



Full wwPDB NMR Structure Validation Report ⓘ

Aug 4, 2026 – 08:41 AM EDT

PDB ID : 2K78 / pdb_00002k78
Title : Solution Structure of the IsdC NEAT domain bound to Zinc Protoporphyrin
Authors : Villareal, V.A.; Pilpa, R.M.; Robson, S.A.; Fadeev, E.A.; Clubb, R.T.
Deposited on : 2008-08-06

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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

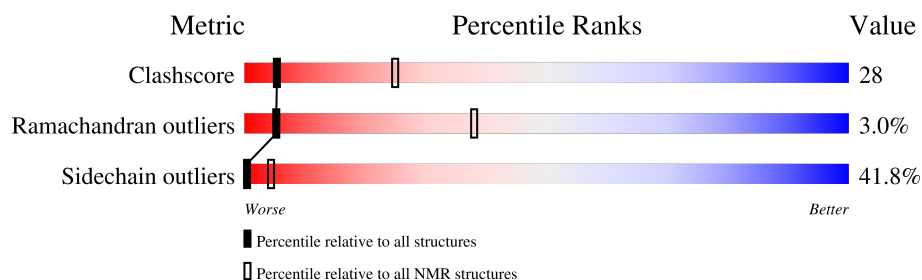
1 Overall quality at a glance

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	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	147	24% 39% 15% 7% 14%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA and RNA chains that are outliers for geometric criteria:

Mol	Chain	Compound	Res	Total models with violations	
				Chirality	Geometry
2	A	ZNH	151	30	-

2 Ensemble composition and analysis

This entry contains 30 models. Model 4 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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:30-A:145 (116)	0.66	4

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

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

The models can be grouped into 5 clusters. No single-model clusters were found.

Cluster number	Models
1	1, 4, 5, 9, 11, 17, 25, 27, 28, 29, 30
2	2, 3, 6, 13, 14, 18, 20
3	7, 15, 16, 21, 22, 26
4	8, 12, 23, 24
5	10, 19

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2057 atoms, of which 1004 are hydrogens and 0 are deuteriums.

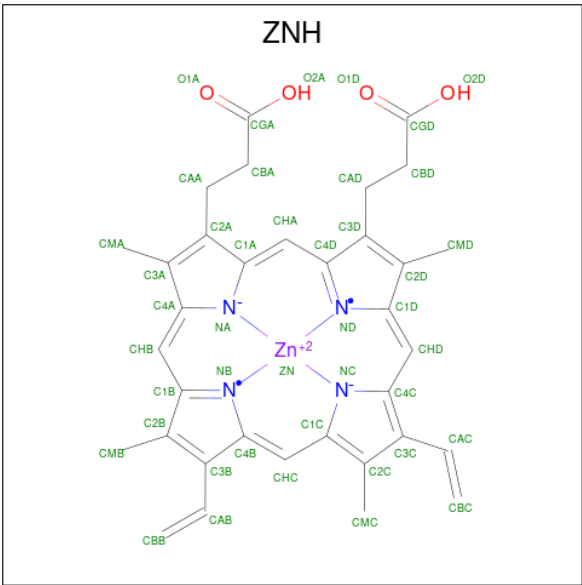
- Molecule 1 is a protein called Iron-regulated surface determinant protein C.

Mol	Chain	Residues	Atoms						Trace
1	A	126	Total	C	H	N	O	S	0
			1984	639	974	171	199	1	

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4	MET	-	expression tag	UNP Q7A151
A	5	GLY	-	expression tag	UNP Q7A151
A	6	SER	-	expression tag	UNP Q7A151
A	7	SER	-	expression tag	UNP Q7A151
A	8	HIS	-	expression tag	UNP Q7A151
A	9	HIS	-	expression tag	UNP Q7A151
A	10	HIS	-	expression tag	UNP Q7A151
A	11	HIS	-	expression tag	UNP Q7A151
A	12	HIS	-	expression tag	UNP Q7A151
A	13	HIS	-	expression tag	UNP Q7A151
A	14	SER	-	expression tag	UNP Q7A151
A	15	SER	-	expression tag	UNP Q7A151
A	16	GLY	-	expression tag	UNP Q7A151
A	17	LEU	-	expression tag	UNP Q7A151
A	18	VAL	-	expression tag	UNP Q7A151
A	19	PRO	-	expression tag	UNP Q7A151
A	20	ARG	-	expression tag	UNP Q7A151
A	21	GLY	-	expression tag	UNP Q7A151
A	22	SER	-	expression tag	UNP Q7A151
A	23	HIS	-	expression tag	UNP Q7A151
A	24	MET	-	expression tag	UNP Q7A151

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING ZN (CCD ID: ZNH) (formula: $C_{34}H_{32}N_4O_4Zn$).



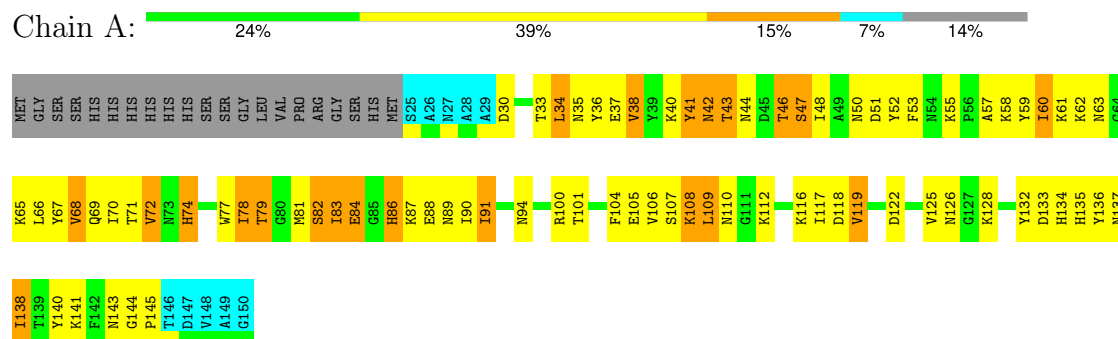
Mol	Chain	Residues	Atoms					
			Total	C	H	N	O	Zn
2	A	1	73	34	30	4	4	1

4 Residue-property plots

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: Iron-regulated surface determinant protein C

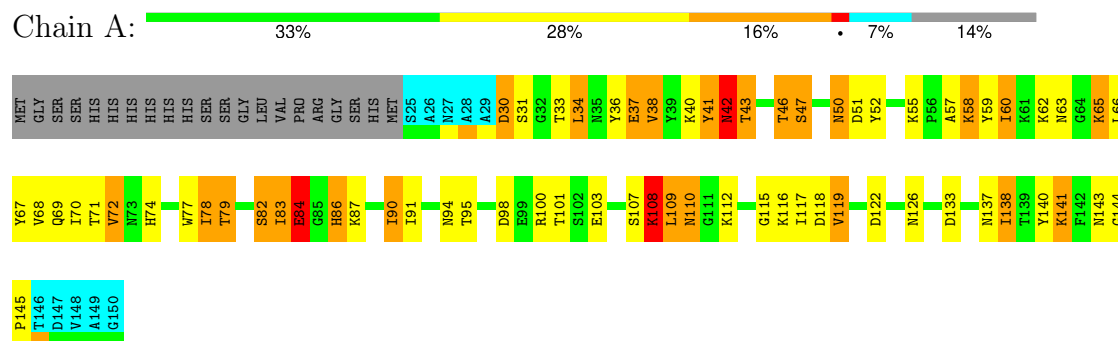


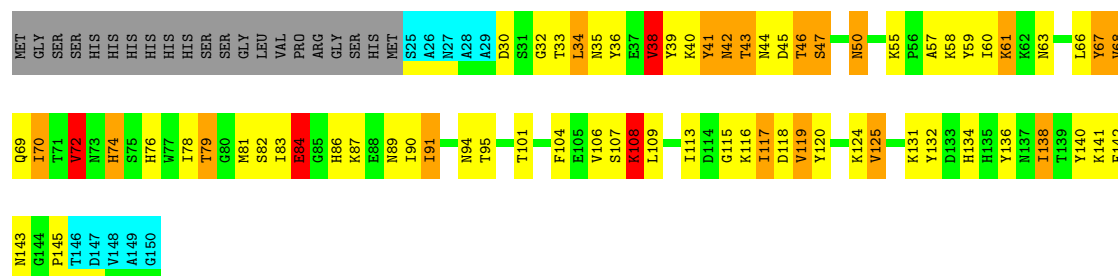
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: Iron-regulated surface determinant protein C

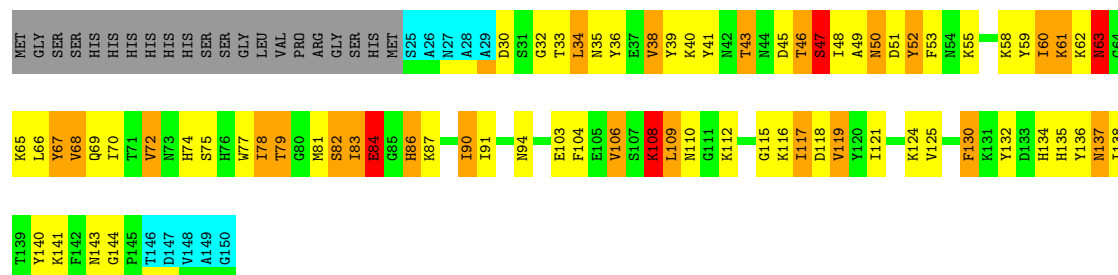




4.2.11 Score per residue for model 11

- Molecule 1: Iron-regulated surface determinant protein C

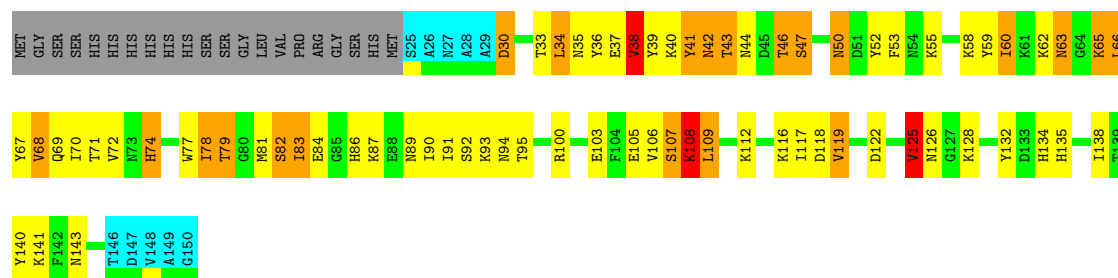
Chain A: 29% 32% 16% 7% 14%



4.2.12 Score per residue for model 12

- Molecule 1: Iron-regulated surface determinant protein C

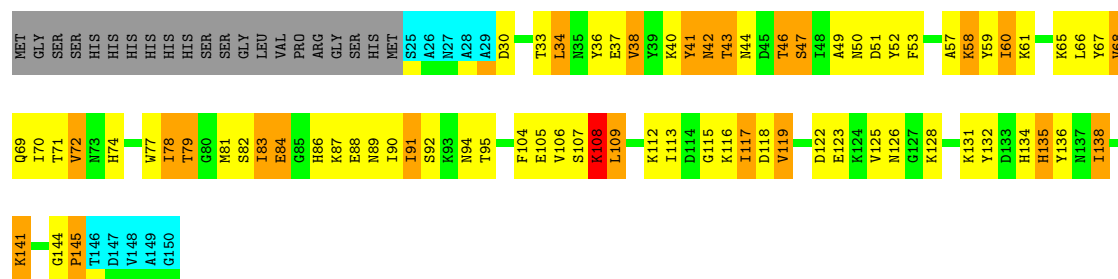
Chain A: 30% 33% 14% 7% 14%



4.2.13 Score per residue for model 13

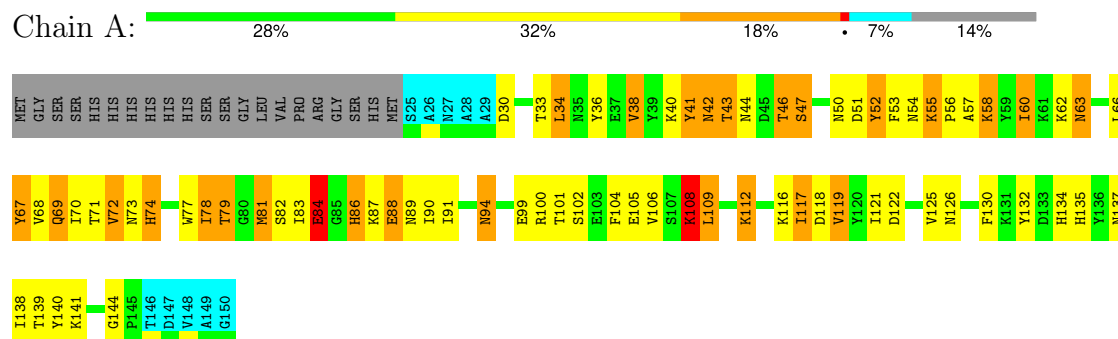
- Molecule 1: Iron-regulated surface determinant protein C

Chain A: 28% 35% 16% 7% 14%



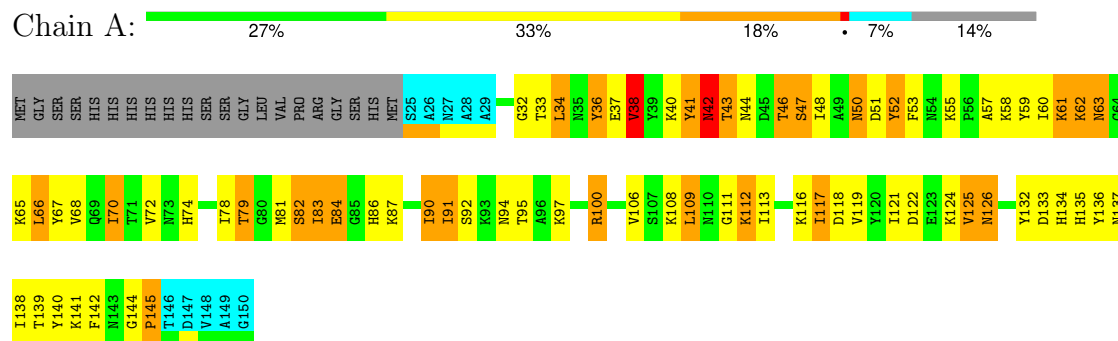
4.2.14 Score per residue for model 14

- Molecule 1: Iron-regulated surface determinant protein C



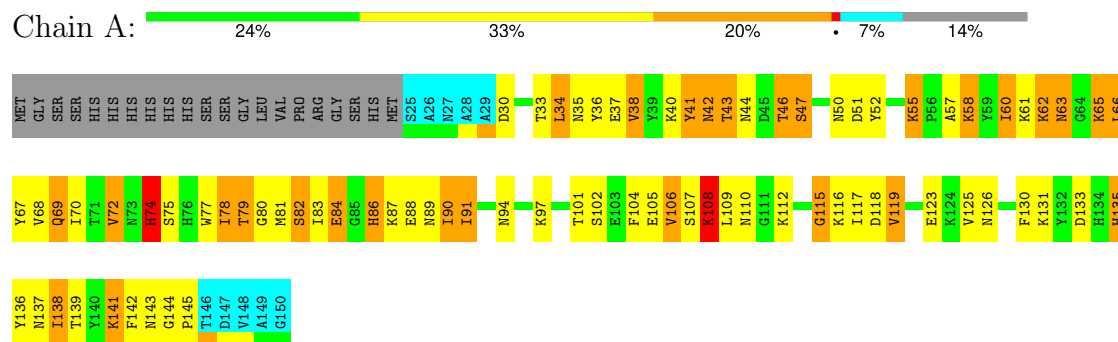
4.2.15 Score per residue for model 15

- Molecule 1: Iron-regulated surface determinant protein C



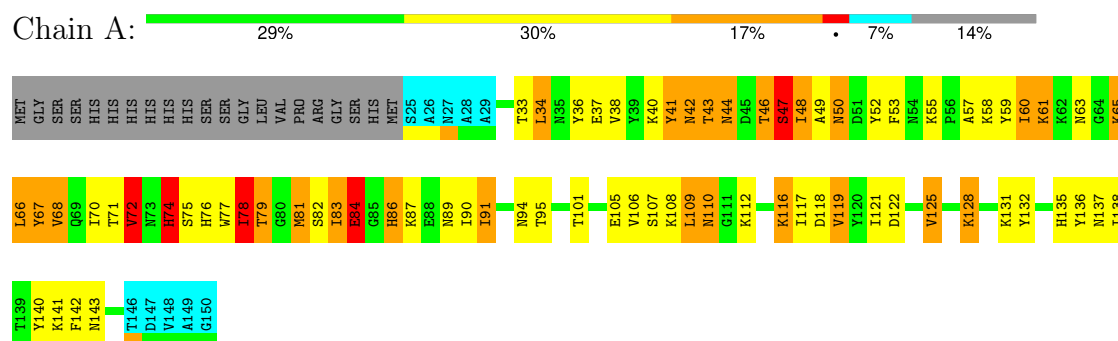
4.2.16 Score per residue for model 16

- Molecule 1: Iron-regulated surface determinant protein C



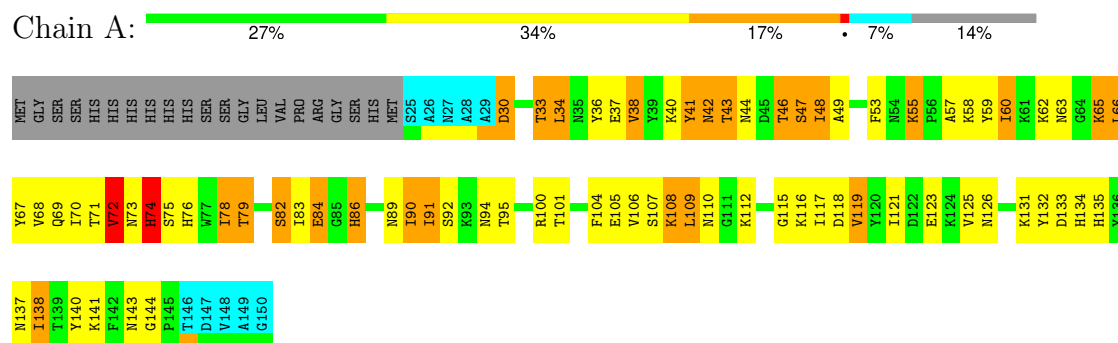
4.2.17 Score per residue for model 17

- Molecule 1: Iron-regulated surface determinant protein C



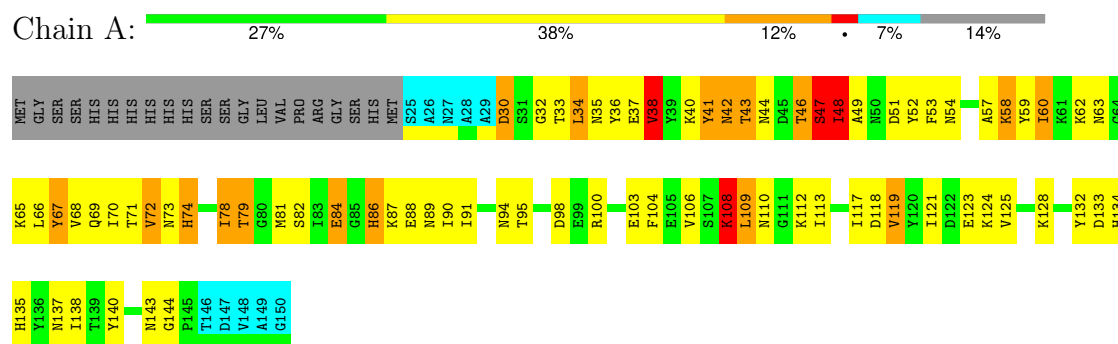
4.2.18 Score per residue for model 18

- Molecule 1: Iron-regulated surface determinant protein C



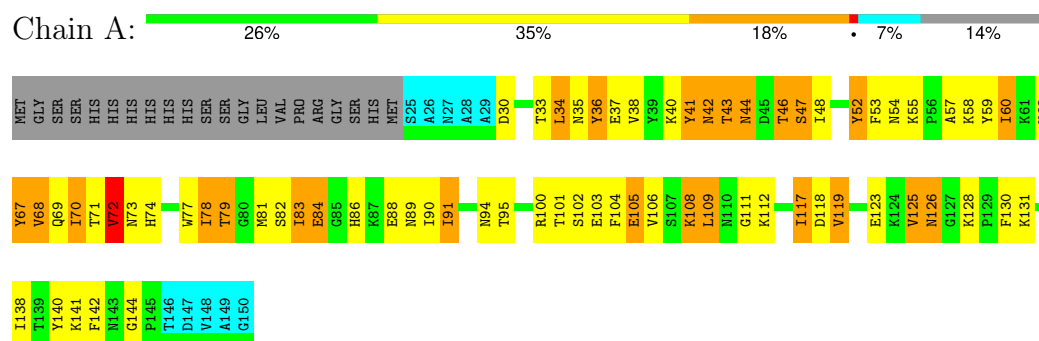
4.2.19 Score per residue for model 19

- Molecule 1: Iron-regulated surface determinant protein C



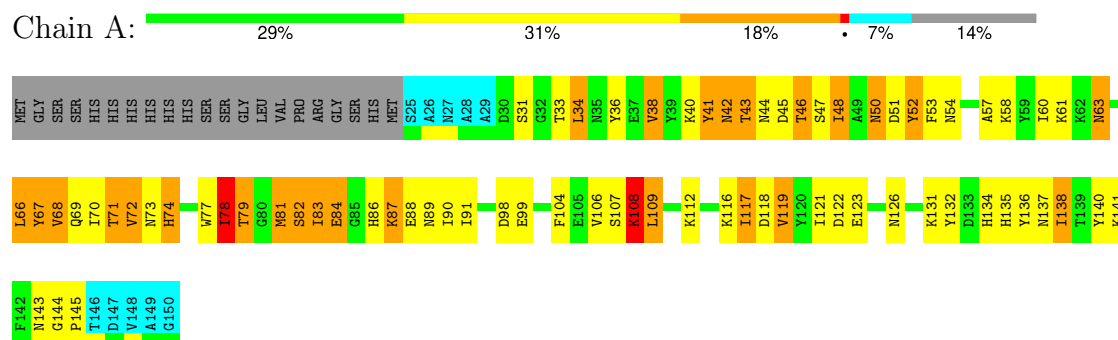
4.2.20 Score per residue for model 20

- Molecule 1: Iron-regulated surface determinant protein C



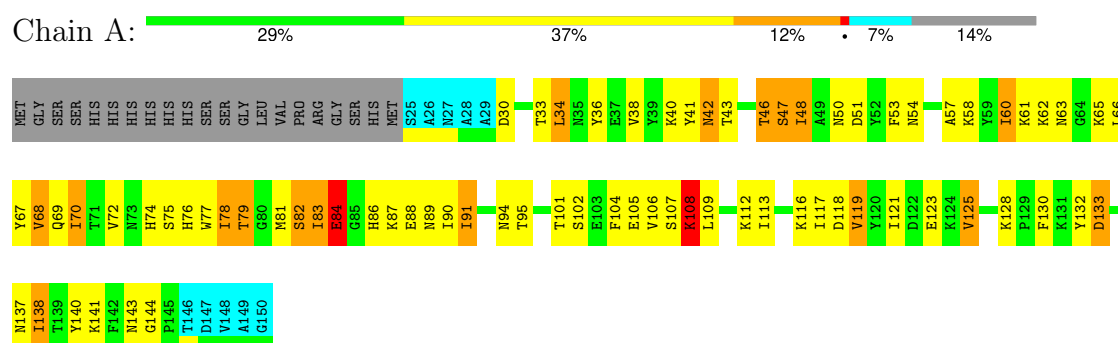
4.2.21 Score per residue for model 21

- Molecule 1: Iron-regulated surface determinant protein C



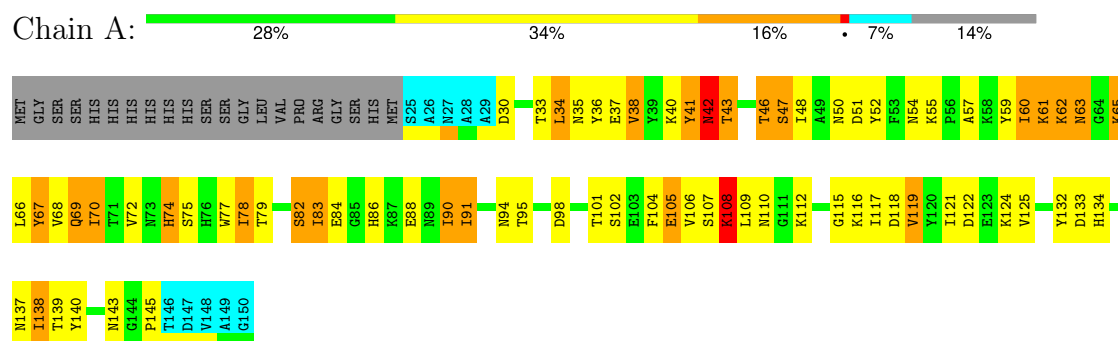
4.2.22 Score per residue for model 22

- Molecule 1: Iron-regulated surface determinant protein C



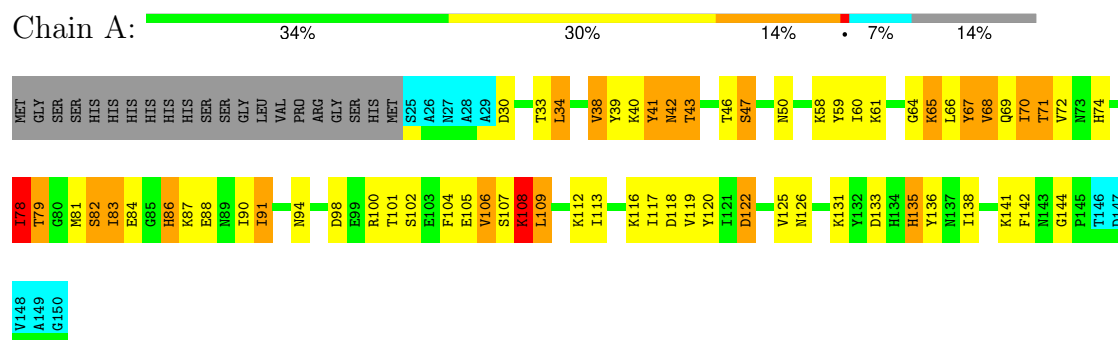
4.2.23 Score per residue for model 23

- Molecule 1: Iron-regulated surface determinant protein C



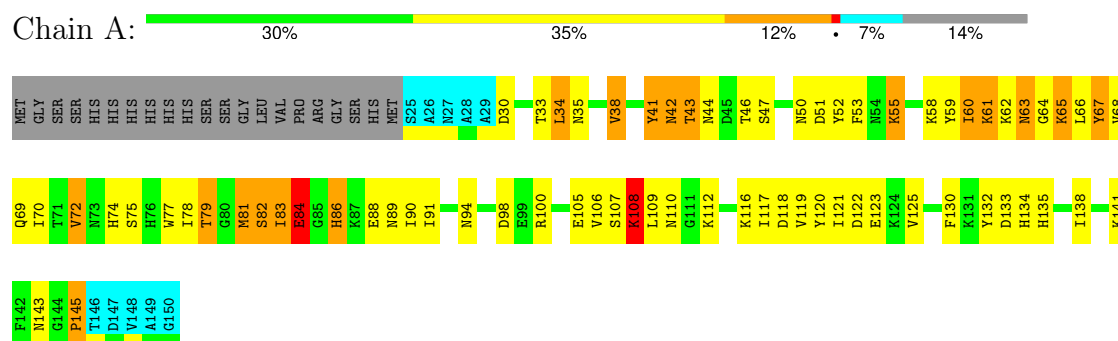
4.2.24 Score per residue for model 24

- Molecule 1: Iron-regulated surface determinant protein C



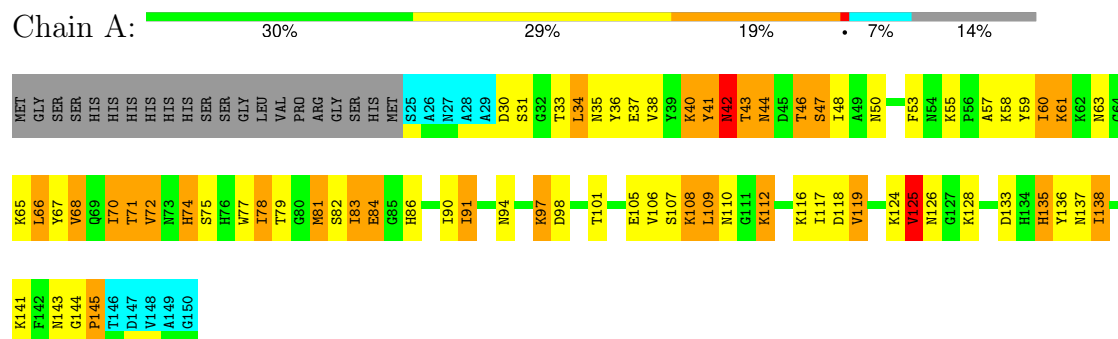
4.2.25 Score per residue for model 25

- Molecule 1: Iron-regulated surface determinant protein C



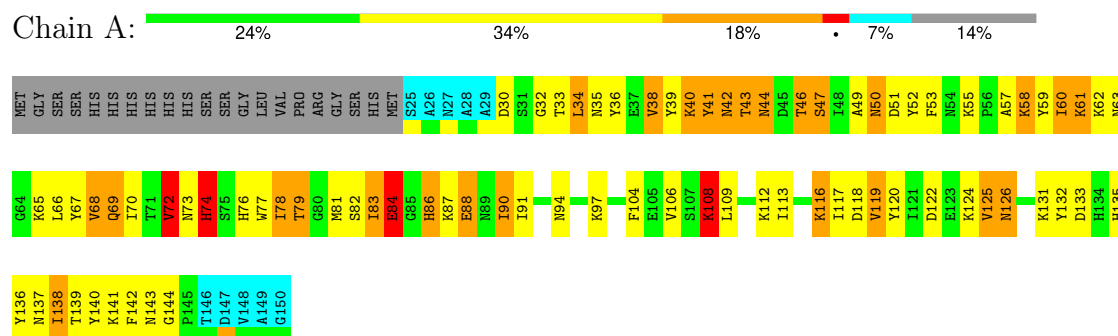
4.2.26 Score per residue for model 26

- Molecule 1: Iron-regulated surface determinant protein C



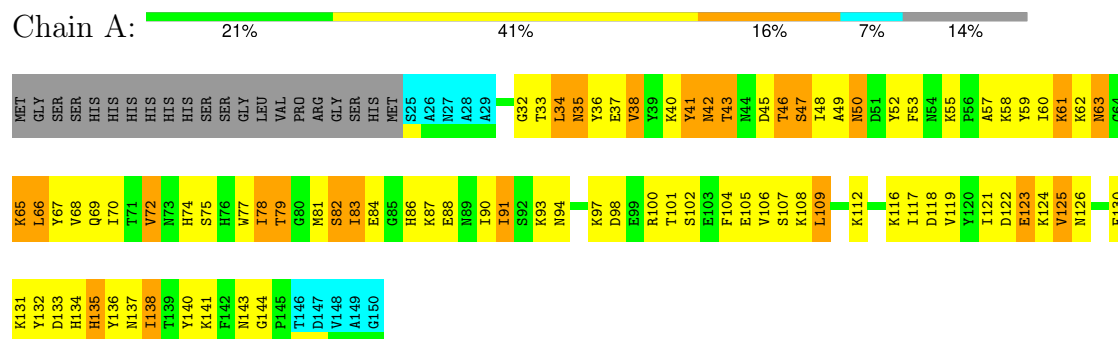
4.2.27 Score per residue for model 27

- Molecule 1: Iron-regulated surface determinant protein C



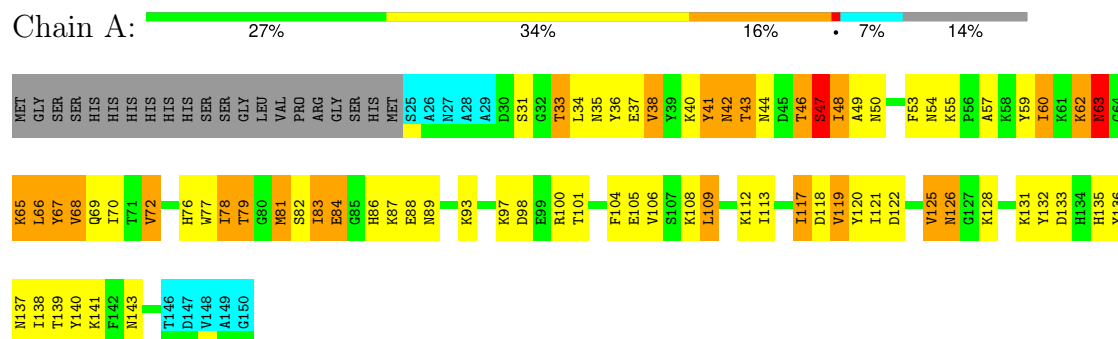
4.2.28 Score per residue for model 28

- Molecule 1: Iron-regulated surface determinant protein C



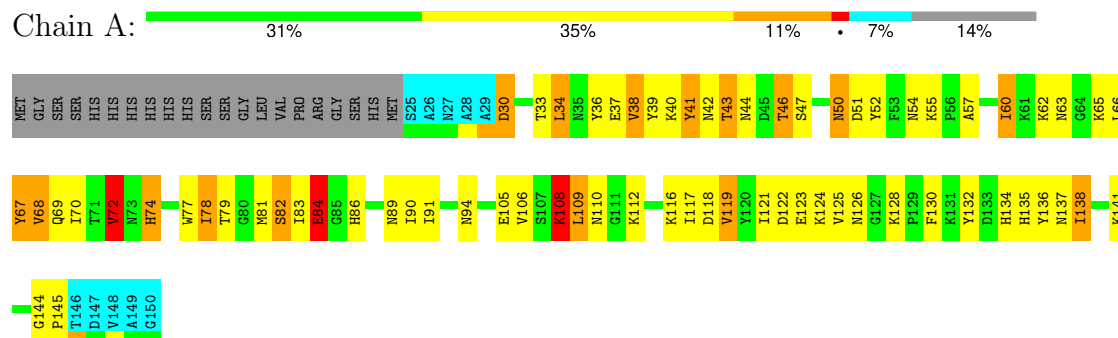
4.2.29 Score per residue for model 29

- Molecule 1: Iron-regulated surface determinant protein C



4.2.30 Score per residue for model 30

- Molecule 1: Iron-regulated surface determinant protein C



5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing*.

Of the 200 calculated structures, 30 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
X-PLOR NIH	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: ZNH

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 (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	1.42±0.01	0±0/975 (0.0± 0.0%)	1.66±0.01	5±1/1316 (0.4± 0.1%)
All	All	1.42	9/29250 (0.0%)	1.66	151/39480 (0.4%)

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
1	A	67	TYR	C-N	-5.46	1.30	1.33	5	1
1	A	115	GLY	N-CA	5.35	1.49	1.44	16	8

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(°)	Ideal(°)	Models	
								Worst	Total
1	A	125	VAL	N-CA-CB	-6.50	104.88	111.90	23	17
1	A	38	VAL	N-CA-CB	-6.45	104.91	112.32	12	10
1	A	110	ASN	N-CA-CB	-5.71	104.94	111.79	19	3
1	A	84	GLU	N-CA-CB	-5.58	105.45	113.65	16	19
1	A	72	VAL	N-CA-CB	-5.57	104.69	111.21	14	16
1	A	47	SER	N-CA-CB	-5.55	103.61	111.00	12	30
1	A	42	ASN	N-CA-CB	-5.54	105.50	113.65	13	27
1	A	78	ILE	N-CA-CB	-5.41	104.83	111.00	21	26
1	A	55	LYS	N-CA-CB	-5.14	104.82	111.35	25	2
1	A	106	VAL	N-CA-CB	-5.07	103.95	112.06	24	1

There are no chirality outliers.

There are no planarity outliers.

6.2 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 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	951	920	920	55±7
2	A	43	30	30	2±1
All	All	29820	28500	28500	1646

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 28.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:34:LEU:HD13	1:A:109:LEU:HD22	0.92	1.39	15	9
1:A:125:VAL:HG23	1:A:130:PHE:CE1	0.86	2.06	3	5
1:A:83:ILE:HD11	1:A:140:TYR:CD2	0.82	2.10	17	9
1:A:36:TYR:CZ	1:A:57:ALA:HB2	0.82	2.10	14	15
1:A:36:TYR:CE1	1:A:57:ALA:HB2	0.81	2.10	10	20
1:A:83:ILE:HD12	1:A:104:PHE:CD2	0.81	2.11	6	8
1:A:34:LEU:HD23	1:A:59:TYR:HB3	0.81	1.51	29	1
1:A:74:HIS:O	1:A:78:ILE:HG22	0.80	1.77	10	13
1:A:66:LEU:HD11	1:A:108:LYS:HA	0.77	1.56	20	11
1:A:66:LEU:O	1:A:66:LEU:HD22	0.77	1.80	17	8
1:A:38:VAL:O	1:A:46:THR:HG22	0.76	1.80	20	5
1:A:34:LEU:HD11	1:A:144:GLY:HA3	0.76	1.57	26	17
1:A:83:ILE:HD11	1:A:140:TYR:CE2	0.75	2.16	3	8
1:A:66:LEU:HD13	1:A:66:LEU:N	0.74	1.97	7	6
1:A:78:ILE:HD11	1:A:81:MET:SD	0.74	2.23	5	8
1:A:72:VAL:HG22	1:A:81:MET:CE	0.74	2.12	16	2
1:A:72:VAL:HG22	1:A:81:MET:HE1	0.74	1.60	21	3
1:A:65:LYS:C	1:A:66:LEU:HD13	0.72	2.09	12	6
1:A:66:LEU:C	1:A:66:LEU:HD22	0.72	2.09	5	5
1:A:38:VAL:HG13	1:A:139:THR:O	0.72	1.84	23	1
1:A:79:THR:HG22	1:A:118:ASP:C	0.72	2.09	23	30
1:A:66:LEU:HD22	1:A:66:LEU:O	0.71	1.86	15	5
1:A:79:THR:HG22	1:A:118:ASP:O	0.71	1.85	23	29
1:A:34:LEU:HD11	1:A:109:LEU:HB2	0.71	1.62	17	1
1:A:34:LEU:HG	1:A:109:LEU:HD22	0.70	1.61	9	2
1:A:72:VAL:HG12	1:A:78:ILE:HG21	0.70	1.63	27	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:38:VAL:C	1:A:46:THR:HG23	0.70	2.11	29	14
1:A:67:TYR:O	1:A:68:VAL:HG23	0.69	1.87	27	2
1:A:66:LEU:HD22	1:A:66:LEU:C	0.68	2.13	29	2
1:A:69:GLN:C	1:A:70:ILE:HD12	0.68	2.14	5	1
1:A:72:VAL:HG12	1:A:81:MET:HE1	0.68	1.65	4	1
1:A:38:VAL:O	1:A:46:THR:HG23	0.68	1.89	15	13
1:A:78:ILE:HD11	1:A:81:MET:CG	0.67	2.19	3	6
1:A:77:TRP:O	1:A:119:VAL:HG12	0.67	1.90	22	17
1:A:34:LEU:HD11	1:A:109:LEU:CB	0.67	2.20	17	1
1:A:66:LEU:HD21	1:A:106:VAL:O	0.66	1.90	27	14
1:A:38:VAL:HG23	1:A:46:THR:OG1	0.66	1.91	13	1
1:A:53:PHE:CZ	1:A:81:MET:HE3	0.66	2.25	17	1
1:A:38:VAL:HG12	1:A:139:THR:O	0.65	1.90	14	3
1:A:125:VAL:HG21	1:A:130:PHE:CD2	0.65	2.26	22	1
1:A:36:TYR:CE2	1:A:57:ALA:HB2	0.65	2.26	20	2
1:A:34:LEU:HD23	1:A:59:TYR:CE1	0.64	2.27	26	2
1:A:66:LEU:HD21	1:A:108:LYS:H	0.64	1.53	14	6
1:A:34:LEU:HD23	1:A:59:TYR:CB	0.64	2.22	29	1
1:A:34:LEU:HD23	1:A:59:TYR:CD1	0.64	2.27	26	2
1:A:83:ILE:HD11	1:A:140:TYR:CE1	0.64	2.27	23	1
1:A:72:VAL:HG13	1:A:78:ILE:HD13	0.64	1.68	2	3
1:A:48:ILE:HG22	1:A:49:ALA:N	0.63	2.07	4	8
1:A:47:SER:HB3	1:A:48:ILE:HD12	0.63	1.71	19	3
1:A:83:ILE:HD12	1:A:104:PHE:CG	0.62	2.30	24	6
1:A:78:ILE:HD11	1:A:81:MET:HG3	0.61	1.70	29	5
1:A:121:ILE:HD13	2:A:151:ZNH:CBB	0.61	2.25	18	13
1:A:66:LEU:HD11	1:A:108:LYS:N	0.61	2.09	25	6
1:A:125:VAL:HG12	1:A:125:VAL:O	0.61	1.96	3	7
1:A:38:VAL:HG23	1:A:46:THR:HG21	0.60	1.73	3	2
1:A:66:LEU:HD11	1:A:106:VAL:O	0.60	1.96	26	1
1:A:34:LEU:C	1:A:34:LEU:HD12	0.60	2.21	29	1
1:A:70:ILE:HD12	1:A:70:ILE:N	0.60	2.11	5	1
1:A:66:LEU:HD23	1:A:66:LEU:N	0.60	2.12	20	11
1:A:77:TRP:O	1:A:119:VAL:HG13	0.60	1.97	4	3
1:A:59:TYR:CD2	1:A:109:LEU:HD21	0.59	2.32	9	4
1:A:113:ILE:HD12	1:A:142:PHE:CD2	0.59	2.32	27	2
1:A:34:LEU:HD12	1:A:144:GLY:HA3	0.59	1.75	20	4
1:A:67:TYR:CD1	1:A:67:TYR:N	0.58	2.71	10	30
1:A:52:TYR:CD1	1:A:78:ILE:HD13	0.58	2.33	21	1
1:A:47:SER:C	1:A:48:ILE:HD12	0.58	2.23	17	2
1:A:91:ILE:HG22	1:A:101:THR:C	0.58	2.24	24	8

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:66:LEU:CD1	1:A:66:LEU:N	0.57	2.67	17	7
1:A:117:ILE:HG21	1:A:138:ILE:HD11	0.57	1.76	13	2
1:A:137:ASN:O	1:A:138:ILE:HG23	0.57	1.99	29	19
1:A:83:ILE:HD12	1:A:104:PHE:CD1	0.57	2.35	11	2
1:A:124:LYS:O	1:A:125:VAL:HG23	0.57	1.98	27	2
1:A:77:TRP:CD2	2:A:151:ZNH:HBB2	0.57	2.35	1	4
1:A:125:VAL:HG23	1:A:130:PHE:CZ	0.57	2.34	3	1
1:A:72:VAL:CG1	1:A:78:ILE:HG21	0.57	2.29	1	8
1:A:38:VAL:HG12	1:A:140:TYR:HA	0.57	1.75	11	4
1:A:67:TYR:C	1:A:68:VAL:HG12	0.57	2.23	24	6
1:A:66:LEU:HD11	1:A:108:LYS:CA	0.56	2.30	11	12
1:A:78:ILE:O	1:A:78:ILE:HG23	0.56	1.99	14	29
1:A:34:LEU:HD11	1:A:144:GLY:CA	0.56	2.30	9	7
1:A:46:THR:HG21	1:A:50:ASN:CB	0.56	2.30	25	2
1:A:78:ILE:HD11	1:A:117:ILE:HG13	0.56	1.77	21	5
1:A:38:VAL:O	1:A:38:VAL:HG23	0.56	1.98	8	1
1:A:106:VAL:HG21	1:A:113:ILE:HD13	0.56	1.76	27	3
1:A:59:TYR:CA	1:A:68:VAL:HG13	0.56	2.30	13	2
1:A:78:ILE:HA	1:A:119:VAL:HG12	0.56	1.77	30	1
1:A:59:TYR:HA	1:A:68:VAL:HG13	0.56	1.76	13	2
1:A:34:LEU:HD11	1:A:145:PRO:HD2	0.56	1.76	15	3
1:A:48:ILE:HG21	2:A:151:ZNH:C1D	0.55	2.31	29	3
1:A:90:ILE:HD12	1:A:90:ILE:O	0.55	2.01	25	2
1:A:34:LEU:HD12	1:A:35:ASN:O	0.55	2.02	29	1
1:A:66:LEU:HD13	1:A:108:LYS:CE	0.55	2.31	20	2
1:A:91:ILE:HG22	1:A:101:THR:O	0.55	2.02	24	12
1:A:117:ILE:HG21	1:A:138:ILE:CD1	0.55	2.31	13	2
1:A:34:LEU:HD13	1:A:35:ASN:N	0.55	2.17	9	1
1:A:72:VAL:CG2	1:A:81:MET:HE1	0.54	2.31	30	1
1:A:91:ILE:HG21	1:A:101:THR:HG22	0.54	1.79	24	9
1:A:53:PHE:CG	1:A:53:PHE:O	0.54	2.60	28	3
1:A:66:LEU:N	1:A:66:LEU:HD13	0.54	2.18	17	7
1:A:125:VAL:HG23	1:A:130:PHE:CD1	0.54	2.37	25	1
1:A:34:LEU:HD13	1:A:35:ASN:H	0.53	1.63	9	1
1:A:38:VAL:CG2	1:A:53:PHE:CE1	0.53	2.91	14	2
1:A:84:GLU:CB	1:A:86:HIS:CE1	0.53	2.92	2	10
1:A:66:LEU:HD13	1:A:108:LYS:HE2	0.53	1.80	9	1
1:A:34:LEU:C	1:A:34:LEU:HD23	0.53	2.29	12	4
1:A:34:LEU:HD13	1:A:59:TYR:CD2	0.53	2.39	8	1
1:A:88:GLU:CB	1:A:104:PHE:CE1	0.52	2.92	7	6
1:A:38:VAL:CG1	1:A:140:TYR:CD1	0.52	2.92	14	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:121:ILE:HD13	2:A:151:ZNH:HBB1	0.52	1.79	21	6
1:A:34:LEU:HD21	1:A:143:ASN:C	0.52	2.28	8	7
1:A:88:GLU:CB	1:A:104:PHE:CZ	0.52	2.92	2	3
1:A:74:HIS:CE1	1:A:77:TRP:CE2	0.52	2.97	30	1
1:A:34:LEU:HD22	1:A:109:LEU:HD22	0.52	1.80	14	3
1:A:66:LEU:N	1:A:66:LEU:CD1	0.52	2.69	7	6
1:A:53:PHE:CZ	1:A:81:MET:CE	0.52	2.92	17	1
1:A:34:LEU:CB	1:A:59:TYR:CD1	0.52	2.92	3	5
1:A:86:HIS:CD2	1:A:86:HIS:C	0.52	2.88	24	24
1:A:59:TYR:CD1	1:A:61:LYS:CE	0.52	2.92	17	1
1:A:72:VAL:HG22	1:A:100:ARG:O	0.51	2.05	6	2
1:A:66:LEU:C	1:A:66:LEU:CD2	0.51	2.83	5	7
1:A:41:TYR:CE2	1:A:137:ASN:ND2	0.51	2.78	9	5
1:A:41:TYR:CD2	1:A:137:ASN:ND2	0.51	2.78	20	4
1:A:49:ALA:HB2	1:A:138:ILE:HG21	0.51	1.83	8	1
1:A:36:TYR:CD2	1:A:37:GLU:N	0.51	2.79	7	1
1:A:132:TYR:OH	2:A:151:ZNH:NA	0.51	2.44	11	20
1:A:83:ILE:O	1:A:86:HIS:CD2	0.51	2.64	22	14
1:A:49:ALA:HB1	1:A:53:PHE:CE2	0.50	2.41	29	1
1:A:66:LEU:HD11	1:A:108:LYS:H	0.50	1.65	6	1
1:A:77:TRP:HA	1:A:121:ILE:HD11	0.50	1.82	11	5
1:A:86:HIS:CD2	1:A:86:HIS:N	0.50	2.79	12	3
1:A:71:THR:O	1:A:71:THR:HG22	0.50	2.06	26	1
1:A:53:PHE:O	1:A:53:PHE:CG	0.50	2.65	20	3
1:A:83:ILE:HD11	1:A:140:TYR:CZ	0.50	2.42	12	1
1:A:60:ILE:HG23	1:A:67:TYR:HB2	0.50	1.82	7	1
1:A:117:ILE:C	1:A:117:ILE:HD13	0.50	2.31	15	3
1:A:105:GLU:C	1:A:106:VAL:HG13	0.50	2.32	25	16
1:A:36:TYR:CE1	1:A:57:ALA:CB	0.50	2.93	10	2
1:A:83:ILE:CD1	1:A:140:TYR:CD2	0.50	2.93	17	1
1:A:36:TYR:CE2	1:A:70:ILE:HD11	0.49	2.42	23	2
1:A:34:LEU:HD23	1:A:34:LEU:O	0.49	2.07	12	3
1:A:77:TRP:CE2	2:A:151:ZNH:HBB2	0.49	2.42	20	1
1:A:106:VAL:HG22	1:A:107:SER:N	0.49	2.22	21	1
1:A:41:TYR:O	1:A:42:ASN:CB	0.49	2.61	22	29
1:A:32:GLY:O	1:A:59:TYR:CD2	0.49	2.65	11	2
1:A:83:ILE:O	1:A:84:GLU:CB	0.49	2.61	27	16
1:A:117:ILE:HD13	1:A:117:ILE:O	0.49	2.08	15	2
1:A:83:ILE:HD11	1:A:140:TYR:CG	0.49	2.43	15	1
1:A:86:HIS:CD2	1:A:86:HIS:O	0.48	2.66	9	15
1:A:52:TYR:CD1	1:A:52:TYR:N	0.48	2.80	17	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:66:LEU:O	1:A:66:LEU:CD2	0.48	2.61	16	6
1:A:66:LEU:N	1:A:66:LEU:CD2	0.48	2.76	20	5
1:A:125:VAL:O	1:A:125:VAL:HG12	0.48	2.08	26	6
1:A:83:ILE:HD11	1:A:140:TYR:CD1	0.48	2.43	15	1
1:A:84:GLU:HB2	1:A:86:HIS:CE1	0.48	2.43	10	8
1:A:70:ILE:CG2	1:A:81:MET:HE1	0.48	2.38	8	2
1:A:34:LEU:HD13	1:A:59:TYR:CG	0.48	2.43	8	1
1:A:59:TYR:CD1	1:A:59:TYR:C	0.48	2.90	11	6
1:A:36:TYR:CD1	1:A:36:TYR:O	0.48	2.67	22	14
1:A:84:GLU:CB	1:A:86:HIS:NE2	0.48	2.76	13	5
1:A:71:THR:HG22	1:A:71:THR:O	0.48	2.08	24	1
1:A:121:ILE:HD12	1:A:132:TYR:CB	0.48	2.39	30	1
1:A:38:VAL:O	1:A:39:TYR:CD1	0.48	2.66	11	8
1:A:41:TYR:CD1	1:A:138:ILE:C	0.48	2.92	6	3
1:A:36:TYR:C	1:A:36:TYR:CD1	0.48	2.91	14	1
1:A:49:ALA:O	1:A:53:PHE:CD2	0.48	2.67	13	4
1:A:88:GLU:HB3	1:A:104:PHE:CZ	0.48	2.44	2	3
1:A:53:PHE:O	1:A:53:PHE:CD2	0.48	2.66	21	4
1:A:41:TYR:O	1:A:43:THR:HG23	0.48	2.08	5	21
1:A:112:LYS:CG	1:A:139:THR:CG2	0.48	2.91	16	2
1:A:90:ILE:HG22	1:A:100:ARG:NH2	0.48	2.24	15	1
1:A:60:ILE:O	1:A:67:TYR:CD1	0.48	2.66	17	2
1:A:117:ILE:HG21	1:A:138:ILE:HD13	0.48	1.85	29	1
1:A:58:LYS:CB	1:A:69:GLN:CG	0.48	2.92	11	6
1:A:38:VAL:HG23	1:A:46:THR:CG2	0.48	2.39	3	1
1:A:41:TYR:CD2	1:A:137:ASN:CG	0.48	2.91	17	1
1:A:36:TYR:CE1	1:A:57:ALA:N	0.48	2.81	19	1
1:A:51:ASP:C	1:A:52:TYR:CG	0.48	2.92	6	8
1:A:66:LEU:C	1:A:67:TYR:CG	0.48	2.92	18	23
1:A:84:GLU:HB3	1:A:86:HIS:CE1	0.48	2.43	29	5
1:A:123:GLU:O	1:A:130:PHE:CD1	0.48	2.67	3	2
1:A:100:ARG:CG	1:A:101:THR:N	0.47	2.77	1	4
1:A:78:ILE:HD12	2:A:151:ZNH:HBC2	0.47	1.85	16	2
1:A:77:TRP:O	1:A:121:ILE:HD11	0.47	2.09	17	1
1:A:34:LEU:HB2	1:A:59:TYR:CD1	0.47	2.45	10	8
1:A:46:THR:CG2	1:A:50:ASN:CB	0.47	2.92	25	8
1:A:70:ILE:N	1:A:70:ILE:CD1	0.47	2.77	5	1
1:A:41:TYR:CD2	1:A:137:ASN:OD1	0.47	2.68	19	2
1:A:32:GLY:O	1:A:59:TYR:CD1	0.47	2.67	2	5
1:A:66:LEU:HD23	1:A:66:LEU:O	0.47	2.09	6	3
1:A:52:TYR:CD2	2:A:151:ZNH:HMC1	0.47	2.45	11	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:135:HIS:O	1:A:136:TYR:CD1	0.47	2.67	2	4
1:A:135:HIS:O	1:A:136:TYR:CG	0.47	2.68	20	14
1:A:72:VAL:CG2	1:A:100:ARG:CB	0.47	2.93	7	2
1:A:72:VAL:HG13	1:A:81:MET:HE1	0.47	1.86	30	1
1:A:66:LEU:CD1	1:A:108:LYS:CA	0.47	2.92	23	3
1:A:51:ASP:C	1:A:52:TYR:CD1	0.47	2.92	19	3
1:A:61:LYS:N	1:A:61:LYS:CD	0.47	2.78	10	1
1:A:135:HIS:CD2	1:A:135:HIS:O	0.47	2.68	21	2
1:A:106:VAL:HG21	1:A:113:ILE:CD1	0.47	2.40	27	3
1:A:55:LYS:CD	1:A:55:LYS:N	0.47	2.78	18	1
1:A:82:SER:HG	1:A:116:LYS:C	0.47	2.17	1	2
1:A:65:LYS:CD	1:A:65:LYS:N	0.47	2.78	17	4
1:A:59:TYR:CD1	1:A:61:LYS:HD2	0.46	2.45	11	4
1:A:119:VAL:HG11	1:A:132:TYR:CE2	0.46	2.44	18	1
1:A:38:VAL:CG1	1:A:140:TYR:CE1	0.46	2.98	27	2
1:A:54:ASN:HB3	1:A:71:THR:HG21	0.46	1.87	14	2
1:A:124:LYS:C	1:A:125:VAL:HG23	0.46	2.36	15	1
1:A:52:TYR:O	1:A:52:TYR:CD1	0.46	2.68	20	3
1:A:66:LEU:C	1:A:67:TYR:CD1	0.46	2.93	18	2
1:A:106:VAL:CG2	1:A:107:SER:N	0.46	2.79	21	2
1:A:59:TYR:CD2	1:A:109:LEU:CD2	0.46	2.98	12	5
1:A:130:PHE:CD1	1:A:130:PHE:N	0.46	2.83	3	1
1:A:137:ASN:C	1:A:138:ILE:HG23	0.46	2.35	11	2
1:A:50:ASN:O	1:A:53:PHE:CE1	0.46	2.68	14	1
1:A:143:ASN:ND2	1:A:143:ASN:N	0.46	2.64	19	3
1:A:106:VAL:CG2	1:A:113:ILE:HD13	0.46	2.40	27	2
1:A:83:ILE:O	1:A:86:HIS:CE1	0.46	2.69	10	7
1:A:66:LEU:CD2	1:A:66:LEU:C	0.46	2.89	12	6
1:A:66:LEU:HD13	1:A:66:LEU:H	0.46	1.69	15	5
1:A:36:TYR:CE2	1:A:53:PHE:CE2	0.46	3.04	20	1
1:A:120:TYR:CZ	1:A:133:ASP:OD2	0.45	2.68	25	2
1:A:82:SER:N	1:A:116:LYS:O	0.45	2.50	11	26
1:A:125:VAL:O	1:A:126:ASN:CB	0.45	2.65	15	5
1:A:41:TYR:CE1	1:A:138:ILE:C	0.45	2.95	16	3
1:A:52:TYR:CE1	1:A:78:ILE:HD13	0.45	2.47	21	2
1:A:90:ILE:CG1	1:A:90:ILE:O	0.45	2.65	18	2
1:A:79:THR:CG2	1:A:118:ASP:CB	0.45	2.94	23	9
1:A:88:GLU:HB2	1:A:104:PHE:CZ	0.45	2.47	7	8
1:A:81:MET:HG3	1:A:117:ILE:HD12	0.45	1.89	5	1
1:A:68:VAL:CG2	1:A:70:ILE:HD12	0.45	2.42	22	1
1:A:61:LYS:HD3	1:A:109:LEU:HD21	0.45	1.87	4	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:74:HIS:CD2	1:A:77:TRP:CE2	0.45	3.05	7	1
1:A:130:PHE:N	1:A:130:PHE:CD1	0.45	2.85	14	2
1:A:135:HIS:O	1:A:136:TYR:CD2	0.45	2.70	17	1
1:A:119:VAL:HG11	1:A:132:TYR:HE2	0.45	1.72	18	2
1:A:74:HIS:CD2	1:A:77:TRP:CZ3	0.45	3.05	26	1
1:A:112:LYS:CD	1:A:112:LYS:C	0.45	2.90	2	2
1:A:34:LEU:CD1	1:A:143:ASN:O	0.45	2.65	26	2
1:A:54:ASN:CB	1:A:71:THR:HG21	0.45	2.42	21	1
1:A:34:LEU:HB2	1:A:59:TYR:CD2	0.45	2.47	23	2
1:A:34:LEU:C	1:A:34:LEU:CD1	0.45	2.89	29	1
1:A:62:LYS:CB	1:A:65:LYS:CD	0.45	2.95	3	3
1:A:43:THR:OG1	1:A:44:ASN:N	0.45	2.50	21	20
1:A:34:LEU:C	1:A:34:LEU:CD2	0.44	2.90	12	4
1:A:74:HIS:CD2	1:A:77:TRP:CE3	0.44	3.04	26	1
1:A:72:VAL:HG13	1:A:81:MET:CE	0.44	2.42	30	1
1:A:53:PHE:CE1	1:A:81:MET:SD	0.44	3.10	8	1
1:A:53:PHE:N	1:A:53:PHE:CD1	0.44	2.83	12	1
1:A:78:ILE:O	1:A:78:ILE:CG2	0.44	2.66	22	20
1:A:113:ILE:HD12	1:A:142:PHE:CE2	0.44	2.47	8	1
1:A:53:PHE:CZ	1:A:81:MET:SD	0.44	3.11	29	5
1:A:59:TYR:HA	1:A:68:VAL:HG23	0.44	1.89	12	1
1:A:41:TYR:CG	1:A:137:ASN:OD1	0.44	2.71	17	2
1:A:82:SER:HG	1:A:86:HIS:C	0.44	2.20	20	1
1:A:30:ASP:O	1:A:60:ILE:HA	0.44	2.13	27	14
1:A:66:LEU:CD1	1:A:108:LYS:HA	0.44	2.42	23	7
1:A:46:THR:HG21	1:A:50:ASN:HB2	0.44	1.90	27	2
1:A:124:LYS:O	1:A:125:VAL:CG2	0.44	2.65	15	1
1:A:41:TYR:CE1	1:A:138:ILE:O	0.44	2.71	23	1
1:A:70:ILE:O	1:A:70:ILE:CG2	0.44	2.66	26	5
1:A:130:PHE:O	1:A:130:PHE:CG	0.44	2.70	25	1
1:A:81:MET:HE2	1:A:138:ILE:HD11	0.44	1.90	28	1
1:A:38:VAL:O	1:A:46:THR:CG2	0.44	2.65	6	11
1:A:118:ASP:C	1:A:119:VAL:CG2	0.44	2.91	28	5
1:A:135:HIS:CD2	1:A:135:HIS:C	0.44	2.96	3	1
1:A:115:GLY:C	1:A:116:LYS:CG	0.44	2.91	16	4
1:A:109:LEU:H	1:A:109:LEU:HD12	0.44	1.73	11	1
1:A:81:MET:HE2	1:A:138:ILE:CD1	0.44	2.43	28	1
1:A:69:GLN:O	1:A:69:GLN:CG	0.44	2.66	18	12
1:A:35:ASN:O	1:A:143:ASN:N	0.44	2.51	11	10
1:A:60:ILE:O	1:A:66:LEU:HA	0.43	2.13	22	3
1:A:130:PHE:CD2	1:A:130:PHE:O	0.43	2.71	11	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:58:LYS:HB2	1:A:69:GLN:CG	0.43	2.43	6	5
1:A:112:LYS:HG3	1:A:139:THR:CG2	0.43	2.43	3	3
1:A:68:VAL:O	1:A:68:VAL:HG12	0.43	2.13	11	3
1:A:36:TYR:CD1	1:A:36:TYR:C	0.43	2.91	8	8
1:A:33:THR:HG22	1:A:57:ALA:O	0.43	2.12	29	1
1:A:68:VAL:HG12	1:A:68:VAL:O	0.43	2.11	6	2
1:A:110:ASN:ND2	1:A:145:PRO:O	0.43	2.52	8	1
1:A:119:VAL:HB	1:A:132:TYR:CE2	0.43	2.49	11	2
1:A:142:PHE:C	1:A:143:ASN:ND2	0.43	2.76	10	2
1:A:64:GLY:C	1:A:65:LYS:CD	0.43	2.92	24	2
1:A:120:TYR:CD1	1:A:120:TYR:C	0.43	2.95	29	1
1:A:126:ASN:ND2	1:A:126:ASN:N	0.43	2.64	2	3
1:A:83:ILE:O	1:A:86:HIS:NE2	0.43	2.51	23	6
1:A:84:GLU:HB3	1:A:86:HIS:NE2	0.43	2.28	12	5
1:A:73:ASN:C	1:A:74:HIS:CG	0.43	2.93	21	1
1:A:36:TYR:HE2	1:A:70:ILE:HD11	0.43	1.73	4	1
1:A:70:ILE:O	1:A:70:ILE:HG22	0.43	2.13	15	3
1:A:137:ASN:O	1:A:138:ILE:CG2	0.43	2.67	11	3
1:A:141:LYS:C	1:A:142:PHE:CG	0.43	2.97	16	1
1:A:110:ASN:CB	1:A:145:PRO:O	0.43	2.67	26	1
1:A:121:ILE:HD12	1:A:132:TYR:HB3	0.43	1.90	30	1
1:A:68:VAL:HG11	1:A:142:PHE:CE1	0.43	2.49	5	1
1:A:34:LEU:CD1	1:A:109:LEU:HD22	0.43	2.29	15	3
1:A:34:LEU:HD12	1:A:35:ASN:C	0.43	2.39	29	1
1:A:72:VAL:HG11	1:A:81:MET:CE	0.43	2.44	7	1
1:A:109:LEU:HD12	1:A:109:LEU:N	0.43	2.29	11	1
1:A:46:THR:OG1	1:A:50:ASN:CB	0.43	2.66	25	1
1:A:119:VAL:HG12	1:A:121:ILE:CD1	0.43	2.43	28	1
1:A:90:ILE:O	1:A:90:ILE:CG1	0.43	2.67	1	9
1:A:105:GLU:C	1:A:106:VAL:CG1	0.43	2.92	4	5
1:A:63:ASN:ND2	1:A:63:ASN:N	0.43	2.64	11	1
1:A:82:SER:O	1:A:83:ILE:CG1	0.43	2.67	25	2
1:A:40:LYS:HZ2	1:A:138:ILE:HG22	0.43	1.73	26	1
1:A:88:GLU:HB2	1:A:104:PHE:CE1	0.43	2.48	19	3
1:A:111:GLY:N	1:A:142:PHE:O	0.43	2.52	20	2
1:A:62:LYS:O	1:A:65:LYS:CG	0.42	2.67	2	7
1:A:130:PHE:CE2	2:A:151:ZNH:HMB3	0.42	2.49	2	2
1:A:102:SER:HB2	1:A:104:PHE:CZ	0.42	2.49	4	3
1:A:109:LEU:C	1:A:109:LEU:HD12	0.42	2.38	7	1
1:A:34:LEU:CD2	1:A:143:ASN:O	0.42	2.67	22	2
1:A:72:VAL:HG22	1:A:100:ARG:HB3	0.42	1.91	25	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:37:GLU:O	1:A:141:LYS:N	0.42	2.52	9	4
1:A:58:LYS:CB	1:A:69:GLN:CD	0.42	2.91	1	3
1:A:77:TRP:CG	2:A:151:ZNH:HBB2	0.42	2.49	5	1
1:A:109:LEU:O	1:A:109:LEU:CD1	0.42	2.67	14	5
1:A:123:GLU:HG2	1:A:130:PHE:CE1	0.42	2.48	25	1
1:A:112:LYS:CD	1:A:112:LYS:O	0.42	2.67	26	1
1:A:125:VAL:N	1:A:128:LYS:O	0.42	2.52	4	3
1:A:59:TYR:HB2	1:A:109:LEU:HD23	0.42	1.91	5	1
1:A:60:ILE:CG2	1:A:67:TYR:O	0.42	2.67	14	4
1:A:41:TYR:O	1:A:43:THR:CG2	0.42	2.68	17	3
1:A:135:HIS:C	1:A:136:TYR:CD2	0.42	2.97	17	1
1:A:90:ILE:O	1:A:90:ILE:CD1	0.42	2.68	25	1
1:A:62:LYS:HB2	1:A:65:LYS:CD	0.42	2.45	3	3
1:A:106:VAL:CG1	1:A:108:LYS:O	0.42	2.67	26	1
1:A:118:ASP:OD1	1:A:119:VAL:N	0.42	2.52	10	1
1:A:120:TYR:CE1	1:A:122:ASP:OD2	0.42	2.72	24	1
1:A:34:LEU:CD1	1:A:35:ASN:O	0.42	2.67	29	1
1:A:50:ASN:CG	1:A:51:ASP:N	0.42	2.78	14	2
1:A:88:GLU:OE1	1:A:102:SER:CB	0.42	2.68	20	1
1:A:50:ASN:HA	1:A:53:PHE:CE1	0.42	2.50	22	2
1:A:107:SER:O	1:A:108:LYS:C	0.42	2.63	2	7
1:A:125:VAL:O	1:A:125:VAL:CG1	0.42	2.68	6	1
1:A:117:ILE:C	1:A:117:ILE:CD1	0.42	2.92	15	1
1:A:119:VAL:HB	1:A:132:TYR:CD2	0.42	2.50	21	3
1:A:77:TRP:O	1:A:119:VAL:CG1	0.42	2.67	28	1
1:A:79:THR:CG2	1:A:80:GLY:N	0.42	2.81	4	2
1:A:102:SER:O	1:A:104:PHE:CE2	0.42	2.73	4	1
1:A:55:LYS:HA	1:A:56:PRO:C	0.42	2.40	6	3
1:A:36:TYR:CD1	1:A:57:ALA:HB2	0.42	2.50	16	1
1:A:130:PHE:CD1	1:A:130:PHE:C	0.42	2.95	16	2
1:A:84:GLU:OE1	1:A:84:GLU:CA	0.42	2.67	19	1
1:A:138:ILE:CD1	1:A:140:TYR:OH	0.42	2.68	7	1
1:A:32:GLY:O	1:A:59:TYR:N	0.42	2.52	11	2
1:A:40:LYS:CG	1:A:43:THR:OG1	0.42	2.68	11	1
1:A:81:MET:HE3	1:A:102:SER:HB3	0.42	1.91	14	1
1:A:69:GLN:NE2	1:A:101:THR:HG21	0.42	2.29	23	1
1:A:136:TYR:CD1	2:A:151:ZNH:HBD2	0.41	2.50	10	3
1:A:48:ILE:HG21	2:A:151:ZNH:CHD	0.41	2.44	21	1
1:A:50:ASN:OD1	1:A:51:ASP:N	0.41	2.53	21	1
1:A:46:THR:CG2	1:A:50:ASN:HB2	0.41	2.45	27	6
1:A:46:THR:CG2	1:A:50:ASN:HB3	0.41	2.45	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:38:VAL:O	1:A:38:VAL:CG2	0.41	2.67	8	1
1:A:110:ASN:O	1:A:110:ASN:ND2	0.41	2.53	11	1
1:A:73:ASN:O	1:A:74:HIS:CB	0.41	2.67	18	1
1:A:88:GLU:HB3	1:A:104:PHE:CE1	0.41	2.49	29	1
1:A:106:VAL:HG23	1:A:113:ILE:CD1	0.41	2.45	2	2
1:A:81:MET:CE	1:A:140:TYR:OH	0.41	2.68	3	1
1:A:105:GLU:CD	1:A:105:GLU:N	0.41	2.79	12	1
1:A:72:VAL:HG13	1:A:81:MET:HE2	0.41	1.91	16	1
1:A:72:VAL:O	1:A:72:VAL:CG2	0.41	2.66	20	2
1:A:100:ARG:HG3	1:A:101:THR:N	0.41	2.31	29	2
1:A:133:ASP:OD1	1:A:133:ASP:N	0.41	2.53	24	1
1:A:59:TYR:CD2	1:A:59:TYR:C	0.41	2.98	2	1
1:A:91:ILE:CG2	1:A:92:SER:N	0.41	2.84	7	4
1:A:113:ILE:N	1:A:140:TYR:O	0.41	2.53	19	2
1:A:121:ILE:HD12	1:A:132:TYR:CD2	0.41	2.50	18	2
1:A:50:ASN:ND2	1:A:51:ASP:OD1	0.41	2.54	21	1
1:A:67:TYR:C	1:A:68:VAL:CG1	0.41	2.91	24	1
1:A:52:TYR:O	1:A:53:PHE:CD1	0.41	2.73	27	1
1:A:34:LEU:HG	1:A:35:ASN:N	0.41	2.29	3	1
1:A:50:ASN:ND2	1:A:51:ASP:OD2	0.41	2.53	4	1
1:A:72:VAL:HG22	1:A:100:ARG:CB	0.41	2.45	7	1
1:A:62:LYS:HB3	1:A:65:LYS:CD	0.41	2.46	9	2
1:A:130:PHE:CZ	2:A:151:ZNH:HMB1	0.41	2.50	11	1
1:A:121:ILE:HG21	2:A:151:ZNH:CBB	0.41	2.46	22	1
1:A:61:LYS:CD	1:A:66:LEU:HB2	0.41	2.46	26	1
1:A:49:ALA:CB	1:A:138:ILE:HG21	0.41	2.46	8	1
1:A:79:THR:CG2	1:A:118:ASP:HB2	0.41	2.46	28	4
1:A:135:HIS:C	1:A:136:TYR:CG	0.41	2.99	16	1
1:A:122:ASP:OD1	1:A:122:ASP:N	0.41	2.53	13	3
1:A:83:ILE:HG13	1:A:140:TYR:CE2	0.41	2.51	29	2
1:A:36:TYR:CZ	1:A:53:PHE:CE2	0.41	3.09	20	1
1:A:40:LYS:HG3	1:A:43:THR:CG2	0.41	2.46	21	1
1:A:52:TYR:CD1	2:A:151:ZNH:HBC1	0.41	2.51	28	1
1:A:51:ASP:O	1:A:52:TYR:CD1	0.41	2.73	2	1
1:A:94:ASN:N	1:A:99:GLU:O	0.41	2.53	14	2
1:A:135:HIS:O	1:A:135:HIS:CD2	0.41	2.74	2	1
1:A:106:VAL:HG23	1:A:108:LYS:O	0.41	2.16	4	1
1:A:59:TYR:CB	1:A:68:VAL:HG23	0.41	2.45	12	1
1:A:62:LYS:O	1:A:63:ASN:O	0.41	2.39	29	5
1:A:135:HIS:ND1	1:A:135:HIS:N	0.41	2.68	19	1
1:A:38:VAL:HG22	1:A:139:THR:O	0.41	2.16	23	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:38:VAL:O	1:A:46:THR:CB	0.41	2.68	25	1
1:A:72:VAL:HB	1:A:78:ILE:HD13	0.41	1.93	29	1
1:A:132:TYR:OH	1:A:136:TYR:CZ	0.41	2.67	30	1
1:A:35:ASN:O	1:A:143:ASN:ND2	0.41	2.51	25	1
1:A:55:LYS:CB	1:A:56:PRO:HA	0.40	2.46	5	1
1:A:51:ASP:HB2	1:A:52:TYR:CE1	0.40	2.51	6	1
1:A:34:LEU:HD11	1:A:145:PRO:CD	0.40	2.46	13	1
1:A:60:ILE:HG22	1:A:67:TYR:O	0.40	2.16	18	1
1:A:61:LYS:HE2	1:A:109:LEU:HD21	0.40	1.92	28	1
1:A:78:ILE:CD1	1:A:81:MET:SD	0.40	3.09	30	2
1:A:67:TYR:C	1:A:68:VAL:CG2	0.40	2.91	20	2
1:A:61:LYS:NZ	1:A:145:PRO:HG2	0.40	2.31	15	1
1:A:80:GLY:C	1:A:81:MET:CG	0.40	2.94	16	1
1:A:61:LYS:CG	1:A:66:LEU:HB2	0.40	2.46	17	1
1:A:133:ASP:O	1:A:134:HIS:CG	0.40	2.74	18	1
1:A:113:ILE:HD12	1:A:142:PHE:HD2	0.40	1.74	24	1
1:A:66:LEU:HD21	1:A:106:VAL:HG13	0.40	1.92	26	1
1:A:77:TRP:CD1	1:A:77:TRP:N	0.40	2.88	26	1
1:A:105:GLU:OE2	1:A:106:VAL:N	0.40	2.55	26	1
1:A:83:ILE:O	1:A:83:ILE:HG22	0.40	2.15	12	1
1:A:36:TYR:HA	1:A:143:ASN:ND2	0.40	2.31	29	1
1:A:90:ILE:O	1:A:90:ILE:HD12	0.40	2.16	1	1
1:A:118:ASP:OD1	1:A:135:HIS:ND1	0.40	2.54	8	1
1:A:79:THR:OG1	1:A:120:TYR:CB	0.40	2.70	10	1
1:A:33:THR:HG23	1:A:58:LYS:HD2	0.40	1.93	18	1
1:A:82:SER:O	1:A:116:LYS:N	0.40	2.53	24	1
1:A:112:LYS:O	1:A:112:LYS:CE	0.40	2.69	26	1
1:A:65:LYS:N	1:A:65:LYS:HD3	0.40	2.32	28	1
1:A:40:LYS:HE3	1:A:47:SER:CB	0.40	2.47	11	1
1:A:52:TYR:C	1:A:53:PHE:CD1	0.40	2.99	12	1
1:A:65:LYS:C	1:A:66:LEU:CD1	0.40	2.92	16	1
1:A:97:LYS:O	1:A:98:ASP:CB	0.40	2.69	26	1
1:A:40:LYS:HG2	1:A:43:THR:HG21	0.40	1.94	27	1

6.3 Torsion angles ⓘ

6.3.1 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 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	116/147 (79%)	95±2 (82±2%)	17±2 (15±2%)	4±1 (3±1%)	5	38
All	All	3480/4410 (79%)	2856 (82%)	519 (15%)	105 (3%)	5	38

All 10 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	63	ASN	26
1	A	74	HIS	25
1	A	108	LYS	21
1	A	145	PRO	12
1	A	133	ASP	8
1	A	48	ILE	8
1	A	125	VAL	2
1	A	106	VAL	1
1	A	87	LYS	1
1	A	123	GLU	1

6.3.2 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 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	105/128 (82%)	61±3 (58±3%)	44±3 (42±3%)	0	4
All	All	3150/3840 (82%)	1833 (58%)	1317 (42%)	0	4

All 87 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	33	THR	30
1	A	43	THR	30
1	A	60	ILE	30
1	A	68	VAL	30
1	A	108	LYS	30
1	A	109	LEU	30
1	A	117	ILE	30

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Mol	Chain	Res	Type	Models (Total)
1	A	41	TYR	29
1	A	46	THR	29
1	A	70	ILE	29
1	A	91	ILE	29
1	A	34	LEU	28
1	A	90	ILE	28
1	A	40	LYS	27
1	A	79	THR	27
1	A	112	LYS	27
1	A	141	LYS	27
1	A	94	ASN	26
1	A	119	VAL	26
1	A	47	SER	25
1	A	84	GLU	24
1	A	50	ASN	23
1	A	55	LYS	23
1	A	58	LYS	23
1	A	138	ILE	23
1	A	38	VAL	22
1	A	82	SER	22
1	A	126	ASN	22
1	A	65	LYS	21
1	A	83	ILE	21
1	A	87	LYS	21
1	A	72	VAL	20
1	A	89	ASN	19
1	A	122	ASP	18
1	A	61	LYS	18
1	A	37	GLU	17
1	A	134	HIS	17
1	A	86	HIS	15
1	A	95	THR	15
1	A	128	LYS	15
1	A	30	ASP	14
1	A	62	LYS	14
1	A	98	ASP	13
1	A	74	HIS	13
1	A	66	LEU	13
1	A	67	TYR	13
1	A	107	SER	13
1	A	131	LYS	13
1	A	71	THR	12

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Mol	Chain	Res	Type	Models (Total)
1	A	75	SER	12
1	A	81	MET	12
1	A	135	HIS	11
1	A	69	GLN	10
1	A	123	GLU	10
1	A	44	ASN	9
1	A	48	ILE	9
1	A	63	ASN	9
1	A	97	LYS	9
1	A	100	ARG	9
1	A	76	HIS	8
1	A	124	LYS	8
1	A	42	ASN	7
1	A	103	GLU	7
1	A	105	GLU	7
1	A	54	ASN	7
1	A	92	SER	6
1	A	35	ASN	6
1	A	88	GLU	6
1	A	102	SER	6
1	A	73	ASN	6
1	A	31	SER	5
1	A	106	VAL	5
1	A	137	ASN	5
1	A	125	VAL	5
1	A	45	ASP	5
1	A	52	TYR	5
1	A	133	ASP	4
1	A	113	ILE	4
1	A	116	LYS	4
1	A	78	ILE	4
1	A	110	ASN	3
1	A	93	LYS	3
1	A	53	PHE	2
1	A	36	TYR	2
1	A	130	PHE	1
1	A	99	GLU	1
1	A	51	ASP	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no oligosaccharides 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	ZNH	A	151	1	50,50,50	2.18±0.00	21±0 (42±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
2	ZNH	A	151	1	65,82,82	1.29±0.02	5±0 (7±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	ZNH	A	151	1	1±0,1,9,9	0±0,14,54,54	-

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	151	ZNH	CHC-C1C	4.75	1.28	1.39	25	30
2	A	151	ZNH	CHC-C4B	4.67	1.29	1.38	29	30
2	A	151	ZNH	C1A-C2A	4.46	1.37	1.45	27	30
2	A	151	ZNH	CHD-C4C	4.16	1.30	1.39	9	30
2	A	151	ZNH	ZN-NA	3.85	1.89	2.03	24	30
2	A	151	ZNH	C4A-C3A	3.79	1.37	1.45	1	30
2	A	151	ZNH	CHB-C4A	3.71	1.31	1.38	14	30
2	A	151	ZNH	ZN-ND	3.66	1.91	2.03	15	30
2	A	151	ZNH	CHA-C1A	3.36	1.31	1.38	14	30
2	A	151	ZNH	C1B-C2B	3.26	1.37	1.44	6	30
2	A	151	ZNH	ZN-NC	3.22	1.91	2.03	16	30
2	A	151	ZNH	ZN-NB	3.16	1.92	2.03	8	30
2	A	151	ZNH	CHD-C1D	2.77	1.32	1.38	14	30
2	A	151	ZNH	CHA-C4D	2.75	1.32	1.39	13	30
2	A	151	ZNH	CHB-C1B	2.67	1.33	1.39	28	30
2	A	151	ZNH	C3D-C2D	2.53	1.31	1.36	12	30
2	A	151	ZNH	C4A-NA	2.30	1.43	1.38	8	30
2	A	151	ZNH	C1C-C2C	2.28	1.37	1.43	27	30
2	A	151	ZNH	C1A-NA	2.18	1.42	1.38	19	30
2	A	151	ZNH	C3B-C2B	2.17	1.32	1.37	11	30
2	A	151	ZNH	C2A-C3A	2.14	1.32	1.36	18	30
2	A	151	ZNH	C4C-NC	2.03	1.42	1.38	27	11

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(°)	Ideal(°)	Models	
								Worst	Total
2	A	151	ZNH	C1C-CHC-C4B	3.86	135.32	126.25	25	30
2	A	151	ZNH	C3B-C2B-C1B	2.92	108.60	106.41	14	30
2	A	151	ZNH	C4A-NA-C1A	2.91	102.19	106.52	4	30
2	A	151	ZNH	C3C-C2C-C1C	2.34	109.93	107.17	2	30
2	A	151	ZNH	CHD-C1D-ND	2.33	127.75	124.75	23	30

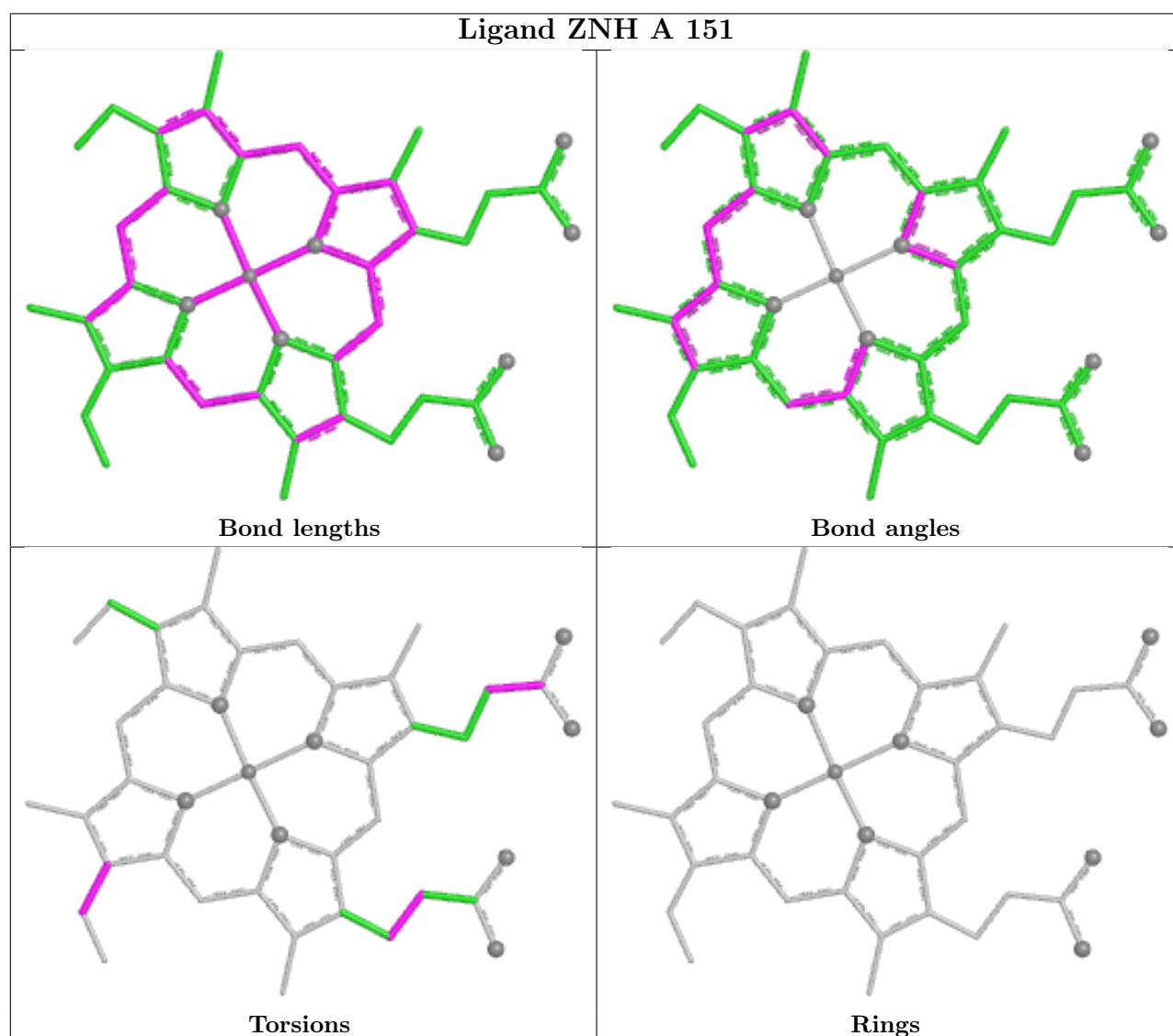
All unique chiral outliers are listed below.

Mol	Chain	Res	Type	Atoms	Models (Total)
2	A	151	ZNH	NA	30

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 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.



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

No chemical shift data were provided