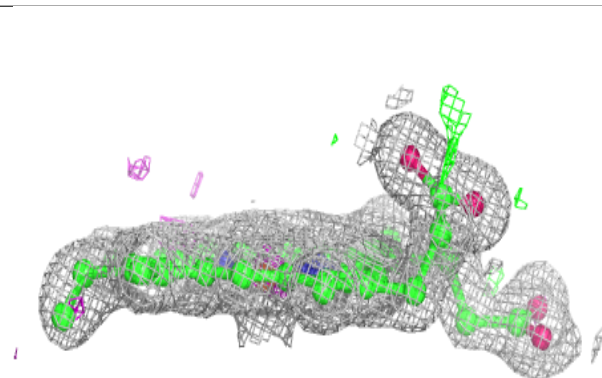
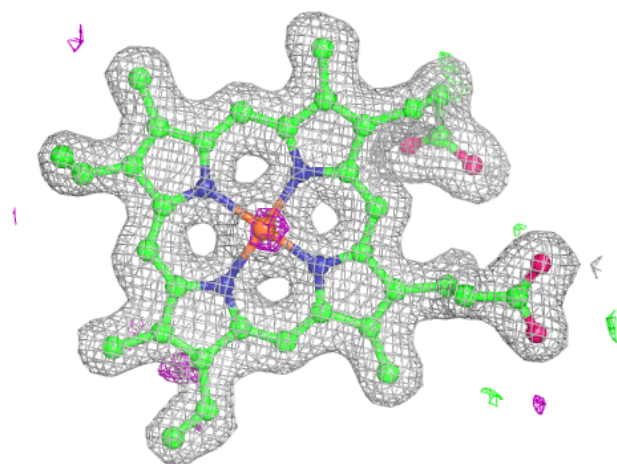
Electron density around HEM F 401:

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



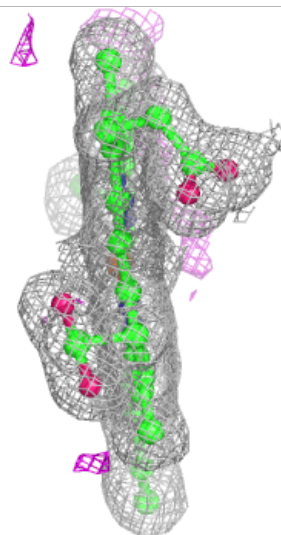
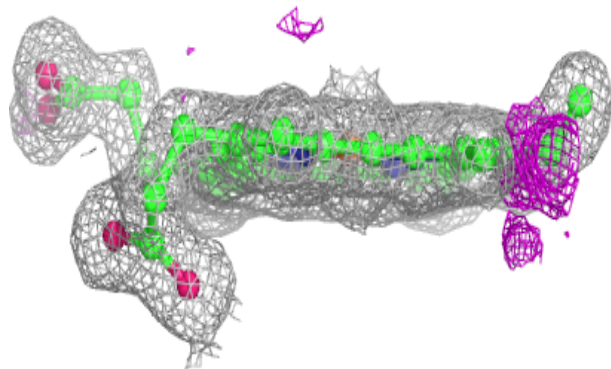
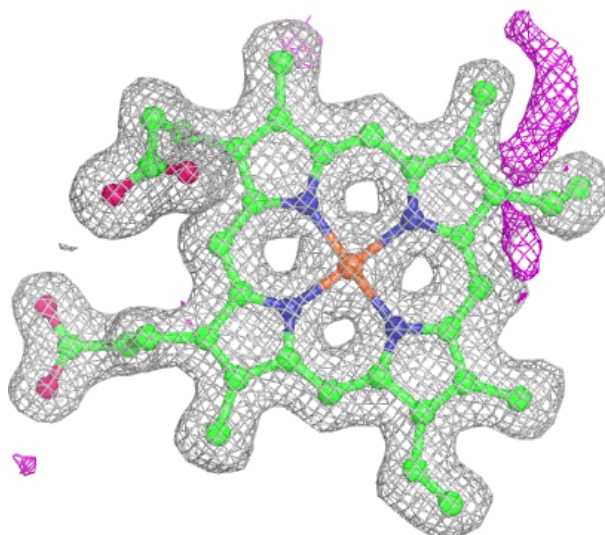
Electron density around HEM C 401:

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and green (positive)



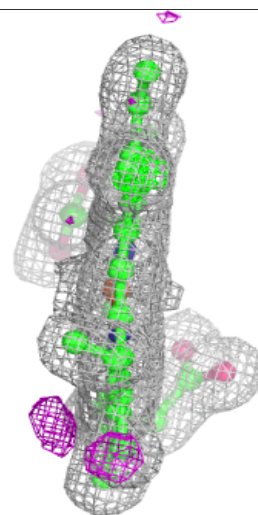
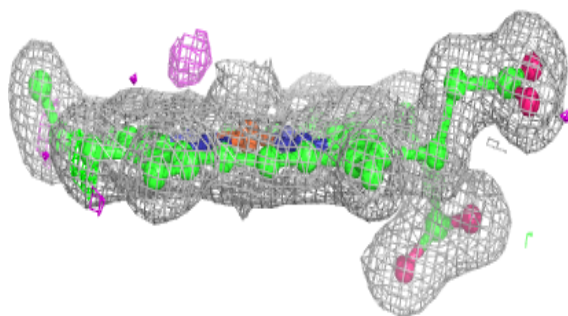
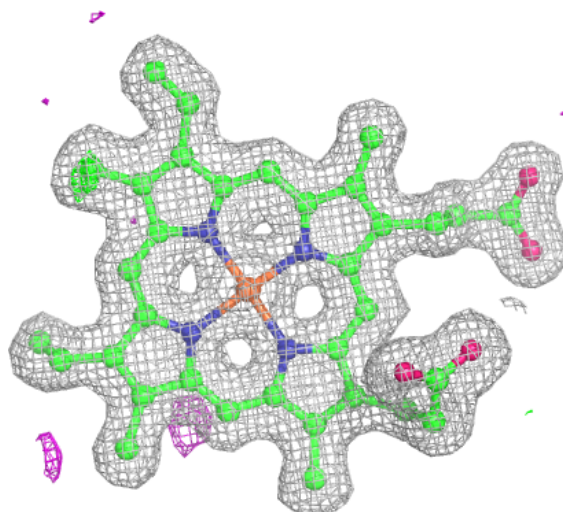
Electron density around HEM D 401:

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



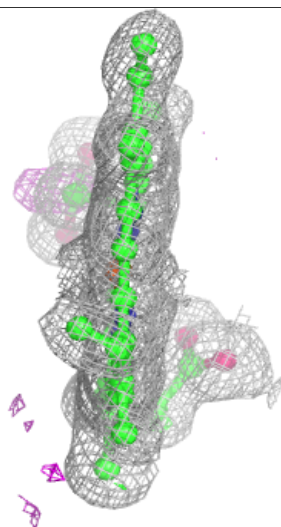
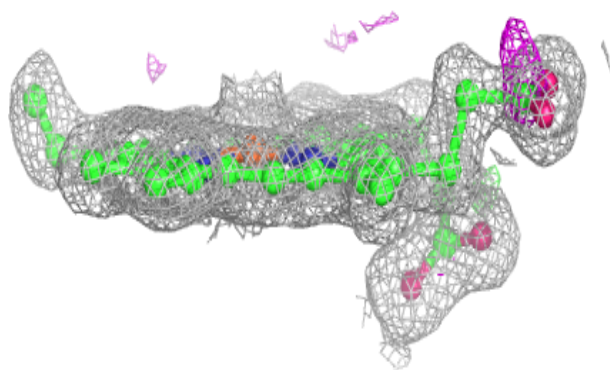
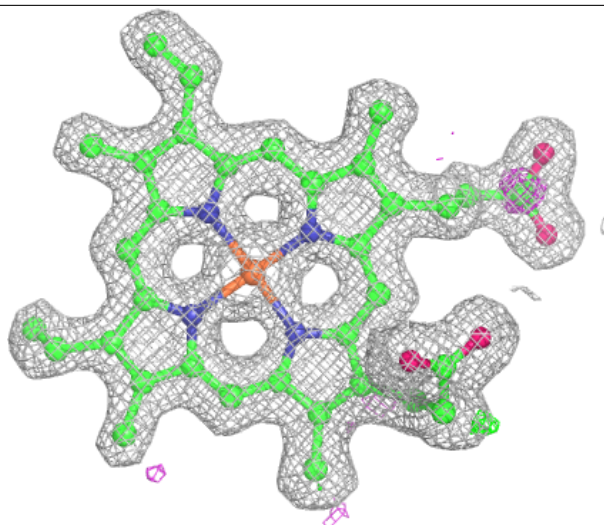
Electron density around HEM E 401:

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



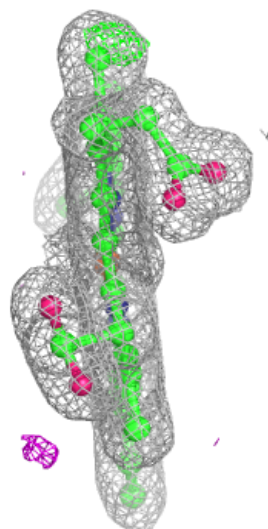
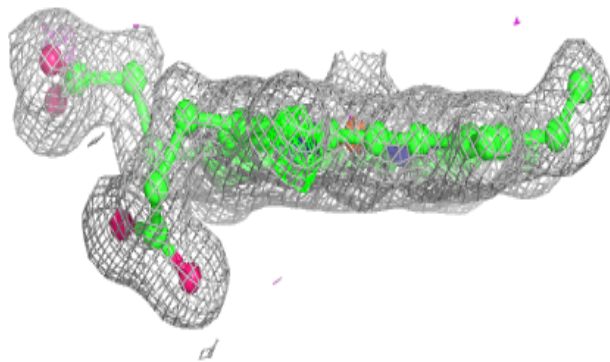
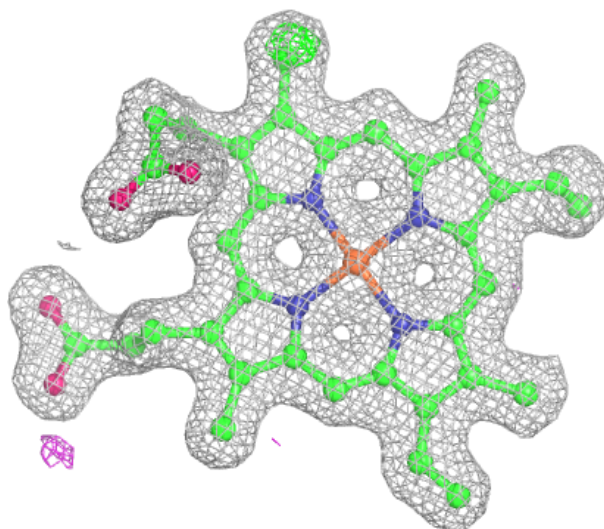
Electron density around HEM A 401:

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



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)



6.5 Other polymers [i](#)

There are no such residues in this entry.