



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:35 AM UTC

PDB ID : 1DW2 / pdb_00001dw2
Title : STRUCTURE OF THE NITRIC OXIDE COMPLEX OF REDUCED SHP,
AN OXYGEN BINDING CYTOCHROME C
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Beeumen, J.J.
Deposited on : 2000-01-24
Resolution : 2.20 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 Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

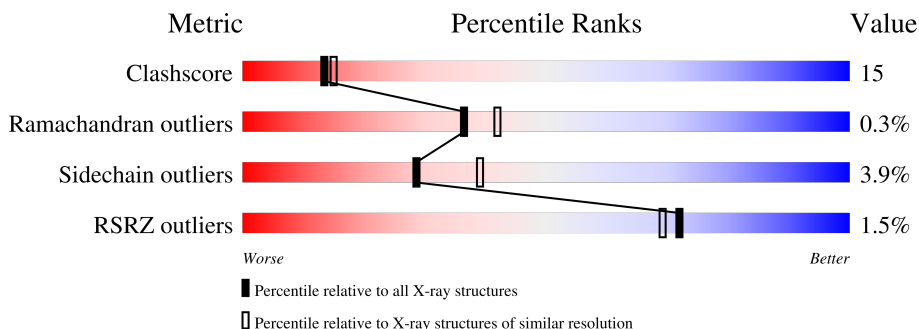
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	112	2% 54% 38% 6%
1	B	112	3% 62% 29% 7%
1	C	112	55% 35% 8%

2 Entry composition [i](#)

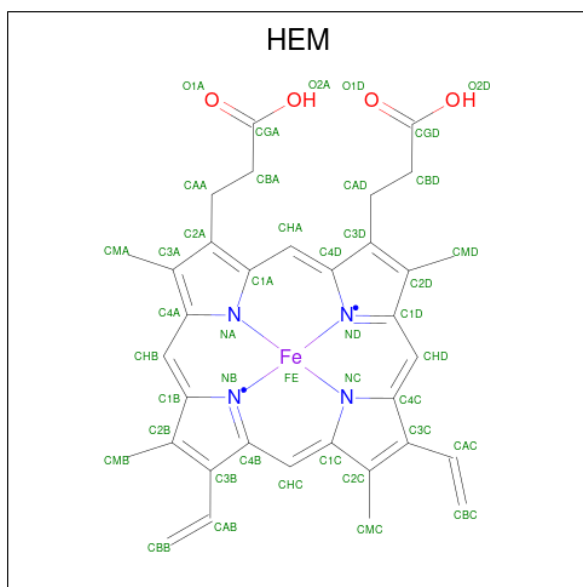
There are 4 unique types of molecules in this entry. The entry contains 2807 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 CYTOCHROME C.

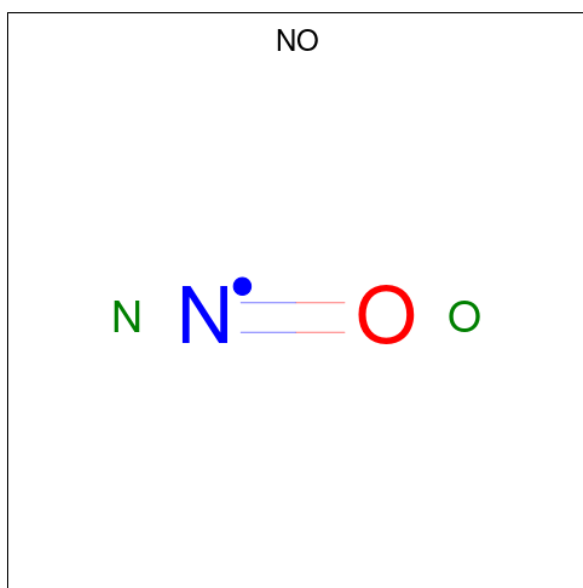
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	112	Total	C	N	O	S	0	1	0
			802	488	149	161	4			
1	B	112	Total	C	N	O	S	0	0	0
			794	483	143	164	4			
1	C	112	Total	C	N	O	S	0	1	0
			802	486	150	162	4			

- 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	0	1
			45	34	1	4	6		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is NITRIC OXIDE (CCD ID: NO) (formula: NO).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	N	O	0	0
			2	1	1		
3	B	1	Total	N	O	0	0
			2	1	1		
3	C	1	Total	N	O	0	0
			2	1	1		

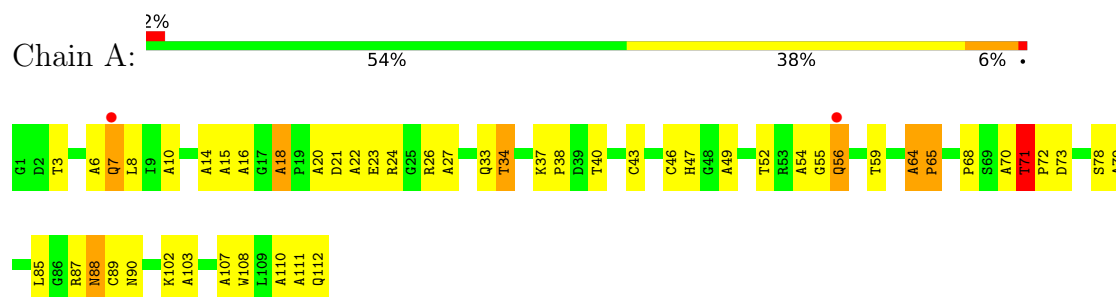
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	82	Total	O	0	0
			82	82		
4	B	81	Total	O	0	0
			81	81		
4	C	109	Total	O	0	0
			109	109		

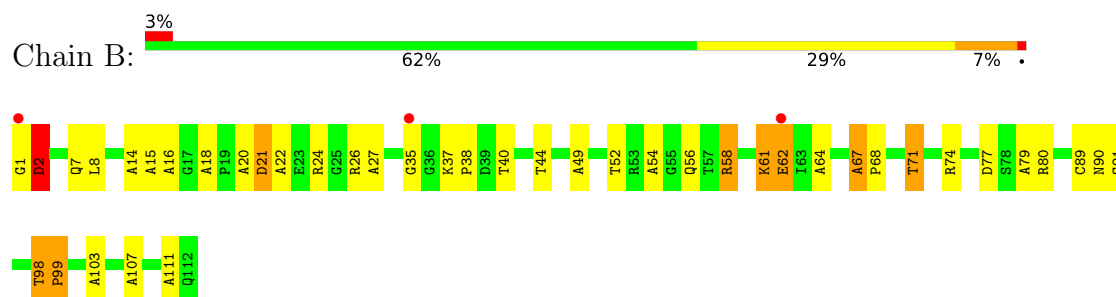
3 Residue-property plots [i](#)

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.

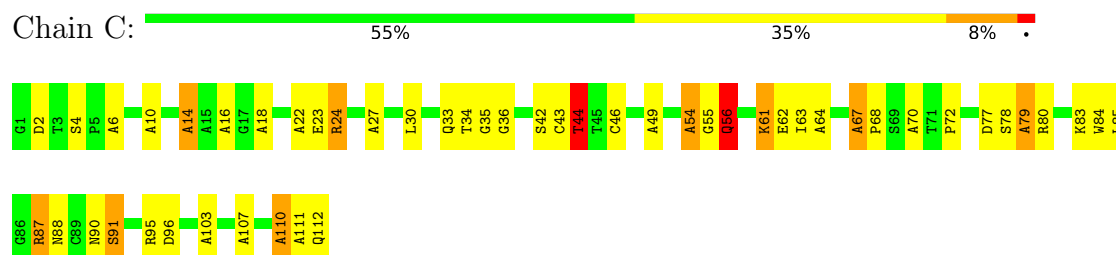
• Molecule 1: CYTOCHROME C



• Molecule 1: CYTOCHROME C



• Molecule 1: CYTOCHROME C



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	83.34Å 103.50Å 113.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 20.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.20) 83.4 (20.00-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 2.09Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.167 , 0.229 0.153 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	28.8	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 62.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2807	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.55% 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: NO, HEM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	2.69	29/824 (3.5%)	2.45	57/1121 (5.1%)
1	B	2.54	31/809 (3.8%)	2.41	46/1105 (4.2%)
1	C	2.85	32/824 (3.9%)	2.41	46/1123 (4.1%)
All	All	2.70	92/2457 (3.7%)	2.42	149/3349 (4.4%)

All (92) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	34	THR	CA-CB	37.13	2.10	1.53
1	B	61	LYS	CA-C	32.25	1.94	1.52
1	C	54	ALA	CA-CB	32.05	2.06	1.53
1	C	79	ALA	CA-CB	31.34	2.03	1.53
1	A	18	ALA	CA-CB	23.44	1.90	1.53
1	A	107	ALA	C-O	-21.45	0.99	1.24
1	C	27	ALA	C-O	19.44	1.47	1.24
1	B	58	ARG	C-O	18.84	1.48	1.24
1	C	54	ALA	C-O	17.59	1.46	1.23
1	C	61	LYS	CA-CB	16.68	1.79	1.53
1	B	15	ALA	C-O	16.34	1.45	1.24
1	B	58	ARG	CA-C	-16.31	1.29	1.52
1	A	64	ALA	CA-CB	16.10	1.77	1.53
1	C	111	ALA	C-O	-16.04	1.02	1.24
1	B	58	ARG	CA-CB	15.75	1.80	1.53
1	C	49	ALA	CA-C	13.96	1.71	1.52
1	C	14	ALA	CA-CB	13.71	1.75	1.53
1	B	27	ALA	C-O	13.33	1.39	1.24
1	B	107	ALA	CA-CB	13.00	1.73	1.53
1	A	6	ALA	CA-C	12.86	1.69	1.52
1	B	15	ALA	CA-C	-12.65	1.35	1.52
1	C	6	ALA	CA-CB	12.47	1.72	1.53
1	C	14	ALA	C-O	12.37	1.38	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	14	ALA	CA-C	-12.35	1.36	1.52
1	C	111	ALA	CA-C	12.20	1.70	1.52
1	A	14	ALA	CA-C	11.65	1.68	1.52
1	A	54	ALA	CA-CB	11.52	1.73	1.53
1	A	34	THR	CA-C	-11.45	1.37	1.52
1	A	49	ALA	C-O	11.31	1.39	1.24
1	B	16	ALA	CA-C	-11.08	1.37	1.52
1	B	54	ALA	CA-CB	11.05	1.71	1.53
1	B	27	ALA	CA-CB	10.98	1.70	1.53
1	C	110	ALA	CA-CB	10.73	1.72	1.53
1	A	6	ALA	C-O	-10.71	1.11	1.24
1	B	18	ALA	CA-C	10.62	1.63	1.52
1	C	110	ALA	CA-C	-10.51	1.38	1.52
1	A	110	ALA	C-O	-10.50	1.10	1.24
1	C	27	ALA	CA-C	-10.45	1.39	1.52
1	B	22	ALA	C-O	10.41	1.37	1.24
1	A	70	ALA	CA-C	-10.33	1.39	1.52
1	A	107	ALA	CA-C	10.32	1.66	1.52
1	A	14	ALA	C-O	-9.79	1.12	1.24
1	C	16	ALA	CA-C	-9.66	1.39	1.52
1	B	22	ALA	CA-CB	9.20	1.68	1.53
1	A	22	ALA	CA-CB	-9.04	1.38	1.53
1	B	64	ALA	CA-CB	9.00	1.66	1.53
1	C	67	ALA	CA-CB	-8.97	1.36	1.53
1	B	103	ALA	C-O	-8.89	1.13	1.24
1	A	103	ALA	CA-C	8.75	1.64	1.52
1	C	79	ALA	CA-C	8.74	1.65	1.52
1	C	18	ALA	CA-CB	8.74	1.67	1.54
1	B	79	ALA	CA-CB	8.61	1.67	1.53
1	C	79	ALA	C-O	-8.47	1.13	1.24
1	A	10	ALA	C-O	-8.14	1.14	1.24
1	A	110	ALA	CA-CB	8.08	1.67	1.53
1	A	49	ALA	CA-C	-8.08	1.41	1.52
1	C	18	ALA	C-O	-8.04	1.14	1.24
1	A	18	ALA	CA-C	-8.04	1.42	1.52
1	A	16	ALA	C-O	-7.80	1.14	1.24
1	B	14	ALA	C-O	-7.75	1.15	1.24
1	B	16	ALA	C-O	7.69	1.34	1.24
1	C	110	ALA	C-O	7.67	1.33	1.24
1	C	64	ALA	CA-C	-7.54	1.45	1.53
1	C	103	ALA	CA-C	7.40	1.62	1.52
1	A	16	ALA	CA-CB	7.16	1.65	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	64	ALA	CA-C	7.14	1.61	1.53
1	C	70	ALA	CA-C	-7.14	1.43	1.52
1	B	14	ALA	CA-C	6.98	1.61	1.52
1	A	10	ALA	CA-C	6.91	1.62	1.52
1	A	64	ALA	C-O	6.88	1.32	1.23
1	A	15	ALA	CA-CB	6.78	1.64	1.53
1	C	103	ALA	CA-CB	6.63	1.63	1.53
1	B	20	ALA	CA-CB	-6.56	1.42	1.53
1	C	70	ALA	CA-CB	6.49	1.63	1.53
1	B	49	ALA	CA-CB	-6.38	1.43	1.53
1	B	107	ALA	C-O	-6.13	1.17	1.24
1	B	18	ALA	C-O	-6.08	1.17	1.24
1	C	6	ALA	CA-C	-5.96	1.45	1.52
1	A	10	ALA	CA-CB	5.88	1.63	1.53
1	B	22	ALA	CA-C	-5.87	1.44	1.52
1	B	61	LYS	C-O	-5.87	1.16	1.23
1	B	27	ALA	CA-C	-5.83	1.45	1.52
1	A	34	THR	C-O	5.75	1.31	1.24
1	B	79	ALA	C-O	5.64	1.30	1.24
1	A	111	ALA	CA-CB	5.62	1.63	1.53
1	C	6	ALA	C-O	5.58	1.30	1.24
1	B	54	ALA	C-O	5.54	1.30	1.23
1	C	22	ALA	CA-CB	5.48	1.62	1.53
1	C	54	ALA	CA-C	-5.37	1.45	1.52
1	C	96	ASP	CG-OD2	5.11	1.35	1.25
1	B	103	ALA	CA-C	5.08	1.59	1.52
1	A	23	GLU	CA-CB	5.07	1.61	1.53

All (149) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	61	LYS	CA-C-O	19.69	143.49	121.05
1	A	34	THR	CB-CA-C	15.27	136.56	109.99
1	C	54	ALA	N-CA-CB	13.94	130.25	110.17
1	A	34	THR	CA-C-O	-13.32	103.46	119.18
1	C	54	ALA	O-C-N	-13.08	107.42	122.86
1	A	107	ALA	O-C-N	12.49	135.36	122.12
1	A	70	ALA	CA-C-O	-12.30	104.90	119.78
1	C	27	ALA	O-C-N	-12.29	108.14	122.15
1	B	58	ARG	O-C-N	-12.28	105.98	122.43
1	A	56	GLN	CB-CG-CD	12.16	133.27	112.60
1	C	49	ALA	N-CA-CB	11.89	129.12	110.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	54	ALA	CA-C-O	11.68	134.29	121.56
1	B	58	ARG	N-CA-CB	-11.39	91.48	110.39
1	C	49	ALA	CB-CA-C	-11.03	88.55	109.72
1	B	2	ASP	CA-CB-CG	10.88	123.48	112.60
1	C	88	ASN	CA-CB-CG	-10.76	101.84	112.60
1	A	7	GLN	CB-CG-CD	10.58	130.58	112.60
1	B	15	ALA	O-C-N	-10.10	108.53	122.46
1	C	79	ALA	N-CA-CB	10.06	126.82	109.72
1	C	111	ALA	O-C-N	9.94	136.08	122.37
1	B	62	GLU	CA-C-O	-9.54	106.87	120.51
1	A	7	GLN	OE1-CD-NE2	-9.50	113.10	122.60
1	C	14	ALA	O-C-N	-9.38	111.45	122.15
1	A	64	ALA	CA-C-O	9.18	131.07	120.69
1	A	111	ALA	CA-C-O	-9.13	108.23	119.28
1	B	90	ASN	CA-CB-CG	-9.04	103.56	112.60
1	A	88	ASN	CA-CB-CG	-8.96	103.64	112.60
1	B	26	ARG	NE-CZ-NH2	8.81	127.12	119.20
1	B	27	ALA	O-C-N	-8.70	112.89	122.12
1	C	34	THR	CA-C-O	-8.22	109.89	119.59
1	A	7	GLN	CG-CD-NE2	8.19	128.68	116.40
1	A	6	ALA	O-C-N	8.07	130.68	122.12
1	B	71	THR	CA-CB-OG1	8.03	121.65	109.60
1	B	111	ALA	CA-C-O	-7.99	110.21	119.56
1	B	62	GLU	CB-CG-CD	7.97	126.14	112.60
1	A	34	THR	N-CA-CB	-7.92	99.01	110.80
1	B	26	ARG	CD-NE-CZ	7.90	135.46	124.40
1	C	90	ASN	CA-C-O	-7.86	112.57	120.82
1	B	16	ALA	O-C-N	-7.82	111.30	122.41
1	B	58	ARG	CB-CA-C	7.74	126.49	110.31
1	B	14	ALA	O-C-N	7.74	130.32	122.12
1	C	44	THR	N-CA-CB	-7.67	97.29	110.32
1	A	71	THR	CB-CA-C	7.55	120.71	110.13
1	A	90	ASN	CA-C-O	-7.50	112.97	120.70
1	B	14	ALA	CA-C-O	-7.50	112.60	120.55
1	A	103	ALA	N-CA-CB	7.40	120.73	110.01
1	B	15	ALA	CA-C-O	-7.38	110.16	119.31
1	C	95	ARG	CA-C-O	-7.23	113.54	121.28
1	A	49	ALA	O-C-N	-7.22	112.40	122.37
1	B	89	CYS	N-CA-C	-7.17	103.40	111.14
1	B	26	ARG	NE-CZ-NH1	-7.04	114.46	121.50
1	B	61	LYS	N-CA-CB	6.96	120.56	110.06
1	B	62	GLU	CA-C-N	-6.96	114.03	123.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	62	GLU	C-N-CA	-6.96	114.03	123.14
1	A	70	ALA	N-CA-CB	-6.92	99.84	110.42
1	C	110	ALA	O-C-N	-6.91	113.17	122.43
1	C	27	ALA	CA-C-N	6.85	129.76	120.44
1	C	27	ALA	C-N-CA	6.85	129.76	120.44
1	A	64	ALA	O-C-N	-6.76	113.86	121.43
1	A	65	PRO	O-C-N	-6.70	114.38	122.89
1	B	67	ALA	CA-C-N	6.68	126.38	119.56
1	B	67	ALA	C-N-CA	6.68	126.38	119.56
1	C	2	ASP	CA-C-O	-6.65	111.27	119.18
1	A	55	GLY	N-CA-C	-6.64	103.64	112.81
1	A	89	CYS	N-CA-C	-6.64	104.13	111.36
1	C	10	ALA	CA-C-O	-6.59	112.39	119.97
1	A	14	ALA	O-C-N	6.56	129.07	122.12
1	A	87	ARG	NE-CZ-NH2	-6.55	113.30	119.20
1	A	72	PRO	N-CA-C	6.39	122.42	113.53
1	A	34	THR	N-CA-C	-6.39	104.03	113.61
1	A	107	ALA	CA-C-O	-6.30	113.88	120.55
1	C	79	ALA	CB-CA-C	-6.29	100.40	110.84
1	A	111	ALA	N-CA-CB	-6.25	100.88	110.44
1	C	56	GLN	CA-C-O	6.21	127.93	121.10
1	C	87	ARG	CD-NE-CZ	6.15	133.01	124.40
1	B	111	ALA	N-CA-C	-6.13	105.38	112.92
1	A	18	ALA	CA-C-N	6.12	126.08	119.78
1	A	18	ALA	C-N-CA	6.12	126.08	119.78
1	A	79	ALA	O-C-N	-6.12	115.64	122.12
1	B	35	GLY	N-CA-C	6.08	125.40	115.73
1	C	49	ALA	CA-C-O	6.06	126.39	119.18
1	C	85	LEU	CA-C-O	6.06	126.94	120.10
1	B	7	GLN	CG-CD-NE2	6.05	125.48	116.40
1	B	64	ALA	CA-C-O	6.05	127.53	120.69
1	C	35	GLY	CA-C-N	-6.05	109.56	121.41
1	C	35	GLY	C-N-CA	-6.05	109.56	121.41
1	A	40	THR	N-CA-CB	6.04	118.64	110.79
1	C	78	SER	CB-CA-C	-5.99	101.22	110.81
1	C	43	CYS	CA-C-O	5.98	126.89	120.55
1	A	103	ALA	CB-CA-C	-5.97	101.50	110.88
1	C	88	ASN	N-CA-C	5.97	120.63	113.23
1	C	84	TRP	CA-CB-CG	-5.94	102.31	113.60
1	A	88	ASN	OD1-CG-ND2	5.91	128.51	122.60
1	B	71	THR	CA-C-N	5.89	126.31	119.47
1	B	71	THR	C-N-CA	5.89	126.31	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	ALA	N-CA-C	-5.89	105.76	113.12
1	A	27	ALA	N-CA-CB	5.84	119.34	110.28
1	A	47	HIS	CA-C-O	-5.81	112.26	119.18
1	A	10	ALA	N-CA-C	5.81	119.18	111.75
1	A	14	ALA	N-CA-CB	5.76	118.59	110.12
1	B	103	ALA	O-C-N	5.73	127.97	122.07
1	A	6	ALA	CB-CA-C	-5.70	101.33	110.79
1	B	62	GLU	CA-CB-CG	5.69	125.49	114.10
1	C	6	ALA	O-C-N	-5.67	116.11	122.12
1	B	40	THR	O-C-N	-5.66	116.28	121.37
1	A	22	ALA	O-C-N	5.64	128.82	122.22
1	C	91	SER	CB-CA-C	-5.63	101.45	110.79
1	A	24	ARG	NE-CZ-NH2	5.62	124.25	119.20
1	B	99	PRO	CA-C-N	5.60	126.20	119.98
1	B	99	PRO	C-N-CA	5.60	126.20	119.98
1	A	23	GLU	N-CA-CB	5.60	118.35	110.12
1	A	59	THR	N-CA-CB	5.59	119.87	110.92
1	C	44	THR	OG1-CB-CG2	5.59	120.49	109.30
1	C	107	ALA	CA-C-O	-5.59	114.63	120.55
1	A	54	ALA	CB-CA-C	5.58	119.60	109.62
1	A	6	ALA	N-CA-CB	5.58	118.32	110.12
1	A	52	THR	N-CA-C	-5.56	106.45	113.18
1	A	79	ALA	CA-C-N	5.56	127.73	120.28
1	A	79	ALA	C-N-CA	5.56	127.73	120.28
1	B	54	ALA	CB-CA-C	-5.54	99.34	109.54
1	A	33	GLN	OE1-CD-NE2	-5.51	117.09	122.60
1	A	6	ALA	N-CA-C	5.48	117.26	111.28
1	C	33	GLN	OE1-CD-NE2	-5.47	117.13	122.60
1	A	20	ALA	CA-C-O	-5.45	115.54	121.87
1	C	87	ARG	CA-C-O	-5.44	115.11	120.82
1	C	23	GLU	CB-CG-CD	5.39	121.77	112.60
1	C	79	ALA	O-C-N	5.38	128.71	122.20
1	B	21	ASP	CB-CG-OD1	5.36	130.72	118.40
1	C	62	GLU	CA-CB-CG	5.31	124.73	114.10
1	A	21	ASP	CA-C-N	5.26	127.85	120.28
1	A	21	ASP	C-N-CA	5.26	127.85	120.28
1	C	61	LYS	CB-CA-C	5.23	118.26	109.89
1	B	74	ARG	NE-CZ-NH2	5.22	123.90	119.20
1	B	79	ALA	CA-C-O	5.22	125.95	120.42
1	C	30	LEU	CA-C-O	-5.22	112.84	119.31
1	C	90	ASN	OD1-CG-ND2	5.17	127.77	122.60
1	A	110	ALA	CA-C-O	-5.16	113.24	119.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	74	ARG	NE-CZ-NH1	-5.16	116.34	121.50
1	A	71	THR	CA-CB-OG1	5.13	117.29	109.60
1	C	24	ARG	CB-CA-C	5.13	119.57	110.85
1	B	61	LYS	N-CA-C	-5.12	102.12	110.20
1	B	52	THR	N-CA-C	-5.11	106.82	113.16
1	B	58	ARG	N-CA-C	-5.11	106.31	112.54
1	C	2	ASP	CA-C-N	-5.11	112.69	122.08
1	C	2	ASP	C-N-CA	-5.11	112.69	122.08
1	C	4	SER	N-CA-C	-5.09	102.35	108.25
1	B	98	THR	CA-C-N	5.05	125.08	119.32
1	B	98	THR	C-N-CA	5.05	125.08	119.32
1	A	110	ALA	O-C-N	5.01	128.90	122.39

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	802	0	762	26	1
1	B	794	0	740	18	1
1	C	802	0	756	29	0
2	A	45	0	0	0	0
2	B	43	0	30	1	0
2	C	43	0	30	7	0
3	A	2	0	0	1	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
4	A	82	0	0	1	1
4	B	81	0	0	2	0
4	C	109	0	0	0	1
All	All	2807	0	2318	72	2

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 15.

All (72) 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:A:64:ALA:CB	1:A:64:ALA:CA	1.77	1.60
1:B:58:ARG:CA	1:B:58:ARG:CB	1.80	1.57
1:C:61:LYS:CA	1:C:61:LYS:CB	1.79	1.56
1:C:14:ALA:CA	1:C:14:ALA:CB	1.75	1.54
1:A:18:ALA:CA	1:A:18:ALA:CB	1.90	1.49
1:C:46:CYS:SG	2:C:113:HEM:CAC	2.06	1.44
1:B:61:LYS:CA	1:B:61:LYS:C	1.94	1.40
1:C:46:CYS:SG	2:C:113:HEM:HAC	1.62	1.38
1:C:79:ALA:CA	1:C:79:ALA:CB	2.03	1.34
1:C:54:ALA:CA	1:C:54:ALA:CB	2.06	1.33
1:A:34:THR:CA	1:A:34:THR:CB	2.10	1.30
1:C:46:CYS:HG	2:C:113:HEM:CAC	1.45	1.25
1:B:58:ARG:CB	1:B:58:ARG:N	2.34	0.90
1:B:61:LYS:CA	1:B:62:GLU:N	2.39	0.85
1:C:80[A]:ARG:NH2	2:C:113:HEM:O2D	2.12	0.83
1:C:14:ALA:CB	1:C:14:ALA:C	2.50	0.83
1:A:78:SER:HB2	4:A:127:HOH:O	1.79	0.82
1:B:62:GLU:HB3	4:B:194:HOH:O	1.78	0.81
1:A:18:ALA:CB	1:A:18:ALA:C	2.62	0.73
1:B:56:GLN:HA	1:B:61:LYS:O	1.92	0.70
1:A:64:ALA:CB	1:A:64:ALA:C	2.69	0.64
1:B:58:ARG:CB	1:B:58:ARG:H	2.13	0.61
1:C:68:PRO:HG2	1:C:110:ALA:HA	1.82	0.61
1:C:46:CYS:SG	2:C:113:HEM:C3C	2.90	0.60
1:C:54:ALA:CB	1:C:54:ALA:C	2.74	0.60
1:B:62:GLU:HG2	4:B:193:HOH:O	2.00	0.60
1:C:36:GLY:HA3	1:C:91:SER:O	2.02	0.59
1:C:56:GLN:HA	1:C:63:ILE:HG13	1.85	0.58
1:A:88:ASN:HB2	3:A:114:NO:N	2.20	0.57
1:B:98:THR:HB	1:B:99:PRO:HD2	1.88	0.55
1:C:46:CYS:SG	2:C:113:HEM:CBC	2.91	0.53
1:A:26:ARG:HG3	1:A:108:TRP:CE2	2.44	0.52
1:C:80[A]:ARG:CZ	2:C:113:HEM:O2D	2.58	0.51
1:B:61:LYS:C	1:B:61:LYS:N	2.66	0.51
1:B:80:ARG:NH2	2:B:113:HEM:O2D	2.45	0.50
1:C:14:ALA:CB	1:C:14:ALA:N	2.67	0.49
1:A:65:PRO:O	1:A:71:THR:HG23	2.12	0.48
1:A:73:ASP:OD1	1:A:73:ASP:N	2.46	0.48
1:C:79:ALA:CB	1:C:79:ALA:C	2.83	0.48
1:A:34:THR:CB	1:A:34:THR:N	2.74	0.47
1:A:8:LEU:HD21	1:A:78:SER:HB3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:LYS:CB	1:C:61:LYS:N	2.69	0.46
1:A:26:ARG:HG3	1:A:108:TRP:CZ2	2.52	0.45
1:B:21:ASP:OD2	1:B:24:ARG:HG3	2.17	0.45
1:C:61:LYS:CB	1:C:61:LYS:HA	2.22	0.44
1:B:61:LYS:CA	1:B:62:GLU:H	2.25	0.44
1:C:42:SER:OG	1:C:44:THR:HB	2.18	0.44
1:C:83:LYS:O	1:C:87:ARG:HG3	2.18	0.44
1:B:67:ALA:HA	1:B:68:PRO:HD3	1.83	0.43
1:B:1:GLY:O	1:B:2:ASP:HB3	2.17	0.43
1:A:3:THR:OG1	1:A:7:GLN:HG2	2.19	0.43
1:B:37:LYS:HA	1:B:38:PRO:HD3	1.85	0.43
1:C:67:ALA:HA	1:C:68:PRO:HD3	1.89	0.43
1:A:85:LEU:HD13	1:A:102:LYS:HG2	2.00	0.43
1:B:44:THR:OG1	1:C:24:ARG:HD2	2.18	0.43
1:C:55:GLY:O	1:C:56:GLN:HB3	2.18	0.42
1:C:42:SER:HB3	1:C:44:THR:HG22	2.00	0.42
1:A:68:PRO:HD2	1:A:112:GLN:OXT	2.20	0.42
1:B:77:ASP:HB3	1:B:80:ARG:HB3	2.00	0.42
1:C:56:GLN:HB2	1:C:61:LYS:O	2.19	0.42
1:A:37:LYS:HB3	1:A:38:PRO:HD2	2.03	0.41
1:C:68:PRO:HD2	1:C:112:GLN:OXT	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:GLN:OE1	4:C:177:HOH:O[3_656]	2.19	0.01
1:B:62:GLU:CA	4:A:182:HOH:O[4_556]	2.19	0.01

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	111/112 (99%)	111 (100%)	0	0	100	100
1	B	110/112 (98%)	108 (98%)	1 (1%)	1 (1%)	14	14
1	C	111/112 (99%)	110 (99%)	1 (1%)	0	100	100
All	All	332/336 (99%)	329 (99%)	2 (1%)	1 (0%)	36	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2	ASP

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	78/81 (96%)	76 (97%)	2 (3%)	40	55
1	B	77/81 (95%)	73 (95%)	4 (5%)	21	26
1	C	78/81 (96%)	75 (96%)	3 (4%)	29	40
All	All	233/243 (96%)	224 (96%)	9 (4%)	28	39

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	71	THR
1	B	2	ASP
1	B	8	LEU
1	B	71	THR
1	B	91	SER
1	C	44	THR
1	C	56	GLN
1	C	72	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	ASN
1	A	112	GLN
1	B	88	ASN
1	B	90	ASN
1	C	7	GLN
1	C	56	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 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	C	113	1,3	50,50,50	1.59	10 (20%)	67,82,82	1.40	10 (14%)
2	HEM	A	113[A]	-	50,50,50	1.64	9 (18%)	67,82,82	1.43	9 (13%)
2	HEM	B	113	1,3	50,50,50	1.64	14 (28%)	67,82,82	1.74	15 (22%)
3	NO	B	114	2	0,1,1	-	-	-	-	-
3	NO	A	114	2	0,1,1	-	-	-	-	-
2	HEM	A	113[B]	-	50,50,50	1.67	10 (20%)	67,82,82	1.39	8 (11%)
3	NO	C	114	2	0,1,1	-	-	-	-	-

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	C	113	1,3	-	7/14/54/54	-
2	HEM	A	113[B]	-	-	7/14/54/54	-
2	HEM	A	113[A]	-	-	9/14/54/54	-
2	HEM	B	113	1,3	-	9/14/54/54	-

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	113[A]	HEM	CAC-C3C	4.32	1.58	1.47
2	A	113[B]	HEM	CAC-C3C	4.32	1.58	1.47
2	C	113	HEM	CAC-C3C	4.31	1.58	1.47
2	B	113	HEM	CAB-C3B	4.15	1.58	1.47
2	A	113[A]	HEM	CAB-C3B	3.93	1.57	1.47
2	A	113[B]	HEM	CAB-C3B	3.93	1.57	1.47
2	B	113	HEM	CMC-C2C	3.51	1.58	1.50
2	C	113	HEM	CAB-C3B	3.46	1.56	1.47
2	A	113[A]	HEM	C3C-C4C	-3.32	1.40	1.46
2	A	113[B]	HEM	C3C-C4C	-3.32	1.40	1.46
2	B	113	HEM	CAC-C3C	3.20	1.55	1.47
2	B	113	HEM	CMB-C2B	3.20	1.57	1.50
2	A	113[A]	HEM	CBB-CAB	2.84	1.44	1.30
2	A	113[B]	HEM	CBB-CAB	2.84	1.44	1.30
2	A	113[A]	HEM	CMA-C3A	2.76	1.56	1.50
2	A	113[B]	HEM	CMA-C3A	2.76	1.56	1.50
2	C	113	HEM	O2A-CGA	-2.69	1.21	1.30
2	C	113	HEM	C3C-C4C	-2.64	1.41	1.46
2	C	113	HEM	CBB-CAB	2.63	1.43	1.30
2	C	113	HEM	CMB-C2B	2.63	1.56	1.50
2	B	113	HEM	C2A-C3A	-2.57	1.32	1.38
2	B	113	HEM	CBB-CAB	2.56	1.42	1.30
2	A	113[A]	HEM	CMB-C2B	2.55	1.56	1.50
2	A	113[B]	HEM	CMB-C2B	2.55	1.56	1.50
2	C	113	HEM	CMD-C2D	2.47	1.55	1.50
2	B	113	HEM	CMD-C2D	2.45	1.55	1.50
2	A	113[A]	HEM	CMC-C2C	2.41	1.55	1.50
2	A	113[B]	HEM	CMC-C2C	2.41	1.55	1.50
2	B	113	HEM	CMA-C3A	2.34	1.55	1.50
2	B	113	HEM	C3C-C2C	-2.26	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	113	HEM	FE-NC	2.24	2.02	1.95
2	B	113	HEM	CBC-CAC	2.23	1.41	1.30
2	A	113[A]	HEM	C3B-C2B	-2.23	1.32	1.37
2	A	113[B]	HEM	C3B-C2B	-2.23	1.32	1.37
2	A	113[B]	HEM	O2A-CGA	-2.17	1.23	1.30
2	B	113	HEM	C3B-C2B	-2.16	1.32	1.37
2	A	113[A]	HEM	FE-ND	2.12	2.01	1.94
2	A	113[B]	HEM	FE-ND	2.12	2.01	1.94
2	B	113	HEM	O1D-CGD	2.07	1.28	1.22
2	C	113	HEM	FE-ND	2.05	2.01	1.94
2	C	113	HEM	CBC-CAC	2.04	1.40	1.30
2	C	113	HEM	FE-NB	2.03	2.01	1.94
2	B	113	HEM	CAD-C3D	2.03	1.56	1.51

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	113	HEM	CAD-CBD-CGD	6.25	130.24	113.67
2	A	113[A]	HEM	CBB-CAB-C3B	-4.05	107.27	127.53
2	A	113[B]	HEM	CBB-CAB-C3B	-4.05	107.27	127.53
2	C	113	HEM	CHC-C4B-NB	3.98	128.71	124.42
2	B	113	HEM	CHC-C1C-NC	3.87	128.67	124.45
2	C	113	HEM	CBB-CAB-C3B	-3.67	109.19	127.53
2	C	113	HEM	O2A-CGA-O1A	-3.52	114.27	123.33
2	A	113[A]	HEM	O1A-CGA-CBA	-3.50	111.99	123.09
2	A	113[A]	HEM	C3B-C4B-NB	-3.34	107.07	109.47
2	A	113[B]	HEM	C3B-C4B-NB	-3.34	107.07	109.47
2	B	113	HEM	O2D-CGD-CBD	3.31	124.47	114.00
2	A	113[A]	HEM	C4B-C3B-C2B	2.96	110.00	107.28
2	A	113[B]	HEM	C4B-C3B-C2B	2.96	110.00	107.28
2	B	113	HEM	O2A-CGA-CBA	2.95	123.31	114.00
2	B	113	HEM	O2D-CGD-O1D	-2.93	115.80	123.33
2	A	113[A]	HEM	CHC-C4B-NB	2.87	127.51	124.42
2	A	113[B]	HEM	CHC-C4B-NB	2.87	127.51	124.42
2	C	113	HEM	O1D-CGD-CBD	-2.81	114.18	123.09
2	B	113	HEM	CMA-C3A-C2A	2.80	131.57	125.62
2	A	113[B]	HEM	O2A-CGA-O1A	2.80	130.53	123.33
2	B	113	HEM	CMA-C3A-C4A	-2.72	121.28	125.42
2	B	113	HEM	CHD-C4C-NC	2.68	127.38	124.45
2	B	113	HEM	CBB-CAB-C3B	-2.65	114.29	127.53
2	C	113	HEM	C3B-C4B-NB	-2.58	107.61	109.47
2	A	113[A]	HEM	CBC-CAC-C3C	-2.57	114.69	127.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	113[B]	HEM	CBC-CAC-C3C	-2.57	114.69	127.53
2	C	113	HEM	C4B-C3B-C2B	2.56	109.64	107.28
2	A	113[A]	HEM	C2A-C1A-NA	-2.50	107.38	110.15
2	A	113[B]	HEM	C2A-C1A-NA	-2.50	107.38	110.15
2	C	113	HEM	O2A-CGA-CBA	2.48	121.84	114.00
2	B	113	HEM	CHD-C1D-ND	2.36	126.96	124.42
2	A	113[A]	HEM	CHA-C4D-ND	2.35	127.28	124.37
2	A	113[B]	HEM	CHA-C4D-ND	2.35	127.28	124.37
2	B	113	HEM	CBA-CAA-C2A	2.34	119.00	112.53
2	C	113	HEM	CMC-C2C-C1C	-2.32	120.65	124.73
2	C	113	HEM	CHA-C1A-NA	2.27	127.97	123.86
2	A	113[A]	HEM	O2A-CGA-CBA	2.14	120.75	114.00
2	B	113	HEM	C3B-C4B-NB	-2.13	107.94	109.47
2	B	113	HEM	C4C-CHD-C1D	-2.05	121.66	126.02
2	B	113	HEM	CBC-CAC-C3C	-2.05	117.30	127.53
2	B	113	HEM	CHC-C4B-NB	2.03	126.61	124.42
2	C	113	HEM	C1C-CHC-C4B	-2.00	121.76	126.02

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	113[A]	HEM	C2B-C3B-CAB-CBB
2	A	113[A]	HEM	C4B-C3B-CAB-CBB
2	A	113[B]	HEM	C2B-C3B-CAB-CBB
2	A	113[B]	HEM	C4B-C3B-CAB-CBB
2	B	113	HEM	C2B-C3B-CAB-CBB
2	B	113	HEM	C4B-C3B-CAB-CBB
2	C	113	HEM	C2B-C3B-CAB-CBB
2	C	113	HEM	C4B-C3B-CAB-CBB
2	C	113	HEM	C2C-C3C-CAC-CBC
2	A	113[A]	HEM	C3D-CAD-CBD-CGD
2	A	113[B]	HEM	C3D-CAD-CBD-CGD
2	B	113	HEM	C3D-CAD-CBD-CGD
2	A	113[A]	HEM	C2C-C3C-CAC-CBC
2	A	113[B]	HEM	C2C-C3C-CAC-CBC
2	B	113	HEM	C2C-C3C-CAC-CBC
2	A	113[A]	HEM	CAA-CBA-CGA-O1A
2	A	113[A]	HEM	CAA-CBA-CGA-O2A
2	C	113	HEM	C2D-C3D-CAD-CBD
2	A	113[A]	HEM	C4C-C3C-CAC-CBC
2	A	113[B]	HEM	C4C-C3C-CAC-CBC

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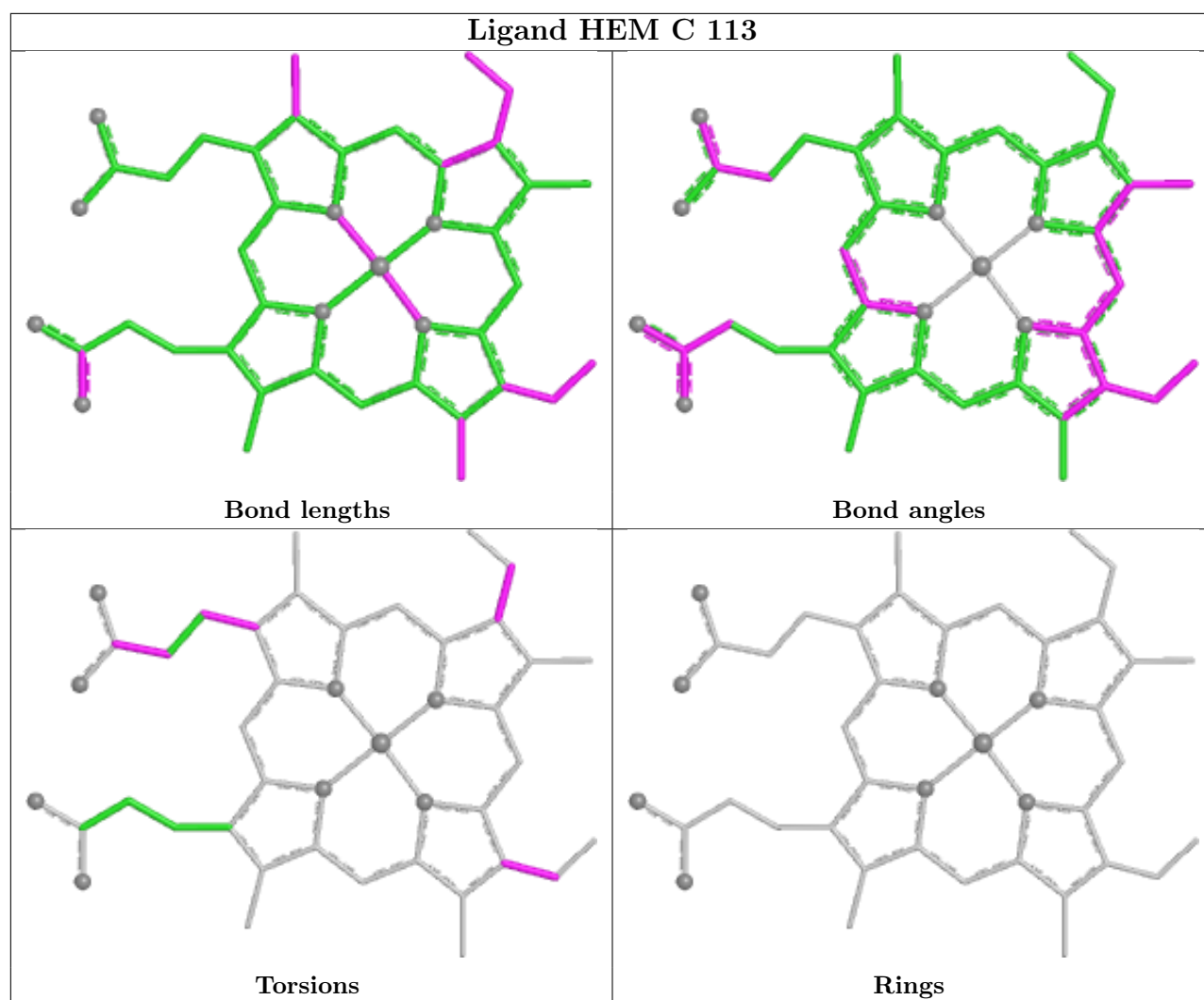
Mol	Chain	Res	Type	Atoms
2	B	113	HEM	C4C-C3C-CAC-CBC
2	C	113	HEM	C4C-C3C-CAC-CBC
2	B	113	HEM	CAD-CBD-CGD-O1D
2	B	113	HEM	CAA-CBA-CGA-O1A
2	B	113	HEM	CAD-CBD-CGD-O2D
2	C	113	HEM	CAD-CBD-CGD-O1D
2	C	113	HEM	CAD-CBD-CGD-O2D
2	B	113	HEM	CAA-CBA-CGA-O2A
2	A	113[A]	HEM	CAD-CBD-CGD-O2D
2	A	113[B]	HEM	CAD-CBD-CGD-O2D
2	A	113[A]	HEM	CAD-CBD-CGD-O1D
2	A	113[B]	HEM	CAD-CBD-CGD-O1D

There are no ring outliers.

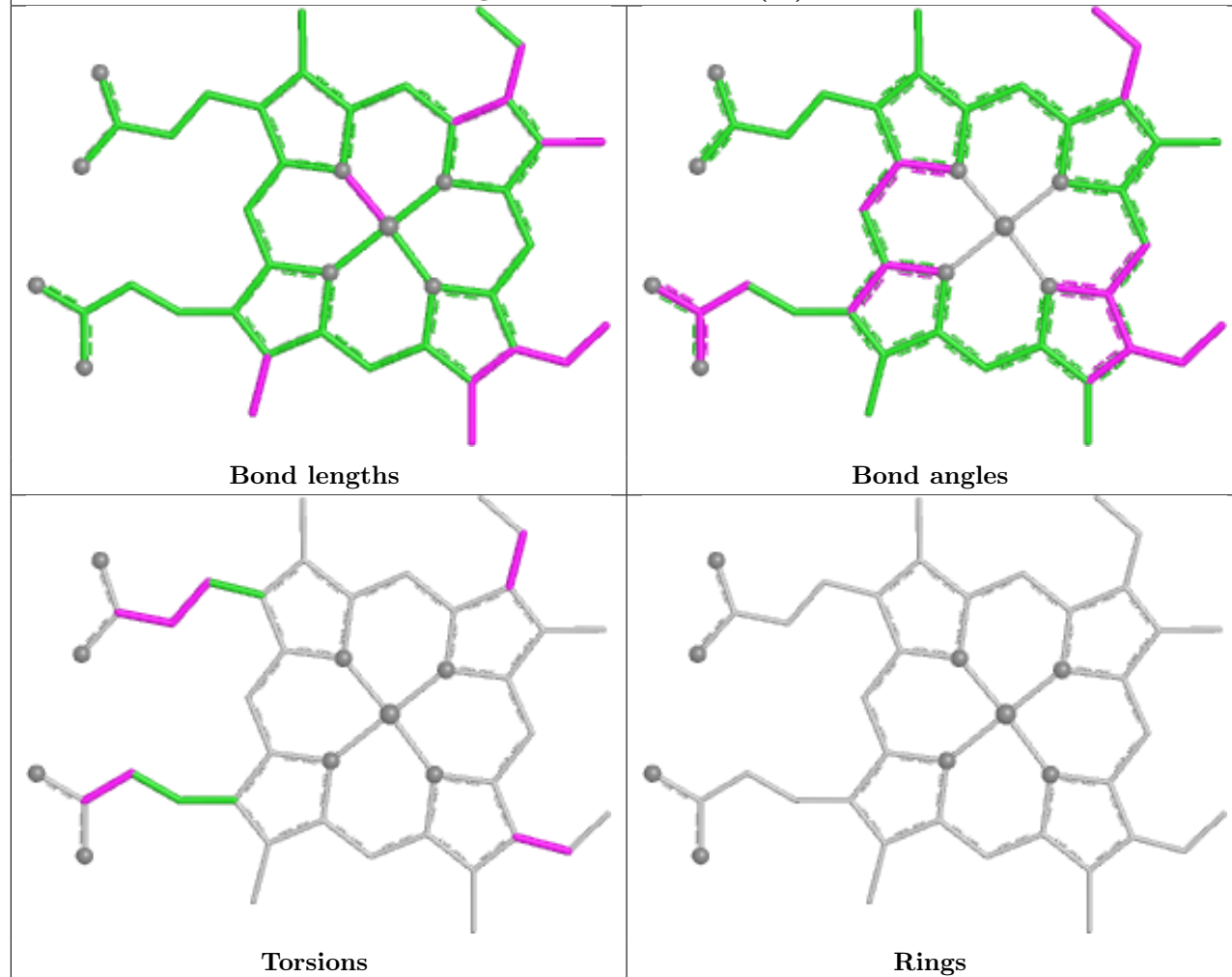
3 monomers are involved in 9 short contacts:

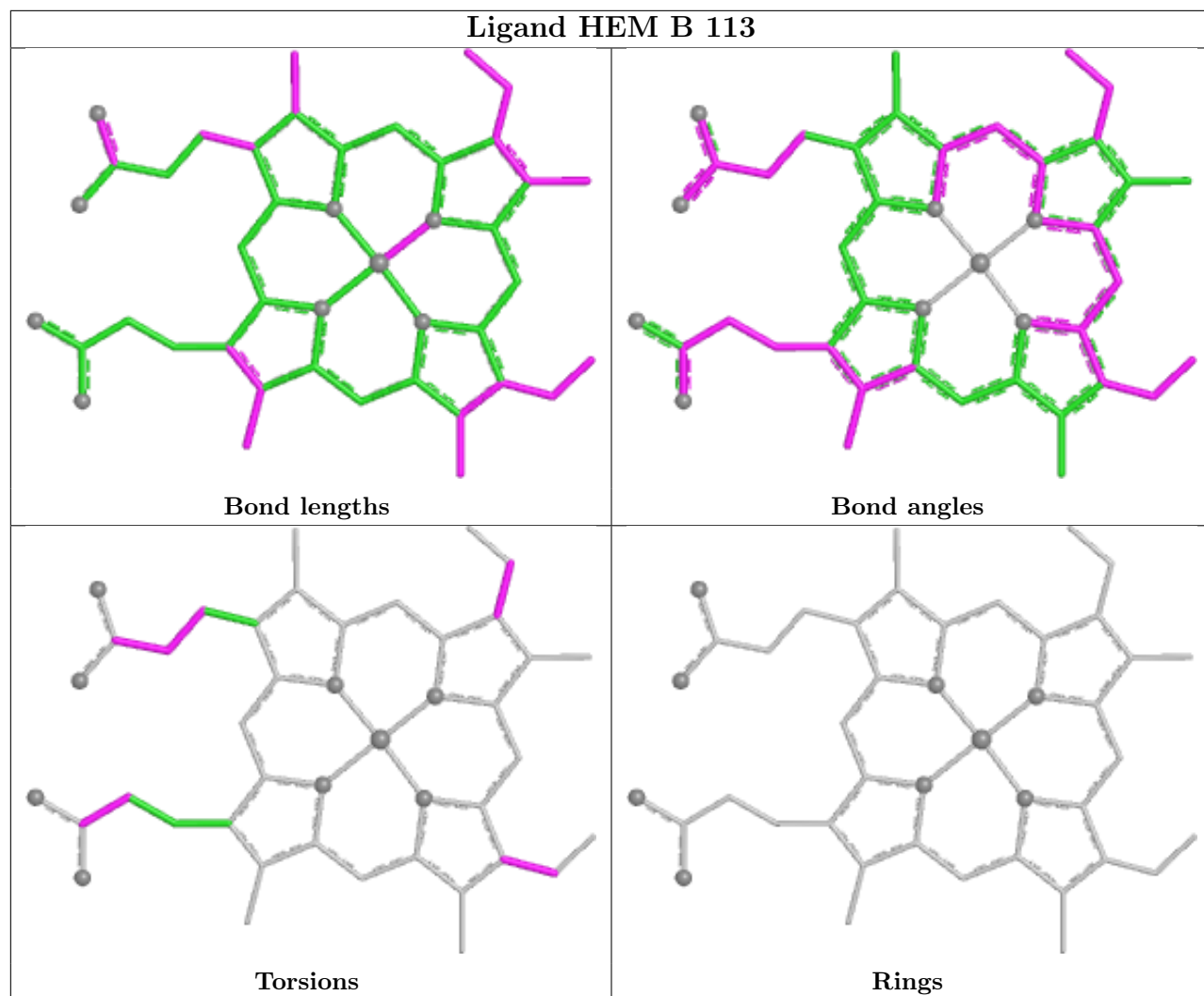
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	113	HEM	7	0
2	B	113	HEM	1	0
3	A	114	NO	1	0

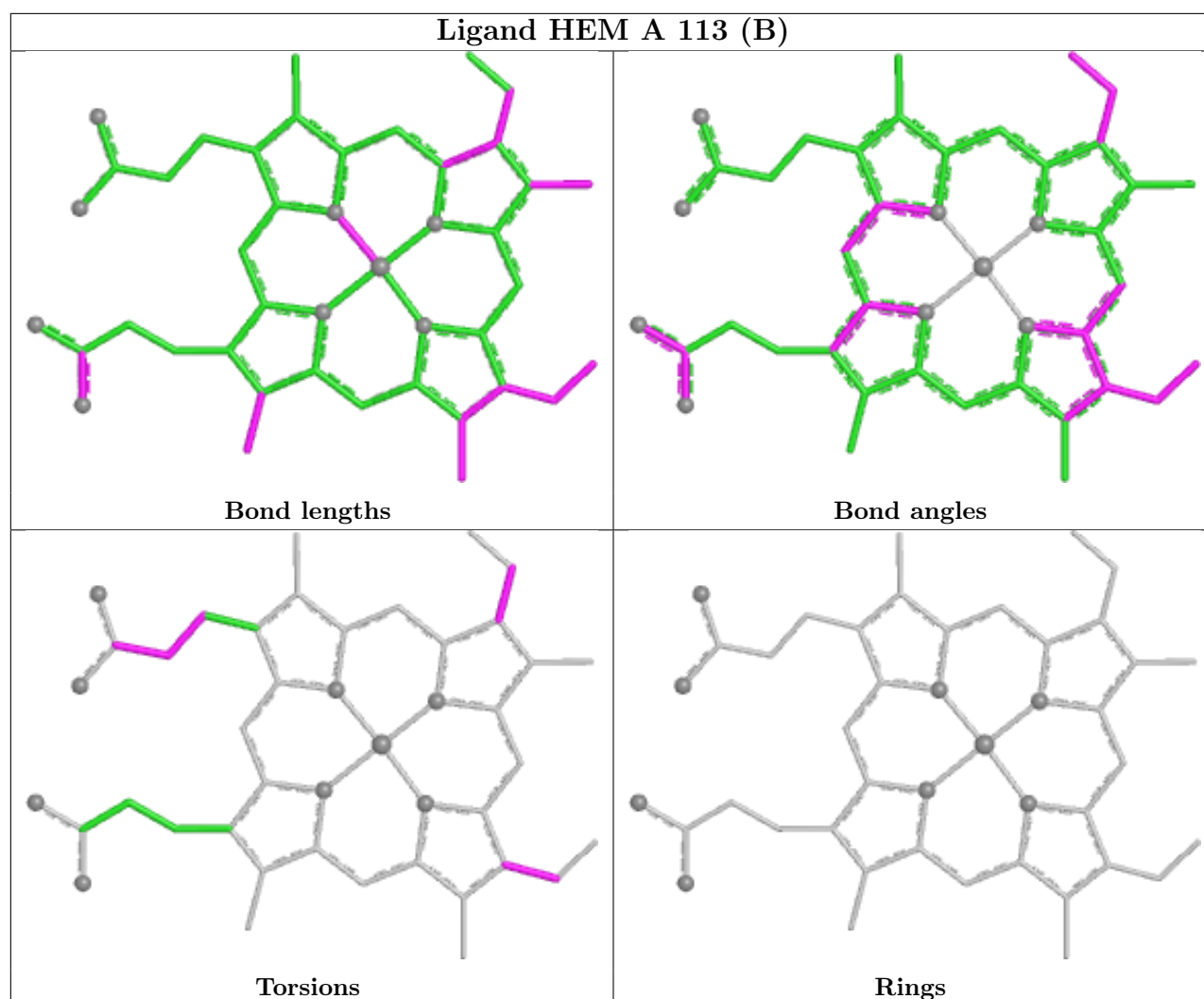
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 HEM A 113 (A)







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	112/112 (100%)	-0.37	2 (1%) 67 64	21, 31, 49, 63	1 (0%)
1	B	112/112 (100%)	-0.43	3 (2%) 56 53	23, 31, 48, 59	0
1	C	112/112 (100%)	-0.40	0 100 100	20, 32, 52, 62	1 (0%)
All	All	336/336 (100%)	-0.40	5 (1%) 72 69	20, 31, 51, 63	2 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	GLY	4.3
1	B	62	GLU	3.2
1	A	7	GLN	2.6
1	B	35	GLY	2.3
1	A	56	GLN	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

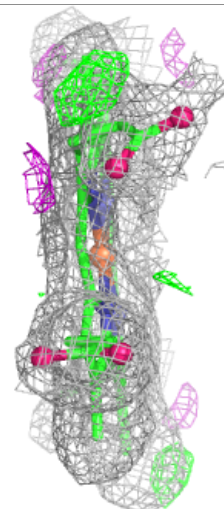
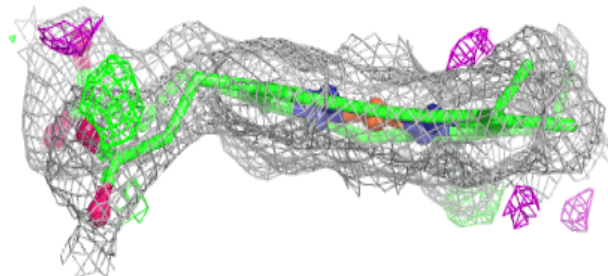
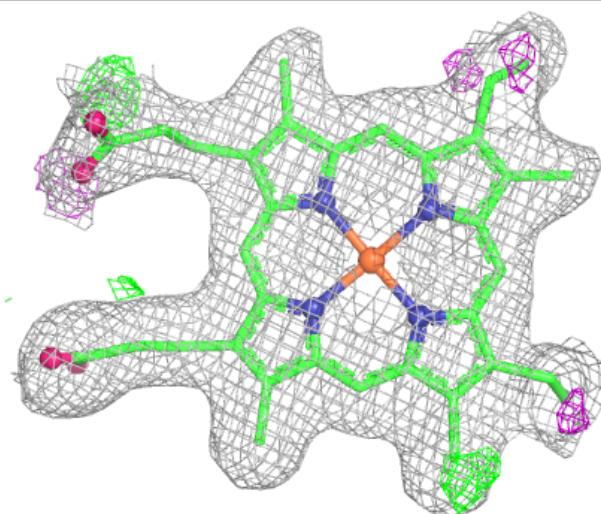
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NO	A	114	2/2	0.97	0.09	36,36,36,44	0
3	NO	C	114	2/2	0.97	0.13	37,37,37,37	0
2	HEM	B	113	43/43	0.98	0.06	20,24,41,50	0
2	HEM	C	113	43/43	0.98	0.06	25,31,43,47	0
2	HEM	A	113[A]	43/43	0.98	0.05	21,25,36,43	2
2	HEM	A	113[B]	43/43	0.98	0.05	21,25,36,43	2
3	NO	B	114	2/2	0.99	0.10	28,28,28,29	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.

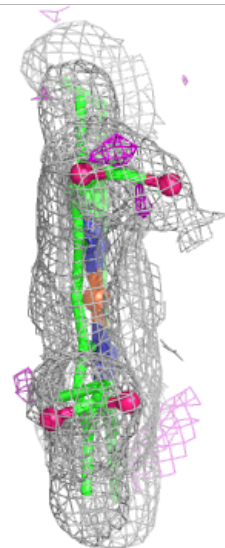
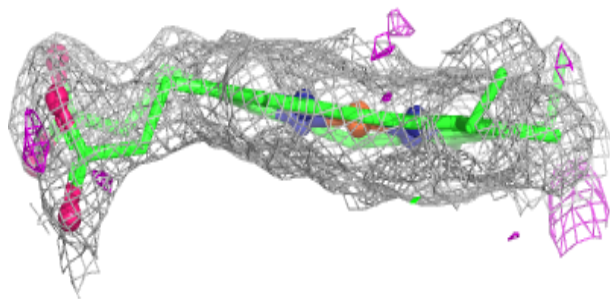
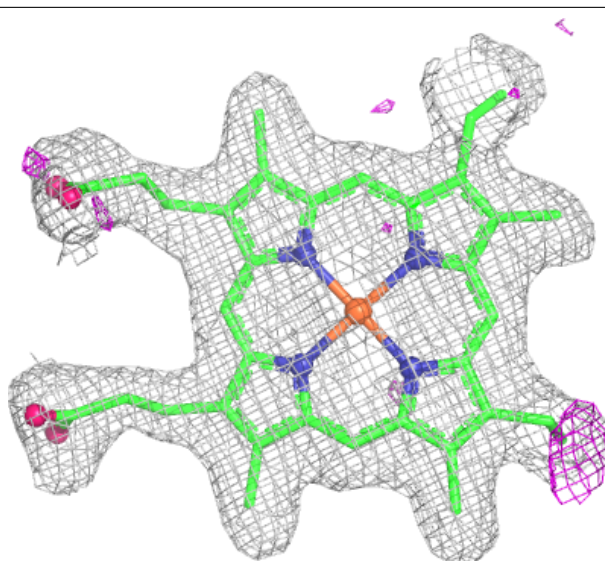
Electron density around HEM B 113:

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



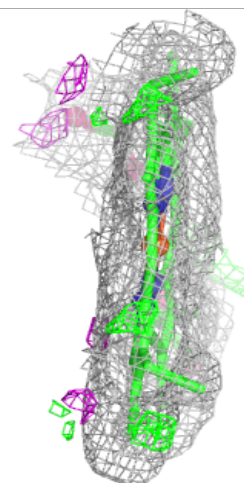
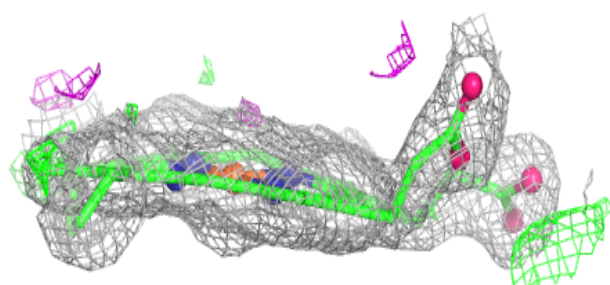
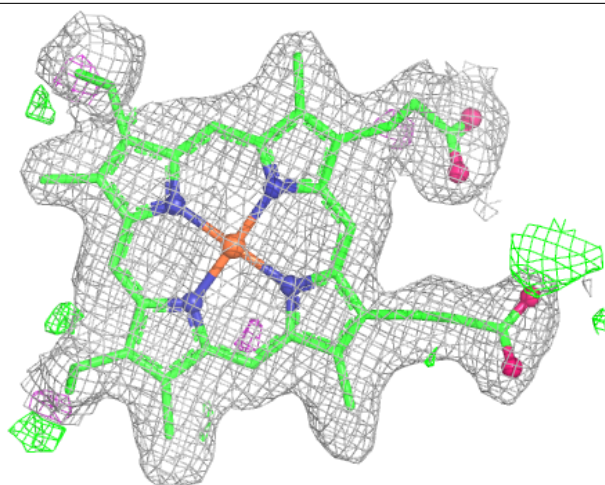
Electron density around HEM C 113:

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



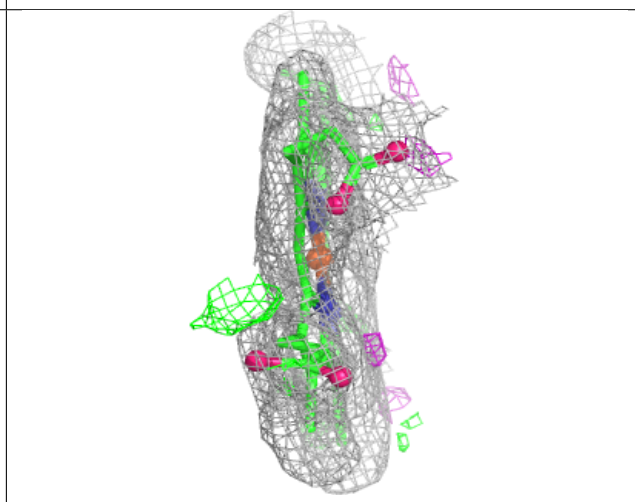
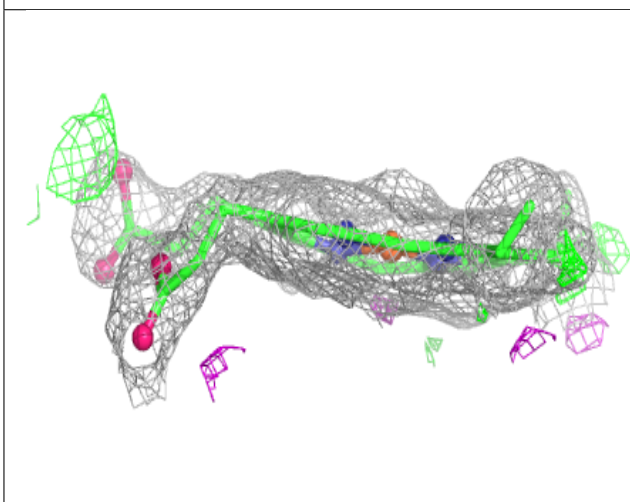
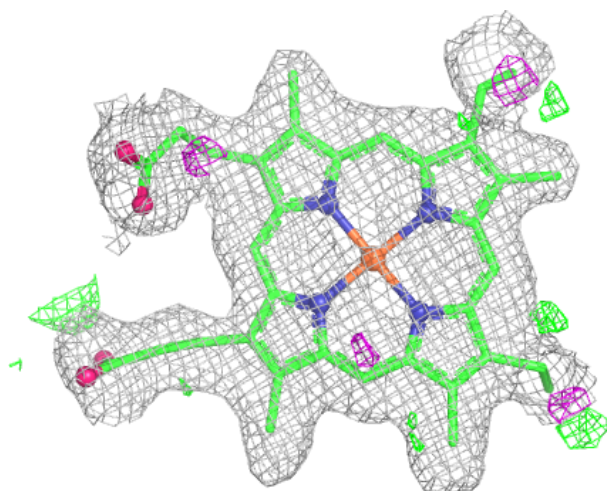
Electron density around HEM A 113 (A):

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



Electron density around HEM A 113 (B):

$2mF_o-DF_c$ (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.