



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 03:11 pm BST

PDB ID : 4L1Q
Title : Crystal Structure of the E113Q-MauG/pre-Methylamine Dehydrogenase Complex
Authors : Yukl, E.Y.; Wilmot, C.M.
Deposited on : 2013-06-03
Resolution : 1.92 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

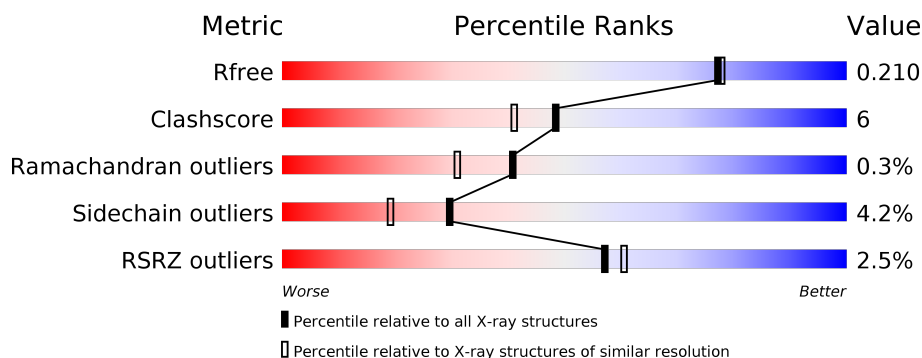
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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 on 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	373	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>5%</div> </div> </div>
1	B	373	<div> <div>%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>..</div> </div> </div>
2	C	137	<div> <div>5%</div> <div> <div></div> <div>77%</div> <div>15%</div> <div>..</div> </div> </div>
2	E	137	<div> <div>%</div> <div> <div></div> <div>73%</div> <div>17%</div> <div>9%</div> </div> </div>
3	D	385	<div> <div>4%</div> <div> <div></div> <div>82%</div> <div>15%</div> <div>..</div> </div> </div>
3	F	385	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div>..</div> </div> </div>

2 Entry composition i

There are 9 unique types of molecules in this entry. The entry contains 15249 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 Methylamine utilization protein MauG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	354	Total	C	N	O	S	0	6	0
			2795	1742	509	533	11			
1	B	357	Total	C	N	O	S	0	4	0
			2801	1745	510	535	11			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	113	GLN	GLU	ENGINEERED MUTATION	UNP Q51658
A	368	HIS	-	EXPRESSION TAG	UNP Q51658
A	369	HIS	-	EXPRESSION TAG	UNP Q51658
A	370	HIS	-	EXPRESSION TAG	UNP Q51658
A	371	HIS	-	EXPRESSION TAG	UNP Q51658
A	372	HIS	-	EXPRESSION TAG	UNP Q51658
A	373	HIS	-	EXPRESSION TAG	UNP Q51658
B	113	GLN	GLU	ENGINEERED MUTATION	UNP Q51658
B	368	HIS	-	EXPRESSION TAG	UNP Q51658
B	369	HIS	-	EXPRESSION TAG	UNP Q51658
B	370	HIS	-	EXPRESSION TAG	UNP Q51658
B	371	HIS	-	EXPRESSION TAG	UNP Q51658
B	372	HIS	-	EXPRESSION TAG	UNP Q51658
B	373	HIS	-	EXPRESSION TAG	UNP Q51658

- Molecule 2 is a protein called Methylamine dehydrogenase light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	131	Total	C	N	O	S	0	1	0
			1017	628	179	196	14			
2	E	125	Total	C	N	O	S	0	4	0
			976	604	163	194	15			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	132	HIS	-	EXPRESSION TAG	UNP A1BBA0
C	133	HIS	-	EXPRESSION TAG	UNP A1BBA0
C	134	HIS	-	EXPRESSION TAG	UNP A1BBA0
C	135	HIS	-	EXPRESSION TAG	UNP A1BBA0
C	136	HIS	-	EXPRESSION TAG	UNP A1BBA0
C	137	HIS	-	EXPRESSION TAG	UNP A1BBA0
E	132	HIS	-	EXPRESSION TAG	UNP A1BBA0
E	133	HIS	-	EXPRESSION TAG	UNP A1BBA0
E	134	HIS	-	EXPRESSION TAG	UNP A1BBA0
E	135	HIS	-	EXPRESSION TAG	UNP A1BBA0
E	136	HIS	-	EXPRESSION TAG	UNP A1BBA0
E	137	HIS	-	EXPRESSION TAG	UNP A1BBA0

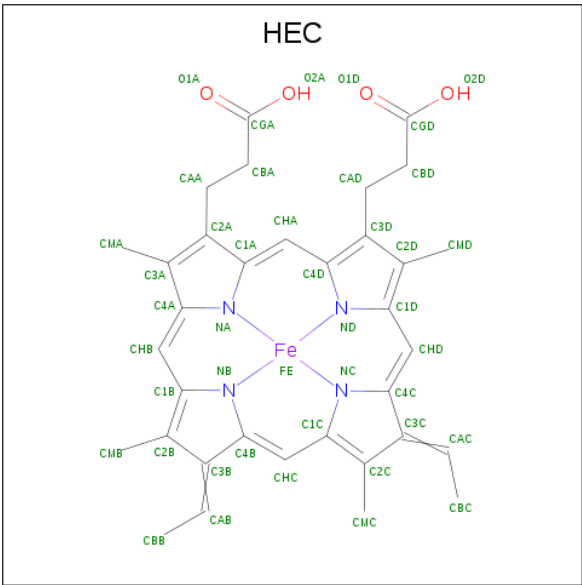
- Molecule 3 is a protein called Methylamine dehydrogenase heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	376	Total	C	N	O	S	0	2	0
			2943	1864	507	564	8			
3	F	376	Total	C	N	O	S	0	5	0
			2963	1875	511	568	9			

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

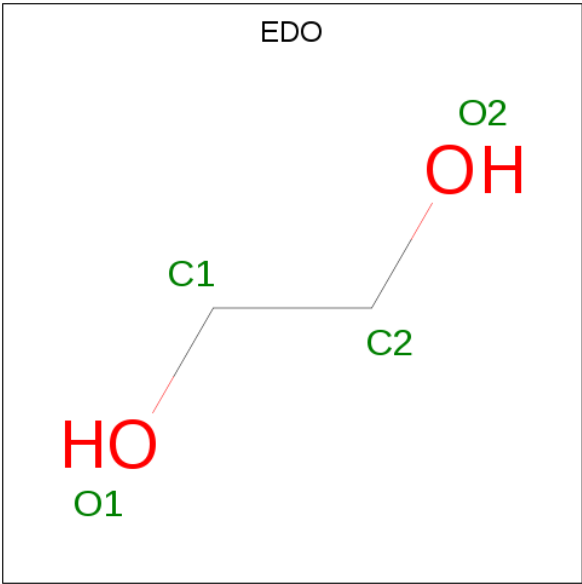
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca	0	0
			1	1		
4	A	1	Total	Ca	0	0
			1	1		

- Molecule 5 is HEME C (three-letter code: HEC) (formula: C₃₄H₃₄FeN₄O₄).



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

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).

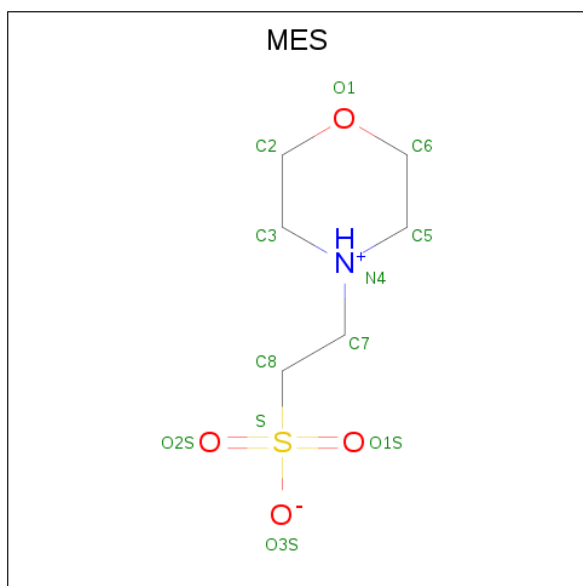


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	2	Total	Na	0	0
			2	2		
7	A	1	Total	Na	0	0
			1	1		

- Molecule 8 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
8	F	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

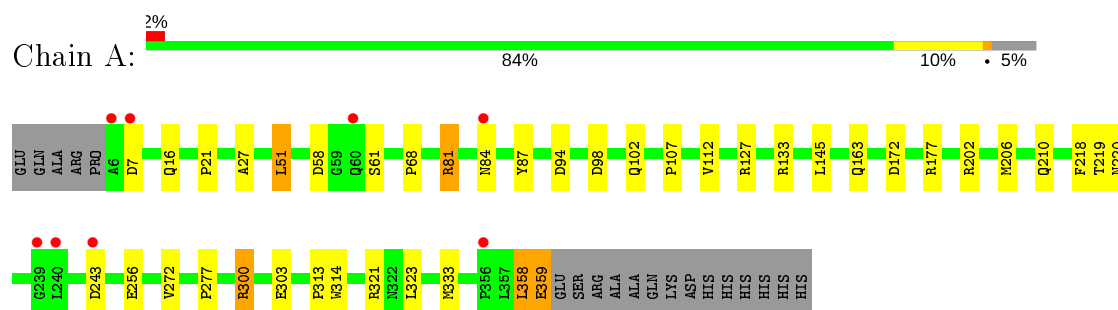
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	270	Total 274	O 274	0	4
9	B	384	Total 390	O 390	0	6
9	C	81	Total 81	O 81	0	0
9	D	272	Total 277	O 277	0	5
9	E	125	Total 125	O 125	0	0
9	F	389	Total 394	O 394	0	5

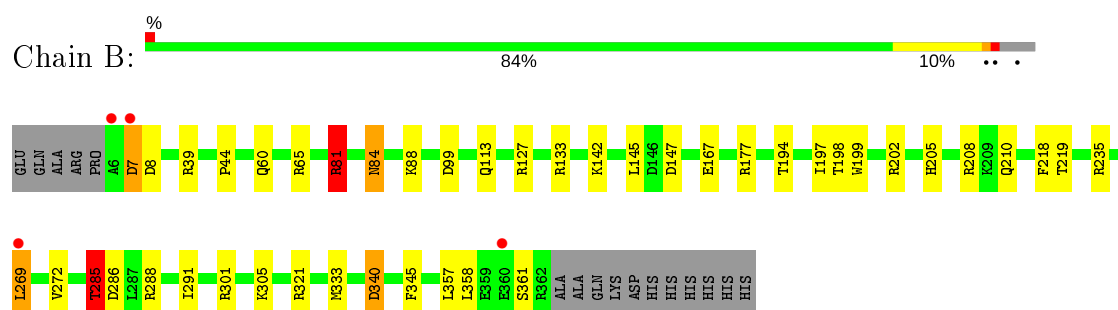
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

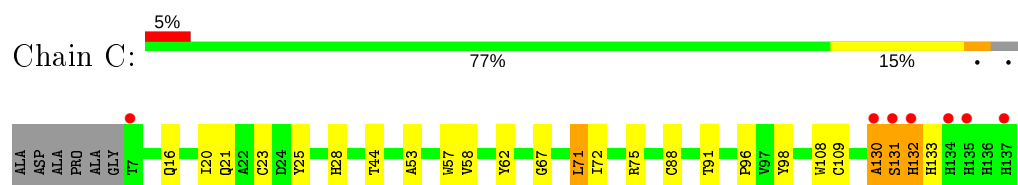
- Molecule 1: Methylamine utilization protein MauG



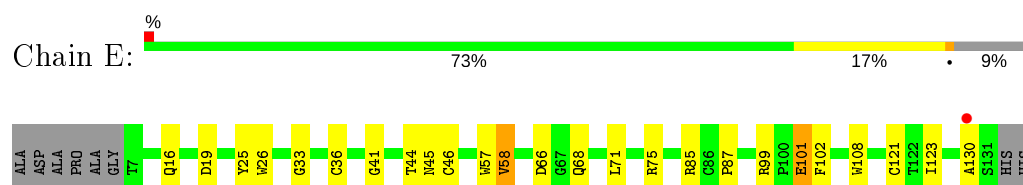
- Molecule 1: Methylamine utilization protein MauG



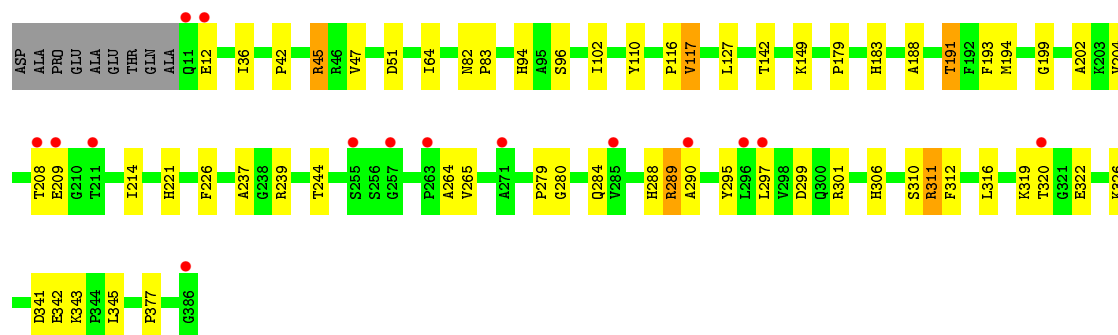
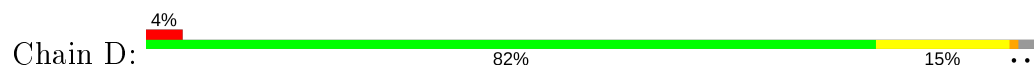
- Molecule 2: Methylamine dehydrogenase light chain



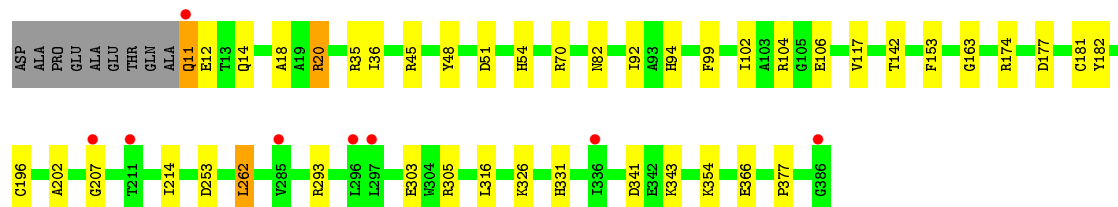
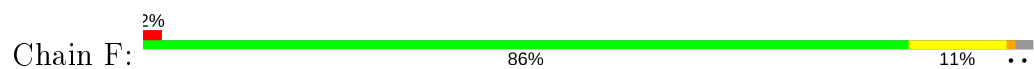
- Molecule 2: Methylamine dehydrogenase light chain



- Molecule 3: Methylamine dehydrogenase heavy chain



• Molecule 3: Methylamine dehydrogenase heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.53Å 83.52Å 107.78Å 109.94° 91.54° 105.78°	Depositor
Resolution (Å)	24.35 – 1.92 24.34 – 1.92	Depositor EDS
% Data completeness (in resolution range)	97.3 (24.35-1.92) 97.4 (24.34-1.92)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.49 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.160 , 0.207 0.169 , 0.210	Depositor DCC
R_{free} test set	6447 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	26.6	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15249	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 5.77% 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: 0AF, NA, CA, EDO, MES, HEC

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	0.91	0/2859	0.94	2/3876 (0.1%)
1	B	1.06	1/2865 (0.0%)	1.02	11/3884 (0.3%)
2	C	0.95	0/1035	0.96	2/1413 (0.1%)
2	E	1.20	3/991 (0.3%)	1.06	3/1353 (0.2%)
3	D	0.89	0/3020	0.93	1/4114 (0.0%)
3	F	1.08	1/3043 (0.0%)	1.06	3/4145 (0.1%)
All	All	1.00	5/13813 (0.0%)	0.99	22/18785 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	26	TRP	CB-CG	-5.96	1.39	1.50
2	E	101[A]	GLU	CD-OE1	-5.52	1.19	1.25
2	E	101[B]	GLU	CD-OE1	-5.52	1.19	1.25
1	B	345	PHE	CB-CG	-5.42	1.42	1.51
3	F	303	GLU	CD-OE2	-5.18	1.20	1.25

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	177	ARG	NE-CZ-NH1	5.97	123.29	120.30
2	C	75	ARG	NE-CZ-NH1	-5.85	117.38	120.30
3	F	35	ARG	NE-CZ-NH1	5.75	123.17	120.30
2	E	85	ARG	NE-CZ-NH1	5.68	123.14	120.30
3	F	305	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	B	99	ASP	CB-CG-OD1	5.68	123.41	118.30
1	B	235	ARG	NE-CZ-NH1	5.65	123.13	120.30
2	E	99	ARG	NE-CZ-NH2	-5.55	117.53	120.30
1	B	208	ARG	NE-CZ-NH1	-5.47	117.56	120.30
1	A	98	ASP	CB-CG-OD2	-5.30	113.53	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	341	ASP	CB-CG-OD1	5.29	123.06	118.30
1	B	285	THR	N-CA-CB	-5.29	100.26	110.30
1	B	65	ARG	NE-CZ-NH1	-5.25	117.68	120.30
1	B	81	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	A	94	ASP	CB-CG-OD1	5.22	123.00	118.30
1	B	81	ARG	NE-CZ-NH1	5.21	122.91	120.30
2	C	132	HIS	N-CA-C	5.12	124.84	111.00
1	B	333	MET	CG-SD-CE	-5.11	92.02	100.20
3	D	239	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	B	340	ASP	CB-CG-OD2	-5.09	113.72	118.30
1	B	142	LYS	CD-CE-NZ	-5.07	100.05	111.70
2	E	66	ASP	CB-CG-OD1	5.00	122.81	118.30

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	2795	0	2673	28	0
1	B	2801	0	2679	34	0
2	C	1017	0	906	31	0
2	E	976	0	883	19	0
3	D	2943	0	2825	36	0
3	F	2963	0	2842	25	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	86	0	60	5	0
5	B	86	0	60	3	0
6	A	4	0	6	0	0
6	B	8	0	12	0	0
7	A	1	0	0	0	0
7	B	2	0	0	0	0
8	D	12	0	13	0	0
8	F	12	0	13	0	0
9	A	274	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	B	390	0	0	5	0
9	C	81	0	0	3	0
9	D	277	0	0	3	0
9	E	125	0	0	2	0
9	F	394	0	0	4	0
All	All	15249	0	12972	160	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 6.

All (160) 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:C:131:SER:H	2:C:132:HIS:HA	0.91	1.08
2:C:131:SER:N	2:C:132:HIS:HA	1.55	1.06
2:E:130:ALA:HB3	9:E:293:HOH:O	1.69	0.93
2:C:131:SER:H	2:C:132:HIS:CA	1.81	0.92
1:B:198:THR:HG22	2:E:58[B]:VAL:HG23	1.50	0.91
1:B:198:THR:HG22	2:E:58[B]:VAL:CG2	2.04	0.86
3:F:181[B]:CYS:HB3	3:F:196:CYS:SG	2.17	0.84
2:C:131:SER:N	2:C:132:HIS:CA	2.39	0.82
1:A:206:MET:HE2	1:A:206:MET:HA	1.62	0.80
1:B:205:HIS:ND1	1:B:269[B]:LEU:HD11	1.96	0.80
9:C:262:HOH:O	3:F:36:ILE:HD11	1.84	0.77
2:C:131:SER:HA	2:C:132:HIS:HB2	1.65	0.77
1:A:300[A]:ARG:HD2	9:A:715:HOH:O	1.86	0.76
1:B:197:ILE:O	1:B:202:ARG:HD2	1.85	0.75
2:C:131:SER:HA	2:C:132:HIS:CB	2.15	0.75
3:D:342:GLU:HA	9:D:596:HOH:O	1.86	0.75
2:E:36[B]:CYS:SG	2:E:45:ASN:O	2.44	0.75
1:B:198:THR:CG2	2:E:58[B]:VAL:HG23	2.17	0.74
2:C:71:LEU:HD13	2:C:130:ALA:HA	1.68	0.73
1:B:210:GLN:HE22	2:E:44:THR:HG21	1.53	0.72
2:E:36[B]:CYS:SG	2:E:41:GLY:HA3	2.30	0.72
1:A:133[A]:ARG:NH1	9:A:764:HOH:O	2.20	0.71
1:B:198:THR:CG2	2:E:58[B]:VAL:CG2	2.70	0.69
1:B:269[B]:LEU:HD12	1:B:269[B]:LEU:N	2.07	0.69
3:F:12:GLU:OE1	3:F:20:ARG:NH1	2.25	0.69
2:C:57:OAF:CE3	2:C:108:TRP:CD1	2.76	0.69
2:C:130:ALA:O	2:C:131:SER:HB2	1.93	0.67
1:A:127[A]:ARG:NH2	9:A:761:HOH:O	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:HIS:CE1	1:B:269[B]:LEU:HD11	2.30	0.66
1:A:210:GLN:HE22	2:C:44:THR:HG21	1.61	0.65
1:B:301:ARG:NH2	3:F:177[B]:ASP:OD1	2.30	0.65
1:B:269[B]:LEU:N	1:B:269[B]:LEU:CD1	2.61	0.64
1:B:81:ARG:HD3	9:B:713:HOH:O	1.98	0.64
1:A:206:MET:CE	1:A:218:PHE:CD1	2.81	0.64
1:B:81:ARG:NH2	9:B:821:HOH:O	2.34	0.61
2:C:131:SER:CA	2:C:132:HIS:CB	2.77	0.61
3:F:14[A]:GLN:OE1	3:F:70:ARG:HD2	2.01	0.61
2:C:130:ALA:HB3	9:C:272:HOH:O	2.01	0.60
2:E:101[B]:GLU:HG2	2:E:102:PHE:CD2	2.36	0.60
1:A:206:MET:HE3	1:A:218:PHE:CD1	2.37	0.60
1:A:303[A]:GLU:OE1	1:A:303[A]:GLU:N	2.33	0.60
3:F:82:ASN:HB3	3:F:142:THR:HB	1.85	0.59
3:F:253:ASP:HB2	3:F:262:LEU:HD21	1.85	0.59
1:B:285:THR:HG22	1:B:286:ASP:OD1	2.03	0.58
3:D:82:ASN:HB3	3:D:142:THR:HB	1.85	0.58
2:C:132:HIS:H	2:C:133:HIS:HA	1.69	0.58
3:D:280:GLY:HA3	3:D:301:ARG:CZ	2.35	0.57
3:F:11:GLN:HG2	3:F:11:GLN:O	2.05	0.57
2:C:130:ALA:HB1	2:C:132:HIS:O	2.05	0.57
3:F:196:CYS:SG	3:F:202:ALA:HB2	2.44	0.57
3:D:237:ALA:HB2	3:D:289:ARG:HD2	1.86	0.56
1:B:194:THR:HG21	2:E:101[B]:GLU:HG3	1.88	0.56
3:D:341:ASP:OD1	3:D:341:ASP:N	2.32	0.56
1:B:305:LYS:HD2	9:B:882:HOH:O	2.07	0.55
1:A:303[A]:GLU:H	1:A:303[A]:GLU:CD	2.08	0.54
2:E:57:0AF:CE3	2:E:108:TRP:CD1	2.90	0.54
1:A:206:MET:HE1	1:A:218:PHE:CE1	2.42	0.54
3:D:96:SER:HB3	3:D:110:TYR:CZ	2.42	0.54
1:B:218:PHE:O	1:B:269[B]:LEU:HD13	2.08	0.54
3:D:188:ALA:HB3	3:D:191:THR:HG23	1.90	0.53
3:D:42:PRO:HG3	3:D:116:PRO:HB2	1.91	0.52
3:F:54:HIS:HE1	9:F:699:HOH:O	1.92	0.52
3:D:188:ALA:HB3	3:D:191:THR:CG2	2.40	0.52
3:D:204:VAL:HG22	3:D:214:ILE:HG12	1.91	0.52
1:A:163:GLN:HE22	5:A:402:HEC:HMA1	1.76	0.51
3:D:299:ASP:OD2	3:D:311[B]:ARG:CD	2.58	0.51
1:B:198:THR:CG2	2:E:58[B]:VAL:HG21	2.40	0.51
3:D:299:ASP:OD2	3:D:311[A]:ARG:HG2	2.10	0.51
3:F:331:HIS:HE1	3:F:366:GLU:OE1	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:ASP:C	1:A:58:ASP:OD1	2.49	0.50
2:C:131:SER:CA	2:C:132:HIS:HB2	2.40	0.50
2:C:132:HIS:N	2:C:133:HIS:HA	2.26	0.50
3:F:45:ARG:NH2	3:F:343:LYS:O	2.45	0.50
1:B:127:ARG:HG2	1:B:145:LEU:HB3	1.94	0.49
3:D:179:PRO:HD3	3:D:214:ILE:HD13	1.94	0.49
3:D:36:ILE:HD13	2:E:46:CYS:HB2	1.95	0.49
1:A:206:MET:HE3	1:A:218:PHE:HD1	1.78	0.49
1:A:313:PRO:O	1:A:314:TRP:C	2.50	0.49
2:E:36[B]:CYS:HA	2:E:121:CYS:SG	2.53	0.49
5:A:402:HEC:HBC3	5:A:402:HEC:HMC1	1.95	0.48
2:C:57:0AF:HE3	2:C:108:TRP:CD1	2.49	0.48
5:A:403:HEC:HBC3	5:A:403:HEC:HMC1	1.96	0.48
3:D:295:TYR:CD1	3:D:295:TYR:N	2.80	0.48
1:B:288:ARG:NH1	1:B:340:ASP:OD1	2.43	0.48
3:D:320:THR:OG1	3:D:322:GLU:HB3	2.14	0.48
1:A:206:MET:HE1	1:A:218:PHE:CD1	2.49	0.48
1:B:272:VAL:HG21	5:B:403:HEC:HMA3	1.94	0.48
3:D:42:PRO:HD3	3:D:117:VAL:HG12	1.96	0.47
1:B:198:THR:HG22	2:E:58[B]:VAL:HG21	1.92	0.47
3:F:153:PHE:CZ	3:F:163:GLY:HA3	2.49	0.47
1:A:51:LEU:HB3	1:A:277:PRO:HD3	1.98	0.46
3:D:279:PRO:HA	3:D:297:LEU:O	2.15	0.46
3:D:45:ARG:HD3	3:D:345:LEU:HD11	1.98	0.46
3:D:51:ASP:HA	3:D:377:PRO:HA	1.97	0.46
1:A:272:VAL:HG21	5:A:403:HEC:HMA3	1.97	0.46
3:F:51:ASP:HA	3:F:377:PRO:HA	1.98	0.46
1:B:147:ASP:C	1:B:147:ASP:OD1	2.54	0.46
1:B:197:ILE:O	1:B:202:ARG:CD	2.61	0.46
1:B:88:LYS:HE3	9:B:869:HOH:O	2.15	0.46
3:D:284:GLN:NE2	9:D:688:HOH:O	2.47	0.46
2:E:75:ARG:HA	9:E:287:HOH:O	2.16	0.46
1:A:107:PRO:HG3	9:A:757:HOH:O	2.16	0.46
3:D:226:PHE:O	3:D:244:THR:HA	2.16	0.45
3:D:264:ALA:HA	9:D:745:HOH:O	2.16	0.45
1:A:68:PRO:HG2	5:A:402:HEC:HBA1	1.99	0.45
2:C:20:ILE:HG22	2:C:25:TYR:CZ	2.51	0.45
1:B:113:GLN:HG2	5:B:402:HEC:HBC2	1.99	0.45
1:A:206:MET:O	1:A:220:ASN:HB3	2.17	0.45
1:A:112:VAL:HG23	1:A:112:VAL:O	2.17	0.45
1:A:21:PRO:O	1:A:27:ALA:HA	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:GLU:HA	9:A:551:HOH:O	2.16	0.44
3:D:288:HIS:HE1	3:D:290:ALA:HB3	1.82	0.44
2:C:91:THR:HB	3:D:306:HIS:CE1	2.52	0.44
1:A:133[A]:ARG:HG2	1:A:145:LEU:HD12	2.00	0.44
1:B:84:ASN:N	1:B:84:ASN:OD1	2.50	0.44
3:D:299:ASP:OD2	3:D:311[B]:ARG:HD2	2.17	0.44
3:D:193:PHE:HA	3:D:202:ALA:O	2.17	0.44
2:E:33:GLY:O	2:E:87:PRO:HA	2.18	0.44
3:F:153:PHE:CE2	3:F:163:GLY:HA3	2.53	0.43
3:D:47:VAL:HG13	3:D:64:ILE:HB	2.00	0.43
1:A:172:ASP:O	1:A:177:ARG:NH1	2.51	0.43
2:C:62:TYR:OH	2:C:67:GLY:HA2	2.18	0.43
1:A:81:ARG:HB2	1:A:87:TYR:CE1	2.53	0.43
2:C:58:VAL:HA	2:C:72:ILE:O	2.18	0.43
1:A:210:GLN:NE2	2:C:44:THR:HG21	2.30	0.43
2:C:23:CYS:SG	2:C:88[B]:CYS:SG	3.16	0.43
3:D:45:ARG:NH1	3:D:343:LYS:O	2.51	0.43
2:C:62:TYR:CZ	2:C:67:GLY:HA2	2.54	0.42
2:C:20:ILE:HD13	3:F:18:ALA:HB1	2.01	0.42
1:B:39:ARG:HD2	1:B:44:PRO:HB3	2.00	0.42
3:D:312:PHE:CD2	3:D:326:LYS:HE3	2.54	0.42
1:B:81:ARG:NH2	9:B:700:HOH:O	2.48	0.42
3:D:199:GLY:HA2	3:D:221:HIS:NE2	2.34	0.42
2:E:44:THR:HA	2:E:123:ILE:HB	2.01	0.42
2:C:23:CYS:SG	2:C:88[B]:CYS:HB2	2.60	0.42
1:A:61:SER:HB3	1:A:112:VAL:HB	2.02	0.41
3:D:297:LEU:HD22	3:D:310:SER:HB2	2.01	0.41
2:E:19:ASP:O	2:E:25:TYR:HB2	2.20	0.41
3:F:326:LYS:NZ	9:F:724:HOH:O	2.50	0.41
2:C:23:CYS:SG	2:C:88[B]:CYS:CB	3.08	0.41
3:D:191:THR:HA	3:D:204:VAL:O	2.21	0.41
3:F:214:ILE:HD12	9:F:539:HOH:O	2.20	0.41
2:C:96:PRO:HB2	2:C:98:TYR:CE1	2.56	0.41
2:C:21:GLN:HE22	3:F:14[A]:GLN:NE2	2.18	0.41
3:F:99:PHE:HA	3:F:106:GLU:O	2.21	0.41
2:C:25:TYR:HB3	2:C:28:HIS:CD2	2.55	0.41
3:D:320:THR:HG1	3:D:322:GLU:HB3	1.85	0.41
3:D:82:ASN:HA	3:D:83:PRO:HD2	1.93	0.41
3:F:366:GLU:OE2	9:F:743:HOH:O	2.22	0.41
3:D:183:HIS:O	3:D:194:MET:HG2	2.21	0.41
3:F:174[B]:ARG:NH1	3:F:207:GLY:O	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:ILE:HG12	5:B:403:HEC:HMB2	2.02	0.40
9:C:275:HOH:O	3:F:36:ILE:CD1	2.69	0.40
1:B:205:HIS:HD1	1:B:269[B]:LEU:HD11	1.83	0.40
1:B:7:ASP:OD1	1:B:7:ASP:N	2.54	0.40
2:C:53:ALA:HB2	2:C:109:CYS:HA	2.02	0.40
1:B:199:TRP:CD1	1:B:199:TRP:N	2.89	0.40
1:B:321:ARG:HG2	1:B:321:ARG:H	1.70	0.40
3:F:48:TYR:CZ	3:F:92:ILE:HG21	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	358/373 (96%)	345 (96%)	12 (3%)	1 (0%)	41	31
1	B	359/373 (96%)	350 (98%)	9 (2%)	0	100	100
2	C	129/137 (94%)	119 (92%)	8 (6%)	2 (2%)	9	2
2	E	126/137 (92%)	124 (98%)	2 (2%)	0	100	100
3	D	376/385 (98%)	360 (96%)	15 (4%)	1 (0%)	41	31
3	F	379/385 (98%)	366 (97%)	12 (3%)	1 (0%)	41	31
All	All	1727/1790 (96%)	1664 (96%)	58 (3%)	5 (0%)	41	31

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	130	ALA
2	C	131	SER
1	A	358	LEU
3	F	102	ILE
3	D	102	ILE

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	282/292 (97%)	265 (94%)	17 (6%)	19	9
1	B	283/292 (97%)	268 (95%)	15 (5%)	22	12
2	C	111/112 (99%)	109 (98%)	2 (2%)	59	53
2	E	108/112 (96%)	103 (95%)	5 (5%)	27	16
3	D	306/310 (99%)	291 (95%)	15 (5%)	25	14
3	F	309/310 (100%)	300 (97%)	9 (3%)	42	33
All	All	1399/1428 (98%)	1336 (96%)	63 (4%)	30	17

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASP
1	A	16	GLN
1	A	51	LEU
1	A	81	ARG
1	A	84	ASN
1	A	102	GLN
1	A	202	ARG
1	A	219	THR
1	A	243	ASP
1	A	256	GLU
1	A	300[A]	ARG
1	A	300[B]	ARG
1	A	321	ARG
1	A	323	LEU
1	A	333	MET
1	A	358	LEU
1	A	359	GLU
1	B	7	ASP
1	B	8	ASP
1	B	60	GLN
1	B	81	ARG
1	B	84	ASN

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Mol	Chain	Res	Type
1	B	133[A]	ARG
1	B	133[B]	ARG
1	B	167	GLU
1	B	219	THR
1	B	269[A]	LEU
1	B	269[B]	LEU
1	B	285	THR
1	B	357	LEU
1	B	358	LEU
1	B	361	SER
2	C	16	GLN
2	C	71	LEU
3	D	12	GLU
3	D	45	ARG
3	D	94	HIS
3	D	117	VAL
3	D	127	LEU
3	D	149	LYS
3	D	191	THR
3	D	208	THR
3	D	209	GLU
3	D	265	VAL
3	D	289	ARG
3	D	311[A]	ARG
3	D	311[B]	ARG
3	D	316	LEU
3	D	319	LYS
2	E	16	GLN
2	E	58[A]	VAL
2	E	58[B]	VAL
2	E	68	GLN
2	E	71	LEU
3	F	11	GLN
3	F	20	ARG
3	F	94	HIS
3	F	104	ARG
3	F	117	VAL
3	F	262	LEU
3	F	293	ARG
3	F	316	LEU
3	F	354	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	91	GLN
1	A	163	GLN
1	A	210	GLN
1	B	16	GLN
1	B	29	GLN
1	B	60	GLN
1	B	163	GLN
1	B	210	GLN
2	C	136	HIS
3	D	30	GLN
3	D	54	HIS
3	D	331	HIS
3	F	11	GLN
3	F	54	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	0AF	E	57	2	13,16,17	1.50	4 (30%)	11,22,24	1.40	1 (9%)
2	0AF	C	57	2	13,16,17	1.50	2 (15%)	11,22,24	1.37	1 (9%)

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	0AF	E	57	2	-	0/4/6/8	0/2/2/2
2	0AF	C	57	2	-	0/4/6/8	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	57	0AF	CZ2-CE2	-3.01	1.37	1.42
2	E	57	0AF	CB-CA	-2.53	1.48	1.53
2	E	57	0AF	CZ2-CE2	-2.50	1.38	1.42
2	C	57	0AF	CA-N	-2.38	1.40	1.48
2	E	57	0AF	CZ3-CE3	2.04	1.41	1.36
2	E	57	0AF	CD1-NE1	-2.04	1.32	1.36

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	57	0AF	CG-CB-CA	3.32	119.66	114.53
2	C	57	0AF	CG-CB-CA	3.08	119.29	114.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	57	0AF	1	0
2	C	57	0AF	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 5 are monoatomic - leaving 9 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	EDO	B	405	-	3,3,3	0.48	0	2,2,2	0.59	0
5	HEC	B	402	1	26,50,50	1.39	3 (11%)	18,82,82	2.39	8 (44%)
6	EDO	A	404	-	3,3,3	0.56	0	2,2,2	0.19	0
8	MES	F	401	-	12,12,12	1.38	1 (8%)	14,16,16	6.75	7 (50%)
8	MES	D	401	-	12,12,12	1.48	2 (16%)	14,16,16	5.92	8 (57%)
5	HEC	A	402	1	26,50,50	1.86	7 (26%)	18,82,82	3.32	9 (50%)
6	EDO	B	404	-	3,3,3	0.45	0	2,2,2	0.11	0
5	HEC	A	403	1	26,50,50	1.56	4 (15%)	18,82,82	3.00	8 (44%)
5	HEC	B	403	1	26,50,50	1.67	8 (30%)	18,82,82	2.50	6 (33%)

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
6	EDO	B	405	-	-	0/1/1/1	-
5	HEC	B	402	1	-	0/6/54/54	-
6	EDO	A	404	-	-	1/1/1/1	-
8	MES	F	401	-	-	3/6/14/14	0/1/1/1
8	MES	D	401	-	-	2/6/14/14	0/1/1/1
5	HEC	A	402	1	-	0/6/54/54	-
6	EDO	B	404	-	-	1/1/1/1	-
5	HEC	A	403	1	-	0/6/54/54	-
5	HEC	B	403	1	-	0/6/54/54	-

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	402	HEC	C3C-C4C	4.47	1.51	1.43
8	F	401	MES	C8-S	-4.45	1.71	1.77
5	B	403	HEC	C3C-C2C	-4.28	1.36	1.40
5	A	402	HEC	C1A-C2A	3.73	1.51	1.42
5	A	403	HEC	C3C-C2C	-3.70	1.36	1.40
8	D	401	MES	C8-S	-3.57	1.72	1.77
5	A	402	HEC	C3C-C2C	-3.39	1.37	1.40
5	A	402	HEC	C4A-C3A	3.22	1.49	1.42
5	A	402	HEC	C3B-C2B	-2.96	1.37	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	403	HEC	C4A-C3A	2.91	1.49	1.42
5	B	402	HEC	C3B-C4B	2.85	1.48	1.43
5	B	403	HEC	C3B-C2B	2.82	1.43	1.40
5	B	402	HEC	C1B-NB	-2.80	1.30	1.36
8	D	401	MES	O2S-S	2.69	1.53	1.45
5	B	403	HEC	C4A-C3A	2.66	1.48	1.42
5	B	403	HEC	C3C-C4C	2.58	1.47	1.43
5	A	403	HEC	C3C-C4C	2.34	1.47	1.43
5	B	402	HEC	C1C-NC	-2.32	1.31	1.36
5	B	403	HEC	CMD-C2D	-2.28	1.46	1.51
5	A	402	HEC	C1B-CHB	2.23	1.47	1.41
5	B	403	HEC	CBC-CAC	-2.22	1.41	1.49
5	B	403	HEC	C4D-ND	-2.20	1.31	1.36
5	A	403	HEC	CAA-C2A	-2.08	1.48	1.52
5	B	403	HEC	CAD-C3D	-2.08	1.49	1.52
5	A	402	HEC	C3B-C4B	2.00	1.46	1.43

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	F	401	MES	O3S-S-C8	-14.04	83.06	105.77
8	D	401	MES	O3S-S-O1S	-13.19	79.04	111.27
8	F	401	MES	O1S-S-C8	10.60	119.68	106.92
8	D	401	MES	O3S-S-O2S	-10.60	85.38	111.27
8	F	401	MES	O3S-S-O1S	-10.59	85.39	111.27
8	F	401	MES	O2S-S-C8	10.25	119.26	106.92
8	F	401	MES	O3S-S-O2S	-9.57	87.90	111.27
8	D	401	MES	O3S-S-C8	-8.95	91.29	105.77
8	D	401	MES	O2S-S-C8	8.14	116.72	106.92
5	A	403	HEC	CBD-CAD-C3D	-6.41	100.67	112.49
5	A	402	HEC	CMC-C2C-C3C	6.33	133.26	125.82
5	A	402	HEC	C1D-C2D-C3D	-6.33	102.59	107.00
5	A	403	HEC	CMB-C2B-C3B	6.33	133.26	125.82
5	B	403	HEC	CMC-C2C-C3C	6.21	133.12	125.82
5	B	403	HEC	CBD-CAD-C3D	-5.90	101.60	112.49
5	B	402	HEC	C1D-C2D-C3D	-5.44	103.21	107.00
5	A	402	HEC	CBD-CAD-C3D	-5.25	102.81	112.49
5	A	402	HEC	CMC-C2C-C1C	-4.73	121.19	128.46
5	A	403	HEC	CMC-C2C-C3C	4.70	131.35	125.82
8	D	401	MES	O2S-S-O1S	4.36	129.05	113.95
8	D	401	MES	O1S-S-C8	4.12	111.88	106.92
5	A	402	HEC	C4C-C3C-C2C	-4.11	101.92	106.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	403	HEC	CMA-C3A-C2A	3.66	131.84	124.94
5	B	402	HEC	CBD-CAD-C3D	-3.43	106.16	112.49
5	B	402	HEC	CMB-C2B-C3B	3.36	129.77	125.82
8	D	401	MES	C2-C3-N4	3.32	115.14	110.10
5	A	403	HEC	CBA-CAA-C2A	-3.29	106.42	112.48
5	A	402	HEC	C4B-C3B-C2B	-3.23	102.87	106.35
5	A	402	HEC	CMA-C3A-C2A	3.18	130.93	124.94
5	B	402	HEC	C4B-C3B-C2B	-3.14	102.96	106.35
5	B	403	HEC	CMC-C2C-C1C	-3.11	123.68	128.46
5	A	403	HEC	CMC-C2C-C1C	-3.08	123.74	128.46
5	B	402	HEC	C4C-C3C-C2C	-3.06	103.05	106.35
8	F	401	MES	C2-C3-N4	-2.76	105.91	110.10
5	A	402	HEC	C3B-C4B-NB	2.75	116.14	110.94
5	B	402	HEC	CMC-C2C-C3C	2.74	129.04	125.82
5	A	403	HEC	C1D-C2D-C3D	-2.66	105.14	107.00
5	B	403	HEC	CMB-C2B-C3B	2.55	128.82	125.82
5	A	403	HEC	CMB-C2B-C1B	-2.53	124.57	128.46
5	B	403	HEC	CAD-CBD-CGD	-2.38	108.67	112.67
5	A	402	HEC	CBA-CAA-C2A	-2.29	108.26	112.48
5	B	402	HEC	CMD-C2D-C3D	2.26	129.20	124.94
5	B	403	HEC	C3B-C4B-NB	2.21	115.12	110.94
8	D	401	MES	C5-N4-C3	2.12	113.60	108.83
8	F	401	MES	C6-O1-C2	2.08	116.83	109.89
5	B	402	HEC	CMA-C3A-C2A	2.06	128.82	124.94

There are no chirality outliers.

All (7) torsion outliers are listed below:

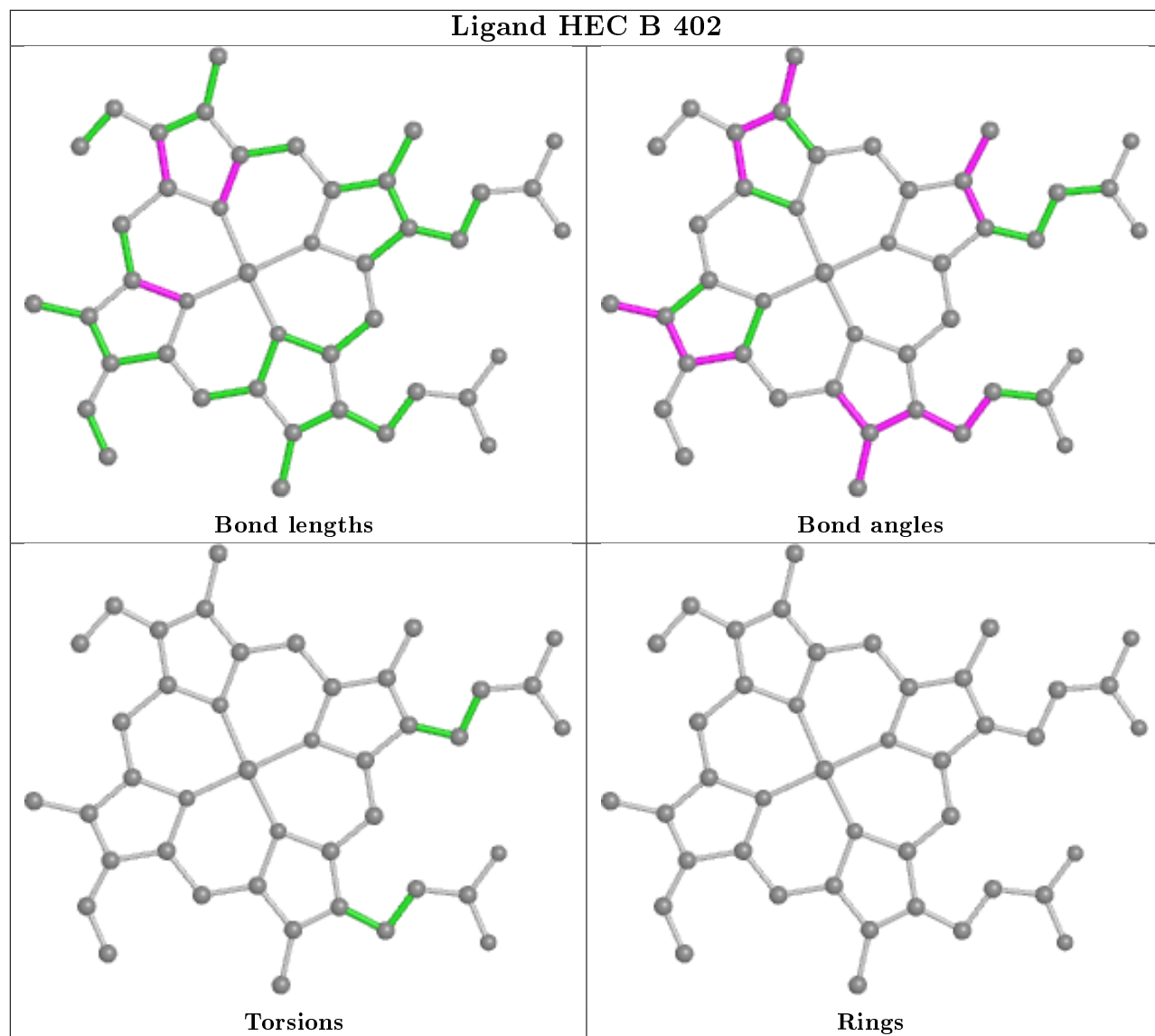
Mol	Chain	Res	Type	Atoms
8	F	401	MES	N4-C7-C8-S
8	F	401	MES	C7-C8-S-O3S
8	D	401	MES	C7-C8-S-O1S
6	B	404	EDO	O1-C1-C2-O2
6	A	404	EDO	O1-C1-C2-O2
8	F	401	MES	C7-C8-S-O1S
8	D	401	MES	C7-C8-S-O2S

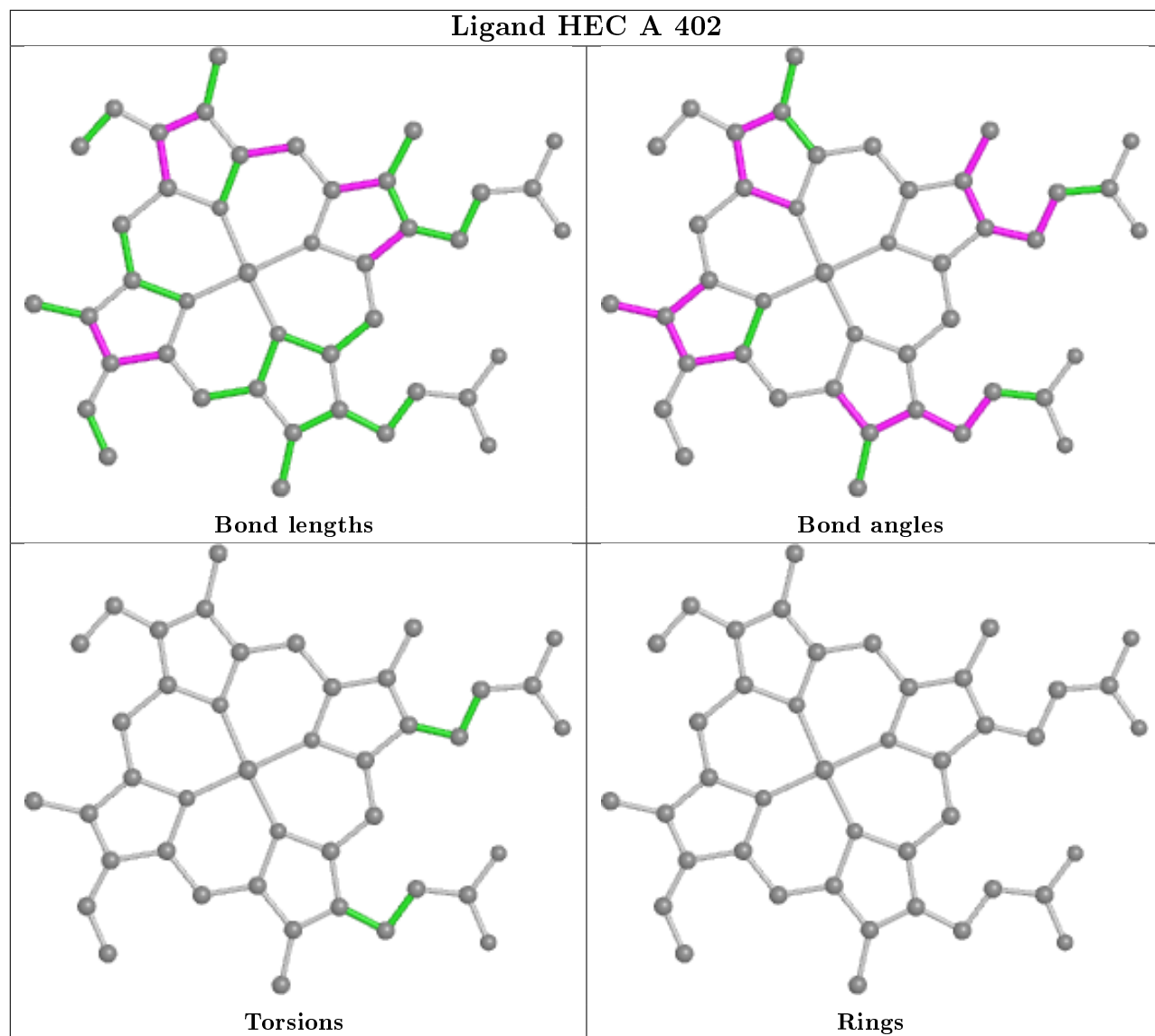
There are no ring outliers.

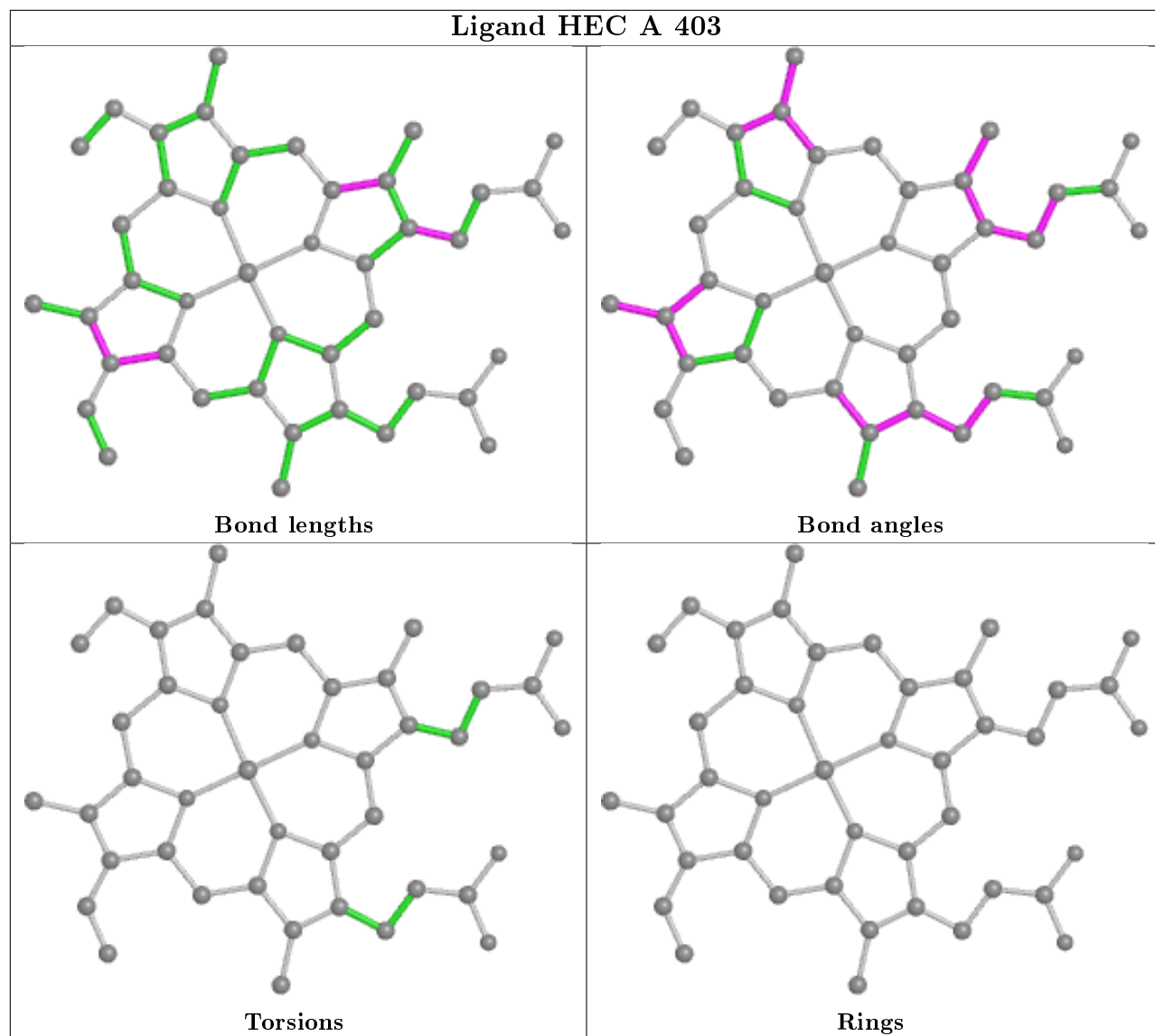
4 monomers are involved in 8 short contacts:

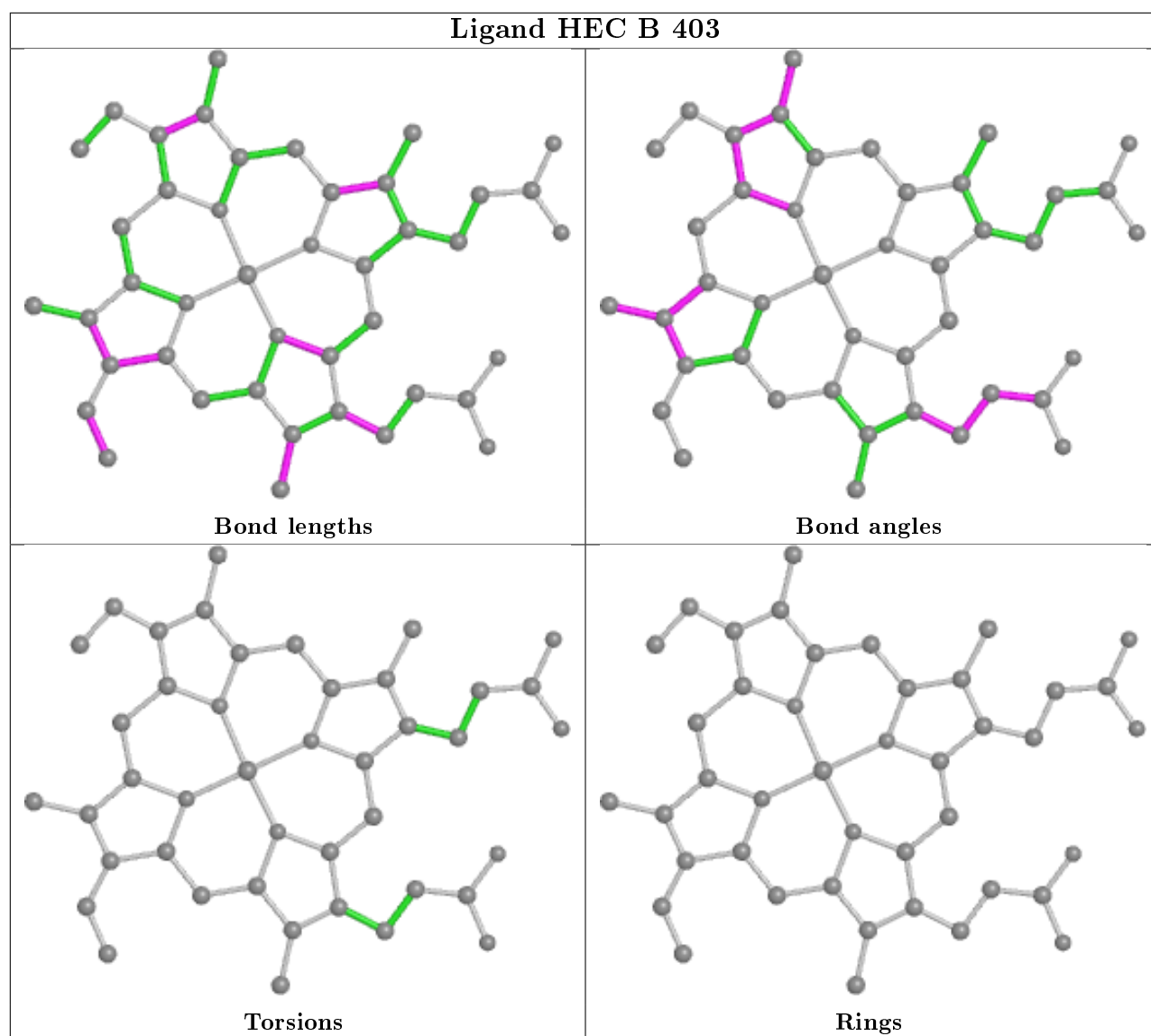
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	402	HEC	1	0
5	A	402	HEC	3	0
5	A	403	HEC	2	0
5	B	403	HEC	2	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	354/373 (94%)	-0.15	8 (2%)	60 63	26, 38, 56, 73	0
1	B	357/373 (95%)	-0.37	4 (1%)	80 82	16, 29, 47, 76	0
2	C	130/137 (94%)	0.11	7 (5%)	25 29	23, 36, 64, 90	2 (1%)
2	E	124/137 (90%)	-0.42	1 (0%)	86 87	20, 26, 38, 71	0
3	D	376/385 (97%)	0.10	15 (3%)	38 41	24, 41, 66, 85	0
3	F	376/385 (97%)	-0.28	8 (2%)	63 66	18, 28, 47, 68	0
All	All	1717/1790 (95%)	-0.17	43 (2%)	57 60	16, 33, 59, 90	2 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	131	SER	7.0
1	B	6	ALA	5.4
1	A	6	ALA	5.2
3	D	386	GLY	4.9
3	F	386	GLY	4.5
3	D	11	GLN	3.9
1	B	7	ASP	3.9
2	C	134	HIS	3.8
3	D	208	THR	3.3
2	C	132	HIS	3.3
3	D	257	GLY	3.3
2	C	7	THR	3.3
3	D	263	PRO	3.2
3	D	285	VAL	3.2
3	D	209	GLU	3.2
3	D	271	ALA	3.1
3	D	255	SER	2.8
3	D	290	ALA	2.7
3	D	296	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	7	ASP	2.6
1	A	243	ASP	2.5
1	A	240	LEU	2.5
1	A	60	GLN	2.5
1	B	360	GLU	2.5
3	F	211	THR	2.4
3	F	297	LEU	2.4
3	D	320	THR	2.4
2	C	137	HIS	2.3
3	F	207	GLY	2.2
3	D	211	THR	2.2
3	F	296	LEU	2.2
1	A	356	PRO	2.2
2	C	135	HIS	2.2
3	F	11	GLN	2.2
3	F	285	VAL	2.2
1	A	239	GLY	2.1
3	D	297	LEU	2.1
2	C	130	ALA	2.1
2	E	130	ALA	2.1
3	F	336	ILE	2.1
3	D	12	GLU	2.1
1	B	269[A]	LEU	2.0
1	A	84	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	0AF	C	57	15/16	0.96	0.10	37,40,45,46	0
2	0AF	E	57	15/16	0.97	0.11	24,27,32,33	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands

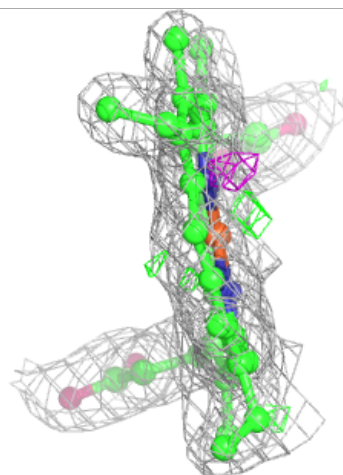
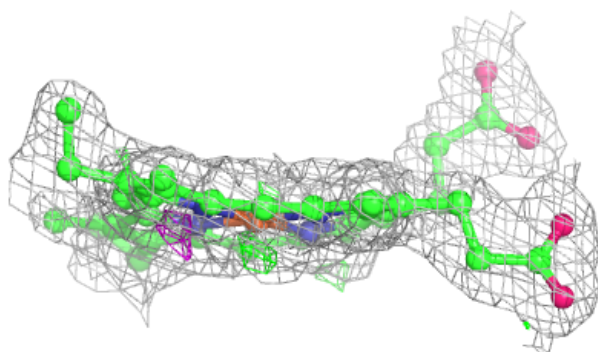
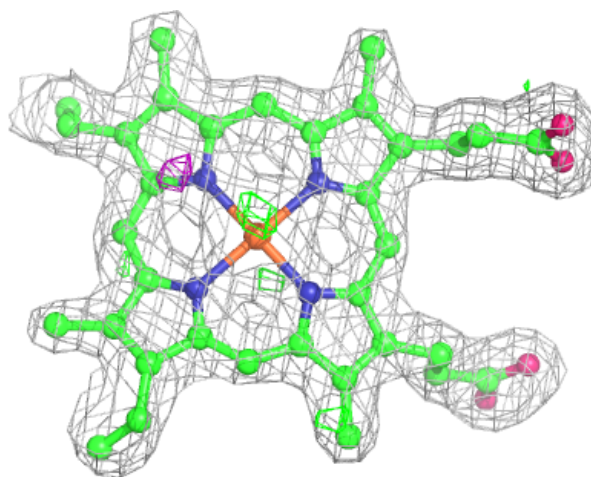
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	EDO	A	404	4/4	0.84	0.10	57,57,59,59	0
8	MES	D	401	12/12	0.90	0.21	29,54,60,61	12
6	EDO	B	405	4/4	0.91	0.18	37,50,50,60	0
8	MES	F	401	12/12	0.94	0.18	44,66,68,70	0
6	EDO	B	404	4/4	0.95	0.07	46,47,50,53	0
5	HEC	A	402	43/43	0.97	0.10	28,31,34,36	0
7	NA	A	405	1/1	0.97	0.17	47,47,47,47	0
7	NA	B	407	1/1	0.98	0.15	31,31,31,31	0
5	HEC	B	402	43/43	0.98	0.08	19,23,26,29	0
5	HEC	A	403	43/43	0.98	0.13	26,29,32,32	0
7	NA	B	406	1/1	0.98	0.13	30,30,30,30	0
5	HEC	B	403	43/43	0.99	0.10	15,17,20,21	0
4	CA	A	401	1/1	0.99	0.02	30,30,30,30	0
4	CA	B	401	1/1	1.00	0.06	20,20,20,20	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.

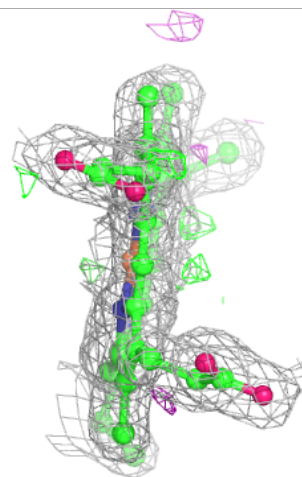
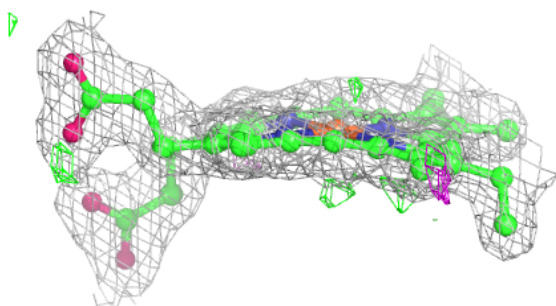
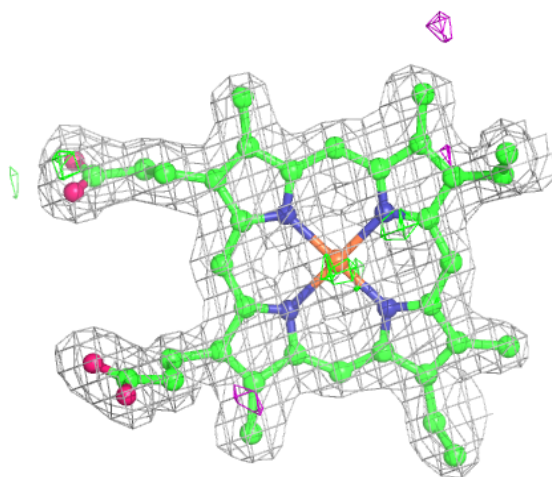
Electron density around HEC A 402:

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



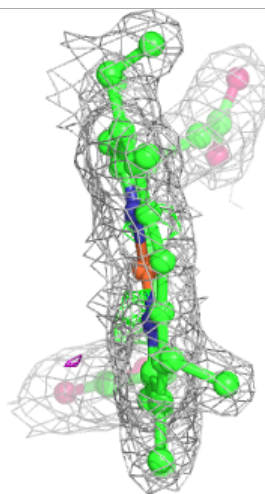
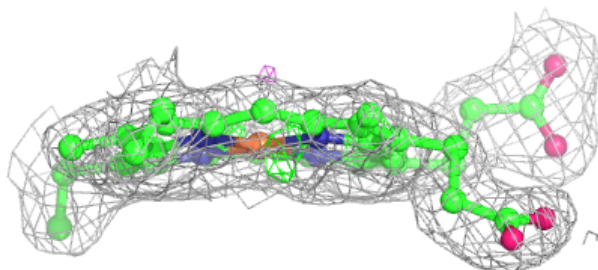
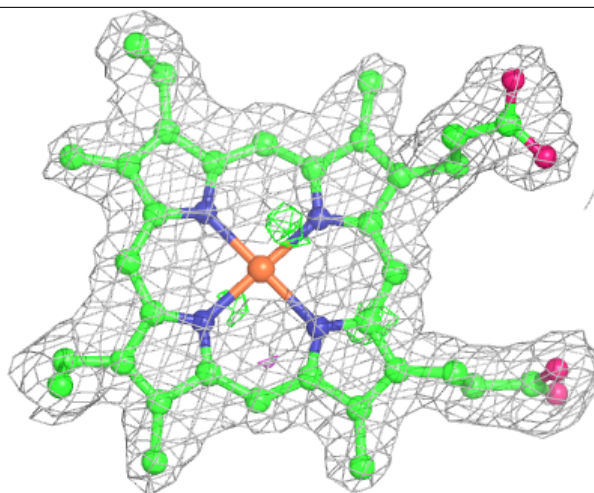
Electron density around HEC B 402:

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



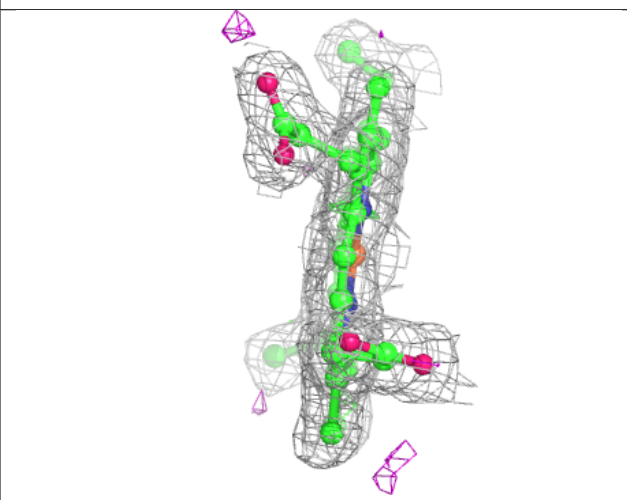
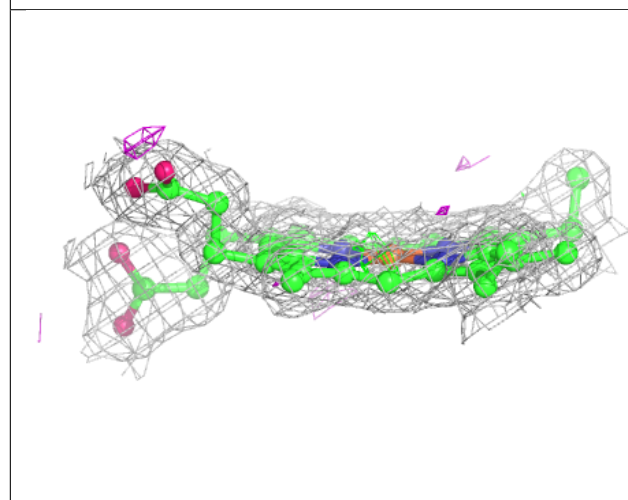
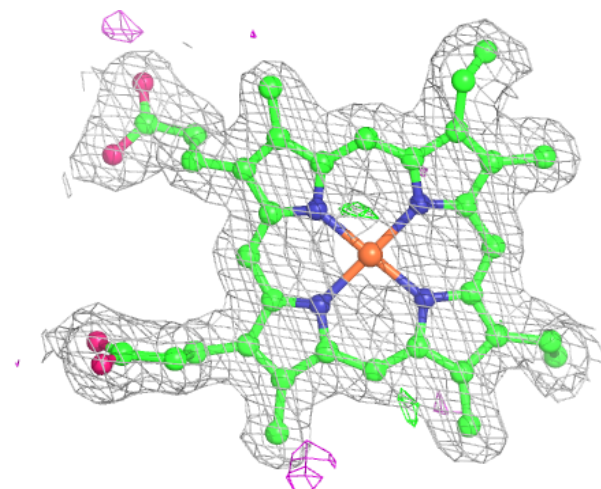
Electron density around HEC A 403:

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



Electron density around HEC B 403:

$2mF_o-DF_c$ (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.