wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	2KI4
Title	:	FGF1-S100A13 complex structure: key component in non-classical path way of FGF1
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Deposited on	:	2009-04-27

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

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

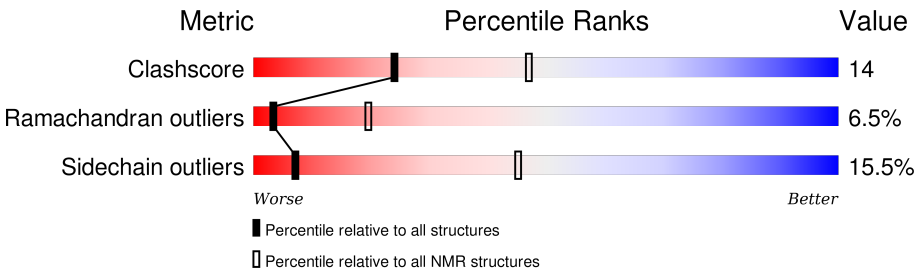
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	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 (i)

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

The overall completeness of chemical shifts assignment was not calculated.

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



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	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 >=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	133	51% 39% 9% •
1	D	133	53% 37% 9% •
2	B	98	68% 20% • 9%
3	C	98	61% 26% • 12%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 7 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:1-A:133, B:8-B:96, C:12-C:93, C:95-C:98, D:1-D:133 (441)	0.16	7

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 and 1 single-model cluster was found.

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

3 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7505 atoms, of which 3764 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Heparin-binding growth factor 1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	133	2113	671	1049	185	204	4	0
1	D	133	2113	671	1049	185	204	4	0

- Molecule 2 is a protein called Protein S100-A13.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
2	B	98	1640	512	834	136	156	2	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	97	DLY	LYS	SEE REMARK 999	UNP Q99584
B	98	DLY	LYS	SEE REMARK 999	UNP Q99584

- Molecule 3 is a protein called Protein S100-A13.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
3	C	98	1639	512	832	136	157	2	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	94	DLY	LYS	SEE REMARK 999	UNP Q99584

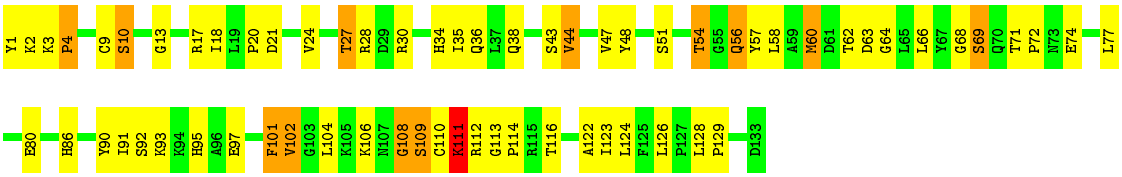
4 Residue-property plots (i)

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.

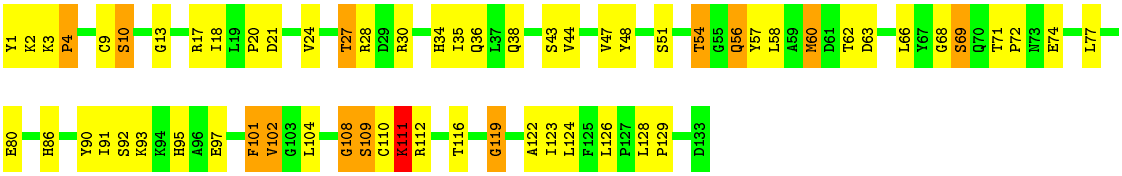
- Molecule 1: Heparin-binding growth factor 1

Chain A:



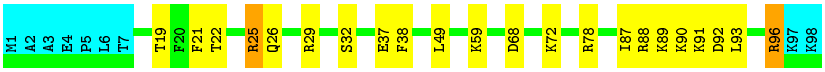
- Molecule 1: Heparin-binding growth factor 1

Chain D:



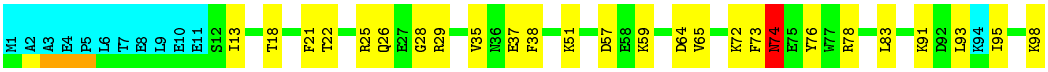
- Molecule 2: Protein S100-A13

Chain B:



- Molecule 3: Protein S100-A13

Chain C:

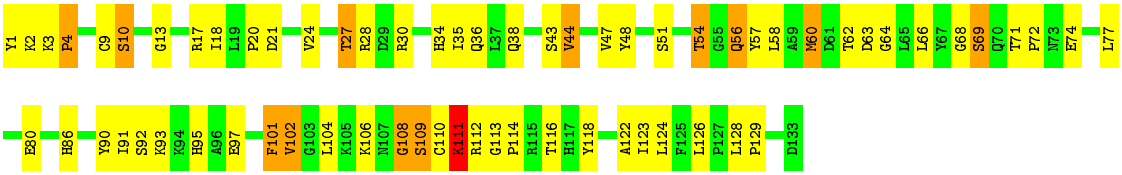


4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 7. Colouring as in section 4.1 above.

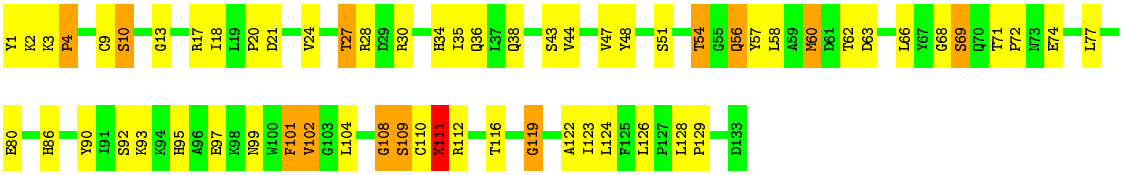
- Molecule 1: Heparin-binding growth factor 1

Chain A:



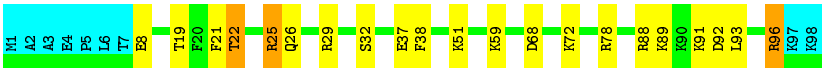
- Molecule 1: Heparin-binding growth factor 1

Chain D:



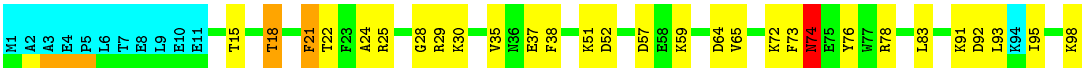
- Molecule 2: Protein S100-A13

Chain B:



- Molecule 3: Protein S100-A13

Chain C:



5 Refinement protocol and experimental data overview (i)

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

Of the 5000 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	structure solution	1.2 & 2.2
CNS	structure solution	1.2
CNS	refinement	1.2

No chemical shift data was provided. 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: DLY

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	0.30±0.00	0±0/1088 (0.0±0.0%)	0.39±0.00	0±0/1466 (0.0±0.0%)
1	D	0.30±0.00	0±0/1088 (0.0±0.0%)	0.39±0.00	0±0/1466 (0.0±0.0%)
2	B	0.50±0.00	0±0/748 (0.0±0.0%)	0.93±0.00	2±0/1000 (0.2±0.0%)
3	C	0.51±0.00	0±0/722 (0.0±0.0%)	1.00±0.01	0±0/961 (0.0±0.0%)
All	All	0.40	0/72920 (0.0%)	0.68	40/97860 (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
2	B	0.0±0.0	4.0±0.0
3	C	0.0±0.0	7.0±0.0
All	All	0	220

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(^{o})	Ideal(^{o})	Models	
								Worst	Total
2	B	37	GLU	OE1-CD-OE2	-5.93	116.18	123.30	17	20
2	B	68	ASP	CB-CG-OD2	5.37	123.13	118.30	14	20

There are no chirality outliers.

5 of 11 unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
3	C	91	LYS	Peptide	20
3	C	74	ASN	Peptide	20
3	C	64	ASP	Peptide	20
2	B	89	LYS	Peptide	20
3	C	35	VAL	Peptide,Mainchain	20

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	1064	1049	1046	49±3
1	D	1064	1049	1046	47±2
2	B	739	756	755	4±1
3	C	714	740	739	4±2
All	All	71620	71880	71720	2036

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.

5 of 150 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:66:LEU:HD12	1:A:110:CYS:HB2	0.74	1.59	16	20
1:D:66:LEU:HD12	1:D:110:CYS:HB2	0.69	1.63	15	20
3:C:22:THR:HG21	1:D:119:GLY:HA3	0.68	1.66	20	12
1:A:3:LYS:CB	1:A:4:PRO:HD2	0.67	2.20	9	20
3:C:21:PHE:O	3:C:25:ARG:HG2	0.67	1.89	6	10

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	131/133 (98%)	95±1 (73±0%)	26±1 (20±1%)	10±1 (7±0%)	3	17
1	D	131/133 (98%)	96±0 (73±0%)	25±1 (19±0%)	10±0 (8±0%)	2	16
2	B	89/98 (91%)	68±1 (76±1%)	17±1 (19±1%)	4±0 (5±0%)	5	28
3	C	85/98 (87%)	65±1 (76±1%)	16±1 (18±1%)	5±1 (5±1%)	4	24
All	All	8720/9240 (94%)	6482 (74%)	1673 (19%)	565 (6%)	3	19

5 of 32 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
2	B	96	ARG	20
1	D	10	SER	20
1	A	4	PRO	20
1	D	111	LYS	20
1	A	102	VAL	20

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	117/117 (100%)	98±0 (83±0%)	19±0 (17±0%)	6	43
1	D	117/117 (100%)	97±1 (83±1%)	20±1 (17±1%)	6	42
2	B	84/89 (94%)	75±1 (90±1%)	9±1 (10±1%)	13	57
3	C	81/90 (90%)	67±1 (83±2%)	14±1 (17±2%)	6	42
All	All	7980/8260 (97%)	6747 (85%)	1233 (15%)	7	45

5 of 74 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	D	62	THR	20
1	A	35	ILE	20
1	A	101	PHE	20
1	A	18	ILE	20
1	D	111	LYS	20

6.3.3 RNA (i)

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues are 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	DLY	B	97	2	6,8,9	0.68±0.01	0±0 (0±0%)
2	DLY	B	98	2	6,8,9	0.65±0.01	0±0 (0±0%)
3	DLY	C	94	3	6,8,9	0.57±0.01	0±0 (0±0%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
2	DLY	B	97	2	6,8,10	1.63±0.01	0±0 (0±0%)
2	DLY	B	98	2	6,8,10	1.26±0.01	0±0 (0±0%)
3	DLY	C	94	3	6,8,10	1.07±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	DLY	B	97	2	-	0±0,5,7,9	0±0,0,0,0
2	DLY	B	98	2	-	0±0,5,7,9	0±0,0,0,0
3	DLY	C	94	3	-	0±0,5,7,9	0±0,0,0,0

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates (i)

There are no carbohydrates in this entry.

6.6 Ligand geometry (i)

There are no ligands in this entry.

6.7 Other polymers (i)

There are no such molecules in this entry.

6.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

7 Chemical shift validation (i)

No chemical shift data were provided