Full wwPDB X-ray Structure Validation Report (i)

Nov 1, 2023 – 03:52 AM EDT

PDB ID	:	3HRW
Title	:	Crystal structure of hemoglobin from mouse (Mus musculus)at 2.8
Authors	:	Sundaresan, S.S.; Ramesh, P.; Ponnuswamy, M.N.
Deposited on	:	2009-06-10
Resolution	:	2.80 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 (i) 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 (i)) were used in the production of this report:

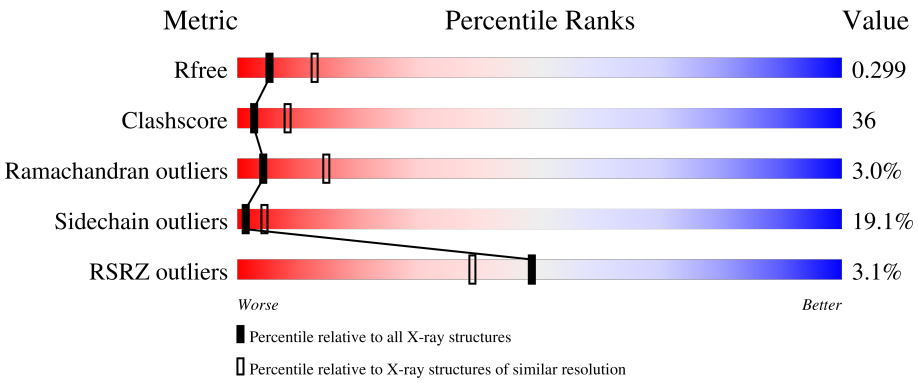
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 <=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	141	3% 47% 41% 11% •
1	C	141	4% 47% 43% 10%
2	B	146	3% 40% 46% 12% •
2	D	146	3% 43% 45% 11% •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4548 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.

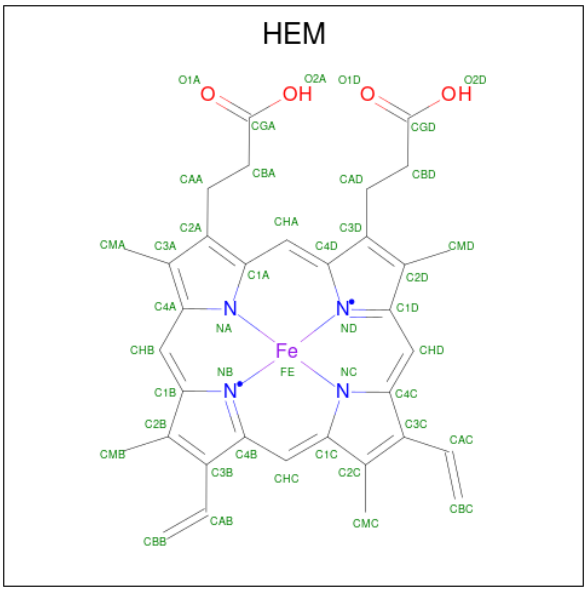
- Molecule 1 is a protein called Hemoglobin subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	141	1057	674	187	194	2	0	0	0
1	C	141	1057	674	187	194	2	0	0	0

- Molecule 2 is a protein called Hemoglobin subunit beta-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	146	1108	708	195	201	4	0	0	0
2	D	146	1108	708	195	201	4	0	0	0

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
3	A	1	43	34	1	4	4	0	0
3	B	1	43	34	1	4	4	0	0
3	C	1	43	34	1	4	4	0	0
3	D	1	43	34	1	4	4	0	0

- Molecule 4 is water.

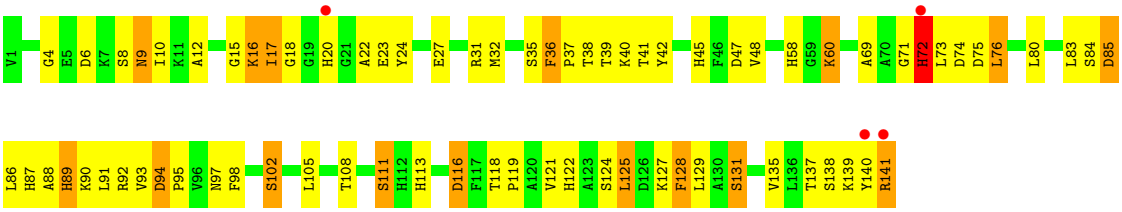
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	15	15	15	0	0
4	B	9	9	9	0	0
4	C	12	12	12	0	0
4	D	10	10	10	0	0

3 Residue-property plots (i)

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

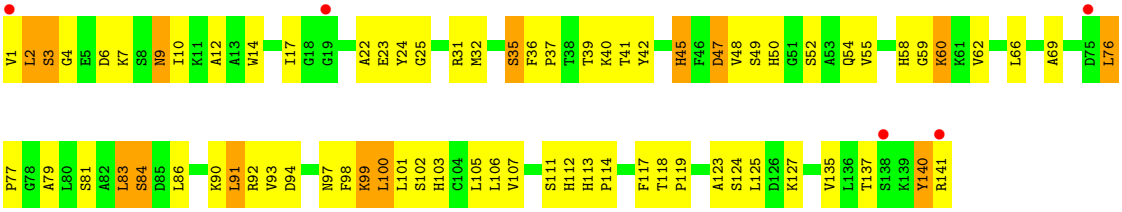
- Molecule 1: Hemoglobin subunit alpha

Chain A:



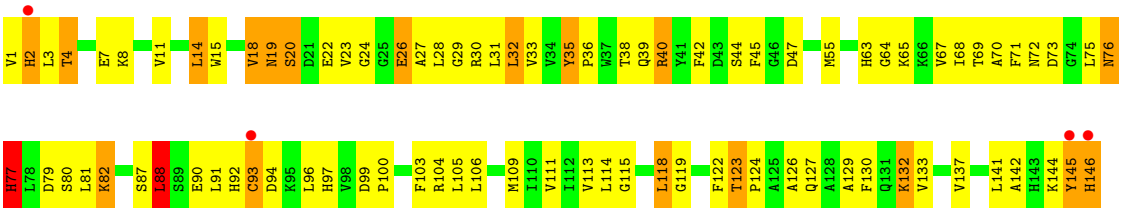
- Molecule 1: Hemoglobin subunit alpha

Chain C:



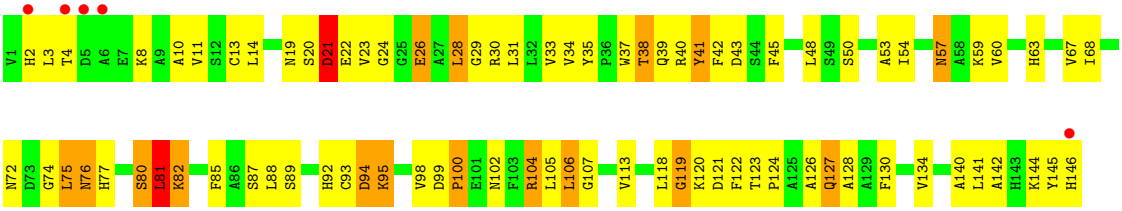
- Molecule 2: Hemoglobin subunit beta-1

Chain B:



- Molecule 2: Hemoglobin subunit beta-1

Chain D:


V1
H2
L3
T4
D5
A6
E7
K8
A9
A10
V11
S12
C13
L14
N19
S20
D21
E22
V23
G24
G25
E26
A27
L28
G29
R30
L31
L32
V33
V34
Y35
P36
W37
T38
Q39
R40
Y41
F42
D43
S44
F45
L48
S49
S50
A53
I54
N57
A58
K59
V60
H63
V67
I68
N72
D73
G74
L75
N76
H77
S80
L81
K82
F85
A86
S87
L88
S89
H92
C93
D94
K95
V98
D99
P100
E101
N102
F103
R104
L105
L106
G107
V113
L118
G119
K120
D121
F122
T123
P124
A125
A126
Q127
A128
A129
F130
V134
A140
L141
A142
H143
K144
Y145
H146

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	53.08Å 65.94Å 150.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.94 – 2.80 19.94 – 2.80	Depositor EDS
% Data completeness (in resolution range)	92.9 (19.94-2.80) 92.9 (19.94-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$< I/\sigma(I) >$ [1]	1.59 (at 2.79Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.243 , 0.306 0.240 , 0.299	Depositor DCC
R_{free} test set	616 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å^2)	47.1	Xtriage
Anisotropy	0.236	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 51.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	4548	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.29% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.49	0/1085	0.76	5/1471 (0.3%)
1	C	0.49	0/1085	0.70	0/1471
2	B	0.53	0/1133	0.85	4/1534 (0.3%)
2	D	0.50	0/1133	0.79	3/1534 (0.2%)
All	All	0.50	0/4436	0.78	12/6010 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	88	LEU	CB-CA-C	-9.06	92.98	110.20
2	D	81	LEU	CB-CA-C	-8.01	94.99	110.20
2	D	81	LEU	N-CA-C	6.17	127.65	111.00
2	D	19	ASN	CB-CA-C	5.76	121.92	110.40
1	A	88	ALA	CB-CA-C	-5.62	101.66	110.10
2	B	96	LEU	CB-CA-C	-5.62	99.52	110.20
1	A	36	PHE	N-CA-C	-5.61	95.86	111.00
1	A	89	HIS	CB-CA-C	-5.35	99.70	110.40
1	A	105	LEU	CA-CB-CG	5.28	127.45	115.30
2	B	35	TYR	N-CA-C	-5.24	96.85	111.00
2	B	2	HIS	CB-CA-C	5.19	120.78	110.40
1	A	75	ASP	N-CA-C	-5.14	97.11	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1057	0	1048	83	0
1	C	1057	0	1048	83	0
2	B	1108	0	1105	95	0
2	D	1108	0	1105	83	0
3	A	43	0	30	8	0
3	B	43	0	30	7	0
3	C	43	0	30	9	0
3	D	43	0	30	7	0
4	A	15	0	0	0	0
4	B	9	0	0	2	0
4	C	12	0	0	1	0
4	D	10	0	0	2	0
All	All	4548	0	4426	318	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 36.

All (318) 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)
2:B:36:PRO:O	2:B:39:GLN:HG2	1.48	1.12
1:A:83:LEU:HD13	3:A:200:HEM:HHB	1.36	1.08
2:B:32:LEU:H	2:B:32:LEU:HD23	1.10	1.06
2:B:87:SER:O	2:B:88:LEU:HB2	1.44	1.06
1:C:3:SER:HB2	1:C:4:GLY:HA2	1.36	1.04
1:C:3:SER:CB	1:C:4:GLY:HA2	1.89	1.02
2:B:20:SER:O	2:B:68:ILE:HD12	1.60	1.01
2:B:35:TYR:O	2:B:38:THR:HG22	1.61	1.00
1:A:6:ASP:OD2	1:A:127:LYS:HE2	1.60	0.99
2:D:63:HIS:HE1	4:D:147:HOH:O	1.47	0.96
2:B:32:LEU:H	2:B:32:LEU:CD2	1.78	0.95
1:C:32:MET:SD	1:C:39:THR:HG21	2.07	0.94
1:A:95:PRO:HG3	1:A:140:TYR:OH	1.69	0.92
1:A:141:ARG:CZ	1:C:1:VAL:HG21	2.01	0.91
2:B:68:ILE:HA	2:B:71:PHE:HB3	1.52	0.89

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:24:TYR:HE1	1:C:112:HIS:HD1	1.22	0.87
2:D:100:PRO:HD3	2:D:145:TYR:CE1	2.09	0.86
2:D:63:HIS:CE1	4:D:147:HOH:O	2.22	0.86
1:A:16:LYS:HD2	1:A:116:ASP:OD2	1.77	0.84
1:A:83:LEU:CD1	3:A:200:HEM:HHB	2.09	0.83
2:B:32:LEU:HD23	2:B:32:LEU:N	1.94	0.82
2:D:34:VAL:HG23	2:D:35:TYR:CD2	2.17	0.80
1:C:24:TYR:HE1	1:C:112:HIS:ND1	1.81	0.79
2:B:72:ASN:HA	2:B:75:LEU:HD12	1.62	0.79
1:A:83:LEU:HD13	3:A:200:HEM:CHB	2.11	0.78
1:A:119:PRO:HG3	2:B:55:MET:HE3	1.65	0.78
2:B:137:VAL:O	2:B:141:LEU:HD23	1.83	0.78
2:D:24:GLY:N	2:D:68:ILE:HD11	2.00	0.77
2:D:22:GLU:O	2:D:26:GLU:HB2	1.85	0.77
1:C:103:HIS:O	1:C:107:VAL:HG23	1.85	0.76
2:B:19:ASN:O	2:B:20:SER:CB	2.32	0.76
3:D:200:HEM:HHC	3:D:200:HEM:HBB2	1.69	0.74
2:B:68:ILE:HA	2:B:71:PHE:CB	2.17	0.74
2:B:68:ILE:O	2:B:71:PHE:N	2.21	0.73
1:C:58:HIS:HE1	3:C:200:HEM:C1A	2.06	0.73
1:C:99:LYS:HB2	1:C:99:LYS:NZ	2.04	0.73
2:B:19:ASN:O	2:B:20:SER:OG	2.06	0.73
2:B:32:LEU:CD2	2:B:32:LEU:N	2.49	0.72
2:D:41:TYR:HE1	3:D:200:HEM:HAC	1.55	0.72
2:B:2:HIS:O	2:B:3:LEU:HD12	1.90	0.72
1:A:138:SER:HA	1:A:141:ARG:HH21	1.53	0.72
2:D:24:GLY:CA	2:D:68:ILE:HD11	2.19	0.72
1:C:58:HIS:HE1	3:C:200:HEM:C2A	2.09	0.71
2:B:100:PRO:HB3	2:B:142:ALA:CB	2.20	0.71
2:B:76:ASN:N	2:B:76:ASN:HD22	1.89	0.70
1:C:31:ARG:HG2	2:D:124:PRO:HG3	1.72	0.70
1:C:102:SER:O	1:C:106:LEU:HG	1.90	0.70
1:A:138:SER:HB3	1:A:139:LYS:HE2	1.72	0.70
2:B:87:SER:O	2:B:88:LEU:CB	2.32	0.69
1:A:113:HIS:HB3	1:A:116:ASP:OD1	1.92	0.69
2:B:104:ARG:HH11	2:B:104:ARG:HB2	1.57	0.69
2:D:104:ARG:HH11	2:D:104:ARG:HB2	1.57	0.69
2:B:4:THR:O	2:B:7:GLU:N	2.24	0.69
1:A:9:ASN:N	1:A:9:ASN:HD22	1.91	0.68
1:A:22:ALA:HB1	1:A:60:LYS:HG2	1.75	0.68
2:B:68:ILE:O	2:B:72:ASN:N	2.27	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:ILE:HG21	1:A:73:LEU:HD21	1.75	0.68
1:A:140:TYR:CE1	2:D:37:TRP:CE3	2.82	0.68
1:C:3:SER:CB	1:C:4:GLY:CA	2.67	0.68
2:D:85:PHE:HB3	2:D:141:LEU:HD23	1.75	0.68
1:A:127:LYS:CD	1:C:141:ARG:HG3	2.25	0.67
2:D:57:ASN:HD21	2:D:59:LYS:HB2	1.60	0.67
2:D:107:GLY:HA3	2:D:134:VAL:CG1	2.25	0.67
1:A:111:SER:HA	2:B:115:GLY:O	1.96	0.66
2:D:31:LEU:HD13	2:D:106:LEU:HA	1.76	0.66
2:D:29:GLY:O	2:D:33:VAL:HG23	1.94	0.66
2:B:100:PRO:HB3	2:B:142:ALA:HB1	1.77	0.66
2:D:94:ASP:O	2:D:95:LYS:HB3	1.96	0.65
1:C:58:HIS:CE1	3:C:200:HEM:C1A	2.84	0.65
1:C:24:TYR:CE1	1:C:112:HIS:ND1	2.61	0.65
1:C:40:LYS:HE2	1:C:48:VAL:HG21	1.76	0.65
2:B:29:GLY:O	2:B:33:VAL:HG23	1.97	0.65
2:B:35:TYR:O	2:B:38:THR:CG2	2.43	0.64
2:B:36:PRO:O	2:B:39:GLN:CG	2.37	0.64
2:B:105:LEU:O	2:B:109:MET:HG2	1.98	0.64
2:B:123:THR:HB	2:B:126:ALA:H	1.63	0.64
2:B:64:GLY:O	2:B:68:ILE:HG12	1.98	0.64
1:C:76:LEU:N	1:C:77:PRO:HD2	2.12	0.64
1:A:128:PHE:HA	1:A:131:SER:OG	1.98	0.63
1:C:52:SER:HB3	1:C:55:VAL:HG22	1.80	0.63
1:A:119:PRO:HG3	2:B:55:MET:CE	2.27	0.63
1:C:76:LEU:H	1:C:77:PRO:HD2	1.64	0.63
2:B:76:ASN:O	2:B:77:HIS:ND1	2.32	0.63
2:B:92:HIS:O	2:B:93:CYS:SG	2.56	0.63
2:D:3:LEU:HD12	2:D:4:THR:H	1.65	0.62
2:D:50:SER:O	2:D:53:ALA:N	2.32	0.62
2:D:93:CYS:SG	2:D:145:TYR:HA	2.40	0.62
2:D:107:GLY:HA3	2:D:134:VAL:HG13	1.82	0.62
1:A:84:SER:OG	1:A:139:LYS:HG2	2.00	0.61
2:D:142:ALA:HA	2:D:145:TYR:HD2	1.66	0.61
2:B:76:ASN:N	2:B:76:ASN:ND2	2.49	0.61
1:A:71:GLY:O	1:A:73:LEU:N	2.30	0.61
2:B:1:VAL:O	2:B:2:HIS:C	2.39	0.61
1:A:20:HIS:O	1:A:23:GLU:HG2	2.00	0.60
2:B:20:SER:O	2:B:68:ILE:CD1	2.44	0.60
2:B:28:LEU:O	2:B:32:LEU:CD2	2.50	0.60
2:D:142:ALA:HA	2:D:145:TYR:CD2	2.36	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:ASN:N	1:A:9:ASN:ND2	2.48	0.60
2:B:104:ARG:HB2	2:B:104:ARG:NH1	2.16	0.60
1:A:141:ARG:NH1	1:C:1:VAL:HG21	2.16	0.59
1:C:17:ILE:HG23	1:C:24:TYR:CE2	2.38	0.59
1:C:6:ASP:O	1:C:10:ILE:HG13	2.03	0.59
2:D:24:GLY:HA3	2:D:68:ILE:CD1	2.32	0.59
1:A:122:HIS:ND1	2:B:30:ARG:HD3	2.17	0.58
1:C:83:LEU:HD21	3:C:200:HEM:HMA1	1.84	0.58
1:A:140:TYR:CE1	2:D:37:TRP:CZ3	2.91	0.58
2:B:76:ASN:HD22	2:B:76:ASN:H	1.48	0.58
1:C:10:ILE:O	1:C:14:TRP:HB2	2.02	0.58
1:A:90:LYS:O	1:A:90:LYS:HG2	2.03	0.58
1:C:99:LYS:HB2	1:C:99:LYS:HZ2	1.68	0.58
2:B:18:VAL:O	2:B:18:VAL:CG1	2.50	0.58
2:B:35:TYR:CE2	2:B:105:LEU:HG	2.39	0.58
2:B:92:HIS:CE1	3:B:200:HEM:C1A	2.92	0.58
1:A:89:HIS:CE1	1:A:139:LYS:HB3	2.39	0.57
2:B:146:HIS:HD2	2:B:146:HIS:O	1.88	0.56
2:B:90:GLU:O	2:B:94:ASP:HB3	2.06	0.56
2:B:100:PRO:HA	2:B:103:PHE:CD2	2.41	0.56
2:D:11:VAL:HG13	2:D:130:PHE:CZ	2.41	0.56
1:A:140:TYR:CE1	2:D:37:TRP:HE3	2.24	0.56
3:B:200:HEM:HBC2	3:B:200:HEM:HHD	1.86	0.56
1:A:6:ASP:HA	1:A:124:SER:HB3	1.88	0.55
2:D:40:ARG:C	2:D:41:TYR:CD2	2.80	0.55
2:B:18:VAL:O	2:B:19:ASN:C	2.45	0.55
1:A:86:LEU:O	1:A:91:LEU:N	2.40	0.55
1:A:93:VAL:HG11	1:A:98:PHE:HZ	1.72	0.55
2:B:15:TRP:HZ3	2:B:114:LEU:HD11	1.73	0.54
1:C:118:THR:HB	1:C:119:PRO:HD2	1.90	0.54
1:C:77:PRO:HA	1:C:135:VAL:HG13	1.89	0.54
2:D:31:LEU:HD22	2:D:106:LEU:HG	1.90	0.54
2:D:81:LEU:O	2:D:140:ALA:HB1	2.08	0.54
1:A:95:PRO:CG	1:A:140:TYR:OH	2.47	0.54
2:B:97:HIS:ND1	1:C:41:THR:HG21	2.23	0.54
2:B:67:VAL:HG12	2:B:68:ILE:N	2.23	0.54
2:D:57:ASN:ND2	2:D:59:LYS:HB2	2.22	0.54
1:C:36:PHE:CD2	1:C:100:LEU:HG	2.43	0.54
2:B:92:HIS:HE1	3:B:200:HEM:C1A	2.26	0.53
2:D:11:VAL:HG13	2:D:130:PHE:CE2	2.42	0.53
2:D:23:VAL:HG13	2:D:113:VAL:HG12	1.88	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:VAL:O	1:A:138:SER:HB2	2.08	0.53
2:B:19:ASN:O	2:B:20:SER:HB3	2.09	0.53
2:D:37:TRP:CD1	2:D:37:TRP:C	2.81	0.53
1:A:72:HIS:HD1	1:A:72:HIS:H	1.57	0.53
2:D:24:GLY:CA	2:D:68:ILE:CD1	2.86	0.53
1:A:15:GLY:C	1:A:17:ILE:N	2.59	0.53
2:D:39:GLN:HA	2:D:42:PHE:HD2	1.72	0.53
2:B:40:ARG:HB3	1:C:92:ARG:CD	2.39	0.53
1:A:58:HIS:HE1	3:A:200:HEM:C1A	2.27	0.52
1:C:77:PRO:HA	1:C:135:VAL:CG1	2.39	0.52
1:C:111:SER:HA	2:D:119:GLY:HA2	1.91	0.52
1:A:121:VAL:HG12	1:A:125:LEU:HD22	1.90	0.52
1:A:127:LYS:HD3	1:C:141:ARG:HG3	1.90	0.52
2:B:144:LYS:O	2:B:145:TYR:CD2	2.62	0.52
2:D:28:LEU:HD23	2:D:60:VAL:HG13	1.90	0.52
1:A:83:LEU:HD11	3:A:200:HEM:HMA3	1.90	0.52
1:A:86:LEU:HD12	1:A:90:LYS:HD3	1.91	0.52
2:D:92:HIS:HE1	3:D:200:HEM:NA	2.05	0.52
1:A:15:GLY:C	1:A:17:ILE:H	2.13	0.52
1:A:32:MET:HG2	1:A:39:THR:HG21	1.91	0.51
2:B:63:HIS:HE2	3:B:200:HEM:CHA	2.24	0.51
1:C:90:LYS:HG3	1:C:91:LEU:HD13	1.93	0.51
2:D:102:ASN:HA	2:D:105:LEU:HB2	1.91	0.51
1:A:17:ILE:O	1:A:20:HIS:HB2	2.10	0.51
1:A:27:GLU:OE1	1:A:108:THR:HG23	2.10	0.51
1:A:111:SER:O	2:B:119:GLY:HA2	2.11	0.51
2:D:40:ARG:O	2:D:41:TYR:CG	2.63	0.51
2:D:10:ALA:O	2:D:14:LEU:HB2	2.11	0.51
1:C:2:LEU:HD21	1:C:127:LYS:HB3	1.93	0.51
2:D:80:SER:O	2:D:82:LYS:N	2.38	0.51
1:C:94:ASP:HB3	1:C:97:ASN:OD1	2.10	0.51
2:B:67:VAL:HG22	3:B:200:HEM:C1B	2.45	0.50
2:D:72:ASN:O	2:D:75:LEU:N	2.43	0.50
2:D:35:TYR:O	2:D:38:THR:HB	2.11	0.50
1:A:24:TYR:OH	1:A:113:HIS:CE1	2.65	0.50
2:D:122:PHE:CZ	2:D:127:GLN:HA	2.46	0.50
2:B:122:PHE:CE2	2:B:127:GLN:HG3	2.47	0.50
1:C:3:SER:CA	1:C:6:ASP:HB2	2.42	0.50
2:D:21:ASP:OD1	2:D:22:GLU:N	2.45	0.49
2:D:57:ASN:HD22	2:D:59:LYS:H	1.60	0.49
2:B:129:ALA:O	2:B:133:VAL:HG23	2.11	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:31:LEU:HG	2:B:38:THR:HG21	1.94	0.49
2:B:100:PRO:HB3	2:B:142:ALA:HB2	1.93	0.49
1:C:103:HIS:HA	1:C:106:LEU:HD12	1.95	0.49
1:A:17:ILE:HG12	1:A:24:TYR:CD2	2.48	0.49
1:A:85:ASP:C	1:A:85:ASP:OD2	2.50	0.49
2:D:10:ALA:O	2:D:14:LEU:N	2.42	0.49
1:A:102:SER:HB2	1:A:129:LEU:HD13	1.94	0.49
2:D:89:SER:OG	2:D:144:LYS:HB2	2.13	0.48
2:B:82:LYS:H	2:B:82:LYS:HD2	1.78	0.48
1:A:15:GLY:O	1:A:18:GLY:N	2.44	0.48
1:C:79:ALA:C	1:C:81:SER:H	2.16	0.48
1:A:127:LYS:CG	1:C:141:ARG:HG3	2.43	0.48
2:B:40:ARG:HB3	1:C:92:ARG:HD3	1.95	0.48
2:D:123:THR:O	2:D:126:ALA:N	2.44	0.48
1:A:102:SER:HB2	1:A:129:LEU:HB3	1.96	0.48
2:B:146:HIS:O	2:B:146:HIS:CD2	2.66	0.47
1:C:24:TYR:HE1	1:C:112:HIS:CG	2.32	0.47
2:B:14:LEU:HD12	2:B:14:LEU:HA	1.68	0.47
2:B:70:ALA:HA	2:B:73:ASP:OD2	2.14	0.47
2:B:97:HIS:CG	1:C:41:THR:HG21	2.50	0.47
2:D:123:THR:O	2:D:127:GLN:N	2.39	0.47
1:A:22:ALA:CB	1:A:60:LYS:HE2	2.44	0.47
1:C:45:HIS:ND1	1:C:45:HIS:N	2.55	0.47
1:A:94:ASP:O	1:A:97:ASN:OD1	2.33	0.47
1:A:140:TYR:HE1	2:D:37:TRP:CZ3	2.32	0.47
1:A:10:ILE:HD13	1:A:128:PHE:HB2	1.97	0.47
1:C:3:SER:H	1:C:6:ASP:HB2	1.80	0.47
1:C:107:VAL:HG11	2:D:127:GLN:OE1	2.15	0.47
2:B:65:LYS:O	2:B:68:ILE:HG13	2.15	0.47
1:C:66:LEU:HD21	3:C:200:HEM:HBB2	1.97	0.46
1:A:9:ASN:ND2	1:A:9:ASN:H	2.13	0.46
2:B:28:LEU:O	2:B:32:LEU:HD23	2.15	0.46
1:C:137:THR:HA	1:C:140:TYR:CE2	2.50	0.46
2:D:48:LEU:O	2:D:54:ILE:HG12	2.15	0.46
2:B:77:HIS:HE1	4:B:155:HOH:O	1.97	0.46
1:A:17:ILE:HG12	1:A:24:TYR:CE2	2.51	0.46
1:A:31:ARG:HD3	2:B:127:GLN:OE1	2.14	0.46
1:C:76:LEU:N	1:C:77:PRO:CD	2.79	0.46
2:D:40:ARG:O	2:D:41:TYR:CD2	2.68	0.46
2:D:76:ASN:OD1	2:D:76:ASN:N	2.48	0.46
2:D:127:GLN:CG	2:D:128:ALA:N	2.78	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:VAL:HG12	1:A:48:VAL:O	2.15	0.46
2:B:68:ILE:C	2:B:71:PHE:H	2.19	0.46
2:D:45:PHE:CE1	2:D:59:LYS:HB3	2.50	0.46
1:A:83:LEU:HD11	3:A:200:HEM:CMA	2.45	0.46
2:B:26:GLU:OE1	2:B:113:VAL:HG13	2.16	0.46
2:B:146:HIS:CD2	2:B:146:HIS:C	2.89	0.46
1:C:17:ILE:HG23	1:C:24:TYR:HE2	1.80	0.46
1:C:22:ALA:HB1	1:C:60:LYS:HB2	1.98	0.46
1:C:35:SER:HA	2:D:128:ALA:CB	2.46	0.46
2:B:23:VAL:O	2:B:27:ALA:N	2.43	0.45
1:C:37:PRO:O	1:C:40:LYS:HB2	2.17	0.45
1:C:42:TYR:O	3:C:200:HEM:CMD	2.65	0.45
1:A:121:VAL:O	1:A:122:HIS:C	2.54	0.45
1:C:113:HIS:N	1:C:114:PRO:HD3	2.31	0.45
2:D:41:TYR:CE1	3:D:200:HEM:HAC	2.42	0.45
2:D:45:PHE:CD1	2:D:59:LYS:HB3	2.52	0.45
1:A:80:LEU:HB2	1:A:135:VAL:HG11	1.99	0.45
2:B:18:VAL:O	2:B:18:VAL:HG12	2.17	0.45
2:D:34:VAL:HG23	2:D:35:TYR:CE2	2.50	0.45
1:A:72:HIS:O	1:A:76:LEU:HD13	2.16	0.45
1:A:140:TYR:HE1	2:D:37:TRP:HZ3	1.63	0.45
2:B:114:LEU:O	2:B:118:LEU:HB2	2.18	0.44
2:D:3:LEU:HB3	2:D:8:LYS:HB2	1.99	0.44
1:C:3:SER:OG	1:C:4:GLY:HA2	2.16	0.44
2:D:107:GLY:CA	2:D:134:VAL:HG13	2.46	0.44
1:A:15:GLY:O	1:A:17:ILE:N	2.51	0.44
1:A:116:ASP:O	1:A:118:THR:N	2.49	0.44
1:C:98:PHE:CE2	1:C:137:THR:CG2	3.01	0.44
2:D:82:LYS:N	2:D:82:LYS:HD2	2.33	0.44
1:A:42:TYR:OH	1:A:94:ASP:HB2	2.18	0.44
1:A:31:ARG:HA	2:B:124:PRO:HB3	1.99	0.44
1:C:140:TYR:N	1:C:140:TYR:CD2	2.86	0.44
2:D:24:GLY:H	2:D:68:ILE:HD11	1.79	0.44
2:D:100:PRO:HD3	2:D:145:TYR:CZ	2.50	0.44
1:A:8:SER:O	1:A:12:ALA:CB	2.66	0.44
1:A:35:SER:HB2	1:A:36:PHE:CE2	2.52	0.44
2:D:33:VAL:HG22	2:D:54:ILE:HD12	2.00	0.44
2:D:94:ASP:O	2:D:95:LYS:CB	2.65	0.44
2:D:106:LEU:HD12	3:D:200:HEM:CHC	2.48	0.44
3:D:200:HEM:HBA1	3:D:200:HEM:CMA	2.48	0.44
2:B:28:LEU:O	2:B:31:LEU:HB3	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:24:TYR:O	1:C:25:GLY:C	2.57	0.43
1:C:86:LEU:O	1:C:91:LEU:HB2	2.18	0.43
2:B:73:ASP:O	2:B:77:HIS:HB2	2.18	0.43
1:C:35:SER:HA	2:D:128:ALA:HB1	2.00	0.43
1:C:54:GLN:HB2	4:C:142:HOH:O	2.19	0.43
2:B:22:GLU:O	2:B:26:GLU:HB2	2.18	0.43
1:C:93:VAL:CG1	3:C:200:HEM:HBC2	2.49	0.43
2:B:68:ILE:O	2:B:69:THR:C	2.56	0.43
1:C:3:SER:N	1:C:6:ASP:HB2	2.34	0.43
2:B:20:SER:HB3	4:B:149:HOH:O	2.18	0.43
1:C:24:TYR:HH	1:C:113:HIS:HE2	1.65	0.43
1:C:47:ASP:OD2	1:C:47:ASP:C	2.57	0.43
1:A:87:HIS:HA	1:A:91:LEU:HB2	2.01	0.43
1:A:37:PRO:O	1:A:40:LYS:HB2	2.19	0.43
2:D:98:VAL:O	2:D:145:TYR:HE1	2.02	0.42
2:B:40:ARG:HB3	1:C:92:ARG:HD2	2.01	0.42
1:C:3:SER:HA	1:C:6:ASP:HB2	2.00	0.42
2:D:57:ASN:ND2	2:D:59:LYS:H	2.16	0.42
1:C:42:TYR:CE1	1:C:93:VAL:HA	2.55	0.42
2:D:41:TYR:CD2	2:D:41:TYR:N	2.88	0.42
1:A:69:ALA:HB2	1:A:80:LEU:HD21	2.00	0.42
2:B:99:ASP:HA	2:B:100:PRO:HD3	1.77	0.42
2:B:92:HIS:O	2:B:93:CYS:CB	2.68	0.42
1:C:9:ASN:HA	1:C:12:ALA:HB3	2.01	0.42
2:B:39:GLN:HA	2:B:42:PHE:HD2	1.84	0.42
2:B:103:PHE:HD1	3:B:200:HEM:HBB2	1.85	0.42
2:B:24:GLY:H	2:B:68:ILE:HD13	1.85	0.41
2:D:107:GLY:CA	2:D:134:VAL:CG1	2.96	0.41
2:B:36:PRO:C	2:B:38:THR:H	2.23	0.41
2:B:92:HIS:HE1	3:B:200:HEM:NA	2.01	0.41
2:B:130:PHE:HA	2:B:133:VAL:HB	2.02	0.41
1:C:81:SER:HA	1:C:84:SER:HB2	2.01	0.41
2:D:23:VAL:HG13	2:D:113:VAL:CG1	2.51	0.41
2:D:74:GLY:O	2:D:77:HIS:N	2.53	0.41
1:A:58:HIS:HE1	3:A:200:HEM:CHA	2.33	0.41
1:C:66:LEU:O	1:C:69:ALA:HB3	2.20	0.41
1:A:137:THR:HG22	1:A:137:THR:O	2.20	0.41
2:D:99:ASP:HA	2:D:100:PRO:HD2	1.79	0.41
2:B:45:PHE:N	2:B:45:PHE:CD2	2.89	0.41
1:C:117:PHE:O	2:D:30:ARG:NH2	2.51	0.41
1:A:27:GLU:O	1:A:31:ARG:HG3	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:HIS:CE1	3:A:200:HEM:CHA	3.04	0.41
2:D:63:HIS:CE1	2:D:67:VAL:HG23	2.56	0.41
1:C:97:ASN:HA	1:C:100:LEU:HB2	2.03	0.41
1:A:95:PRO:O	1:A:98:PHE:HD2	2.03	0.40
1:A:137:THR:HG23	1:A:140:TYR:OH	2.21	0.40
2:B:132:LYS:HE3	2:B:132:LYS:HB2	1.58	0.40
1:C:42:TYR:CD2	3:C:200:HEM:HBC1	2.57	0.40
1:C:49:SER:O	1:C:50:HIS:C	2.59	0.40
1:C:66:LEU:HD21	3:C:200:HEM:CBB	2.50	0.40
1:C:123:ALA:HA	2:D:34:VAL:HG12	2.03	0.40
1:A:141:ARG:NH2	1:C:1:VAL:HG11	2.37	0.40
1:C:97:ASN:O	1:C:100:LEU:HB2	2.22	0.40
3:D:200:HEM:HBA1	3:D:200:HEM:HMA1	2.03	0.40
1:A:6:ASP:OD2	1:A:127:LYS:CE	2.50	0.40
1:C:59:GLY:O	1:C:62:VAL:HB	2.21	0.40
1:C:97:ASN:HA	1:C:100:LEU:HD22	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	139/141 (99%)	123 (88%)	13 (9%)	3 (2%)	6 22
1	C	139/141 (99%)	127 (91%)	10 (7%)	2 (1%)	11 34
2	B	144/146 (99%)	121 (84%)	17 (12%)	6 (4%)	3 9
2	D	144/146 (99%)	124 (86%)	14 (10%)	6 (4%)	3 9
All	All	566/574 (99%)	495 (88%)	54 (10%)	17 (3%)	4 15

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	GLY
2	B	77	HIS
2	B	93	CYS
2	D	81	LEU
2	D	95	LYS
1	A	72	HIS
2	B	19	ASN
2	B	20	SER
2	B	47	ASP
1	A	16	LYS
2	D	21	ASP
2	D	121	ASP
1	C	3	SER
1	C	2	LEU
2	B	88	LEU
2	D	119	GLY
2	D	100	PRO

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	110/110 (100%)	90 (82%)	20 (18%)	1 5
1	C	110/110 (100%)	92 (84%)	18 (16%)	2 7
2	B	115/115 (100%)	91 (79%)	24 (21%)	1 3
2	D	115/115 (100%)	91 (79%)	24 (21%)	1 3
All	All	450/450 (100%)	364 (81%)	86 (19%)	1 4

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	17	ILE
1	A	38	THR
1	A	41	THR
1	A	45	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	47	ASP
1	A	60	LYS
1	A	72	HIS
1	A	74	ASP
1	A	76	LEU
1	A	85	ASP
1	A	92	ARG
1	A	94	ASP
1	A	102	SER
1	A	111	SER
1	A	116	ASP
1	A	125	LEU
1	A	128	PHE
1	A	131	SER
1	A	141	ARG
2	B	4	THR
2	B	8	LYS
2	B	11	VAL
2	B	14	LEU
2	B	18	VAL
2	B	26	GLU
2	B	32	LEU
2	B	40	ARG
2	B	44	SER
2	B	76	ASN
2	B	77	HIS
2	B	79	ASP
2	B	80	SER
2	B	81	LEU
2	B	82	LYS
2	B	88	LEU
2	B	91	LEU
2	B	106	LEU
2	B	111	VAL
2	B	118	LEU
2	B	123	THR
2	B	132	LYS
2	B	145	TYR
2	B	146	HIS
1	C	7	LYS
1	C	9	ASN
1	C	23	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	35	SER
1	C	45	HIS
1	C	47	ASP
1	C	60	LYS
1	C	76	LEU
1	C	83	LEU
1	C	84	SER
1	C	91	LEU
1	C	99	LYS
1	C	100	LEU
1	C	101	LEU
1	C	105	LEU
1	C	124	SER
1	C	125	LEU
1	C	140	TYR
2	D	2	HIS
2	D	13	CYS
2	D	20	SER
2	D	21	ASP
2	D	26	GLU
2	D	28	LEU
2	D	38	THR
2	D	41	TYR
2	D	43	ASP
2	D	57	ASN
2	D	75	LEU
2	D	76	ASN
2	D	80	SER
2	D	81	LEU
2	D	82	LYS
2	D	87	SER
2	D	88	LEU
2	D	94	ASP
2	D	104	ARG
2	D	106	LEU
2	D	118	LEU
2	D	120	LYS
2	D	127	GLN
2	D	146	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	50	HIS
1	A	58	HIS
1	A	103	HIS
1	A	113	HIS
2	B	39	GLN
2	B	76	ASN
2	B	102	ASN
2	B	108	ASN
2	B	131	GLN
2	B	146	HIS
1	C	9	ASN
1	C	58	HIS
2	D	57	ASN
2	D	63	HIS
2	D	116	HIS
2	D	131	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

4 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	$\#\|Z\| > 2$
3	HEM	D	200	4,2	41,50,50	2.05	7 (17%)	45,82,82	2.06	11 (24%)
3	HEM	B	200	2	41,50,50	2.01	6 (14%)	45,82,82	1.86	15 (33%)
3	HEM	C	200	1,4	41,50,50	1.93	7 (17%)	45,82,82	1.70	10 (22%)
3	HEM	A	200	1,4	41,50,50	1.95	7 (17%)	45,82,82	1.75	9 (20%)

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	HEM	D	200	4,2	-	10/12/54/54	-
3	HEM	B	200	2	-	6/12/54/54	-
3	HEM	C	200	1,4	-	9/12/54/54	-
3	HEM	A	200	1,4	-	6/12/54/54	-

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	200	HEM	C3D-C2D	7.81	1.53	1.36
3	A	200	HEM	C3D-C2D	7.76	1.53	1.36
3	B	200	HEM	C3D-C2D	7.64	1.53	1.36
3	C	200	HEM	C3D-C2D	7.33	1.52	1.36
3	A	200	HEM	C3C-C2C	-5.07	1.33	1.40
3	D	200	HEM	C3C-CAC	4.68	1.57	1.47
3	B	200	HEM	C3C-C2C	-4.60	1.34	1.40
3	D	200	HEM	C3C-C2C	-4.28	1.34	1.40
3	C	200	HEM	C3C-C2C	-4.24	1.34	1.40
3	B	200	HEM	FE-ND	4.16	2.17	1.96
3	B	200	HEM	C3C-CAC	4.02	1.56	1.47
3	C	200	HEM	C3C-CAC	3.67	1.55	1.47
3	D	200	HEM	FE-ND	3.58	2.14	1.96
3	A	200	HEM	C3C-CAC	3.53	1.55	1.47
3	C	200	HEM	CAA-C2A	3.37	1.57	1.52
3	C	200	HEM	CAB-C3B	2.82	1.55	1.47
3	D	200	HEM	CAB-C3B	2.78	1.55	1.47
3	B	200	HEM	CAB-C3B	2.67	1.54	1.47
3	A	200	HEM	CAB-C3B	2.43	1.54	1.47
3	A	200	HEM	CAA-C2A	2.42	1.55	1.52
3	C	200	HEM	FE-NB	2.35	2.08	1.96

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	200	HEM	CMD-C2D	2.32	1.55	1.50
3	B	200	HEM	C3B-C2B	-2.10	1.33	1.37
3	A	200	HEM	C3B-C2B	-2.09	1.33	1.37
3	D	200	HEM	O1D-CGD	2.05	1.28	1.22
3	C	200	HEM	CMB-C2B	2.03	1.55	1.50
3	D	200	HEM	FE-NB	2.01	2.06	1.96

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	200	HEM	C4D-ND-C1D	6.22	111.49	105.07
3	B	200	HEM	C4D-ND-C1D	6.11	111.38	105.07
3	D	200	HEM	C4D-ND-C1D	6.06	111.33	105.07
3	A	200	HEM	C4D-ND-C1D	5.97	111.24	105.07
3	D	200	HEM	C4C-CHD-C1D	4.88	129.00	122.56
3	D	200	HEM	CAD-C3D-C4D	4.72	132.91	124.66
3	C	200	HEM	CMA-C3A-C4A	-4.16	122.08	128.46
3	D	200	HEM	CMA-C3A-C4A	-4.04	122.25	128.46
3	A	200	HEM	C4C-CHD-C1D	3.81	127.58	122.56
3	B	200	HEM	C4C-CHD-C1D	3.50	127.18	122.56
3	D	200	HEM	C3B-C2B-C1B	3.33	108.96	106.49
3	B	200	HEM	CHC-C4B-C3B	3.20	129.47	124.57
3	C	200	HEM	CAD-C3D-C4D	2.98	129.87	124.66
3	B	200	HEM	CMA-C3A-C4A	-2.86	124.06	128.46
3	B	200	HEM	CMD-C2D-C1D	2.85	129.38	125.04
3	B	200	HEM	CBD-CAD-C3D	-2.82	104.79	112.63
3	A	200	HEM	CBA-CAA-C2A	2.78	117.37	112.62
3	B	200	HEM	C3B-C2B-C1B	2.70	108.49	106.49
3	C	200	HEM	C4A-C3A-C2A	2.63	108.83	107.00
3	D	200	HEM	CAD-C3D-C2D	-2.59	123.06	127.88
3	A	200	HEM	CMD-C2D-C1D	2.57	128.96	125.04
3	C	200	HEM	C4C-CHD-C1D	2.57	125.95	122.56
3	A	200	HEM	C1B-NB-C4B	2.56	107.72	105.07
3	D	200	HEM	C1B-NB-C4B	2.56	107.71	105.07
3	D	200	HEM	CMA-C3A-C2A	2.55	129.75	124.94
3	B	200	HEM	C4B-CHC-C1C	2.44	125.78	122.56
3	B	200	HEM	CHC-C4B-NB	-2.23	122.01	124.43
3	C	200	HEM	CHD-C1D-ND	2.20	126.83	124.43
3	C	200	HEM	CMA-C3A-C2A	2.20	129.09	124.94
3	D	200	HEM	CHC-C4B-C3B	2.20	127.94	124.57
3	D	200	HEM	O1A-CGA-CBA	-2.19	116.03	123.08
3	C	200	HEM	O1D-CGD-CBD	-2.19	116.04	123.08

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	200	HEM	C1B-NB-C4B	2.18	107.33	105.07
3	B	200	HEM	CAA-CBA-CGA	-2.15	107.75	113.76
3	A	200	HEM	CBD-CAD-C3D	-2.14	106.69	112.63
3	B	200	HEM	CAD-C3D-C4D	2.12	128.36	124.66
3	A	200	HEM	C3B-C2B-C1B	2.11	108.05	106.49
3	C	200	HEM	CBD-CAD-C3D	-2.10	106.80	112.63
3	C	200	HEM	CMD-C2D-C1D	2.09	128.22	125.04
3	A	200	HEM	CAD-CBD-CGD	-2.08	109.12	113.60
3	A	200	HEM	CBB-CAB-C3B	-2.05	117.42	127.62
3	B	200	HEM	C3C-C4C-NC	-2.04	107.09	110.94
3	D	200	HEM	O2A-CGA-CBA	2.03	120.56	114.03
3	B	200	HEM	CBA-CAA-C2A	-2.02	109.18	112.62
3	B	200	HEM	CHD-C1D-C2D	2.01	128.11	124.98

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	200	HEM	C2B-C3B-CAB-CBB
3	A	200	HEM	C4B-C3B-CAB-CBB
3	B	200	HEM	C2A-CAA-CBA-CGA
3	C	200	HEM	C1A-C2A-CAA-CBA
3	C	200	HEM	C3A-C2A-CAA-CBA
3	D	200	HEM	C1A-C2A-CAA-CBA
3	D	200	HEM	C3A-C2A-CAA-CBA
3	D	200	HEM	C2D-C3D-CAD-CBD
3	D	200	HEM	C4D-C3D-CAD-CBD
3	C	200	HEM	C3D-CAD-CBD-CGD
3	D	200	HEM	C3D-CAD-CBD-CGD
3	C	200	HEM	C2A-CAA-CBA-CGA
3	B	200	HEM	C3D-CAD-CBD-CGD
3	D	200	HEM	C4B-C3B-CAB-CBB
3	C	200	HEM	C2D-C3D-CAD-CBD
3	D	200	HEM	CAA-CBA-CGA-O2A
3	D	200	HEM	CAA-CBA-CGA-O1A
3	D	200	HEM	CAD-CBD-CGD-O1D
3	A	200	HEM	C3D-CAD-CBD-CGD
3	B	200	HEM	CAD-CBD-CGD-O2D
3	C	200	HEM	CAA-CBA-CGA-O2A
3	B	200	HEM	CAA-CBA-CGA-O2A
3	A	200	HEM	CAD-CBD-CGD-O2D
3	D	200	HEM	CAD-CBD-CGD-O2D

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	C	200	HEM	C4D-C3D-CAD-CBD
3	C	200	HEM	CAA-CBA-CGA-O1A
3	B	200	HEM	CAD-CBD-CGD-O1D
3	B	200	HEM	CAA-CBA-CGA-O1A
3	A	200	HEM	CAD-CBD-CGD-O1D
3	C	200	HEM	CAD-CBD-CGD-O2D
3	A	200	HEM	CAA-CBA-CGA-O2A

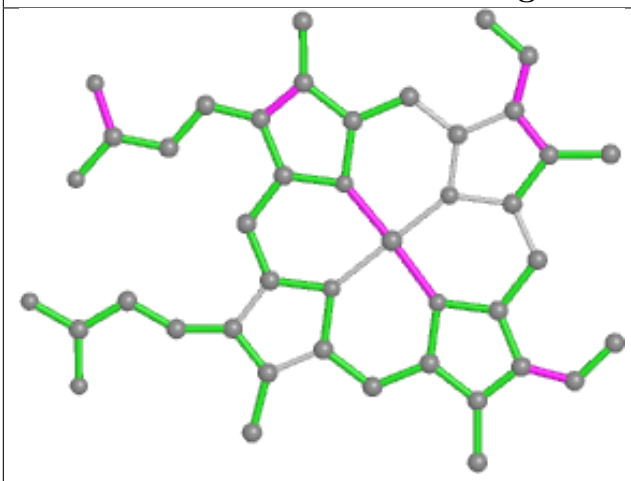
There are no ring outliers.

4 monomers are involved in 31 short contacts:

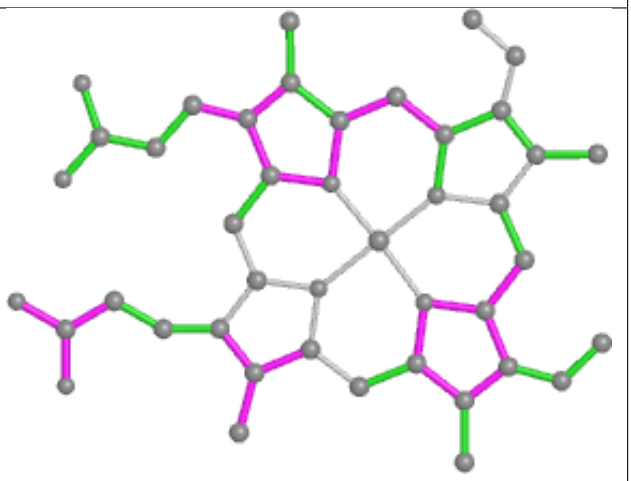
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	200	HEM	7	0
3	B	200	HEM	7	0
3	C	200	HEM	9	0
3	A	200	HEM	8	0

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 then 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.

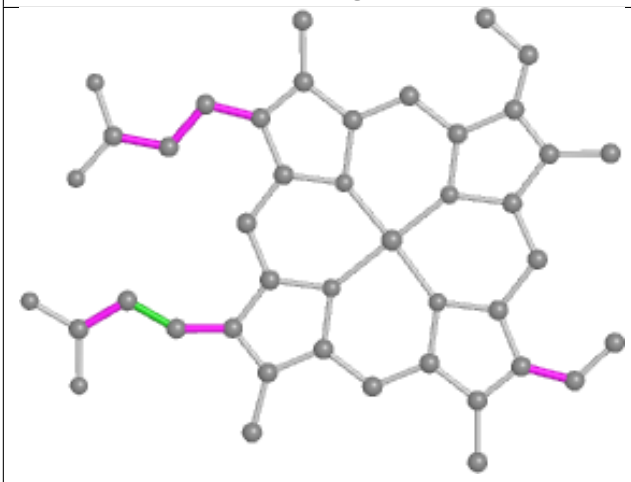
Ligand HEM D 200



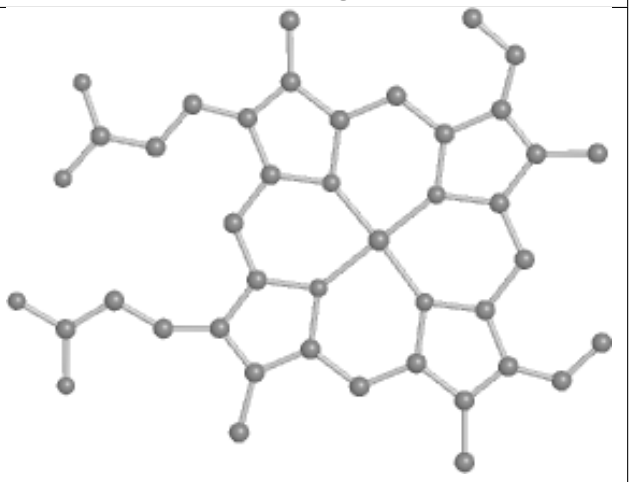
Bond lengths



Bond angles

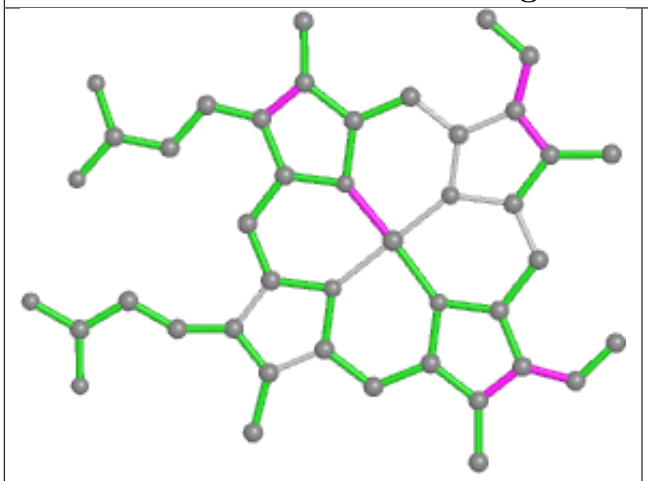


Torsions

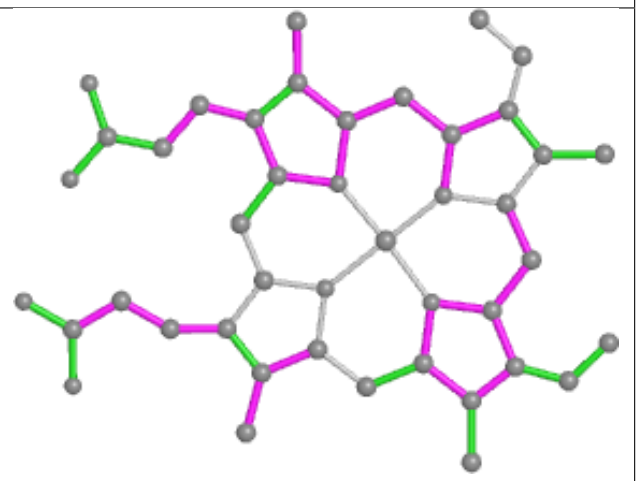


Rings

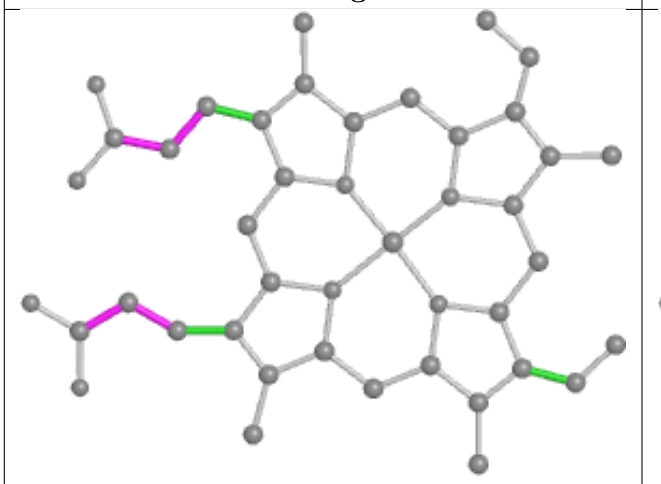
Ligand HEM B 200



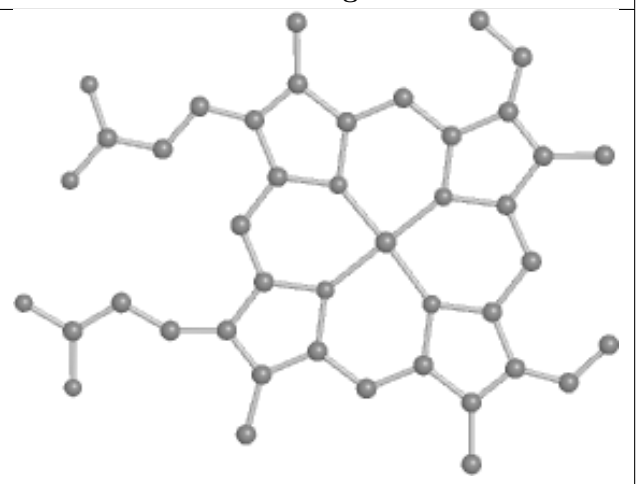
Bond lengths



Bond angles

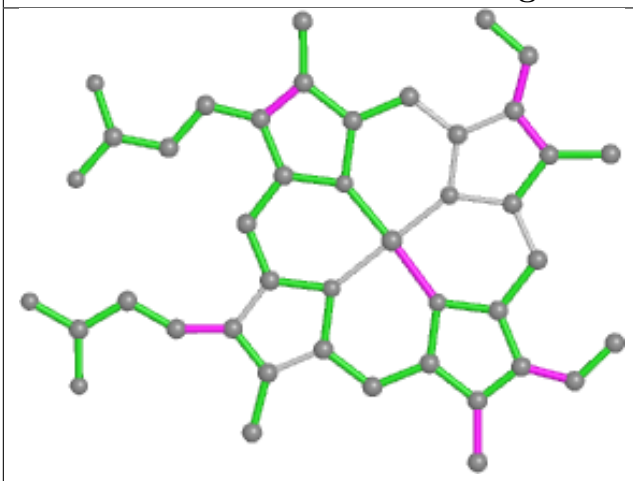


Torsions

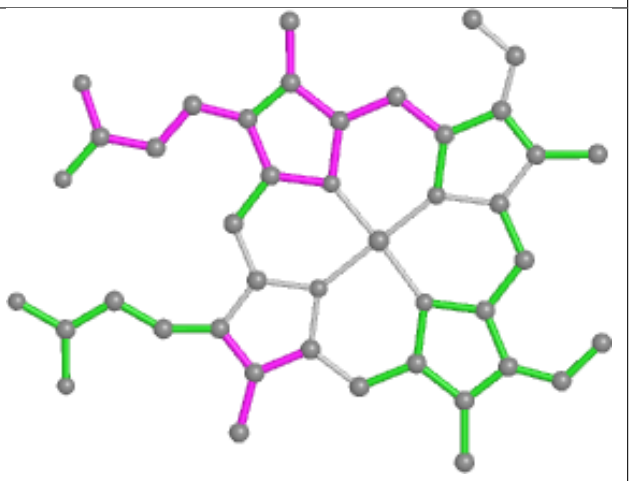


Rings

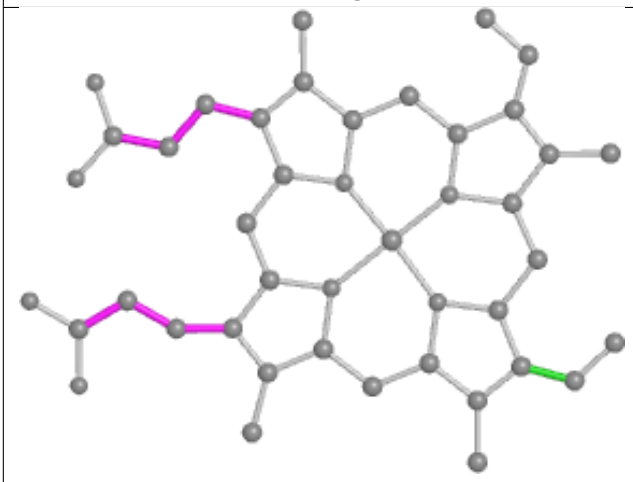
Ligand HEM C 200



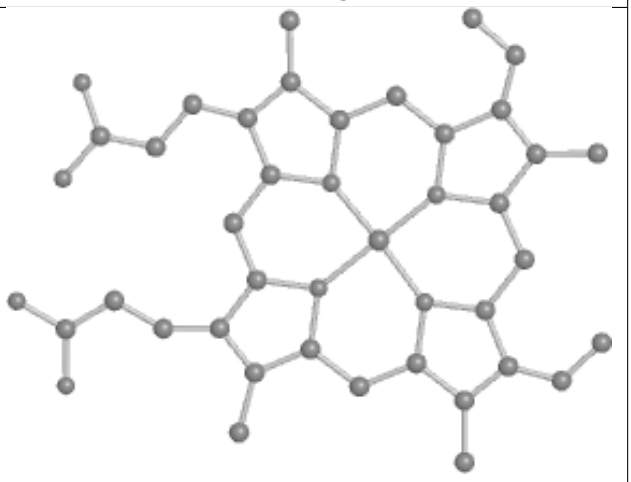
Bond lengths



Bond angles

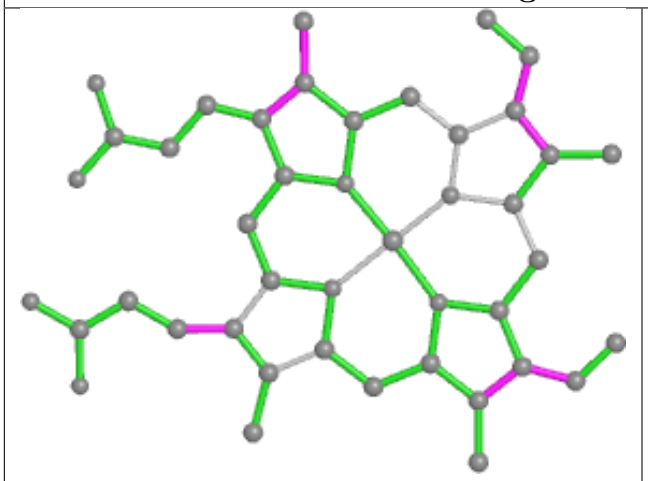


Torsions

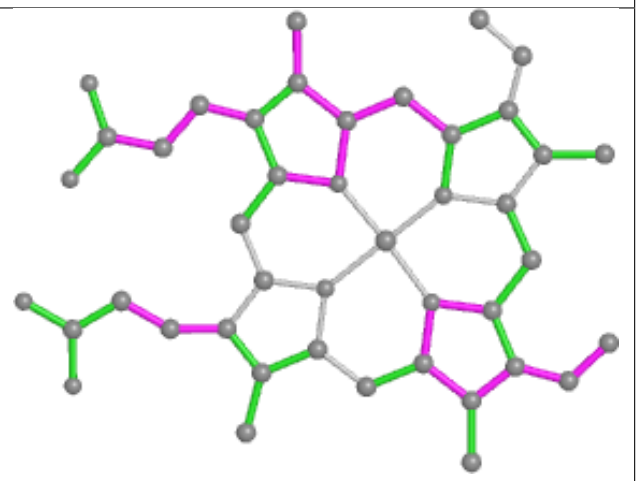


Rings

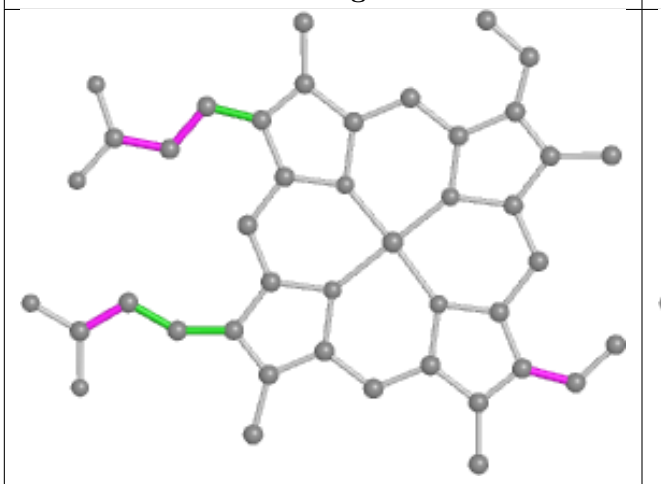
Ligand HEM A 200



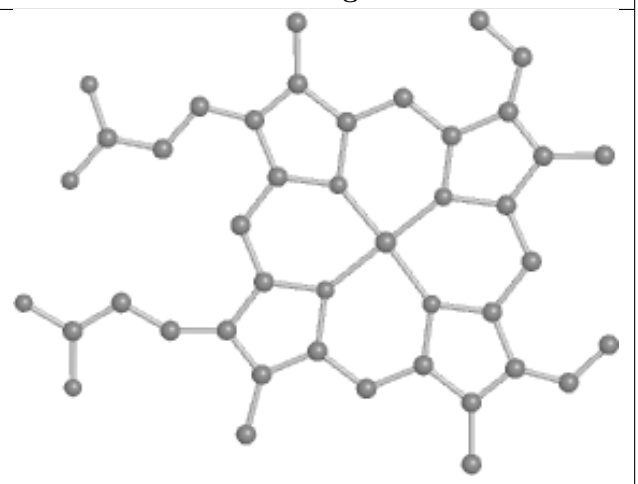
Bond lengths



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 (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	141/141 (100%)	-0.22	4 (2%)	53	43	22, 34, 49, 53	0
1	C	141/141 (100%)	-0.08	5 (3%)	44	34	20, 38, 64, 68	0
2	B	146/146 (100%)	-0.20	4 (2%)	54	44	15, 34, 51, 56	0
2	D	146/146 (100%)	-0.20	5 (3%)	45	35	21, 33, 53, 65	0
All	All	574/574 (100%)	-0.17	18 (3%)	49	39	15, 34, 56, 68	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	146	HIS	7.5
1	A	141	ARG	4.6
2	B	2	HIS	4.2
2	B	145	TYR	3.9
2	D	2	HIS	3.7
1	A	20	HIS	3.5
1	C	19	GLY	3.3
1	A	140	TYR	3.0
2	D	5	ASP	3.0
2	B	93	CYS	3.0
2	D	4	THR	3.0
1	C	75	ASP	2.6
1	C	1	VAL	2.5
2	D	6	ALA	2.5
2	D	146	HIS	2.4
1	C	138	SER	2.3
1	C	141	ARG	2.3
1	A	72	HIS	2.2

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

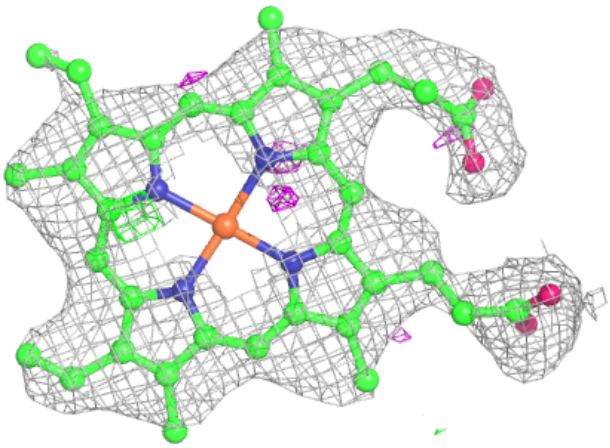
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

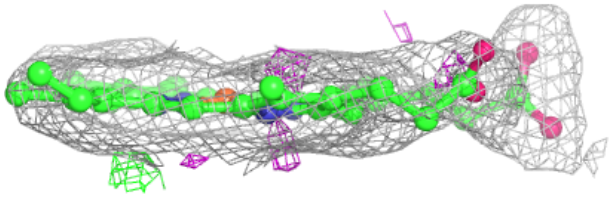
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	HEM	D	200	43/43	0.87	0.23	20,23,30,32	0
3	HEM	A	200	43/43	0.91	0.22	27,28,32,33	0
3	HEM	B	200	43/43	0.92	0.18	32,34,36,40	0
3	HEM	C	200	43/43	0.93	0.17	27,33,42,44	0

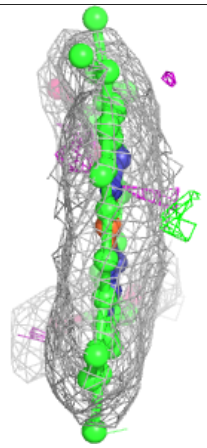
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HEM D 200:

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

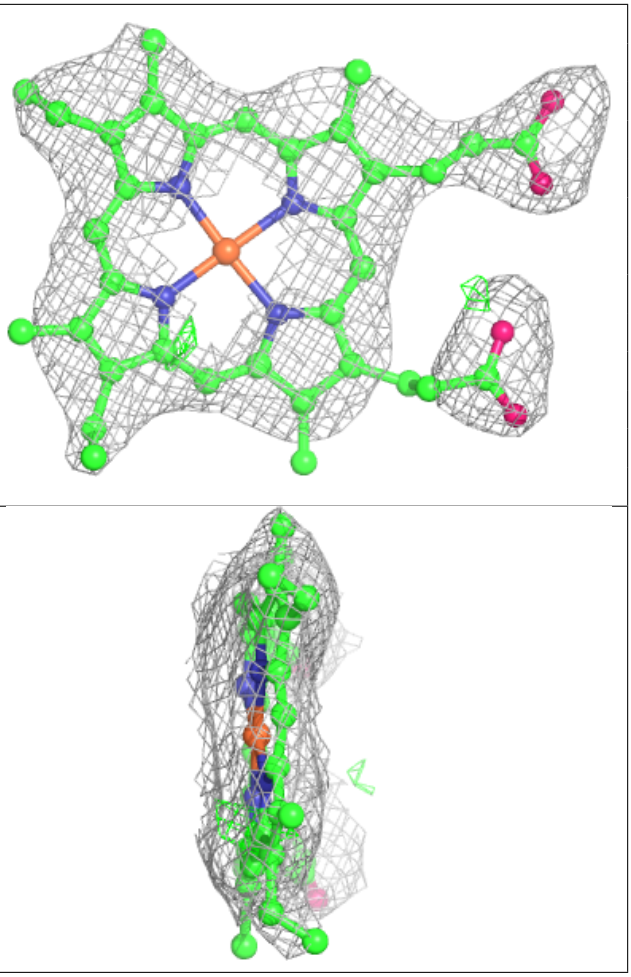






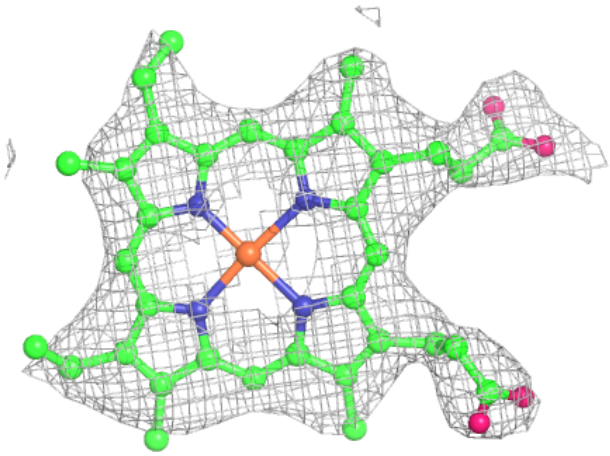
Electron density around HEM A 200:

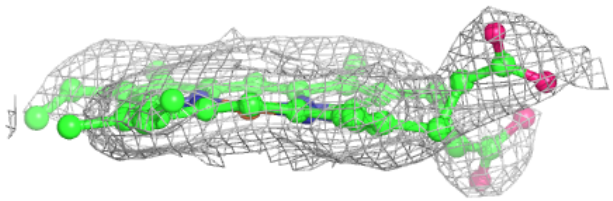
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

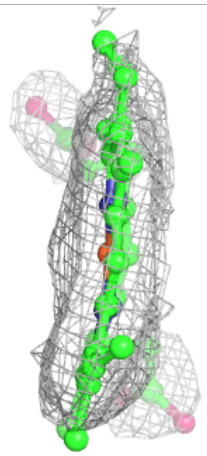


Electron density around HEM B 200:

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

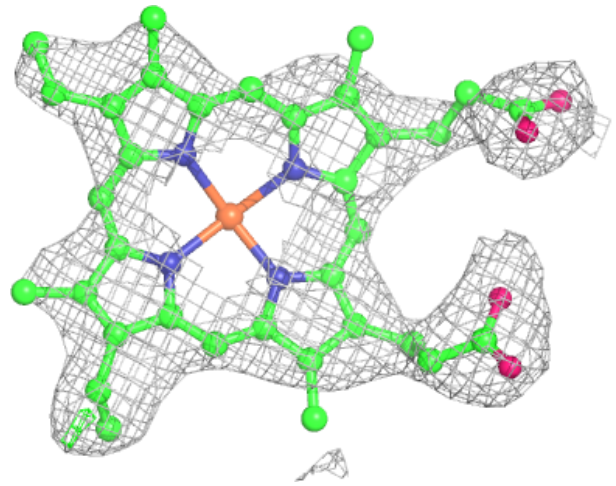


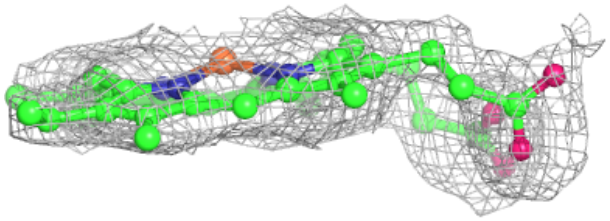


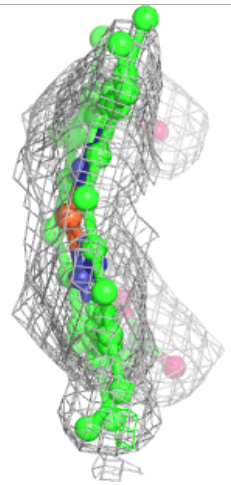


Electron density around HEM C 200:

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.