



# Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 11:23 PM BST

PDB ID : 2KOR  
Title : NMR solution structures of 2-octenoyl-ACP from *Streptomyces coelicolor*  
Fatty Acid Synthase  
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Deposited on : 2009-09-29

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

Cyrange : Kirchner and Güntert (2011)  
NmrClust : Kelley et al. (1996)  
MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
RCI : v\_1n\_11\_5\_13\_A (Berjanski et al., 2005)  
PANAV : Wang et al. (2010)  
ShiftChecker : rb-20027457  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027457

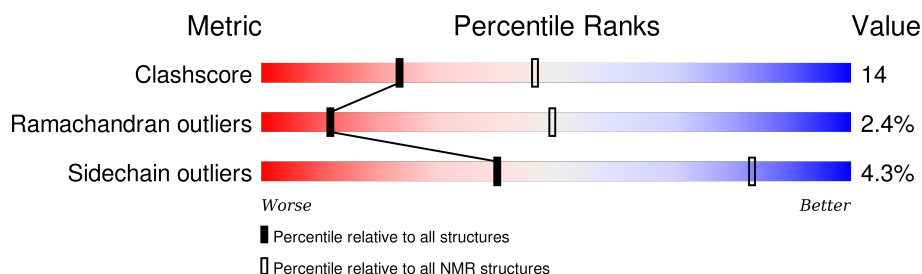
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 84%.

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	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

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  $\geq 3$ , 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	81	<div> <div>62%</div> <div>36%</div> <div>.</div> </div>

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	SOO	101	20	-

## 2 Ensemble composition and analysis

This entry contains 20 models. Model 8 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

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:2-A:80 (79)	0.32	8

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 4 clusters and 1 single-model cluster was found.

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

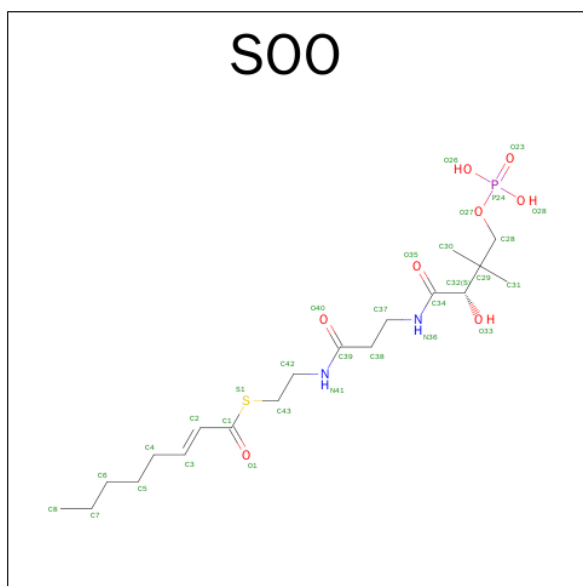
### 3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 1291 atoms, of which 644 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Acyl carrier protein.

Mol	Chain	Residues	Atoms						Trace
1	A	81	Total	C	H	N	O	S	0
			1228	386	611	96	134	1	

- Molecule 2 is S-[2-[3-[[[(2S)-2-HYDROXY-3,3-DIMETHYL-4-PHOSPHONOOXY-BUTANOYL]AMINO]PROPANOYLAMINO]ETHYL] (E)-OCT-2-ENETHIOATE (three-letter code: SOO) (formula: C<sub>19</sub>H<sub>35</sub>N<sub>2</sub>O<sub>8</sub>PS).

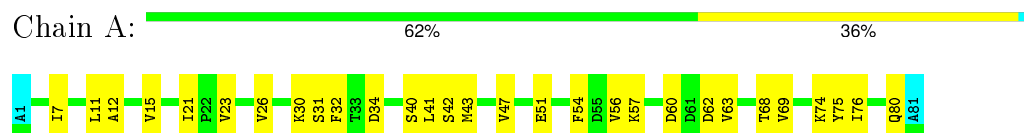


## 4 Residue-property plots

### 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA 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: Acyl carrier protein

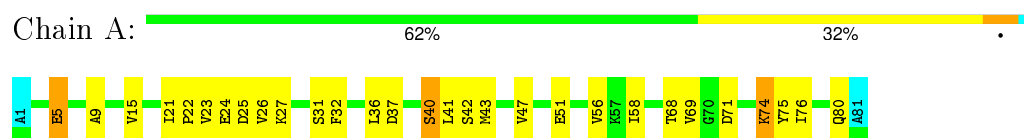


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

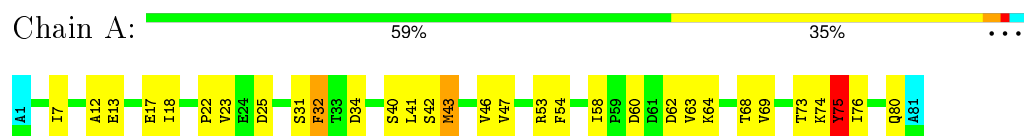
#### 4.2.1 Score per residue for model 1

- Molecule 1: Acyl carrier protein



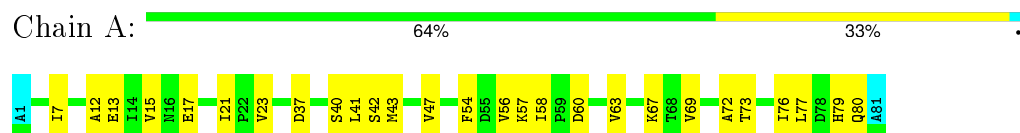
#### 4.2.2 Score per residue for model 2

- Molecule 1: Acyl carrier protein



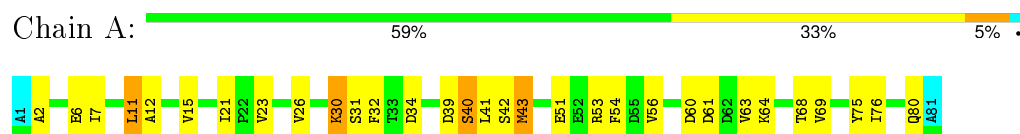
### 4.2.3 Score per residue for model 3

- Molecule 1: Acyl carrier protein



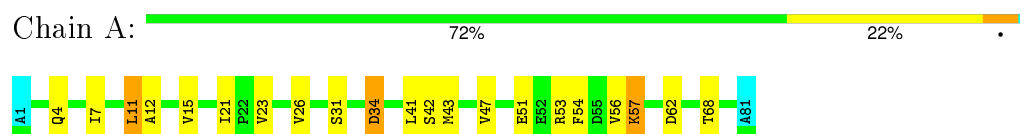
### 4.2.4 Score per residue for model 4

- Molecule 1: Acyl carrier protein



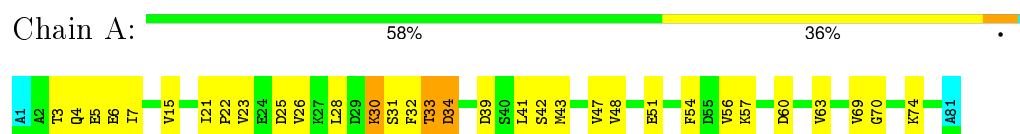
### 4.2.5 Score per residue for model 5

- Molecule 1: Acyl carrier protein



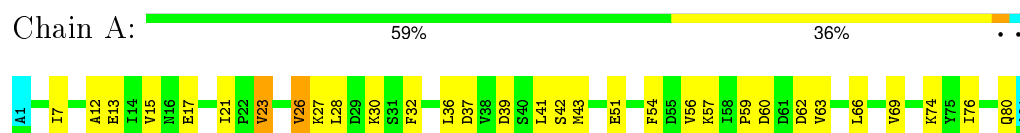
### 4.2.6 Score per residue for model 6

- Molecule 1: Acyl carrier protein



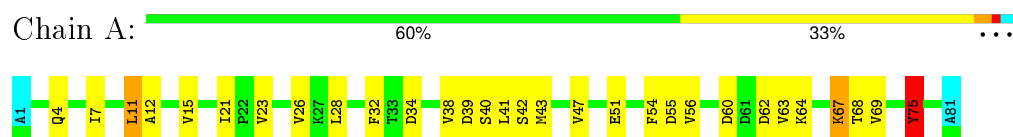
### 4.2.7 Score per residue for model 7

- Molecule 1: Acyl carrier protein



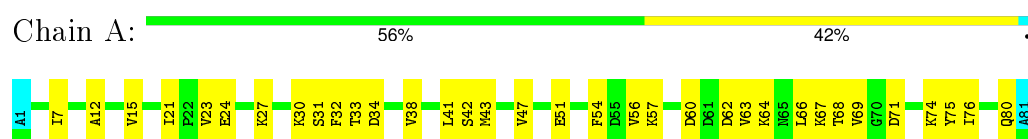
#### 4.2.8 Score per residue for model 8 (medoid)

- Molecule 1: Acyl carrier protein



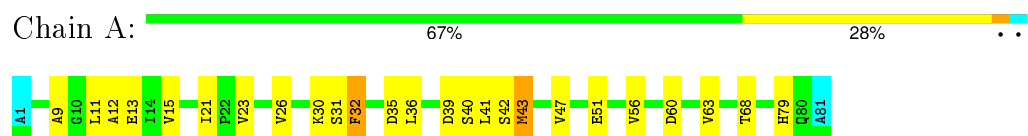
#### 4.2.9 Score per residue for model 9

- Molecule 1: Acyl carrier protein



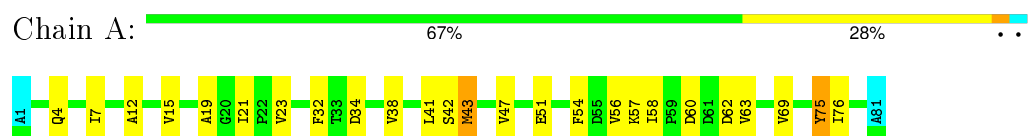
#### 4.2.10 Score per residue for model 10

- Molecule 1: Acyl carrier protein



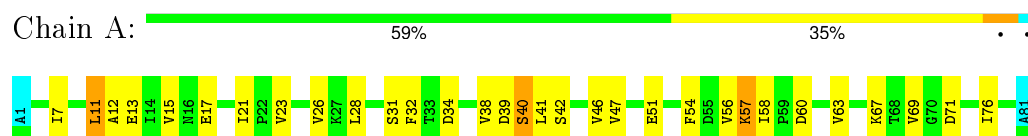
#### 4.2.11 Score per residue for model 11

- Molecule 1: Acyl carrier protein



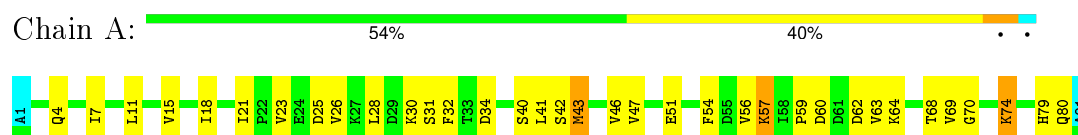
#### 4.2.12 Score per residue for model 12

- Molecule 1: Acyl carrier protein



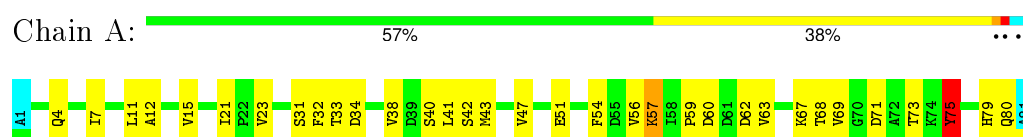
### 4.2.13 Score per residue for model 13

- Molecule 1: Acyl carrier protein



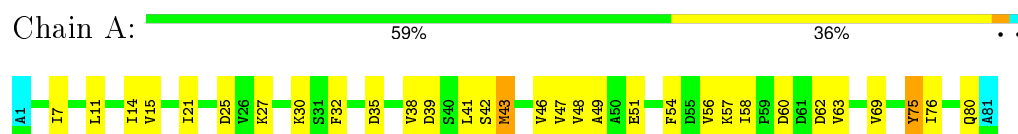
### 4.2.14 Score per residue for model 14

- Molecule 1: Acyl carrier protein



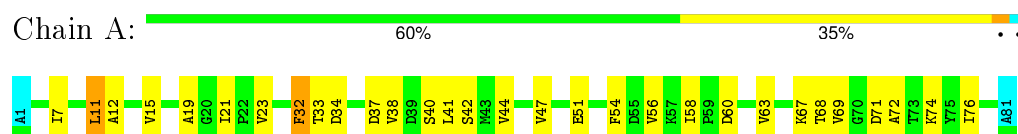
### 4.2.15 Score per residue for model 15

- Molecule 1: Acyl carrier protein



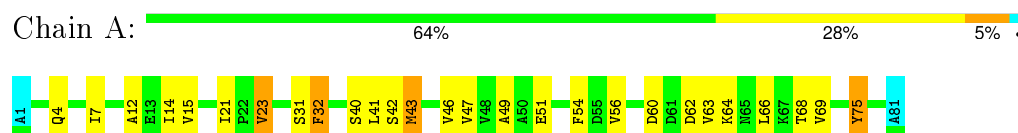
### 4.2.16 Score per residue for model 16

- Molecule 1: Acyl carrier protein



### 4.2.17 Score per residue for model 17

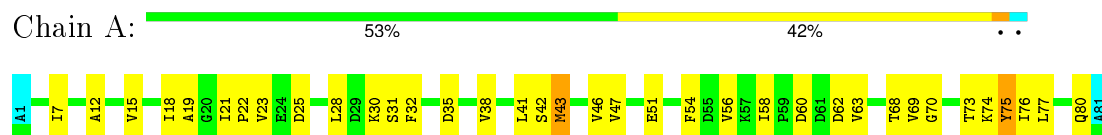
- Molecule 1: Acyl carrier protein





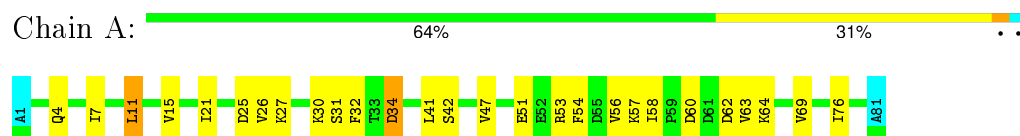
#### 4.2.18 Score per residue for model 18

- Molecule 1: Acyl carrier protein



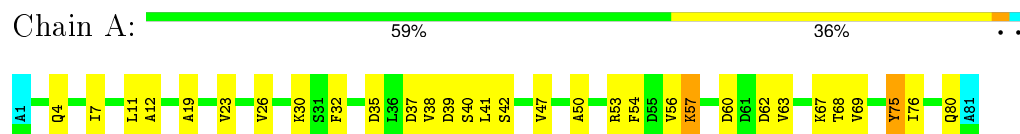
#### 4.2.19 Score per residue for model 19

- Molecule 1: Acyl carrier protein



#### 4.2.20 Score per residue for model 20

- Molecule 1: Acyl carrier protein



## 5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 20 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
ARIA	refinement	1.2
ARIA	structure solution	1.2

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)	BMRB entry 16527
Number of chemical shift lists	1
Total number of shifts	949
Number of shifts mapped to atoms	917
Number of unparsed shifts	0
Number of shifts with mapping errors	32
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	84%

No validations of the models with respect to experimental NMR restraints is performed at this time.

## 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: SOO

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	0.79±0.10	1±1/612 (0.1±0.2%)	0.73±0.03	0±0/833 (0.0±0.0%)
All	All	0.79	15/12240 (0.1%)	0.73	2/16660 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	0.2±0.4
All	All	0	5

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	75	TYR	CE1-CZ	13.52	1.56	1.38	17	8
1	A	75	TYR	CE2-CZ	-12.54	1.22	1.38	17	7

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	75	TYR	CE1-CZ-OH	-5.46	105.35	120.10	17	1
1	A	75	TYR	OH-CZ-CE2	-5.12	106.28	120.10	14	1

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	75	TYR	Sidechain	4
1	A	53	ARG	Sidechain	1

## 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	606	599	598	18±3
2	A	30	33	33	2±1
All	All	12720	12640	12618	367

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 14.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:56:VAL:HG11	1:A:76:ILE:HG23	0.77	1.56	4	2
1:A:30:LYS:HE3	1:A:30:LYS:HA	0.75	1.58	4	2
1:A:7:ILE:HA	1:A:54:PHE:CE1	0.74	2.17	15	18
1:A:19:ALA:HB2	1:A:38:VAL:HG13	0.71	1.61	18	4
1:A:33:THR:HG23	2:A:101:SOO:S1	0.67	2.28	6	1
1:A:12:ALA:HB1	1:A:23:VAL:HA	0.67	1.64	11	15
1:A:56:VAL:HG12	1:A:80:GLN:HB3	0.65	1.66	4	1
1:A:47:VAL:HG22	2:A:101:SOO:H6	0.64	1.70	11	5
1:A:39:ASP:HB3	1:A:42:SER:HB3	0.64	1.68	7	2
1:A:15:VAL:HG12	1:A:21:ILE:HB	0.63	1.71	18	17
1:A:58:ILE:HD11	1:A:76:ILE:HG12	0.60	1.73	3	9
1:A:67:LYS:HG3	1:A:68:THR:HG23	0.60	1.72	16	1
1:A:56:VAL:HG23	1:A:80:GLN:HB3	0.60	1.74	3	1
1:A:28:LEU:HA	1:A:69:VAL:HG22	0.60	1.74	12	3
1:A:47:VAL:HG21	1:A:63:VAL:HG21	0.60	1.74	12	2
1:A:62:ASP:OD2	1:A:75:TYR:OH	0.59	2.19	8	5
1:A:43:MET:SD	2:A:101:SOO:H37	0.59	2.37	15	4
1:A:18:ILE:HG13	1:A:46:VAL:HG22	0.59	1.74	2	2
1:A:31:SER:HA	1:A:68:THR:HA	0.59	1.73	4	9
1:A:12:ALA:HB2	1:A:26:VAL:HB	0.59	1.75	5	3
1:A:51:GLU:HA	1:A:56:VAL:HG12	0.58	1.75	18	9

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:40:SER:O	1:A:43:MET:HG3	0.58	1.99	4	4
1:A:39:ASP:HB3	1:A:42:SER:HB2	0.58	1.76	10	5
1:A:25:ASP:O	1:A:30:LYS:HE2	0.58	1.99	13	1
1:A:30:LYS:HE3	1:A:35:ASP:HB3	0.57	1.75	10	3
1:A:14:ILE:HG22	1:A:46:VAL:HG23	0.57	1.76	15	2
1:A:26:VAL:HG22	1:A:69:VAL:HG11	0.57	1.75	7	1
1:A:43:MET:HB2	2:A:101:SOO:H4	0.57	1.74	13	2
1:A:43:MET:HB2	2:A:101:SOO:H4A	0.56	1.76	9	1
1:A:60:ASP:HA	1:A:63:VAL:HG12	0.56	1.77	15	17
1:A:48:VAL:HA	1:A:51:GLU:OE1	0.56	2.00	15	2
1:A:67:LYS:HG2	1:A:71:ASP:OD2	0.56	2.00	16	1
1:A:43:MET:O	1:A:47:VAL:HG23	0.56	2.00	15	7
1:A:32:PHE:HA	1:A:36:LEU:HD12	0.56	1.77	1	3
1:A:30:LYS:HE2	1:A:35:ASP:HB3	0.55	1.77	18	1
1:A:51:GLU:HA	1:A:56:VAL:CG2	0.55	2.30	4	1
1:A:56:VAL:HG12	1:A:57:LYS:H	0.55	1.60	20	1
1:A:76:ILE:O	1:A:80:GLN:HG2	0.55	2.01	15	2
1:A:26:VAL:HG12	1:A:69:VAL:HG11	0.55	1.79	6	7
1:A:71:ASP:HA	1:A:74:LYS:HE2	0.55	1.78	1	1
1:A:28:LEU:O	1:A:68:THR:HB	0.54	2.02	13	1
1:A:22:PRO:HB2	1:A:25:ASP:HB2	0.54	1.79	6	4
1:A:56:VAL:HG22	1:A:57:LYS:H	0.54	1.63	15	7
1:A:56:VAL:HG12	1:A:80:GLN:CB	0.53	2.33	4	1
1:A:62:ASP:OD1	1:A:75:TYR:OH	0.53	2.25	15	2
1:A:47:VAL:HG21	1:A:63:VAL:HG11	0.53	1.80	9	1
1:A:56:VAL:HG22	1:A:80:GLN:HB3	0.53	1.81	20	1
1:A:24:GLU:O	1:A:27:LYS:HE3	0.52	2.05	1	2
1:A:41:LEU:HD12	1:A:42:SER:N	0.52	2.20	3	20
2:A:101:SOO:HN36	2:A:101:SOO:H31	0.51	1.65	10	1
1:A:7:ILE:HG21	1:A:73:THR:HA	0.51	1.82	2	2
1:A:62:ASP:HB3	1:A:75:TYR:OH	0.51	2.05	20	2
1:A:28:LEU:HA	1:A:69:VAL:CG2	0.51	2.35	7	3
1:A:15:VAL:CG1	1:A:21:ILE:HB	0.51	2.36	8	9
1:A:60:ASP:O	1:A:64:LYS:HG3	0.51	2.05	9	6
1:A:31:SER:HB2	1:A:34:ASP:HB3	0.50	1.83	5	3
1:A:64:LYS:HG2	2:A:101:SOO:O40	0.50	2.06	17	1
1:A:32:PHE:CZ	1:A:69:VAL:HG12	0.50	2.41	8	16
1:A:51:GLU:O	1:A:55:ASP:HA	0.50	2.06	8	1
1:A:38:VAL:HG11	1:A:46:VAL:HG21	0.50	1.83	12	1
1:A:41:LEU:HA	2:A:101:SOO:H38	0.50	1.84	4	1
1:A:38:VAL:HB	1:A:43:MET:HB3	0.50	1.83	8	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:28:LEU:HA	1:A:69:VAL:HG23	0.49	1.82	7	1
1:A:56:VAL:HG21	1:A:76:ILE:HG23	0.49	1.84	20	2
1:A:40:SER:HB2	2:A:101:SOO:H28	0.49	1.84	12	1
1:A:25:ASP:HA	1:A:27:LYS:NZ	0.48	2.23	19	1
1:A:67:LYS:HG2	1:A:68:THR:H	0.48	1.69	8	1
1:A:56:VAL:HA	1:A:80:GLN:HA	0.48	1.84	13	4
1:A:13:GLU:O	1:A:17:GLU:HG2	0.48	2.08	7	4
1:A:43:MET:HG3	2:A:101:SOO:C2	0.48	2.38	2	2
1:A:3:THR:H	1:A:6:GLU:HB3	0.48	1.69	6	1
1:A:32:PHE:CZ	2:A:101:SOO:H6	0.48	2.43	20	1
1:A:67:LYS:HG2	1:A:68:THR:HG23	0.48	1.85	20	1
1:A:27:LYS:HB2	1:A:30:LYS:HG2	0.48	1.86	19	1
1:A:31:SER:CB	1:A:34:ASP:HB3	0.47	2.38	5	1
1:A:47:VAL:O	1:A:51:GLU:HG3	0.47	2.09	19	7
1:A:43:MET:HE3	2:A:101:SOO:H2	0.47	1.85	18	1
1:A:7:ILE:O	1:A:11:LEU:HB2	0.47	2.10	12	5
1:A:11:LEU:HD21	2:A:101:SOO:H8B	0.47	1.87	16	1
1:A:31:SER:HB3	1:A:34:ASP:HB3	0.47	1.85	13	2
1:A:7:ILE:HA	1:A:54:PHE:CZ	0.47	2.44	12	2
1:A:2:ALA:HB1	1:A:6:GLU:CB	0.47	2.40	4	1
1:A:14:ILE:HG21	1:A:49:ALA:HB3	0.47	1.87	15	2
1:A:73:THR:O	1:A:77:LEU:HG	0.47	2.10	18	2
1:A:11:LEU:HD21	2:A:101:SOO:H8A	0.47	1.87	8	1
1:A:5:GLU:O	1:A:9:ALA:HB3	0.46	2.10	1	1
1:A:27:LYS:HB2	1:A:30:LYS:HG3	0.46	1.87	7	1
1:A:51:GLU:HA	1:A:56:VAL:CG1	0.46	2.41	13	4
1:A:63:VAL:O	1:A:66:LEU:HG	0.45	2.11	7	1
1:A:70:GLY:O	1:A:74:LYS:HD3	0.45	2.12	6	2
1:A:38:VAL:HG13	1:A:43:MET:HB2	0.45	1.86	14	1
1:A:56:VAL:C	1:A:57:LYS:HD2	0.45	2.32	11	4
1:A:43:MET:HG2	2:A:101:SOO:H2	0.44	1.89	4	1
1:A:2:ALA:HB1	1:A:6:GLU:HB2	0.44	1.89	4	1
1:A:47:VAL:HG22	2:A:101:SOO:H6A	0.44	1.90	1	1
1:A:43:MET:O	1:A:46:VAL:HG12	0.44	2.12	17	1
1:A:43:MET:C	1:A:43:MET:SD	0.43	2.97	4	2
2:A:101:SOO:HN41	2:A:101:SOO:HN36	0.43	1.56	1	1
1:A:11:LEU:HD21	2:A:101:SOO:H7	0.43	1.89	4	1
1:A:69:VAL:HA	1:A:72:ALA:HB3	0.43	1.90	16	2
1:A:44:VAL:O	1:A:47:VAL:HB	0.43	2.14	16	1
1:A:32:PHE:CD1	2:A:101:SOO:H4	0.43	2.49	12	1
1:A:33:THR:HG22	2:A:101:SOO:S1	0.43	2.53	16	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:59:PRO:HG2	1:A:62:ASP:HB2	0.43	1.90	14	3
1:A:38:VAL:HG11	1:A:46:VAL:CG2	0.43	2.44	12	1
1:A:51:GLU:HA	1:A:56:VAL:HG22	0.43	1.91	4	1
1:A:50:ALA:O	1:A:53:ARG:HB3	0.42	2.14	20	1
1:A:43:MET:HE2	2:A:101:SOO:H2	0.42	1.90	13	1
1:A:66:LEU:HA	1:A:71:ASP:OD2	0.42	2.15	9	1
1:A:30:LYS:HE2	1:A:30:LYS:HA	0.42	1.90	9	1
1:A:71:ASP:O	1:A:75:TYR:HB2	0.42	2.14	1	2
1:A:54:PHE:HB2	1:A:56:VAL:HG23	0.42	1.90	20	1
1:A:25:ASP:HA	1:A:27:LYS:HE3	0.42	1.92	15	1
1:A:43:MET:SD	2:A:101:SOO:H4A	0.42	2.54	4	1
1:A:25:ASP:O	1:A:30:LYS:HG3	0.42	2.15	18	1
1:A:39:ASP:HB3	1:A:42:SER:CB	0.41	2.45	20	1
1:A:56:VAL:HG13	1:A:80:GLN:HA	0.41	1.92	20	1
1:A:76:ILE:HG22	1:A:80:GLN:HE21	0.41	1.75	7	1
1:A:61:ASP:HA	1:A:64:LYS:HD3	0.41	1.93	4	1
1:A:63:VAL:HG23	1:A:66:LEU:HD12	0.41	1.90	17	1
1:A:56:VAL:HB	1:A:80:GLN:HB3	0.41	1.92	9	1
1:A:70:GLY:O	1:A:74:LYS:HG2	0.41	2.16	18	1
1:A:32:PHE:CG	2:A:101:SOO:H4	0.41	2.50	7	1
1:A:9:ALA:O	1:A:13:GLU:HG2	0.40	2.17	10	1
1:A:54:PHE:HB2	1:A:56:VAL:HG13	0.40	1.92	4	1
1:A:67:LYS:HB3	1:A:71:ASP:OD1	0.40	2.16	12	1
1:A:71:ASP:O	1:A:74:LYS:HD3	0.40	2.16	1	1
1:A:18:ILE:HD12	1:A:46:VAL:HA	0.40	1.92	18	1
1:A:5:GLU:CD	1:A:5:GLU:H	0.40	2.19	6	1
1:A:32:PHE:CE1	1:A:69:VAL:HG12	0.40	2.52	17	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	79/81 (98%)	66±2 (83±2%)	11±2 (14±3%)	2±1 (2±1%)	12	49
All	All	1580/1620 (98%)	1319 (83%)	223 (14%)	38 (2%)	12	49

All 9 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	40	SER	10
1	A	34	ASP	10
1	A	37	ASP	5
1	A	23	VAL	5
1	A	32	PHE	4
1	A	26	VAL	1
1	A	57	LYS	1
1	A	28	LEU	1
1	A	39	ASP	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	69/69 (100%)	66±1 (96±2%)	3±1 (4±2%)	40	83
All	All	1380/1380 (100%)	1320 (96%)	60 (4%)	40	83

All 13 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	11	LEU	11
1	A	43	MET	10
1	A	4	GLN	9
1	A	74	LYS	6
1	A	79	HIS	4
1	A	57	LYS	4
1	A	67	LYS	4
1	A	53	ARG	3
1	A	33	THR	3
1	A	62	ASP	2
1	A	30	LYS	2
1	A	51	GLU	1
1	A	5	GLU	1



### 6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 6.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.6 Ligand geometry ⓘ

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	SOO	A	101	1	20,29,30	0.77±0.13	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
2	SOO	A	101	1	26,36,39	1.09±0.09	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	SOO	A	101	1	1±0,1,10,10	2±0,33,36,37	0±0,0,0,0

There are no bond-length outliers.

There are no bond-angle outliers.

All unique chiral outliers are listed below.

Mol	Chain	Res	Type	Atoms	Models (Total)
2	A	101	SOO	C32	20

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

Mol	Chain	Res	Type	Atoms	Models (Total)
2	A	101	SOO	O1-C1-S1-C43	17
2	A	101	SOO	C2-C1-S1-C43	14

There are no ring outliers.

## 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 84% for the well-defined parts and 84% for the entire structure.

### 7.1 Chemical shift list 1

File name: BMRB entry 16527

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	949
Number of shifts mapped to atoms	917
Number of unparsed shifts	0
Number of shifts with mapping errors	32
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

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

- Chain not found in structure. All 32 occurrences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
UNMAPPED	1	SOD	H4B	1.58	0.0	1
UNMAPPED	1	SOD	H36A	8.07	0.0	1
UNMAPPED	1	SOD	H6B	0.81	0.0	1
UNMAPPED	1	SOD	H28B	3.3	0.0	2
UNMAPPED	1	SOD	H28A	3.74	0.0	2
UNMAPPED	1	SOD	H5A	1.02	0.0	1
UNMAPPED	1	SOD	H8A	0.63	0.0	1
UNMAPPED	1	SOD	H31A	0.87	0.0	2
UNMAPPED	1	SOD	H37B	3.6	0.0	1
UNMAPPED	1	SOD	H42B	3.29	0.0	1
UNMAPPED	1	SOD	H3A	6.02	0.0	1
UNMAPPED	1	SOD	H37A	3.39	0.0	1
UNMAPPED	1	SOD	H42A	3.11	0.0	1
UNMAPPED	1	SOD	H38B	2.46	0.0	1

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Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
UNMAPPED	1	SOD	H41A	8.24	0.0	1
UNMAPPED	1	SOD	H31C	0.87	0.0	2
UNMAPPED	1	SOD	H4A	1.76	0.0	1
UNMAPPED	1	SOD	H5B	0.79	0.0	1
UNMAPPED	1	SOD	H30A	0.87	0.0	2
UNMAPPED	1	SOD	H8B	0.63	0.0	1
UNMAPPED	1	SOD	H43A	2.84	0.0	1
UNMAPPED	1	SOD	H31B	0.87	0.0	2
UNMAPPED	1	SOD	H7B	0.96	0.0	1
UNMAPPED	1	SOD	H32A	4.04	0.0	1
UNMAPPED	1	SOD	H8C	0.63	0.0	1
UNMAPPED	1	SOD	H30C	0.87	0.0	2
UNMAPPED	1	SOD	H2A	6.63	0.0	2
UNMAPPED	1	SOD	H43B	2.95	0.0	1
UNMAPPED	1	SOD	H7A	0.96	0.0	1
UNMAPPED	1	SOD	H30B	0.87	0.0	2
UNMAPPED	1	SOD	H38A	2.46	0.0	1
UNMAPPED	1	SOD	H6A	0.75	0.0	1

### 7.1.2 Chemical shift referencing ⓘ

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	76	$-0.58 \pm 0.22$	Should be applied
$^{13}\text{C}_\beta$	78	$-0.16 \pm 0.26$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}'$	0	—	—
$^{15}\text{N}$	78	$0.86 \pm 0.48$	None needed (imprecise)

### 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 84%, i.e. 771 atoms were assigned a chemical shift out of a possible 916. 0 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	307/391 (79%)	156/156 (100%)	74/158 (47%)	77/77 (100%)
Sidechain	436/491 (89%)	268/277 (97%)	165/201 (82%)	3/13 (23%)
Aromatic	28/34 (82%)	15/18 (83%)	13/14 (93%)	0/2 (0%)
Overall	771/916 (84%)	439/451 (97%)	252/373 (68%)	80/92 (87%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 84%, i.e. 781 atoms were assigned a chemical shift out of a possible 930. 0 out of 18 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	313/401 (78%)	159/160 (99%)	76/162 (47%)	78/79 (99%)
Sidechain	440/495 (89%)	270/279 (97%)	167/203 (82%)	3/13 (23%)
Aromatic	28/34 (82%)	15/18 (83%)	13/14 (93%)	0/2 (0%)
Overall	781/930 (84%)	444/457 (97%)	256/379 (68%)	81/94 (86%)

#### 7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

#### 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.

Random coil index (RCI) for chain A:

