Full wwPDB NMR Structure Validation Report (i)

May 28, 2020 – 10:51 pm BST

PDB ID	:	2L22
Title	:	Mupirocin didomain ACP
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Deposited on	:	2010-08-10

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

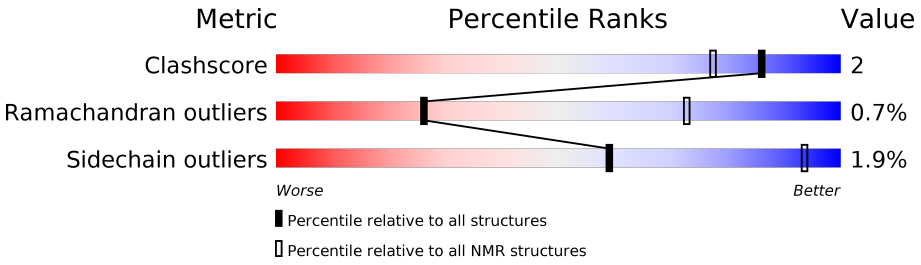
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance

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

The overall completeness of chemical shifts assignment was not calculated.

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



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

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

Mol	Chain	Length	Quality of chain
1	A	212	72% • 13% 14%

2 Ensemble composition and analysis

This entry contains 20 models. Model 5 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:4-A:78 (75)	0.65	9
2	A:104-A:183 (80)	0.99	5

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 2 clusters. No single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2850 atoms, of which 1414 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Mupirocin didomain Acyl Carrier Protein.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	183	2850	910	1414	250	271	5	0

There are 29 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	EXPRESSION TAG	UNP Q8RL76
A	-19	GLY	-	EXPRESSION TAG	UNP Q8RL76
A	-18	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-17	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-16	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-15	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-14	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-13	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-12	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-11	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	-10	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-9	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-8	GLY	-	EXPRESSION TAG	UNP Q8RL76
A	-7	LEU	-	EXPRESSION TAG	UNP Q8RL76
A	-6	VAL	-	EXPRESSION TAG	UNP Q8RL76
A	-5	PRO	-	EXPRESSION TAG	UNP Q8RL76
A	-4	ARG	-	EXPRESSION TAG	UNP Q8RL76
A	-3	GLY	-	EXPRESSION TAG	UNP Q8RL76
A	-2	SER	-	EXPRESSION TAG	UNP Q8RL76
A	-1	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	0	MET	-	EXPRESSION TAG	UNP Q8RL76
A	184	LEU	-	EXPRESSION TAG	UNP Q8RL76
A	185	GLU	-	EXPRESSION TAG	UNP Q8RL76
A	186	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	187	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	188	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	189	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	190	HIS	-	EXPRESSION TAG	UNP Q8RL76
A	191	HIS	-	EXPRESSION TAG	UNP Q8RL76

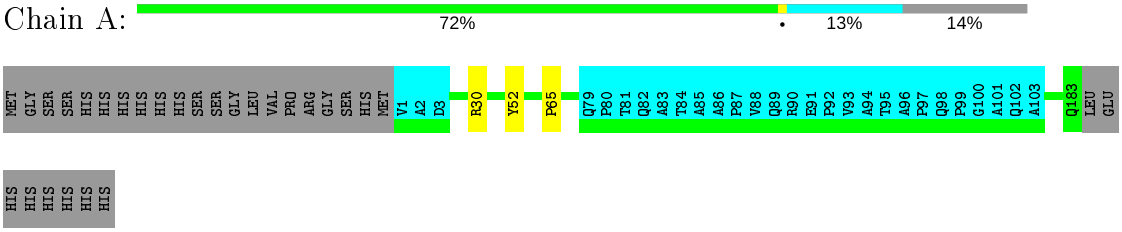
4 Residue-property plots

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: Mupirocin didomain Acyl Carrier Protein

Chain A:



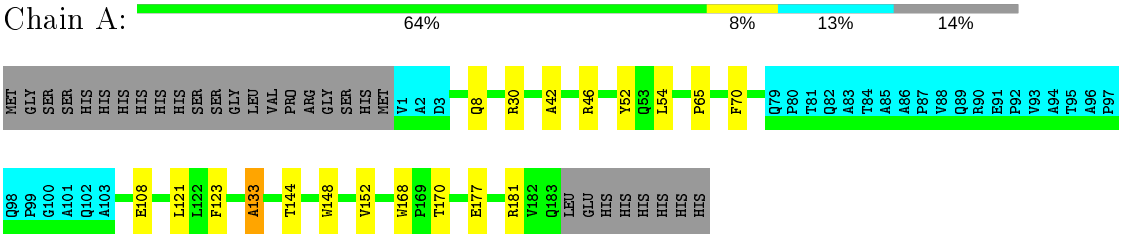
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: Mupirocin didomain Acyl Carrier Protein

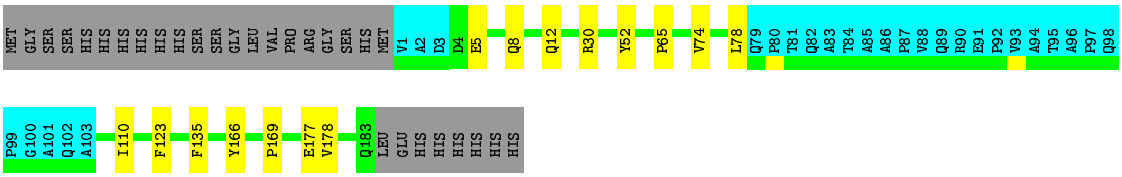
Chain A:



4.2.2 Score per residue for model 2

- Molecule 1: Mupirocin didomain Acyl Carrier Protein

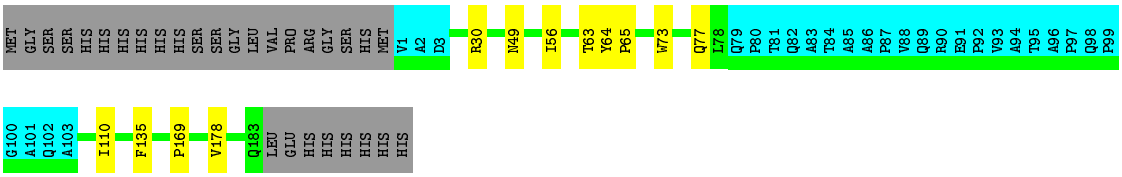
Chain A:



4.2.3 Score per residue for model 3

- Molecule 1: Mupirocin didomain Acyl Carrier Protein

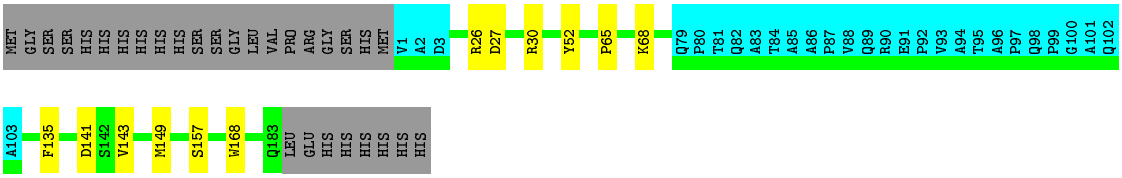
Chain A:



4.2.4 Score per residue for model 4

- Molecule 1: Mupirocin didomain Acyl Carrier Protein

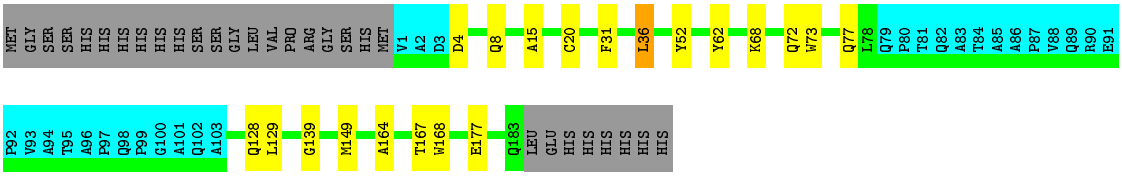
Chain A:



4.2.5 Score per residue for model 5 (medoid)

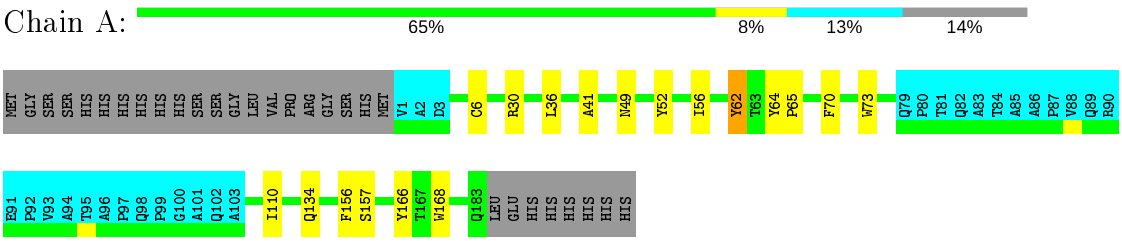
- Molecule 1: Mupirocin didomain Acyl Carrier Protein

Chain A:



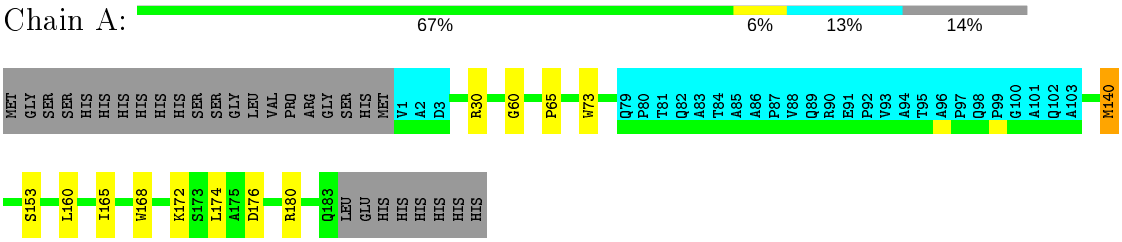
4.2.6 Score per residue for model 6

- Molecule 1: Mupirocin didomain Acyl Carrier Protein



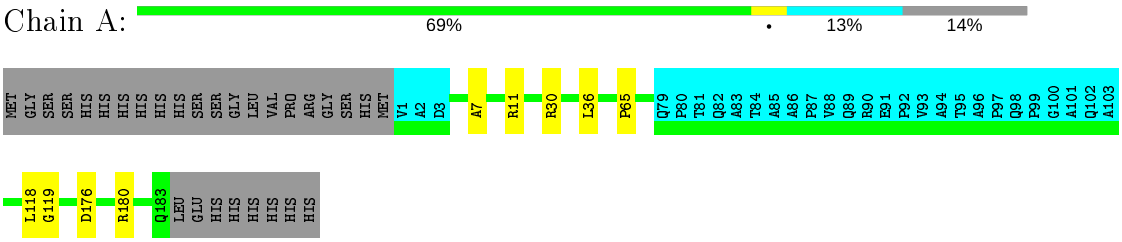
4.2.7 Score per residue for model 7

- Molecule 1: Mupirocin didomain Acyl Carrier Protein



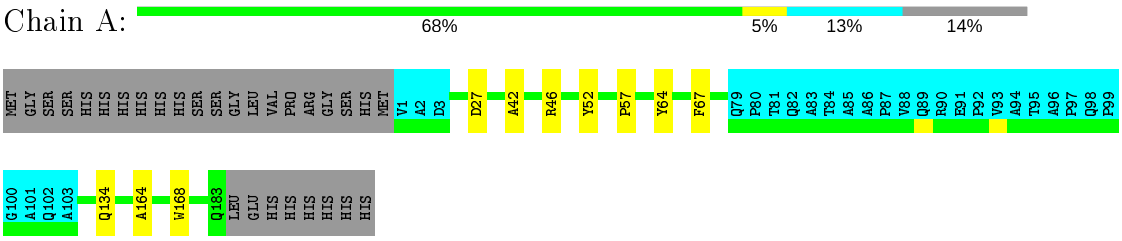
4.2.8 Score per residue for model 8

- Molecule 1: Mupirocin didomain Acyl Carrier Protein



4.2.9 Score per residue for model 9

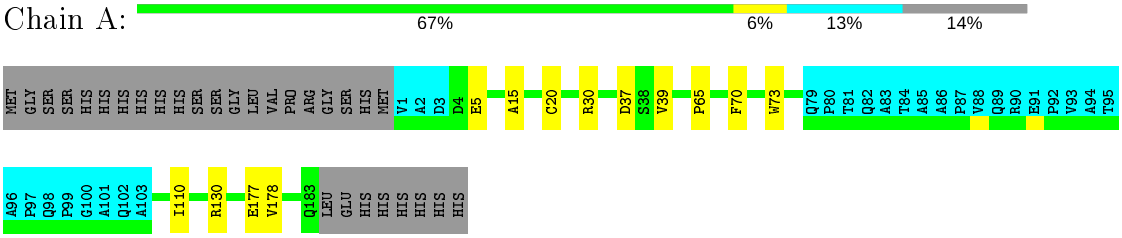
- Molecule 1: Mupirocin didomain Acyl Carrier Protein



4.2.10 Score per residue for model 10

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

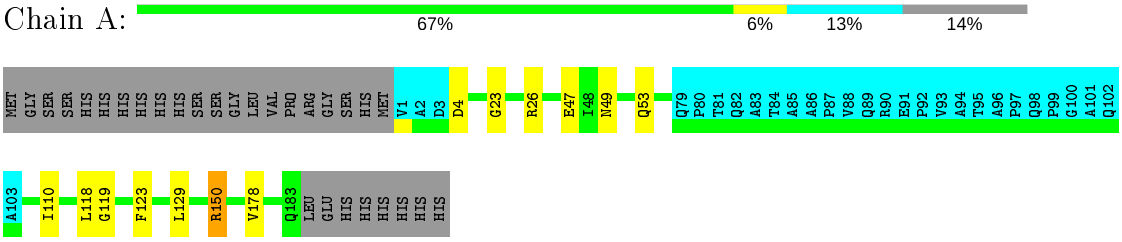
Chain A:



4.2.11 Score per residue for model 11

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

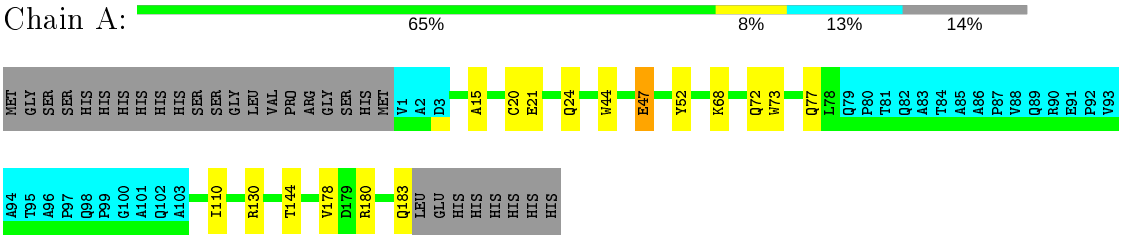
Chain A:



4.2.12 Score per residue for model 12

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

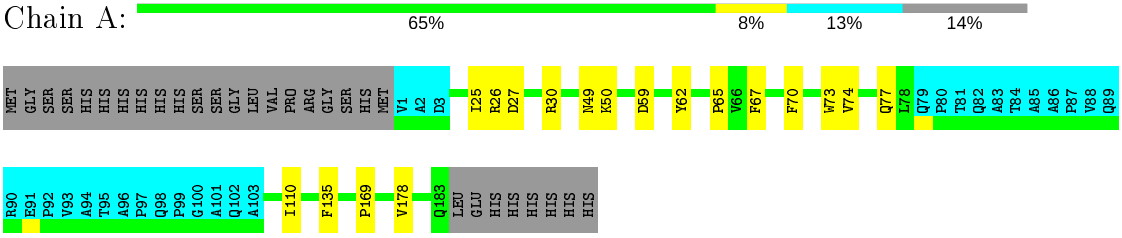
Chain A:



4.2.13 Score per residue for model 13

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

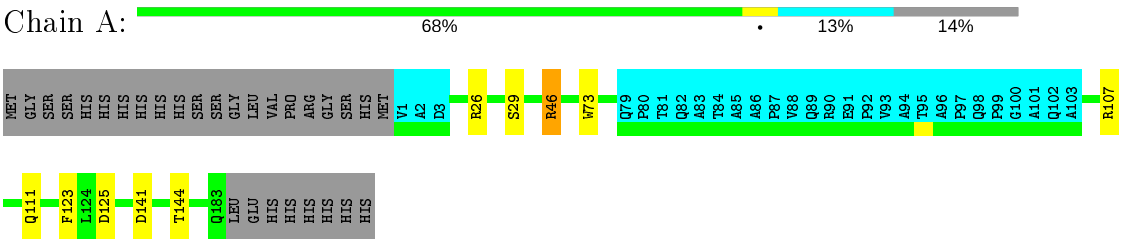
Chain A:



4.2.14 Score per residue for model 14

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

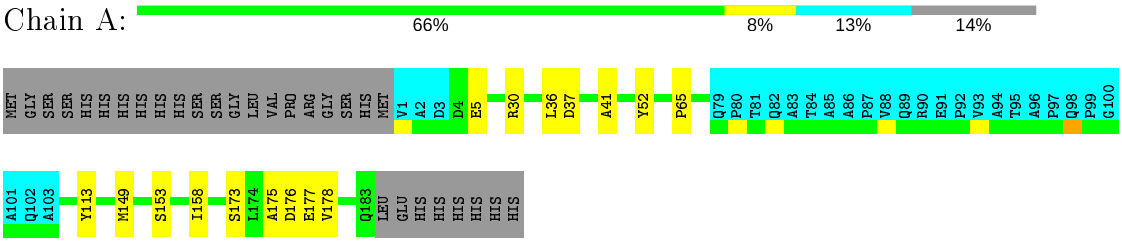
Chain A:



4.2.15 Score per residue for model 15

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

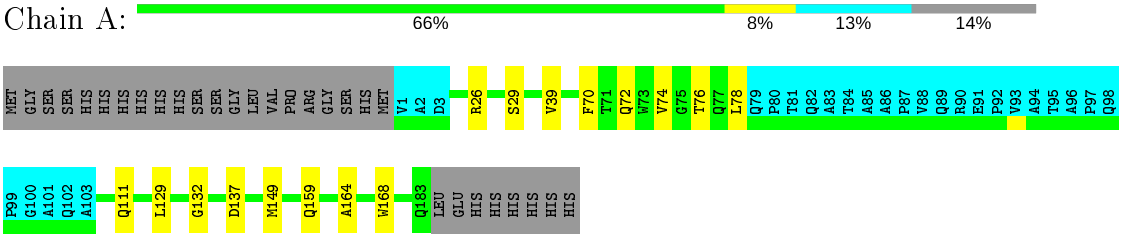
Chain A:



4.2.16 Score per residue for model 16

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

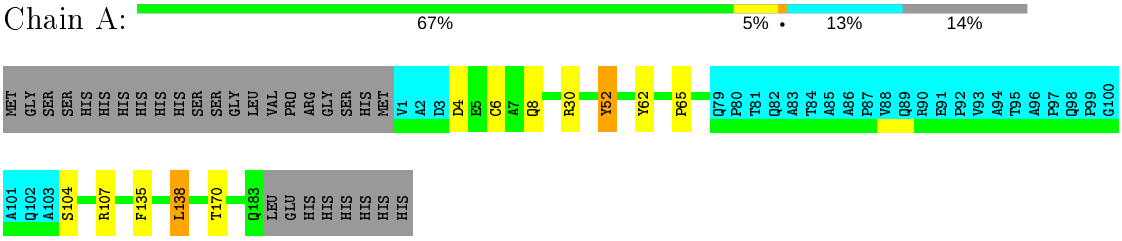
Chain A:



4.2.17 Score per residue for model 17

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

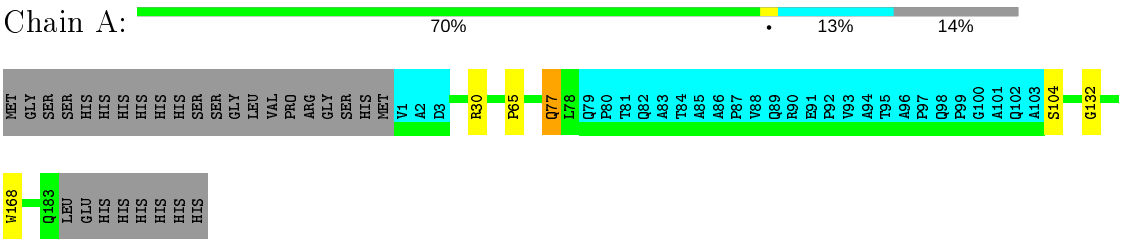
Chain A:



4.2.18 Score per residue for model 18

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

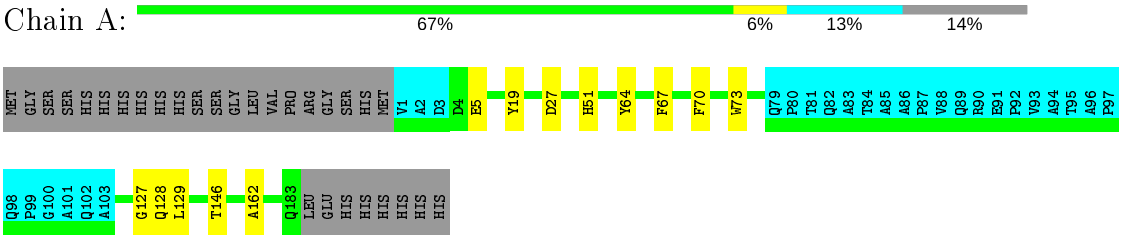
Chain A:



4.2.19 Score per residue for model 19

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

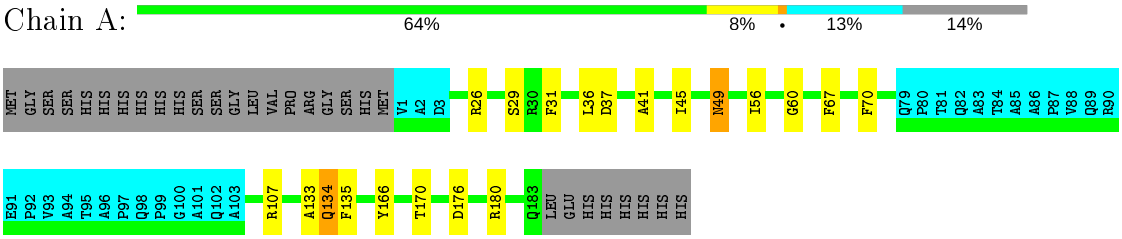
Chain A:



4.2.20 Score per residue for model 20

• Molecule 1: Mupirocin didomain Acyl Carrier Protein

Chain A:



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	2.2

No chemical shift data was provided. No validations of the models with respect to experimental NMR restraints is performed at this time.

COVALENT-GEOMETRY INFOmissingINFO

5.1 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	1242	1225	1223	5±2
All	All	24840	24500	24460	103

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 2.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:164:ALA:HA	1:A:168:TRP:CZ3	0.62	2.29	16	1
1:A:110:ILE:HG21	1:A:178:VAL:HG11	0.60	1.73	3	6
1:A:60:GLY:HA2	1:A:73:TRP:CZ3	0.60	2.32	7	1
1:A:30:ARG:HA	1:A:65:PRO:O	0.59	1.97	8	12
1:A:27:ASP:HA	1:A:67:PHE:CZ	0.56	2.36	13	2
1:A:150:ARG:HA	1:A:150:ARG:NE	0.56	2.16	11	1
1:A:15:ALA:HA	1:A:20:CYS:SG	0.55	2.41	10	3
1:A:73:TRP:O	1:A:77:GLN:HG2	0.53	2.04	12	2
1:A:7:ALA:O	1:A:11:ARG:HG3	0.53	2.03	8	1
1:A:49:ASN:OD1	1:A:56:ILE:HB	0.51	2.04	6	1
1:A:164:ALA:HA	1:A:168:TRP:CE3	0.49	2.42	16	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:134:GLN:NE2	1:A:135:PHE:H	0.48	2.06	20	1
1:A:36:LEU:HD21	1:A:41:ALA:HB2	0.48	1.85	15	3
1:A:74:VAL:O	1:A:78:LEU:HG	0.48	2.08	16	2
1:A:168:TRP:CZ2	1:A:177:GLU:HG2	0.48	2.44	1	2
1:A:173:SER:HA	1:A:176:ASP:OD2	0.47	2.09	15	1
1:A:176:ASP:O	1:A:180:ARG:HG2	0.47	2.08	7	2
1:A:49:ASN:O	1:A:53:GLN:HA	0.47	2.09	11	1
1:A:148:TRP:O	1:A:152:VAL:HG23	0.47	2.10	1	1
1:A:135:PHE:CD1	1:A:169:PRO:HA	0.47	2.45	13	2
1:A:175:ALA:O	1:A:178:VAL:HG12	0.46	2.10	15	1
1:A:8:GLN:O	1:A:12:GLN:HG3	0.46	2.11	2	1
1:A:141:ASP:OD2	1:A:143:VAL:HG12	0.46	2.11	4	1
1:A:133:ALA:O	1:A:170:THR:HA	0.46	2.11	20	2
1:A:46:ARG:HA	1:A:46:ARG:NE	0.46	2.25	14	1
1:A:4:ASP:O	1:A:8:GLN:HG3	0.45	2.11	5	2
1:A:135:PHE:HB3	1:A:168:TRP:O	0.45	2.12	4	1
1:A:6:CYS:HA	1:A:52:TYR:OH	0.45	2.11	17	2
1:A:176:ASP:O	1:A:180:ARG:HG3	0.45	2.11	8	1
1:A:70:PHE:O	1:A:73:TRP:HB3	0.45	2.12	19	4
1:A:118:LEU:HD12	1:A:119:GLY:N	0.44	2.27	8	2
1:A:110:ILE:HG12	1:A:156:PHE:CZ	0.44	2.47	6	1
1:A:181:ARG:HA	1:A:181:ARG:NE	0.44	2.27	1	1
1:A:42:ALA:O	1:A:46:ARG:HG2	0.44	2.12	1	1
1:A:164:ALA:HA	1:A:167:THR:OG1	0.44	2.12	5	1
1:A:140:MET:SD	1:A:165:ILE:HD12	0.44	2.52	7	1
1:A:168:TRP:CZ3	1:A:174:LEU:HD13	0.44	2.48	7	1
1:A:74:VAL:O	1:A:77:GLN:HG2	0.44	2.12	13	1
1:A:49:ASN:ND2	1:A:56:ILE:HG22	0.44	2.28	20	1
1:A:153:SER:OG	1:A:160:LEU:HB2	0.44	2.12	7	1
1:A:26:ARG:HD2	1:A:26:ARG:N	0.44	2.28	13	1
1:A:68:LYS:O	1:A:72:GLN:HG2	0.43	2.13	12	1
1:A:49:ASN:ND2	1:A:56:ILE:HB	0.43	2.29	3	1
1:A:135:PHE:HA	1:A:169:PRO:O	0.43	2.14	2	1
1:A:73:TRP:CZ3	1:A:77:GLN:HG3	0.43	2.48	3	1
1:A:164:ALA:HB1	1:A:168:TRP:CD1	0.43	2.48	9	1
1:A:44:TRP:O	1:A:47:GLU:HG3	0.43	2.13	12	1
1:A:121:LEU:HD21	1:A:144:THR:HB	0.42	1.90	1	1
1:A:107:ARG:O	1:A:111:GLN:HG2	0.42	2.14	14	1
1:A:26:ARG:HB3	1:A:29:SER:OG	0.42	2.14	20	1
1:A:42:ALA:O	1:A:46:ARG:HG3	0.42	2.14	9	1
1:A:27:ASP:HA	1:A:67:PHE:CE2	0.41	2.49	9	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:72:GLN:O	1:A:76:THR:HG22	0.41	2.16	16	1
1:A:26:ARG:HG2	1:A:29:SER:OG	0.41	2.15	16	1
1:A:104:SER:O	1:A:107:ARG:HB3	0.41	2.16	17	1
1:A:60:GLY:HA3	1:A:70:PHE:CE1	0.41	2.51	20	1
1:A:63:THR:HG23	1:A:64:TYR:CD2	0.41	2.50	3	1
1:A:153:SER:HB2	1:A:158:ILE:O	0.41	2.16	15	1
1:A:31:PHE:CZ	1:A:67:PHE:HA	0.41	2.51	20	1
1:A:27:ASP:O	1:A:68:LYS:HB2	0.41	2.16	4	1
1:A:68:LYS:O	1:A:72:GLN:HG3	0.41	2.16	5	1
1:A:77:GLN:HE21	1:A:77:GLN:HA	0.41	1.76	18	1
1:A:21:GLU:O	1:A:24:GLN:HG2	0.40	2.16	12	1
1:A:41:ALA:O	1:A:45:ILE:HG12	0.40	2.16	20	1
1:A:146:THR:HB	1:A:162:ALA:HB1	0.40	1.92	19	1
1:A:37:ASP:OD2	1:A:39:VAL:HG22	0.40	2.17	10	1
1:A:26:ARG:HB2	1:A:29:SER:OG	0.40	2.16	14	1
1:A:135:PHE:O	1:A:138:LEU:HD22	0.40	2.17	17	1
1:A:36:LEU:N	1:A:36:LEU:HD23	0.40	2.32	5	1
1:A:59:ASP:HA	1:A:62:TYR:CD2	0.40	2.52	13	1

5.2 Torsion angles (i)

5.2.1 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	154/212 (73%)	146±2 (95±1%)	7±2 (5±1%)	1±1 (1±1%)	26	73
All	All	3080/4240 (73%)	2914 (95%)	144 (5%)	22 (1%)	26	73

All 14 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	123	PHE	4
1	A	129	LEU	3
1	A	130	ARG	2
1	A	157	SER	2

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Mol	Chain	Res	Type	Models (Total)
1	A	132	GLY	2
1	A	128	GLN	1
1	A	139	GLY	1
1	A	127	GLY	1
1	A	25	ILE	1
1	A	104	SER	1
1	A	23	GLY	1
1	A	26	ARG	1
1	A	133	ALA	1
1	A	140	MET	1

5.2.2 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	131/176 (74%)	129±1 (98±1%)	2±1 (2±1%)	59 93
All	All	2620/3520 (74%)	2571 (98%)	49 (2%)	59 93

All 34 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	149	MET	4
1	A	5	GLU	3
1	A	134	GLN	3
1	A	177	GLU	3
1	A	47	GLU	2
1	A	168	TRP	2
1	A	49	ASN	2
1	A	36	LEU	2
1	A	37	ASP	2
1	A	144	THR	2
1	A	39	VAL	1
1	A	159	GLN	1
1	A	26	ARG	1
1	A	73	TRP	1
1	A	54	LEU	1

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Mol	Chain	Res	Type	Models (Total)
1	A	150	ARG	1
1	A	138	LEU	1
1	A	128	GLN	1
1	A	51	HIS	1
1	A	137	ASP	1
1	A	183	GLN	1
1	A	108	GLU	1
1	A	50	LYS	1
1	A	125	ASP	1
1	A	141	ASP	1
1	A	107	ARG	1
1	A	170	THR	1
1	A	46	ARG	1
1	A	129	LEU	1
1	A	172	LYS	1
1	A	111	GLN	1
1	A	8	GLN	1
1	A	77	GLN	1
1	A	57	PRO	1

5.2.3 RNA

There are no RNA molecules in this entry.

5.3 Non-standard residues in protein, DNA, RNA chains

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

5.4 Carbohydrates

There are no carbohydrates in this entry.

5.5 Ligand geometry

There are no ligands in this entry.

5.6 Other polymers

There are no such molecules in this entry.

5.7 Polymer linkage issues

There are no chain breaks in this entry.

6 Chemical shift validation

No chemical shift data were provided