



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 12:57 pm BST

PDB ID : 4FA1
Title : Crystal Structure of WT MauG in Complex with Pre-Methylamine Dehydrogenase Aged 130 Days.
Authors : Yukl, E.T.; Wilmot, C.M.
Deposited on : 2012-05-21
Resolution : 2.18 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

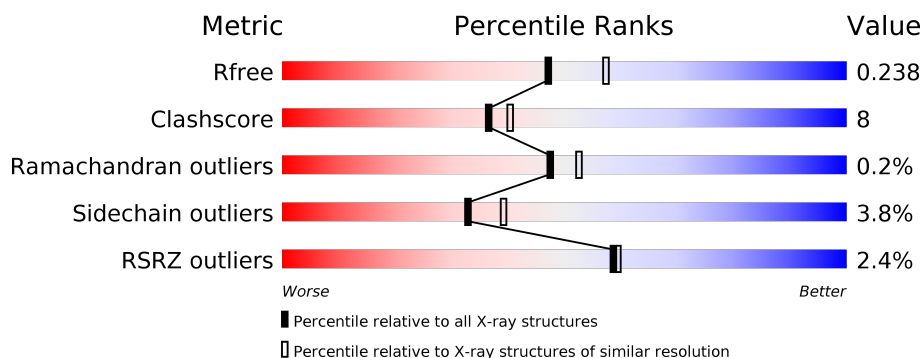
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.18 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	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

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>76%</div> <div>17%</div> <div>• 5%</div> </div>
1	B	373	<div> <div>81%</div> <div>13%</div> <div>• 5%</div> </div>
2	C	137	<div> <div>4%</div> <div>69%</div> <div>20%</div> <div>• 9%</div> </div>
2	E	137	<div> <div>3%</div> <div>73%</div> <div>17%</div> <div>• 9%</div> </div>
3	D	385	<div> <div>5%</div> <div>80%</div> <div>17%</div> <div>• •</div> </div>
3	F	385	<div> <div>80%</div> <div>17%</div> <div>• •</div> </div>

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 14647 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	5	0
			2772	1731	499	531	11			
1	B	355	Total	C	N	O	S	0	7	0
			2809	1751	512	534	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
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	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	125	Total	C	N	O	S	0	1	0
			958	592	161	191	14			
2	E	125	Total	C	N	O	S	0	2	0
			964	596	161	193	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	132	HIS	-	EXPRESSION TAG	UNP P22619
C	133	HIS	-	EXPRESSION TAG	UNP P22619
C	134	HIS	-	EXPRESSION TAG	UNP P22619
C	135	HIS	-	EXPRESSION TAG	UNP P22619
C	136	HIS	-	EXPRESSION TAG	UNP P22619
C	137	HIS	-	EXPRESSION TAG	UNP P22619
E	132	HIS	-	EXPRESSION TAG	UNP P22619
E	133	HIS	-	EXPRESSION TAG	UNP P22619
E	134	HIS	-	EXPRESSION TAG	UNP P22619
E	135	HIS	-	EXPRESSION TAG	UNP P22619
E	136	HIS	-	EXPRESSION TAG	UNP P22619
E	137	HIS	-	EXPRESSION TAG	UNP P22619

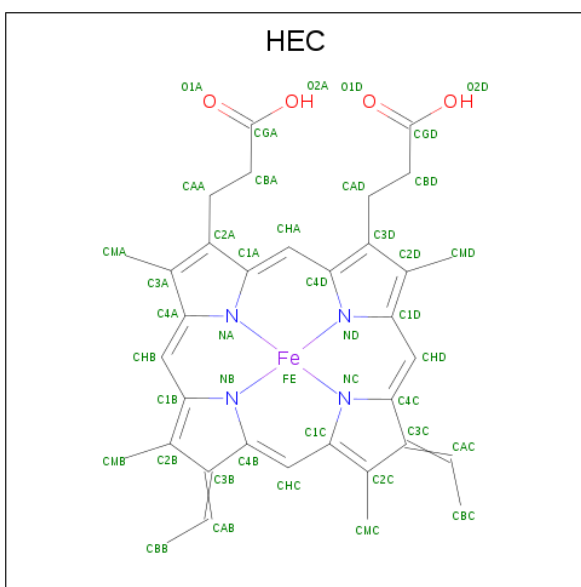
- Molecule 3 is a protein called Methylamine dehydrogenase heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	376	Total	C	N	O	S	0	4	0
			2945	1867	505	564	9			
3	F	376	Total	C	N	O	S	0	8	0
			2983	1891	514	570	8			

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

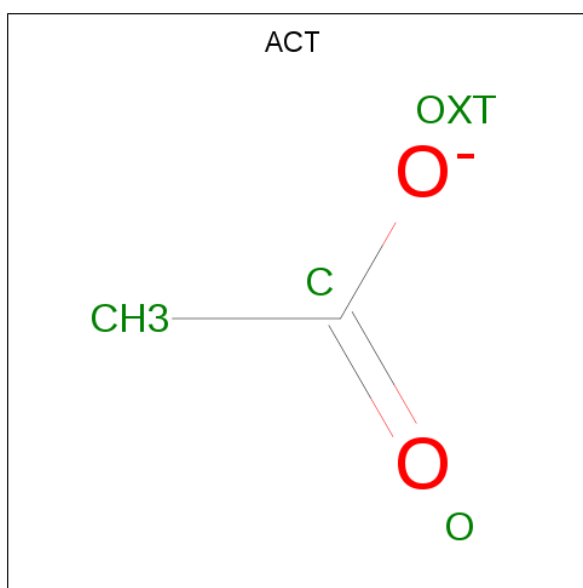
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total Na 1 1	0	0
6	A	1	Total Na 1 1	0	0

- Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



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

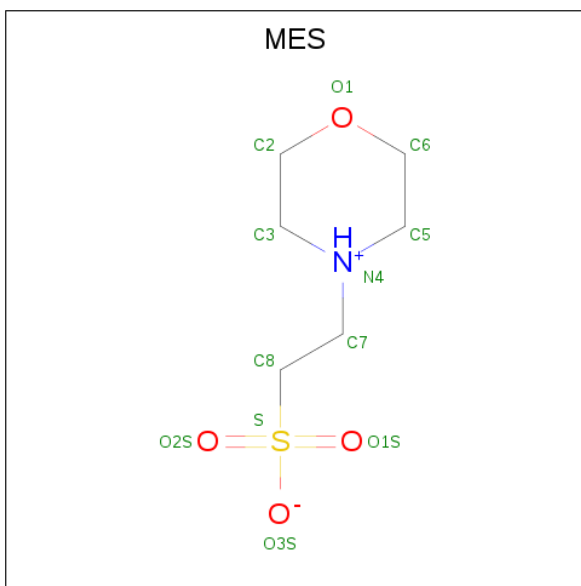
- Molecule 8 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 9 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES)

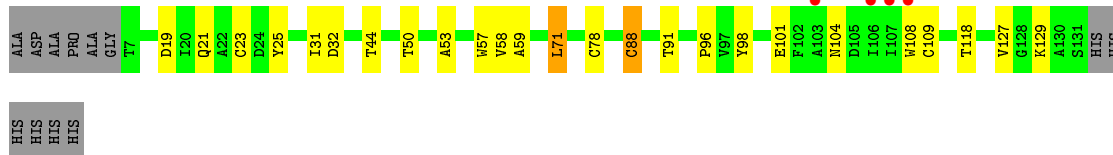
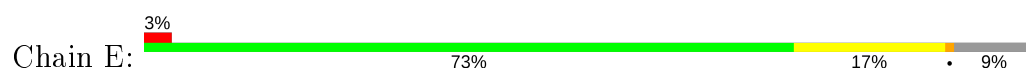
(formula: C₆H₁₃NO₄S).



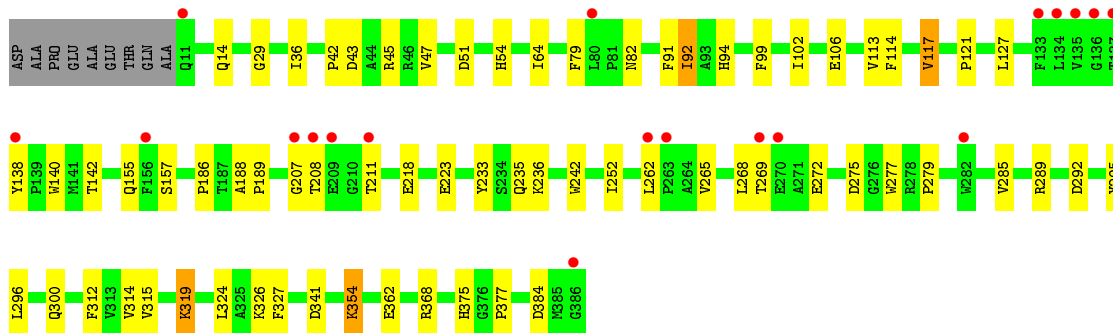
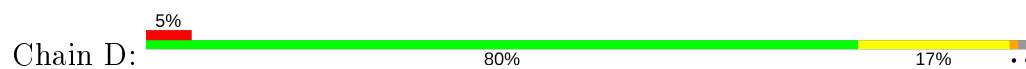
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	F	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 10 is water.

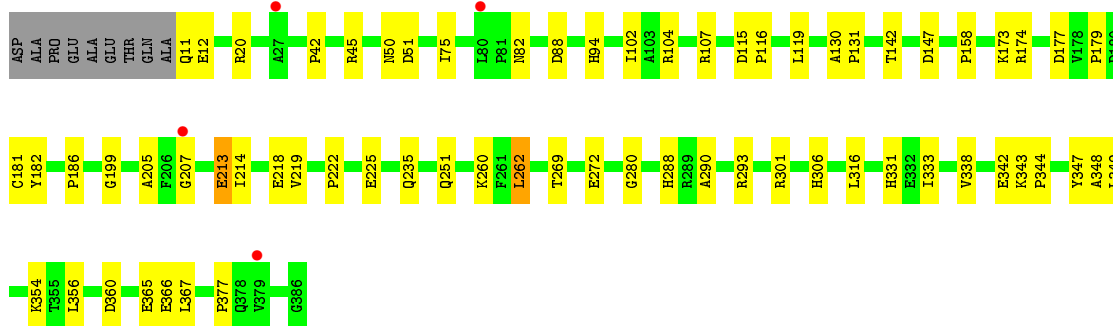
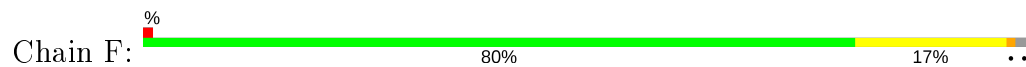
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	142	Total	O	0	2
			144	144		
10	B	230	Total	O	0	2
			232	232		
10	C	61	Total	O	0	0
			61	61		
10	D	188	Total	O	0	0
			188	188		
10	E	77	Total	O	0	0
			77	77		
10	F	313	Total	O	0	1
			314	314		



- 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)	37.17 – 2.18 37.17 – 2.18	Depositor EDS
% Data completeness (in resolution range)	93.7 (37.17-2.18) 93.7 (37.17-2.18)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.18Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.157 , 0.227 0.171 , 0.238	Depositor DCC
R_{free} test set	4239 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	31.4	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14647	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% 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: NA, CA, HEC, EDO, MES, ACT, TRQ

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	0.80	0/2845	0.84	1/3859 (0.0%)
1	B	0.91	0/2879	0.85	3/3900 (0.1%)
2	C	0.84	0/969	0.86	0/1323
2	E	1.04	2/978 (0.2%)	0.95	0/1335
3	D	0.83	1/3035 (0.0%)	0.83	0/4134
3	F	1.01	2/3075 (0.1%)	0.92	0/4185
All	All	0.90	5/13781 (0.0%)	0.87	4/18736 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	88[A]	CYS	CB-SG	-5.60	1.72	1.81
2	E	88[B]	CYS	CB-SG	-5.60	1.72	1.81
3	F	338	VAL	CB-CG1	5.49	1.64	1.52
3	F	213	GLU	CG-CD	5.10	1.59	1.51
3	D	79	PHE	CE2-CZ	5.01	1.46	1.37

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	252	ARG	NE-CZ-NH2	-5.74	117.43	120.30
1	A	252	ARG	NE-CZ-NH2	-5.37	117.62	120.30
1	B	321	ARG	NE-CZ-NH1	5.17	122.88	120.30
1	B	269	LEU	CB-CG-CD1	5.10	119.67	111.00

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	2772	0	2655	53	0
1	B	2809	0	2696	42	0
2	C	958	0	862	23	0
2	E	964	0	868	22	0
3	D	2945	0	2834	44	0
3	F	2983	0	2880	44	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	86	0	60	7	0
5	B	86	0	60	6	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	B	8	0	12	2	0
8	D	4	0	3	0	0
9	F	12	0	12	0	0
10	A	144	0	0	4	0
10	B	232	0	0	9	0
10	C	61	0	0	2	0
10	D	188	0	0	6	0
10	E	77	0	0	2	0
10	F	314	0	0	8	0
All	All	14647	0	12942	208	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 8.

All (208) 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)
3:D:289[B]:ARG:HH11	3:D:289[B]:ARG:HG2	0.85	0.99
3:D:289[B]:ARG:CG	3:D:289[B]:ARG:HH11	1.76	0.98
3:D:289[B]:ARG:NH1	3:D:289[B]:ARG:HG2	1.67	0.97
1:A:299:SER:HB2	1:A:333:MET:HG3	1.48	0.95
1:A:48:LYS:H	1:A:62:HIS:HE1	1.22	0.87
3:D:121:PRO:HG3	3:F:104[A]:ARG:NH1	1.91	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:57:TRQ:HB2	2:C:108:TRP:NE1	1.93	0.83
1:B:198:THR:HG22	2:E:58:VAL:HG13	1.61	0.82
3:D:14:GLN:HE21	2:E:21:GLN:HE22	1.30	0.80
1:B:210:GLN:HE22	2:E:44:THR:HG21	1.47	0.79
1:A:210:GLN:HE22	2:C:44:THR:HG21	1.48	0.78
1:B:197:ILE:O	1:B:202:ARG:HD2	1.83	0.78
1:B:39[A]:ARG:HD2	10:B:710:HOH:O	1.82	0.78
1:A:197:ILE:O	1:A:202[B]:ARG:HD2	1.83	0.77
1:B:300[A]:ARG:HG3	3:F:158:PRO:HG2	1.66	0.77
3:F:11[B]:GLN:HB2	10:F:730:HOH:O	1.85	0.75
1:B:200:ASN:O	5:B:403:HEC:HMC3	1.89	0.73
1:A:358:LEU:O	1:A:359:GLU:HB2	1.87	0.73
3:F:342:GLU:HA	10:F:794:HOH:O	1.89	0.72
1:B:198:THR:CG2	2:E:58:VAL:HG13	2.20	0.72
3:F:11[A]:GLN:HA	3:F:11[A]:GLN:NE2	2.05	0.72
1:B:177[B]:ARG:NH2	1:B:355:GLU:OE2	2.23	0.71
2:C:71:LEU:HD13	2:C:130:ALA:HA	1.73	0.71
1:A:202[B]:ARG:HH12	2:C:127:VAL:HG11	1.54	0.71
1:B:208[A]:ARG:NH1	3:D:29:GLY:O	2.20	0.70
2:E:57:TRQ:HB2	2:E:108:TRP:NE1	2.08	0.69
1:A:197:ILE:HD13	2:C:71:LEU:HD23	1.75	0.68
5:B:402:HEC:HBB3	5:B:402:HEC:HMB1	1.75	0.67
1:B:86:LYS:HD2	10:B:661:HOH:O	1.94	0.67
3:D:268:LEU:HD22	3:D:277:TRP:HB3	1.77	0.67
3:F:12:GLU:OE1	3:F:20:ARG:NH1	2.29	0.66
2:E:23:CYS:SG	2:E:88[B]:CYS:HB2	2.35	0.66
1:A:70:LEU:HD13	1:A:163:GLN:NE2	2.12	0.65
1:B:39[A]:ARG:CD	10:B:710:HOH:O	2.43	0.64
1:B:198:THR:HG22	2:E:58:VAL:CG1	2.28	0.64
1:A:206:MET:HE2	1:A:206:MET:HA	1.80	0.63
3:D:272:GLU:HA	10:D:665:HOH:O	1.98	0.63
1:B:215:ARG:HH22	7:B:405:EDO:H22	1.64	0.62
3:D:121:PRO:HG3	3:F:104[A]:ARG:HH12	1.62	0.62
1:A:202[B]:ARG:HH21	2:C:75:ARG:HD2	1.63	0.62
1:B:197:ILE:O	1:B:202:ARG:CD	2.48	0.62
5:B:403:HEC:HBC3	5:B:403:HEC:HMC1	1.82	0.62
1:A:48:LYS:H	1:A:62:HIS:CE1	2.12	0.61
1:B:107:PRO:HG3	10:B:723[A]:HOH:O	2.01	0.60
3:F:222:PRO:HG2	3:F:225:GLU:HB2	1.83	0.60
1:A:202[B]:ARG:NH1	2:C:127:VAL:HG11	2.16	0.60
3:F:82:ASN:HB3	3:F:142:THR:HB	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:53:ALA:HB2	2:C:109:CYS:HA	1.84	0.59
2:C:105:ASP:HB2	3:D:138:TYR:OH	2.03	0.59
1:A:7:ASP:OD1	1:A:7:ASP:N	2.35	0.59
3:D:279:PRO:HB3	3:D:296:LEU:HD21	1.85	0.58
3:F:11[A]:GLN:HE21	3:F:11[A]:GLN:CA	2.16	0.58
1:A:172:ASP:O	1:A:177:ARG:NH1	2.36	0.58
2:E:96:PRO:HB2	2:E:98:TYR:CE1	2.39	0.58
3:D:188:ALA:HB1	3:D:189:PRO:HD2	1.87	0.57
1:A:197:ILE:O	1:A:202[B]:ARG:NH1	2.35	0.56
2:C:101:GLU:HB2	10:C:232:HOH:O	2.05	0.56
1:A:163:GLN:HE22	5:A:402:HEC:HMA1	1.71	0.56
1:A:60:GLN:O	1:A:62:HIS:HD2	1.89	0.56
3:D:289[A]:ARG:NH1	3:D:384:ASP:OD1	2.39	0.55
2:E:101[A]:GLU:HA	10:E:274:HOH:O	2.07	0.55
1:A:202[A]:ARG:HB2	1:A:206:MET:HG3	1.88	0.54
1:B:210:GLN:NE2	2:E:44:THR:HG21	2.20	0.54
5:A:403:HEC:HBC3	5:A:403:HEC:HMC1	1.89	0.54
1:A:119:ARG:HG2	1:A:152:PHE:CG	2.43	0.54
3:F:186:PRO:HB2	3:F:235:GLN:NE2	2.23	0.54
1:B:299:SER:HB3	1:B:304:ALA:CB	2.38	0.54
3:D:82:ASN:HB3	3:D:142:THR:HB	1.90	0.54
1:A:210:GLN:NE2	10:A:605:HOH:O	2.26	0.53
1:B:272:VAL:HG21	5:B:403:HEC:HMA3	1.89	0.53
1:B:81:ARG:NH1	10:B:681:HOH:O	2.38	0.53
3:D:47:VAL:HG13	3:D:64:ILE:HB	1.91	0.53
2:C:58:VAL:HG22	2:C:59:ALA:N	2.24	0.53
1:B:198:THR:CG2	2:E:58:VAL:CG1	2.86	0.53
2:E:31:ILE:CG2	2:E:88[B]:CYS:SG	2.97	0.53
3:F:179:PRO:HD3	3:F:214:ILE:HD13	1.90	0.53
2:C:23:CYS:SG	2:C:88[B]:CYS:HB2	2.48	0.53
1:B:95:GLY:HA3	1:B:223:TYR:OH	2.08	0.53
1:B:91:GLN:O	1:B:92:PHE:HB2	2.09	0.52
3:F:331:HIS:HE1	3:F:366:GLU:OE1	1.92	0.52
1:B:212:VAL:HG11	2:E:129:LYS:HD2	1.92	0.52
1:B:285:THR:HG22	1:B:286:ASP:OD1	2.11	0.51
3:F:181:CYS:C	3:F:182:TYR:CD1	2.84	0.51
3:F:347:TYR:HB3	3:F:356:LEU:HD11	1.93	0.51
1:A:136:PHE:CG	1:A:145:LEU:HD21	2.46	0.51
2:C:110:PHE:CE1	3:D:157:SER:HB2	2.46	0.50
1:A:91:GLN:O	1:A:92:PHE:HB2	2.11	0.50
1:A:65:ARG:HH12	1:A:94:ASP:CG	2.14	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:LEU:HD22	1:A:261:SER:HB2	1.93	0.50
1:B:81:ARG:NH1	10:B:707:HOH:O	2.37	0.50
1:A:113:GLU:HG2	5:A:402:HEC:HBC2	1.92	0.50
3:D:54:HIS:HD2	10:D:677:HOH:O	1.94	0.50
3:D:277:TRP:CE2	3:D:300:GLN:HG3	2.47	0.49
1:A:288:ARG:NH1	1:A:340:ASP:OD1	2.44	0.49
3:D:362:GLU:OE1	3:D:362:GLU:HA	2.12	0.49
3:F:45:ARG:NH2	3:F:343:LYS:O	2.45	0.49
1:A:70:LEU:HD13	1:A:163:GLN:HE21	1.76	0.49
1:A:312:ALA:HB1	1:A:313:PRO:HD2	1.95	0.48
3:F:174:ARG:NH2	3:F:207:GLY:O	2.41	0.48
1:A:21:PRO:O	1:A:27:ALA:HA	2.13	0.48
2:C:57:TRQ:HB2	2:C:108:TRP:HE1	1.73	0.48
3:F:205:ALA:HB3	3:F:213:GLU:HB3	1.95	0.48
3:D:275:ASP:HB2	10:D:665:HOH:O	2.13	0.48
10:B:706:HOH:O	2:E:127:VAL:HG12	2.13	0.48
1:A:40:ALA:HA	1:A:354:TYR:CZ	2.49	0.48
3:D:92:ILE:HG12	3:D:114:PHE:HB2	1.96	0.48
1:A:48:LYS:N	1:A:62:HIS:HE1	2.01	0.48
3:D:51:ASP:HA	3:D:377:PRO:HA	1.96	0.48
1:B:215:ARG:HH22	7:B:405:EDO:C2	2.26	0.47
2:C:51:LYS:HD2	2:C:113:GLU:HB3	1.96	0.47
3:F:42:PRO:HG3	3:F:116:PRO:HB2	1.96	0.47
3:F:115:ASP:O	3:F:119:LEU:HA	2.13	0.47
1:A:277:PRO:HB2	1:A:282:GLY:O	2.15	0.47
1:B:299:SER:HB2	1:B:333:MET:HG3	1.96	0.47
3:F:365[A]:GLU:OE1	10:F:738:HOH:O	2.20	0.47
3:D:269:THR:OG1	3:D:272:GLU:HB2	2.15	0.47
3:F:11[A]:GLN:HA	3:F:11[A]:GLN:HE21	1.75	0.47
1:A:223:TYR:CE2	1:A:265:LYS:HB2	2.49	0.47
2:C:19:ASP:O	2:C:25:TYR:HB2	2.14	0.47
3:F:199:GLY:HA3	3:F:219:VAL:HG13	1.97	0.47
1:B:42:THR:HG21	1:B:275:THR:HB	1.97	0.46
1:A:312:ALA:HB1	1:A:313:PRO:CD	2.45	0.46
1:B:247:HIS:HB3	1:B:250:LEU:HB3	1.98	0.46
3:F:343:LYS:N	3:F:344:PRO:CD	2.78	0.46
2:C:68:GLN:HB3	2:C:129:LYS:HE3	1.98	0.46
3:F:269:THR:OG1	3:F:272:GLU:HG3	2.16	0.46
1:A:46[A]:GLU:HG3	1:A:47:GLY:N	2.30	0.46
1:A:48:LYS:HB2	1:A:62:HIS:CE1	2.51	0.46
3:F:104[B]:ARG:HD3	10:F:530:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:ILE:HG12	5:B:403:HEC:HMB2	1.97	0.46
3:D:315:VAL:O	3:D:324:LEU:N	2.44	0.46
3:D:368:ARG:HA	10:D:625:HOH:O	2.15	0.46
1:A:192:GLY:HA3	1:A:341:ALA:O	2.17	0.45
1:B:163:GLN:HE22	5:B:402:HEC:HMA1	1.81	0.45
3:D:252:ILE:HD12	3:D:252:ILE:N	2.31	0.45
2:C:25:TYR:CE2	2:C:27:ARG:HB2	2.51	0.45
3:D:236:LYS:NZ	3:D:384:ASP:O	2.44	0.45
1:B:223:TYR:CZ	1:B:265:LYS:HB2	2.51	0.45
3:D:140:TRP:CE2	3:D:233:TYR:HB3	2.52	0.45
1:B:223:TYR:CE2	1:B:265:LYS:HB2	2.51	0.45
2:E:101[B]:GLU:HA	10:E:274:HOH:O	2.17	0.45
2:E:91:THR:HB	3:F:306:HIS:CE1	2.51	0.45
1:B:252:ARG:HD3	10:B:636:HOH:O	2.17	0.45
1:B:305:LYS:NZ	10:B:638:HOH:O	2.48	0.45
1:A:353:ARG:NE	10:A:583:HOH:O	2.49	0.44
3:F:349:LEU:HD21	3:F:377:PRO:HB2	1.99	0.44
1:A:68:PRO:HG2	5:A:402:HEC:HBA1	1.99	0.44
2:C:57:TRQ:HB2	2:C:108:TRP:CE2	2.52	0.44
3:D:186:PRO:HB2	3:D:235:GLN:NE2	2.32	0.44
3:D:289[B]:ARG:NH1	3:D:289[B]:ARG:CG	2.49	0.44
1:B:301:ARG:NH2	3:F:177[B]:ASP:OD1	2.50	0.44
1:A:197:ILE:HA	1:A:202[A]:ARG:HB3	2.00	0.44
3:F:280:GLY:HA3	3:F:301:ARG:CZ	2.48	0.44
1:B:186:THR:OG1	1:B:189:GLU:HG3	2.18	0.44
2:E:53:ALA:HB2	2:E:109:CYS:HA	1.99	0.44
3:F:288:HIS:HE1	3:F:290:ALA:HB3	1.83	0.44
1:A:35:HIS:CE1	1:A:70:LEU:HD21	2.53	0.44
1:A:71:GLY:HA2	1:A:168:PHE:O	2.17	0.44
3:F:288:HIS:CE1	3:F:290:ALA:HB3	2.53	0.43
3:F:130:ALA:N	3:F:131:PRO:HD3	2.32	0.43
1:A:29:GLN:HE22	5:A:402:HEC:HBC3	1.83	0.43
2:C:80:TYR:HB2	2:C:120:HIS:HB2	2.00	0.43
3:D:188:ALA:HB1	3:D:189:PRO:CD	2.49	0.43
3:D:314:VAL:HG12	3:D:326:LYS:HG3	2.00	0.43
2:C:96:PRO:HB2	2:C:98:TYR:CE1	2.53	0.43
2:C:131:SER:HA	10:C:255:HOH:O	2.19	0.43
3:D:312:PHE:HA	3:D:327:PHE:O	2.19	0.43
3:F:333:ILE:HD12	3:F:348:ALA:HB1	2.00	0.43
2:E:32:ASP:OD2	2:E:104:ASN:O	2.37	0.43
1:A:40:ALA:HA	1:A:354:TYR:CE1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:43:ASP:HB2	10:D:639:HOH:O	2.17	0.42
2:C:118:THR:HG23	3:D:99:PHE:CE1	2.54	0.42
1:A:188:LEU:HG	1:A:341:ALA:HA	2.01	0.42
1:A:267:PRO:HD3	5:A:403:HEC:CAD	2.48	0.42
3:D:285:VAL:HA	3:D:295:TYR:O	2.19	0.42
3:D:42:PRO:HD3	3:D:117:VAL:HG12	2.01	0.42
3:F:260[A]:LYS:HE3	3:F:262:LEU:CD1	2.49	0.42
1:A:112[A]:VAL:HG22	10:A:613:HOH:O	2.19	0.42
1:A:206:MET:O	1:A:220:ASN:HB3	2.19	0.42
3:D:341:ASP:N	3:D:341:ASP:OD1	2.49	0.42
1:B:20:ASP:HA	1:B:21:PRO:HD2	1.92	0.42
3:F:50:ASN:HB3	10:F:536:HOH:O	2.20	0.42
3:F:251:GLN:O	3:F:262:LEU:HB2	2.20	0.42
1:A:49:ALA:CB	1:A:234:ALA:HA	2.50	0.41
2:E:78:CYS:HA	2:E:118:THR:O	2.20	0.41
3:F:147:ASP:OD1	3:F:147:ASP:C	2.59	0.41
3:F:173[A]:LYS:HE2	10:F:692:HOH:O	2.20	0.41
1:B:81:ARG:HB2	1:B:87:TYR:CE2	2.55	0.41
3:D:289[B]:ARG:HD3	3:D:384:ASP:OD2	2.20	0.41
2:E:19:ASP:O	2:E:25:TYR:HB2	2.20	0.41
3:F:104[B]:ARG:NH1	10:F:520:HOH:O	2.38	0.41
3:F:107:ARG:NE	3:F:130:ALA:HB1	2.35	0.41
3:F:51:ASP:HA	3:F:377:PRO:HA	2.02	0.41
1:A:114:MET:HG3	5:A:402:HEC:HMC2	2.01	0.41
3:D:354:LYS:HE2	3:D:375:HIS:O	2.20	0.41
1:A:107:PRO:HG3	10:A:638[A]:HOH:O	2.20	0.41
3:F:360:ASP:HB2	3:F:367:LEU:HD11	2.01	0.41
1:B:105:GLY:O	1:B:108:MET:HB2	2.21	0.41
3:D:155:GLN:NE2	10:D:666:HOH:O	2.50	0.41
3:D:36:ILE:HD12	2:E:50:THR:O	2.21	0.41
1:A:303:GLU:H	1:A:303:GLU:HG3	1.33	0.41
3:D:292:ASP:OD2	3:D:319:LYS:HD3	2.21	0.41
2:E:59:ALA:O	2:E:71:LEU:HA	2.21	0.41
1:B:70:LEU:HD13	1:B:163:GLN:NE2	2.35	0.40
3:F:88:ASP:HA	10:F:676:HOH:O	2.20	0.40
1:B:296:LYS:O	1:B:305:LYS:HE3	2.21	0.40
1:A:231:ASN:ND2	1:A:283:VAL:HA	2.36	0.40
3:D:91:PHE:CD1	3:D:113:VAL:HG13	2.56	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	357/373 (96%)	345 (97%)	12 (3%)	0	100	100
1	B	360/373 (96%)	351 (98%)	9 (2%)	0	100	100
2	C	123/137 (90%)	119 (97%)	4 (3%)	0	100	100
2	E	124/137 (90%)	116 (94%)	8 (6%)	0	100	100
3	D	378/385 (98%)	358 (95%)	17 (4%)	3 (1%)	19	17
3	F	381/385 (99%)	367 (96%)	13 (3%)	1 (0%)	41	43
All	All	1723/1790 (96%)	1656 (96%)	63 (4%)	4 (0%)	47	52

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	D	207	GLY
3	D	102	ILE
3	D	223	GLU
3	F	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	281/292 (96%)	263 (94%)	18 (6%)	17	18
1	B	284/292 (97%)	271 (95%)	13 (5%)	27	31
2	C	105/112 (94%)	102 (97%)	3 (3%)	42	51

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	106/112 (95%)	105 (99%)	1 (1%)	78	87
3	D	308/310 (99%)	293 (95%)	15 (5%)	25	28
3	F	312/310 (101%)	305 (98%)	7 (2%)	52	62
All	All	1396/1428 (98%)	1339 (96%)	57 (4%)	33	36

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASP
1	A	23	LEU
1	A	75	LEU
1	A	112[A]	VAL
1	A	112[B]	VAL
1	A	190	GLU
1	A	202[A]	ARG
1	A	202[B]	ARG
1	A	209	LYS
1	A	232	GLU
1	A	256	GLU
1	A	300	ARG
1	A	303	GLU
1	A	306	ILE
1	A	321	ARG
1	A	357	LEU
1	A	358	LEU
1	A	359	GLU
1	B	7	ASP
1	B	60	GLN
1	B	102	GLN
1	B	157	GLU
1	B	167	GLU
1	B	208[A]	ARG
1	B	208[B]	ARG
1	B	219	THR
1	B	236	GLU
1	B	269	LEU
1	B	300[A]	ARG
1	B	300[B]	ARG
1	B	357	LEU
2	C	16	GLN
2	C	71	LEU

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Mol	Chain	Res	Type
2	C	131	SER
3	D	45	ARG
3	D	92	ILE
3	D	94	HIS
3	D	106	GLU
3	D	117	VAL
3	D	127	LEU
3	D	208	THR
3	D	211	THR
3	D	218[A]	GLU
3	D	218[B]	GLU
3	D	242	TRP
3	D	262	LEU
3	D	265	VAL
3	D	319	LYS
3	D	354	LYS
2	E	71	LEU
3	F	75	ILE
3	F	94	HIS
3	F	218	GLU
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 (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	60	GLN
1	A	62	HIS
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	84	ASN
1	B	91	GLN
1	B	163	GLN
1	B	210	GLN
3	D	30	GLN

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Mol	Chain	Res	Type
3	D	155	GLN
2	E	21	GLN
3	F	14	GLN
3	F	331	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	TRQ	C	57	2	13,17,18	2.47	4 (30%)	14,24,26	3.64	7 (50%)
2	TRQ	E	57	2	13,17,18	2.36	5 (38%)	14,24,26	2.85	6 (42%)

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	TRQ	C	57	2	-	0/4/19/21	0/2/2/2
2	TRQ	E	57	2	-	0/4/19/21	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	57	TRQ	CH2-CZ2	-6.12	1.47	1.54
2	E	57	TRQ	CH2-CZ2	-5.05	1.48	1.54
2	E	57	TRQ	CD2-CE3	4.35	1.54	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	57	TRQ	CD2-CE3	3.80	1.53	1.44
2	E	57	TRQ	CZ3-CE3	3.14	1.40	1.34
2	C	57	TRQ	CB-CA	3.05	1.60	1.53
2	E	57	TRQ	CB-CA	2.89	1.59	1.53
2	C	57	TRQ	CZ3-CE3	2.56	1.39	1.34
2	E	57	TRQ	CZ3-CH2	-2.02	1.40	1.45

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	57	TRQ	CB-CA-C	7.54	125.61	111.47
2	E	57	TRQ	CB-CG-CD1	-6.04	120.50	127.97
2	C	57	TRQ	CZ2-CE2-NE1	5.79	129.18	119.94
2	E	57	TRQ	CZ2-CE2-NE1	5.69	129.03	119.94
2	C	57	TRQ	O6-CH2-CZ2	-5.53	114.75	118.51
2	C	57	TRQ	CB-CG-CD1	-4.97	121.82	127.97
2	C	57	TRQ	CG-CB-CA	-4.93	106.91	114.53
2	E	57	TRQ	CG-CB-CA	-3.23	109.54	114.53
2	E	57	TRQ	CB-CA-C	3.14	117.35	111.47
2	E	57	TRQ	O7-CZ2-CE2	-2.93	118.73	121.84
2	C	57	TRQ	CE3-CZ3-CH2	-2.23	119.39	121.08
2	C	57	TRQ	CZ3-CH2-CZ2	2.20	121.73	118.72
2	E	57	TRQ	O7-CZ2-CH2	2.13	121.50	119.00

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	57	TRQ	3	0
2	E	57	TRQ	1	0

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
7	EDO	B	405	-	3,3,3	0.29	0	2,2,2	0.30	0
8	ACT	D	401	-	1,3,3	1.47	0	0,3,3	0.00	-
7	EDO	B	406	-	3,3,3	0.65	0	2,2,2	0.16	0
5	HEC	B	402	1,10	26,50,50	1.71	7 (26%)	18,82,82	3.35	12 (66%)
5	HEC	A	402	1,10	26,50,50	1.69	8 (30%)	18,82,82	2.45	8 (44%)
9	MES	F	401	-	12,12,12	1.77	1 (8%)	14,16,16	2.64	5 (35%)
5	HEC	B	403	1	26,50,50	1.53	7 (26%)	18,82,82	3.12	7 (38%)
5	HEC	A	403	1	26,50,50	1.55	6 (23%)	18,82,82	3.31	11 (61%)

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
7	EDO	B	405	-	-	1/1/1/1	-
7	EDO	B	406	-	-	0/1/1/1	-
5	HEC	B	402	1,10	-	0/6/54/54	-
5	HEC	A	402	1,10	-	0/6/54/54	-
9	MES	F	401	-	-	5/6/14/14	0/1/1/1
5	HEC	B	403	1	-	0/6/54/54	-
5	HEC	A	403	1	-	0/6/54/54	-

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	F	401	MES	C8-S	-5.63	1.69	1.77
5	B	402	HEC	C3C-C4C	3.78	1.49	1.43
5	A	402	HEC	C3C-C4C	3.56	1.49	1.43
5	A	403	HEC	C3C-C4C	3.47	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	402	HEC	C4A-C3A	3.23	1.49	1.42
5	B	403	HEC	C3C-C4C	3.22	1.48	1.43
5	A	402	HEC	C1D-CHD	3.09	1.49	1.41
5	B	403	HEC	C4D-CHA	3.04	1.49	1.41
5	A	402	HEC	C1A-C2A	3.02	1.49	1.42
5	A	402	HEC	C4A-C3A	3.01	1.49	1.42
5	A	403	HEC	C4A-C3A	2.85	1.49	1.42
5	B	403	HEC	C1A-C2A	2.85	1.49	1.42
5	A	403	HEC	C4D-CHA	2.75	1.48	1.41
5	A	402	HEC	C4D-CHA	2.67	1.48	1.41
5	B	402	HEC	C1A-C2A	2.60	1.48	1.42
5	A	402	HEC	C3B-C4B	2.59	1.47	1.43
5	B	402	HEC	C3B-C4B	2.58	1.47	1.43
5	A	402	HEC	C1B-CHB	2.52	1.48	1.41
5	A	402	HEC	C1C-CHC	2.51	1.48	1.41
5	B	402	HEC	C1D-CHD	2.45	1.47	1.41
5	A	403	HEC	C3B-C4B	2.37	1.47	1.43
5	B	402	HEC	C4D-CHA	2.27	1.47	1.41
5	B	403	HEC	C3B-C2B	2.26	1.43	1.40
5	B	403	HEC	C1C-CHC	2.25	1.47	1.41
5	B	402	HEC	C1B-CHB	2.22	1.47	1.41
5	B	403	HEC	C4A-C3A	2.22	1.47	1.42
5	A	403	HEC	C1C-CHC	2.16	1.47	1.41
5	A	403	HEC	C1D-CHD	2.07	1.46	1.41
5	B	403	HEC	C4D-ND	-2.03	1.32	1.36

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	403	HEC	CBD-CAD-C3D	-7.63	98.41	112.49
5	B	403	HEC	CMB-C2B-C3B	7.36	134.48	125.82
5	B	402	HEC	CMB-C2B-C3B	6.71	133.71	125.82
5	B	402	HEC	CMC-C2C-C3C	6.30	133.23	125.82
5	A	403	HEC	C1D-C2D-C3D	-6.15	102.71	107.00
5	A	403	HEC	CMB-C2B-C3B	5.87	132.72	125.82
5	B	403	HEC	CBD-CAD-C3D	-5.39	102.54	112.49
9	F	401	MES	C5-N4-C3	5.30	120.75	108.83
5	B	403	HEC	C1D-C2D-C3D	-5.27	103.33	107.00
5	A	402	HEC	CMC-C2C-C3C	5.01	131.71	125.82
9	F	401	MES	O3S-S-C8	4.99	113.84	105.77
5	A	402	HEC	C1D-C2D-C3D	-4.46	103.89	107.00
9	F	401	MES	C7-N4-C5	4.19	121.96	111.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	403	HEC	CBA-CAA-C2A	-4.17	104.79	112.48
5	B	403	HEC	CMC-C2C-C3C	4.04	130.57	125.82
5	A	402	HEC	CBD-CAD-C3D	-3.99	105.12	112.49
5	B	402	HEC	CMB-C2B-C1B	-3.94	122.40	128.46
5	B	402	HEC	CMC-C2C-C1C	-3.86	122.53	128.46
5	B	402	HEC	C4B-C3B-C2B	-3.72	102.33	106.35
5	B	402	HEC	CMD-C2D-C3D	3.67	131.86	124.94
9	F	401	MES	C2-C3-N4	-3.62	104.62	110.10
5	B	402	HEC	CMD-C2D-C1D	-3.61	122.91	128.46
5	B	403	HEC	CMB-C2B-C1B	-3.38	123.28	128.46
5	B	402	HEC	CBD-CAD-C3D	-3.26	106.48	112.49
5	A	403	HEC	CBA-CAA-C2A	-3.17	106.63	112.48
5	A	403	HEC	CMC-C2C-C3C	3.08	129.45	125.82
5	A	403	HEC	CMB-C2B-C1B	-2.94	123.95	128.46
5	B	402	HEC	C3B-C4B-NB	2.93	116.48	110.94
5	A	402	HEC	CBA-CAA-C2A	-2.86	107.21	112.48
5	A	403	HEC	C4C-C3C-C2C	-2.74	103.39	106.35
5	A	402	HEC	C4B-C3B-C2B	-2.68	103.46	106.35
5	B	402	HEC	C4C-C3C-C2C	-2.64	103.50	106.35
5	A	403	HEC	CMD-C2D-C3D	2.62	129.89	124.94
5	B	403	HEC	CAD-CBD-CGD	-2.58	108.34	112.67
5	B	402	HEC	C1D-C2D-C3D	-2.56	105.21	107.00
5	A	402	HEC	CMB-C2B-C3B	2.55	128.82	125.82
5	A	402	HEC	CMC-C2C-C1C	-2.50	124.62	128.46
5	B	402	HEC	CBA-CAA-C2A	-2.41	108.05	112.48
5	A	403	HEC	CMA-C3A-C2A	2.38	129.42	124.94
9	F	401	MES	C7-N4-C3	2.36	117.27	111.23
5	A	403	HEC	CAA-CBA-CGA	2.06	116.13	112.67
5	A	403	HEC	C4B-C3B-C2B	-2.03	104.16	106.35
5	A	402	HEC	CAA-CBA-CGA	-2.02	109.28	112.67

There are no chirality outliers.

All (6) torsion outliers are listed below:

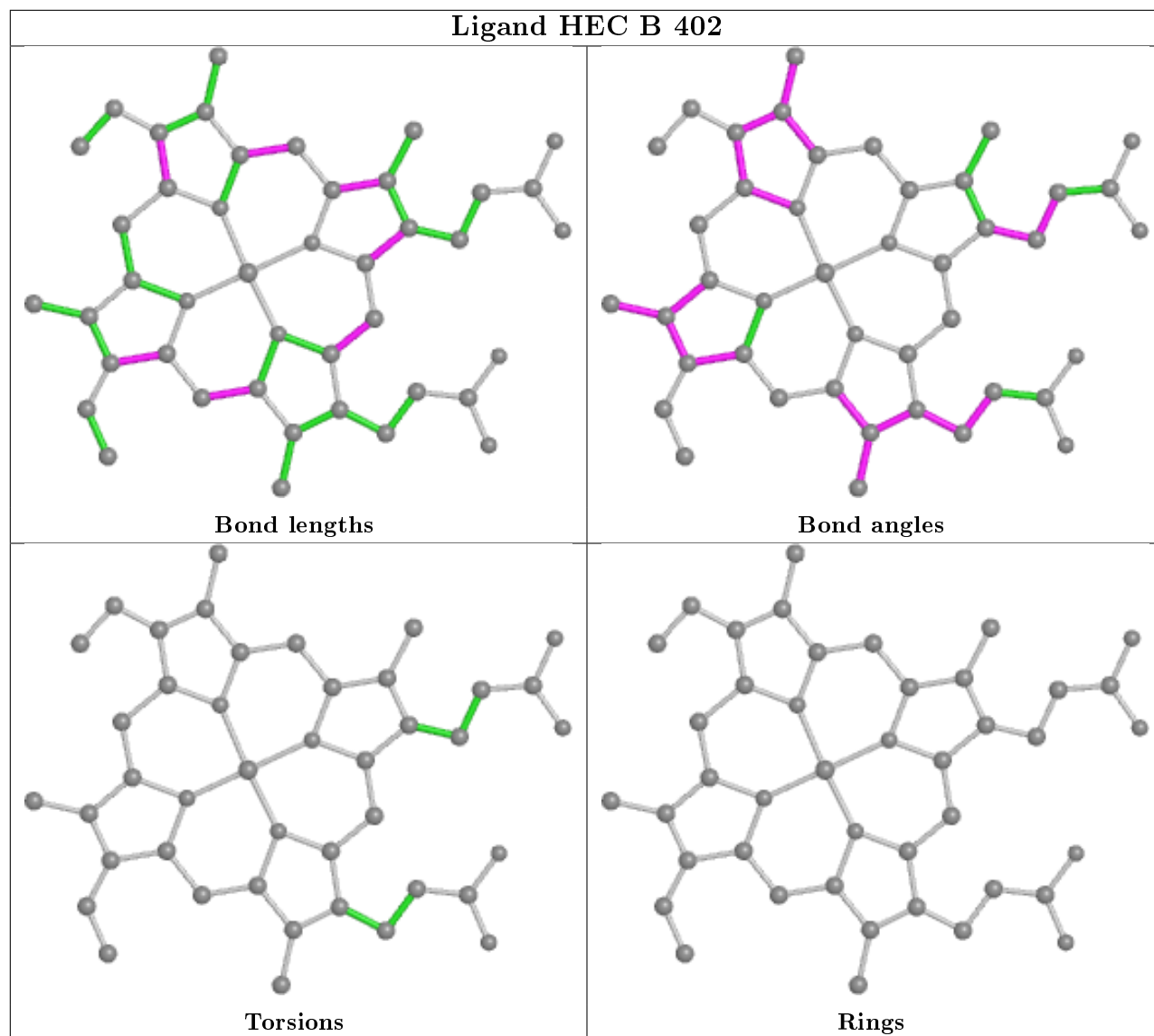
Mol	Chain	Res	Type	Atoms
9	F	401	MES	C8-C7-N4-C5
9	F	401	MES	N4-C7-C8-S
9	F	401	MES	C7-C8-S-O1S
9	F	401	MES	C7-C8-S-O3S
7	B	405	EDO	O1-C1-C2-O2
9	F	401	MES	C7-C8-S-O2S

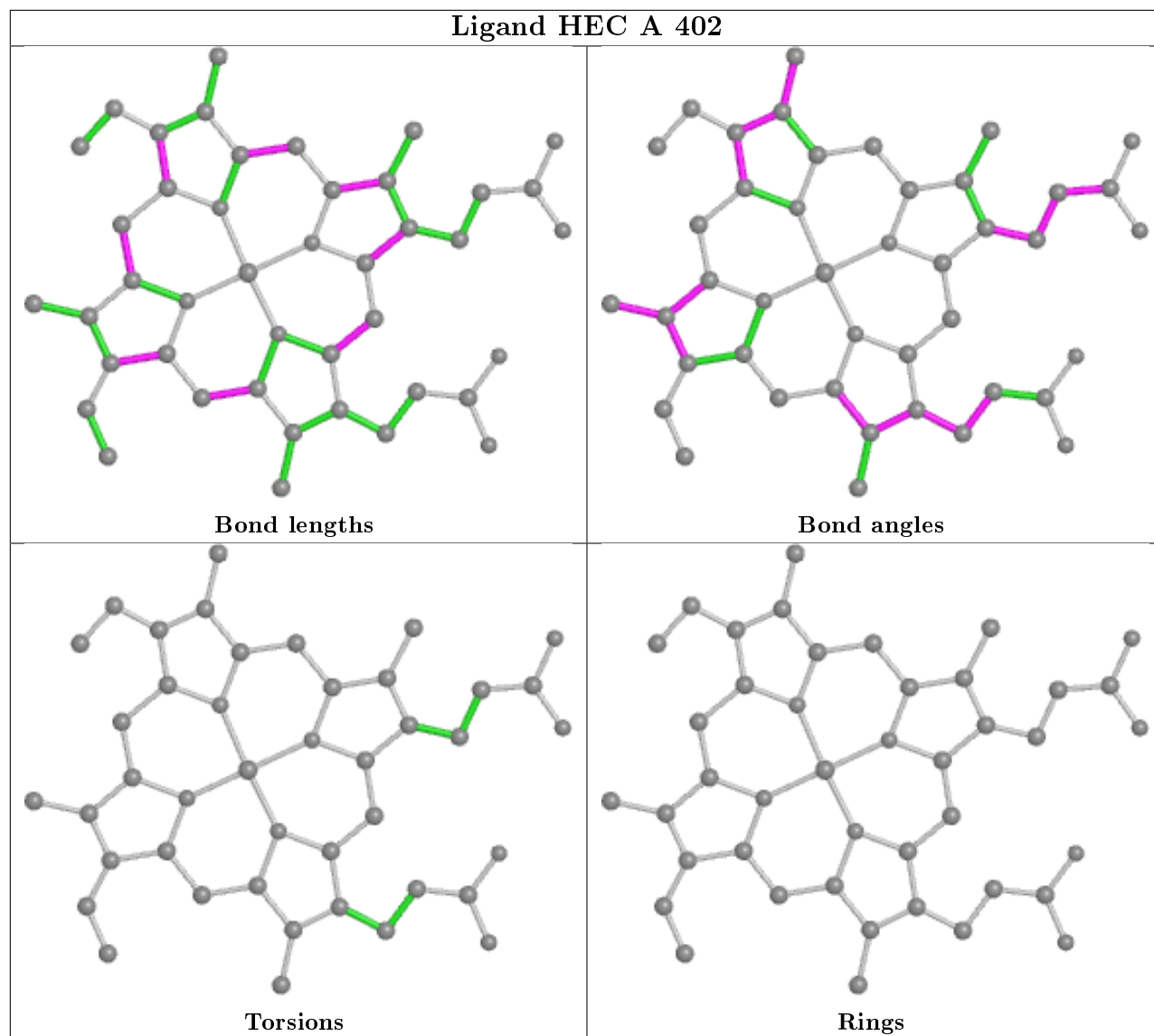
There are no ring outliers.

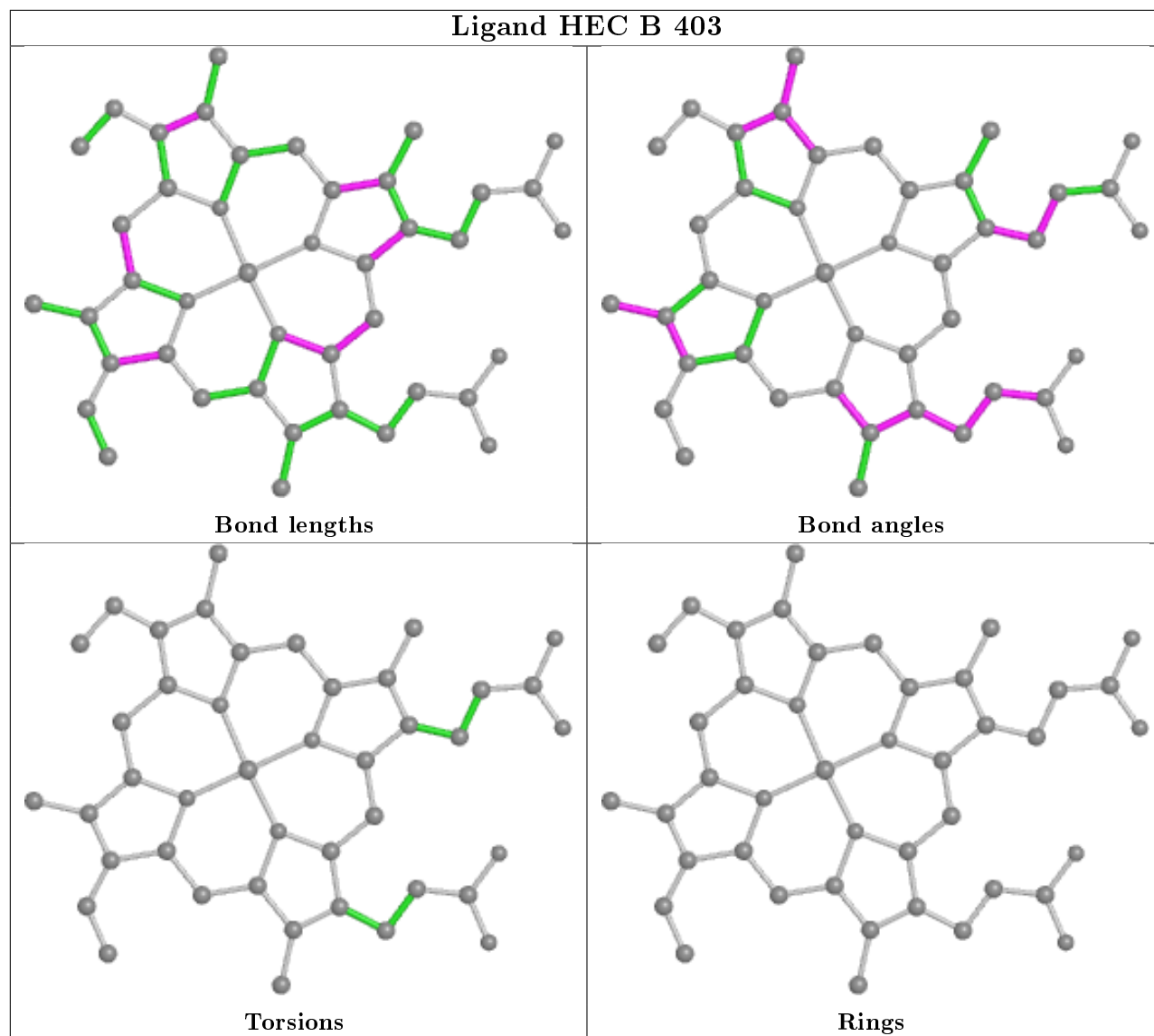
5 monomers are involved in 15 short contacts:

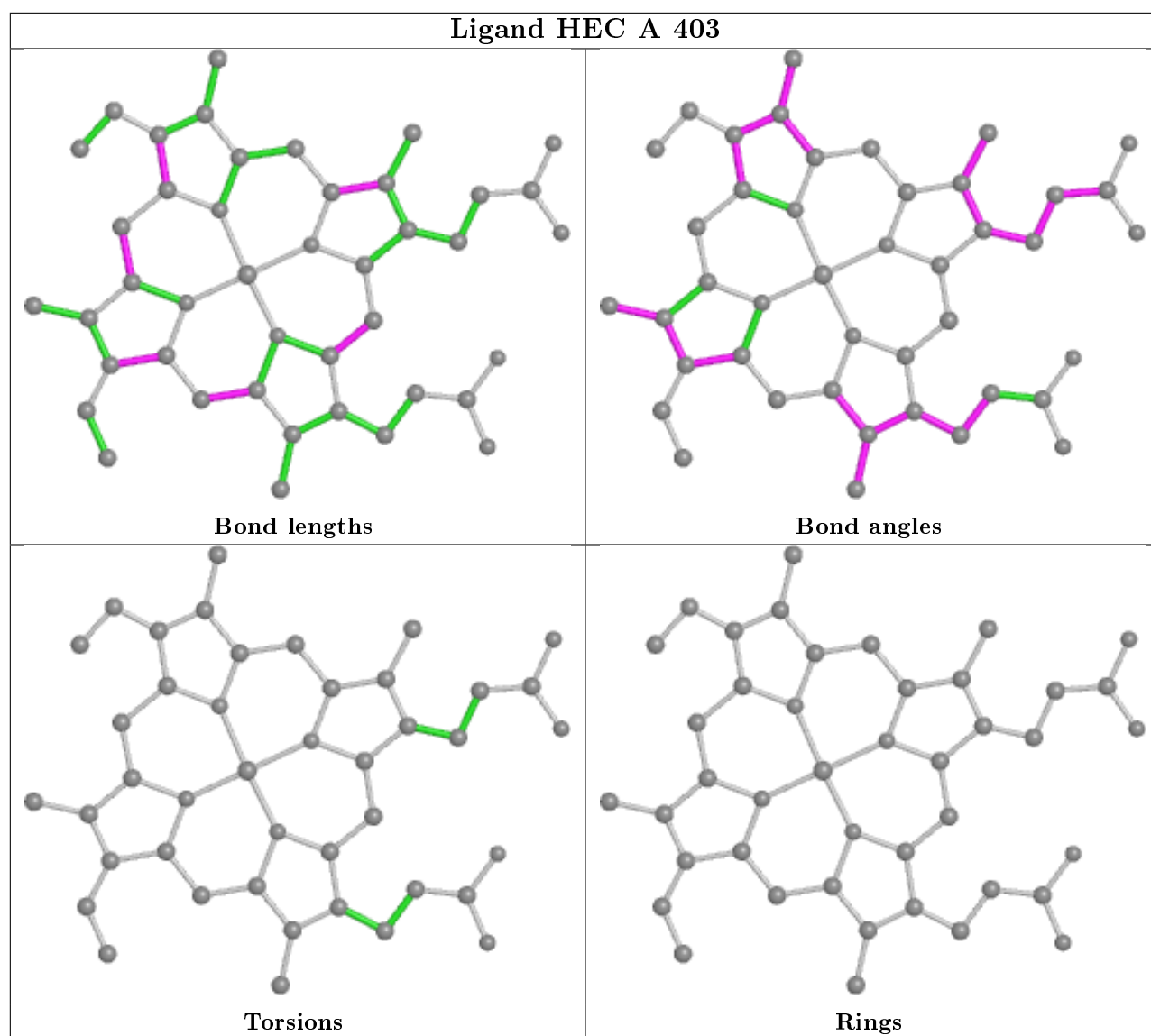
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	405	EDO	2	0
5	B	402	HEC	2	0
5	A	402	HEC	5	0
5	B	403	HEC	4	0
5	A	403	HEC	2	0

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%)	-0.06	5 (1%) 75 75	29, 42, 61, 72	0
1	B	355/373 (95%)	-0.20	3 (0%) 86 86	22, 35, 53, 75	0
2	C	124/137 (90%)	0.20	6 (4%) 30 32	25, 36, 57, 74	0
2	E	124/137 (90%)	-0.08	4 (3%) 47 48	23, 27, 39, 65	0
3	D	376/385 (97%)	0.13	19 (5%) 28 29	24, 43, 71, 83	0
3	F	376/385 (97%)	-0.21	4 (1%) 80 80	20, 30, 47, 61	0
All	All	1709/1790 (95%)	-0.07	41 (2%) 59 59	20, 36, 61, 83	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	6	ALA	6.6
1	A	6	ALA	4.5
2	C	106	ILE	4.3
3	D	207	GLY	4.2
1	B	7	ASP	3.8
3	D	208	THR	3.3
3	D	269	THR	3.2
2	C	108	TRP	3.2
3	F	207	GLY	3.2
1	A	7	ASP	3.2
3	D	133	PHE	3.1
2	C	107	ILE	3.1
2	E	106	ILE	2.8
3	D	135	VAL	2.8
3	D	11	GLN	2.8
3	D	134	LEU	2.7
3	D	270	GLU	2.7
2	E	108	TRP	2.6
2	C	56	SER	2.5

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Mol	Chain	Res	Type	RSRZ
3	D	209	GLU	2.5
2	E	103	ALA	2.4
3	F	27	ALA	2.4
3	D	262	LEU	2.4
1	A	84	ASN	2.4
1	B	84	ASN	2.3
2	E	107	ILE	2.3
3	D	263	PRO	2.2
1	A	243	ASP	2.2
3	D	386	GLY	2.2
3	D	138	TYR	2.2
3	D	137	THR	2.2
3	F	379	VAL	2.2
2	C	7	THR	2.2
3	D	211	THR	2.2
3	D	156	PHE	2.1
2	C	103	ALA	2.1
3	D	136	GLY	2.1
3	D	80	LEU	2.1
3	F	80	LEU	2.1
3	D	282	TRP	2.1
1	A	241	GLY	2.0

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

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	TRQ	C	57	16/17	0.93	0.28	41,48,53,58	0
2	TRQ	E	57	16/17	0.94	0.23	33,38,45,47	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

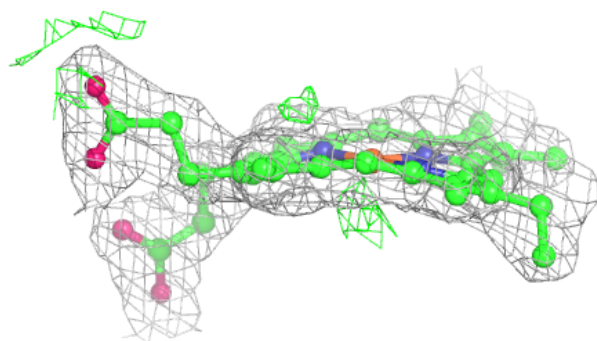
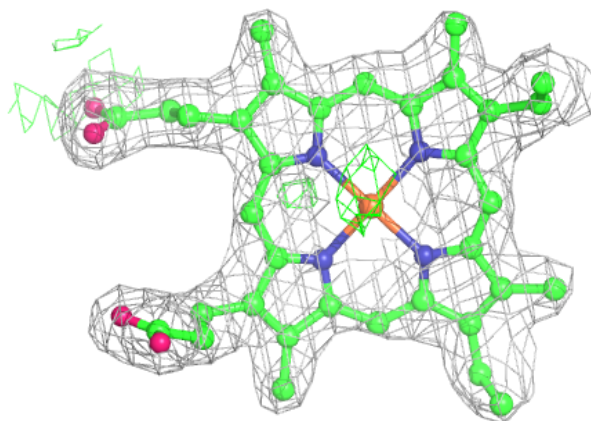
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	EDO	B	406	4/4	0.85	0.12	59,60,60,60	0
6	NA	A	404	1/1	0.91	0.13	58,58,58,58	0
9	MES	F	401	12/12	0.92	0.25	73,78,79,79	0
6	NA	B	404	1/1	0.94	0.06	42,42,42,42	0
8	ACT	D	401	4/4	0.95	0.12	48,50,51,51	0
5	HEC	A	402	43/43	0.97	0.14	27,36,38,39	0
5	HEC	B	402	43/43	0.98	0.13	22,29,33,34	0
4	CA	A	401	1/1	0.98	0.10	40,40,40,40	0
7	EDO	B	405	4/4	0.98	0.09	44,45,46,48	0
5	HEC	A	403	43/43	0.98	0.15	30,34,39,43	0
5	HEC	B	403	43/43	0.99	0.15	15,21,26,31	0
4	CA	B	401	1/1	0.99	0.12	26,26,26,26	0

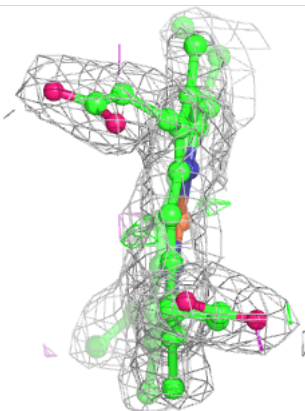
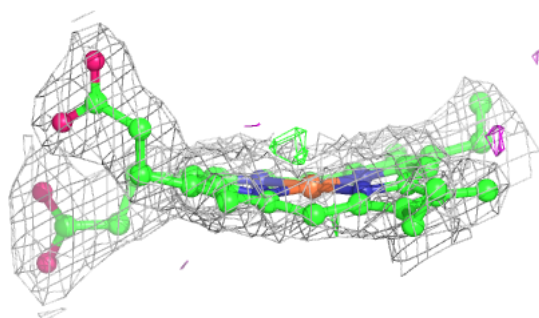
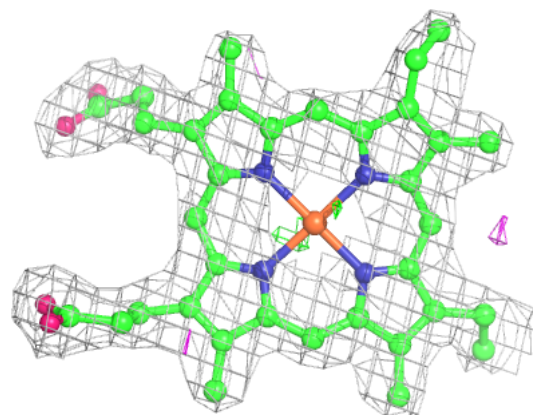
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

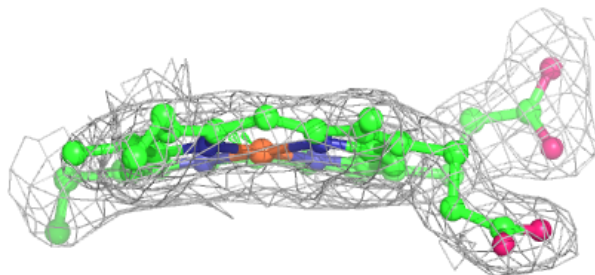
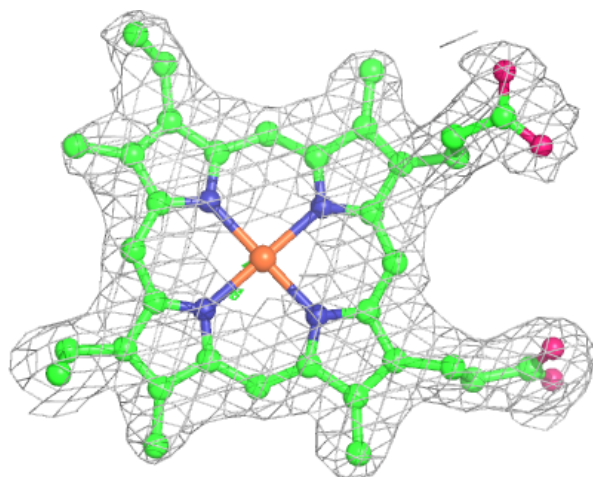
**Electron density around HEC B 402:**

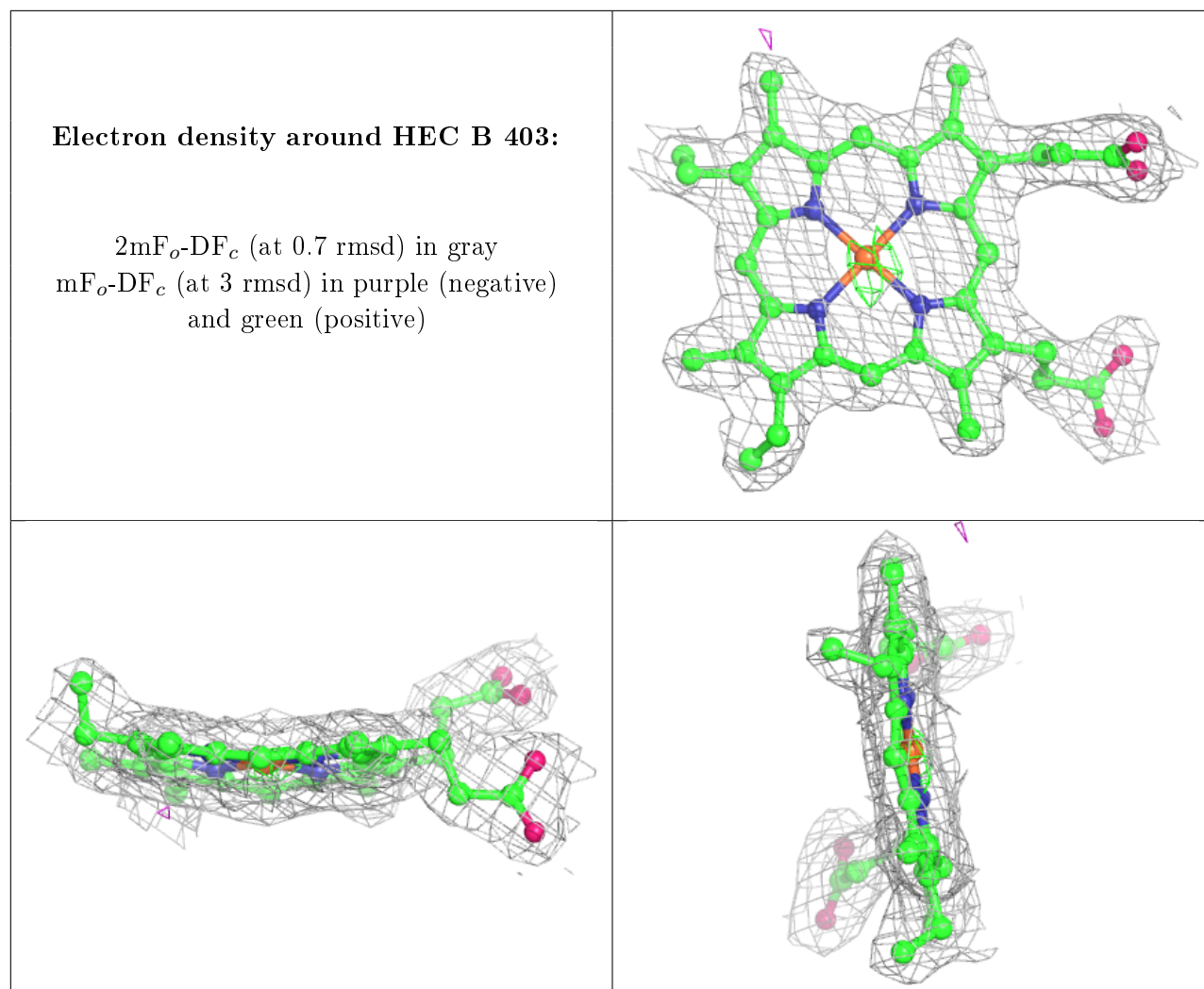
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





6.5 Other polymers ⓘ

There are no such residues in this entry.