



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 02:25 PM BST

PDB ID : 1E1G
Title : HUMAN PRION PROTEIN VARIANT M166V
Authors : Calzolari, L.; Lysek, D.A.; Guntert, P.; Von Schroetter, C.; Zahn, R.; Riek, R.; Wuthrich, K.
Deposited on : 2000-05-04

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

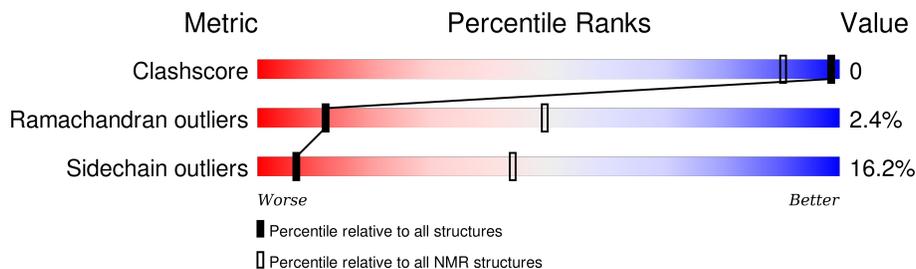
1 Overall quality at a glance i

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	104	

2 Ensemble composition and analysis

This entry contains 20 models. Model 12 is the overall representative, medoid model (most similar to other models).

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:125-A:228 (104)	0.65	12

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 4 clusters and 2 single-model clusters were found.

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

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 1690 atoms, of which 814 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called PRION PROTEIN.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	104	1690	544	814	153	171	8	0

There is a discrepancy between the modelled and reference sequences:

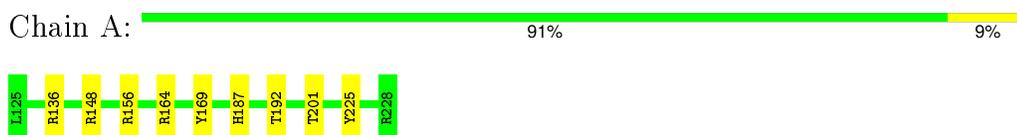
Chain	Residue	Modelled	Actual	Comment	Reference
A	166	VAL	MET	ENGINEERED	UNP P78446

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: PRION PROTEIN

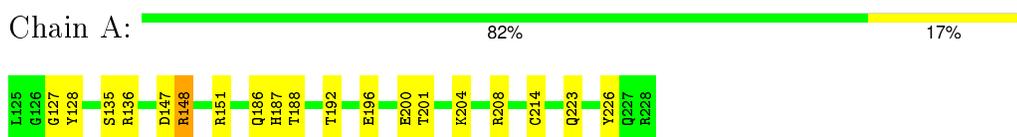


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: PRION PROTEIN



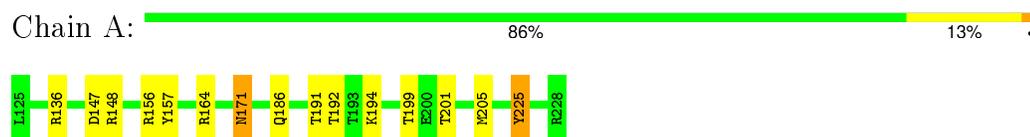
4.2.2 Score per residue for model 2

- Molecule 1: PRION PROTEIN



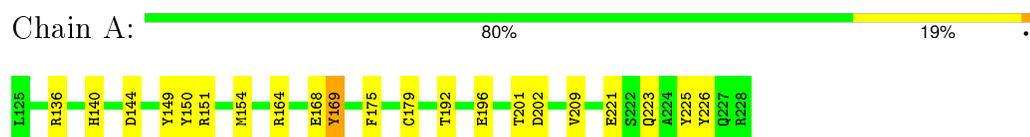
4.2.3 Score per residue for model 3

- Molecule 1: PRION PROTEIN



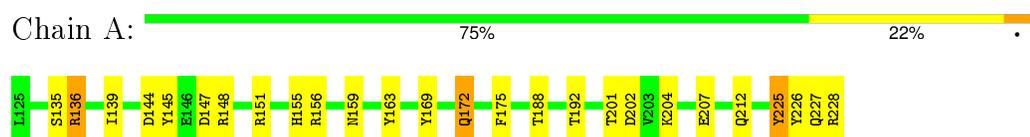
4.2.4 Score per residue for model 4

- Molecule 1: PRION PROTEIN



4.2.5 Score per residue for model 5

- Molecule 1: PRION PROTEIN



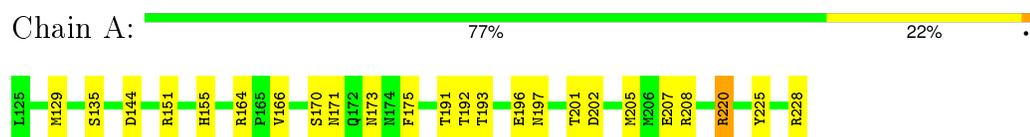
4.2.6 Score per residue for model 6

- Molecule 1: PRION PROTEIN



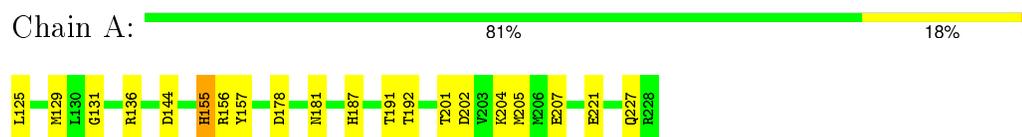
4.2.7 Score per residue for model 7

- Molecule 1: PRION PROTEIN



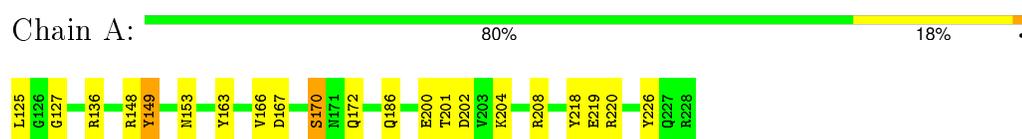
4.2.8 Score per residue for model 8

- Molecule 1: PRION PROTEIN



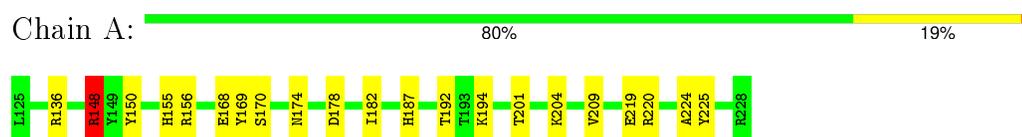
4.2.9 Score per residue for model 9

- Molecule 1: PRION PROTEIN



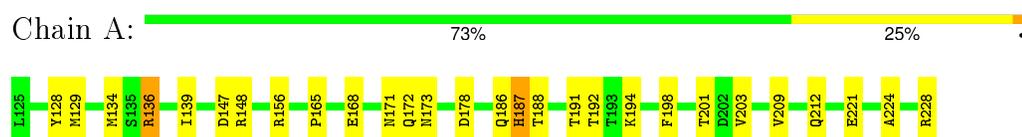
4.2.10 Score per residue for model 10

- Molecule 1: PRION PROTEIN



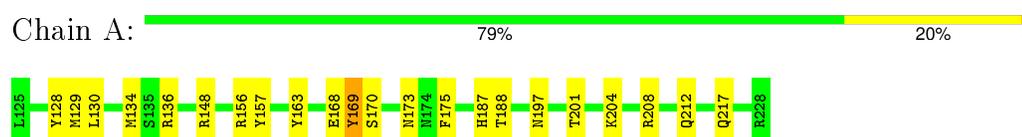
4.2.11 Score per residue for model 11

- Molecule 1: PRION PROTEIN



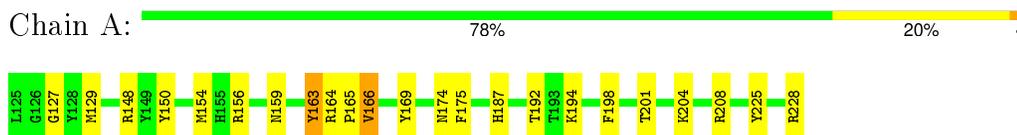
4.2.12 Score per residue for model 12 (medoid)

- Molecule 1: PRION PROTEIN



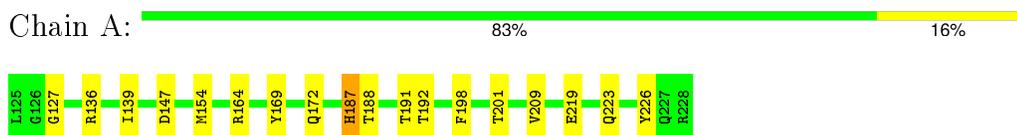
4.2.13 Score per residue for model 13

- Molecule 1: PRION PROTEIN



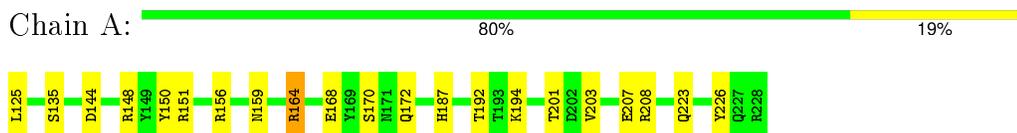
4.2.14 Score per residue for model 14

- Molecule 1: PRION PROTEIN



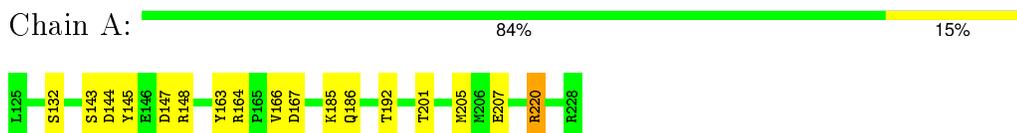
4.2.15 Score per residue for model 15

- Molecule 1: PRION PROTEIN



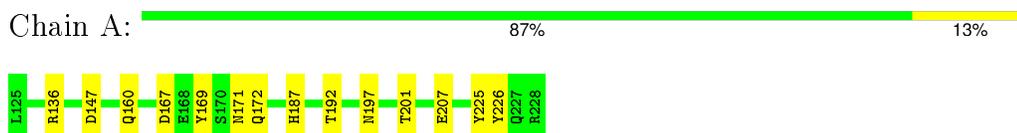
4.2.16 Score per residue for model 16

- Molecule 1: PRION PROTEIN



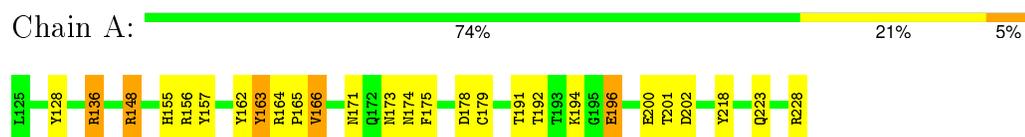
4.2.17 Score per residue for model 17

- Molecule 1: PRION PROTEIN



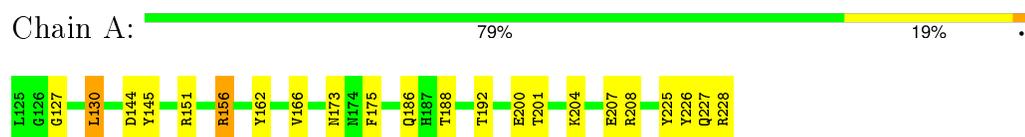
4.2.18 Score per residue for model 18

- Molecule 1: PRION PROTEIN



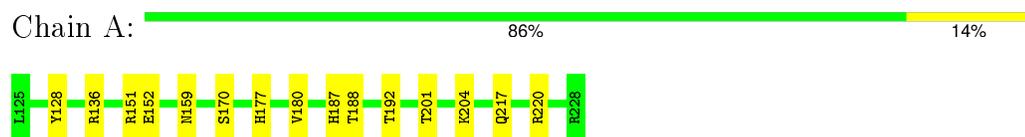
4.2.19 Score per residue for model 19

- Molecule 1: PRION PROTEIN



4.2.20 Score per residue for model 20

- Molecule 1: PRION PROTEIN



5 Refinement protocol and experimental data overview

The models were refined using the following method: *TORSION ANGLE DYNAMICS*.

Of the 20 calculated structures, 20 were deposited, based on the following criterion: ?

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

Software name	Classification	Version
OPALP	refinement	
DYANA	structure solution	

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

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	0.68±0.01	0±0/896 (0.0±0.0%)	1.06±0.04	1±1/1210 (0.1±0.1%)
All	All	0.68	0/17920 (0.0%)	1.06	26/24200 (0.1%)

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	2.2±1.9
All	All	0	45

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(°)	Ideal(°)	Models	
								Worst	Total
1	A	148	ARG	NE-CZ-NH2	-8.22	116.19	120.30	6	3
1	A	169	TYR	CB-CG-CD2	-6.82	116.91	121.00	13	1
1	A	164	ARG	NE-CZ-NH2	-6.68	116.96	120.30	7	3
1	A	149	TYR	CB-CG-CD2	-6.51	117.09	121.00	9	1
1	A	157	TYR	CB-CG-CD2	-6.50	117.10	121.00	3	3
1	A	156	ARG	NE-CZ-NH1	6.41	123.50	120.30	19	2
1	A	148	ARG	NE-CZ-NH1	6.03	123.31	120.30	10	1
1	A	220	ARG	NE-CZ-NH2	-5.68	117.46	120.30	16	2
1	A	180	VAL	CA-CB-CG1	5.64	119.36	110.90	20	1
1	A	136	ARG	CD-NE-CZ	5.62	131.47	123.60	11	1
1	A	136	ARG	NE-CZ-NH1	5.59	123.09	120.30	5	1
1	A	225	TYR	CB-CG-CD2	-5.36	117.78	121.00	5	1
1	A	150	TYR	CB-CG-CD1	-5.35	117.79	121.00	15	1
1	A	202	ASP	CB-CG-OD1	5.29	123.06	118.30	8	1
1	A	151	ARG	NE-CZ-NH2	-5.26	117.67	120.30	20	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	136	ARG	NE-CZ-NH2	-5.07	117.77	120.30	17	1
1	A	228	ARG	NE-CZ-NH2	-5.06	117.77	120.30	19	1
1	A	157	TYR	CB-CG-CD1	5.00	124.00	121.00	3	1

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	128	TYR	Sidechain	5
1	A	148	ARG	Sidechain	4
1	A	164	ARG	Sidechain	4
1	A	225	TYR	Sidechain	4
1	A	220	ARG	Sidechain	4
1	A	150	TYR	Sidechain	3
1	A	151	ARG	Sidechain	3
1	A	218	TYR	Sidechain	3
1	A	169	TYR	Sidechain	2
1	A	163	TYR	Sidechain	2
1	A	156	ARG	Sidechain	2
1	A	228	ARG	Sidechain	1
1	A	167	ASP	Peptide	1
1	A	136	ARG	Sidechain	1
1	A	166	VAL	Peptide	1
1	A	226	TYR	Sidechain	1
1	A	149	TYR	Sidechain	1
1	A	208	ARG	Sidechain	1
1	A	198	PHE	Sidechain	1
1	A	162	TYR	Sidechain	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	876	814	813	0±1
All	All	17520	16280	16260	10

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 0.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:139:ILE:HD11	1:A:209:VAL:HG22	0.53	1.80	14	1
1:A:130:LEU:HD23	1:A:162:TYR:CE1	0.51	2.39	19	1
1:A:191:THR:HG23	1:A:196:GLU:HG2	0.51	1.83	18	1
1:A:187:HIS:CD2	1:A:198:PHE:CD2	0.47	3.03	11	2
1:A:169:TYR:CD2	1:A:175:PHE:CE1	0.46	3.04	6	1
1:A:139:ILE:HD11	1:A:212:GLN:HG2	0.43	1.88	5	1
1:A:139:ILE:HD11	1:A:212:GLN:CG	0.43	2.43	11	2
1:A:196:GLU:HG2	1:A:197:ASN:H	0.40	1.76	7	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	102/104 (98%)	89±3 (87±3%)	11±3 (11±3%)	2±1 (2±1%)	12	49
All	All	2040/2080 (98%)	1771 (87%)	220 (11%)	49 (2%)	12	49

All 19 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	170	SER	6
1	A	127	GLY	6
1	A	172	GLN	6
1	A	166	VAL	5
1	A	165	PRO	4
1	A	171	ASN	3
1	A	169	TYR	3
1	A	198	PHE	2
1	A	135	SER	2
1	A	155	HIS	2
1	A	224	ALA	2

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Mol	Chain	Res	Type	Models (Total)
1	A	128	TYR	1
1	A	126	GLY	1
1	A	219	GLU	1
1	A	156	ARG	1
1	A	200	GLU	1
1	A	131	GLY	1
1	A	167	ASP	1
1	A	157	TYR	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	97/97 (100%)	81±3 (84±3%)	16±3 (16±3%)	7 44
All	All	1940/1940 (100%)	1625 (84%)	315 (16%)	7 44

All 70 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	201	THR	20
1	A	192	THR	18
1	A	136	ARG	12
1	A	204	LYS	10
1	A	187	HIS	10
1	A	226	TYR	9
1	A	147	ASP	8
1	A	207	GLU	8
1	A	208	ARG	8
1	A	188	THR	8
1	A	148	ARG	8
1	A	194	LYS	7
1	A	175	PHE	7
1	A	225	TYR	7
1	A	144	ASP	7
1	A	156	ARG	7
1	A	186	GLN	7

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Mol	Chain	Res	Type	Models (Total)
1	A	173	ASN	7
1	A	163	TYR	7
1	A	168	GLU	6
1	A	191	THR	6
1	A	202	ASP	6
1	A	155	HIS	5
1	A	223	GLN	5
1	A	129	MET	5
1	A	164	ARG	5
1	A	151	ARG	4
1	A	205	MET	4
1	A	228	ARG	4
1	A	159	ASN	4
1	A	221	GLU	4
1	A	178	ASP	4
1	A	169	TYR	4
1	A	196	GLU	4
1	A	171	ASN	3
1	A	166	VAL	3
1	A	154	MET	3
1	A	197	ASN	3
1	A	217	GLN	3
1	A	200	GLU	3
1	A	145	TYR	3
1	A	134	MET	3
1	A	227	GLN	3
1	A	174	ASN	3
1	A	135	SER	3
1	A	125	LEU	3
1	A	209	VAL	3
1	A	219	GLU	2
1	A	185	LYS	2
1	A	179	CYS	2
1	A	220	ARG	2
1	A	167	ASP	2
1	A	203	VAL	2
1	A	181	ASN	2
1	A	130	LEU	2
1	A	153	ASN	1
1	A	170	SER	1
1	A	193	THR	1
1	A	172	GLN	1

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Mol	Chain	Res	Type	Models (Total)
1	A	152	GLU	1
1	A	199	THR	1
1	A	177	HIS	1
1	A	182	ILE	1
1	A	160	GLN	1
1	A	214	CYS	1
1	A	149	TYR	1
1	A	212	GLN	1
1	A	132	SER	1
1	A	140	HIS	1
1	A	143	SER	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

No chemical shift data were provided