

Full wwPDB NMR Structure Validation Report (i)

Nov 2, 2022 – 09:53 AM EDT

PDB ID : 8DI2 BMRB ID : 31030

Title : Site 2 insulin receptor binding peptide IM459N21

Authors: Lawrence, M.C.; Hu, H.; Martinez, F.J.; Espinosa, J.F.

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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
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

Mol Probity : 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

RCI : v 1n 11 5 13 A (Berjanski et al., 2005)

PANAV : Wang et al. (2010)

ShiftChecker : 2.31.2

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

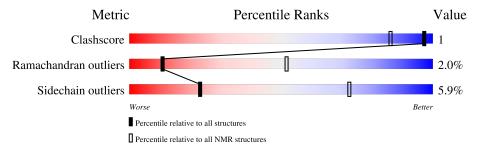
Validation Pipeline (wwPDB-VP) : 2.31.2

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 is 81%.

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	NMR archive		
Metric	$(\# \mathrm{Entries})$	$(\# \mathrm{Entries})$		
Clashscore	158937	12864		
Ramachandran outliers	154571	11451		
Sidechain outliers	154315	11428		

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	22	55%	27%	5%	14%			



2 Ensemble composition and analysis (i)

This entry contains 16 models. Model 10 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: fewest violations.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues							
Well-defined core	Well-defined core Residue range (total) Backbone RMSD (Å) Medoid model						
1	A:2-A:7, A:9-A:21 (19)	0.95	10				

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. No single-model clusters were found.

Cluster number	Models
1	1, 3, 6, 7, 8, 9, 10, 12, 13, 14, 15, 16
2	2, 4
3	5, 11



3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 368 atoms, of which 184 are hydrogens and 0 are deuteriums.

• Molecule 1 is a protein called Site 2 binding peptide IM459N21.

Mol	Chain	Residues	Atoms					Trace	
1	Λ	29	Total	С	Н	N	О	S	0
1	A	22	368	120	184	29	33	2	U



4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide 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: Site 2 binding peptide IM459N21



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: Site 2 binding peptide IM459N21



4.2.2 Score per residue for model 2

• Molecule 1: Site 2 binding peptide IM459N21





4.2.3 Score per residue for model 3

• Molecule 1: Site 2 binding peptide IM459N21





4.2.4 Score per residue for model 4

• Molecule 1: Site 2 binding peptide IM459N21

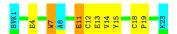




4.2.5 Score per residue for model 5

• Molecule 1: Site 2 binding peptide IM459N21

Chain A: 45% 32% 9% 14%



4.2.6 Score per residue for model 6

• Molecule 1: Site 2 binding peptide IM459N21

Chain A: 68% 5% 14% 14%



4.2.7 Score per residue for model 7

• Molecule 1: Site 2 binding peptide IM459N21

Chain A: 55% 23% 9% 14%





4.2.8 Score per residue for model 8

• Molecule 1: Site 2 binding peptide IM459N21

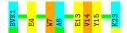




4.2.9 Score per residue for model 9

• Molecule 1: Site 2 binding peptide IM459N21

Chain A: 64% 14% 9% 14%



4.2.10 Score per residue for model 10 (medoid)

• Molecule 1: Site 2 binding peptide IM459N21

Chain A: 55% 27% 5% 14%



4.2.11 Score per residue for model 11

• Molecule 1: Site 2 binding peptide IM459N21

Chain A: 55% 23% 9% 14%



4.2.12 Score per residue for model 12

• Molecule 1: Site 2 binding peptide IM459N21

Chain A: 59% 23% 5% 14%





4.2.13 Score per residue for model 13

• Molecule 1: Site 2 binding peptide IM459N21



4.2.14 Score per residue for model 14

• Molecule 1: Site 2 binding peptide IM459N21



4.2.15 Score per residue for model 15

• Molecule 1: Site 2 binding peptide IM459N21



4.2.16 Score per residue for model 16

• Molecule 1: Site 2 binding peptide IM459N21





Refinement protocol and experimental data overview (i) 5



The models were refined using the following method: molecular dynamics.

Of the 16 calculated structures, 16 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
GROMACS	refinement	5.0.7
GROMACS	structure calculation	5.0.7

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)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	247
Number of shifts mapped to atoms	246
Number of unparsed shifts	1
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	81%



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: AIB, LYN, BVK

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).

Mal	Chain	В	Sond lengths	Bond angles		
MIOI	Chain	RMSZ	#Z>5	RMSZ	#Z>5	
1	A	0.58 ± 0.02	$0\pm0/160~(~0.0\pm~0.0\%)$	1.96 ± 0.10	$4\pm 2/213$ ($1.9\pm~0.8\%$)	
All	All	0.58	0/2560~(~0.0%)	1.97	66/3408 (1.9%)	

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	$4.7{\pm}1.6$
All	All	0	75

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.

Mal	ol Chain H		Trino	Trung Atoms	Z	Observed(0)	Ideal(0)	Mod	dels
Mol	Chain	Res	Type	Atoms	L	$Observed(^o)$	$\operatorname{Ideal}(^{o})$	Worst	Total
1	A	15	TYR	CB-CG-CD1	-12.55	113.47	121.00	16	5
1	A	15	TYR	CG-CD2-CE2	-9.96	113.33	121.30	15	2
1	A	11	GLU	OE1-CD-OE2	-8.83	112.70	123.30	11	3
1	A	7	TRP	CB-CA-C	7.99	126.37	110.40	8	5
1	A	7	TRP	CA-CB-CG	7.82	128.56	113.70	12	6
1	A	7	TRP	NE1-CE2-CD2	-7.81	99.49	107.30	9	2
1	A	3	LEU	CB-CG-CD2	7.73	124.14	111.00	2	1
1	A	15	TYR	CB-CG-CD2	-7.45	116.53	121.00	3	3
1	A	7	TRP	CG-CD2-CE3	-7.22	127.40	133.90	1	1
1	A	15	TYR	CD1-CG-CD2	7.10	125.71	117.90	15	2
1	A	13	GLU	OE1-CD-OE2	7.00	131.70	123.30	11	2
1	A	7	TRP	CE2-CD2-CG	6.98	112.89	107.30	9	3

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Mol	Mol Chain R		Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$	Models	
MIOI	Chain	Res	туре	Atoms	L	Observed()	ideai()	Worst	Total
1	A	12	CYS	CA-CB-SG	6.92	126.45	114.00	8	2
1	A	7	TRP	CD1-NE1-CE2	6.73	115.06	109.00	9	3
1	A	14	VAL	CA-CB-CG2	6.71	120.96	110.90	3	3
1	A	4	GLU	O-C-N	-6.57	112.19	122.70	4	1
1	A	15	TYR	CD1-CE1-CZ	-6.50	113.95	119.80	15	1
1	A	19	PRO	N-CA-CB	6.25	110.80	103.30	5	2
1	A	15	TYR	CG-CD1-CE1	-6.17	116.36	121.30	16	1
1	A	7	TRP	CZ3-CH2-CZ2	-6.02	114.38	121.60	10	2
1	A	14	VAL	CG1-CB-CG2	-6.01	101.28	110.90	9	1
1	A	7	TRP	NE1-CE2-CZ2	6.00	137.00	130.40	1	3
1	A	19	PRO	N-CD-CG	5.94	112.11	103.20	2	1
1	A	7	TRP	CE2-CD2-CE3	-5.93	111.58	118.70	9	1
1	A	7	TRP	CB-CG-CD2	5.81	134.15	126.60	14	1
1	A	15	TYR	CA-CB-CG	5.77	124.36	113.40	10	3
1	A	7	TRP	CD2-CE3-CZ3	5.45	125.88	118.80	9	1
1	A	15	TYR	CB-CA-C	5.35	121.09	110.40	7	1
1	A	14	VAL	CA-CB-CG1	5.23	118.74	110.90	8	1
1	A	7	TRP	CH2-CZ2-CE2	5.20	122.60	117.40	12	1
1	A	3	LEU	C-N-CA	5.12	134.51	121.70	11	1
1	A	17	LYS	N-CA-CB	-5.08	101.45	110.60	2	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	4	GLU	Peptide	14
1	A	19	PRO	Peptide	13
1	A	13	GLU	Peptide	11
1	A	12	CYS	Peptide	10
1	A	14	VAL	Peptide, Mainchain	8
1	A	15	TYR	Sidechain	7
1	A	7	TRP	Mainchain,Peptide	4
1	A	3	LEU	Peptide	2
1	A	11	GLU	Peptide	2
1	A	2	SER	Peptide	1
1	A	16	GLY	Peptide	1
1	A	6	GLU	Sidechain	1



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.

\mathbf{N}	/Iol	Chain	Non-H	H(model)	H(added)	Clashes
	1	A	157	151	151	0±0
A	All	All	2512	2416	2416	3

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 1.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models		
Atom-1	Atom-2	Clash(A)	Distance(A)	Worst	Total	
1:A:18:CYS:SG	1:A:21:LYS:HB2	0.47	2.50	2	1	
1:A:4:GLU:HA	1:A:7:TRP:NE1	0.41	2.30	15	1	
1:A:4:GLU:HA	1:A:7:TRP:CE2	0.41	2.50	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	19/22 (86%)	16±1 (83±6%)	3±1 (15±6%)	0±0 (2±3%)	11	52	
All	All	304/352 (86%)	252 (83%)	46 (15%)	6 (2%)	11	52	

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	14	VAL	4
1	A	16	GLY	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	18/18 (100%)	17±1 (94±5%)	1±1 (6±5%)	23 72		
All	All	288/288 (100%)	271 (94%)	17 (6%)	23 72		

All 4 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	7	TRP	9
1	A	12	CYS	4
1	A	18	CYS	3
1	A	6	GLU	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)

2 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	Tuno	Chain	Pag	Link		gths	
	IVIOI	туре	Chain	nes	Lilik	Counts	RMSZ	#Z>2
Ī	1	LYN	A	23	1	9,9,9	0.34 ± 0.05	0±0 (0±0%)
Ī	1	AIB	A	8	1	1,5,6	1.09 ± 0.10	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.

Mal	Trunc	Chain	Dag	Timle		Bond angles Counts RMSZ #Z>2		
MIOI	туре	Chain	nes	Lilik	Counts	RMSZ	$\#Z{>}2$	
1	LYN	A	23	1	9,10,10	1.20 ± 0.44	1±1 (8±10%)	
1	AIB	A	8	1	2,7,9	1.10 ± 0.37	0±0 (3±12%)	

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
1	AIB	A	8	1	-	$0\pm0,2,3,6$	-
1	LYN	A	23	1	-	$0\pm0,9,9,9$	-

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	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$	Models	
WIOI	Chain	nes	Type	Atoms		Observed()	Ideal()	Worst	Total
1	A	23	LYN	CA-C-NT	4.99	125.24	116.68	4	4
1	A	23	LYN	O-C-NT	4.93	114.43	123.00	4	2
1	A	23	LYN	O-C-CA	2.91	116.03	120.30	7	4
1	A	8	AIB	CB2-CA-C	2.48	118.38	108.99	5	1
1	A	23	LYN	CB-CA-N	2.39	103.90	110.17	1	1
1	A	23	LYN	CG-CB-CA	2.09	120.08	113.35	6	1

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates (i)

There are no monosaccharides 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)

The completeness of assignment taking into account all chemical shift lists is 81% for the well-defined parts and 81% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: assigned_chem_shift_list

7.1.1 Bookkeeping (i)

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	247
Number of shifts mapped to atoms	246
Number of unparsed shifts	1
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

The following errors were found when reading this chemical shift list.

• Chemical shift has been reported more than once. The only occurrence is reported below.

Shift ID	Chain	Pos	Type	Atom	Shift Data Value Uncertainty Ambiguity		
	Chain	nes	туре	Atom	Value	Uncertainty	Ambiguity
247	A	3	LEU	HB2	1.168	0.001	2

7.1.2 Chemical shift referencing (i)

No chemical shift referencing corrections were calculated (not enough data).

7.1.3 Completeness of resonance assignments (i)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 81%, i.e. 199 atoms were assigned a chemical shift out of a possible 245. 0 out of 2 assigned methyl groups (LEU and VAL) were assigned stereospecifically.



	Total	$^{1}\mathrm{H}$	$^{13}\mathbf{C}$	$^{15}{ m N}$
Backbone	55/91 (60%)	36/36 (100%)	19/38 (50%)	0/17 (0%)
Sidechain	125/134 (93%)	81/81 (100%)	44/49 (90%)	0/4 (0%)
Aromatic	19/20 (95%)	10/10 (100%)	9/9 (100%)	0/1 (0%)
Overall	199/245 (81%)	127/127 (100%)	72/96 (75%)	0/22~(0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 81%, i.e. 199 atoms were assigned a chemical shift out of a possible 245. 0 out of 2 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^{1}\mathrm{H}$	$^{13}\mathbf{C}$	$^{15}{ m N}$
Backbone	55/91 (60%)	36/36~(100%)	19/38 (50%)	0/17~(0%)
Sidechain	125/134 (93%)	81/81 (100%)	44/49 (90%)	0/4 (0%)
Aromatic	19/20 (95%)	10/10 (100%)	9/9 (100%)	0/1 (0%)
Overall	199/245 (81%)	127/127 (100%)	72/96 (75%)	0/22~(0%)

7.1.4 Statistically unusual chemical shifts (i)

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	9	LYS	CG	36.46	30.67 - 19.17	10.0

7.1.5 Random Coil Index (RCI) plots (i)

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:



