Full wwPDB NMR Structure Validation Report (i)

Apr 26, 2016 – 09:23 PM BST

PDB ID	:	2JQJ
Title	:	NMR structure of yeast Dun1 FHA domain
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Deposited on	:	2007-06-02

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 (i) symbol.

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

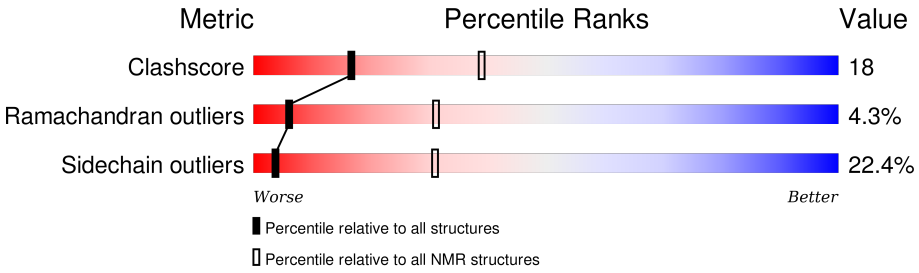
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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 <=5%

Mol	Chain	Length	Quality of chain
1	A	151	37% 31% • 23% 7%

2 Ensemble composition and analysis

This entry contains 20 models. Model 9 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:32-A:138 (107)	0.43	9

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 6 clusters and 6 single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2253 atoms, of which 1123 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called DNA damage response protein kinase DUN1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	141	2253	698	1123	205	223	4	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	MET	-	EXPRESSION TAG	UNP P39009
A	18	ALA	-	EXPRESSION TAG	UNP P39009
A	160	ARG	-	EXPRESSION TAG	UNP P39009
A	161	SER	-	EXPRESSION TAG	UNP P39009
A	162	HIS	-	EXPRESSION TAG	UNP P39009
A	163	HIS	-	EXPRESSION TAG	UNP P39009
A	164	HIS	-	EXPRESSION TAG	UNP P39009
A	165	HIS	-	EXPRESSION TAG	UNP P39009
A	166	HIS	-	EXPRESSION TAG	UNP P39009
A	167	HIS	-	EXPRESSION TAG	UNP P39009

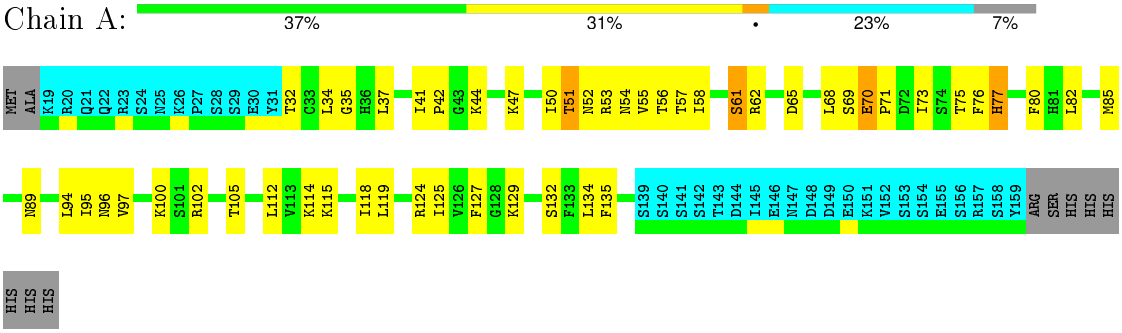
4 Residue-property plots

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: DNA damage response protein kinase DUN1

Chain A:



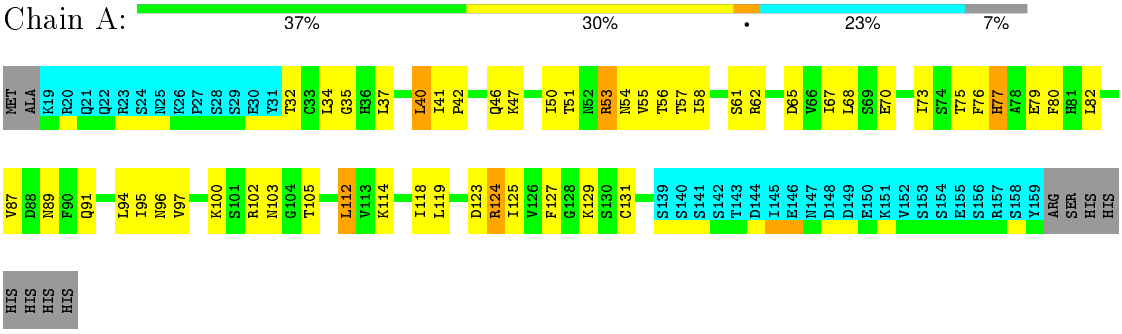
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: DNA damage response protein kinase DUN1

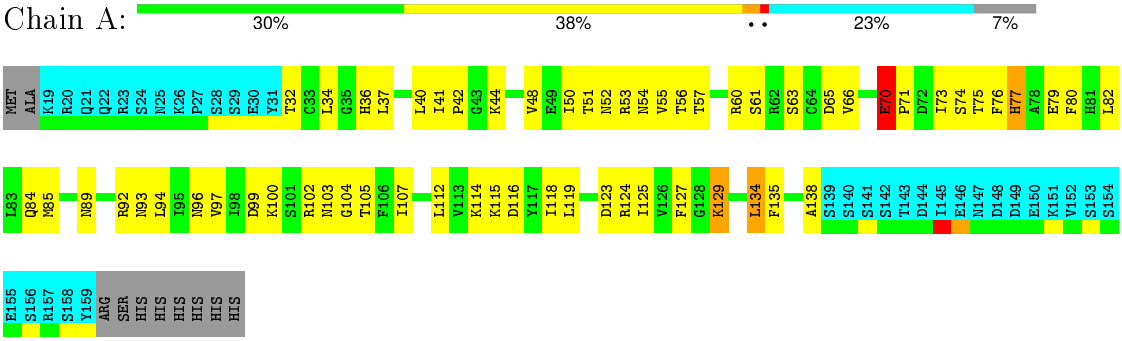
Chain A:



4.2.2 Score per residue for model 2

• Molecule 1: DNA damage response protein kinase DUN1

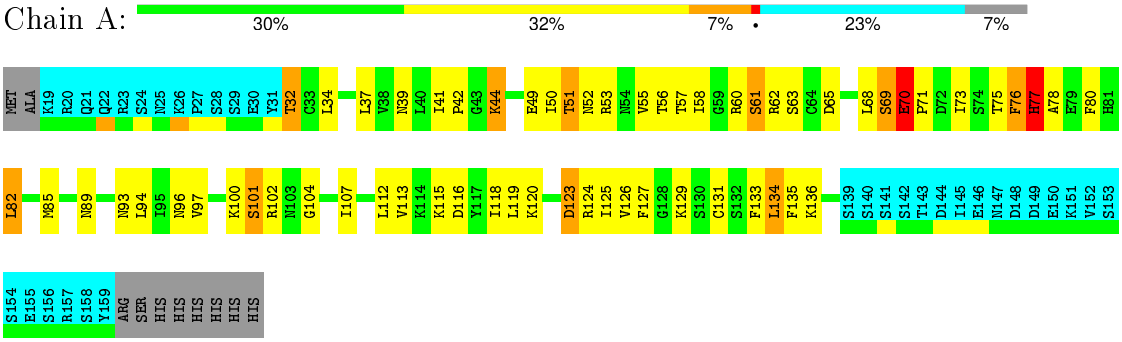
Chain A:



4.2.3 Score per residue for model 3

• Molecule 1: DNA damage response protein kinase DUN1

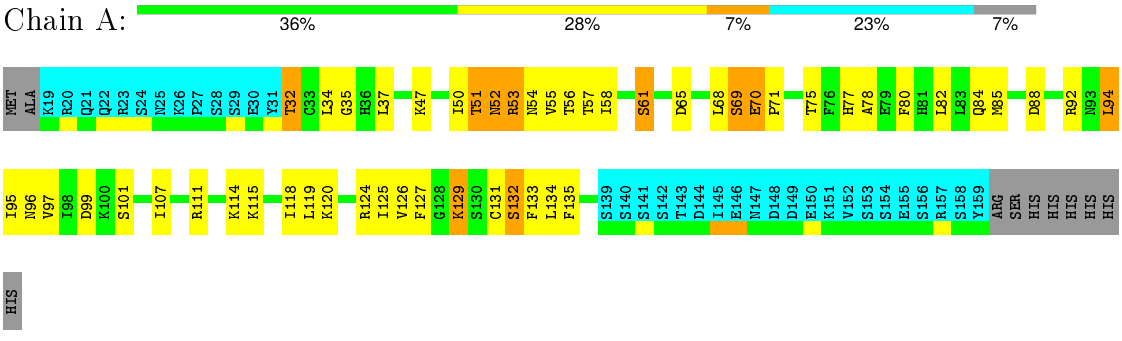
Chain A:



4.2.4 Score per residue for model 4

• Molecule 1: DNA damage response protein kinase DUN1

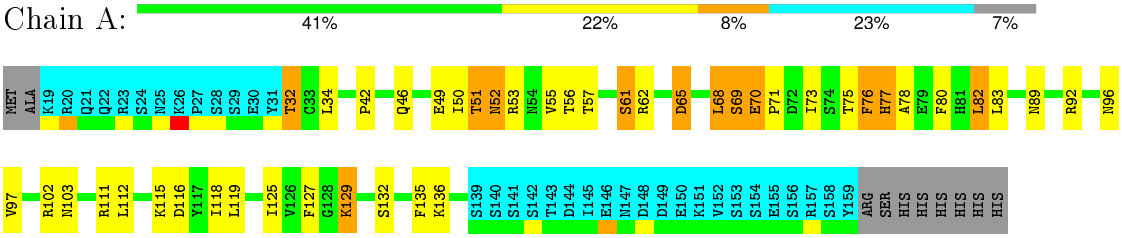
Chain A:



4.2.5 Score per residue for model 5

• Molecule 1: DNA damage response protein kinase DUN1

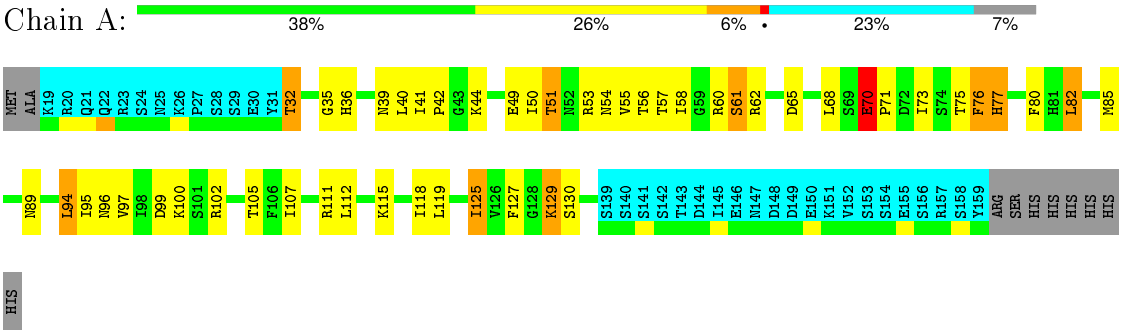
Chain A:



4.2.6 Score per residue for model 6

• Molecule 1: DNA damage response protein kinase DUN1

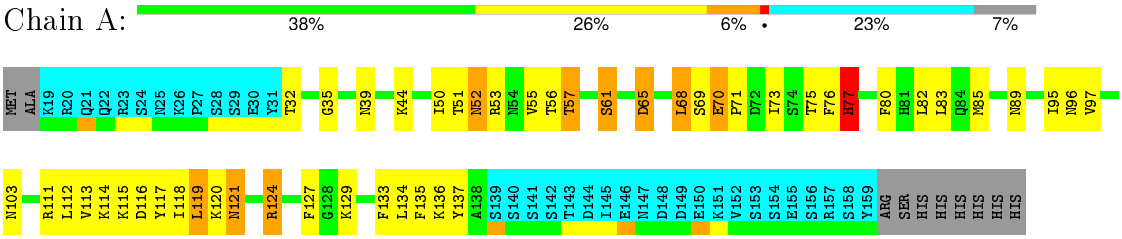
Chain A:



4.2.7 Score per residue for model 7

• Molecule 1: DNA damage response protein kinase DUN1

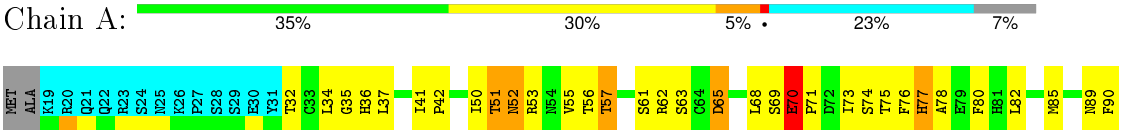
Chain A:

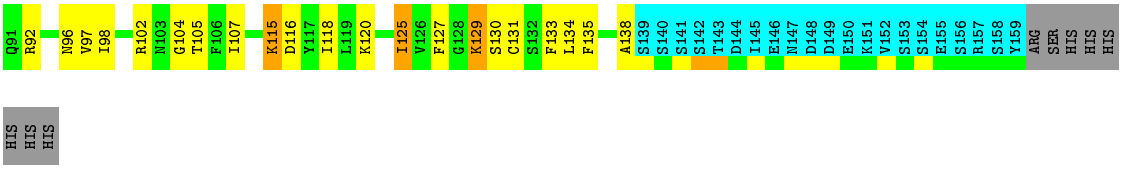


4.2.8 Score per residue for model 8

• Molecule 1: DNA damage response protein kinase DUN1

Chain A:

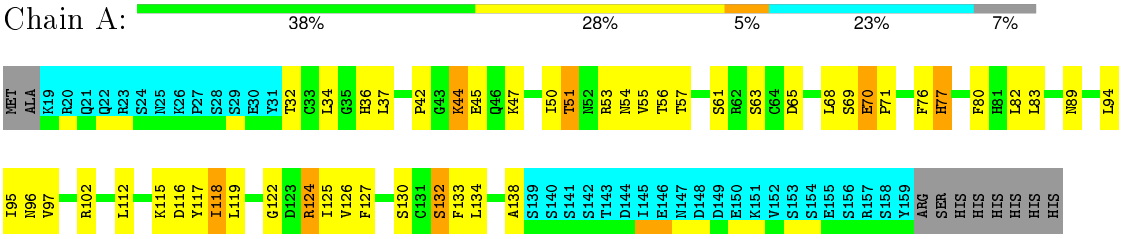




4.2.9 Score per residue for model 9 (medoid)

- Molecule 1: DNA damage response protein kinase DUN1

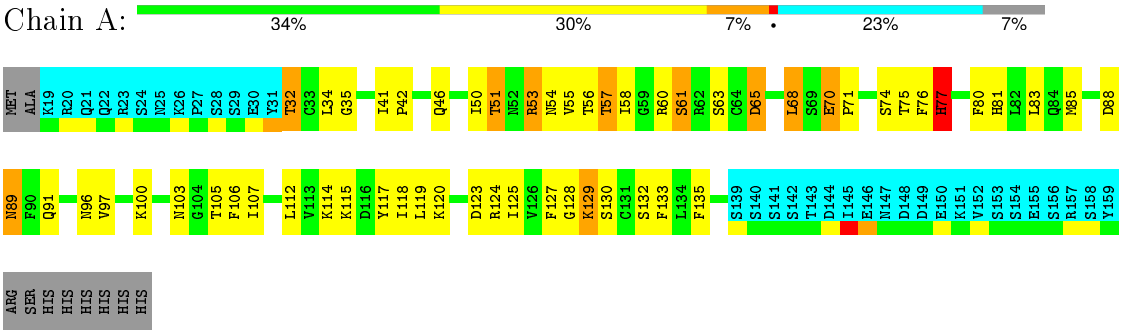
Chain A:



4.2.10 Score per residue for model 10

- Molecule 1: DNA damage response protein kinase DUN1

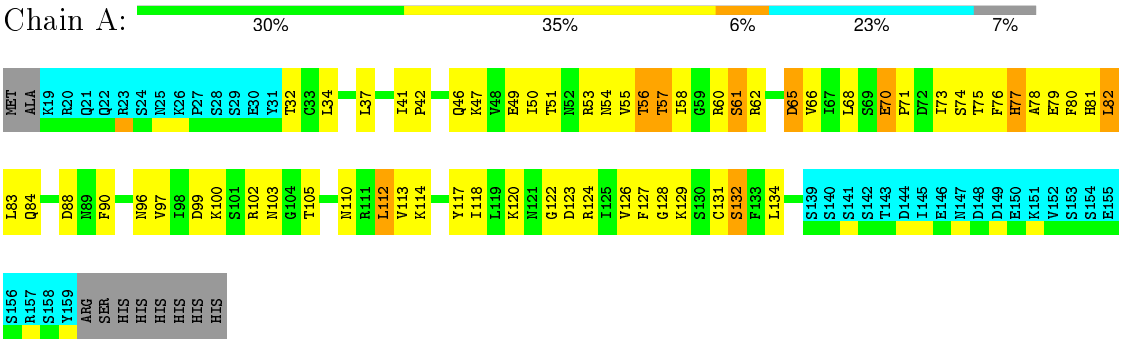
Chain A:



4.2.11 Score per residue for model 11

- Molecule 1: DNA damage response protein kinase DUN1

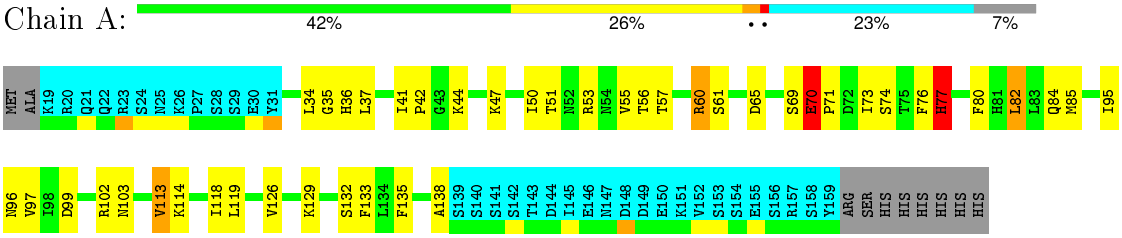
Chain A:



4.2.12 Score per residue for model 12

• Molecule 1: DNA damage response protein kinase DUN1

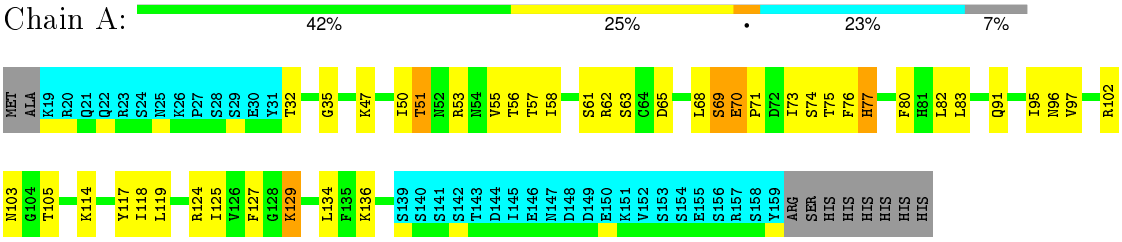
Chain A:



4.2.13 Score per residue for model 13

• Molecule 1: DNA damage response protein kinase DUN1

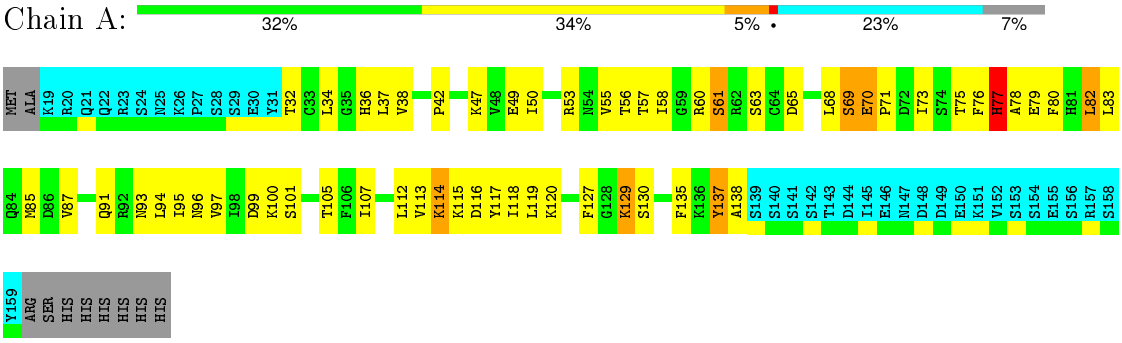
Chain A:



4.2.14 Score per residue for model 14

• Molecule 1: DNA damage response protein kinase DUN1

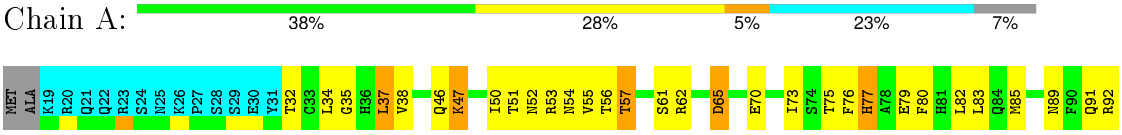
Chain A:

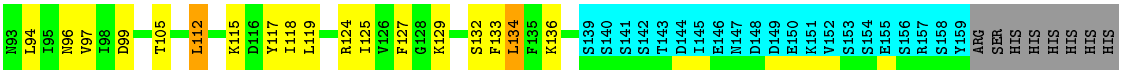


4.2.15 Score per residue for model 15

• Molecule 1: DNA damage response protein kinase DUN1

Chain A:

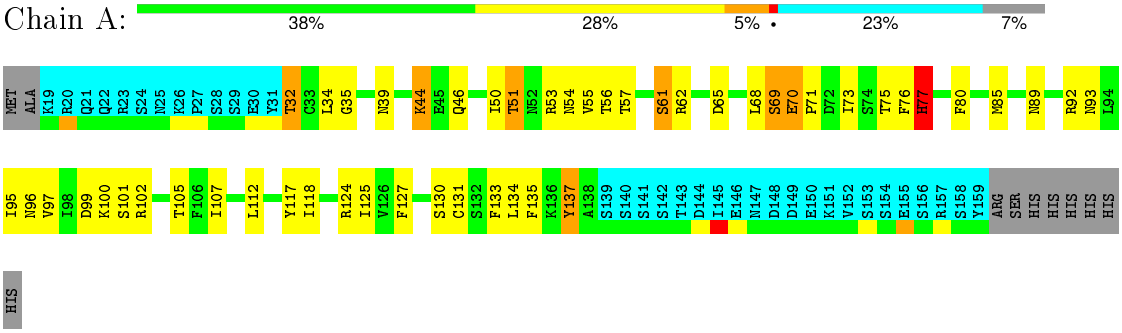




4.2.16 Score per residue for model 16

• Molecule 1: DNA damage response protein kinase DUN1

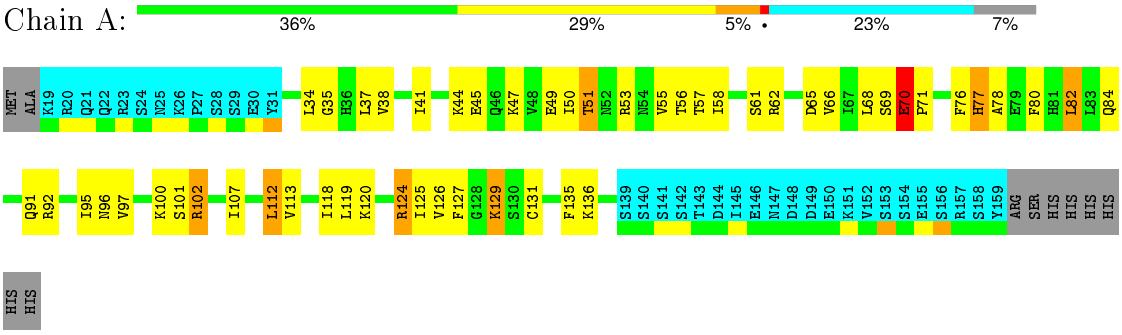
Chain A:



4.2.17 Score per residue for model 17

• Molecule 1: DNA damage response protein kinase DUN1

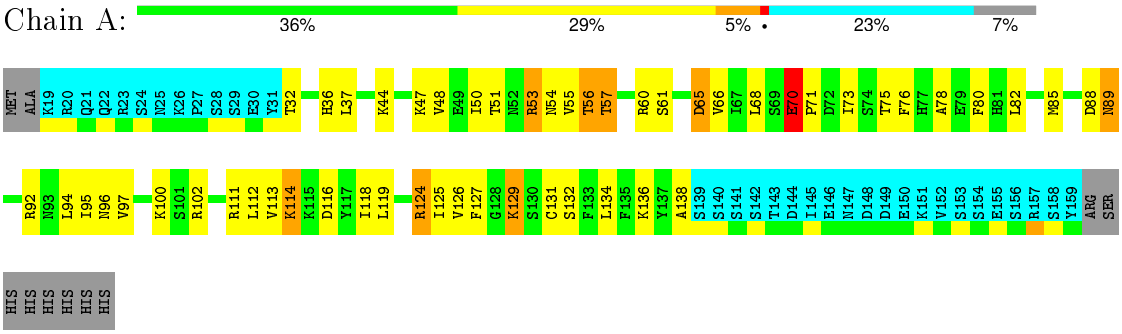
Chain A:



4.2.18 Score per residue for model 18

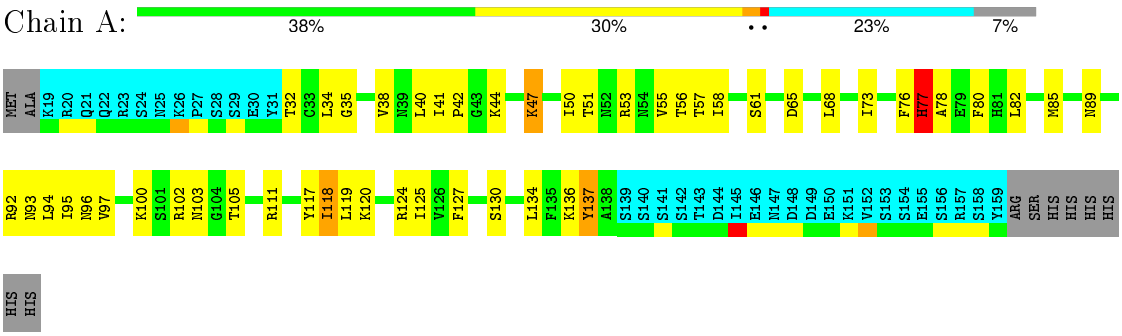
• Molecule 1: DNA damage response protein kinase DUN1

Chain A:



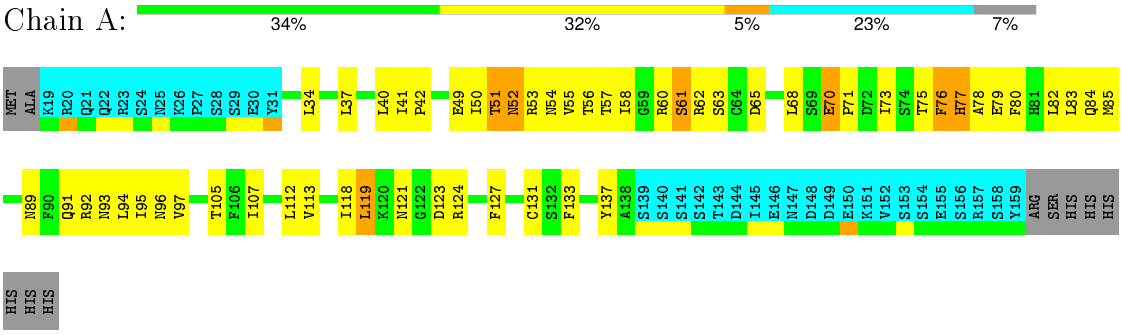
4.2.19 Score per residue for model 19

• Molecule 1: DNA damage response protein kinase DUN1



4.2.20 Score per residue for model 20

• Molecule 1: DNA damage response protein kinase DUN1



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 least restraint violations*.

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

Software name	Classification	Version
CNS	structure solution	1.1
CNS	refinement	1.1

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

6.1 Standard geometry

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

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	857	874	871	32±6
All	All	17140	17480	17420	632

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 18.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:80:PHE:CE2	1:A:97:VAL:HG22	0.85	2.06	14	14
1:A:51:THR:O	1:A:82:LEU:HD13	0.84	1.72	2	4
1:A:80:PHE:CE1	1:A:97:VAL:HG13	0.82	2.10	17	13
1:A:56:THR:OG1	1:A:82:LEU:HD11	0.81	1.76	17	1
1:A:37:LEU:HD11	1:A:50:ILE:HD11	0.80	1.51	2	1
1:A:80:PHE:CZ	1:A:97:VAL:HG13	0.80	2.10	13	14
1:A:37:LEU:HD12	1:A:50:ILE:HD11	0.79	1.54	3	5
1:A:119:LEU:HD11	1:A:125:ILE:HD11	0.76	1.56	4	9
1:A:37:LEU:CD1	1:A:50:ILE:HD11	0.76	2.11	2	5
1:A:94:LEU:HD12	1:A:118:ILE:HD11	0.76	1.56	15	1
1:A:52:ASN:HA	1:A:82:LEU:HD22	0.74	1.60	2	5
1:A:107:ILE:HD12	1:A:124:ARG:O	0.73	1.82	16	4
1:A:34:LEU:HD22	1:A:82:LEU:HD21	0.73	1.61	2	2
1:A:61:SER:HB2	1:A:75:THR:HG23	0.72	1.60	14	12

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:58:ILE:HG23	1:A:68:LEU:HD11	0.71	1.63	3	1
1:A:56:THR:HG23	1:A:65:ASP:CB	0.70	2.16	5	18
1:A:68:LEU:HD23	1:A:133:PHE:CE2	0.70	2.22	10	1
1:A:58:ILE:HG23	1:A:68:LEU:HD13	0.70	1.62	6	1
1:A:119:LEU:CD1	1:A:125:ILE:HD11	0.69	2.18	4	3
1:A:107:ILE:HG22	1:A:112:LEU:CD2	0.69	2.18	14	1
1:A:119:LEU:HD23	1:A:120:LYS:N	0.68	2.03	3	2
1:A:68:LEU:HD13	1:A:73:ILE:HD11	0.68	1.64	1	1
1:A:61:SER:HB3	1:A:75:THR:HG23	0.68	1.64	2	4
1:A:51:THR:O	1:A:82:LEU:HD23	0.68	1.88	9	3
1:A:107:ILE:HG22	1:A:112:LEU:HD21	0.68	1.66	14	1
1:A:68:LEU:HD13	1:A:73:ILE:HG12	0.67	1.66	19	2
1:A:56:THR:HG23	1:A:65:ASP:HB3	0.66	1.66	20	17
1:A:107:ILE:HD13	1:A:125:ILE:HG22	0.66	1.65	6	1
1:A:34:LEU:HD13	1:A:82:LEU:HD21	0.65	1.67	15	1
1:A:58:ILE:CG2	1:A:68:LEU:HD13	0.65	2.21	6	1
1:A:38:VAL:HG22	1:A:47:LYS:HG3	0.65	1.67	14	1
1:A:58:ILE:CG2	1:A:68:LEU:HD23	0.65	2.22	20	1
1:A:119:LEU:HD21	1:A:135:PHE:CD2	0.65	2.27	7	2
1:A:78:ALA:HB3	1:A:127:PHE:CE2	0.64	2.28	3	6
1:A:112:LEU:HD13	1:A:113:VAL:N	0.64	2.07	17	1
1:A:35:GLY:C	1:A:50:ILE:HG22	0.63	2.14	13	10
1:A:78:ALA:HB3	1:A:127:PHE:CZ	0.63	2.29	14	5
1:A:99:ASP:HB2	1:A:112:LEU:HD11	0.62	1.69	15	1
1:A:112:LEU:HD21	1:A:114:LYS:O	0.61	1.94	18	1
1:A:124:ARG:CA	1:A:134:LEU:HD12	0.61	2.25	3	1
1:A:107:ILE:CG2	1:A:125:ILE:HG22	0.61	2.24	3	1
1:A:58:ILE:CG2	1:A:68:LEU:HD11	0.61	2.25	19	4
1:A:68:LEU:CD1	1:A:73:ILE:HD11	0.61	2.26	1	2
1:A:119:LEU:HD21	1:A:123:ASP:HB2	0.61	1.73	3	1
1:A:68:LEU:HD12	1:A:73:ILE:HD11	0.61	1.71	13	2
1:A:107:ILE:HD13	1:A:125:ILE:HD12	0.60	1.74	8	1
1:A:61:SER:CB	1:A:75:THR:HG23	0.60	2.27	8	10
1:A:105:THR:HG23	1:A:127:PHE:CD1	0.60	2.31	6	4
1:A:69:SER:O	1:A:70:GLU:CB	0.60	2.50	9	11
1:A:34:LEU:HD22	1:A:82:LEU:HD23	0.59	1.75	5	1
1:A:118:ILE:C	1:A:118:ILE:HD13	0.58	2.18	9	1
1:A:68:LEU:HD22	1:A:131:CYS:HB3	0.57	1.75	17	4
1:A:80:PHE:CE2	1:A:97:VAL:HG13	0.57	2.35	13	1
1:A:50:ILE:HG21	1:A:135:PHE:CE2	0.57	2.35	8	4
1:A:50:ILE:HG22	1:A:82:LEU:HD22	0.57	1.77	18	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:34:LEU:HD22	1:A:82:LEU:HD11	0.57	1.77	11	2
1:A:95:ILE:HD12	1:A:95:ILE:N	0.57	2.15	1	5
1:A:80:PHE:CZ	1:A:97:VAL:HG22	0.57	2.35	2	6
1:A:34:LEU:HD11	1:A:93:ASN:OD1	0.57	2.00	16	1
1:A:50:ILE:HG22	1:A:82:LEU:HD12	0.56	1.75	2	1
1:A:37:LEU:CD1	1:A:66:VAL:HG21	0.56	2.30	17	1
1:A:113:VAL:O	1:A:113:VAL:HG22	0.56	2.00	3	3
1:A:118:ILE:HD13	1:A:118:ILE:C	0.56	2.20	19	1
1:A:122:GLY:HA2	1:A:134:LEU:HD21	0.56	1.77	11	2
1:A:70:GLU:CB	1:A:71:PRO:HD3	0.56	2.30	17	8
1:A:34:LEU:HD22	1:A:52:ASN:ND2	0.56	2.15	20	1
1:A:40:LEU:HD13	1:A:124:ARG:NH2	0.56	2.16	1	1
1:A:32:THR:N	1:A:51:THR:HG22	0.56	2.16	4	5
1:A:70:GLU:HA	1:A:73:ILE:HG22	0.56	1.78	18	3
1:A:58:ILE:HG22	1:A:68:LEU:HD11	0.56	1.78	19	2
1:A:68:LEU:HD12	1:A:131:CYS:HB3	0.56	1.78	20	1
1:A:68:LEU:HD23	1:A:133:PHE:HE2	0.55	1.60	10	1
1:A:48:VAL:HG11	1:A:66:VAL:HG13	0.55	1.78	2	2
1:A:50:ILE:HG21	1:A:82:LEU:HD13	0.55	1.77	14	1
1:A:35:GLY:O	1:A:50:ILE:HG22	0.55	2.03	19	7
1:A:112:LEU:HD22	1:A:114:LYS:O	0.55	2.02	1	1
1:A:36:HIS:ND1	1:A:138:ALA:HB2	0.54	2.17	8	6
1:A:70:GLU:HB3	1:A:71:PRO:HD3	0.54	1.80	3	3
1:A:94:LEU:HD22	1:A:118:ILE:HD11	0.54	1.79	3	1
1:A:94:LEU:HD23	1:A:118:ILE:HD11	0.53	1.80	6	1
1:A:107:ILE:CD1	1:A:125:ILE:HD13	0.53	2.33	10	1
1:A:41:ILE:HG22	1:A:42:PRO:HD2	0.53	1.80	3	1
1:A:70:GLU:CB	1:A:71:PRO:CD	0.53	2.87	8	7
1:A:34:LEU:HD13	1:A:82:LEU:HD23	0.53	1.80	12	1
1:A:50:ILE:HG22	1:A:82:LEU:CD1	0.53	2.34	2	2
1:A:50:ILE:CG2	1:A:82:LEU:HD13	0.53	2.33	14	1
1:A:67:ILE:O	1:A:67:ILE:HG23	0.53	2.03	1	1
1:A:34:LEU:HD21	1:A:93:ASN:ND2	0.52	2.19	3	1
1:A:50:ILE:HG21	1:A:135:PHE:HE2	0.52	1.65	4	5
1:A:97:VAL:CG2	1:A:119:LEU:HD12	0.52	2.35	3	1
1:A:107:ILE:HG22	1:A:125:ILE:HG22	0.52	1.82	3	1
1:A:37:LEU:N	1:A:37:LEU:HD23	0.51	2.20	15	1
1:A:58:ILE:HD12	1:A:58:ILE:N	0.51	2.20	17	3
1:A:70:GLU:CA	1:A:73:ILE:HG22	0.51	2.34	18	1
1:A:127:PHE:CE1	1:A:133:PHE:CZ	0.51	2.98	9	1
1:A:126:VAL:HG22	1:A:132:SER:OG	0.51	2.06	4	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:124:ARG:CG	1:A:134:LEU:HD12	0.51	2.36	2	1
1:A:113:VAL:HG22	1:A:113:VAL:O	0.51	2.06	12	1
1:A:124:ARG:HD2	1:A:134:LEU:HD12	0.51	1.82	7	1
1:A:124:ARG:CB	1:A:134:LEU:HD12	0.51	2.35	3	1
1:A:53:ARG:C	1:A:55:VAL:H	0.51	2.10	16	20
1:A:69:SER:O	1:A:70:GLU:HB2	0.50	2.05	4	1
1:A:77:HIS:CD2	1:A:127:PHE:CD1	0.50	2.99	20	3
1:A:82:LEU:HD12	1:A:95:ILE:HG13	0.50	1.83	13	1
1:A:68:LEU:HD11	1:A:133:PHE:CE2	0.50	2.41	16	1
1:A:68:LEU:HD13	1:A:73:ILE:CD1	0.50	2.34	1	1
1:A:118:ILE:HG23	1:A:118:ILE:O	0.50	2.05	20	13
1:A:69:SER:C	1:A:71:PRO:HD2	0.50	2.27	4	2
1:A:82:LEU:HD13	1:A:95:ILE:HG13	0.50	1.84	4	1
1:A:118:ILE:O	1:A:118:ILE:HG23	0.50	2.07	7	7
1:A:50:ILE:CG2	1:A:82:LEU:HD12	0.49	2.36	2	1
1:A:69:SER:OG	1:A:73:ILE:HD13	0.49	2.08	7	1
1:A:76:PHE:O	1:A:77:HIS:C	0.49	2.49	13	18
1:A:119:LEU:C	1:A:119:LEU:HD13	0.49	2.28	17	1
1:A:34:LEU:HD21	1:A:93:ASN:OD1	0.49	2.06	2	2
1:A:58:ILE:HG23	1:A:68:LEU:HD23	0.49	1.83	20	1
1:A:125:ILE:N	1:A:125:ILE:HD13	0.49	2.22	6	1
1:A:41:ILE:HD12	1:A:45:GLU:HA	0.49	1.85	17	1
1:A:94:LEU:HD21	1:A:118:ILE:HG12	0.49	1.83	9	1
1:A:98:ILE:HD13	1:A:115:LYS:O	0.49	2.08	8	1
1:A:36:HIS:CE1	1:A:138:ALA:HB2	0.49	2.43	14	1
1:A:34:LEU:HD11	1:A:93:ASN:ND2	0.48	2.22	20	1
1:A:34:LEU:HD23	1:A:34:LEU:O	0.48	2.08	4	1
1:A:50:ILE:HG22	1:A:82:LEU:CD2	0.48	2.37	18	1
1:A:126:VAL:HG13	1:A:131:CYS:O	0.48	2.09	17	2
1:A:50:ILE:HG12	1:A:82:LEU:HD22	0.48	1.84	13	1
1:A:78:ALA:HB3	1:A:127:PHE:HE2	0.48	1.67	3	1
1:A:80:PHE:HE1	1:A:97:VAL:HG13	0.48	1.68	6	1
1:A:137:TYR:N	1:A:137:TYR:CD1	0.48	2.79	19	3
1:A:134:LEU:HD23	1:A:135:PHE:N	0.48	2.24	8	1
1:A:68:LEU:HD23	1:A:131:CYS:HB2	0.47	1.86	11	1
1:A:95:ILE:N	1:A:95:ILE:HD12	0.47	2.24	16	4
1:A:68:LEU:HD22	1:A:131:CYS:CB	0.47	2.40	1	1
1:A:32:THR:HG23	1:A:52:ASN:ND2	0.47	2.25	4	1
1:A:105:THR:HG23	1:A:127:PHE:CE1	0.47	2.44	1	5
1:A:34:LEU:C	1:A:34:LEU:HD23	0.47	2.29	4	1
1:A:41:ILE:CG2	1:A:42:PRO:HD2	0.47	2.40	8	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:34:LEU:HD22	1:A:93:ASN:OD1	0.47	2.10	19	1
1:A:94:LEU:HG	1:A:118:ILE:HD11	0.47	1.86	20	3
1:A:126:VAL:HG22	1:A:132:SER:HB3	0.47	1.86	9	3
1:A:52:ASN:CA	1:A:82:LEU:HD22	0.47	2.39	3	1
1:A:56:THR:N	1:A:80:PHE:O	0.47	2.48	15	18
1:A:54:ASN:O	1:A:81:HIS:HA	0.47	2.10	10	2
1:A:58:ILE:CG2	1:A:68:LEU:HD12	0.47	2.39	4	1
1:A:124:ARG:N	1:A:134:LEU:HD12	0.47	2.25	9	1
1:A:68:LEU:CD2	1:A:73:ILE:HD11	0.47	2.40	16	1
1:A:57:THR:O	1:A:65:ASP:N	0.46	2.48	11	6
1:A:105:THR:CG2	1:A:127:PHE:CE1	0.46	2.98	10	6
1:A:124:ARG:HA	1:A:134:LEU:HD12	0.46	1.86	3	1
1:A:56:THR:HG23	1:A:65:ASP:HB2	0.46	1.84	17	1
1:A:82:LEU:HD23	1:A:95:ILE:HD12	0.46	1.87	17	1
1:A:121:ASN:ND2	1:A:137:TYR:CD2	0.46	2.84	20	1
1:A:37:LEU:HD11	1:A:66:VAL:HG21	0.46	1.87	2	1
1:A:34:LEU:HD11	1:A:93:ASN:HD21	0.46	1.69	20	1
1:A:51:THR:O	1:A:82:LEU:HD22	0.46	2.11	6	1
1:A:133:PHE:CD1	1:A:133:PHE:N	0.46	2.83	8	3
1:A:37:LEU:HD23	1:A:133:PHE:CE2	0.46	2.46	12	1
1:A:38:VAL:HG22	1:A:47:LYS:HD3	0.46	1.87	17	1
1:A:125:ILE:CG2	1:A:127:PHE:CE1	0.46	2.99	9	1
1:A:95:ILE:HD13	1:A:119:LEU:HD23	0.46	1.87	6	2
1:A:68:LEU:HD23	1:A:73:ILE:CD1	0.45	2.41	16	1
1:A:107:ILE:HD13	1:A:125:ILE:HG12	0.45	1.88	17	2
1:A:73:ILE:O	1:A:73:ILE:CG2	0.45	2.64	8	5
1:A:38:VAL:HG22	1:A:47:LYS:HD2	0.45	1.88	15	1
1:A:135:PHE:CD1	1:A:136:LYS:N	0.45	2.85	3	2
1:A:107:ILE:HD13	1:A:125:ILE:CG1	0.45	2.41	16	1
1:A:60:ARG:CD	1:A:70:GLU:O	0.45	2.65	2	5
1:A:82:LEU:HG	1:A:95:ILE:HD12	0.45	1.88	12	1
1:A:126:VAL:HG22	1:A:132:SER:CB	0.45	2.42	12	1
1:A:36:HIS:CE1	1:A:138:ALA:CB	0.45	2.99	2	1
1:A:119:LEU:HD21	1:A:135:PHE:CG	0.45	2.46	2	1
1:A:133:PHE:N	1:A:133:PHE:CD1	0.45	2.84	3	3
1:A:99:ASP:CB	1:A:112:LEU:HD11	0.45	2.41	15	1
1:A:78:ALA:CB	1:A:127:PHE:CZ	0.45	3.00	20	2
1:A:53:ARG:O	1:A:55:VAL:N	0.44	2.50	6	4
1:A:125:ILE:CG2	1:A:133:PHE:CE2	0.44	2.99	8	1
1:A:107:ILE:CD1	1:A:125:ILE:HG22	0.44	2.37	6	1
1:A:125:ILE:CG2	1:A:127:PHE:CE2	0.44	3.00	16	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:125:ILE:CD1	1:A:133:PHE:CE2	0.44	3.00	3	1
1:A:34:LEU:HD13	1:A:34:LEU:O	0.44	2.11	10	1
1:A:60:ARG:HB2	1:A:73:ILE:HG23	0.44	1.88	18	3
1:A:60:ARG:HD2	1:A:70:GLU:O	0.44	2.12	6	1
1:A:38:VAL:HG22	1:A:47:LYS:HG2	0.44	1.88	19	1
1:A:118:ILE:C	1:A:118:ILE:CD1	0.44	2.86	9	1
1:A:68:LEU:HD23	1:A:131:CYS:CB	0.44	2.43	8	1
1:A:99:ASP:OD1	1:A:112:LEU:HD22	0.44	2.12	11	1
1:A:37:LEU:HD13	1:A:133:PHE:CE1	0.44	2.47	9	1
1:A:61:SER:OG	1:A:75:THR:HG23	0.44	2.13	10	1
1:A:94:LEU:N	1:A:94:LEU:HD12	0.44	2.27	4	1
1:A:124:ARG:CZ	1:A:134:LEU:HD23	0.44	2.43	15	1
1:A:119:LEU:HD21	1:A:123:ASP:CB	0.44	2.42	3	1
1:A:32:THR:C	1:A:51:THR:HG22	0.44	2.32	6	1
1:A:119:LEU:HD23	1:A:119:LEU:C	0.44	2.32	1	2
1:A:58:ILE:HG21	1:A:68:LEU:HD11	0.44	1.89	13	1
1:A:68:LEU:O	1:A:69:SER:O	0.43	2.36	5	2
1:A:68:LEU:HD12	1:A:68:LEU:N	0.43	2.28	3	1
1:A:80:PHE:HE2	1:A:119:LEU:HD13	0.43	1.72	6	1
1:A:53:ARG:O	1:A:54:ASN:HB2	0.43	2.13	6	1
1:A:82:LEU:HD23	1:A:82:LEU:C	0.43	2.33	15	1
1:A:52:ASN:HA	1:A:82:LEU:HB3	0.43	1.89	5	1
1:A:113:VAL:HG13	1:A:114:LYS:N	0.43	2.29	11	2
1:A:34:LEU:HD12	1:A:84:GLN:CG	0.43	2.42	17	1
1:A:56:THR:HG22	1:A:65:ASP:CB	0.43	2.44	11	2
1:A:58:ILE:CD1	1:A:80:PHE:CD1	0.43	3.02	20	1
1:A:73:ILE:HG23	1:A:73:ILE:O	0.43	2.13	15	1
1:A:118:ILE:CD1	1:A:118:ILE:C	0.43	2.88	19	1
1:A:68:LEU:HD23	1:A:73:ILE:HD11	0.43	1.91	16	1
1:A:37:LEU:HD12	1:A:66:VAL:HG21	0.42	1.90	11	2
1:A:50:ILE:HG21	1:A:135:PHE:HE1	0.42	1.74	16	1
1:A:118:ILE:HD13	1:A:119:LEU:N	0.42	2.29	19	2
1:A:68:LEU:HD21	1:A:131:CYS:CB	0.42	2.44	16	1
1:A:128:GLY:O	1:A:129:LYS:HB2	0.42	2.13	10	1
1:A:80:PHE:CE2	1:A:119:LEU:HD13	0.42	2.49	6	1
1:A:68:LEU:HD13	1:A:127:PHE:CB	0.42	2.44	7	1
1:A:124:ARG:HG2	1:A:134:LEU:HD12	0.42	1.90	18	1
1:A:68:LEU:HD21	1:A:133:PHE:CE1	0.42	2.50	7	1
1:A:124:ARG:HG3	1:A:134:LEU:HD12	0.42	1.90	4	2
1:A:37:LEU:O	1:A:37:LEU:HD12	0.42	2.14	8	1
1:A:107:ILE:HG22	1:A:112:LEU:CD1	0.42	2.44	6	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:93:ASN:C	1:A:94:LEU:HD22	0.42	2.35	14	1
1:A:58:ILE:HD11	1:A:80:PHE:CD2	0.42	2.49	1	1
1:A:40:LEU:HD13	1:A:124:ARG:NH1	0.42	2.30	20	1
1:A:106:PHE:CD1	1:A:106:PHE:N	0.42	2.87	10	1
1:A:73:ILE:CD1	1:A:77:HIS:CE1	0.42	3.02	5	1
1:A:60:ARG:HB2	1:A:73:ILE:CG2	0.42	2.44	2	1
1:A:70:GLU:HB3	1:A:71:PRO:CD	0.42	2.45	9	1
1:A:107:ILE:HG21	1:A:112:LEU:CD1	0.41	2.45	2	1
1:A:37:LEU:HD22	1:A:133:PHE:HD2	0.41	1.75	20	1
1:A:73:ILE:HD12	1:A:77:HIS:CE1	0.41	2.51	14	1
1:A:87:VAL:O	1:A:87:VAL:HG13	0.41	2.14	14	2
1:A:82:LEU:HD22	1:A:95:ILE:HG23	0.41	1.92	4	1
1:A:41:ILE:HG23	1:A:42:PRO:HD2	0.41	1.93	8	3
1:A:68:LEU:HD11	1:A:133:PHE:HE2	0.41	1.74	16	1
1:A:82:LEU:HA	1:A:95:ILE:HD13	0.41	1.92	17	1
1:A:56:THR:CG2	1:A:65:ASP:CB	0.41	2.98	8	1
1:A:80:PHE:CD2	1:A:97:VAL:HG22	0.41	2.50	8	1
1:A:118:ILE:CG2	1:A:118:ILE:O	0.41	2.68	20	2
1:A:53:ARG:CD	1:A:55:VAL:O	0.41	2.69	1	1
1:A:60:ARG:HB3	1:A:70:GLU:HA	0.41	1.93	12	1
1:A:58:ILE:HG21	1:A:127:PHE:CD2	0.41	2.51	11	1
1:A:73:ILE:CG2	1:A:73:ILE:O	0.41	2.69	7	2
1:A:118:ILE:O	1:A:118:ILE:CG2	0.41	2.69	5	2
1:A:80:PHE:CZ	1:A:97:VAL:CG1	0.41	2.99	20	1
1:A:50:ILE:HG22	1:A:82:LEU:HD13	0.41	1.93	3	1
1:A:124:ARG:HB2	1:A:134:LEU:HD12	0.41	1.93	3	1
1:A:50:ILE:HD11	1:A:56:THR:HG21	0.41	1.92	5	1
1:A:121:ASN:OD1	1:A:137:TYR:CG	0.41	2.74	7	1
1:A:34:LEU:HD13	1:A:52:ASN:OD1	0.41	2.16	2	1
1:A:113:VAL:HG23	1:A:114:LYS:N	0.41	2.31	14	1
1:A:68:LEU:O	1:A:69:SER:C	0.41	2.59	9	1
1:A:53:ARG:CB	1:A:53:ARG:CZ	0.40	2.99	1	1
1:A:58:ILE:HD13	1:A:80:PHE:CD1	0.40	2.51	10	1
1:A:124:ARG:NH1	1:A:134:LEU:HD23	0.40	2.31	15	1
1:A:78:ALA:CB	1:A:127:PHE:CE1	0.40	3.04	11	1
1:A:128:GLY:O	1:A:129:LYS:CB	0.40	2.69	11	1
1:A:41:ILE:HD12	1:A:41:ILE:N	0.40	2.32	6	1
1:A:36:HIS:NE2	1:A:49:GLU:CG	0.40	2.85	6	1
1:A:113:VAL:O	1:A:113:VAL:HG13	0.40	2.14	7	1
1:A:39:ASN:N	1:A:39:ASN:OD1	0.40	2.55	6	1
1:A:37:LEU:HD11	1:A:50:ILE:HB	0.40	1.92	4	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:53:ARG:C	1:A:55:VAL:N	0.40	2.75	16	1
1:A:40:LEU:HD13	1:A:124:ARG:HH12	0.40	1.76	20	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	107/151 (71%)	90±2 (84±2%)	12±3 (11±2%)	5±2 (4±2%)	6	31
All	All	2140/3020 (71%)	1804 (84%)	244 (11%)	92 (4%)	6	31

All 16 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	77	HIS	18
1	A	70	GLU	17
1	A	129	LYS	12
1	A	89	ASN	12
1	A	44	LYS	5
1	A	101	SER	5
1	A	69	SER	4
1	A	52	ASN	4
1	A	42	PRO	3
1	A	104	GLY	3
1	A	90	PHE	2
1	A	102	ARG	2
1	A	68	LEU	2
1	A	116	ASP	1
1	A	113	VAL	1
1	A	130	SER	1

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	98/141 (70%)	76±3 (78±3%)	22±3 (22±3%)	4 31
All	All	1960/2820 (70%)	1521 (78%)	439 (22%)	4 31

All 64 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	96	ASN	20
1	A	57	THR	20
1	A	51	THR	18
1	A	32	THR	17
1	A	61	SER	14
1	A	85	MET	14
1	A	102	ARG	13
1	A	129	LYS	12
1	A	115	LYS	11
1	A	62	ARG	11
1	A	112	LEU	11
1	A	100	LYS	11
1	A	92	ARG	10
1	A	82	LEU	10
1	A	47	LYS	9
1	A	117	TYR	9
1	A	83	LEU	9
1	A	103	ASN	9
1	A	70	GLU	9
1	A	114	LYS	8
1	A	44	LYS	8
1	A	124	ARG	8
1	A	63	SER	8
1	A	77	HIS	8
1	A	54	ASN	8
1	A	116	ASP	7
1	A	91	GLN	7
1	A	65	ASP	7
1	A	120	LYS	7

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Mol	Chain	Res	Type	Models (Total)
1	A	49	GLU	6
1	A	46	GLN	6
1	A	74	SER	6
1	A	99	ASP	6
1	A	79	GLU	6
1	A	123	ASP	6
1	A	132	SER	6
1	A	136	LYS	6
1	A	111	ARG	6
1	A	130	SER	6
1	A	134	LEU	5
1	A	76	PHE	5
1	A	84	GLN	5
1	A	89	ASN	4
1	A	88	ASP	4
1	A	53	ARG	4
1	A	40	LEU	4
1	A	94	LEU	4
1	A	137	TYR	3
1	A	39	ASN	3
1	A	34	LEU	3
1	A	37	LEU	2
1	A	125	ILE	2
1	A	60	ARG	2
1	A	56	THR	2
1	A	52	ASN	2
1	A	118	ILE	2
1	A	69	SER	2
1	A	119	LEU	2
1	A	121	ASN	1
1	A	105	THR	1
1	A	110	ASN	1
1	A	68	LEU	1
1	A	101	SER	1
1	A	45	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

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided