



# Full wwPDB NMR Structure Validation Report ⓘ

Apr 27, 2016 – 01:44 AM BST

PDB ID : 2LSK  
Title : C-terminal domain of human REV1 in complex with DNA-polymerase H (eta)  
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Deposited on : 2012-05-01

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 : 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

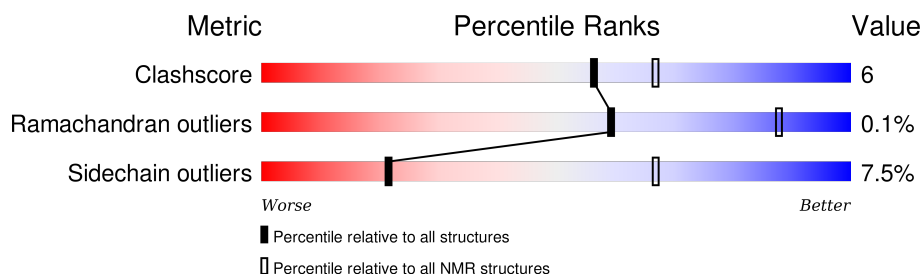
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 76%.

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	95	
2	B	16	

## 2 Ensemble composition and analysis ⓘ

This entry contains 20 models. Model 11 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:1159-A:1178, A:1184-A:1219, A:1227-A:1250, B:531-B:539 (89)	0.37	11

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

NmrClust was unable to cluster the ensemble.

Error message: Inconsistent models

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

There are 2 unique types of molecules in this entry. The entry contains 1814 atoms, of which 914 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called DNA repair protein REV1.

Mol	Chain	Residues	Atoms						Trace
1	A	95	Total	C	H	N	O	S	0
			1552	494	781	120	152	5	

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1157	GLY	-	EXPRESSION TAG	UNP Q9UBZ9

- Molecule 2 is a protein called DNA polymerase eta.

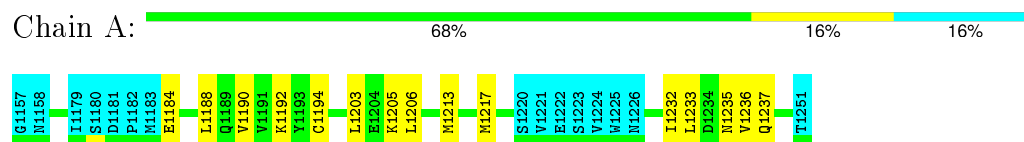
Mol	Chain	Residues	Atoms					Trace
2	B	16	Total	C	H	N	O	0
			262	84	133	20	25	

## 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: DNA repair protein REV1



- Molecule 2: DNA polymerase eta

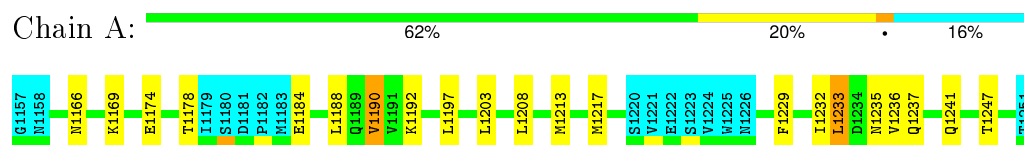


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

#### 4.2.1 Score per residue for model 1

- Molecule 1: DNA repair protein REV1

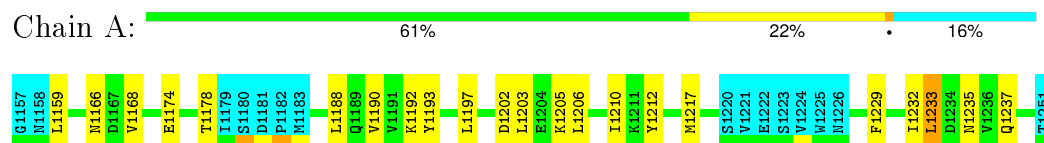


- Molecule 2: DNA polymerase eta



### 4.2.2 Score per residue for model 2

- Molecule 1: DNA repair protein REV1

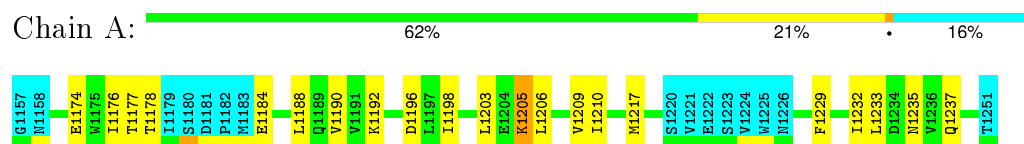


- Molecule 2: DNA polymerase eta

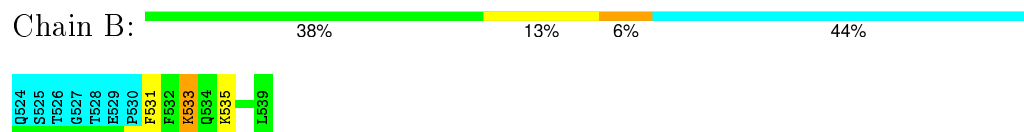


### 4.2.3 Score per residue for model 3

- Molecule 1: DNA repair protein REV1

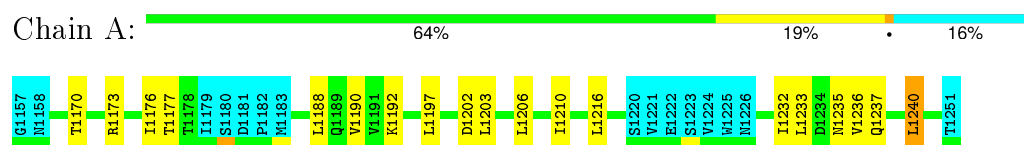


- Molecule 2: DNA polymerase eta



### 4.2.4 Score per residue for model 4

- Molecule 1: DNA repair protein REV1

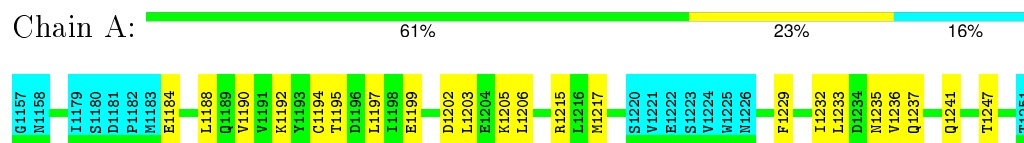


- Molecule 2: DNA polymerase eta



### 4.2.5 Score per residue for model 5

- Molecule 1: DNA repair protein REV1

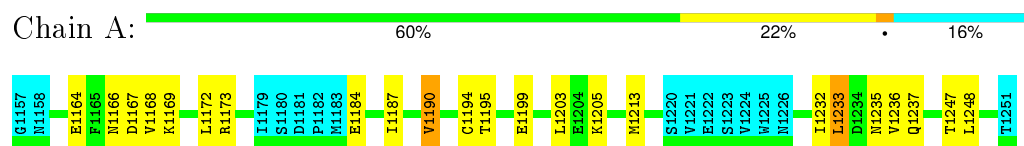


- Molecule 2: DNA polymerase eta

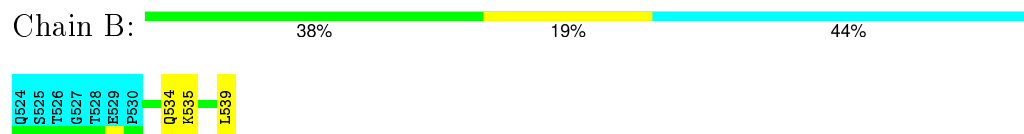


### 4.2.6 Score per residue for model 6

- Molecule 1: DNA repair protein REV1

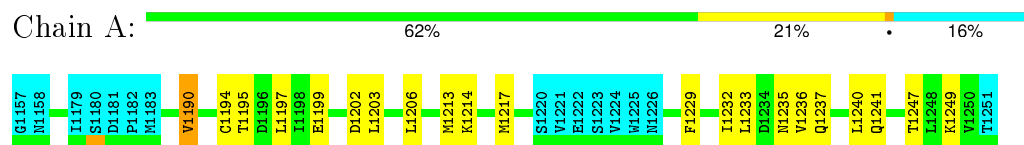


- Molecule 2: DNA polymerase eta



### 4.2.7 Score per residue for model 7

- Molecule 1: DNA repair protein REV1

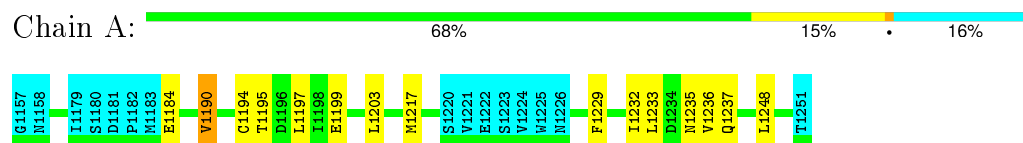


- Molecule 2: DNA polymerase eta



### 4.2.8 Score per residue for model 8

- Molecule 1: DNA repair protein REV1

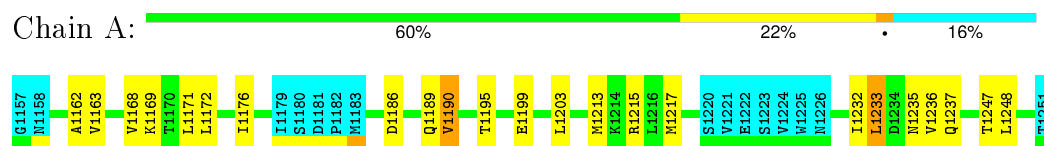


- Molecule 2: DNA polymerase eta

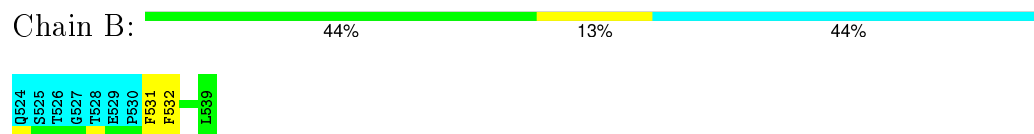


### 4.2.9 Score per residue for model 9

- Molecule 1: DNA repair protein REV1

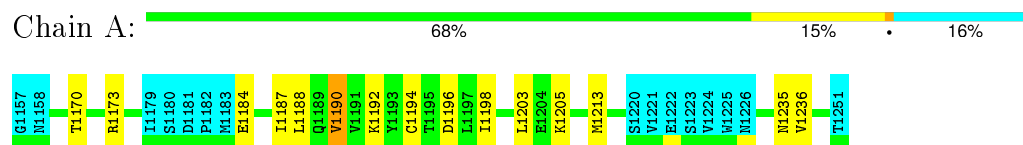


- Molecule 2: DNA polymerase eta

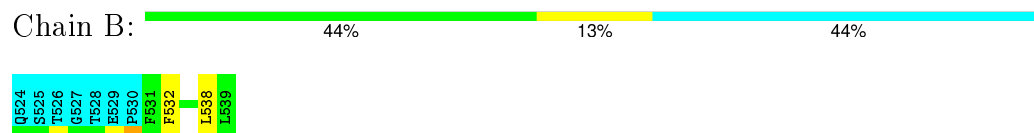


### 4.2.10 Score per residue for model 10

- Molecule 1: DNA repair protein REV1



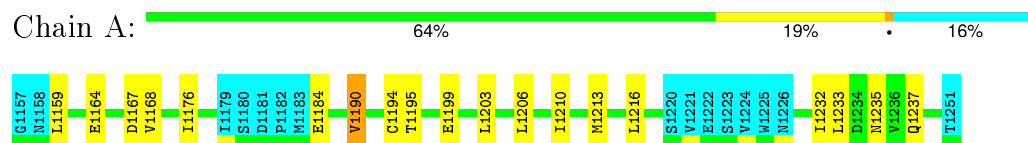
- Molecule 2: DNA polymerase eta



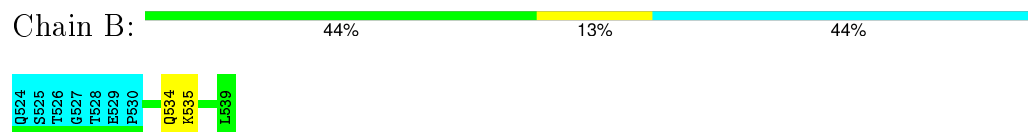


### 4.2.11 Score per residue for model 11 (medoid)

- Molecule 1: DNA repair protein REV1

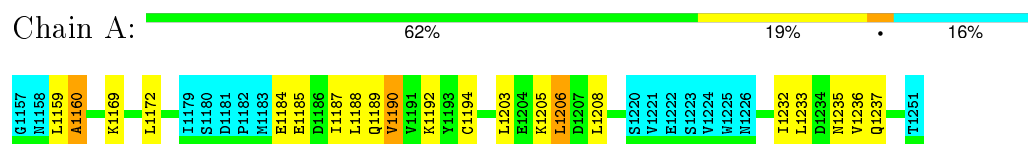


- Molecule 2: DNA polymerase eta

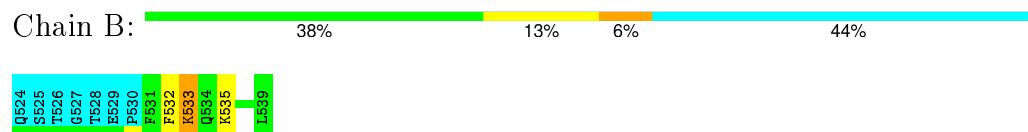


### 4.2.12 Score per residue for model 12

- Molecule 1: DNA repair protein REV1

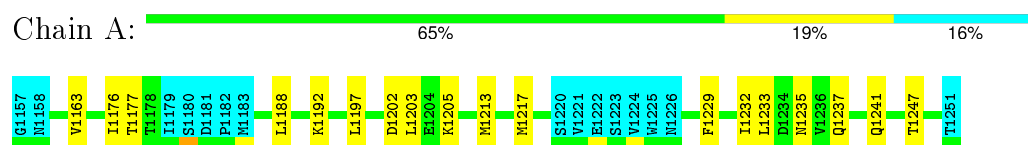


- Molecule 2: DNA polymerase eta



### 4.2.13 Score per residue for model 13

- Molecule 1: DNA repair protein REV1

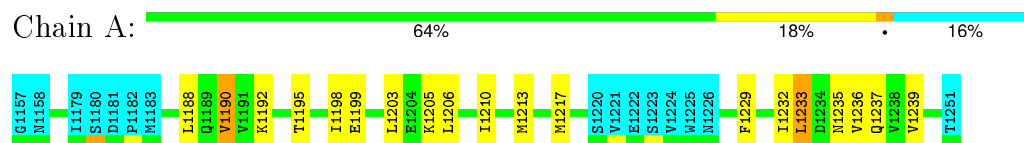


- Molecule 2: DNA polymerase eta



#### 4.2.14 Score per residue for model 14

- Molecule 1: DNA repair protein REV1

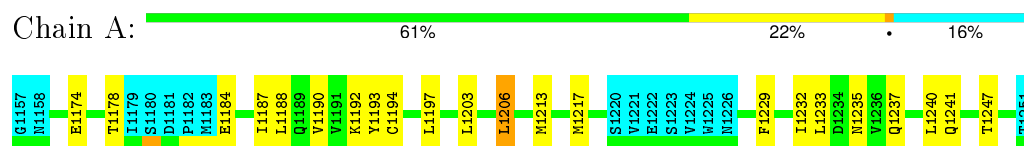


- Molecule 2: DNA polymerase eta



#### 4.2.15 Score per residue for model 15

- Molecule 1: DNA repair protein REV1

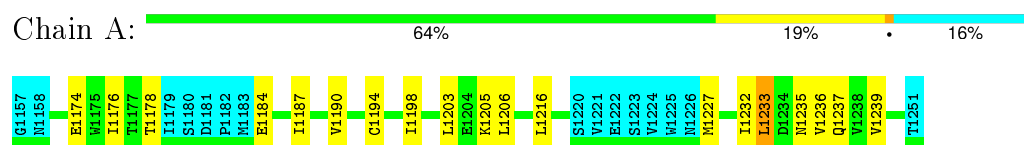


- Molecule 2: DNA polymerase eta



#### 4.2.16 Score per residue for model 16

- Molecule 1: DNA repair protein REV1

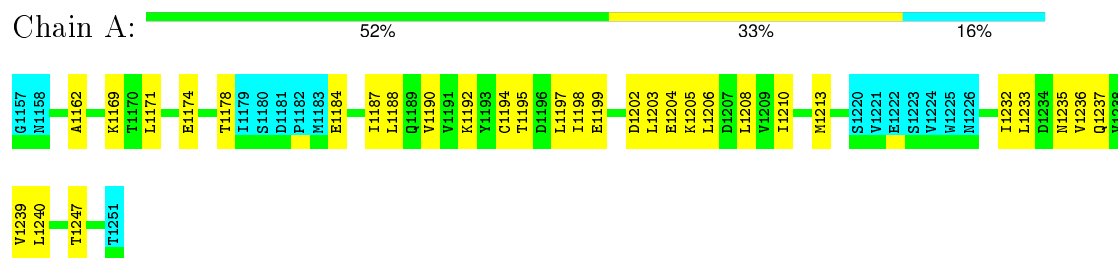


- Molecule 2: DNA polymerase eta

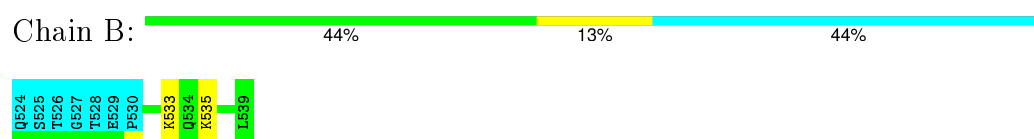


### 4.2.17 Score per residue for model 17

- Molecule 1: DNA repair protein REV1

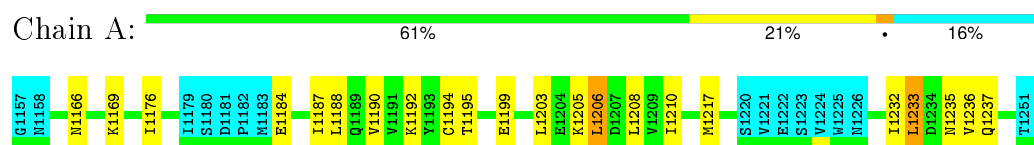


- Molecule 2: DNA polymerase eta



### 4.2.18 Score per residue for model 18

- Molecule 1: DNA repair protein REV1

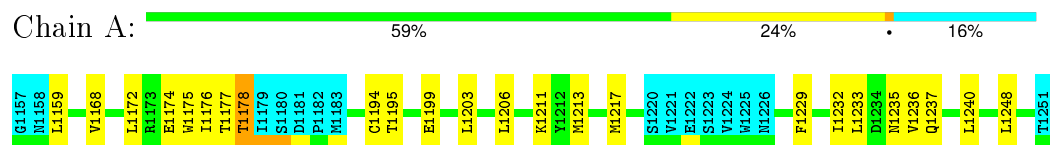


- Molecule 2: DNA polymerase eta



### 4.2.19 Score per residue for model 19

- Molecule 1: DNA repair protein REV1



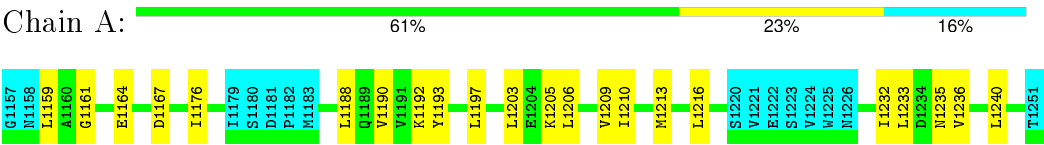
- Molecule 2: DNA polymerase eta





4.2.20 Score per residue for model 20

- Molecule 1: DNA repair protein REV1



- Molecule 2: DNA polymerase eta



## 5 Refinement protocol and experimental data overview

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

Of the 20 calculated structures, 20 were deposited, based on the following criterion: *all calculated structures submitted*.

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

Software name	Classification	Version
CYANA	structure solution	
CNS	refinement	

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)	2lsk_cs.str
Number of chemical shift lists	1
Total number of shifts	1118
Number of shifts mapped to atoms	1118
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	76%

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.

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	0.1±0.2
All	All	0	1

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

All unique planar outliers are listed below.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	1178	THR	Peptide	1

### 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	657	679	679	9±2
2	B	80	90	90	0±1
All	All	14740	15380	15380	191

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 6.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:1213:MET:SD	1:A:1232:ILE:HG12	0.69	2.28	13	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:1176:ILE:HD11	1:A:1217:MET:HG3	0.61	1.72	9	3
1:A:1194:CYS:SG	1:A:1206:LEU:HD21	0.61	2.36	18	8
1:A:1190:VAL:HB	1:A:1232:ILE:HD11	0.59	1.75	11	11
1:A:1198:ILE:HG21	1:A:1239:VAL:HG12	0.57	1.76	14	3
1:A:1190:VAL:HG11	1:A:1213:MET:SD	0.56	2.40	7	5
1:A:1176:ILE:HG21	1:A:1216:LEU:HB3	0.53	1.80	4	2
1:A:1197:LEU:HD12	1:A:1206:LEU:HB2	0.53	1.80	17	2
1:A:1194:CYS:SG	1:A:1236:VAL:HG22	0.52	2.44	19	8
1:A:1184:GLU:HA	1:A:1187:ILE:HD12	0.52	1.81	10	7
1:A:1186:ASP:HA	1:A:1189:GLN:HG2	0.52	1.82	9	1
1:A:1169:LYS:HE2	1:A:1208:LEU:HD21	0.52	1.80	1	1
1:A:1166:ASN:HA	1:A:1169:LYS:HD3	0.52	1.82	6	2
1:A:1176:ILE:HG23	1:A:1177:THR:HG23	0.51	1.82	3	4
1:A:1162:ALA:HB2	1:A:1171:LEU:HD13	0.51	1.83	17	2
1:A:1213:MET:SD	1:A:1232:ILE:HG21	0.51	2.45	6	2
2:B:539:LEU:HD23	2:B:539:LEU:H	0.50	1.67	19	1
1:A:1217:MET:SD	1:A:1229:PHE:HA	0.50	2.47	8	8
1:A:1169:LYS:HG2	1:A:1208:LEU:HD21	0.50	1.84	12	3
1:A:1217:MET:HB3	1:A:1229:PHE:HB2	0.50	1.83	15	2
1:A:1161:GLY:HA2	2:B:539:LEU:HD21	0.49	1.82	20	1
1:A:1197:LEU:HD12	1:A:1202:ASP:HB3	0.48	1.85	4	1
1:A:1241:GLN:HE21	1:A:1247:THR:HG23	0.48	1.68	15	5
1:A:1195:THR:O	1:A:1199:GLU:HG3	0.47	2.09	11	4
1:A:1237:GLN:NE2	1:A:1247:THR:HB	0.47	2.25	1	3
1:A:1233:LEU:O	1:A:1237:GLN:HG2	0.47	2.10	16	18
1:A:1227:MET:SD	1:A:1227:MET:N	0.47	2.88	16	1
1:A:1159:LEU:HG	1:A:1172:LEU:HD11	0.47	1.85	12	1
1:A:1197:LEU:HD22	1:A:1202:ASP:HB2	0.47	1.87	7	2
1:A:1195:THR:O	1:A:1199:GLU:HG2	0.46	2.10	8	6
1:A:1210:ILE:HD11	1:A:1236:VAL:HG21	0.46	1.87	20	2
1:A:1188:LEU:O	1:A:1192:LYS:HG2	0.46	2.11	10	13
1:A:1206:LEU:O	1:A:1210:ILE:HG12	0.45	2.11	3	6
1:A:1205:LYS:HE3	1:A:1205:LYS:HA	0.45	1.88	3	1
1:A:1190:VAL:HG23	1:A:1232:ILE:HD11	0.45	1.87	18	1
2:B:535:LYS:O	2:B:539:LEU:HD23	0.45	2.12	6	1
1:A:1159:LEU:HD11	1:A:1193:TYR:CZ	0.45	2.47	2	1
1:A:1176:ILE:HD13	1:A:1216:LEU:HB3	0.44	1.89	11	2
1:A:1193:TYR:O	1:A:1197:LEU:HG	0.44	2.12	15	2
1:A:1170:THR:HG23	1:A:1173:ARG:NH2	0.44	2.28	4	2
1:A:1197:LEU:HD22	1:A:1202:ASP:HB3	0.44	1.88	5	3
1:A:1164:GLU:HB2	1:A:1167:ASP:HB2	0.43	1.88	20	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:1236:VAL:O	1:A:1240:LEU:HB2	0.43	2.13	4	1
1:A:1232:ILE:O	1:A:1236:VAL:HG23	0.43	2.14	14	9
1:A:1159:LEU:HD23	1:A:1168:VAL:HB	0.43	1.90	11	1
1:A:1174:GLU:O	1:A:1178:THR:HG22	0.43	2.13	16	7
1:A:1209:VAL:O	1:A:1213:MET:HB3	0.43	2.14	20	1
1:A:1205:LYS:O	1:A:1209:VAL:HG23	0.43	2.13	3	1
2:B:533:LYS:HE3	2:B:533:LYS:HA	0.42	1.91	12	1
1:A:1159:LEU:HD12	1:A:1168:VAL:HB	0.42	1.90	19	1
1:A:1172:LEU:HA	1:A:1175:TRP:HB3	0.42	1.90	19	1
1:A:1176:ILE:HD11	1:A:1217:MET:SD	0.42	2.55	18	1
1:A:1190:VAL:HG22	2:B:532:PHE:CE2	0.41	2.51	9	1
1:A:1185:GLU:O	1:A:1189:GLN:HG2	0.41	2.16	12	1
1:A:1193:TYR:CE2	1:A:1197:LEU:HD11	0.41	2.50	20	1
1:A:1160:ALA:HB1	2:B:535:LYS:HB3	0.41	1.93	12	1
1:A:1194:CYS:HA	1:A:1206:LEU:HG	0.41	1.92	19	1
1:A:1159:LEU:HD22	1:A:1168:VAL:HB	0.41	1.91	2	1
1:A:1168:VAL:O	1:A:1172:LEU:HG	0.40	2.16	9	2
1:A:1160:ALA:HB2	2:B:532:PHE:HA	0.40	1.93	12	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	80/95 (84%)	78±1 (98±1%)	2±1 (2±1%)	0±0 (0±0%)	59	88
2	B	8/16 (50%)	7±1 (89±10%)	1±1 (11±10%)	0±0 (0±0%)	100	100
All	All	1760/2220 (79%)	1709 (97%)	49 (3%)	2 (0%)	59	88

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	1163	VAL	1
1	A	1160	ALA	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	75/89 (84%)	70±1 (93±2%)	6±1 (7±2%)	22	68
2	B	9/15 (60%)	8±1 (91±8%)	1±1 (9±8%)	17	62
All	All	1680/2080 (81%)	1554 (92%)	126 (8%)	21	67

All 30 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	1235	ASN	20
1	A	1203	LEU	20
1	A	1190	VAL	13
1	A	1205	LYS	12
1	A	1233	LEU	8
1	A	1240	LEU	6
1	A	1184	GLU	5
1	A	1248	LEU	4
2	B	535	LYS	3
2	B	538	LEU	3
1	A	1206	LEU	3
2	B	533	LYS	3
2	B	531	PHE	2
1	A	1196	ASP	2
2	B	534	GLN	2
1	A	1215	ARG	2
1	A	1166	ASN	2
1	A	1198	ILE	2
2	B	539	LEU	2
1	A	1197	LEU	2
1	A	1159	LEU	1
1	A	1211	LYS	1
1	A	1247	THR	1
1	A	1204	GLU	1
2	B	532	PHE	1
1	A	1214	LYS	1
1	A	1169	LYS	1
1	A	1173	ARG	1

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Mol	Chain	Res	Type	Models (Total)
1	A	1212	TYR	1
1	A	1249	LYS	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](#)

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

The completeness of assignment taking into account all chemical shift lists is 76% for the well-defined parts and 70% for the entire structure.

### 7.1 Chemical shift list 1

File name: 2lsk\_cs.str

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	1118
Number of shifts mapped to atoms	1118
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	3

#### 7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	90	$-0.55 \pm 0.13$	Should be applied
$^{13}\text{C}_\beta$	88	$0.36 \pm 0.09$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}'$	84	$-0.21 \pm 0.08$	None needed ( $< 0.5$ ppm)
$^{15}\text{N}$	85	$0.34 \pm 0.35$	None needed ( $< 0.5$ ppm)

#### 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 76%, i.e. 879 atoms were assigned a chemical shift out of a possible 1164. 0 out of 24 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	386/445 (87%)	155/178 (87%)	154/178 (87%)	77/89 (87%)
Sidechain	467/638 (73%)	284/367 (77%)	175/247 (71%)	8/24 (33%)

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	<b>Total</b>	<b><sup>1</sup>H</b>	<b><sup>13</sup>C</b>	<b><sup>15</sup>N</b>
Aromatic	26/81 (32%)	13/43 (30%)	12/37 (32%)	1/1 (100%)
Overall	879/1164 (76%)	452/588 (77%)	341/462 (74%)	86/114 (75%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 70%, i.e. 982 atoms were assigned a chemical shift out of a possible 1398. 0 out of 26 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	<b>Total</b>	<b><sup>1</sup>H</b>	<b><sup>13</sup>C</b>	<b><sup>15</sup>N</b>
Backbone	433/551 (79%)	174/220 (79%)	174/222 (78%)	85/109 (78%)
Sidechain	521/754 (69%)	318/435 (73%)	193/292 (66%)	10/27 (37%)
Aromatic	28/93 (30%)	14/49 (29%)	12/42 (29%)	2/2 (100%)
Overall	982/1398 (70%)	506/704 (72%)	379/556 (68%)	97/138 (70%)

#### 7.1.4 Statistically unusual chemical shifts ⓘ

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	1182	PRO	CG	20.25	32.66 – 21.76	-6.4
1	A	1182	PRO	HD3	1.66	5.52 – 1.72	-5.1
1	A	1169	LYS	HG3	-0.06	2.76 – -0.04	-5.1

#### 7.1.5 Random Coil Index (RCI) plots ⓘ

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:

