



Full wwPDB NMR Structure Validation Report ⓘ

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PDB ID : 2JYI
Title : Solution structure of MLL CXXC domain
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Deposited on : 2007-12-13

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
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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 : unknown
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

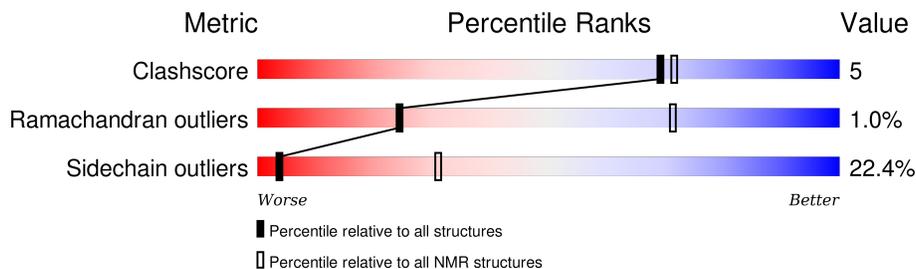
1 Overall quality at a glance i

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	57	

2 Ensemble composition and analysis

This entry contains 14 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:1154-A:1181, A:1189-A:1196 (36)	0.14	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 3 clusters and 2 single-model clusters were found.

Cluster number	Models
1	1, 2, 6, 8, 11
2	5, 7, 9, 10, 12
3	4, 14
Single-model clusters	3; 13

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 905 atoms, of which 457 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Zinc finger protein HRX.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
1	A	57	903	267	457	93	75	11	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

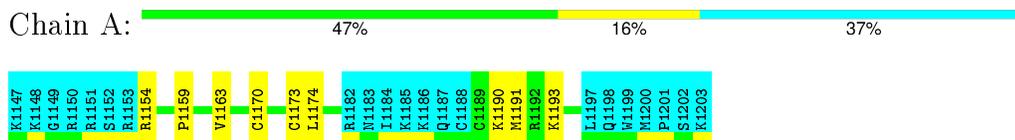
Mol	Chain	Residues	Atoms	
2	A	2	Total	Zn
			2	2

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: Zinc finger protein HRX



4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

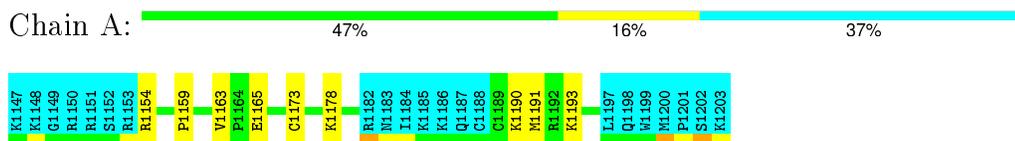
4.2.1 Score per residue for model 1

- Molecule 1: Zinc finger protein HRX



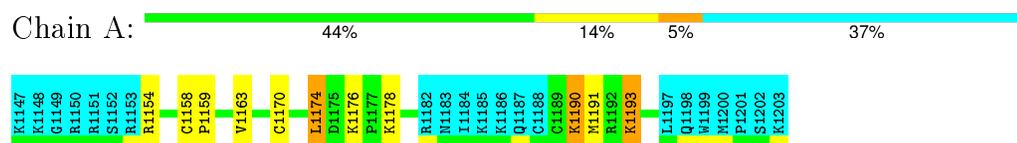
4.2.2 Score per residue for model 2

- Molecule 1: Zinc finger protein HRX



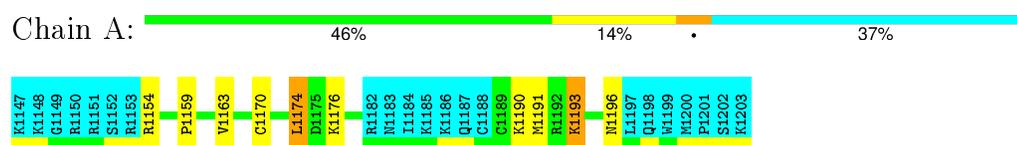
4.2.3 Score per residue for model 3

- Molecule 1: Zinc finger protein HRX



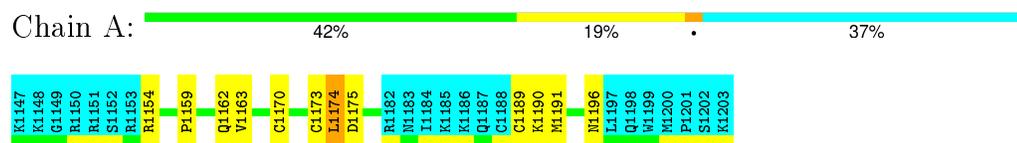
4.2.4 Score per residue for model 4

- Molecule 1: Zinc finger protein HRX



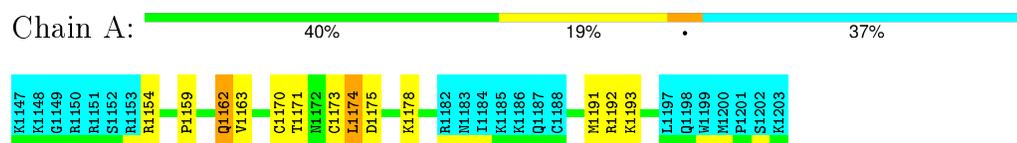
4.2.5 Score per residue for model 5

- Molecule 1: Zinc finger protein HRX



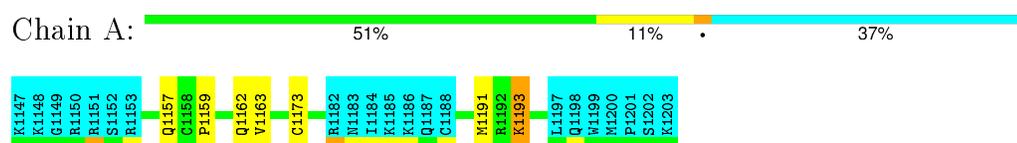
4.2.6 Score per residue for model 6

- Molecule 1: Zinc finger protein HRX



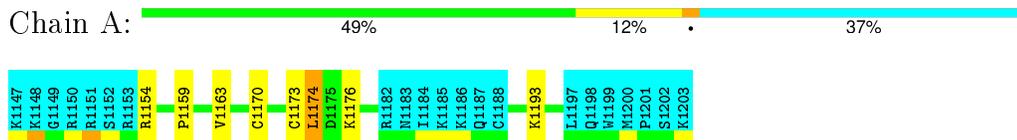
4.2.7 Score per residue for model 7

- Molecule 1: Zinc finger protein HRX



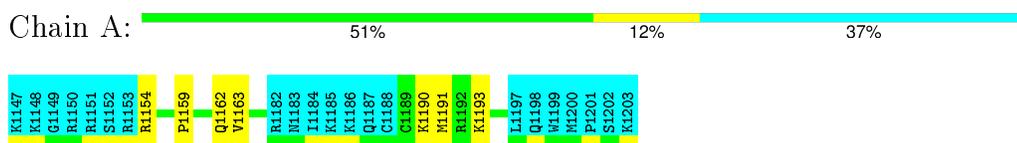
4.2.8 Score per residue for model 8 (medoid)

- Molecule 1: Zinc finger protein HRX



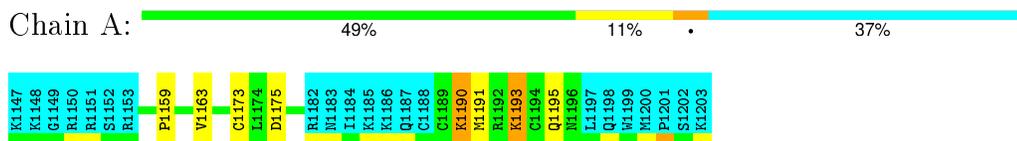
4.2.9 Score per residue for model 9

- Molecule 1: Zinc finger protein HRX



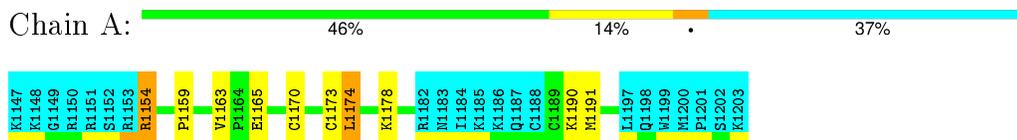
4.2.10 Score per residue for model 10

- Molecule 1: Zinc finger protein HRX



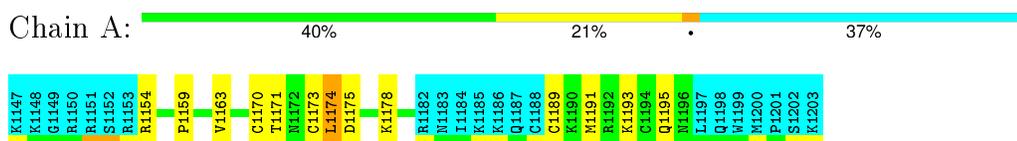
4.2.11 Score per residue for model 11

- Molecule 1: Zinc finger protein HRX



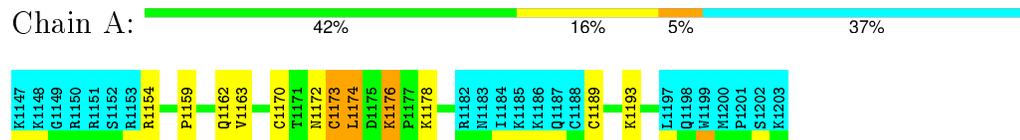
4.2.12 Score per residue for model 12

- Molecule 1: Zinc finger protein HRX



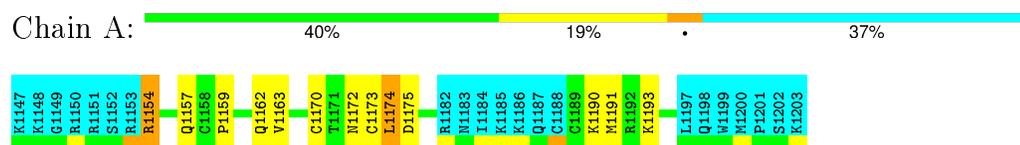
4.2.13 Score per residue for model 13

- Molecule 1: Zinc finger protein HRX



4.2.14 Score per residue for model 14

- Molecule 1: Zinc finger protein HRX



5 Refinement protocol and experimental data overview

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

Of the 200 calculated structures, 14 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
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:
ZN

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

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	263	250	250	2±1
All	All	3710	3500	3500	33

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 5.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:1159:PRO:O	1:A:1163:VAL:HG23	0.79	1.77	13	14
1:A:1170:CYS:O	1:A:1174:LEU:HD12	0.63	1.94	5	1
1:A:1170:CYS:O	1:A:1174:LEU:CD2	0.45	2.64	4	9
1:A:1190:LYS:O	1:A:1193:LYS:HG3	0.44	2.12	10	2
1:A:1193:LYS:CD	1:A:1193:LYS:O	0.43	2.66	4	3
1:A:1173:CYS:O	1:A:1176:LYS:CG	0.41	2.68	13	1
1:A:1159:PRO:HA	1:A:1162:GLN:CG	0.41	2.46	6	2
1:A:1159:PRO:HA	1:A:1162:GLN:HG2	0.40	1.93	6	1

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	36/57 (63%)	34±1 (96±2%)	1±0 (3±1%)	0±0 (1±1%)	24	71
All	All	504/798 (63%)	482 (96%)	17 (3%)	5 (1%)	24	71

All 2 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	1189	CYS	3
1	A	1154	ARG	2

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	31/51 (61%)	24±2 (78±5%)	7±2 (22±5%)	4	31
All	All	434/714 (61%)	337 (78%)	97 (22%)	4	31

All 18 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	1193	LYS	12
1	A	1173	CYS	11
1	A	1191	MET	11
1	A	1174	LEU	10
1	A	1154	ARG	9
1	A	1190	LYS	8
1	A	1178	LYS	6
1	A	1162	GLN	5

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Mol	Chain	Res	Type	Models (Total)
1	A	1176	LYS	5
1	A	1175	ASP	5
1	A	1165	GLU	3
1	A	1172	ASN	2
1	A	1171	THR	2
1	A	1195	GLN	2
1	A	1157	GLN	2
1	A	1196	ASN	2
1	A	1158	CYS	1
1	A	1192	ARG	1

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 carbohydrates in this entry.

6.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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