



wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 1E8E / pdb_00001e8e
BMRB ID : 4837
Title : Solution Structure of Methylophilus methylotrophus Cytochrome c". Insights into the Structural Basis of Haem-Ligand Detachment
Authors : Brennan, L.; Turner, D.L.; Fareleira, P.; Santos, H.
Deposited on : 2000-09-20

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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 : 1.8.4, CSD as541be (2020)
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

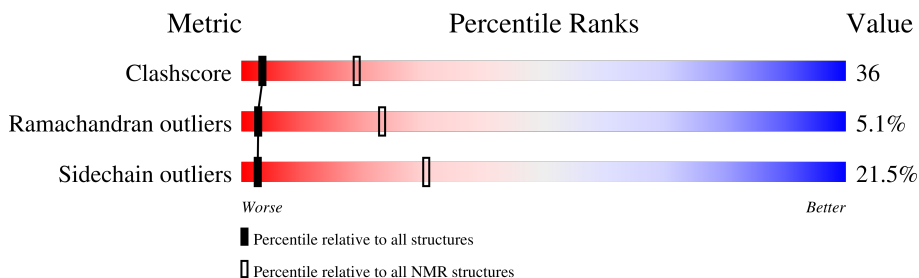
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 52%.

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	124	

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	HEC	125	20	-

2 Ensemble composition and analysis

This entry contains 20 models. Model 10 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative.

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:118 (116)	0.51	10

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 3 clusters and 3 single-model clusters were found.

Cluster number	Models
1	1, 2, 3, 4, 6, 7, 12, 19
2	5, 8, 9, 10, 13, 15, 17
3	11, 18
Single-model clusters	14; 16; 20

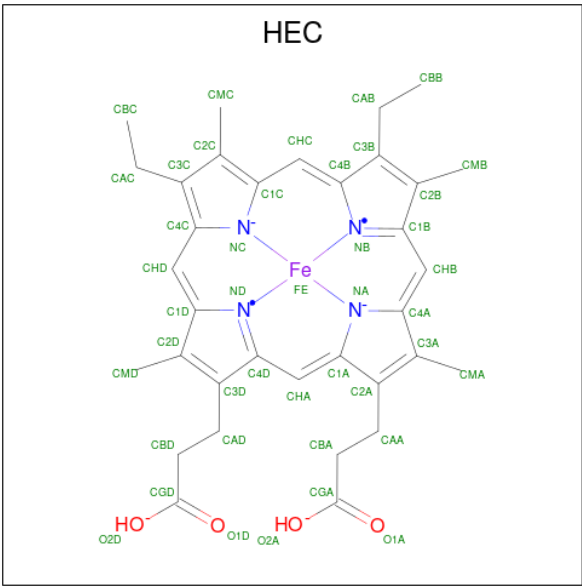
3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 1986 atoms, of which 982 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called CYTOCHROME C''.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	124	1911	604	950	164	188	5	0

- Molecule 2 is HEME C (CCD ID: HEC) (formula: C₃₄H₃₆FeN₄O₄).



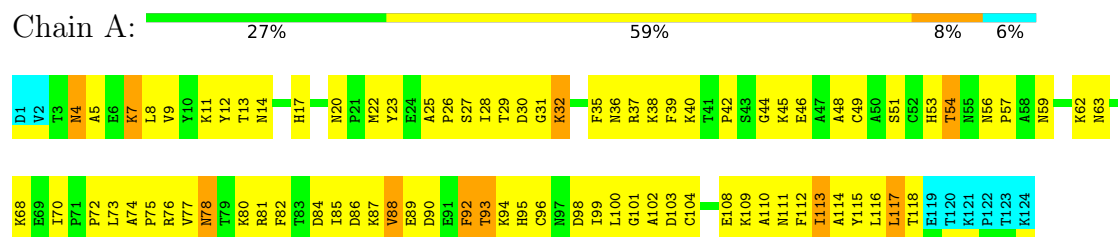
Mol	Chain	Residues	Atoms					
			Total	C	Fe	H	N	O
2	A	1	75	34	1	32	4	4

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.

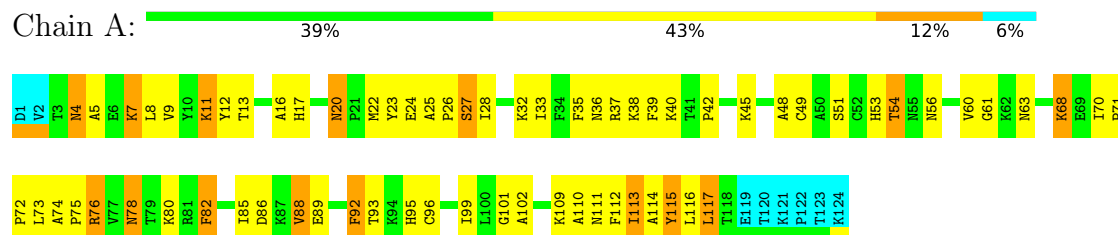
- Molecule 1: CYTOCHROME C''



4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 10. Colouring as in section 4.1 above.

- Molecule 1: CYTOCHROME C''



5 Refinement protocol and experimental data overview

The models were refined using the following method: *TORSION ANGLE DYNAMICS WITH SIMULATED ANNEALING*.

Of the 400 calculated structures, 20 were deposited, based on the following criterion: *LEAST RESTRAINT VIOLATION*.

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

Software name	Classification	Version
PARADYANA	refinement	
PARADYANA	structure solution	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	2
Total number of shifts	884
Number of shifts mapped to atoms	859
Number of unparsed shifts	0
Number of shifts with mapping errors	25
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	52%

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

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	0.66±0.00	0±0/918 (0.0± 0.0%)	1.27±0.01	1±1/1246 (0.1± 0.1%)
All	All	0.66	0/18360 (0.0%)	1.27	28/24920 (0.1%)

There are no bond-length outliers.

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	70	ILE	CA-C-N	6.49	124.33	119.66	20	1
1	A	70	ILE	C-N-CA	6.49	124.33	119.66	20	1
1	A	56	ASN	CA-C-N	5.05	124.35	118.85	7	13
1	A	56	ASN	C-N-CA	5.05	124.35	118.85	7	13

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	897	882	882	66±6
2	A	43	32	30	6±2
All	All	18800	18280	18240	1339

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 36.

5 of 342 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:73:LEU:HD12	1:A:73:LEU:O	0.92	1.65	20	2
1:A:92:PHE:CE1	2:A:125:HEC:HMA3	0.89	2.03	10	13
1:A:5:ALA:HB1	1:A:117:LEU:HD21	0.86	1.48	11	4
1:A:5:ALA:CB	1:A:117:LEU:HD21	0.86	2.00	6	3
1:A:63:ASN:ND2	1:A:70:ILE:HD11	0.86	1.85	8	18

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	116/124 (94%)	86±3 (74±3%)	24±3 (21±2%)	6±1 (5±1%)	3	23
All	All	2320/2480 (94%)	1714 (74%)	487 (21%)	119 (5%)	3	23

5 of 23 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	54	THR	20
1	A	38	LYS	18
1	A	44	GLY	13
1	A	45	LYS	9
1	A	68	LYS	9

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	99/107 (93%)	78±3 (78±3%)	21±3 (22±3%)	3	30
All	All	1980/2140 (93%)	1554 (78%)	426 (22%)	3	30

5 of 52 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	88	VAL	20
1	A	92	PHE	20
1	A	93	THR	20
1	A	113	ILE	20
1	A	7	LYS	19

6.3.3 RNA [i](#)

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	HEC	A	125	1	50,50,50	1.09±0.02	5±0 (10±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	HEC	A	125	1	68,82,82	0.66±0.01	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
2	HEC	A	125	1	1±0,1,7,7	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	125	HEC	CHC-C1C	3.77	1.47	1.39	11	20
2	A	125	HEC	FE-NB	3.25	2.04	1.94	6	20
2	A	125	HEC	FE-NC	2.78	2.04	1.95	7	20
2	A	125	HEC	FE-ND	2.61	2.02	1.94	20	20
2	A	125	HEC	FE-NA	2.21	2.02	1.95	11	20

There are no bond-angle outliers.

All unique chiral outliers are listed below.

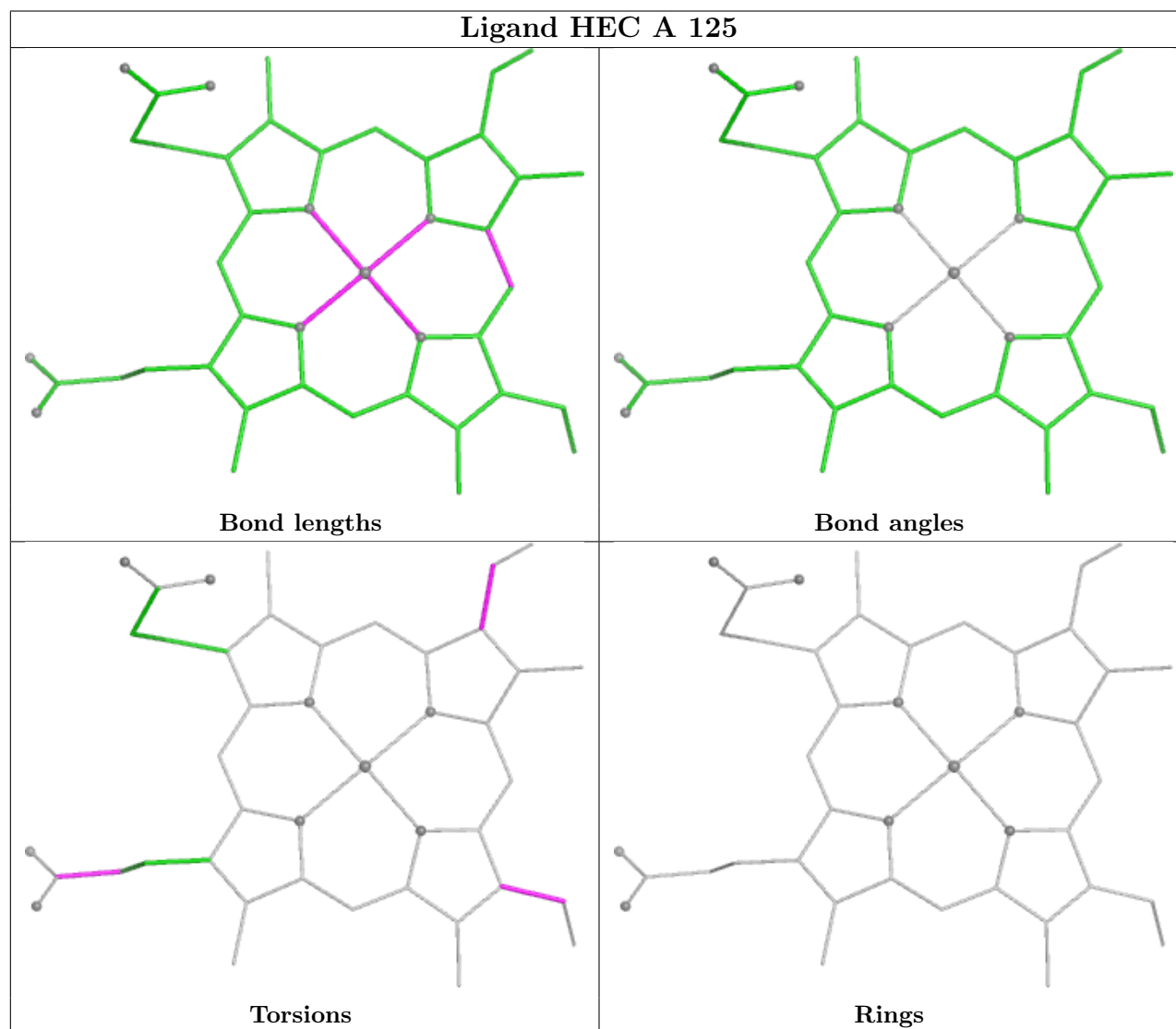
Mol	Chain	Res	Type	Atoms	Models (Total)
2	A	125	HEC	NA	20

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

The completeness of assignment taking into account all chemical shift lists is 52% for the well-defined parts and 52% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	857
Number of shifts mapped to atoms	857
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	52

7.1.2 Chemical shift referencing

No chemical shift referencing corrections were calculated (not enough data).

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 52%, i.e. 797 atoms were assigned a chemical shift out of a possible 1524. 0 out of 10 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	228/569 (40%)	228/229 (100%)	0/232 (0%)	0/108 (0%)
Sidechain	520/838 (62%)	520/542 (96%)	0/264 (0%)	0/32 (0%)
Aromatic	49/117 (42%)	49/58 (84%)	0/56 (0%)	0/3 (0%)
Overall	797/1524 (52%)	797/829 (96%)	0/552 (0%)	0/143 (0%)

7.1.4 Statistically unusual chemical shifts ⓘ

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	95	HIS	HB2	20.64	1.36 – 4.85	50.2
1	A	53	HIS	HB2	15.43	1.36 – 4.85	35.3
1	A	54	THR	HG1	6.87	0.08 – 2.19	27.2
1	A	95	HIS	HB3	11.89	1.18 – 4.91	23.7
1	A	118	THR	HG1	5.61	0.08 – 2.19	21.2
1	A	53	HIS	HB3	10.83	1.18 – 4.91	20.9
1	A	66	THR	HG1	4.99	0.08 – 2.19	18.3
1	A	41	THR	HG1	4.96	0.08 – 2.19	18.1
1	A	13	THR	HG1	4.63	0.08 – 2.19	16.6
1	A	53	HIS	HA	11.38	2.49 – 6.71	16.1
1	A	95	HIS	HA	11.24	2.49 – 6.71	15.7
1	A	73	LEU	HB2	5.90	-0.07 – 3.30	12.7
1	A	98	ASP	HB2	5.56	1.41 – 4.01	11.0
1	A	99	ILE	HG12	5.56	-0.69 – 3.24	10.9
1	A	108	GLU	HG3	0.07	1.20 – 3.30	-10.4
1	A	73	LEU	HD21	3.41	-0.65 – 2.13	9.6
1	A	73	LEU	HD22	3.41	-0.65 – 2.13	9.6
1	A	73	LEU	HD23	3.41	-0.65 – 2.13	9.6
1	A	109	LYS	HD2	-0.30	0.58 – 2.64	-9.3
1	A	109	LYS	HD3	-0.30	0.54 – 2.65	-9.0
1	A	99	ILE	HG21	-1.49	-0.56 – 2.11	-8.5
1	A	99	ILE	HG22	-1.49	-0.56 – 2.11	-8.5
1	A	99	ILE	HG23	-1.49	-0.56 – 2.11	-8.5
1	A	98	ASP	HB3	4.92	1.32 – 4.00	8.4
1	A	57	PRO	HA	6.90	2.78 – 6.00	7.8
1	A	94	LYS	HG2	3.26	0.13 – 2.61	7.6
1	A	94	LYS	HB2	3.58	0.58 – 2.97	7.5
1	A	35	PHE	HZ	9.98	4.94 – 9.06	7.2
1	A	54	THR	HG21	2.62	0.08 – 2.19	7.0
1	A	54	THR	HG22	2.62	0.08 – 2.19	7.0
1	A	54	THR	HG23	2.62	0.08 – 2.19	7.0
1	A	61	GLY	HA2	6.48	2.15 – 5.77	7.0
1	A	53	HIS	H	12.81	4.92 – 11.57	6.9
1	A	49	CYS	HB3	-0.03	0.69 – 5.10	-6.6
1	A	49	CYS	HB2	0.22	0.81 – 5.11	-6.4
1	A	61	GLY	HA3	6.10	2.08 – 5.71	6.1
1	A	95	HIS	H	12.26	4.92 – 11.57	6.0
1	A	81	ARG	HD3	1.57	1.81 – 4.39	-5.9

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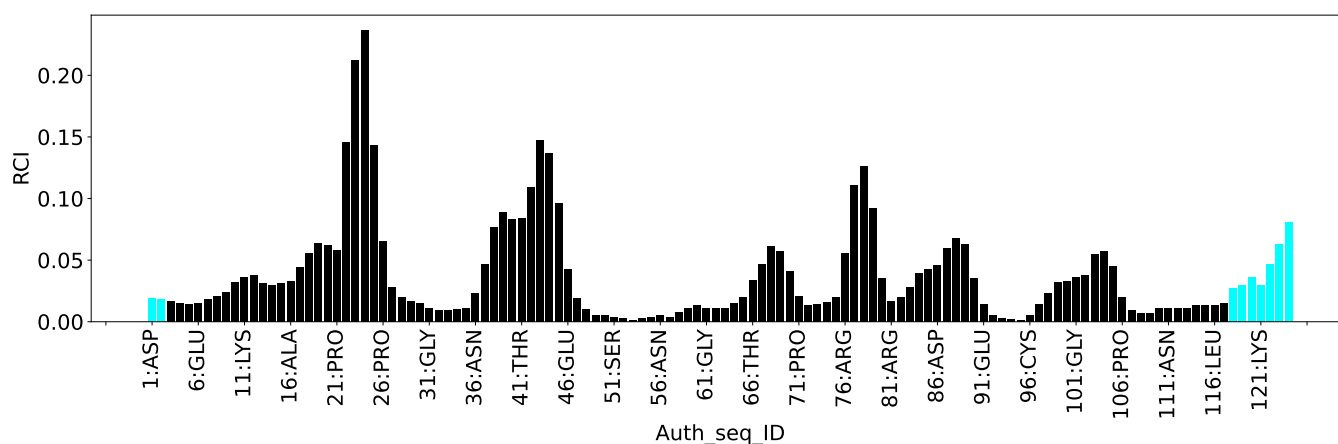
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List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	94	LYS	HB3	3.27	0.46 – 3.04	5.9
1	A	35	PHE	HE1	8.87	5.56 – 8.62	5.8
1	A	35	PHE	HE2	8.87	5.54 – 8.63	5.8
1	A	75	PRO	HG2	0.24	0.41 – 3.45	-5.5
1	A	75	PRO	HB3	0.09	0.25 – 3.76	-5.5
1	A	96	CYS	HB2	5.19	0.81 – 5.11	5.2
1	A	100	LEU	HD21	-0.70	-0.65 – 2.13	-5.2
1	A	100	LEU	HD22	-0.70	-0.65 – 2.13	-5.2
1	A	100	LEU	HD23	-0.70	-0.65 – 2.13	-5.2
1	A	47	ALA	HB1	0.12	0.14 – 2.58	-5.1
1	A	47	ALA	HB2	0.12	0.14 – 2.58	-5.1
1	A	47	ALA	HB3	0.12	0.14 – 2.58	-5.1
1	A	39	PHE	HZ	4.92	4.94 – 9.06	-5.0
1	A	94	LYS	HG3	2.68	0.04 – 2.67	5.0

7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:



7.2 Chemical shift list 2

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_2*

7.2.1 Bookkeeping ⓘ

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	27
Number of shifts mapped to atoms	2
Number of unparsed shifts	0
Number of shifts with mapping errors	25
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	2

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- No matching atom found in the structure. First 5 (of 25) occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
2	A	1	ASP	1HAA	8.11	?	2
2	A	1	ASP	1HAD	9.61	?	1
2	A	1	ASP	1HBC	0.21	?	1
2	A	1	ASP	1HBD	-2.16	?	2
2	A	1	ASP	1HMA	14.56	?	1
2	A	1	ASP	1HMB	8.62	?	1
2	A	1	ASP	1HMC	17.57	?	1
2	A	1	ASP	1HMD	12.8	?	1
2	A	1	ASP	2HAA	4.63	?	2
2	A	1	ASP	2HAD	3.83	?	1
2	A	1	ASP	2HBC	0.21	?	1
2	A	1	ASP	2HBD	0.62	?	2
2	A	1	ASP	2HMA	14.56	?	1
2	A	1	ASP	2HMB	8.62	?	1
2	A	1	ASP	2HMC	17.57	?	1
2	A	1	ASP	2HMD	12.8	?	1
2	A	1	ASP	3HBC	0.21	?	1
2	A	1	ASP	3HMA	14.56	?	1
2	A	1	ASP	3HMB	8.62	?	1
2	A	1	ASP	3HMC	17.57	?	1
2	A	1	ASP	3HMD	12.8	?	1
2	A	1	ASP	HAC	-0.42	?	1
2	A	1	ASP	HHA	1.57	?	1
2	A	1	ASP	HHB	-2.38	?	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
2	A	1	ASP	HHC	2.43	?	1

7.2.2 Chemical shift referencing [i](#)

No chemical shift referencing corrections were calculated (not enough data).

7.2.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 0%, i.e. 0 atoms were assigned a chemical shift out of a possible 1524. 0 out of 10 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	0/569 (0%)	0/229 (0%)	0/232 (0%)	0/108 (0%)
Sidechain	0/838 (0%)	0/542 (0%)	0/264 (0%)	0/32 (0%)
Aromatic	0/117 (0%)	0/58 (0%)	0/56 (0%)	0/3 (0%)
Overall	0/1524 (0%)	0/829 (0%)	0/552 (0%)	0/143 (0%)

7.2.4 Statistically unusual chemical shifts [i](#)

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
2	A	1	ASP	HA	-2.27	3.04 – 6.12	-22.2
2	A	1	ASP	HB2	-0.09	1.41 – 4.01	-10.8

7.2.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:

