Full wwPDB NMR Structure Validation Report (i)

Oct 23, 2021 – 05:35 PM EDT

PDB ID	:	1YFC
Title	:	Solution nmr structure of a yeast iso-1-ferrocytochrome C
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Deposited on	:	1996-08-08

This is a Full wwPDB NMR 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/NMRValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (i)) were used in the production of this report:

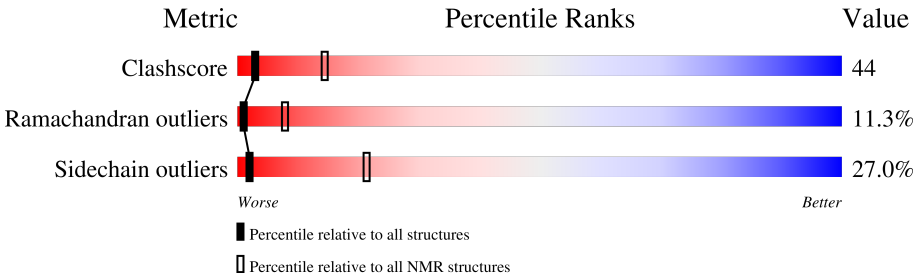
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.23.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

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)	NMR archive (#Entries)
Clashscore	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Mol	Chain	Length	Quality of chain
1	A	108	26% 54% 18% •

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 17 is the overall representative, medoid model (most similar to other models).

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	**Residue range (total)**	**Backbone RMSD (Å)**	**Medoid model**
1	A:-3-A:-1, A:1-A:71, A:73-A:103 (105)	0.56	17

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

NmrClust was unable to cluster the ensemble.

Error message: 1 Unexpected character

3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1777 atoms, of which 885 are hydrogens and 0 are deuteriums.

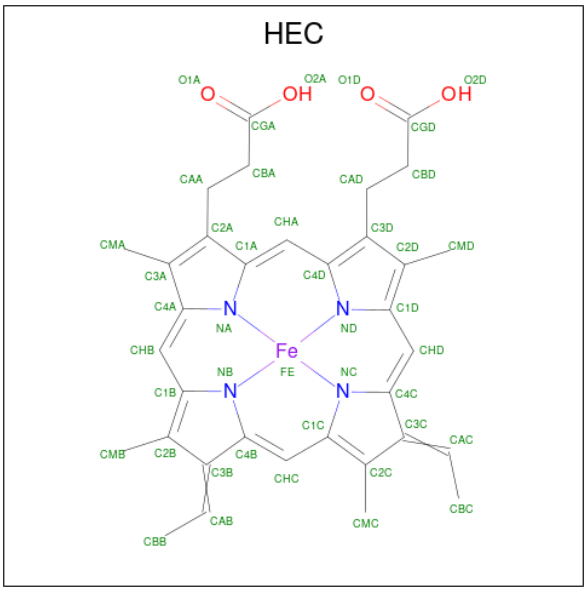
- Molecule 1 is a protein called YEAST ISO-1-FERROCYTOCHROME C.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	108	1702	537	853	151	157	4	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	72	M3L	LYS	modified residue	UNP P00044
A	102	SER	CYS	engineered mutation	UNP P00044

- Molecule 2 is HEME C (three-letter code: HEC) (formula: $C_{34}H_{34}FeN_4O_4$).



Mol	Chain	Residues	Atoms					
			Total	C	Fe	H	N	O
2	A	1	75	34	1	32	4	4

4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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 outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

• Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

Chain A: 26% 54% 18%

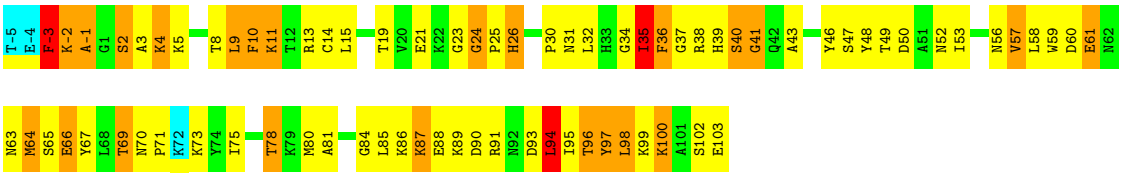
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

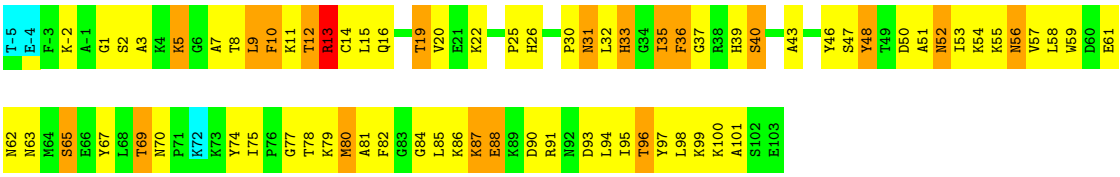
Chain A: 27% 46% 21%



4.2.2 Score per residue for model 2

• Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

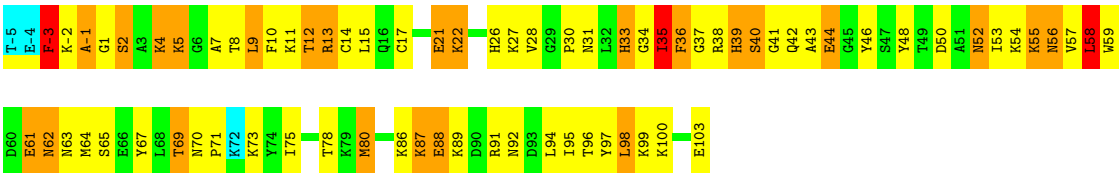
Chain A: 29% 50% 18%



4.2.3 Score per residue for model 3

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

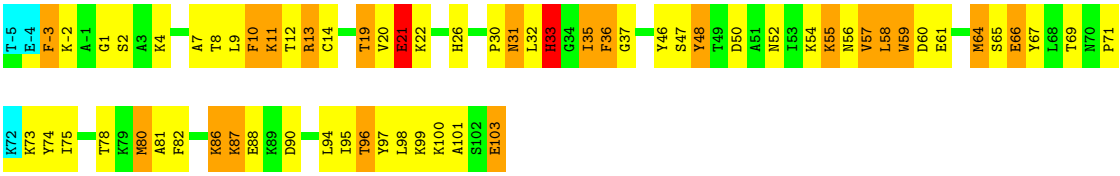
Chain A: 29% 44% 22%



4.2.4 Score per residue for model 4

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

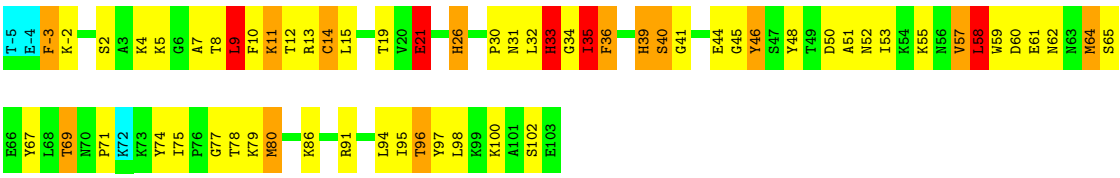
Chain A: 38% 39% 19%



4.2.5 Score per residue for model 5

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

Chain A: 40% 41% 12% 5%



4.2.6 Score per residue for model 6

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

Chain A: 35% 38% 22%

T-5 E-4 F-3 K-2 A-1 K4 K5 G6 A7 T8 L9 F10 K11 T12 R13 C14 L15 T19 V20 E21 K22 H26 K27 V28 G29 P30 N31 L32 H33 G34 I35 F36 G37 R38 H39 S40 G41 Q42 Y46 S47 Y48 T49 D50 A51 N52 K55 N56 V57 L58 W59 D60 E61 N62 N63 M64

S65 E66 Y67 L68 T69 N70 P71 K72 G77 T78 K79 M80 A81 K86 K89 D90 R91 L94 I95 T96 Y97 L98 K99 K100 A101 S102 E103

4.2.7 Score per residue for model 7

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

Chain A: 29% 49% 18%

T-5 E-4 F-3 G1 S2 A3 K4 K5 G6 A7 T8 L9 F10 K11 T12 R13 C14 L15 T19 V20 E21 K22 H26 P30 N31 L32 H33 G34 I35 F36 G37 R38 H39 S40 G41 Q42 A43 Y46 S47 Y48 T49 D50 A51 N52 I53 K54 K55 N56 V57 L58 W59 D60 E61 N62 N63

M64 S65 E66 Y67 N70 P71 K72 K73 Y74 I75 P76 G77 T78 K79 M80 A81 F82 G83 G84 K87 E88 K89 D90 R91 N92 D93 L94 I95 T96 Y97 L98 K99 K100 A101 S102 E103

4.2.8 Score per residue for model 8

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

Chain A: 35% 46% 14%

T-5 E-4 F-3 K-2 A-1 G1 S2 A3 K4 K5 G6 A7 T8 L9 F10 K11 T12 R13 C14 L15 T19 V20 E21 K22 H26 K27 V28 G29 P30 N31 L32 H33 G34 I35 F36 G37 R38 H39 S40 G41 Q42 A43 Y46 S47 Y48 T49 D50 A51 N52 K55 N56 V57 L58 W59 D60 E61

N62 N63 M64 S65 E66 Y67 P71 K72 I75 T78 K79 M80 A81 L85 K86 K87 E88 K89 D90 R91 N92 D93 L94 I95 T96 Y97 L98 K99 K100 A101 S102 E103

4.2.9 Score per residue for model 9

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

Chain A: 39% 37% 20%

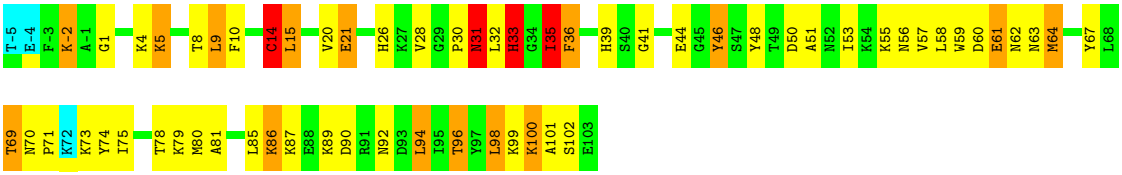
T-5 E-4 G1 S2 A3 K4 K5 T8 L9 F10 K11 T12 R13 C14 L15 Q16 C17 H18 T19 V20 E21 K22 H26 K27 V28 G29 P30 N31 L32 H33 G34 I35 F36 H39 S40 G41 Y46 S47 Y48 T49 D50 A51 N52 I53 K54 K55 N56 V57 L58 W59 D60 E61 M64 S65

L68 T69 N70 P71 K72 K73 T78 K79 M80 A81 F82 G83 G84 L85 K86 K87 E88 R91 L94 I95 T96 Y97 L98 K99 K100 A101 S102 E103

4.2.10 Score per residue for model 10

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

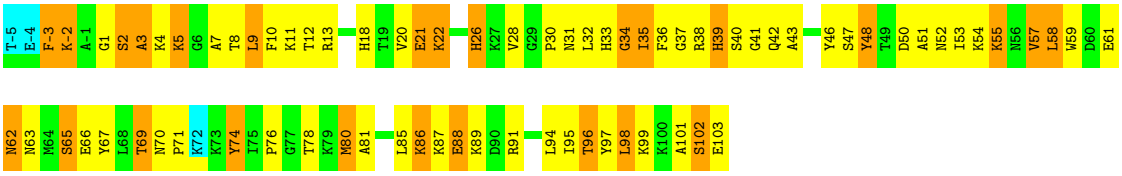
Chain A: 41% 39% 14%



4.2.11 Score per residue for model 11

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

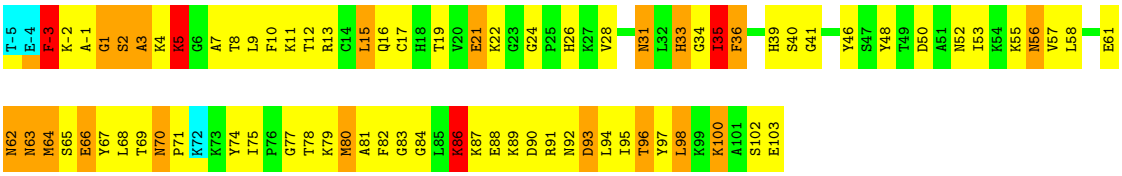
Chain A: 28% 45% 24%



4.2.12 Score per residue for model 12

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

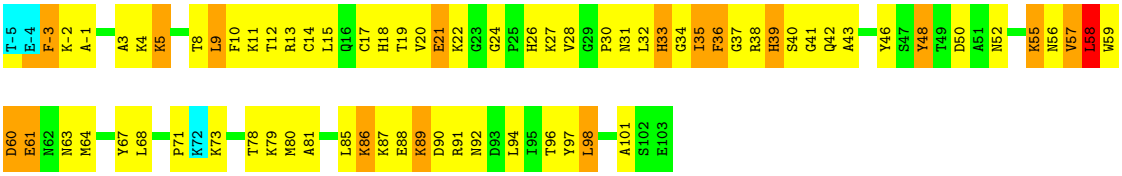
Chain A: 25% 51% 18%



4.2.13 Score per residue for model 13

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

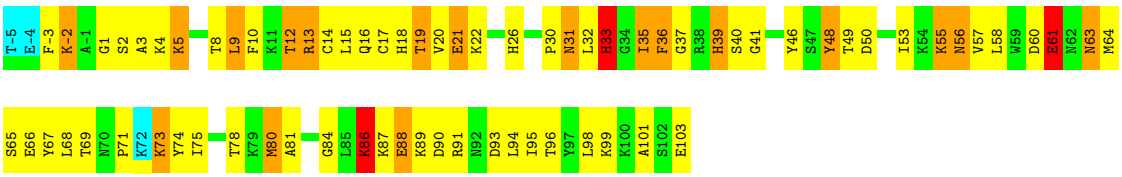
Chain A: 31% 51% 15%



4.2.14 Score per residue for model 14

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

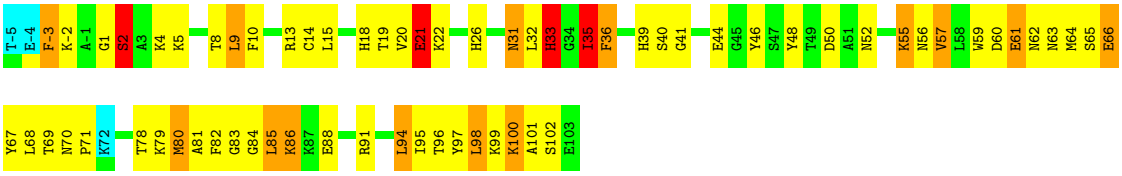
Chain A: 31% 47% 17%



4.2.15 Score per residue for model 15

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

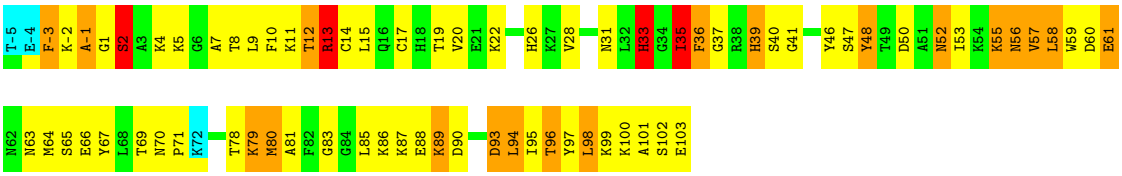
Chain A: 35% 45% 13%



4.2.16 Score per residue for model 16

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

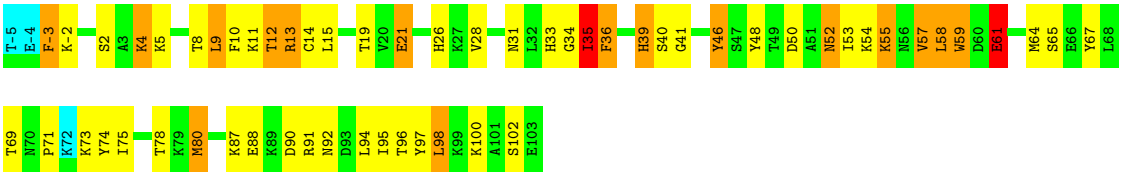
Chain A: 30% 46% 18%



4.2.17 Score per residue for model 17 (medoid)

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

Chain A: 44% 37% 15%



4.2.18 Score per residue for model 18

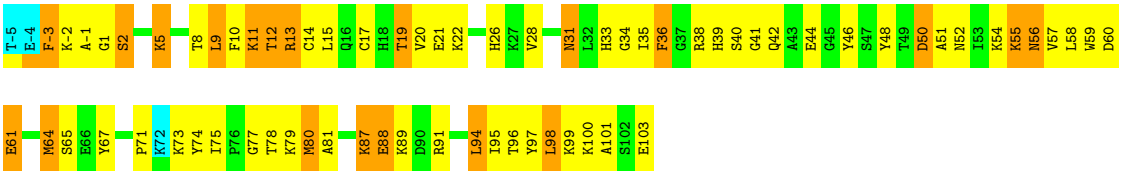
- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

Chain A: 28% 51% 17%

4.2.19 Score per residue for model 19

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

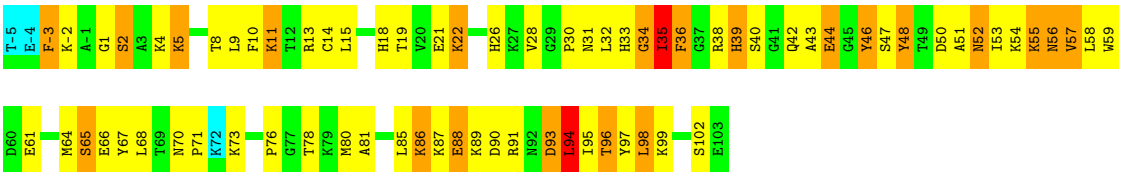
Chain A: 32% 46% 19%



4.2.20 Score per residue for model 20

- Molecule 1: YEAST ISO-1-FERROCYTOCHROME C

Chain A: 30% 46% 19%



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *distance geometry*.

Of the 500 calculated structures, 20 were deposited, based on the following criterion: *TARGET FUNCTION <= 0.82 A2*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
DIANA	structure solution	
DIANA	refinement	

No chemical shift data was provided.

6 Model quality (i)

6.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: HEC, M3L

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	821	821	821	76±11
2	A	43	32	30	9±3
All	All	17280	17060	17026	1521

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 44.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:14:CYS:SG	2:A:104:HEC:HAB	1.69	1.28	13	3
1:A:14:CYS:SG	2:A:104:HEC:C3B	1.20	2.27	5	1
1:A:67:TYR:CE1	1:A:75:ILE:HD12	1.07	1.85	17	7
1:A:9:LEU:HD22	1:A:94:LEU:HD22	0.99	1.33	2	7
1:A:9:LEU:HD22	1:A:94:LEU:HD13	0.95	1.37	17	5
1:A:-3:PHE:CE2	1:A:-1:ALA:HB2	0.91	2.01	16	1
1:A:20:VAL:HG21	1:A:101:ALA:HB1	0.90	1.41	18	9
1:A:-3:PHE:CZ	1:A:-1:ALA:HB2	0.90	2.02	16	1
1:A:26:HIS:CE1	1:A:43:ALA:HB1	0.89	2.01	20	1
1:A:28:VAL:O	2:A:104:HEC:HMD3	0.88	1.68	17	12
1:A:38:ARG:O	1:A:58:LEU:HD12	0.87	1.68	7	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:39:HIS:CD2	1:A:58:LEU:HD13	0.86	2.05	7	6
1:A:98:LEU:HD11	2:A:104:HEC:CMB	0.85	2.02	3	1
1:A:9:LEU:HD21	1:A:85:LEU:HD21	0.85	1.48	2	1
1:A:9:LEU:HD22	1:A:94:LEU:CD1	0.85	2.01	9	10
1:A:9:LEU:HD22	1:A:94:LEU:CD2	0.84	2.02	1	5
1:A:9:LEU:HD22	1:A:94:LEU:HD23	0.84	1.47	18	1
1:A:9:LEU:HD22	1:A:94:LEU:HD12	0.83	1.50	15	4
1:A:94:LEU:HD13	1:A:97:TYR:CE2	0.83	2.08	1	1
1:A:20:VAL:HG22	1:A:102:SER:OG	0.82	1.72	18	1
1:A:9:LEU:HD22	1:A:94:LEU:HG	0.82	1.50	6	2
1:A:94:LEU:CD1	1:A:98:LEU:HD12	0.82	2.04	18	2
1:A:14:CYS:HG	2:A:104:HEC:HAB	0.81	1.28	13	2
1:A:57:VAL:HG11	1:A:67:TYR:OH	0.80	1.76	11	2
1:A:57:VAL:HG11	1:A:59:TRP:CH2	0.80	2.12	3	5
1:A:67:TYR:CE1	1:A:74:TYR:CD2	0.79	2.70	11	1
1:A:14:CYS:CB	2:A:104:HEC:HAB	0.79	2.06	13	2
1:A:94:LEU:HD12	1:A:98:LEU:HD12	0.78	1.56	18	1
1:A:81:ALA:O	2:A:104:HEC:HBC2	0.76	1.81	1	7
1:A:10:PHE:CD2	1:A:15:LEU:HD13	0.76	2.16	14	9
1:A:94:LEU:HD21	2:A:104:HEC:CBB	0.76	2.11	5	6
1:A:20:VAL:HG23	1:A:101:ALA:HB1	0.76	1.55	4	3
1:A:32:LEU:HD11	2:A:104:HEC:CHB	0.75	2.11	1	7
1:A:14:CYS:SG	2:A:104:HEC:CBB	0.75	2.74	16	4
1:A:97:TYR:CE1	1:A:101:ALA:HB2	0.74	2.18	13	1
1:A:9:LEU:HD11	1:A:85:LEU:HD22	0.73	1.61	20	1
1:A:64:MET:SD	1:A:68:LEU:HD11	0.72	2.23	12	3
1:A:35:ILE:O	1:A:36:PHE:CG	0.72	2.42	7	1
1:A:67:TYR:CE1	1:A:74:TYR:CE2	0.72	2.77	11	1
1:A:67:TYR:CZ	1:A:75:ILE:HD12	0.71	2.21	17	1
1:A:12:THR:HG22	1:A:12:THR:O	0.71	1.84	4	1
1:A:9:LEU:HD21	1:A:13:ARG:NH1	0.71	2.01	4	1
1:A:91:ARG:O	1:A:95:ILE:HD12	0.71	1.85	17	8
1:A:7:ALA:HB2	1:A:97:TYR:CE2	0.70	2.21	7	2
1:A:19:THR:HG23	1:A:31:ASN:ND2	0.70	2.00	19	5
1:A:-3:PHE:CZ	1:A:-1:ALA:CB	0.70	2.74	16	1
1:A:48:TYR:CE2	2:A:104:HEC:CGA	0.70	2.74	7	6
1:A:94:LEU:HD23	1:A:98:LEU:CD1	0.70	2.17	9	1
1:A:65:SER:OG	1:A:95:ILE:HD11	0.70	1.87	14	7
1:A:10:PHE:CE2	1:A:15:LEU:CD1	0.69	2.75	14	14
1:A:35:ILE:CD1	1:A:36:PHE:CD2	0.69	2.75	9	5
1:A:13:ARG:O	2:A:104:HEC:HMC3	0.69	1.87	1	7

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:35:ILE:HD12	1:A:36:PHE:N	0.69	2.03	2	6
1:A:61:GLU:HG2	1:A:95:ILE:HG21	0.69	1.63	3	1
1:A:9:LEU:CD2	1:A:94:LEU:HD13	0.69	2.18	14	3
1:A:9:LEU:HD11	1:A:90:ASP:CG	0.69	2.09	18	3
1:A:35:ILE:CD1	1:A:36:PHE:CD1	0.69	2.76	18	1
1:A:43:ALA:HB3	1:A:48:TYR:CE1	0.69	2.23	1	6
1:A:85:LEU:HD11	1:A:91:ARG:HG2	0.68	1.66	9	1
1:A:36:PHE:CZ	1:A:64:MET:CE	0.68	2.76	20	5
1:A:10:PHE:CD2	1:A:15:LEU:CD1	0.67	2.77	17	5
1:A:94:LEU:HD22	1:A:98:LEU:HD12	0.67	1.64	8	2
1:A:57:VAL:CG1	1:A:59:TRP:CZ2	0.67	2.78	3	8
1:A:9:LEU:HD22	1:A:94:LEU:CG	0.67	2.19	6	2
1:A:57:VAL:CG1	1:A:59:TRP:CH2	0.67	2.77	1	4
1:A:98:LEU:HD11	2:A:104:HEC:HMB1	0.66	1.65	3	1
1:A:20:VAL:CG2	1:A:101:ALA:HB1	0.66	2.21	8	5
1:A:94:LEU:HD23	1:A:98:LEU:HD12	0.66	1.67	9	1
1:A:61:GLU:HG3	1:A:95:ILE:HG21	0.66	1.66	19	2
1:A:39:HIS:CE1	1:A:58:LEU:HD13	0.65	2.25	10	6
1:A:49:THR:HG23	2:A:104:HEC:O2D	0.65	1.91	1	1
1:A:9:LEU:HD13	1:A:94:LEU:HD23	0.65	1.68	20	2
1:A:97:TYR:C	1:A:97:TYR:CD1	0.65	2.69	1	1
1:A:97:TYR:CD1	1:A:98:LEU:N	0.65	2.65	1	1
1:A:51:ALA:HB1	1:A:75:ILE:CG2	0.64	2.23	19	1
1:A:22:LYS:CD	1:A:33:HIS:CE1	0.64	2.80	8	1
1:A:35:ILE:HD11	1:A:36:PHE:CD2	0.64	2.27	8	5
1:A:57:VAL:O	1:A:57:VAL:HG13	0.64	1.90	1	4
1:A:18:HIS:CG	1:A:32:LEU:HD11	0.64	2.28	11	2
1:A:9:LEU:CD2	1:A:94:LEU:HD22	0.63	2.24	14	6
1:A:20:VAL:HG21	1:A:101:ALA:CB	0.63	2.23	7	4
1:A:39:HIS:CD2	1:A:58:LEU:HD12	0.63	2.29	3	2
1:A:14:CYS:SG	2:A:104:HEC:C2B	0.62	2.87	5	1
1:A:7:ALA:CB	1:A:97:TYR:CE1	0.62	2.82	11	1
1:A:94:LEU:HD21	2:A:104:HEC:HBB1	0.62	1.70	17	3
1:A:36:PHE:CE1	1:A:64:MET:HE2	0.62	2.30	6	2
1:A:7:ALA:HB2	1:A:97:TYR:CE1	0.62	2.30	11	2
1:A:67:TYR:CD1	1:A:75:ILE:HD12	0.62	2.29	18	2
1:A:48:TYR:CE2	2:A:104:HEC:O2A	0.61	2.53	9	7
1:A:19:THR:HG23	1:A:31:ASN:OD1	0.61	1.96	14	1
1:A:9:LEU:HD11	1:A:90:ASP:OD2	0.61	1.95	13	1
1:A:39:HIS:CD2	1:A:56:ASN:O	0.61	2.54	3	2
1:A:12:THR:O	1:A:13:ARG:CB	0.61	2.49	16	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:33:HIS:CG	1:A:33:HIS:O	0.61	2.54	9	3
1:A:36:PHE:CZ	1:A:64:MET:SD	0.60	2.94	8	3
1:A:59:TRP:CE3	1:A:64:MET:SD	0.60	2.94	15	5
1:A:94:LEU:HD12	1:A:98:LEU:CD1	0.60	2.26	18	1
1:A:61:GLU:CG	1:A:95:ILE:HG21	0.60	2.27	3	1
1:A:36:PHE:CD1	1:A:60:ASP:O	0.60	2.54	4	8
1:A:94:LEU:HD11	1:A:98:LEU:HD12	0.60	1.71	16	1
1:A:33:HIS:CE1	1:A:102:SER:OG	0.60	2.55	10	1
1:A:94:LEU:O	1:A:97:TYR:CD2	0.60	2.55	1	1
1:A:48:TYR:CE2	2:A:104:HEC:O1A	0.60	2.55	12	6
1:A:-3:PHE:CZ	1:A:-2:LYS:O	0.60	2.55	8	2
1:A:36:PHE:CE1	1:A:60:ASP:O	0.60	2.55	19	5
1:A:39:HIS:CE1	1:A:56:ASN:OD1	0.60	2.55	14	2
1:A:65:SER:HB2	1:A:95:ILE:HD11	0.60	1.72	17	5
1:A:36:PHE:CE1	1:A:64:MET:CE	0.60	2.85	6	2
1:A:-3:PHE:CZ	1:A:61:GLU:OE2	0.60	2.55	1	1
1:A:97:TYR:CG	1:A:98:LEU:N	0.59	2.70	1	1
1:A:36:PHE:CZ	1:A:64:MET:HE3	0.59	2.32	9	1
1:A:63:ASN:O	1:A:67:TYR:CD2	0.59	2.55	16	1
1:A:26:HIS:HE1	1:A:43:ALA:HB1	0.59	1.56	20	1
1:A:38:ARG:C	1:A:58:LEU:HD12	0.59	2.16	7	2
1:A:35:ILE:CD1	1:A:36:PHE:CG	0.59	2.85	4	6
1:A:36:PHE:CZ	1:A:64:MET:HE1	0.59	2.32	20	2
1:A:36:PHE:CE1	1:A:64:MET:SD	0.59	2.96	3	2
1:A:33:HIS:O	1:A:33:HIS:CD2	0.59	2.56	4	2
1:A:-3:PHE:CD2	1:A:61:GLU:OE2	0.59	2.56	5	2
1:A:-3:PHE:CG	1:A:-2:LYS:N	0.59	2.70	18	2
1:A:-3:PHE:CE2	1:A:61:GLU:OE2	0.59	2.55	1	1
1:A:100:LYS:O	1:A:101:ALA:HB3	0.59	1.97	15	2
1:A:43:ALA:HB3	1:A:48:TYR:HE1	0.58	1.58	7	4
1:A:10:PHE:CE2	1:A:15:LEU:HD13	0.58	2.33	14	6
1:A:62:ASN:N	1:A:62:ASN:HD22	0.58	1.96	18	2
1:A:57:VAL:HG13	1:A:59:TRP:CZ2	0.58	2.34	7	3
1:A:39:HIS:CE1	1:A:58:LEU:CD1	0.58	2.87	5	2
1:A:65:SER:HB3	1:A:95:ILE:HD11	0.58	1.74	19	1
1:A:48:TYR:CB	1:A:53:ILE:HD12	0.58	2.29	5	4
1:A:39:HIS:NE2	1:A:58:LEU:HD13	0.58	2.13	7	8
1:A:57:VAL:HG22	1:A:63:ASN:ND2	0.57	2.14	1	3
1:A:6:GLY:O	1:A:97:TYR:CD2	0.57	2.57	18	1
1:A:57:VAL:HG22	1:A:63:ASN:HD21	0.57	1.59	1	3
1:A:22:LYS:HA	1:A:33:HIS:CE1	0.57	2.34	4	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:17:CYS:SG	1:A:28:VAL:HG11	0.56	2.39	9	5
1:A:-3:PHE:CE1	1:A:-1:ALA:CA	0.56	2.88	16	1
1:A:9:LEU:HB3	1:A:94:LEU:HD23	0.56	1.77	10	1
1:A:10:PHE:CE2	1:A:15:LEU:HD12	0.56	2.35	12	2
1:A:6:GLY:O	1:A:97:TYR:CE2	0.56	2.58	18	1
1:A:35:ILE:H	1:A:35:ILE:HD13	0.56	1.60	7	2
1:A:35:ILE:CD1	1:A:36:PHE:N	0.56	2.69	2	4
1:A:67:TYR:CE1	1:A:75:ILE:CD1	0.56	2.76	17	1
1:A:46:TYR:CD1	1:A:47:SER:O	0.56	2.59	18	5
1:A:12:THR:O	1:A:12:THR:CG2	0.55	2.53	4	1
1:A:26:HIS:CB	1:A:46:TYR:CD2	0.55	2.89	13	8
1:A:67:TYR:OH	2:A:104:HEC:C1A	0.55	2.55	2	1
1:A:59:TRP:CE3	1:A:64:MET:CE	0.55	2.90	13	2
1:A:22:LYS:CD	1:A:33:HIS:NE2	0.55	2.70	3	1
1:A:-3:PHE:CE1	1:A:-2:LYS:O	0.55	2.59	8	2
1:A:22:LYS:CG	1:A:33:HIS:NE2	0.55	2.69	7	1
1:A:10:PHE:CE2	1:A:20:VAL:HG12	0.55	2.37	10	1
1:A:48:TYR:CZ	2:A:104:HEC:O2A	0.55	2.59	13	5
1:A:33:HIS:O	1:A:34:GLY:C	0.55	2.45	7	1
1:A:67:TYR:CD2	1:A:71:PRO:HA	0.55	2.36	15	1
1:A:6:GLY:O	1:A:97:TYR:CZ	0.55	2.60	18	1
1:A:46:TYR:CE1	2:A:104:HEC:O1D	0.54	2.60	5	1
1:A:22:LYS:CE	1:A:33:HIS:ND1	0.54	2.70	20	1
1:A:40:SER:HB3	1:A:57:VAL:HG12	0.54	1.77	12	6
1:A:66:GLU:OE2	1:A:74:TYR:CD1	0.54	2.59	12	1
1:A:67:TYR:CD1	1:A:71:PRO:HA	0.54	2.37	13	3
1:A:9:LEU:HB3	1:A:94:LEU:HD13	0.54	1.79	11	4
1:A:48:TYR:CZ	2:A:104:HEC:O1A	0.54	2.60	2	4
1:A:10:PHE:O	1:A:11:LYS:C	0.54	2.46	4	7
1:A:65:SER:CB	1:A:95:ILE:HD11	0.54	2.32	6	4
1:A:22:LYS:HD2	1:A:33:HIS:CE1	0.54	2.38	8	1
1:A:94:LEU:CD1	1:A:98:LEU:CD1	0.54	2.85	16	1
1:A:6:GLY:O	1:A:97:TYR:CG	0.54	2.60	18	1
1:A:9:LEU:HD11	1:A:85:LEU:CD2	0.54	2.30	20	1
1:A:10:PHE:HB2	1:A:94:LEU:HD21	0.54	1.80	9	1
1:A:30:PRO:CB	1:A:48:TYR:CE1	0.54	2.90	14	1
1:A:35:ILE:HD12	1:A:35:ILE:C	0.54	2.23	8	4
2:A:104:HEC:CHA	2:A:104:HEC:O2D	0.54	2.56	20	1
1:A:66:GLU:O	1:A:69:THR:HG22	0.54	2.03	11	5
1:A:59:TRP:CE3	1:A:64:MET:HE3	0.54	2.38	13	1
2:A:104:HEC:HMD1	2:A:104:HEC:CBD	0.54	2.33	16	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:49:THR:CG2	2:A:104:HEC:CGD	0.53	2.86	1	1
1:A:58:LEU:O	1:A:59:TRP:C	0.53	2.46	16	7
1:A:19:THR:CG2	1:A:31:ASN:ND2	0.53	2.72	4	3
1:A:35:ILE:HD12	1:A:36:PHE:CG	0.53	2.39	19	6
1:A:14:CYS:HG	2:A:104:HEC:CAB	0.53	1.96	13	1
1:A:30:PRO:CA	1:A:46:TYR:CE2	0.53	2.92	8	1
1:A:1:GLY:O	1:A:2:SER:CB	0.53	2.57	16	6
1:A:66:GLU:OE2	1:A:74:TYR:CG	0.53	2.62	12	1
1:A:20:VAL:CG2	1:A:101:ALA:CB	0.53	2.87	7	1
1:A:32:LEU:CD1	2:A:104:HEC:CHB	0.53	2.87	7	3
1:A:36:PHE:CZ	1:A:64:MET:HE2	0.53	2.39	19	3
1:A:20:VAL:CG2	1:A:101:ALA:O	0.53	2.57	10	2
1:A:57:VAL:O	1:A:59:TRP:N	0.52	2.41	5	7
1:A:26:HIS:HB3	1:A:46:TYR:CE1	0.52	2.39	14	1
1:A:32:LEU:O	1:A:34:GLY:N	0.52	2.43	7	2
1:A:2:SER:HG	1:A:93:ASP:N	0.52	2.02	12	1
1:A:6:GLY:O	1:A:97:TYR:CD1	0.52	2.62	18	1
1:A:33:HIS:O	1:A:102:SER:CB	0.52	2.58	16	2
1:A:30:PRO:HB3	1:A:48:TYR:CE1	0.52	2.38	14	1
1:A:26:HIS:CB	1:A:46:TYR:CE2	0.52	2.93	4	8
1:A:39:HIS:CD2	1:A:58:LEU:CD1	0.52	2.93	3	6
1:A:11:LYS:HA	1:A:15:LEU:HD22	0.52	1.82	8	3
1:A:82:PHE:HZ	1:A:85:LEU:HD23	0.52	1.64	9	1
1:A:6:GLY:O	1:A:97:TYR:CE1	0.52	2.63	18	1
1:A:26:HIS:HB3	1:A:46:TYR:CE2	0.52	2.40	1	15
1:A:30:PRO:HA	1:A:46:TYR:CE2	0.52	2.40	8	1
1:A:9:LEU:CD2	1:A:94:LEU:CD2	0.52	2.84	1	1
1:A:67:TYR:CE1	1:A:71:PRO:CB	0.52	2.93	1	1
1:A:48:TYR:CD2	1:A:53:ILE:CD1	0.52	2.93	2	1
1:A:-3:PHE:CD2	1:A:-3:PHE:O	0.52	2.63	5	2
1:A:57:VAL:CG1	1:A:57:VAL:O	0.52	2.58	1	2
1:A:26:HIS:HB3	1:A:46:TYR:CD2	0.52	2.40	6	9
1:A:48:TYR:OH	2:A:104:HEC:CGA	0.52	2.58	19	5
1:A:39:HIS:CG	1:A:58:LEU:HD12	0.52	2.40	14	1
1:A:35:ILE:O	1:A:37:GLY:N	0.51	2.43	13	3
1:A:-3:PHE:CE1	1:A:-1:ALA:HA	0.51	2.41	16	1
1:A:67:TYR:O	1:A:71:PRO:CA	0.51	2.59	14	4
1:A:30:PRO:CG	2:A:104:HEC:O1A	0.51	2.59	1	6
1:A:9:LEU:CD2	1:A:94:LEU:CD1	0.51	2.84	9	2
1:A:4:LYS:O	1:A:5:LYS:CB	0.51	2.58	12	2
1:A:10:PHE:CD1	1:A:14:CYS:HB2	0.51	2.41	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:5:LYS:O	1:A:5:LYS:CG	0.51	2.59	3	2
1:A:49:THR:HG23	2:A:104:HEC:CGD	0.51	2.36	1	1
1:A:59:TRP:CE3	1:A:64:MET:HE2	0.51	2.40	7	1
1:A:14:CYS:SG	2:A:104:HEC:HBB1	0.51	2.45	16	2
1:A:12:THR:O	1:A:13:ARG:HB2	0.51	2.06	14	8
1:A:59:TRP:CE3	1:A:64:MET:HG2	0.51	2.41	6	2
1:A:35:ILE:HD13	1:A:36:PHE:CD1	0.51	2.41	18	1
1:A:26:HIS:CG	1:A:46:TYR:CD2	0.51	2.98	8	9
1:A:78:THR:CG2	1:A:78:THR:O	0.51	2.59	10	12
1:A:35:ILE:HD12	1:A:36:PHE:CB	0.51	2.36	8	4
1:A:59:TRP:CZ2	2:A:104:HEC:HAA1	0.51	2.40	19	4
1:A:55:LYS:O	1:A:56:ASN:CB	0.51	2.58	8	4
1:A:41:GLY:HA3	1:A:48:TYR:CE2	0.50	2.41	1	4
1:A:85:LEU:O	1:A:86:LYS:CG	0.50	2.60	9	3
1:A:57:VAL:HG11	1:A:74:TYR:CE2	0.50	2.40	4	2
1:A:20:VAL:O	1:A:33:HIS:CB	0.50	2.59	9	1
1:A:25:PRO:O	1:A:26:HIS:CB	0.50	2.60	1	3
1:A:30:PRO:HA	1:A:46:TYR:CD2	0.50	2.40	5	5
1:A:98:LEU:O	1:A:102:SER:CB	0.50	2.60	12	3
1:A:35:ILE:HD13	1:A:35:ILE:N	0.50	2.21	7	1
1:A:51:ALA:HB3	1:A:77:GLY:O	0.50	2.06	5	4
1:A:30:PRO:HB2	1:A:43:ALA:HB2	0.50	1.81	20	1
1:A:35:ILE:HD12	1:A:36:PHE:CD2	0.50	2.41	9	1
2:A:104:HEC:HHA	2:A:104:HEC:CGD	0.50	2.37	20	1
1:A:-3:PHE:C	1:A:-3:PHE:CD1	0.50	2.85	7	1
1:A:9:LEU:O	1:A:13:ARG:CG	0.50	2.59	4	3
1:A:30:PRO:HB3	1:A:46:TYR:CD2	0.50	2.42	8	2
1:A:9:LEU:CD2	1:A:94:LEU:HD12	0.50	2.31	8	1
1:A:32:LEU:O	1:A:35:ILE:CG2	0.50	2.59	1	1
1:A:69:THR:HG22	1:A:70:ASN:N	0.50	2.20	1	6
1:A:48:TYR:CD2	1:A:53:ILE:HG13	0.50	2.41	2	3
1:A:17:CYS:SG	1:A:28:VAL:CG1	0.50	3.00	9	4
1:A:78:THR:HG23	1:A:78:THR:O	0.50	2.07	16	2
1:A:50:ASP:OD1	1:A:54:LYS:CE	0.50	2.60	19	1
1:A:9:LEU:HD11	1:A:90:ASP:HB3	0.50	1.84	2	1
1:A:94:LEU:HD21	2:A:104:HEC:HBB3	0.50	1.80	5	2
2:A:104:HEC:O1A	2:A:104:HEC:CBD	0.50	2.60	6	1
1:A:26:HIS:CB	1:A:46:TYR:CD1	0.50	2.94	20	1
1:A:40:SER:O	1:A:41:GLY:C	0.50	2.51	5	3
1:A:20:VAL:HG22	1:A:101:ALA:O	0.50	2.07	10	1
1:A:57:VAL:HG22	1:A:63:ASN:HD22	0.50	1.67	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:65:SER:OG	1:A:95:ILE:CD1	0.49	2.60	20	2
1:A:97:TYR:O	1:A:101:ALA:N	0.49	2.45	18	2
1:A:78:THR:OG1	2:A:104:HEC:CGD	0.49	2.59	18	2
1:A:97:TYR:CE1	1:A:101:ALA:CB	0.49	2.95	13	1
1:A:57:VAL:CG1	1:A:74:TYR:OH	0.49	2.60	18	2
1:A:51:ALA:CB	1:A:77:GLY:O	0.49	2.61	5	1
1:A:10:PHE:CE1	1:A:14:CYS:HB3	0.49	2.42	10	1
1:A:38:ARG:CB	1:A:42:GLN:CD	0.49	2.80	20	5
1:A:67:TYR:CE1	1:A:74:TYR:CG	0.49	3.00	11	1
1:A:9:LEU:HD21	1:A:85:LEU:CD1	0.49	2.38	16	1
1:A:49:THR:CG2	2:A:104:HEC:O2D	0.49	2.60	14	2
1:A:9:LEU:CD1	1:A:90:ASP:CG	0.49	2.81	10	2
1:A:9:LEU:O	1:A:12:THR:N	0.49	2.46	6	8
1:A:91:ARG:O	1:A:95:ILE:CG1	0.49	2.61	8	1
1:A:34:GLY:CA	1:A:102:SER:O	0.49	2.61	20	1
1:A:55:LYS:HD3	1:A:75:ILE:HD13	0.49	1.84	3	1
1:A:85:LEU:O	1:A:86:LYS:CB	0.49	2.60	15	4
1:A:61:GLU:CG	1:A:95:ILE:CG2	0.49	2.90	3	2
1:A:94:LEU:CD2	1:A:98:LEU:HD23	0.49	2.38	6	1
1:A:10:PHE:CE2	1:A:20:VAL:HB	0.49	2.43	19	4
1:A:36:PHE:CD2	1:A:99:LYS:O	0.49	2.66	15	1
2:A:104:HEC:HHA	2:A:104:HEC:HBD2	0.48	1.84	18	2
1:A:30:PRO:CG	2:A:104:HEC:O2A	0.48	2.61	18	6
1:A:59:TRP:CE3	1:A:64:MET:HG3	0.48	2.43	5	3
1:A:78:THR:O	1:A:78:THR:HG23	0.48	2.08	11	7
1:A:22:LYS:HD3	1:A:33:HIS:CE1	0.48	2.43	8	2
1:A:10:PHE:CD1	1:A:14:CYS:CB	0.48	2.96	10	2
1:A:41:GLY:HA2	1:A:48:TYR:CE2	0.48	2.44	14	13
1:A:-3:PHE:CZ	1:A:-1:ALA:HA	0.48	2.43	1	2
1:A:46:TYR:CE1	1:A:47:SER:O	0.48	2.67	2	3
1:A:55:LYS:NZ	1:A:74:TYR:CZ	0.48	2.81	10	1
1:A:98:LEU:O	1:A:99:LYS:C	0.48	2.52	15	2
1:A:98:LEU:HA	1:A:101:ALA:HB3	0.48	1.83	18	2
1:A:9:LEU:O	1:A:12:THR:O	0.48	2.31	2	14
1:A:35:ILE:O	1:A:36:PHE:CD1	0.48	2.65	7	1
1:A:94:LEU:HD22	1:A:98:LEU:HD23	0.48	1.85	6	3
1:A:85:LEU:HD11	1:A:91:ARG:CG	0.48	2.35	9	1
1:A:91:ARG:O	1:A:95:ILE:CD1	0.48	2.60	14	3
1:A:69:THR:CB	1:A:91:ARG:NH1	0.48	2.76	18	1
1:A:7:ALA:HA	1:A:97:TYR:CE1	0.48	2.44	5	3
1:A:36:PHE:CD2	1:A:60:ASP:O	0.48	2.66	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:57:VAL:CG1	1:A:67:TYR:OH	0.48	2.62	16	1
1:A:-3:PHE:CD1	1:A:-3:PHE:N	0.48	2.81	3	2
1:A:38:ARG:CB	1:A:42:GLN:NE2	0.48	2.76	19	2
1:A:36:PHE:CE1	1:A:64:MET:HE3	0.48	2.44	20	1
1:A:97:TYR:CE1	1:A:98:LEU:HD22	0.48	2.43	1	1
1:A:59:TRP:CE3	1:A:64:MET:CG	0.48	2.97	6	2
1:A:1:GLY:CA	1:A:96:THR:HG21	0.48	2.39	8	1
1:A:51:ALA:CB	1:A:76:PRO:O	0.48	2.62	20	2
1:A:82:PHE:CE2	1:A:84:GLY:HA2	0.47	2.44	18	4
1:A:-3:PHE:CD1	1:A:-3:PHE:C	0.47	2.87	3	2
1:A:-3:PHE:CG	1:A:-3:PHE:O	0.47	2.65	12	1
1:A:2:SER:O	1:A:3:ALA:HB3	0.47	2.09	14	1
1:A:69:THR:CG2	1:A:70:ASN:N	0.47	2.78	9	3
1:A:22:LYS:HG3	1:A:33:HIS:CD2	0.47	2.45	12	5
1:A:25:PRO:O	1:A:26:HIS:CD2	0.47	2.67	1	3
1:A:32:LEU:HB3	1:A:35:ILE:HD12	0.47	1.86	7	2
1:A:94:LEU:HD13	1:A:98:LEU:HD23	0.47	1.85	10	1
1:A:63:ASN:O	1:A:67:TYR:CE2	0.47	2.67	16	1
1:A:88:GLU:O	1:A:91:ARG:N	0.47	2.47	11	4
1:A:78:THR:O	1:A:78:THR:CG2	0.47	2.62	13	8
1:A:9:LEU:HD12	1:A:90:ASP:CG	0.47	2.29	10	1
1:A:33:HIS:CE1	1:A:102:SER:HB2	0.47	2.44	15	1
1:A:57:VAL:O	1:A:57:VAL:CG1	0.47	2.62	19	1
1:A:36:PHE:CE2	1:A:64:MET:HE1	0.47	2.45	20	1
1:A:67:TYR:O	1:A:71:PRO:N	0.47	2.47	16	16
1:A:57:VAL:HG11	1:A:74:TYR:HE2	0.47	1.69	5	2
1:A:79:LYS:C	2:A:104:HEC:HMD2	0.47	2.30	2	1
1:A:43:ALA:O	1:A:44:GLU:C	0.47	2.53	3	2
1:A:11:LYS:HA	1:A:15:LEU:HD12	0.47	1.87	5	1
1:A:94:LEU:HD11	2:A:104:HEC:HBB1	0.47	1.86	6	1
1:A:22:LYS:NZ	1:A:33:HIS:NE2	0.47	2.60	19	2
1:A:39:HIS:CD2	1:A:58:LEU:N	0.47	2.83	18	1
1:A:48:TYR:CB	1:A:53:ILE:CD1	0.47	2.92	7	3
1:A:63:ASN:O	1:A:66:GLU:N	0.47	2.48	12	3
1:A:87:LYS:O	1:A:88:GLU:C	0.47	2.54	18	9
1:A:41:GLY:CA	1:A:48:TYR:CE2	0.47	2.98	5	1
1:A:36:PHE:CE1	1:A:64:MET:HG3	0.47	2.45	10	2
1:A:9:LEU:CD2	1:A:94:LEU:HD23	0.47	2.33	18	1
1:A:10:PHE:CZ	1:A:20:VAL:HB	0.47	2.45	7	3
1:A:82:PHE:CZ	1:A:84:GLY:HA2	0.47	2.45	9	1
1:A:-3:PHE:CE2	1:A:-2:LYS:O	0.47	2.68	20	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:56:ASN:O	1:A:58:LEU:N	0.47	2.48	7	3
1:A:-3:PHE:CD2	1:A:-1:ALA:HB2	0.47	2.45	16	1
1:A:67:TYR:CE1	1:A:75:ILE:HG13	0.46	2.45	8	2
1:A:67:TYR:OH	2:A:104:HEC:C2A	0.46	2.63	2	1
1:A:9:LEU:HD22	1:A:94:LEU:HD21	0.46	1.85	20	1
1:A:35:ILE:HG12	1:A:36:PHE:CD2	0.46	2.44	12	2
1:A:74:TYR:CD2	1:A:75:ILE:HG13	0.46	2.45	2	1
1:A:52:ASN:O	1:A:55:LYS:N	0.46	2.48	20	13
1:A:94:LEU:O	1:A:97:TYR:N	0.46	2.49	15	9
1:A:98:LEU:O	1:A:100:LYS:N	0.46	2.49	15	2
1:A:69:THR:O	1:A:84:GLY:CA	0.46	2.64	14	2
1:A:100:LYS:O	1:A:101:ALA:CB	0.46	2.64	15	1
1:A:30:PRO:CG	1:A:48:TYR:OH	0.46	2.64	10	1
1:A:67:TYR:HE1	1:A:75:ILE:HD12	0.46	1.67	12	1
1:A:63:ASN:OD1	1:A:64:MET:N	0.46	2.48	16	2
1:A:93:ASP:OD1	1:A:93:ASP:N	0.46	2.49	20	1
1:A:93:ASP:O	1:A:96:THR:N	0.46	2.48	16	4
1:A:21:GLU:O	1:A:31:ASN:ND2	0.46	2.49	9	11
1:A:48:TYR:HB3	1:A:53:ILE:HD12	0.46	1.87	7	3
1:A:32:LEU:O	1:A:33:HIS:O	0.46	2.33	10	3
1:A:38:ARG:NE	1:A:42:GLN:OE1	0.46	2.49	3	1
1:A:62:ASN:O	1:A:65:SER:CB	0.46	2.64	11	1
1:A:64:MET:HG2	1:A:68:LEU:HD11	0.46	1.88	13	1
2:A:104:HEC:CHA	2:A:104:HEC:CGD	0.46	2.94	20	1
1:A:90:ASP:O	1:A:93:ASP:N	0.46	2.49	16	3
1:A:62:ASN:O	1:A:65:SER:N	0.46	2.49	3	2
1:A:89:LYS:O	1:A:92:ASN:N	0.46	2.49	3	2
1:A:51:ALA:HB1	1:A:75:ILE:HG22	0.46	1.88	19	2
1:A:22:LYS:HG3	1:A:33:HIS:CG	0.46	2.46	11	2
1:A:48:TYR:HH	2:A:104:HEC:CGA	0.46	2.23	12	1
1:A:-3:PHE:O	1:A:-3:PHE:CG	0.46	2.69	17	2
1:A:48:TYR:CB	1:A:53:ILE:HD11	0.46	2.41	14	1
1:A:97:TYR:O	1:A:100:LYS:N	0.46	2.49	12	6
1:A:39:HIS:ND1	1:A:56:ASN:OD1	0.46	2.49	2	2
1:A:37:GLY:N	1:A:59:TRP:O	0.45	2.49	2	6
1:A:58:LEU:O	1:A:63:ASN:ND2	0.45	2.49	6	3
1:A:19:THR:O	1:A:31:ASN:ND2	0.45	2.50	20	6
1:A:22:LYS:HG2	1:A:33:HIS:CD2	0.45	2.47	7	1
1:A:10:PHE:CE2	1:A:20:VAL:CG1	0.45	2.99	10	1
1:A:57:VAL:CG2	1:A:63:ASN:OD1	0.45	2.64	12	1
1:A:18:HIS:CG	1:A:32:LEU:CD1	0.45	3.00	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:30:PRO:HG3	1:A:48:TYR:CZ	0.45	2.46	10	1
1:A:67:TYR:O	1:A:70:ASN:N	0.45	2.49	16	4
1:A:2:SER:OG	1:A:3:ALA:N	0.45	2.48	2	1
1:A:1:GLY:N	1:A:96:THR:OG1	0.45	2.50	10	4
1:A:78:THR:C	1:A:80:MET:N	0.45	2.70	16	6
1:A:86:LYS:O	1:A:88:GLU:N	0.45	2.50	14	3
1:A:33:HIS:O	1:A:35:ILE:N	0.45	2.49	7	1
1:A:97:TYR:CE2	1:A:101:ALA:HB2	0.45	2.47	18	3
1:A:36:PHE:CE2	1:A:61:GLU:HA	0.45	2.46	10	1
1:A:82:PHE:CE1	1:A:83:GLY:O	0.45	2.69	12	1
1:A:39:HIS:CG	1:A:58:LEU:CD1	0.45	2.99	14	1
1:A:39:HIS:NE2	1:A:56:ASN:O	0.45	2.50	14	1
1:A:69:THR:OG1	1:A:91:ARG:NH1	0.45	2.50	1	4
1:A:10:PHE:CE1	1:A:98:LEU:HD21	0.45	2.46	20	1
1:A:13:ARG:O	2:A:104:HEC:CMC	0.45	2.63	1	2
1:A:39:HIS:O	1:A:40:SER:C	0.45	2.55	5	11
1:A:67:TYR:CZ	1:A:71:PRO:HB3	0.45	2.45	1	4
1:A:94:LEU:HA	1:A:97:TYR:CE2	0.45	2.47	1	1
1:A:12:THR:O	1:A:13:ARG:CG	0.45	2.65	16	4
1:A:63:ASN:CG	1:A:64:MET:N	0.45	2.70	7	5
1:A:100:LYS:O	1:A:103:GLU:N	0.45	2.50	8	7
1:A:52:ASN:ND2	2:A:104:HEC:O2A	0.45	2.50	8	2
1:A:67:TYR:OH	2:A:104:HEC:CHA	0.45	2.64	8	1
1:A:2:SER:OG	1:A:93:ASP:N	0.45	2.50	12	1
1:A:64:MET:O	1:A:67:TYR:N	0.45	2.50	18	2
1:A:52:ASN:OD1	1:A:53:ILE:N	0.45	2.49	20	3
1:A:88:GLU:O	1:A:92:ASN:ND2	0.45	2.50	13	3
1:A:7:ALA:CB	1:A:97:TYR:CZ	0.45	3.00	12	1
1:A:87:LYS:O	1:A:89:LYS:N	0.45	2.50	16	2
1:A:5:LYS:NZ	1:A:93:ASP:OD2	0.45	2.50	14	1
1:A:-3:PHE:CG	1:A:61:GLU:OE2	0.45	2.70	19	1
1:A:36:PHE:CE1	1:A:64:MET:HE1	0.45	2.46	19	1
1:A:82:PHE:CD1	2:A:104:HEC:HMC2	0.45	2.46	4	1
1:A:93:ASP:N	1:A:93:ASP:OD1	0.45	2.49	12	1
1:A:9:LEU:HD11	1:A:90:ASP:OD1	0.45	2.10	14	1
1:A:65:SER:OG	1:A:91:ARG:NH1	0.45	2.50	20	1
1:A:10:PHE:C	1:A:12:THR:N	0.45	2.70	18	3
1:A:30:PRO:CG	1:A:48:TYR:CZ	0.45	3.00	20	1
1:A:34:GLY:N	1:A:102:SER:O	0.45	2.50	20	1
1:A:10:PHE:CE1	1:A:15:LEU:CD1	0.45	2.99	2	2
1:A:34:GLY:O	1:A:35:ILE:O	0.45	2.35	6	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:-2:LYS:O	1:A:-1:ALA:C	0.45	2.55	16	1
1:A:67:TYR:O	1:A:71:PRO:CB	0.45	2.65	4	2
1:A:19:THR:OG1	1:A:31:ASN:ND2	0.45	2.50	5	4
1:A:65:SER:O	1:A:91:ARG:NH1	0.45	2.50	9	1
1:A:7:ALA:HA	1:A:97:TYR:CE2	0.45	2.47	12	1
1:A:35:ILE:CG1	1:A:36:PHE:N	0.45	2.80	16	2
1:A:67:TYR:CE1	1:A:71:PRO:HB3	0.44	2.46	17	2
1:A:95:ILE:O	1:A:96:THR:C	0.44	2.56	1	2
1:A:94:LEU:C	1:A:96:THR:N	0.44	2.70	7	7
1:A:15:LEU:O	1:A:17:CYS:N	0.44	2.51	12	2
1:A:22:LYS:CD	1:A:33:HIS:ND1	0.44	2.80	11	1
1:A:36:PHE:CE1	1:A:61:GLU:HA	0.44	2.48	15	2
2:A:104:HEC:O2D	2:A:104:HEC:C4D	0.44	2.65	20	1
1:A:7:ALA:C	1:A:9:LEU:N	0.44	2.71	18	3
1:A:1:GLY:O	1:A:96:THR:HG21	0.44	2.12	8	1
1:A:67:TYR:CZ	1:A:74:TYR:CE2	0.44	3.05	11	1
1:A:94:LEU:CD2	1:A:98:LEU:CD1	0.44	2.93	9	1
1:A:64:MET:SD	1:A:68:LEU:CD1	0.44	3.01	12	1
1:A:93:ASP:O	1:A:96:THR:CB	0.44	2.66	20	1
1:A:39:HIS:O	1:A:40:SER:O	0.44	2.36	2	6
1:A:88:GLU:OE2	1:A:91:ARG:NH2	0.44	2.51	9	1
1:A:-3:PHE:CE1	1:A:-1:ALA:HB2	0.44	2.44	16	1
1:A:7:ALA:O	1:A:9:LEU:N	0.44	2.50	18	2
1:A:34:GLY:N	1:A:102:SER:OG	0.44	2.51	11	1
1:A:65:SER:O	1:A:91:ARG:NH2	0.44	2.50	14	1
1:A:99:LYS:O	1:A:103:GLU:N	0.44	2.50	14	1
1:A:36:PHE:CE2	1:A:99:LYS:HA	0.44	2.47	10	1
1:A:57:VAL:HG23	1:A:63:ASN:OD1	0.44	2.12	12	1
1:A:-3:PHE:CD1	1:A:-1:ALA:N	0.44	2.86	16	1
1:A:39:HIS:NE2	1:A:56:ASN:OD1	0.44	2.50	20	1
1:A:15:LEU:C	1:A:17:CYS:N	0.44	2.71	12	5
1:A:36:PHE:CE1	1:A:60:ASP:C	0.44	2.90	9	2
1:A:10:PHE:O	1:A:12:THR:N	0.44	2.51	18	3
1:A:52:ASN:O	1:A:53:ILE:C	0.44	2.57	20	10
1:A:38:ARG:CD	1:A:42:GLN:OE1	0.44	2.66	3	1
1:A:67:TYR:CD1	1:A:75:ILE:HG13	0.44	2.47	8	1
1:A:48:TYR:HB2	1:A:53:ILE:HD11	0.44	1.87	10	1
1:A:98:LEU:C	1:A:100:LYS:N	0.44	2.71	10	2
1:A:39:HIS:CG	1:A:56:ASN:OD1	0.43	2.71	2	1
1:A:61:GLU:HG3	1:A:95:ILE:CG2	0.43	2.44	3	3
1:A:99:LYS:O	1:A:103:GLU:CB	0.43	2.66	19	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:52:ASN:C	1:A:54:LYS:N	0.43	2.71	18	8
1:A:80:MET:HB2	2:A:104:HEC:C4D	0.43	2.43	14	12
1:A:31:ASN:OD1	1:A:33:HIS:ND1	0.43	2.49	4	1
1:A:67:TYR:O	1:A:71:PRO:CD	0.43	2.66	16	2
1:A:95:ILE:O	1:A:99:LYS:CB	0.43	2.66	20	1
1:A:80:MET:HA	2:A:104:HEC:C2D	0.43	2.44	2	3
1:A:87:LYS:C	1:A:89:LYS:N	0.43	2.71	16	2
1:A:48:TYR:HB2	1:A:53:ILE:HD12	0.43	1.90	5	2
1:A:94:LEU:O	1:A:96:THR:N	0.43	2.51	8	5
1:A:24:GLY:O	1:A:31:ASN:ND2	0.43	2.51	12	1
1:A:7:ALA:HA	1:A:97:TYR:CZ	0.43	2.49	16	1
1:A:93:ASP:O	1:A:97:TYR:N	0.43	2.47	16	2
2:A:104:HEC:HMD1	2:A:104:HEC:HBD1	0.43	1.90	16	1
1:A:40:SER:CB	1:A:57:VAL:HG12	0.43	2.43	20	1
1:A:2:SER:O	1:A:3:ALA:C	0.43	2.56	12	4
1:A:80:MET:HE3	2:A:104:HEC:CHC	0.43	2.44	8	2
1:A:86:LYS:O	1:A:91:ARG:NH2	0.43	2.51	8	2
1:A:18:HIS:CD2	1:A:32:LEU:HD11	0.43	2.48	15	3
1:A:18:HIS:CE1	1:A:32:LEU:HD21	0.43	2.49	13	1
1:A:66:GLU:O	1:A:69:THR:CG2	0.43	2.66	16	1
1:A:94:LEU:HA	1:A:97:TYR:CD2	0.43	2.49	1	1
1:A:32:LEU:O	1:A:33:HIS:C	0.43	2.57	9	4
1:A:9:LEU:O	1:A:10:PHE:C	0.43	2.57	1	5
1:A:52:ASN:ND2	2:A:104:HEC:O1A	0.43	2.51	4	2
1:A:51:ALA:CB	1:A:77:GLY:C	0.43	2.87	5	1
1:A:102:SER:O	1:A:103:GLU:C	0.43	2.57	18	1
1:A:2:SER:OG	1:A:96:THR:HG22	0.43	2.14	2	1
1:A:35:ILE:O	1:A:36:PHE:O	0.42	2.37	10	7
1:A:39:HIS:CE1	1:A:58:LEU:HD12	0.42	2.49	5	1
1:A:48:TYR:CD2	2:A:104:HEC:O2A	0.42	2.72	7	1
1:A:9:LEU:C	1:A:11:LYS:N	0.42	2.73	19	2
1:A:2:SER:O	1:A:5:LYS:CG	0.42	2.67	20	1
1:A:21:GLU:O	1:A:33:HIS:ND1	0.42	2.52	4	1
1:A:19:THR:HG23	1:A:31:ASN:CG	0.42	2.34	8	3
1:A:88:GLU:O	1:A:92:ASN:CB	0.42	2.67	7	1
1:A:18:HIS:ND1	1:A:30:PRO:O	0.42	2.49	20	1
1:A:4:LYS:O	1:A:5:LYS:C	0.42	2.57	3	1
1:A:68:LEU:O	1:A:69:THR:CB	0.42	2.67	9	1
1:A:22:LYS:O	1:A:24:GLY:N	0.42	2.50	13	1
1:A:10:PHE:O	1:A:15:LEU:HD13	0.42	2.14	17	1
1:A:57:VAL:HG13	1:A:59:TRP:CE2	0.42	2.50	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:55:LYS:NZ	1:A:74:TYR:CE2	0.42	2.87	10	1
1:A:61:GLU:HG2	1:A:95:ILE:HG23	0.42	1.90	14	1
1:A:52:ASN:O	1:A:54:LYS:N	0.42	2.52	18	2
1:A:34:GLY:O	1:A:35:ILE:C	0.42	2.57	7	1
1:A:57:VAL:HG23	1:A:74:TYR:OH	0.42	2.15	7	1
1:A:63:ASN:C	1:A:65:SER:N	0.42	2.73	11	1
1:A:26:HIS:CB	1:A:46:TYR:CE1	0.42	3.03	14	1
1:A:2:SER:O	1:A:4:LYS:N	0.42	2.49	16	1
1:A:65:SER:O	1:A:66:GLU:C	0.42	2.57	16	1
1:A:78:THR:O	1:A:80:MET:N	0.42	2.52	16	2
1:A:64:MET:SD	2:A:104:HEC:CMA	0.42	3.08	17	1
1:A:55:LYS:NZ	1:A:74:TYR:O	0.42	2.50	18	1
1:A:35:ILE:HD13	1:A:64:MET:CE	0.42	2.44	8	2
1:A:35:ILE:HD13	1:A:64:MET:HE3	0.42	1.92	8	1
1:A:10:PHE:CD1	1:A:15:LEU:HD13	0.42	2.50	1	1
1:A:80:MET:HA	2:A:104:HEC:C1D	0.42	2.45	7	2
1:A:59:TRP:CZ3	1:A:67:TYR:CD2	0.42	3.08	8	1
1:A:3:ALA:O	1:A:5:LYS:N	0.42	2.51	11	1
1:A:13:ARG:O	1:A:14:CYS:SG	0.42	2.78	14	5
1:A:39:HIS:CD2	1:A:56:ASN:OD1	0.42	2.73	8	1
1:A:39:HIS:ND1	1:A:58:LEU:CD1	0.42	2.83	16	1
1:A:23:GLY:O	1:A:24:GLY:C	0.42	2.58	1	1
1:A:-2:LYS:CB	1:A:92:ASN:ND2	0.42	2.82	10	1
1:A:64:MET:CE	1:A:67:TYR:CD2	0.42	3.03	10	1
1:A:-1:ALA:O	1:A:1:GLY:O	0.42	2.38	12	1
1:A:7:ALA:HB2	1:A:97:TYR:CZ	0.42	2.50	12	1
1:A:20:VAL:O	1:A:21:GLU:O	0.41	2.38	15	3
1:A:82:PHE:HD2	2:A:104:HEC:HHC	0.41	1.75	12	1
1:A:65:SER:HA	1:A:95:ILE:HD11	0.41	1.93	2	1
1:A:94:LEU:O	1:A:95:ILE:C	0.41	2.59	16	7
1:A:30:PRO:O	1:A:31:ASN:C	0.41	2.59	10	1
1:A:63:ASN:O	1:A:65:SER:N	0.41	2.52	11	1
1:A:80:MET:CB	2:A:104:HEC:O1D	0.41	2.68	13	1
1:A:22:LYS:CE	1:A:33:HIS:NE2	0.41	2.83	15	2
1:A:33:HIS:O	1:A:101:ALA:HB3	0.41	2.15	15	1
1:A:62:ASN:C	1:A:64:MET:N	0.41	2.74	15	1
1:A:12:THR:CG2	1:A:13:ARG:HG2	0.41	2.45	3	5
1:A:36:PHE:CE2	1:A:99:LYS:HB2	0.41	2.50	7	1
1:A:36:PHE:CE2	1:A:64:MET:CE	0.41	3.03	9	1
1:A:80:MET:HG3	2:A:104:HEC:C4C	0.41	2.45	10	1
1:A:64:MET:HE1	2:A:104:HEC:HMA1	0.41	1.92	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:41:GLY:N	2:A:104:HEC:O2A	0.41	2.53	16	1
1:A:-3:PHE:CD2	1:A:61:GLU:CD	0.41	2.94	17	1
1:A:73:LYS:O	1:A:74:TYR:C	0.41	2.59	7	2
1:A:9:LEU:CD1	1:A:90:ASP:HB3	0.41	2.46	10	1
1:A:61:GLU:O	1:A:95:ILE:HG21	0.41	2.15	16	1
1:A:47:SER:O	1:A:48:TYR:O	0.41	2.38	20	1
1:A:-2:LYS:O	1:A:-1:ALA:O	0.41	2.39	1	2
1:A:14:CYS:C	1:A:16:GLN:N	0.41	2.73	2	1
1:A:19:THR:O	1:A:31:ASN:CG	0.41	2.58	7	3
1:A:19:THR:O	1:A:31:ASN:OD1	0.41	2.39	2	3
1:A:9:LEU:O	1:A:11:LYS:N	0.41	2.54	8	1
1:A:39:HIS:ND1	1:A:58:LEU:HD12	0.41	2.30	11	2
1:A:48:TYR:HB3	1:A:53:ILE:HD11	0.41	1.92	14	1
1:A:57:VAL:HG11	1:A:74:TYR:OH	0.41	2.15	17	1
1:A:62:ASN:N	1:A:62:ASN:ND2	0.41	2.68	18	1
1:A:1:GLY:C	1:A:2:SER:HG	0.41	2.19	2	1
1:A:52:ASN:CG	2:A:104:HEC:O2A	0.41	2.59	6	3
1:A:59:TRP:CE3	1:A:63:ASN:ND2	0.41	2.89	2	1
1:A:86:LYS:O	1:A:91:ARG:NE	0.41	2.53	6	1
1:A:21:GLU:O	1:A:31:ASN:OD1	0.41	2.38	11	2
1:A:2:SER:O	1:A:93:ASP:CG	0.41	2.59	12	1
1:A:82:PHE:CD2	2:A:104:HEC:HHC	0.41	2.51	12	1
1:A:87:LYS:N	1:A:91:ARG:HE	0.41	2.14	13	1
1:A:64:MET:CE	2:A:104:HEC:HMA1	0.41	2.46	14	1
1:A:34:GLY:O	1:A:42:GLN:OE1	0.41	2.39	18	1
1:A:69:THR:OG1	1:A:91:ARG:CZ	0.41	2.69	1	1
1:A:93:ASP:O	1:A:94:LEU:C	0.41	2.59	2	1
1:A:62:ASN:O	1:A:63:ASN:C	0.41	2.59	15	3
1:A:100:LYS:O	1:A:101:ALA:C	0.41	2.60	8	1
1:A:57:VAL:HG21	1:A:59:TRP:CZ3	0.41	2.51	15	1
1:A:88:GLU:O	1:A:92:ASN:N	0.41	2.52	18	1
1:A:30:PRO:CB	1:A:43:ALA:HB2	0.41	2.45	20	1
1:A:10:PHE:CE1	1:A:15:LEU:HD13	0.40	2.51	1	1
1:A:3:ALA:O	1:A:4:LYS:C	0.40	2.59	8	1
1:A:22:LYS:HB2	1:A:33:HIS:CE1	0.40	2.51	13	1
1:A:63:ASN:O	1:A:64:MET:C	0.40	2.58	14	1
1:A:64:MET:O	1:A:68:LEU:N	0.40	2.49	14	1
1:A:2:SER:OG	1:A:93:ASP:CG	0.40	2.59	16	1
1:A:35:ILE:HD12	1:A:36:PHE:CD1	0.40	2.46	18	1
1:A:64:MET:HG2	1:A:68:LEU:CD1	0.40	2.46	18	1
1:A:57:VAL:HG11	1:A:59:TRP:CZ2	0.40	2.50	19	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:10:PHE:HB2	1:A:97:TYR:CE2	0.40	2.51	6	1
1:A:1:GLY:O	1:A:2:SER:C	0.40	2.60	14	1
1:A:67:TYR:O	1:A:68:LEU:C	0.40	2.60	15	2
1:A:67:TYR:O	1:A:70:ASN:C	0.40	2.60	20	2
1:A:38:ARG:HB2	1:A:42:GLN:NE2	0.40	2.32	20	1
1:A:51:ALA:HB2	1:A:76:PRO:O	0.40	2.16	20	1
1:A:32:LEU:HD11	2:A:104:HEC:HHB	0.40	1.88	1	1
1:A:43:ALA:O	1:A:44:GLU:O	0.40	2.39	3	1
1:A:22:LYS:HG3	1:A:33:HIS:NE2	0.40	2.31	7	1
1:A:69:THR:O	1:A:83:GLY:O	0.40	2.40	15	1
1:A:78:THR:O	1:A:79:LYS:C	0.40	2.59	16	1
1:A:-3:PHE:CE2	1:A:-1:ALA:HA	0.40	2.52	3	1
1:A:94:LEU:HG	1:A:98:LEU:HD13	0.40	1.93	11	1
1:A:94:LEU:HD23	1:A:94:LEU:HA	0.40	1.79	15	1
1:A:9:LEU:HD21	1:A:85:LEU:HD11	0.40	1.93	16	1
1:A:90:ASP:O	1:A:91:ARG:C	0.40	2.60	17	1
1:A:33:HIS:CG	1:A:102:SER:HB3	0.40	2.51	18	1
1:A:88:GLU:O	1:A:89:LYS:C	0.40	2.60	18	1
1:A:3:ALA:O	1:A:4:LYS:O	0.40	2.40	1	1
1:A:25:PRO:C	1:A:26:HIS:CG	0.40	2.93	1	1
1:A:94:LEU:CD1	1:A:98:LEU:HD23	0.40	2.46	1	1
1:A:44:GLU:O	1:A:45:GLY:C	0.40	2.60	5	1
1:A:70:ASN:OD1	1:A:73:LYS:CB	0.40	2.70	7	1
1:A:2:SER:O	1:A:3:ALA:O	0.40	2.40	8	1
1:A:67:TYR:CE1	1:A:74:TYR:CZ	0.40	3.09	11	1
1:A:13:ARG:C	1:A:14:CYS:SG	0.40	3.00	13	1
1:A:88:GLU:HG2	1:A:91:ARG:CZ	0.40	2.46	15	1
1:A:89:LYS:O	1:A:93:ASP:OD1	0.40	2.40	16	1

6.3 Torsion angles (i)

6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	104/108 (96%)	61±3 (59±3%)	31±4 (30±4%)	12±2 (11±2%)	1 8

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	2080/2160 (96%)	1219 (59%)	626 (30%)	235 (11%)	1 8

All 46 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	36	PHE	18
1	A	61	GLU	18
1	A	80	MET	16
1	A	35	ILE	14
1	A	5	LYS	14
1	A	-3	PHE	12
1	A	33	HIS	12
1	A	21	GLU	11
1	A	58	LEU	9
1	A	2	SER	8
1	A	-1	ALA	7
1	A	87	LYS	7
1	A	40	SER	6
1	A	13	ARG	6
1	A	52	ASN	6
1	A	86	LYS	6
1	A	4	LYS	5
1	A	26	HIS	5
1	A	57	VAL	5
1	A	34	GLY	5
1	A	3	ALA	4
1	A	1	GLY	4
1	A	11	LYS	3
1	A	77	GLY	3
1	A	85	LEU	3
1	A	41	GLY	2
1	A	94	LEU	2
1	A	44	GLU	2
1	A	59	TRP	2
1	A	8	THR	2
1	A	56	ASN	2
1	A	16	GLN	2
1	A	24	GLY	1
1	A	60	ASP	1
1	A	9	LEU	1
1	A	47	SER	1

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Mol	Chain	Res	Type	Models (Total)
1	A	14	CYS	1
1	A	31	ASN	1
1	A	84	GLY	1
1	A	63	ASN	1
1	A	83	GLY	1
1	A	88	GLU	1
1	A	38	ARG	1
1	A	45	GLY	1
1	A	102	SER	1
1	A	48	TYR	1

6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	86/88 (98%)	63±3 (73±3%)	23±3 (27±3%)	2 21
All	All	1720/1760 (98%)	1255 (73%)	465 (27%)	2 21

All 58 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	8	THR	20
1	A	50	ASP	20
1	A	96	THR	20
1	A	98	LEU	20
1	A	9	LEU	16
1	A	55	LYS	16
1	A	56	ASN	15
1	A	35	ILE	14
1	A	86	LYS	14
1	A	-2	LYS	13
1	A	4	LYS	12
1	A	5	LYS	11
1	A	94	LEU	11
1	A	33	HIS	11
1	A	73	LYS	10

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Mol	Chain	Res	Type	Models (Total)
1	A	89	LYS	10
1	A	14	CYS	9
1	A	69	THR	9
1	A	88	GLU	9
1	A	100	LYS	9
1	A	31	ASN	9
1	A	2	SER	9
1	A	11	LYS	9
1	A	21	GLU	9
1	A	39	HIS	9
1	A	64	MET	8
1	A	12	THR	8
1	A	62	ASN	8
1	A	79	LYS	8
1	A	66	GLU	7
1	A	99	LYS	7
1	A	19	THR	7
1	A	57	VAL	7
1	A	46	TYR	7
1	A	-3	PHE	6
1	A	13	ARG	6
1	A	48	TYR	6
1	A	22	LYS	6
1	A	60	ASP	5
1	A	52	ASN	4
1	A	87	LYS	4
1	A	103	GLU	4
1	A	65	SER	4
1	A	61	GLU	4
1	A	93	ASP	4
1	A	10	PHE	3
1	A	78	THR	3
1	A	102	SER	3
1	A	27	LYS	3
1	A	58	LEU	3
1	A	90	ASP	3
1	A	15	LEU	3
1	A	44	GLU	3
1	A	70	ASN	2
1	A	74	TYR	2
1	A	97	TYR	1
1	A	85	LEU	1

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Mol	Chain	Res	Type	Models (Total)
1	A	63	ASN	1

6.3.3 RNA (i)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds 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 is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
1	M3L	A	72	1	10,11,12	0.69±0.01	0±0 (0±0%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of 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 angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
1	M3L	A	72	1	9,14,16	1.03±0.00	0±0 (0±0%)

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
1	M3L	A	72	1	-	0±0,9,10,12	-

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates (i)

There are no monosaccharides in this entry.

6.6 Ligand geometry (i)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds 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 is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
2	HEC	A	104	1	26,50,50	2.13±0.00	8±0 (31±1%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of 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 angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
2	HEC	A	104	1	18,82,82	2.12±0.00	6±0 (33±0%)

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	HEC	A	104	1	-	0±0,6,54,54	-

All unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
2	A	104	HEC	C3C-C2C	5.23	1.35	1.40	10	20
2	A	104	HEC	C3B-C2B	4.54	1.36	1.40	13	20
2	A	104	HEC	CMD-C2D	4.10	1.60	1.51	10	20
2	A	104	HEC	CMA-C3A	2.76	1.58	1.51	8	20
2	A	104	HEC	C1A-C2A	2.75	1.48	1.42	15	20
2	A	104	HEC	CMB-C2B	2.59	1.57	1.51	12	20
2	A	104	HEC	CAD-C3D	2.32	1.55	1.52	14	20
2	A	104	HEC	C1D-CHD	2.20	1.34	1.41	13	20
2	A	104	HEC	C1C-CHC	2.03	1.46	1.41	13	4
2	A	104	HEC	C3C-C4C	2.01	1.46	1.43	3	1

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})	Models	
								Worst	Total
2	A	104	HEC	CMB-C2B-C1B	4.71	121.23	128.46	13	20
2	A	104	HEC	C4B-C3B-C2B	3.32	102.77	106.35	3	20
2	A	104	HEC	CAD-CBD-CGD	3.32	118.24	112.67	12	20
2	A	104	HEC	CAA-CBA-CGA	3.19	118.03	112.67	4	20
2	A	104	HEC	C4C-C3C-C2C	2.42	108.97	106.35	15	20
2	A	104	HEC	CMC-C2C-C3C	2.38	128.61	125.82	17	20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

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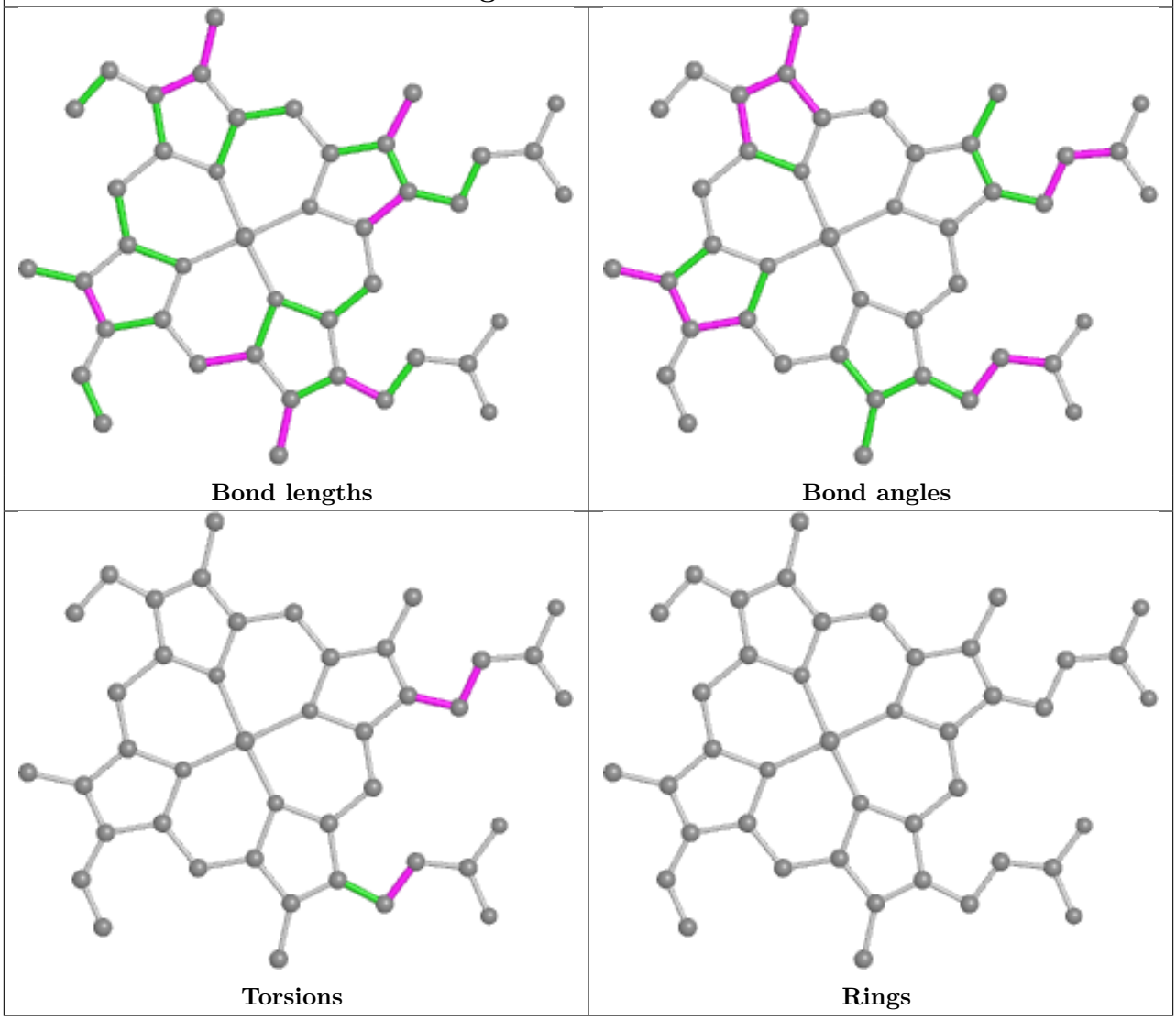

6.7 Other polymers (i)

There are no such molecules in this entry.

6.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

7 Chemical shift validation (i)

No chemical shift data were provided