



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 11:13 am BST

PDB ID : 1IQC  
Title : Crystal structure of Di-Heme Peroxidase from Nitrosomonas europaea  
Authors : Shimizu, H.  
Deposited on : 2001-07-20  
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

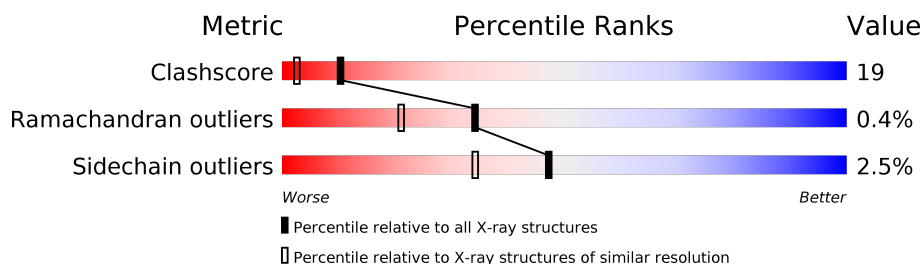
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ .

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	308	
1	B	308	
1	C	308	
1	D	308	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10767 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 di-heme peroxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	308	Total	C	N	O	S	0	0	0
			2388	1497	411	467	13			
1	B	308	Total	C	N	O	S	0	0	0
			2388	1497	411	467	13			
1	C	308	Total	C	N	O	S	0	0	0
			2388	1497	411	467	13			
1	D	308	Total	C	N	O	S	0	0	0
			2388	1497	411	467	13			

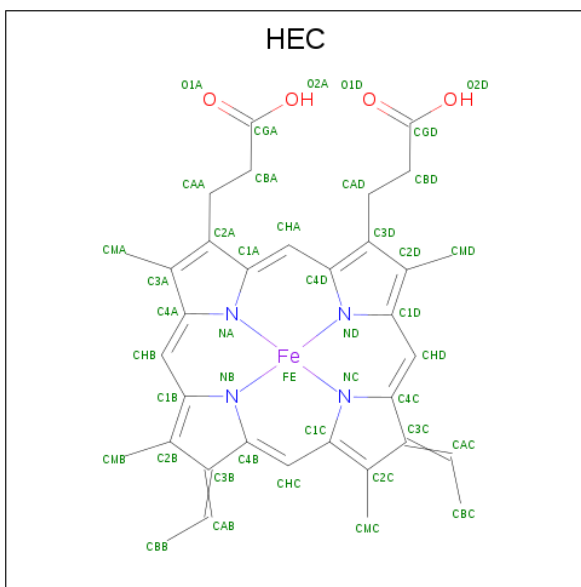
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	2	Total	Ca	0	0
			2	2		
2	D	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		

- Molecule 4 is HEME C (three-letter code: HEC) (formula: C<sub>34</sub>H<sub>34</sub>FeN<sub>4</sub>O<sub>4</sub>).



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

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula:  $\text{C}_3\text{H}_8\text{O}_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

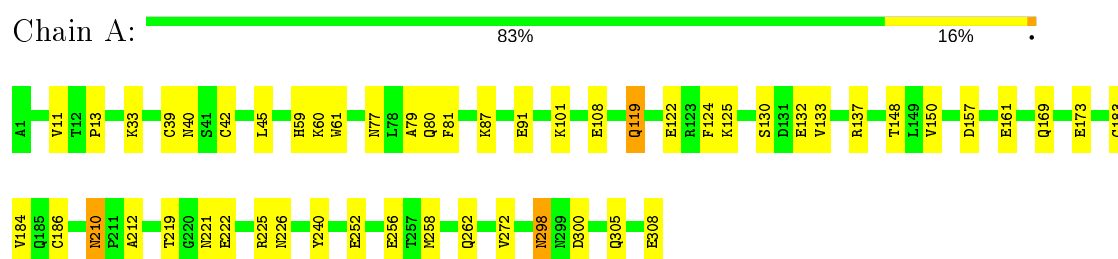
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	260	Total	O	0	0
			260	260		
6	B	199	Total	O	0	0
			199	199		
6	C	279	Total	O	0	0
			279	279		
6	D	120	Total	O	0	0
			120	120		

### 3 Residue-property plots

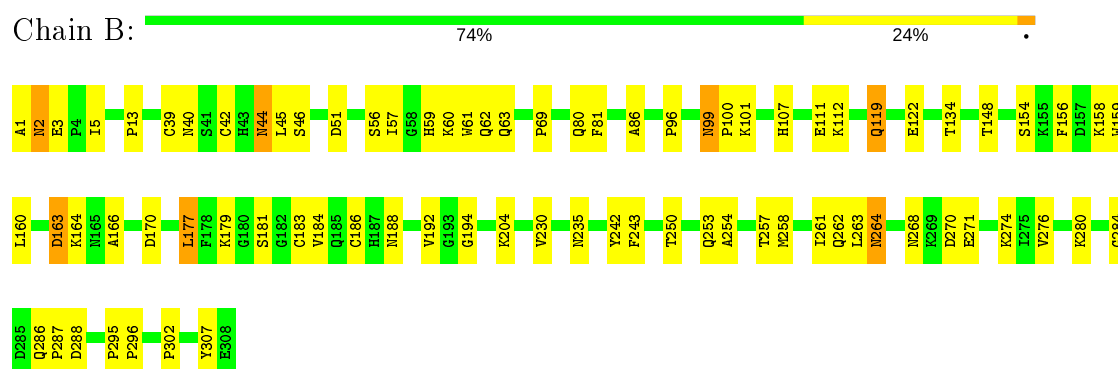
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according 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. 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.

Note EDS was not executed.

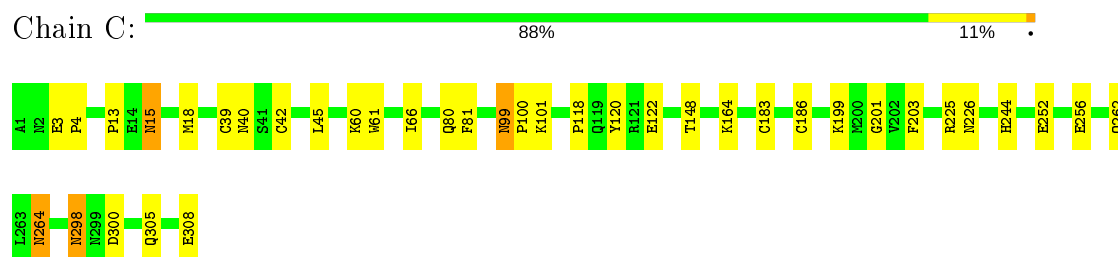
- Molecule 1: di-heme peroxidase



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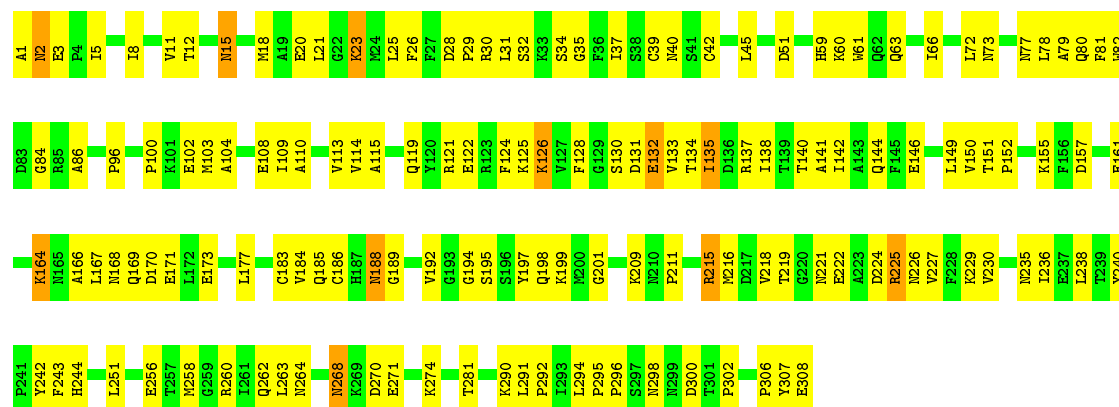


- Molecule 1: di-heme peroxidase



- Molecule 1: di-heme peroxidase





## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.13 Å   55.11 Å   144.00 Å 90.00°   103.60°   90.00°	Depositor
Resolution (Å)	30.00 – 1.80	Depositor
% Data completeness (in resolution range)	95.9 (30.00-1.80)	Depositor
$R_{merge}$	0.03	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS	Depositor
R, $R_{free}$	0.222 , 0.257	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10767	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP



## 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: GOL, CA, MG, 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	0.32	0/2441	0.58	0/3304
1	B	0.29	0/2441	0.56	0/3304
1	C	0.32	0/2441	0.58	0/3304
1	D	0.29	0/2441	0.52	0/3304
All	All	0.30	0/9764	0.56	0/13216

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	240	TYR	Sidechain

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2388	0	2331	53	0
1	B	2388	0	2331	81	0
1	C	2388	0	2331	40	0
1	D	2388	0	2331	179	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	D	1	0	0	0	0
4	A	86	0	64	15	0
4	B	86	0	64	15	0
4	C	86	0	64	14	0
4	D	86	0	64	17	0
5	B	6	0	8	0	0
6	A	260	0	0	4	0
6	B	199	0	0	4	0
6	C	279	0	0	0	0
6	D	120	0	0	6	0
All	All	10767	0	9588	348	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 19.

All (348) 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:186:CYS:SG	4:B:402:HEC:HAC	1.50	1.50
1:D:42:CYS:SG	4:D:401:HEC:HAC	1.70	1.31
1:C:39:CYS:SG	4:C:401:HEC:HAB	1.72	1.25
1:A:183:CYS:SG	4:A:402:HEC:HAB	1.75	1.21
1:A:39:CYS:SG	4:A:401:HEC:HAB	1.77	1.15
1:B:183:CYS:SG	4:B:402:HEC:HAB	1.79	1.15
1:D:183:CYS:SG	4:D:402:HEC:HAB	1.86	1.13
1:B:39:CYS:SG	4:B:401:HEC:HAB	1.86	1.07
1:C:186:CYS:SG	4:C:402:HEC:HAC	1.90	1.06
1:A:42:CYS:SG	4:A:401:HEC:HAC	1.87	1.05
1:D:39:CYS:SG	4:D:401:HEC:HAB	1.91	1.04
1:C:42:CYS:SG	4:C:401:HEC:HAC	1.97	1.03
1:C:183:CYS:SG	4:C:402:HEC:HAB	1.95	1.03
1:A:186:CYS:SG	4:A:402:HEC:HAC	1.98	1.01
1:D:151:THR:HG21	1:D:235:ASN:HD21	1.23	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:125:LYS:HE2	1:D:131:ASP:HB3	1.42	1.00
1:D:73:ASN:HD21	1:D:151:THR:HG22	1.25	0.98
1:D:186:CYS:SG	4:D:402:HEC:HAC	2.02	0.97
1:B:119:GLN:H	1:B:119:GLN:HE21	0.97	0.95
1:A:298:ASN:HD22	1:A:300:ASP:H	1.19	0.90
1:B:42:CYS:SG	4:B:401:HEC:HAC	2.10	0.90
1:D:225:ARG:HH11	1:D:225:ARG:HB2	1.38	0.88
1:B:119:GLN:H	1:B:119:GLN:NE2	1.70	0.88
1:D:224:ASP:O	1:D:227:VAL:HG12	1.74	0.88
1:D:201:GLY:H	1:D:262:GLN:HE22	1.20	0.87
1:D:268:ASN:ND2	1:D:271:GLU:H	1.72	0.87
1:D:201:GLY:H	1:D:262:GLN:NE2	1.72	0.87
1:D:40:ASN:HD21	1:D:45:LEU:H	1.22	0.86
1:A:33:LYS:HD3	6:A:1640:HOH:O	1.74	0.85
1:C:201:GLY:H	1:C:262:GLN:HE22	1.27	0.83
1:B:119:GLN:N	1:B:119:GLN:HE21	1.77	0.82
1:D:122:GLU:O	1:D:126:LYS:HD3	1.79	0.82
1:C:298:ASN:HD22	1:C:300:ASP:H	1.24	0.81
1:D:164:LYS:HA	1:D:164:LYS:HE3	1.61	0.81
1:D:151:THR:HG21	1:D:235:ASN:ND2	1.98	0.79
1:C:42:CYS:SG	4:C:401:HEC:CBC	2.70	0.79
1:C:183:CYS:SG	4:C:402:HEC:CBB	2.71	0.78
1:D:268:ASN:HD21	1:D:271:GLU:H	1.32	0.78
1:D:113:VAL:HG22	1:D:302:PRO:HG2	1.66	0.78
1:B:5:ILE:HD12	1:B:192:VAL:HG22	1.67	0.77
1:C:39:CYS:SG	4:C:401:HEC:CBB	2.71	0.77
1:B:39:CYS:SG	4:B:401:HEC:CBB	2.73	0.77
1:D:31:LEU:HB2	6:D:2502:HOH:O	1.84	0.76
1:A:210:ASN:HD22	1:A:212:ALA:H	1.33	0.76
1:D:40:ASN:ND2	1:D:45:LEU:H	1.83	0.75
1:B:204:LYS:HB2	1:B:261:ILE:HG22	1.69	0.75
1:B:44:ASN:HD22	1:B:46:SER:H	1.33	0.75
1:D:73:ASN:ND2	1:D:151:THR:HG22	2.01	0.74
1:D:298:ASN:HD22	1:D:300:ASP:H	1.35	0.74
1:B:42:CYS:SG	4:B:401:HEC:C3C	2.75	0.74
1:B:42:CYS:SG	4:B:401:HEC:CBC	2.75	0.74
1:A:119:GLN:NE2	1:A:119:GLN:H	1.84	0.74
1:D:66:ILE:HD11	1:D:244:HIS:HB2	1.69	0.73
1:A:186:CYS:SG	4:A:402:HEC:CBC	2.75	0.73
1:D:130:SER:HB2	1:D:137:ARG:HH12	1.54	0.73
1:D:103:MET:HE1	6:D:2502:HOH:O	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:SER:OG	1:D:291:LEU:HD21	1.89	0.73
1:A:42:CYS:SG	4:A:401:HEC:C3C	2.76	0.72
1:D:125:LYS:CE	1:D:131:ASP:HB3	2.19	0.72
1:A:252:GLU:O	1:A:256:GLU:HG3	1.90	0.72
1:B:13:PRO:HG3	1:B:148:THR:CG2	2.19	0.71
1:D:3:GLU:O	1:D:188:ASN:HB2	1.89	0.71
1:D:73:ASN:HD21	1:D:151:THR:CG2	2.00	0.71
1:D:263:LEU:HD21	4:D:402:HEC:HMC2	1.71	0.71
1:B:44:ASN:ND2	1:B:46:SER:H	1.88	0.71
1:D:130:SER:HB2	1:D:137:ARG:HH22	1.54	0.70
1:D:186:CYS:SG	4:D:402:HEC:C3C	2.79	0.70
1:D:221:ASN:HB2	1:D:224:ASP:OD2	1.91	0.70
1:D:216:MET:SD	1:D:225:ARG:NH1	2.65	0.70
1:D:42:CYS:SG	4:D:401:HEC:C3C	2.80	0.70
1:A:308:GLU:OXT	1:A:308:GLU:HG3	1.92	0.70
1:B:186:CYS:SG	4:B:402:HEC:C3C	2.80	0.69
1:A:39:CYS:SG	4:A:401:HEC:CBB	2.79	0.69
1:D:215:ARG:HB3	1:D:227:VAL:HG13	1.73	0.69
1:A:308:GLU:OXT	1:A:308:GLU:CG	2.40	0.69
1:C:186:CYS:SG	4:C:402:HEC:C3C	2.81	0.69
1:D:169:GLN:HG3	1:D:170:ASP:N	2.08	0.69
1:B:170:ASP:OD1	1:B:274:LYS:HE3	1.93	0.69
1:D:183:CYS:SG	4:D:402:HEC:CBB	2.81	0.68
1:C:80:GLN:O	1:C:81:PHE:HB2	1.93	0.68
1:D:108:GLU:OE1	1:D:109:ILE:HG13	1.94	0.68
1:D:298:ASN:HD21	1:D:300:ASP:HB2	1.59	0.67
1:D:3:GLU:HG3	1:D:189:GLY:HA2	1.77	0.67
1:D:114:VAL:HG21	1:D:133:VAL:HG11	1.77	0.67
1:A:184:VAL:HG22	6:A:1464:HOH:O	1.95	0.67
1:A:186:CYS:SG	4:A:402:HEC:C3C	2.81	0.67
1:A:169:GLN:O	1:A:173:GLU:HG3	1.95	0.67
1:B:13:PRO:HG3	1:B:148:THR:HG21	1.77	0.67
1:C:201:GLY:H	1:C:262:GLN:NE2	1.93	0.66
1:D:215:ARG:HB3	1:D:227:VAL:CG1	2.26	0.66
1:A:80:GLN:O	1:A:81:PHE:HB2	1.96	0.66
1:C:42:CYS:SG	4:C:401:HEC:C3C	2.81	0.66
1:B:96:PRO:HG3	6:B:1609:HOH:O	1.96	0.66
1:D:18:MET:HB3	1:D:144:GLN:HG2	1.77	0.66
1:D:110:ALA:O	1:D:114:VAL:HG22	1.96	0.65
1:D:1:ALA:O	1:D:2:ASN:HB2	1.96	0.65
1:D:20:GLU:O	1:D:23:LYS:HG2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:186:CYS:SG	4:D:402:HEC:CBC	2.86	0.64
1:C:186:CYS:SG	4:C:402:HEC:CBC	2.84	0.64
1:D:298:ASN:ND2	1:D:300:ASP:HB2	2.13	0.64
4:A:402:HEC:HBC3	4:A:402:HEC:HMC1	1.80	0.64
1:C:298:ASN:ND2	1:C:300:ASP:H	1.96	0.64
1:C:183:CYS:SG	4:C:402:HEC:C3B	2.85	0.64
1:D:225:ARG:NH1	1:D:225:ARG:HB2	2.12	0.63
1:A:125:LYS:HA	1:A:130:SER:O	1.98	0.63
1:D:268:ASN:HD22	1:D:268:ASN:C	2.01	0.63
1:A:298:ASN:ND2	1:A:300:ASP:H	1.94	0.63
1:B:286:GLN:HB3	1:B:287:PRO:HD2	1.80	0.63
1:B:270:ASP:O	1:B:274:LYS:HD3	1.99	0.62
1:C:252:GLU:O	1:C:256:GLU:HG3	1.99	0.62
1:D:39:CYS:SG	4:D:401:HEC:CBB	2.87	0.62
1:B:80:GLN:O	1:B:81:PHE:HB2	1.99	0.62
1:B:1:ALA:O	1:B:2:ASN:HB2	1.99	0.62
1:D:188:ASN:C	1:D:188:ASN:HD22	2.03	0.62
1:D:66:ILE:HG13	1:D:244:HIS:O	2.00	0.62
1:B:268:ASN:HB2	1:B:271:GLU:HG3	1.82	0.62
1:A:210:ASN:ND2	1:A:212:ALA:H	1.96	0.61
1:D:125:LYS:HA	1:D:130:SER:O	2.00	0.61
1:A:42:CYS:SG	4:A:401:HEC:CBC	2.87	0.61
1:D:80:GLN:O	1:D:81:PHE:HB2	1.99	0.61
1:A:130:SER:HB3	1:A:137:ARG:HH12	1.65	0.60
1:A:183:CYS:SG	4:A:402:HEC:CBB	2.86	0.60
1:B:183:CYS:SG	4:B:402:HEC:CBB	2.88	0.60
1:C:15:ASN:HD22	1:C:15:ASN:C	2.05	0.60
1:D:108:GLU:CD	1:D:109:ILE:HG13	2.22	0.60
1:D:135:ILE:HD13	1:D:135:ILE:O	2.01	0.60
1:D:96:PRO:HG3	6:D:2540:HOH:O	2.02	0.60
1:C:13:PRO:HG3	1:C:148:THR:HG22	1.83	0.59
1:D:3:GLU:HG3	1:D:189:GLY:CA	2.33	0.59
1:B:39:CYS:SG	4:B:401:HEC:C3B	2.89	0.59
1:B:107:HIS:O	1:B:111:GLU:HG3	2.03	0.58
1:D:155:LYS:O	1:D:166:ALA:HB1	2.04	0.58
1:D:29:PRO:O	1:D:35:GLY:HA2	2.04	0.58
1:D:78:LEU:C	1:D:78:LEU:HD12	2.23	0.58
1:A:298:ASN:HD21	1:A:300:ASP:HB2	1.68	0.58
1:A:119:GLN:HE21	1:A:119:GLN:H	1.47	0.58
1:D:130:SER:HB2	1:D:137:ARG:NH2	2.18	0.58
1:D:225:ARG:HH11	1:D:225:ARG:CB	2.14	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:298:ASN:HD21	1:C:300:ASP:HB2	1.68	0.58
1:D:130:SER:HB2	1:D:137:ARG:NH1	2.17	0.58
1:D:39:CYS:SG	4:D:401:HEC:C3B	2.89	0.58
1:D:2:ASN:ND2	1:D:3:GLU:H	2.03	0.57
1:B:163:ASP:OD2	1:B:166:ALA:HB2	2.05	0.57
1:B:61:TRP:HB2	1:D:61:TRP:HB2	1.87	0.57
1:A:157:ASP:O	1:A:161:GLU:HG3	2.05	0.56
1:D:198:GLN:C	1:D:227:VAL:HG23	2.25	0.56
1:C:118:PRO:O	1:C:122:GLU:HG3	2.05	0.56
1:B:288:ASP:HA	6:B:1514:HOH:O	2.06	0.56
1:C:13:PRO:HG3	1:C:148:THR:CG2	2.36	0.56
1:B:13:PRO:HG3	1:B:148:THR:HG22	1.86	0.56
1:B:44:ASN:HD22	1:B:44:ASN:C	2.09	0.56
1:D:15:ASN:OD1	1:D:18:MET:HB2	2.06	0.55
4:D:402:HEC:HMC1	4:D:402:HEC:HBC3	1.88	0.55
1:D:124:PHE:CE2	1:D:133:VAL:HG13	2.41	0.55
1:A:13:PRO:HG3	1:A:148:THR:CG2	2.36	0.55
1:B:1:ALA:O	1:B:2:ASN:CB	2.54	0.55
1:D:21:LEU:O	1:D:25:LEU:HD23	2.06	0.55
1:D:5:ILE:H	1:D:5:ILE:HD12	1.72	0.55
4:B:401:HEC:HMC1	4:B:401:HEC:HBC3	1.88	0.55
1:D:151:THR:O	1:D:151:THR:HG23	2.06	0.55
1:A:305:GLN:HB3	1:A:308:GLU:HG2	1.89	0.54
1:B:264:ASN:HD22	1:B:264:ASN:C	2.10	0.54
1:A:183:CYS:SG	4:A:402:HEC:C3B	2.89	0.54
1:B:258:MET:O	1:B:262:GLN:HB2	2.07	0.54
1:B:242:TYR:O	1:B:243:PHE:HB2	2.08	0.54
1:D:130:SER:CB	1:D:137:ARG:HH12	2.19	0.54
1:D:28:ASP:OD1	1:D:30:ARG:HB2	2.08	0.54
1:C:39:CYS:SG	4:C:401:HEC:C3B	2.91	0.54
1:D:215:ARG:CB	1:D:227:VAL:HG13	2.37	0.54
1:D:298:ASN:ND2	1:D:300:ASP:H	2.02	0.54
4:C:401:HEC:HMC1	4:C:401:HEC:HBC3	1.90	0.54
1:D:5:ILE:N	1:D:5:ILE:HD12	2.22	0.53
1:B:134:THR:HG22	6:B:1575:HOH:O	2.09	0.53
1:B:235:ASN:HD21	1:B:284:GLY:H	1.54	0.53
1:D:209:LYS:O	1:D:211:PRO:HD3	2.09	0.53
1:D:268:ASN:HD21	1:D:271:GLU:HG3	1.73	0.53
1:D:5:ILE:HA	1:D:189:GLY:O	2.09	0.53
1:D:114:VAL:CG2	1:D:133:VAL:HG11	2.38	0.53
1:D:42:CYS:HA	1:D:51:ASP:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:PRO:HG3	1:A:148:THR:HG21	1.91	0.53
1:D:135:ILE:HA	1:D:138:ILE:HD12	1.90	0.53
1:B:194:GLY:HA2	1:B:230:VAL:O	2.08	0.53
1:C:99:ASN:ND2	1:C:101:LYS:H	2.07	0.53
1:D:134:THR:O	1:D:138:ILE:HG13	2.10	0.52
1:D:37:ILE:N	1:D:37:ILE:HD12	2.23	0.52
1:D:82:TRP:HZ3	4:D:402:HEC:HBD2	1.75	0.52
1:D:268:ASN:C	1:D:268:ASN:ND2	2.62	0.52
1:A:39:CYS:SG	4:A:401:HEC:C3B	2.91	0.52
1:D:194:GLY:N	1:D:230:VAL:O	2.43	0.52
1:B:183:CYS:SG	4:B:402:HEC:C3B	2.94	0.52
1:C:40:ASN:HD21	1:C:45:LEU:H	1.57	0.52
1:D:132:GLU:HA	1:D:132:GLU:OE1	2.08	0.51
1:D:169:GLN:O	1:D:173:GLU:HG3	2.10	0.51
1:D:78:LEU:HD11	1:D:197:TYR:OH	2.10	0.51
1:A:130:SER:OG	1:A:132:GLU:HG2	2.10	0.51
1:C:99:ASN:C	1:C:99:ASN:HD22	2.13	0.51
1:D:306:PRO:HG2	1:D:307:TYR:CD1	2.45	0.51
1:A:60:LYS:HA	1:C:61:TRP:CD2	2.45	0.51
1:B:5:ILE:HD11	1:B:188:ASN:HA	1.93	0.51
1:D:229:LYS:NZ	4:D:401:HEC:O1A	2.43	0.51
1:D:291:LEU:HD23	1:D:291:LEU:C	2.31	0.51
1:D:168:ASN:OD1	1:D:171:GLU:HG3	2.11	0.51
1:D:177:LEU:HD21	1:D:271:GLU:HB3	1.92	0.51
1:D:258:MET:O	1:D:262:GLN:HB2	2.10	0.51
1:D:268:ASN:HD21	1:D:271:GLU:N	2.03	0.50
1:D:270:ASP:OD2	1:D:274:LYS:HE2	2.10	0.50
1:D:242:TYR:O	1:D:243:PHE:HB2	2.12	0.50
1:A:40:ASN:HD21	1:A:45:LEU:H	1.60	0.50
1:D:8:ILE:O	1:D:152:PRO:HB3	2.11	0.50
1:B:40:ASN:ND2	1:B:45:LEU:H	2.10	0.50
1:B:44:ASN:HD21	1:B:46:SER:HB2	1.76	0.50
1:D:80:GLN:HB2	1:D:86:ALA:HB3	1.94	0.50
1:B:56:SER:HB3	4:B:401:HEC:HAC	1.94	0.50
1:A:61:TRP:HB2	1:C:61:TRP:HB2	1.93	0.49
1:D:201:GLY:N	1:D:262:GLN:HE22	2.00	0.49
1:B:60:LYS:HA	1:D:61:TRP:CD2	2.47	0.49
1:B:5:ILE:CD1	1:B:188:ASN:HA	2.41	0.49
1:D:2:ASN:CG	1:D:3:GLU:H	2.15	0.49
1:B:59:HIS:O	1:B:62:GLN:HG3	2.12	0.49
1:D:268:ASN:ND2	1:D:271:GLU:HB2	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:77:ASN:HB3	1:D:79:ALA:O	2.12	0.49
1:A:256:GLU:HG2	1:A:272:VAL:HG21	1.95	0.49
1:C:40:ASN:ND2	1:C:45:LEU:H	2.11	0.49
1:D:135:ILE:HD13	1:D:135:ILE:C	2.33	0.49
1:A:61:TRP:CD2	1:C:60:LYS:HA	2.48	0.48
1:D:215:ARG:O	1:D:218:VAL:HG12	2.13	0.48
1:D:34:SER:HB3	1:D:59:HIS:CE1	2.47	0.48
1:D:183:CYS:CB	4:D:402:HEC:HAB	2.42	0.48
1:D:218:VAL:HG13	1:D:219:THR:HG23	1.95	0.48
1:D:222:GLU:OE2	1:D:225:ARG:NH1	2.46	0.48
1:B:112:LYS:HD3	1:B:302:PRO:HG3	1.94	0.48
1:D:78:LEU:HD11	1:D:197:TYR:CZ	2.48	0.48
1:B:40:ASN:HD21	1:B:45:LEU:H	1.62	0.48
1:D:96:PRO:HA	1:D:102:GLU:OE1	2.14	0.48
1:D:291:LEU:HD23	1:D:292:PRO:N	2.28	0.47
1:B:57:ILE:HD11	1:D:306:PRO:HB2	1.97	0.47
1:C:15:ASN:ND2	1:C:18:MET:H	2.13	0.47
1:A:101:LYS:HD3	6:A:1524:HOH:O	2.12	0.47
1:B:177:LEU:HD23	1:B:181:SER:HG	1.79	0.47
1:D:155:LYS:HD3	1:D:166:ALA:O	2.14	0.47
1:B:57:ILE:HD13	1:D:307:TYR:CZ	2.49	0.47
1:C:66:ILE:HD11	1:C:244:HIS:HB2	1.97	0.47
1:C:308:GLU:HG2	1:C:308:GLU:OXT	2.15	0.47
1:D:1:ALA:HB2	1:D:184:VAL:HG23	1.97	0.47
1:D:78:LEU:HD23	1:D:195:SER:C	2.34	0.47
1:D:80:GLN:HA	1:D:80:GLN:NE2	2.30	0.47
1:D:183:CYS:SG	4:D:402:HEC:C3B	2.97	0.47
1:B:296:PRO:HB3	1:D:238:LEU:HD22	1.97	0.46
1:D:113:VAL:HG22	1:D:302:PRO:CG	2.41	0.46
1:D:167:LEU:HD22	1:D:281:THR:HG21	1.98	0.46
1:C:203:PHE:CD2	1:C:264:ASN:HB2	2.51	0.46
1:D:80:GLN:O	1:D:84:GLY:HA2	2.15	0.46
1:B:99:ASN:HD22	1:B:100:PRO:N	2.14	0.46
1:B:99:ASN:HD22	1:B:100:PRO:CD	2.27	0.46
1:B:181:SER:O	1:B:263:LEU:HD13	2.16	0.46
1:D:128:PHE:HE2	1:D:140:THR:HG21	1.81	0.46
1:D:134:THR:OG1	1:D:137:ARG:HG3	2.16	0.46
1:D:188:ASN:C	1:D:188:ASN:ND2	2.69	0.46
1:D:225:ARG:O	1:D:226:ASN:HB2	2.16	0.46
1:D:240:TYR:HB2	6:D:2535:HOH:O	2.15	0.46
1:B:288:ASP:OD1	1:D:290:LYS:HD3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:ALA:HB1	4:B:402:HEC:HMA2	1.98	0.46
1:D:256:GLU:CD	1:D:260:ARG:HH12	2.19	0.46
1:C:99:ASN:HD22	1:C:100:PRO:N	2.13	0.45
1:D:100:PRO:HA	1:D:104:ALA:HA	1.98	0.45
1:D:146:GLU:HA	1:D:149:LEU:HG	1.99	0.45
1:B:177:LEU:HD11	1:B:271:GLU:HB3	1.98	0.45
1:D:23:LYS:NZ	1:D:23:LYS:HB2	2.32	0.45
1:B:63:GLN:HG3	1:D:307:TYR:CZ	2.51	0.45
1:A:40:ASN:ND2	1:A:45:LEU:H	2.14	0.45
1:B:42:CYS:HA	1:B:51:ASP:HB3	1.99	0.45
1:D:78:LEU:HD12	1:D:79:ALA:HB2	1.99	0.45
1:D:150:VAL:HG23	1:D:152:PRO:HD3	1.99	0.45
1:A:59:HIS:CD2	1:A:60:LYS:HG3	2.52	0.45
1:D:142:ILE:O	1:D:146:GLU:HG3	2.16	0.45
1:C:225:ARG:O	1:C:226:ASN:HB2	2.18	0.44
1:D:268:ASN:ND2	1:D:271:GLU:CB	2.80	0.44
1:D:72:LEU:O	1:D:73:ASN:HB2	2.18	0.44
1:A:108:GLU:CD	1:A:108:GLU:H	2.20	0.44
1:A:225:ARG:O	1:A:226:ASN:HB2	2.18	0.44
1:B:99:ASN:HD22	1:B:100:PRO:HD2	1.83	0.44
1:B:99:ASN:C	1:B:99:ASN:HD22	2.19	0.44
1:B:179:LYS:HG2	1:B:184:VAL:HG11	2.00	0.44
1:D:260:ARG:O	1:D:264:ASN:HA	2.18	0.44
1:D:270:ASP:O	1:D:274:LYS:HG3	2.18	0.44
1:C:305:GLN:HB3	1:C:308:GLU:HB3	1.99	0.44
1:B:257:THR:O	1:B:261:ILE:HG12	2.18	0.44
1:C:199:LYS:HE3	1:C:226:ASN:HB2	2.00	0.44
1:D:183:CYS:C	1:D:185:GLN:H	2.21	0.44
1:B:307:TYR:CZ	1:D:63:GLN:HG3	2.53	0.43
1:D:194:GLY:HA2	1:D:230:VAL:O	2.18	0.43
1:B:235:ASN:ND2	1:B:284:GLY:H	2.15	0.43
4:C:402:HEC:HBC3	4:C:402:HEC:HMC1	2.00	0.43
1:B:112:LYS:HE2	6:B:1474:HOH:O	2.18	0.43
1:B:177:LEU:HD23	1:B:181:SER:OG	2.18	0.43
1:D:26:PHE:O	1:D:39:CYS:HB2	2.18	0.43
1:B:99:ASN:ND2	1:B:101:LYS:H	2.16	0.43
1:A:298:ASN:ND2	1:A:300:ASP:HB2	2.32	0.43
1:D:11:VAL:HG12	1:D:12:THR:N	2.33	0.43
1:D:59:HIS:O	1:D:60:LYS:HB2	2.18	0.43
1:D:236:ILE:HA	6:D:2461:HOH:O	2.18	0.43
1:B:250:THR:HG23	1:B:253:GLN:H	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:ASN:ND2	1:B:46:SER:HB2	2.33	0.43
1:D:157:ASP:O	1:D:161:GLU:HG3	2.18	0.42
1:D:229:LYS:HE2	1:D:229:LYS:HB3	1.89	0.42
1:A:87:LYS:HG2	1:A:91:GLU:OE1	2.19	0.42
1:D:164:LYS:HA	1:D:164:LYS:CE	2.42	0.42
1:D:294:LEU:HA	1:D:295:PRO:HD3	1.89	0.42
1:D:199:LYS:HA	1:D:227:VAL:HA	2.02	0.42
1:B:276:VAL:HG12	1:B:280:LYS:HE3	2.01	0.42
1:D:199:LYS:N	1:D:227:VAL:HG23	2.34	0.42
1:B:69:PRO:HG2	4:B:401:HEC:HBA1	2.02	0.42
1:D:258:MET:SD	1:D:262:GLN:HG3	2.60	0.42
1:A:219:THR:HB	1:A:221:ASN:ND2	2.35	0.42
1:D:32:SER:HB3	1:D:37:ILE:O	2.20	0.42
1:D:115:ALA:O	1:D:121:ARG:HD3	2.19	0.42
1:D:192:VAL:HG12	1:D:192:VAL:O	2.20	0.42
1:D:198:GLN:HA	1:D:198:GLN:NE2	2.34	0.42
1:D:268:ASN:HD21	1:D:271:GLU:CG	2.32	0.42
1:A:77:ASN:HB3	1:A:79:ALA:O	2.20	0.41
1:D:5:ILE:CD1	1:D:5:ILE:H	2.33	0.41
1:A:11:VAL:HG23	1:A:150:VAL:HG21	2.01	0.41
1:A:258:MET:O	1:A:262:GLN:HB2	2.20	0.41
1:B:119:GLN:CD	1:B:295:PRO:HB3	2.40	0.41
4:A:401:HEC:HBC3	4:A:401:HEC:HMC1	2.02	0.41
1:B:154:SER:O	1:B:158:LYS:HG3	2.20	0.41
1:B:112:LYS:HD3	1:B:302:PRO:CG	2.50	0.41
1:D:130:SER:HB2	1:D:137:ARG:CZ	2.50	0.41
1:D:183:CYS:CB	4:D:402:HEC:CAB	2.98	0.41
1:A:87:LYS:HE2	6:A:1550:HOH:O	2.21	0.41
1:B:80:GLN:HB2	1:B:86:ALA:HB3	2.02	0.41
1:C:164:LYS:HA	1:C:164:LYS:HD3	1.87	0.41
1:B:156:PHE:CE2	1:B:160:LEU:HD11	2.56	0.41
1:C:3:GLU:HA	1:C:4:PRO:HD3	1.82	0.41
1:D:119:GLN:NE2	1:D:296:PRO:HD2	2.36	0.41
1:D:141:ALA:O	1:D:144:GLN:HB3	2.21	0.41
1:D:236:ILE:HG13	1:D:251:LEU:HD21	2.03	0.41
1:D:201:GLY:N	1:D:262:GLN:NE2	2.54	0.41
1:D:78:LEU:HD12	1:D:79:ALA:CB	2.51	0.41
1:A:13:PRO:HG3	1:A:148:THR:HG22	2.03	0.41
1:B:159:TRP:CD1	1:B:164:LYS:HA	2.56	0.41
1:D:119:GLN:HE22	1:D:296:PRO:HD2	1.86	0.41
1:D:21:LEU:O	1:D:25:LEU:CD2	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:PHE:CE2	1:A:133:VAL:HG13	2.55	0.40
1:A:80:GLN:O	1:A:81:PHE:CB	2.68	0.40
4:A:402:HEC:HBB3	4:A:402:HEC:HMB1	2.02	0.40
1:D:80:GLN:HA	1:D:80:GLN:HE21	1.85	0.40
1:D:151:THR:HG23	6:D:2519:HOH:O	2.20	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	306/308 (99%)	296 (97%)	10 (3%)	0	100	100
1	B	306/308 (99%)	287 (94%)	16 (5%)	3 (1%)	15	5
1	C	306/308 (99%)	298 (97%)	8 (3%)	0	100	100
1	D	306/308 (99%)	285 (93%)	19 (6%)	2 (1%)	22	10
All	All	1224/1232 (99%)	1166 (95%)	53 (4%)	5 (0%)	34	21

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2	ASN
1	D	2	ASN
1	B	163	ASP
1	D	215	ARG
1	B	3	GLU

### 5.3.2 Protein sidechains [i](#)

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

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	260/260 (100%)	255 (98%)	5 (2%)	57	46
1	B	260/260 (100%)	254 (98%)	6 (2%)	50	37
1	C	260/260 (100%)	255 (98%)	5 (2%)	57	46
1	D	260/260 (100%)	250 (96%)	10 (4%)	33	18
All	All	1040/1040 (100%)	1014 (98%)	26 (2%)	47	34

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	GLN
1	A	122	GLU
1	A	210	ASN
1	A	222	GLU
1	A	298	ASN
1	B	44	ASN
1	B	99	ASN
1	B	119	GLN
1	B	122	GLU
1	B	177	LEU
1	B	264	ASN
1	C	15	ASN
1	C	99	ASN
1	C	120	TYR
1	C	264	ASN
1	C	298	ASN
1	D	15	ASN
1	D	23	LYS
1	D	126	LYS
1	D	132	GLU
1	D	135	ILE
1	D	164	LYS
1	D	188	ASN
1	D	225	ARG
1	D	268	ASN
1	D	308	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (56) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	52	ASN
1	A	62	GLN
1	A	119	GLN
1	A	165	ASN
1	A	198	GLN
1	A	210	ASN
1	A	221	ASN
1	A	226	ASN
1	A	253	GLN
1	A	298	ASN
1	A	299	ASN
1	B	40	ASN
1	B	44	ASN
1	B	52	ASN
1	B	59	HIS
1	B	63	GLN
1	B	99	ASN
1	B	119	GLN
1	B	165	ASN
1	B	176	ASN
1	B	198	GLN
1	B	226	ASN
1	B	235	ASN
1	B	264	ASN
1	B	268	ASN
1	B	299	ASN
1	C	2	ASN
1	C	15	ASN
1	C	40	ASN
1	C	52	ASN
1	C	62	GLN
1	C	63	GLN
1	C	99	ASN
1	C	119	GLN
1	C	176	ASN
1	C	185	GLN
1	C	221	ASN
1	C	226	ASN
1	C	253	GLN
1	C	262	GLN
1	C	264	ASN
1	C	298	ASN

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Mol	Chain	Res	Type
1	D	40	ASN
1	D	62	GLN
1	D	63	GLN
1	D	77	ASN
1	D	80	GLN
1	D	119	GLN
1	D	144	GLN
1	D	198	GLN
1	D	226	ASN
1	D	262	GLN
1	D	268	ASN
1	D	298	ASN
1	D	299	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 7 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	$\# Z  > 2$
5	GOL	B	1410	-	5,5,5	0.29	0	5,5,5	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	HEC	B	402	1	26,50,50	1.45	3 (11%)	18,82,82	0.78	0
4	HEC	D	402	1	26,50,50	1.48	4 (15%)	18,82,82	1.80	1 (5%)
4	HEC	B	401	1,6	26,50,50	1.41	6 (23%)	18,82,82	0.85	1 (5%)
4	HEC	A	401	1,6	26,50,50	1.26	3 (11%)	18,82,82	1.03	1 (5%)
4	HEC	D	401	1,6	26,50,50	1.45	4 (15%)	18,82,82	0.78	0
4	HEC	A	402	1	26,50,50	1.56	2 (7%)	18,82,82	1.22	1 (5%)
4	HEC	C	402	1	26,50,50	1.50	3 (11%)	18,82,82	1.29	3 (16%)
4	HEC	C	401	1,6	26,50,50	1.49	4 (15%)	18,82,82	1.08	2 (11%)

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
5	GOL	B	1410	-	-	0/4/4/4	-
4	HEC	B	402	1	-	0/6/54/54	-
4	HEC	D	402	1	-	1/6/54/54	-
4	HEC	B	401	1,6	-	0/6/54/54	-
4	HEC	A	401	1,6	-	0/6/54/54	-
4	HEC	D	401	1,6	-	1/6/54/54	-
4	HEC	A	402	1	-	0/6/54/54	-
4	HEC	C	402	1	-	0/6/54/54	-
4	HEC	C	401	1,6	-	0/6/54/54	-

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	402	HEC	C3C-C2C	-4.63	1.35	1.40
4	A	402	HEC	C3B-C2B	-4.60	1.35	1.40
4	A	402	HEC	C3C-C2C	-4.48	1.36	1.40
4	C	401	HEC	C3B-C2B	-4.30	1.36	1.40
4	D	402	HEC	C3C-C2C	-4.05	1.36	1.40
4	B	401	HEC	C3C-C2C	-4.01	1.36	1.40
4	C	402	HEC	C3B-C2B	-3.96	1.36	1.40
4	A	401	HEC	C3B-C2B	-3.73	1.36	1.40
4	B	402	HEC	C3C-C2C	-3.69	1.36	1.40
4	D	401	HEC	C3C-C2C	-3.57	1.37	1.40
4	B	402	HEC	C3B-C2B	-3.55	1.37	1.40
4	D	401	HEC	C3B-C2B	-3.38	1.37	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	401	HEC	C3C-C2C	-3.37	1.37	1.40
4	D	402	HEC	C3B-C2B	-3.29	1.37	1.40
4	A	401	HEC	C3C-C2C	-2.53	1.38	1.40
4	B	401	HEC	C4D-ND	2.52	1.41	1.36
4	D	402	HEC	C4D-ND	2.48	1.41	1.36
4	B	401	HEC	C3B-C2B	-2.48	1.38	1.40
4	B	401	HEC	C3C-C4C	2.39	1.47	1.43
4	C	401	HEC	C1A-C2A	2.24	1.47	1.42
4	D	402	HEC	C3C-C4C	2.11	1.46	1.43
4	A	401	HEC	C1C-NC	2.10	1.40	1.36
4	D	401	HEC	C1C-NC	2.09	1.40	1.36
4	C	402	HEC	C4D-ND	2.08	1.40	1.36
4	B	402	HEC	C3C-C4C	2.04	1.46	1.43
4	B	401	HEC	C1D-ND	2.03	1.40	1.36
4	D	401	HEC	C1D-ND	2.03	1.40	1.36
4	C	401	HEC	C1B-NB	2.02	1.40	1.36
4	B	401	HEC	C1B-NB	2.01	1.40	1.36

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	402	HEC	CBD-CAD-C3D	6.52	124.51	112.49
4	C	401	HEC	CBD-CAD-C3D	-2.69	107.52	112.49
4	A	402	HEC	CBD-CAD-C3D	-2.67	107.56	112.49
4	A	401	HEC	CBD-CAD-C3D	-2.62	107.66	112.49
4	C	402	HEC	CMB-C2B-C1B	-2.52	124.58	128.46
4	C	401	HEC	CMC-C2C-C3C	2.28	128.50	125.82
4	C	402	HEC	CBD-CAD-C3D	-2.24	108.35	112.49
4	C	402	HEC	CMC-C2C-C1C	-2.24	125.03	128.46
4	B	401	HEC	CMC-C2C-C1C	-2.05	125.32	128.46

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	402	HEC	C4D-C3D-CAD-CBD
4	D	401	HEC	C3A-C2A-CAA-CBA

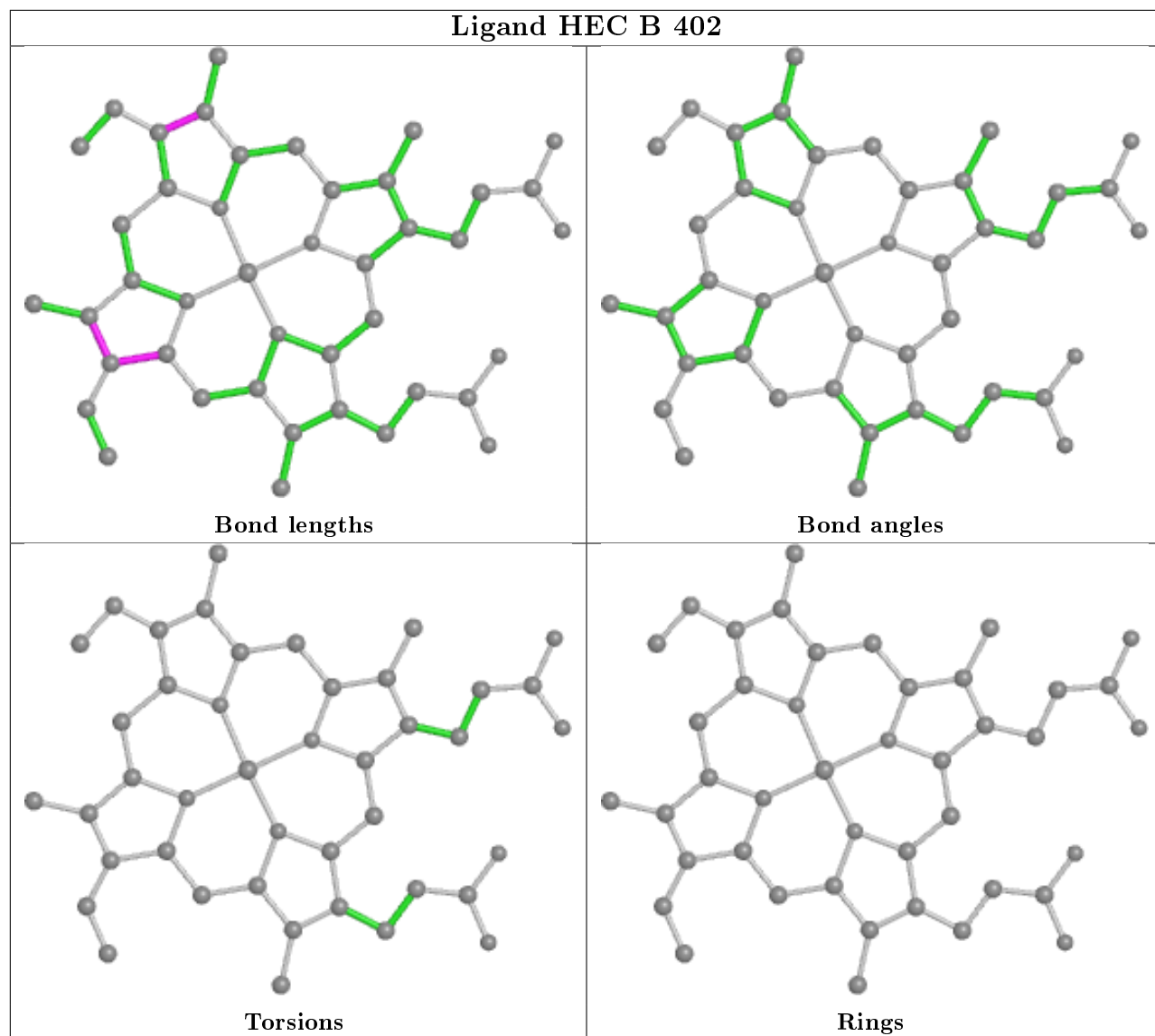
There are no ring outliers.

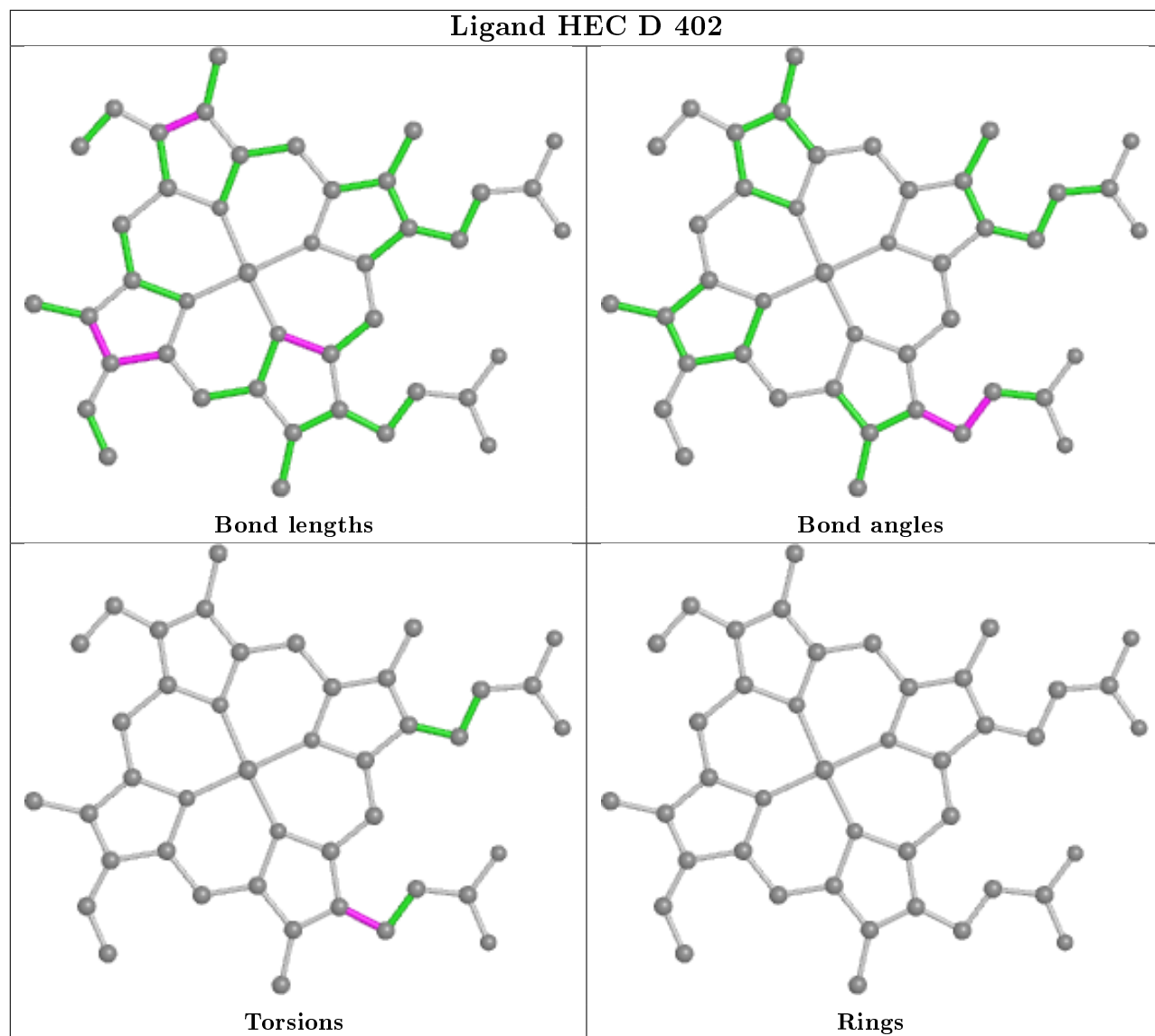
8 monomers are involved in 61 short contacts:

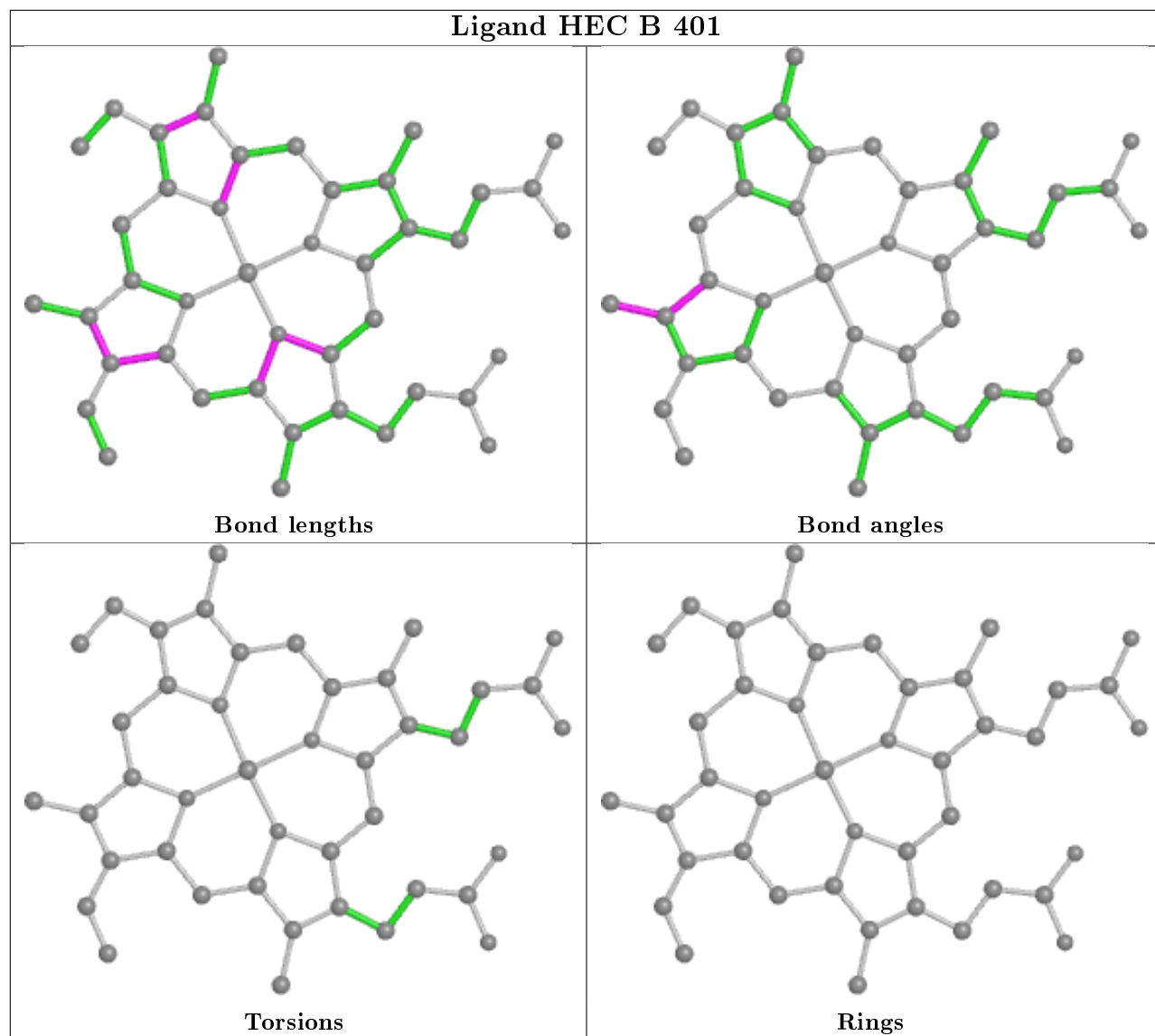


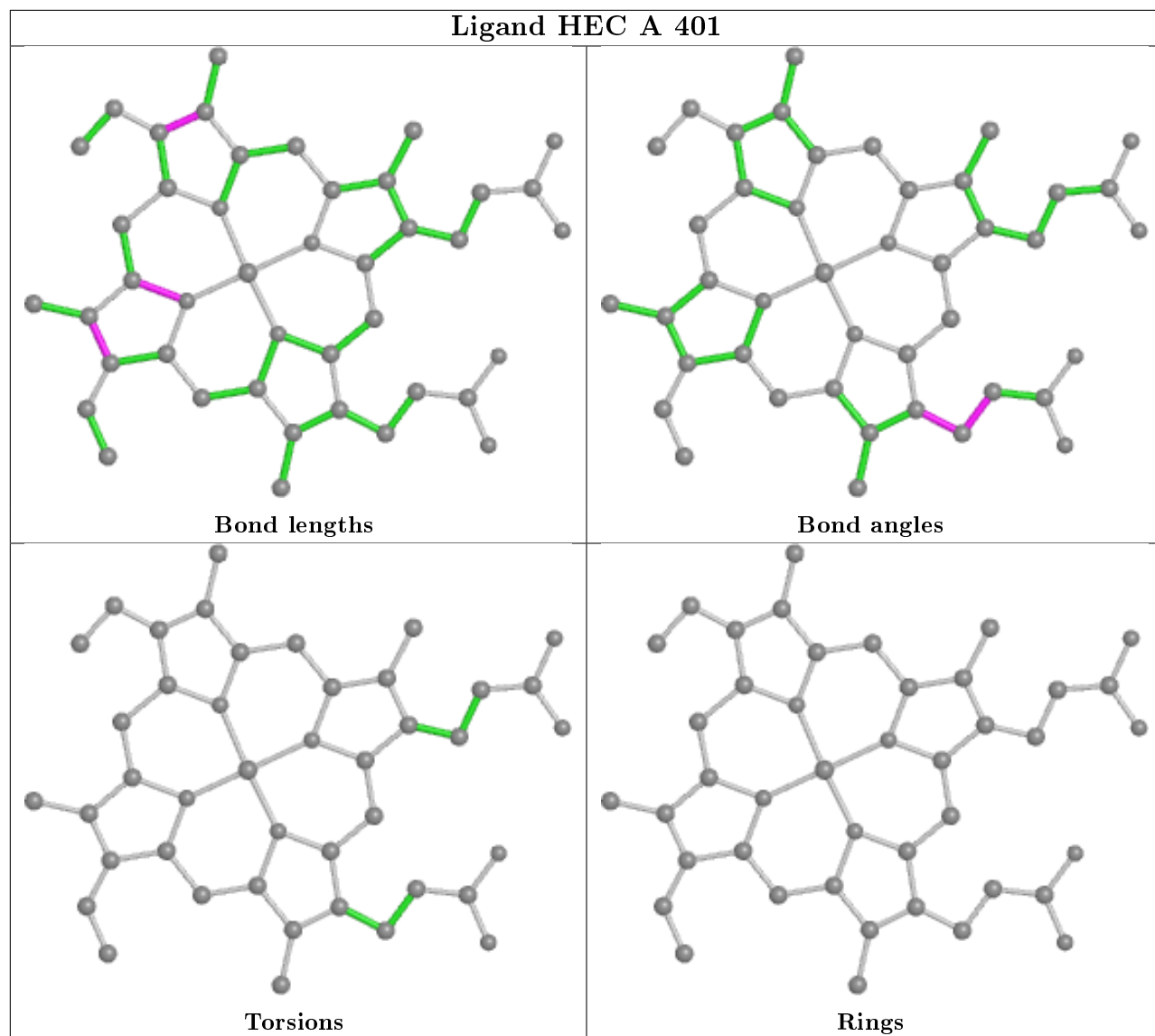
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	402	HEC	6	0
4	D	402	HEC	11	0
4	B	401	HEC	9	0
4	A	401	HEC	7	0
4	D	401	HEC	6	0
4	A	402	HEC	8	0
4	C	402	HEC	7	0
4	C	401	HEC	7	0

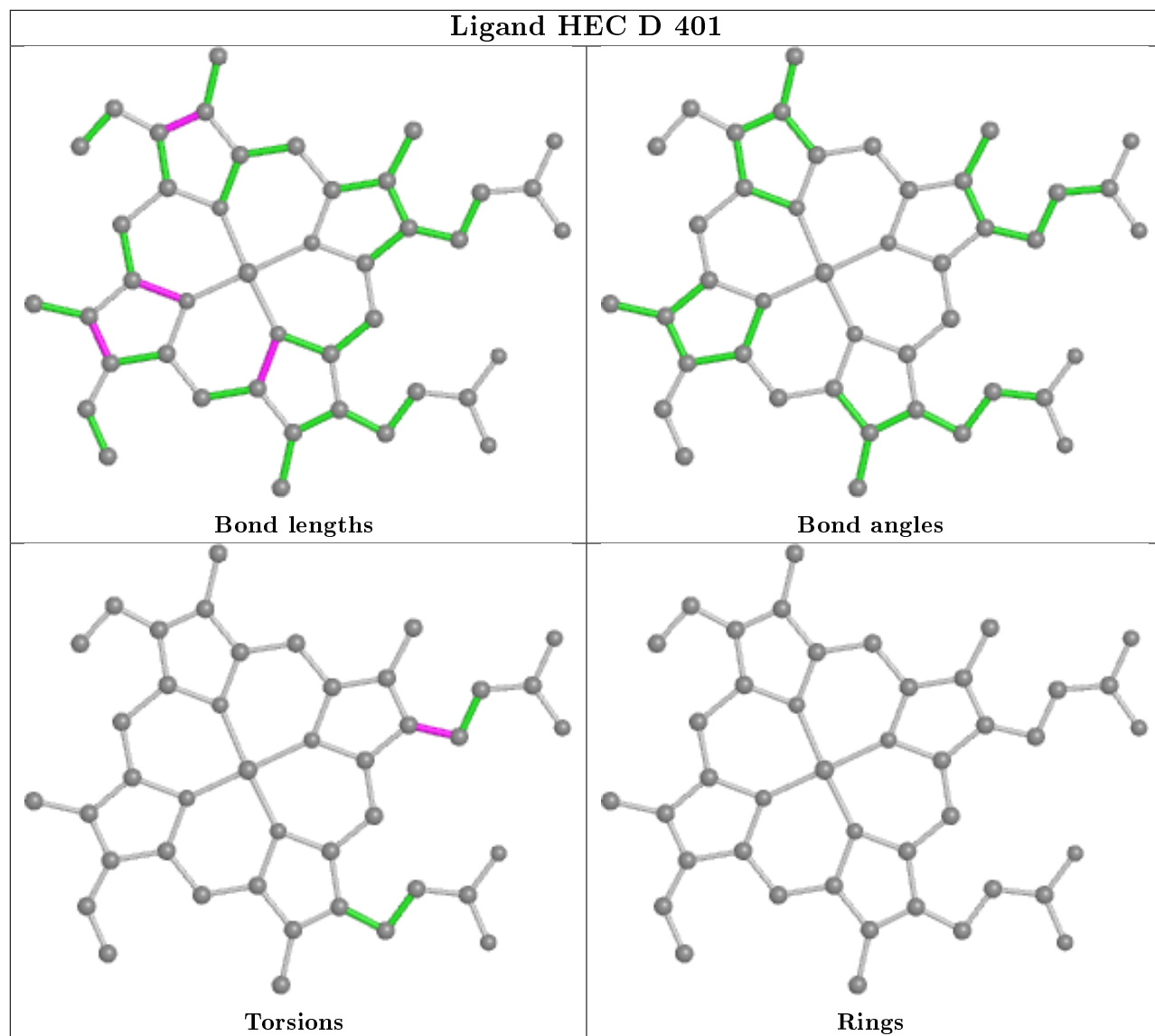
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

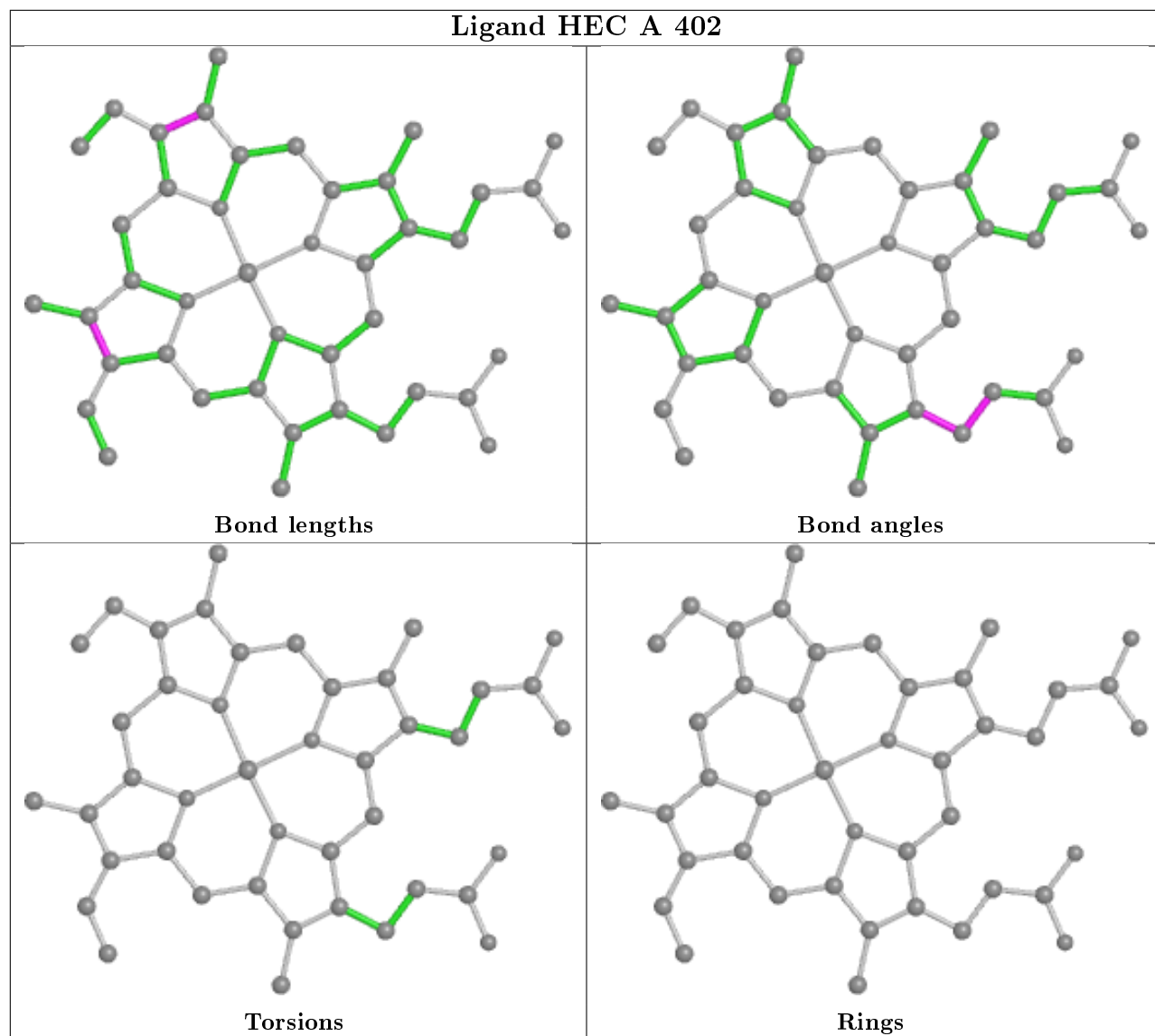


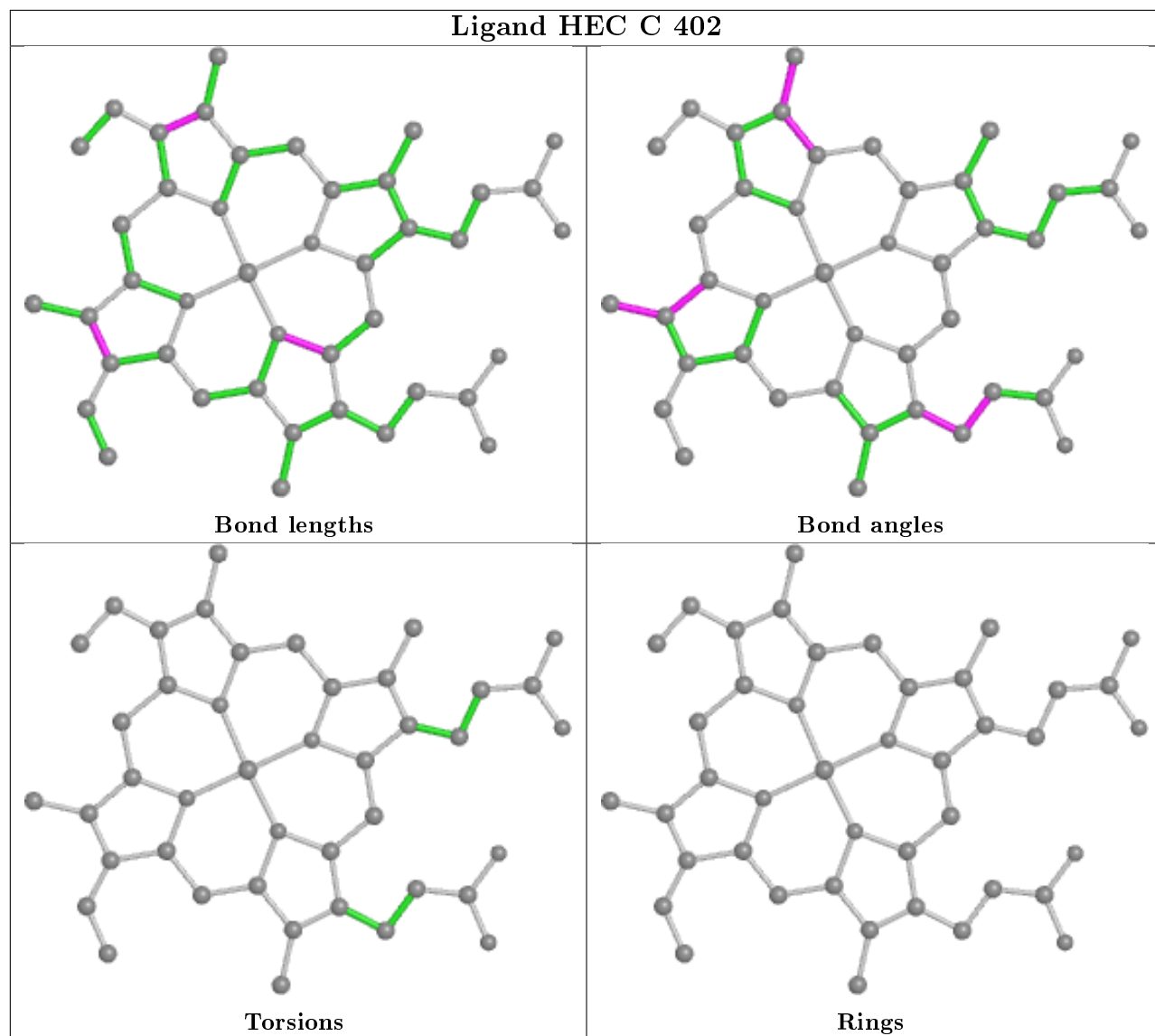




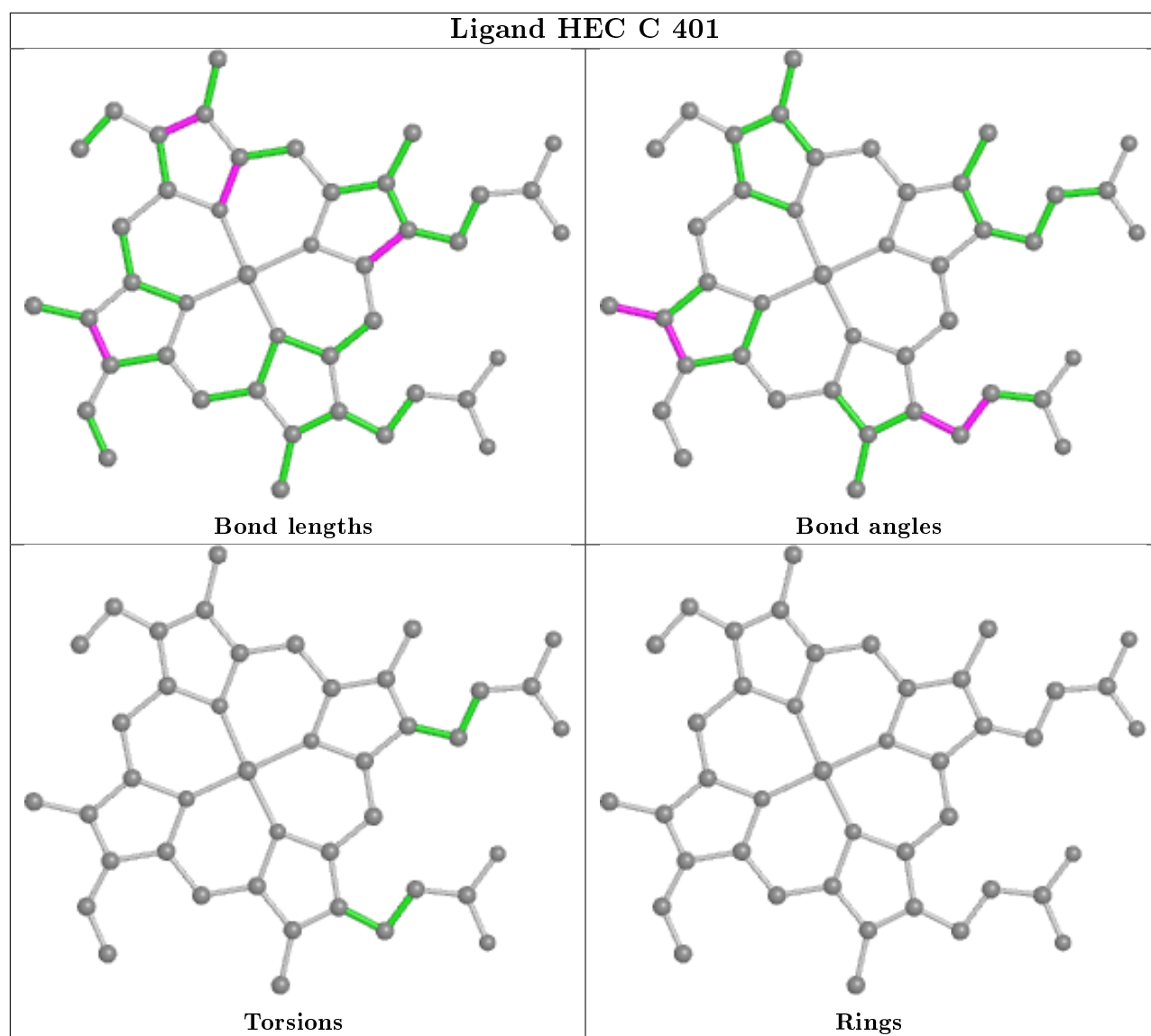












## 5.7 Other polymers [i](#)

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.