



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:21 PM UTC

PDB ID : 3TTU / pdb_00003ttu
Title : Structure of F413Y/H128N double variant of E. coli KatE
Authors : Loewen, P.C.; Jha, V.
Deposited on : 2011-09-15
Resolution : 1.89 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

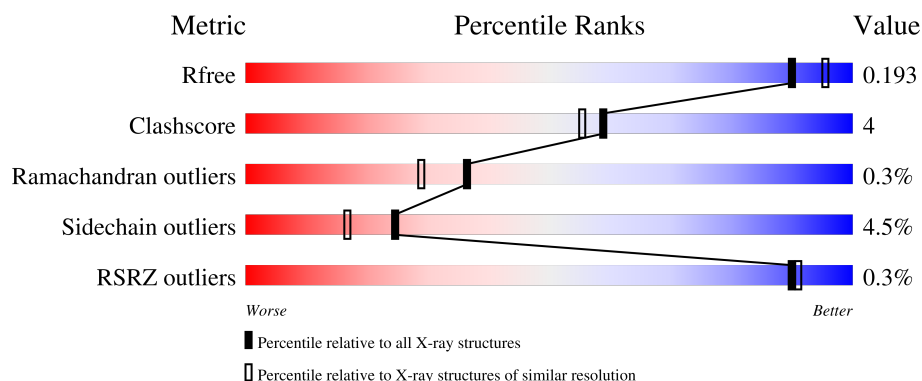
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	753	86% 9% ...
1	B	753	84% 10% ...
1	C	753	83% 12% ...
1	D	753	85% 9% ...

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 25986 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 Catalase HP11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	726	Total	C	N	O	S	0	1	0
			5743	3645	1005	1081	12			
1	B	726	Total	C	N	O	S	0	1	0
			5744	3646	1005	1081	12			
1	C	726	Total	C	N	O	S	0	1	0
			5743	3645	1005	1081	12			
1	D	726	Total	C	N	O	S	0	1	0
			5744	3646	1005	1081	12			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	128	ASN	HIS	engineered mutation	UNP P21179
A	413	TYR	PHE	engineered mutation	UNP P21179
B	128	ASN	HIS	engineered mutation	UNP P21179
B	413	TYR	PHE	engineered mutation	UNP P21179
C	128	ASN	HIS	engineered mutation	UNP P21179
C	413	TYR	PHE	engineered mutation	UNP P21179
D	128	ASN	HIS	engineered mutation	UNP P21179
D	413	TYR	PHE	engineered mutation	UNP P21179

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

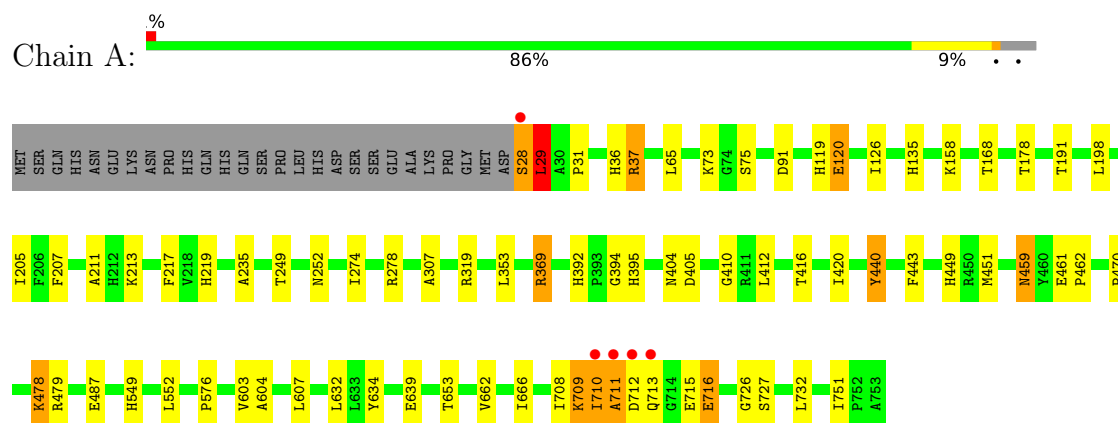
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	772	Total	O	0	0
			772	772		
3	B	639	Total	O	0	0
			639	639		
3	C	680	Total	O	0	0
			680	680		
3	D	749	Total	O	0	0
			749	749		

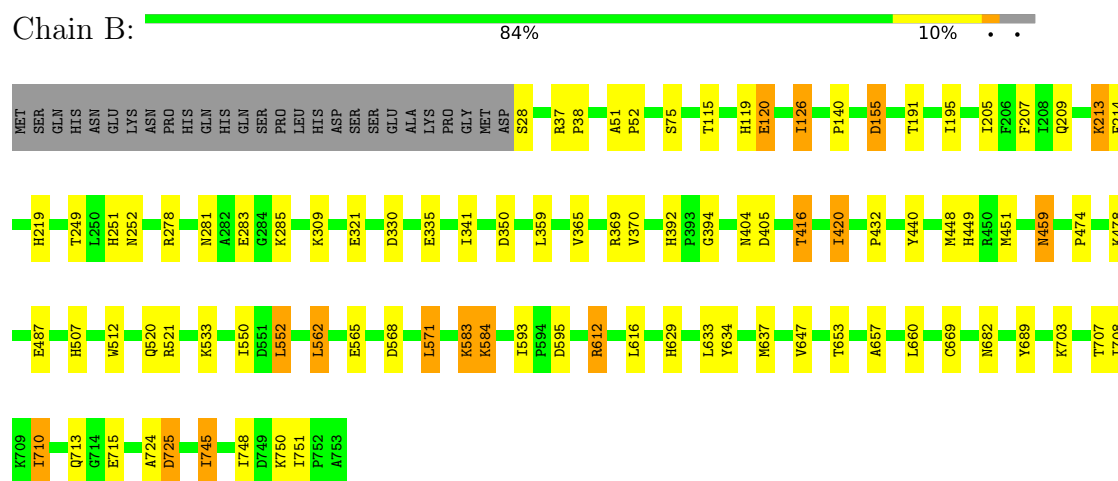
3 Residue-property plots

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.

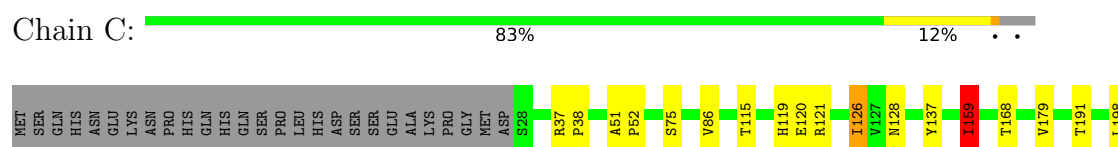
• Molecule 1: Catalase HP11

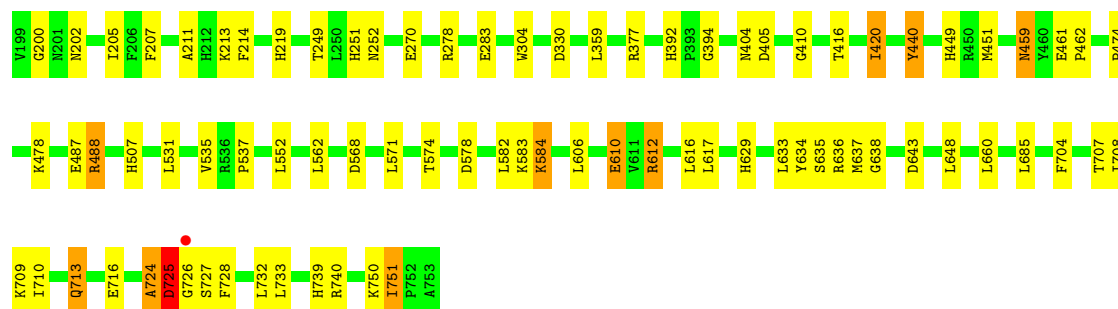


• Molecule 1: Catalase HP11



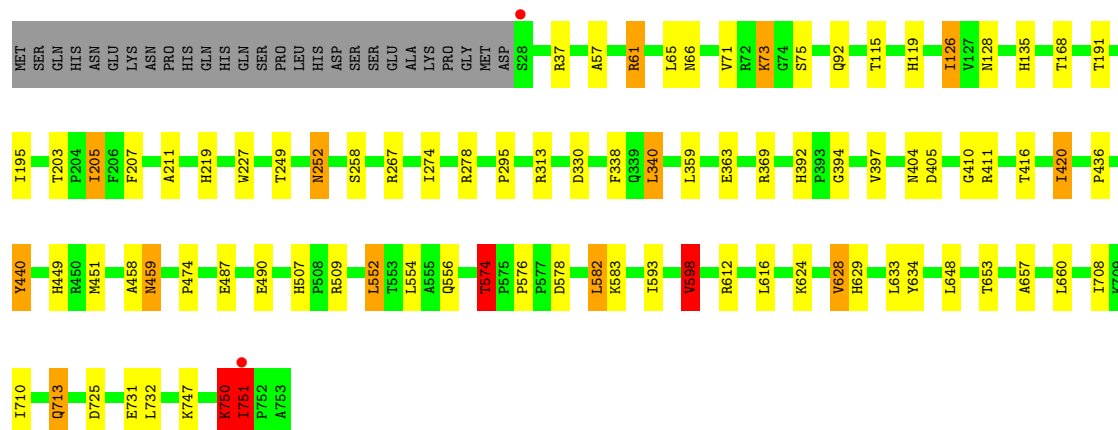
• Molecule 1: Catalase HP11





● Molecule 1: Catalase HP11

Chain D: 85% 9% . . .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	93.62Å 132.96Å 122.67Å 90.00° 109.39° 90.00°	Depositor
Resolution (Å)	35.46 – 1.89 35.46 – 1.90	Depositor EDS
% Data completeness (in resolution range)	84.3 (35.46-1.89) 84.3 (35.46-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.17 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.144 , 0.193 0.144 , 0.193	Depositor DCC
R_{free} test set	9439 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	15.3	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.0	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	25986	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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 [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	1.28	8/5903 (0.1%)	1.11	7/8026 (0.1%)
1	B	1.24	5/5903 (0.1%)	1.12	14/8026 (0.2%)
1	C	1.24	8/5903 (0.1%)	1.10	6/8026 (0.1%)
1	D	1.28	11/5903 (0.2%)	1.12	8/8026 (0.1%)
All	All	1.26	32/23612 (0.1%)	1.11	35/32104 (0.1%)

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	A	0	1
1	C	0	2
All	All	0	3

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	474	PRO	CA-C	8.46	1.56	1.51
1	B	474	PRO	CA-C	7.94	1.56	1.51
1	D	474	PRO	CA-C	7.52	1.56	1.51
1	A	29	LEU	N-CA	-7.27	1.36	1.45
1	B	593	ILE	CA-CB	7.19	1.61	1.53
1	B	647	VAL	CA-CB	6.51	1.61	1.54
1	D	458	ALA	C-O	6.44	1.31	1.23
1	A	412	LEU	N-CA	6.35	1.54	1.46
1	D	203	THR	N-CA	6.30	1.50	1.45
1	A	178	THR	N-CA	5.93	1.54	1.46
1	B	341	ILE	N-CA	5.84	1.50	1.45
1	A	709	LYS	CA-C	-5.80	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	440	TYR	C-O	5.67	1.30	1.23
1	C	86	VAL	N-CA	5.65	1.53	1.46
1	C	179	VAL	CA-CB	5.62	1.60	1.54
1	C	751	ILE	CA-CB	5.54	1.60	1.54
1	D	57	ALA	CA-CB	5.42	1.59	1.53
1	C	214	PHE	C-O	-5.42	1.19	1.24
1	A	37	ARG	CA-C	5.42	1.58	1.52
1	A	353	LEU	N-CA	5.34	1.53	1.46
1	A	307	ALA	CA-CB	5.33	1.61	1.53
1	D	593	ILE	CA-CB	5.27	1.59	1.53
1	D	37	ARG	CA-C	5.22	1.58	1.52
1	C	638	GLY	N-CA	5.18	1.50	1.45
1	D	397	VAL	CA-CB	5.17	1.62	1.54
1	C	200	GLY	N-CA	5.14	1.49	1.44
1	D	628	VAL	CA-CB	5.11	1.60	1.54
1	B	710	ILE	CA-CB	5.10	1.60	1.54
1	A	235	ALA	CA-CB	5.08	1.59	1.52
1	C	420	ILE	CG1-CD1	-5.05	1.32	1.51
1	D	295	PRO	N-CA	5.03	1.53	1.47
1	D	71	VAL	CA-CB	5.01	1.60	1.54

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	61	ARG	NE-CZ-NH2	8.55	126.90	119.20
1	A	120	GLU	N-CA-C	7.37	119.31	111.28
1	B	365	VAL	CA-C-N	-6.48	113.61	120.03
1	B	365	VAL	C-N-CA	-6.48	113.61	120.03
1	B	209	GLN	N-CA-C	6.47	119.64	111.69
1	A	726	GLY	N-CA-C	-6.44	104.70	112.49
1	D	657	ALA	CA-C-N	6.30	126.66	119.92
1	D	657	ALA	C-N-CA	6.30	126.66	119.92
1	A	710	ILE	CA-C-N	6.23	133.43	121.54
1	A	710	ILE	C-N-CA	6.23	133.43	121.54
1	D	61	ARG	NE-CZ-NH1	-6.09	115.41	121.50
1	B	745	ILE	CA-C-N	-6.04	113.40	119.56
1	B	745	ILE	C-N-CA	-6.04	113.40	119.56
1	C	725	ASP	N-CA-C	6.00	123.58	110.80
1	D	574	THR	CB-CA-C	5.70	118.36	109.42
1	C	159	ILE	CB-CG1-CD1	-5.61	102.02	113.80
1	C	214	PHE	CA-C-N	-5.59	113.27	119.19
1	C	214	PHE	C-N-CA	-5.59	113.27	119.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	593	ILE	CA-C-N	5.57	126.81	119.84
1	B	593	ILE	C-N-CA	5.57	126.81	119.84
1	D	258	SER	N-CA-C	-5.38	103.59	110.53
1	C	126	ILE	CB-CG1-CD1	-5.37	102.53	113.80
1	B	120	GLU	N-CA-C	5.29	117.79	111.71
1	D	598	VAL	N-CA-CB	-5.27	100.59	111.76
1	B	214	PHE	CA-C-N	-5.21	113.54	119.28
1	B	214	PHE	C-N-CA	-5.21	113.54	119.28
1	A	711	ALA	CA-C-N	5.16	131.56	122.42
1	A	711	ALA	C-N-CA	5.16	131.56	122.42
1	A	168	THR	N-CA-C	-5.10	101.95	109.86
1	C	304	TRP	N-CA-C	5.07	117.19	111.11
1	B	550	ILE	N-CA-C	-5.04	106.74	111.48
1	D	168	THR	N-CA-C	-5.04	102.05	109.86
1	B	350	ASP	CB-CA-C	5.04	119.07	110.56
1	B	657	ALA	CA-C-N	5.03	125.30	119.92
1	B	657	ALA	C-N-CA	5.03	125.30	119.92

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	708	ILE	Peptide
1	C	724	ALA	Peptide
1	C	725	ASP	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5743	0	5576	55	0
1	B	5744	0	5580	58	0
1	C	5743	0	5576	59	0
1	D	5744	0	5580	54	0
2	A	43	0	30	1	0
2	B	43	0	30	0	0
2	C	43	0	30	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	43	0	30	2	0
3	A	772	0	0	12	0
3	B	639	0	0	11	0
3	C	680	0	0	13	0
3	D	749	0	0	8	1
All	All	25986	0	22432	200	1

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 4.

All (200) 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:B:449[B]:HIS:CE1	3:B:3275:HOH:O	1.90	1.22
1:A:449[A]:HIS:CE1	3:A:3276:HOH:O	1.96	1.16
1:B:449[B]:HIS:NE2	3:B:3275:HOH:O	1.77	1.14
3:B:2705:HOH:O	1:D:73:LYS:HD3	1.47	1.13
1:B:416:THR:HG21	3:B:3233:HOH:O	1.47	1.12
1:B:451:MET:HE2	1:D:451:MET:HE2	1.32	1.10
1:A:449[A]:HIS:CD2	1:C:449[A]:HIS:CD2	2.40	1.08
1:B:521:ARG:HD3	3:B:3047:HOH:O	1.53	1.08
1:A:451:MET:HE2	1:C:451:MET:HE2	1.42	0.98
1:B:612:ARG:HH11	1:B:612:ARG:HB2	1.27	0.95
1:A:639:GLU:HG3	3:A:2414:HOH:O	1.75	0.85
3:B:2705:HOH:O	1:D:73:LYS:CD	2.13	0.84
1:B:583:LYS:H	1:B:583:LYS:NZ	1.76	0.84
1:B:449[B]:HIS:CD2	1:D:449[B]:HIS:CD2	2.67	0.83
1:A:449[A]:HIS:NE2	3:A:3276:HOH:O	2.04	0.83
1:A:449[A]:HIS:CD2	1:C:449[A]:HIS:HD2	1.94	0.81
1:B:392:HIS:CD2	1:B:394:GLY:H	1.99	0.81
1:B:583:LYS:H	1:B:583:LYS:HZ2	1.30	0.78
1:D:267:ARG:HG3	3:D:1920:HOH:O	1.82	0.78
1:A:479:ARG:NH2	3:A:2607:HOH:O	2.16	0.76
1:D:750:LYS:HD2	1:D:751:ILE:H	1.51	0.74
1:B:708:ILE:HG13	1:B:710:ILE:HG12	1.68	0.74
1:C:392:HIS:CD2	1:C:394:GLY:H	2.06	0.74
1:A:392:HIS:CD2	1:A:394:GLY:H	2.06	0.72
1:B:533:LYS:HE3	3:C:2623:HOH:O	1.89	0.71
1:C:612:ARG:HG3	1:C:612:ARG:HH11	1.56	0.70
1:A:28:SER:O	1:A:28:SER:OG	2.10	0.70
1:B:119:HIS:CE1	1:D:420:ILE:HG21	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:750:LYS:HD2	1:D:751:ILE:N	2.07	0.69
1:A:392:HIS:HD2	1:A:394:GLY:H	1.42	0.67
1:B:629:HIS:HD2	3:B:1046:HOH:O	1.77	0.67
1:C:283:GLU:OE1	3:C:2292:HOH:O	2.13	0.67
1:B:392:HIS:HD2	1:B:394:GLY:H	1.43	0.66
1:D:731:GLU:OE2	3:D:3028:HOH:O	2.14	0.65
1:C:629:HIS:HD2	3:C:1129:HOH:O	1.80	0.65
1:B:583:LYS:O	1:B:584:LYS:HB3	1.96	0.65
1:A:449[B]:HIS:CG	1:C:449[B]:HIS:CG	2.67	0.65
1:A:751:ILE:O	1:A:751:ILE:HD12	1.97	0.64
1:C:137:TYR:HB2	1:C:159:ILE:CD1	2.28	0.64
1:C:727:SER:HA	3:C:2714:HOH:O	1.97	0.64
1:B:448:MET:HG3	1:B:449[A]:HIS:CD2	2.32	0.64
1:B:552:LEU:HD21	1:B:571:LEU:HD12	1.80	0.64
3:B:2705:HOH:O	1:D:73:LYS:CE	2.44	0.63
1:A:29:LEU:CB	3:C:2405:HOH:O	2.46	0.63
1:C:392:HIS:HD2	1:C:394:GLY:H	1.44	0.63
1:C:583:LYS:O	1:C:584:LYS:HB3	1.98	0.62
1:D:392:HIS:CD2	1:D:394:GLY:H	2.17	0.62
1:D:629:HIS:HD2	3:D:1554:HOH:O	1.82	0.62
1:B:281:ASN:OD1	1:B:283:GLU:HG3	2.00	0.62
1:C:330:ASP:OD1	1:C:629:HIS:HE1	1.83	0.62
1:B:682:ASN:HB3	1:B:707:THR:HG21	1.82	0.61
1:C:704:PHE:O	1:C:707:THR:HG22	1.99	0.61
1:A:29:LEU:HB2	3:C:2405:HOH:O	2.01	0.61
1:C:612:ARG:HH11	1:C:612:ARG:CG	2.14	0.60
1:C:708:ILE:HG13	1:C:710:ILE:HG12	1.82	0.60
1:B:449[A]:HIS:CG	1:D:449[A]:HIS:CG	2.55	0.60
1:D:330:ASP:OD1	1:D:629:HIS:HE1	1.85	0.59
1:D:747:LYS:O	1:D:750:LYS:HE2	2.03	0.58
1:D:392:HIS:HE1	1:D:416:THR:OG1	1.86	0.58
1:A:449[A]:HIS:HD2	1:C:449[A]:HIS:CD2	2.18	0.58
1:A:36:HIS:CD2	1:A:36:HIS:H	2.22	0.58
1:A:278:ARG:HH12	1:A:487:GLU:CD	2.13	0.57
1:B:724:ALA:O	1:B:725:ASP:O	2.21	0.57
1:D:750:LYS:O	1:D:751:ILE:O	2.23	0.56
1:A:36:HIS:HE1	3:A:1872:HOH:O	1.89	0.56
1:A:392:HIS:HE1	1:A:416:THR:OG1	1.90	0.55
1:B:612:ARG:HB2	1:B:612:ARG:NH1	2.10	0.55
1:B:330:ASP:OD1	1:B:629:HIS:HE1	1.90	0.55
1:D:61:ARG:HH11	1:D:66:ASN:HA	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:LEU:HD22	3:C:2405:HOH:O	2.08	0.54
1:D:359:LEU:H	1:D:507:HIS:HD2	1.54	0.54
1:B:38:PRO:HG2	1:B:51:ALA:HB2	1.89	0.53
1:C:612:ARG:HG3	1:C:612:ARG:NH1	2.21	0.53
1:C:578:ASP:HB2	1:C:582:LEU:O	2.08	0.53
1:A:710:ILE:HG12	1:A:715:GLU:OE1	2.08	0.53
1:D:392:HIS:HD2	1:D:394:GLY:H	1.55	0.53
1:A:603:VAL:HG11	1:A:666:ILE:HD12	1.91	0.53
1:A:120:GLU:HB2	1:D:126:ILE:CD1	2.39	0.53
1:A:478:LYS:HG2	3:A:1955:HOH:O	2.08	0.53
1:A:119:HIS:CE1	1:C:420:ILE:HG21	2.45	0.52
1:C:359:LEU:H	1:C:507:HIS:HD2	1.57	0.52
1:C:416:THR:CG2	3:C:3238:HOH:O	2.58	0.52
1:C:392:HIS:HE1	1:C:416:THR:OG1	1.92	0.52
1:C:724:ALA:O	1:C:725:ASP:HB2	2.10	0.52
1:B:533:LYS:CE	3:C:2623:HOH:O	2.53	0.51
1:B:120:GLU:HB2	1:C:126:ILE:HD11	1.93	0.51
1:C:535:VAL:O	1:C:537:PRO:HD3	2.10	0.51
1:A:607:LEU:HD11	1:A:632:LEU:HB3	1.92	0.51
1:A:710:ILE:CG1	1:A:715:GLU:OE1	2.59	0.51
1:D:509:ARG:HD2	1:D:576:PRO:HD2	1.92	0.51
1:C:38:PRO:HG2	1:C:51:ALA:HB2	1.92	0.51
1:D:708:ILE:HG13	1:D:710:ILE:HG12	1.92	0.50
1:B:359:LEU:H	1:B:507:HIS:HD2	1.58	0.50
1:A:274:ILE:HD12	2:A:760:HEM:HMB1	1.92	0.50
1:A:219:HIS:HB3	1:B:459:ASN:ND2	2.27	0.50
1:C:610:GLU:O	1:C:610:GLU:HG3	2.12	0.50
1:D:115:THR:O	1:D:119:HIS:HD2	1.94	0.50
1:B:612:ARG:NH1	1:B:669:CYS:SG	2.85	0.49
1:A:459:ASN:ND2	1:B:219:HIS:HB3	2.27	0.49
1:C:219:HIS:HB3	1:D:459:ASN:ND2	2.27	0.49
1:D:205:ILE:H	1:D:205:ILE:HD13	1.77	0.49
1:A:449[B]:HIS:CD2	3:A:1077:HOH:O	2.65	0.49
1:D:598:VAL:HG13	1:D:628:VAL:CG2	2.42	0.49
1:B:420:ILE:HG21	1:D:119:HIS:CE1	2.48	0.49
1:B:533:LYS:HE2	3:B:3100:HOH:O	2.11	0.48
1:A:440:TYR:OH	3:A:938:HOH:O	2.18	0.48
1:D:363:GLU:HB2	1:D:582:LEU:HD21	1.95	0.48
1:A:459:ASN:C	1:A:459:ASN:HD22	2.20	0.48
1:B:392:HIS:HE1	1:B:416:THR:OG1	1.96	0.48
1:C:115:THR:O	1:C:119:HIS:HD2	1.97	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:ASP:OD1	1:C:461:GLU:OE1	2.32	0.48
1:A:634:TYR:O	1:A:653:THR:HA	2.13	0.48
1:A:207:PHE:CD1	1:A:217:PHE:CZ	3.02	0.48
1:C:137:TYR:HB2	1:C:159:ILE:HD11	1.97	0.47
1:D:750:LYS:O	1:D:751:ILE:C	2.57	0.47
1:B:126:ILE:CD1	1:C:120:GLU:HB2	2.45	0.47
1:D:65:LEU:HD21	1:D:135:HIS:CG	2.50	0.47
1:B:28:SER:HB2	3:D:2425:HOH:O	2.15	0.47
1:B:309:LYS:HD2	1:B:660:LEU:HD11	1.97	0.47
1:C:610:GLU:OE1	1:C:643:ASP:HA	2.15	0.47
1:D:211:ALA:CB	1:D:410:GLY:HA3	2.45	0.46
1:C:278:ARG:HH12	1:C:487:GLU:CD	2.22	0.46
1:A:603:VAL:HG11	1:A:666:ILE:CD1	2.46	0.46
1:A:716:GLU:HG2	3:A:1656:HOH:O	2.15	0.46
1:D:278:ARG:HH12	1:D:487:GLU:CD	2.24	0.46
1:C:207:PHE:O	1:C:249:THR:HA	2.15	0.46
1:A:65:LEU:HD21	1:A:135:HIS:CG	2.50	0.46
1:B:637:MET:HB2	1:C:562:LEU:HA	1.98	0.46
1:B:725:ASP:OD2	1:B:725:ASP:C	2.58	0.46
1:D:713:GLN:O	1:D:713:GLN:HG2	2.15	0.46
1:A:449[B]:HIS:CD2	1:C:449[B]:HIS:CG	3.01	0.46
1:C:459:ASN:ND2	1:D:219:HIS:HB3	2.31	0.46
1:D:459:ASN:HD22	1:D:459:ASN:C	2.24	0.45
1:D:556:GLN:NE2	3:D:2773:HOH:O	2.48	0.45
1:A:158:LYS:HD3	3:A:2278:HOH:O	2.16	0.45
1:A:213:LYS:HD3	1:D:92:GLN:HA	1.99	0.45
1:A:395:HIS:HE1	3:A:3270:HOH:O	1.99	0.45
1:A:461:GLU:HA	1:A:462:PRO:C	2.41	0.45
1:B:478:LYS:HD2	3:B:2857:HOH:O	2.17	0.45
1:B:583:LYS:H	1:B:583:LYS:HZ3	1.60	0.45
1:D:274:ILE:HD12	2:D:760:HEM:HMB1	1.99	0.45
1:C:51:ALA:HB1	1:C:52:PRO:HD2	1.98	0.45
1:B:335:GLU:OE1	1:B:369:ARG:HD2	2.17	0.44
1:C:488:ARG:HD3	3:C:2379:HOH:O	2.16	0.44
1:C:739:HIS:CD2	1:C:740:ARG:HG2	2.52	0.44
1:B:689:TYR:OH	1:B:715:GLU:OE1	2.35	0.44
1:B:37:ARG:HD3	3:B:2886:HOH:O	2.17	0.44
1:D:725:ASP:OD1	1:D:725:ASP:C	2.61	0.44
1:D:207:PHE:O	1:D:249:THR:HA	2.16	0.44
1:A:207:PHE:O	1:A:249:THR:HA	2.18	0.44
1:A:211:ALA:CB	1:A:410:GLY:HA3	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:338:PHE:HB3	1:D:340:LEU:HD13	2.00	0.44
1:C:202:ASN:HA	1:C:270:GLU:O	2.18	0.43
1:A:31:PRO:HD2	1:A:36:HIS:HB3	1.99	0.43
1:C:416:THR:HG21	3:C:3238:HOH:O	2.18	0.43
1:D:404:ASN:O	1:D:405:ASP:C	2.61	0.43
1:D:552:LEU:HD22	1:D:552:LEU:HA	1.85	0.43
1:A:459:ASN:ND2	1:A:459:ASN:C	2.77	0.43
1:C:725:ASP:HB3	1:C:726:GLY:H	1.50	0.43
1:D:634:TYR:O	1:D:653:THR:HA	2.18	0.43
1:C:128:ASN:HA	1:C:168:THR:O	2.18	0.43
1:B:562:LEU:HA	1:C:637:MET:HB2	1.99	0.43
1:D:411:ARG:HG2	2:D:760:HEM:C2C	2.54	0.43
1:B:634:TYR:O	1:B:653:THR:HA	2.19	0.43
1:B:251:HIS:CE1	1:B:507:HIS:HB3	2.54	0.43
1:C:713:GLN:H	1:C:713:GLN:HG3	1.31	0.43
1:C:716:GLU:OE2	1:C:716:GLU:HA	2.18	0.42
1:D:359:LEU:H	1:D:507:HIS:CD2	2.35	0.42
1:D:578:ASP:HB2	3:D:2129:HOH:O	2.18	0.42
1:C:725:ASP:HA	1:C:728:PHE:HB3	2.00	0.42
1:B:359:LEU:H	1:B:507:HIS:CD2	2.35	0.42
1:D:195:ILE:HD11	1:D:436:PRO:HA	2.02	0.42
1:B:52:PRO:HG3	3:D:1451:HOH:O	2.19	0.42
1:B:278:ARG:HH22	1:B:487:GLU:CD	2.28	0.42
1:A:449[A]:HIS:CG	1:C:449[A]:HIS:CD2	3.03	0.42
1:A:443:PHE:CZ	1:A:470:PRO:HD2	2.55	0.42
1:A:549:HIS:O	1:A:576:PRO:HD3	2.20	0.42
1:B:512:TRP:CH2	1:B:520:GLN:HB3	2.54	0.41
1:B:745:ILE:O	1:B:748:ILE:HG12	2.20	0.41
1:C:211:ALA:CB	1:C:410:GLY:HA3	2.50	0.41
1:A:404:ASN:O	1:A:405:ASP:C	2.63	0.41
1:B:213:LYS:HB3	1:B:213:LYS:HE3	1.78	0.41
1:A:369:ARG:HG2	3:A:1078:HOH:O	2.19	0.41
1:B:115:THR:O	1:B:119:HIS:HD2	2.03	0.41
1:B:207:PHE:O	1:B:249:THR:HA	2.20	0.41
1:C:507:HIS:HE1	3:C:926:HOH:O	2.03	0.41
1:C:461:GLU:HA	1:C:462:PRO:C	2.44	0.41
1:D:574:THR:HG22	3:D:1614:HOH:O	2.21	0.41
1:C:440:TYR:OH	3:C:1028:HOH:O	1.93	0.41
1:C:634:TYR:CG	1:C:635:SER:N	2.89	0.41
1:A:319:ARG:HD3	1:D:227:TRP:O	2.19	0.41
1:A:604:ALA:HB2	1:A:662:VAL:HG11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:ILE:HD12	1:C:121:ARG:CZ	2.51	0.41
1:B:140:PRO:HG2	1:B:155:ASP:O	2.21	0.41
1:C:404:ASN:O	1:C:405:ASP:C	2.64	0.41
1:D:313:ARG:HG3	1:D:660:LEU:HD12	2.02	0.41
1:D:578:ASP:HB3	1:D:582:LEU:O	2.20	0.40
1:B:404:ASN:O	1:B:405:ASP:C	2.65	0.40
1:C:251:HIS:CE1	1:C:507:HIS:HB3	2.56	0.40
1:D:207:PHE:CD2	1:D:252:ASN:HB3	2.56	0.40

All (1) 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)
3:D:2178:HOH:O	3:D:2976:HOH:O[1_655]	2.09	0.11

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	725/753 (96%)	706 (97%)	17 (2%)	2 (0%)	36	29
1	B	725/753 (96%)	703 (97%)	19 (3%)	3 (0%)	30	22
1	C	725/753 (96%)	704 (97%)	19 (3%)	2 (0%)	36	29
1	D	725/753 (96%)	703 (97%)	19 (3%)	3 (0%)	30	22
All	All	2900/3012 (96%)	2816 (97%)	74 (3%)	10 (0%)	36	29

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	711	ALA
1	B	725	ASP

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Mol	Chain	Res	Type
1	C	725	ASP
1	D	750	LYS
1	D	751	ILE
1	B	75	SER
1	C	75	SER
1	D	75	SER
1	A	75	SER
1	B	584	LYS

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	612/636 (96%)	591 (97%)	21 (3%)	32	25
1	B	612/636 (96%)	583 (95%)	29 (5%)	23	15
1	C	612/636 (96%)	578 (94%)	34 (6%)	19	11
1	D	612/636 (96%)	585 (96%)	27 (4%)	25	17
All	All	2448/2544 (96%)	2337 (96%)	111 (4%)	24	17

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	SER
1	A	29	LEU
1	A	37	ARG
1	A	73	LYS
1	A	126	ILE
1	A	191	THR
1	A	198	LEU
1	A	205	ILE
1	A	252	ASN
1	A	369	ARG
1	A	420	ILE
1	A	440	TYR
1	A	459	ASN

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Mol	Chain	Res	Type
1	A	478	LYS
1	A	552	LEU
1	A	709	LYS
1	A	712	ASP
1	A	713	GLN
1	A	716	GLU
1	A	727	SER
1	A	732	LEU
1	B	126	ILE
1	B	155	ASP
1	B	191	THR
1	B	195	ILE
1	B	205	ILE
1	B	213	LYS
1	B	252	ASN
1	B	285	LYS
1	B	321	GLU
1	B	370	VAL
1	B	416	THR
1	B	420	ILE
1	B	432	PRO
1	B	440	TYR
1	B	459	ASN
1	B	552	LEU
1	B	562	LEU
1	B	565	GLU
1	B	568	ASP
1	B	571	LEU
1	B	583	LYS
1	B	595	ASP
1	B	612	ARG
1	B	616	LEU
1	B	633	LEU
1	B	703	LYS
1	B	713	GLN
1	B	750	LYS
1	B	751	ILE
1	C	37	ARG
1	C	159	ILE
1	C	191	THR
1	C	198	LEU
1	C	205	ILE

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Mol	Chain	Res	Type
1	C	213	LYS
1	C	252	ASN
1	C	377	ARG
1	C	440	TYR
1	C	459	ASN
1	C	478	LYS
1	C	488	ARG
1	C	531	LEU
1	C	552	LEU
1	C	568	ASP
1	C	571	LEU
1	C	574	THR
1	C	584	LYS
1	C	606	LEU
1	C	610	GLU
1	C	612	ARG
1	C	616	LEU
1	C	617	LEU
1	C	633	LEU
1	C	636	ARG
1	C	648	LEU
1	C	660	LEU
1	C	685	LEU
1	C	709	LYS
1	C	713	GLN
1	C	732	LEU
1	C	733	LEU
1	C	750	LYS
1	C	751	ILE
1	D	73	LYS
1	D	126	ILE
1	D	128	ASN
1	D	191	THR
1	D	205	ILE
1	D	252	ASN
1	D	340	LEU
1	D	369	ARG
1	D	420	ILE
1	D	440	TYR
1	D	459	ASN
1	D	490	GLU
1	D	552	LEU

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Mol	Chain	Res	Type
1	D	554	LEU
1	D	574	THR
1	D	582	LEU
1	D	583	LYS
1	D	598	VAL
1	D	612	ARG
1	D	616	LEU
1	D	624	LYS
1	D	633	LEU
1	D	648	LEU
1	D	713	GLN
1	D	732	LEU
1	D	750	LYS
1	D	751	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	252	ASN
1	A	392	HIS
1	A	459	ASN
1	B	252	ASN
1	B	392	HIS
1	B	459	ASN
1	B	507	HIS
1	B	629	HIS
1	C	252	ASN
1	C	392	HIS
1	C	459	ASN
1	C	507	HIS
1	C	546	GLN
1	C	629	HIS
1	C	671	ASN
1	D	48	GLN
1	D	128	ASN
1	D	252	ASN
1	D	392	HIS
1	D	459	ASN
1	D	507	HIS
1	D	546	GLN
1	D	556	GLN

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Mol	Chain	Res	Type
1	D	629	HIS
1	D	671	ASN

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](#)

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	$\# Z > 2$
2	HEM	A	760	1	50,50,50	1.87	13 (26%)	67,82,82	1.65	14 (20%)
2	HEM	B	760	1	50,50,50	2.00	13 (26%)	67,82,82	1.71	14 (20%)
2	HEM	D	760	1	50,50,50	2.09	11 (22%)	67,82,82	1.55	10 (14%)
2	HEM	C	760	1	50,50,50	1.88	12 (24%)	67,82,82	1.85	13 (19%)

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	A	760	1	-	2/14/54/54	-
2	HEM	B	760	1	-	3/14/54/54	-
2	HEM	D	760	1	-	3/14/54/54	-
2	HEM	C	760	1	-	4/14/54/54	-

All (49) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	760	HEM	FE-ND	6.91	2.16	1.94
2	D	760	HEM	C3D-C2D	6.21	1.50	1.36
2	B	760	HEM	FE-NB	5.73	2.12	1.94
2	C	760	HEM	C3D-C2D	5.67	1.49	1.36
2	B	760	HEM	C3D-C2D	5.56	1.48	1.36
2	A	760	HEM	C3D-C2D	5.41	1.48	1.36
2	A	760	HEM	FE-ND	5.39	2.11	1.94
2	C	760	HEM	FE-NC	5.12	2.12	1.95
2	D	760	HEM	FE-NB	4.62	2.09	1.94
2	A	760	HEM	FE-NB	4.47	2.08	1.94
2	B	760	HEM	FE-NC	4.34	2.09	1.95
2	D	760	HEM	FE-NA	4.26	2.09	1.95
2	C	760	HEM	FE-NB	4.17	2.07	1.94
2	B	760	HEM	FE-ND	4.06	2.07	1.94
2	C	760	HEM	FE-ND	3.82	2.06	1.94
2	A	760	HEM	CMC-C2C	3.77	1.58	1.50
2	C	760	HEM	FE-NA	3.59	2.07	1.95
2	A	760	HEM	FE-NC	3.52	2.06	1.95
2	D	760	HEM	CAC-C3C	3.45	1.56	1.47
2	B	760	HEM	CMC-C2C	3.32	1.57	1.50
2	D	760	HEM	CMD-C2D	3.24	1.57	1.50
2	B	760	HEM	CMA-C3A	3.15	1.57	1.50
2	D	760	HEM	CMB-C2B	3.06	1.57	1.50
2	B	760	HEM	FE-NA	3.05	2.05	1.95
2	C	760	HEM	CAB-C3B	3.01	1.55	1.47
2	B	760	HEM	CAC-C3C	3.00	1.55	1.47
2	B	760	HEM	C1B-NB	-2.98	1.35	1.40
2	C	760	HEM	CMA-C3A	2.97	1.56	1.50
2	D	760	HEM	CMC-C2C	2.75	1.56	1.50
2	D	760	HEM	CAB-C3B	2.68	1.54	1.47
2	D	760	HEM	FE-NC	2.66	2.03	1.95
2	C	760	HEM	CAC-C3C	2.51	1.54	1.47
2	B	760	HEM	CAB-C3B	2.42	1.53	1.47
2	A	760	HEM	CAC-C3C	2.37	1.53	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	760	HEM	CMD-C2D	2.35	1.55	1.50
2	A	760	HEM	CAA-C2A	2.32	1.57	1.51
2	A	760	HEM	CMA-C3A	2.31	1.55	1.50
2	A	760	HEM	FE-NA	2.27	2.02	1.95
2	A	760	HEM	CMB-C2B	2.22	1.55	1.50
2	A	760	HEM	O1A-CGA	2.22	1.29	1.22
2	D	760	HEM	CMA-C3A	2.20	1.55	1.50
2	B	760	HEM	CMD-C2D	2.17	1.55	1.50
2	C	760	HEM	CBD-CGD	2.15	1.55	1.50
2	B	760	HEM	C4C-NC	-2.15	1.35	1.39
2	C	760	HEM	O1D-CGD	2.12	1.29	1.22
2	A	760	HEM	CMD-C2D	2.09	1.55	1.50
2	A	760	HEM	CAB-C3B	2.08	1.52	1.47
2	B	760	HEM	C2A-C3A	-2.04	1.33	1.38
2	C	760	HEM	O2A-CGA	-2.01	1.24	1.30

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	760	HEM	C4D-ND-C1D	5.72	111.98	105.21
2	B	760	HEM	C4D-ND-C1D	4.98	111.11	105.21
2	C	760	HEM	C3B-C4B-NB	-4.76	106.05	109.47
2	A	760	HEM	C4D-ND-C1D	4.69	110.77	105.21
2	C	760	HEM	CHC-C4B-NB	4.64	129.42	124.42
2	B	760	HEM	C1B-NB-C4B	4.39	110.41	105.21
2	A	760	HEM	C1B-NB-C4B	4.13	110.10	105.21
2	C	760	HEM	C1B-NB-C4B	4.11	110.08	105.21
2	A	760	HEM	C3B-C4B-NB	-4.11	106.52	109.47
2	A	760	HEM	CBD-CAD-C3D	-4.10	101.20	112.53
2	C	760	HEM	CAA-CBA-CGA	-4.01	103.02	113.67
2	C	760	HEM	C2B-C1B-NB	-3.99	105.25	109.84
2	C	760	HEM	C4D-ND-C1D	3.98	109.92	105.21
2	B	760	HEM	CHA-C4D-ND	3.90	129.19	124.37
2	C	760	HEM	CHA-C4D-ND	3.67	128.91	124.37
2	D	760	HEM	CBD-CAD-C3D	-3.66	102.42	112.53
2	B	760	HEM	C3B-C2B-C1B	3.66	109.16	106.41
2	D	760	HEM	CHD-C4C-NC	3.60	128.37	124.45
2	B	760	HEM	C3B-C4B-NB	-3.56	106.91	109.47
2	B	760	HEM	C2B-C1B-NB	-3.39	105.94	109.84
2	A	760	HEM	CAA-CBA-CGA	-3.31	104.89	113.67
2	B	760	HEM	C3D-C4D-ND	-3.10	106.77	110.17
2	D	760	HEM	C3B-C4B-NB	-3.06	107.27	109.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	760	HEM	C3B-C2B-C1B	3.06	108.71	106.41
2	C	760	HEM	CBD-CAD-C3D	-2.96	104.34	112.53
2	D	760	HEM	C2A-C1A-NA	-2.87	106.97	110.15
2	D	760	HEM	C3D-C4D-ND	-2.87	107.03	110.17
2	B	760	HEM	C2A-C1A-NA	-2.84	107.00	110.15
2	B	760	HEM	CAA-CBA-CGA	-2.68	106.56	113.67
2	A	760	HEM	C2B-C1B-NB	-2.67	106.77	109.84
2	D	760	HEM	CAA-CBA-CGA	-2.63	106.70	113.67
2	B	760	HEM	CHD-C4C-NC	2.60	127.29	124.45
2	A	760	HEM	CHD-C4C-NC	2.55	127.23	124.45
2	A	760	HEM	CBC-CAC-C3C	-2.54	114.82	127.53
2	A	760	HEM	C4C-NC-C1C	2.45	109.81	105.82
2	C	760	HEM	CHD-C4C-NC	2.41	127.08	124.45
2	A	760	HEM	C3B-C2B-C1B	2.40	108.21	106.41
2	D	760	HEM	C1B-NB-C4B	2.39	108.04	105.21
2	A	760	HEM	C4C-C3C-C2C	2.36	108.86	106.81
2	C	760	HEM	C2A-C1A-NA	-2.32	107.57	110.15
2	B	760	HEM	CBD-CAD-C3D	-2.29	106.19	112.53
2	B	760	HEM	C4C-NC-C1C	2.29	109.55	105.82
2	A	760	HEM	CAD-C3D-C4D	2.26	128.64	124.70
2	A	760	HEM	C2D-C1D-ND	-2.23	107.32	109.90
2	B	760	HEM	O2A-CGA-O1A	-2.22	117.62	123.33
2	A	760	HEM	CHA-C4D-ND	2.18	127.06	124.37
2	B	760	HEM	CHC-C4B-NB	2.14	126.72	124.42
2	D	760	HEM	C4A-NA-C1A	2.13	109.29	105.82
2	D	760	HEM	CHA-C4D-ND	2.12	126.99	124.37
2	C	760	HEM	C1D-C2D-C3D	-2.10	104.77	106.98
2	C	760	HEM	CBA-CAA-C2A	-2.00	107.00	112.53

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	760	HEM	CAA-CBA-CGA-O1A
2	C	760	HEM	CAA-CBA-CGA-O1A
2	A	760	HEM	CAA-CBA-CGA-O2A
2	D	760	HEM	CAA-CBA-CGA-O2A
2	D	760	HEM	CAA-CBA-CGA-O1A
2	C	760	HEM	CAA-CBA-CGA-O2A
2	B	760	HEM	CAA-CBA-CGA-O2A
2	A	760	HEM	CAA-CBA-CGA-O1A
2	C	760	HEM	CAD-CBD-CGD-O2D

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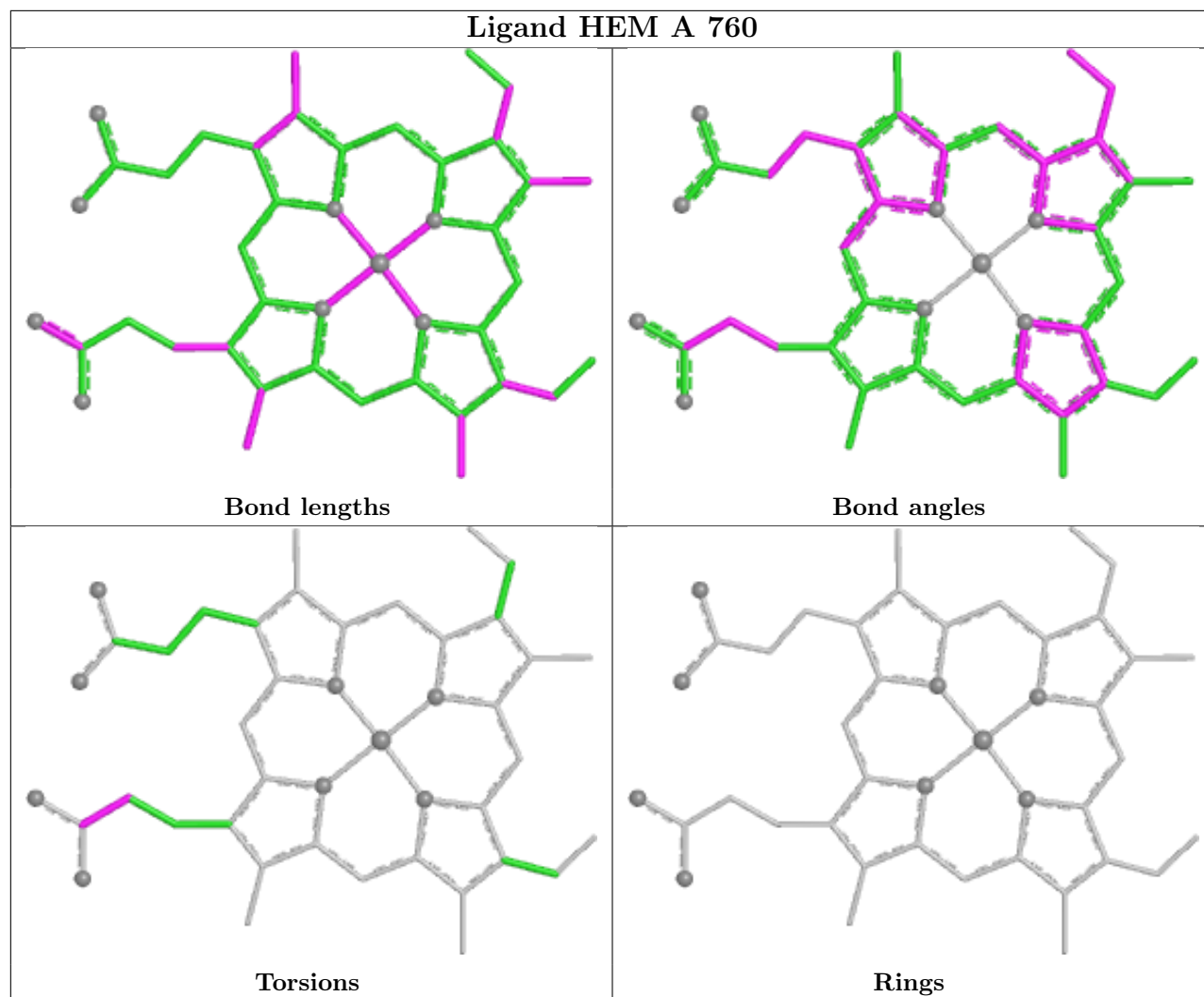
Mol	Chain	Res	Type	Atoms
2	C	760	HEM	CAD-CBD-CGD-O1D
2	D	760	HEM	CAD-CBD-CGD-O2D
2	B	760	HEM	CAD-CBD-CGD-O1D

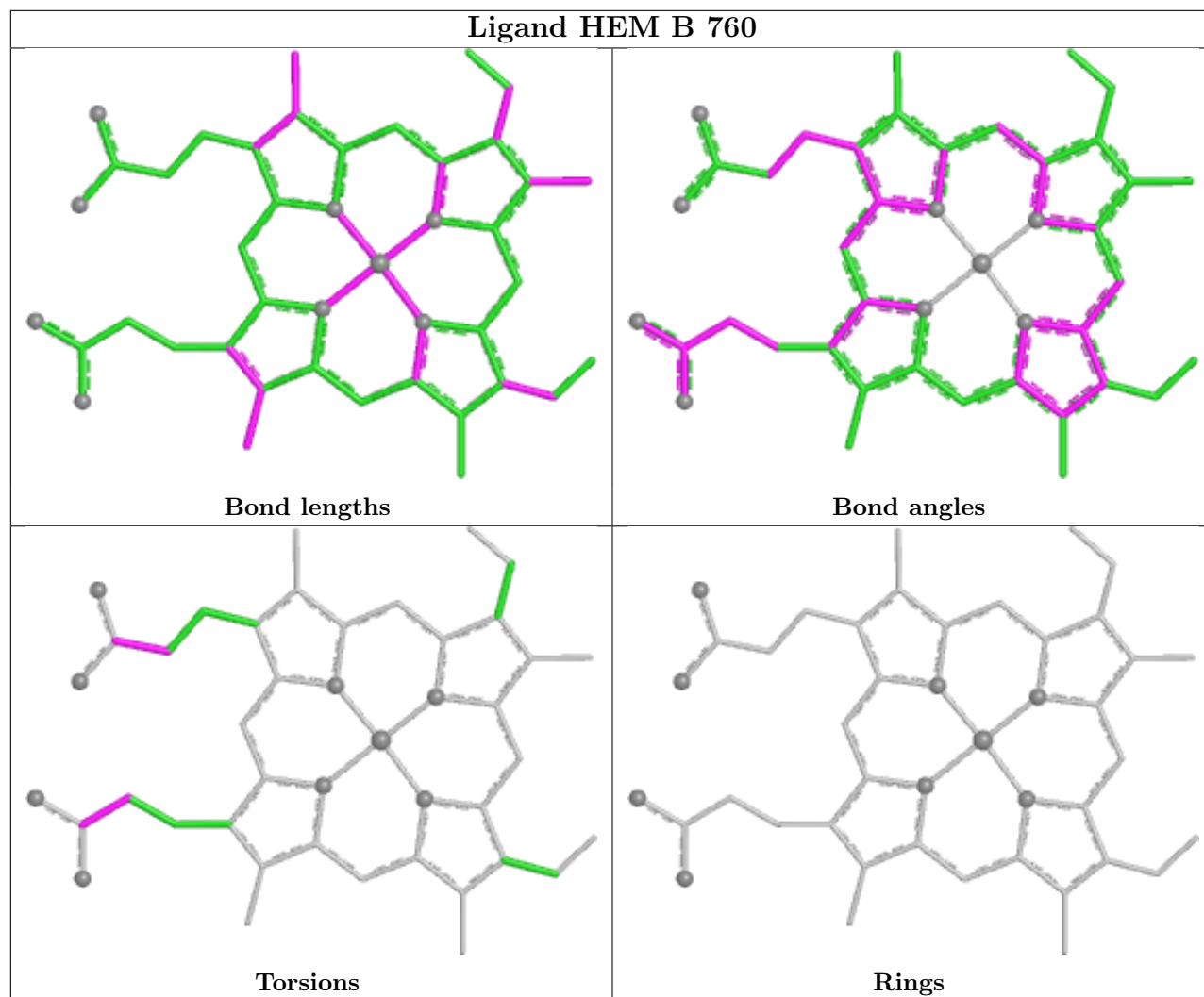
There are no ring outliers.

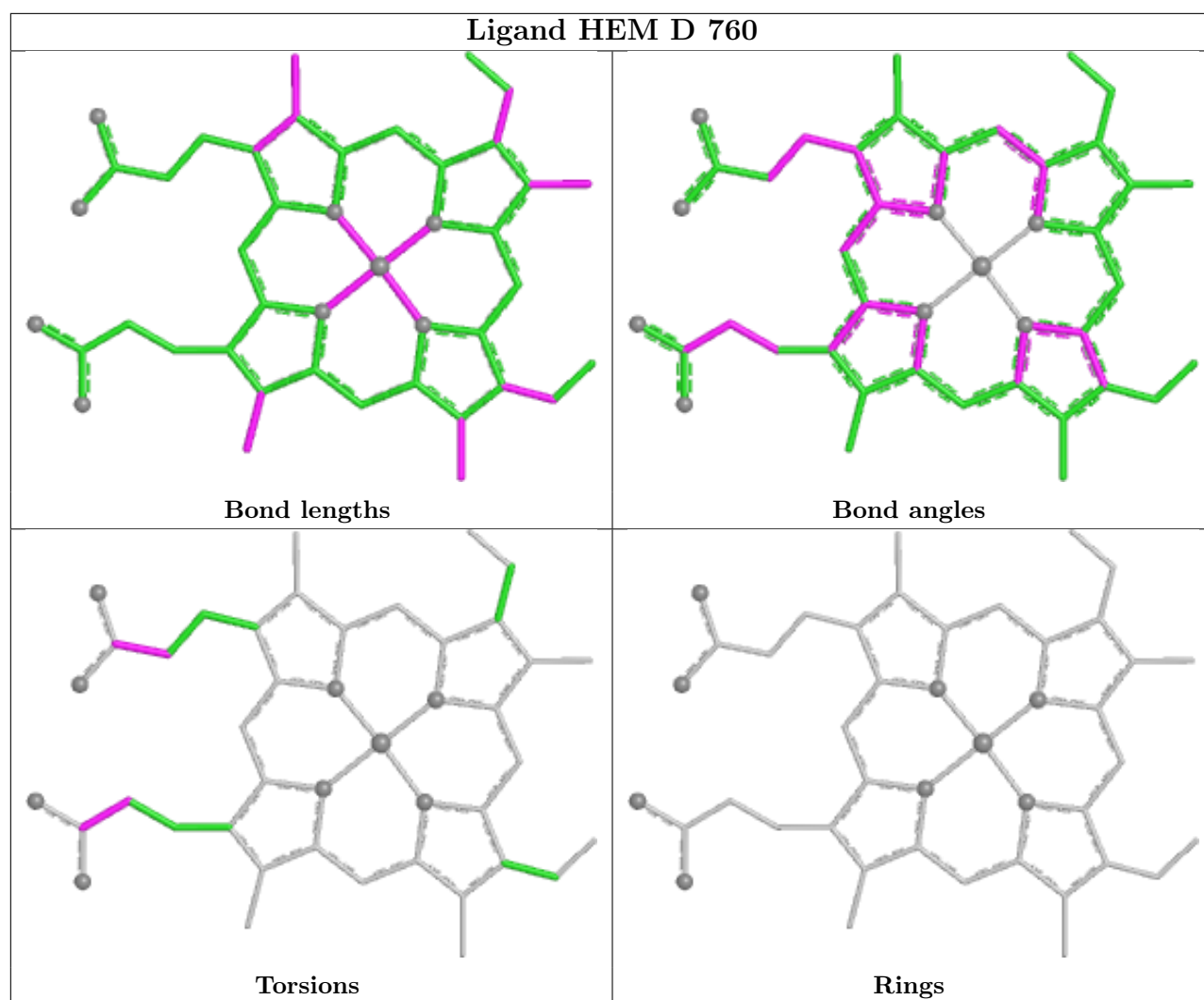
2 monomers are involved in 3 short contacts:

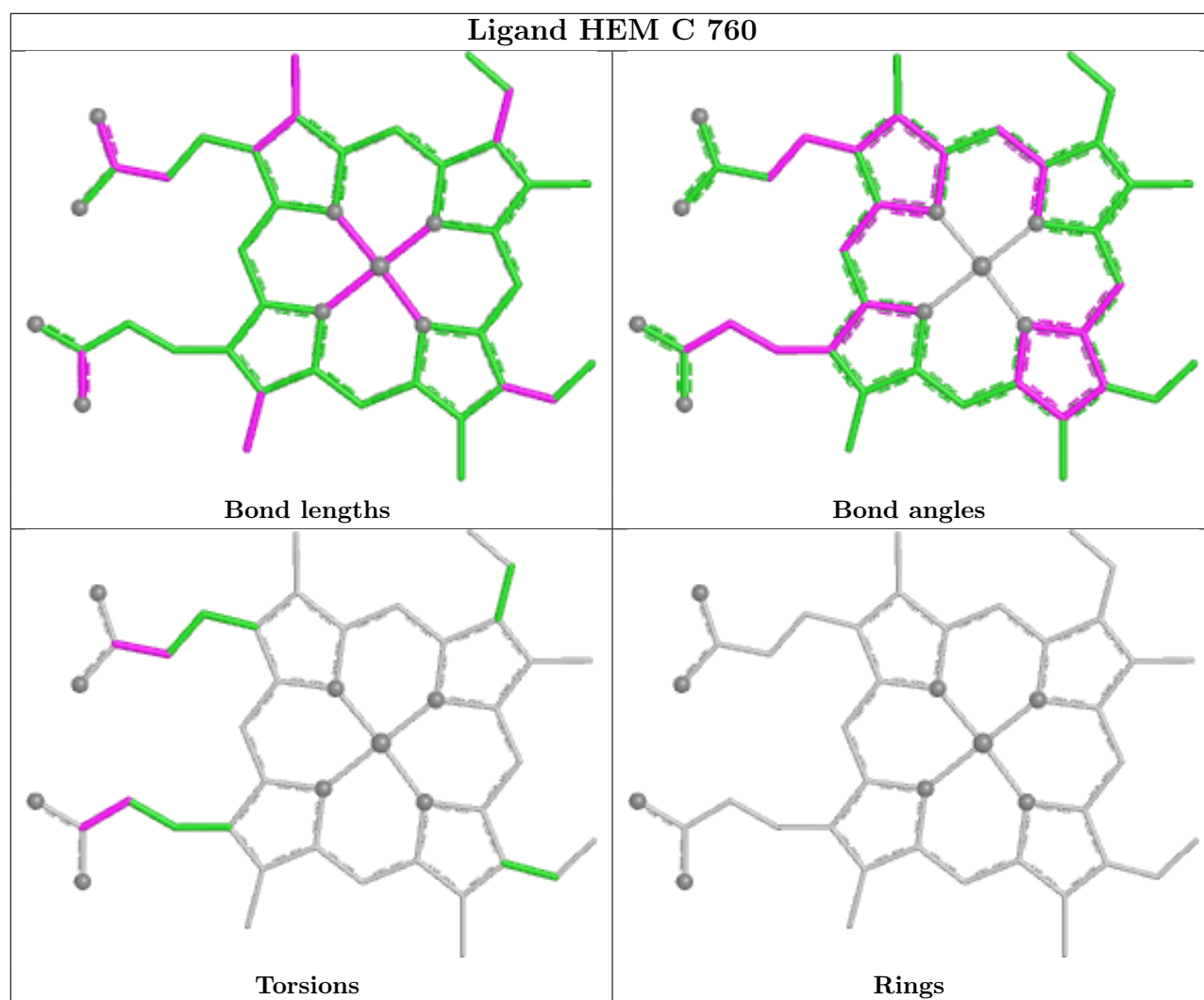
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	760	HEM	1	0
2	D	760	HEM	2	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	726/753 (96%)	-0.77	5 (0%) 84 86	6, 12, 27, 42	2 (0%)
1	B	726/753 (96%)	-0.68	0 100 100	6, 14, 34, 47	2 (0%)
1	C	726/753 (96%)	-0.72	1 (0%) 92 92	6, 14, 34, 46	2 (0%)
1	D	726/753 (96%)	-0.77	2 (0%) 90 91	6, 12, 28, 42	2 (0%)
All	All	2904/3012 (96%)	-0.73	8 (0%) 90 91	6, 13, 31, 47	8 (0%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	711	ALA	4.7
1	A	28	SER	4.0
1	A	710	ILE	3.9
1	A	712	ASP	2.6
1	D	28	SER	2.6
1	C	726	GLY	2.4
1	A	713	GLN	2.2
1	D	751	ILE	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](#)

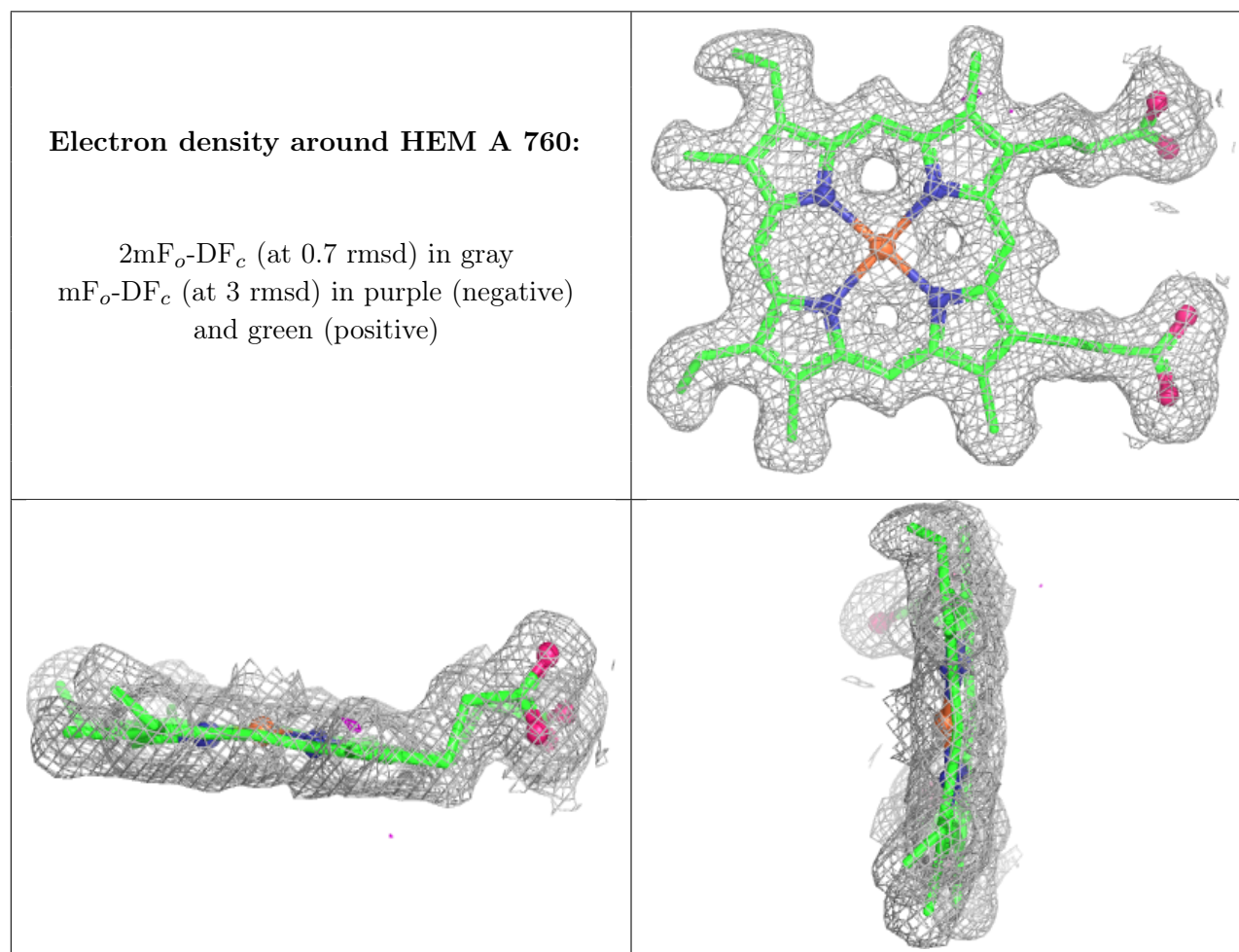
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

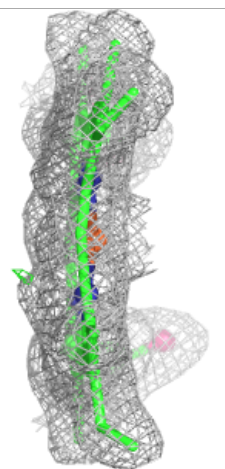
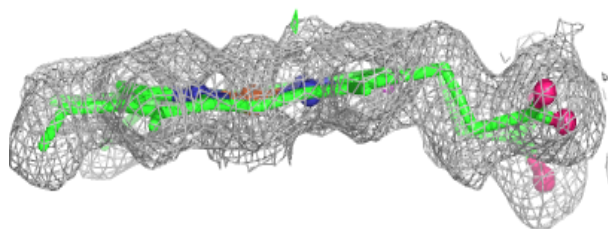
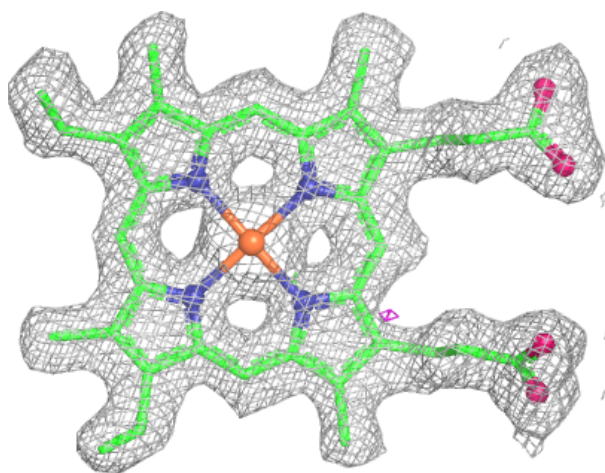
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	HEM	A	760	43/43	0.99	0.04	3,7,10,13	0
2	HEM	B	760	43/43	0.99	0.04	6,8,11,12	0
2	HEM	C	760	43/43	0.99	0.04	5,9,11,12	0
2	HEM	D	760	43/43	0.99	0.04	5,7,9,10	0

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.



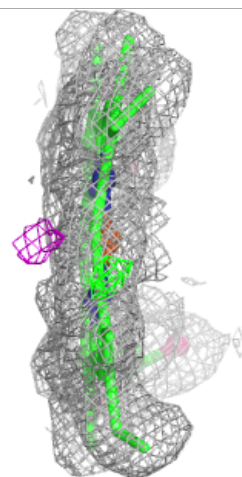
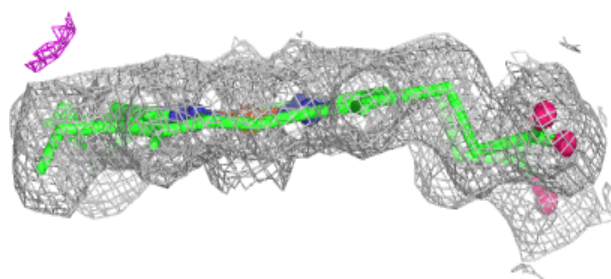
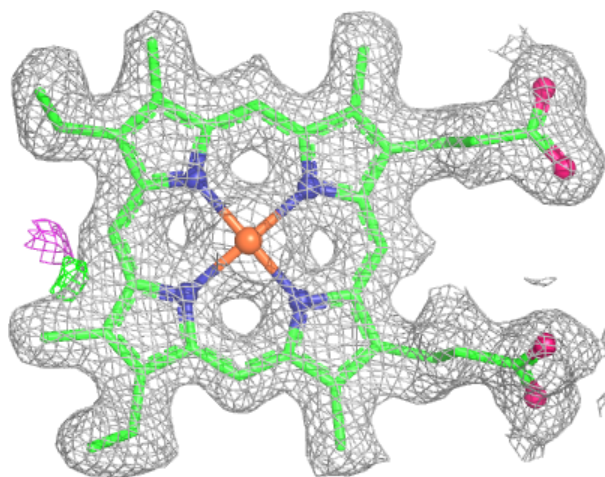
Electron density around HEM B 760:

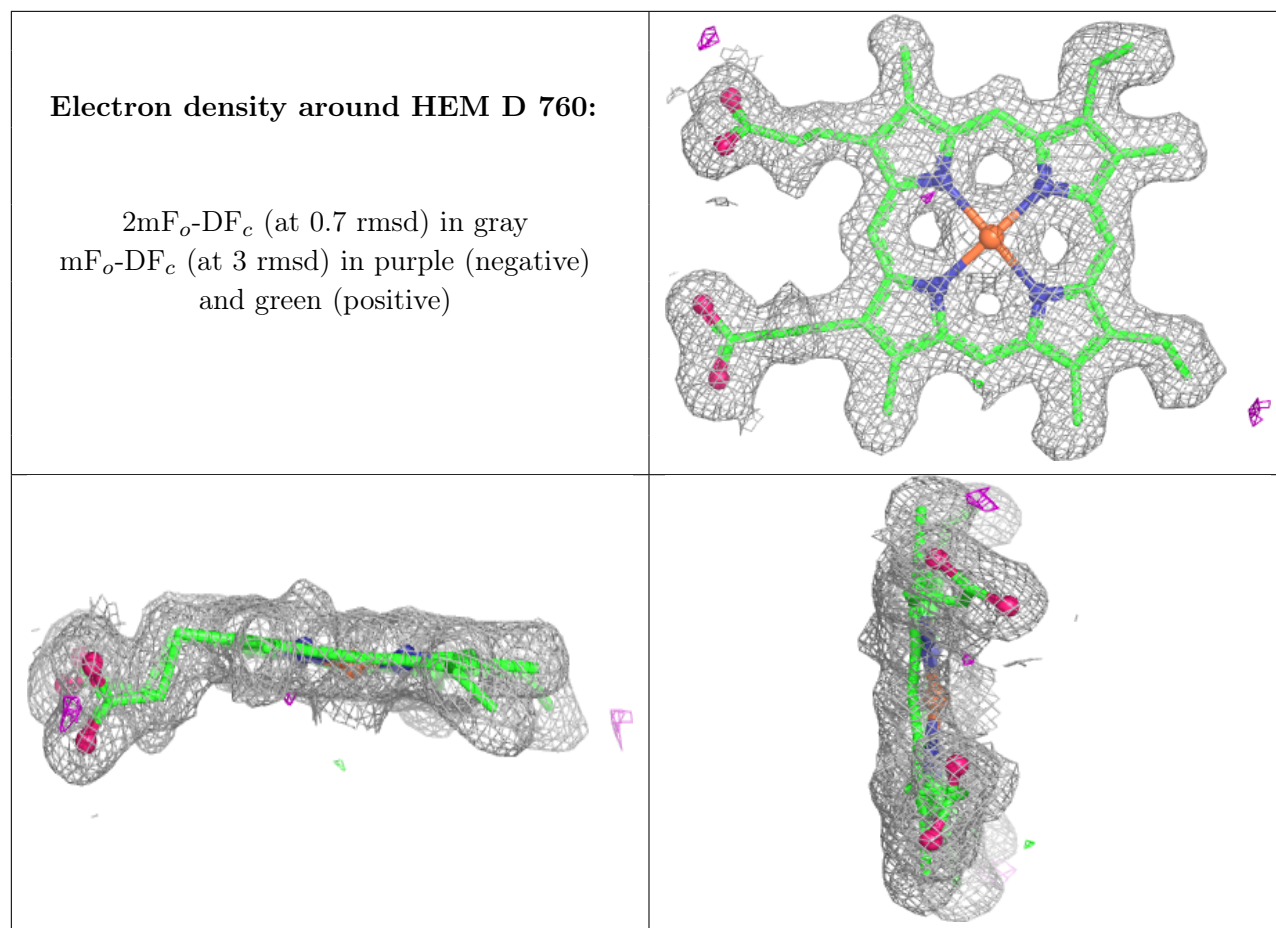
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM C 760:

$2mF_o-DF_c$ (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.