



Full wwPDB NMR Structure Validation Report ⓘ

Feb 20, 2022 – 01:36 AM EST

PDB ID : 1U5S
Title : NMR structure of the complex between Nck-2 SH3 domain and PINCH-1 LIM4 domain
Authors : Vaynberg, J.; Fukuda, T.; Vinogradova, O.; Velyvis, A.; Ng, L.; Wu, C.; Qin, J.
Deposited on : 2004-07-28

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 ⓘ symbol.

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

MolProbity : 4.02b-467
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 : 2.26
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.26

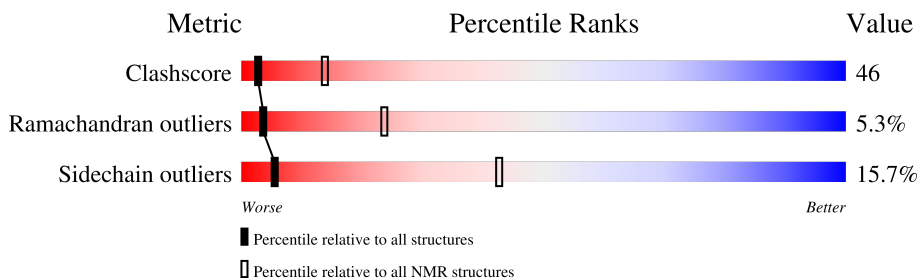
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment 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	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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	71	 39% 42% 17% .
2	B	66	 35% 61% . .

2 Ensemble composition and analysis

This entry contains 18 models. Model 3 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:1-A:71 (71)	0.00	3
2	B:72-B:137 (66)	0.00	7

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

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

The models can be grouped into 3 clusters and 1 single-model cluster was found.

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

3 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2150 atoms, of which 1057 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Cytoplasmic protein NCK2.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
1	A	71	1121	361	555	95	108	2	0

- Molecule 2 is a protein called PINCH protein.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
2	B	66	1027	331	502	100	87	7	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	72	GLY	-	cloning artifact	UNP P48059
B	73	SER	-	cloning artifact	UNP P48059

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

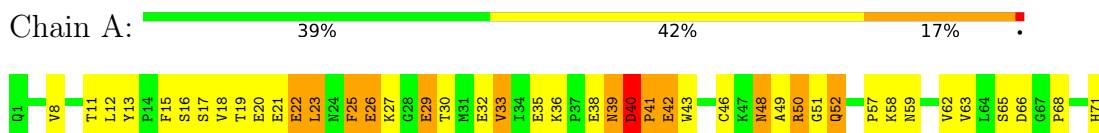
Mol	Chain	Residues	Atoms	
			Total	Zn
3	B	2	2	2

4 Residue-property plots

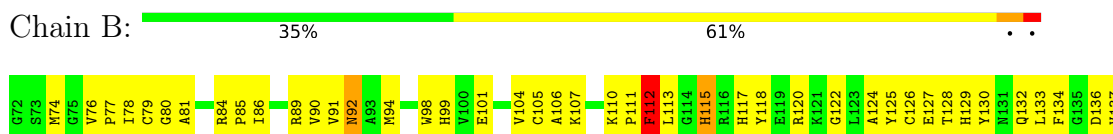
4.1 Average score per residue in the NMR ensemble

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: Cytoplasmic protein NCK2



- Molecule 2: PINCH protein

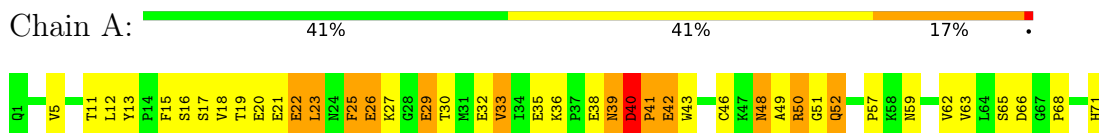


4.2 Scores per residue for each member of the ensemble

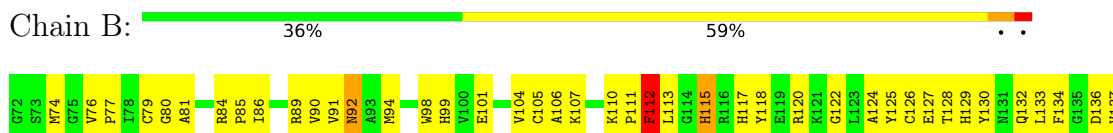
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: Cytoplasmic protein NCK2

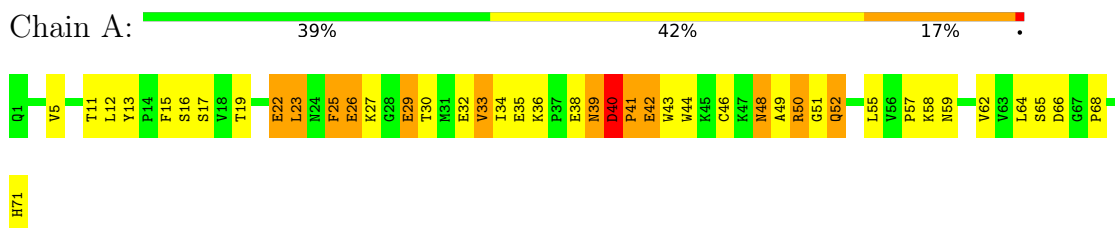


- Molecule 2: PINCH protein

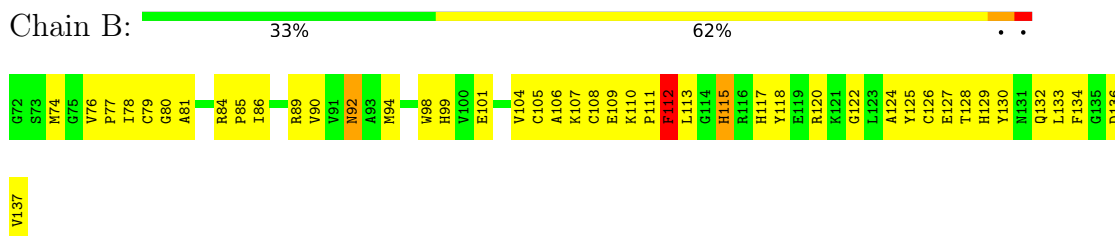


4.2.2 Score per residue for model 2

- Molecule 1: Cytoplasmic protein NCK2

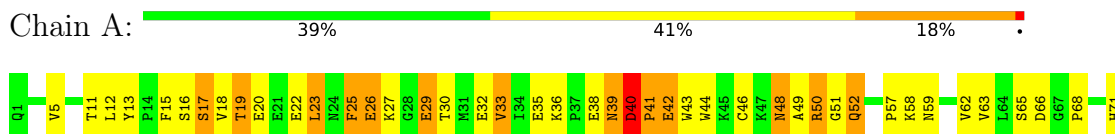


- Molecule 2: PINCH protein

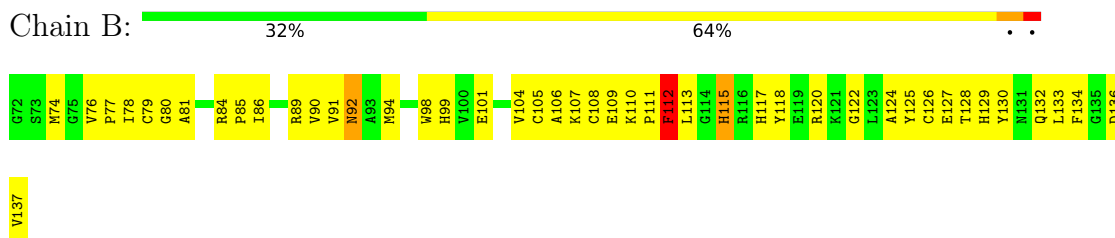


4.2.3 Score per residue for model 3 (medoid)

- Molecule 1: Cytoplasmic protein NCK2



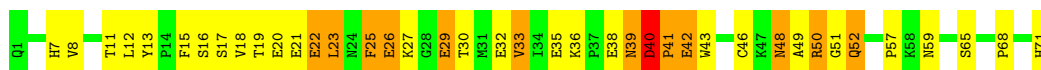
- Molecule 2: PINCH protein



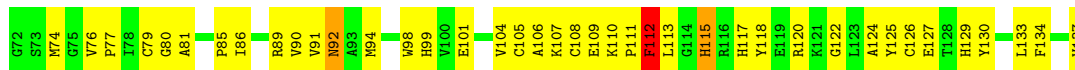
4.2.4 Score per residue for model 4

- Molecule 1: Cytoplasmic protein NCK2



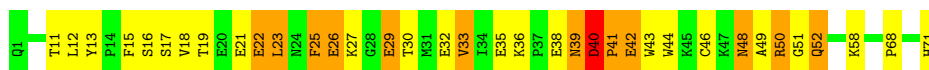


- Molecule 2: PINCH protein

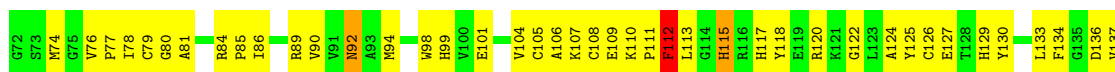


4.2.5 Score per residue for model 5

- Molecule 1: Cytoplasmic protein NCK2

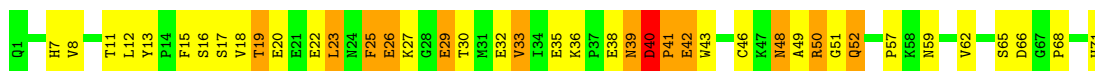


- Molecule 2: PINCH protein



4.2.6 Score per residue for model 6

- Molecule 1: Cytoplasmic protein NCK2



- Molecule 2: PINCH protein



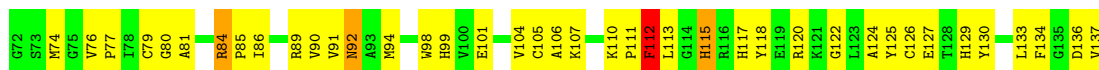
4.2.7 Score per residue for model 7

- Molecule 1: Cytoplasmic protein NCK2





- Molecule 2: PINCH protein

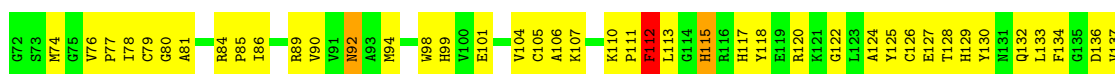


4.2.8 Score per residue for model 8

- Molecule 1: Cytoplasmic protein NCK2

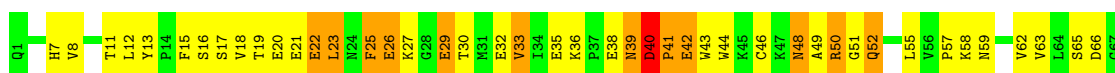


- Molecule 2: PINCH protein

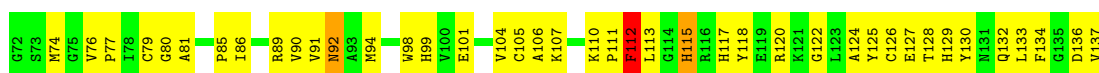
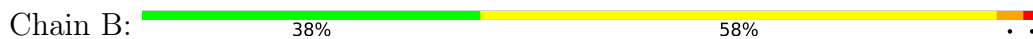


4.2.9 Score per residue for model 9

- Molecule 1: Cytoplasmic protein NCK2

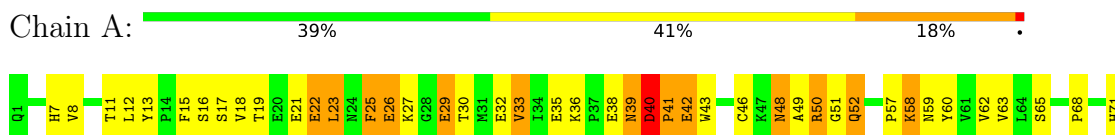


- Molecule 2: PINCH protein

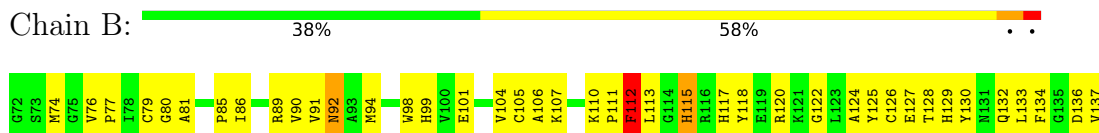


4.2.10 Score per residue for model 10

- Molecule 1: Cytoplasmic protein NCK2

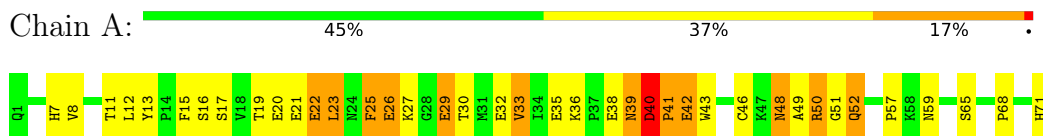


- Molecule 2: PINCH protein

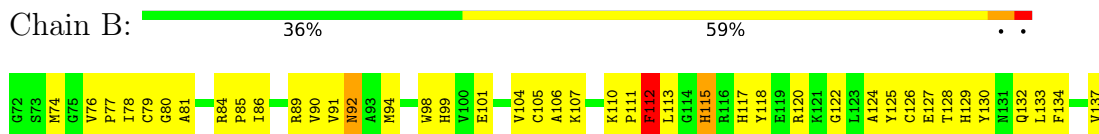


4.2.11 Score per residue for model 11

- Molecule 1: Cytoplasmic protein NCK2

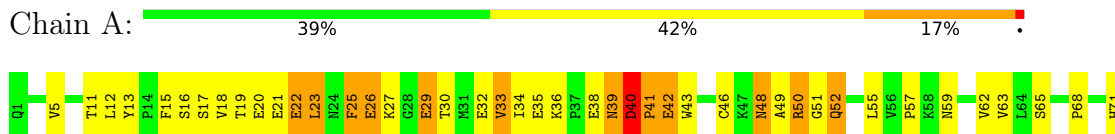


- Molecule 2: PINCH protein

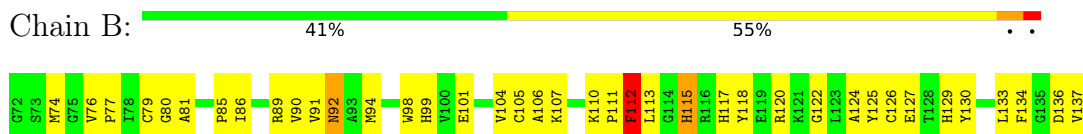


4.2.12 Score per residue for model 12

- Molecule 1: Cytoplasmic protein NCK2

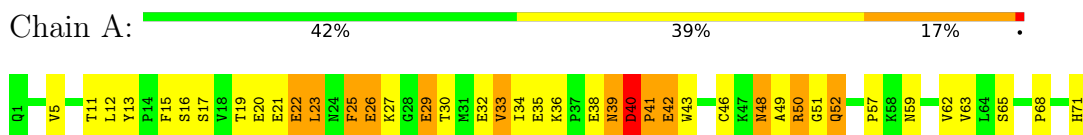


- Molecule 2: PINCH protein

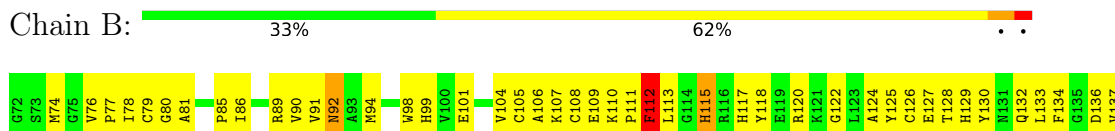


4.2.13 Score per residue for model 13

- Molecule 1: Cytoplasmic protein NCK2

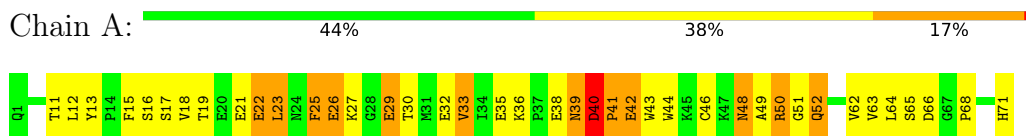


- Molecule 2: PINCH protein

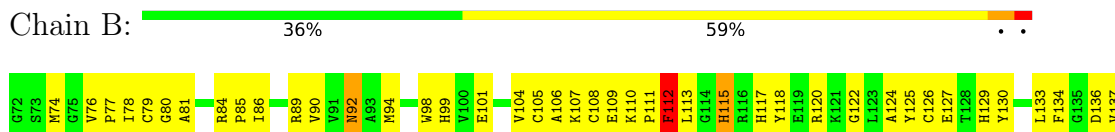


4.2.14 Score per residue for model 14

- Molecule 1: Cytoplasmic protein NCK2

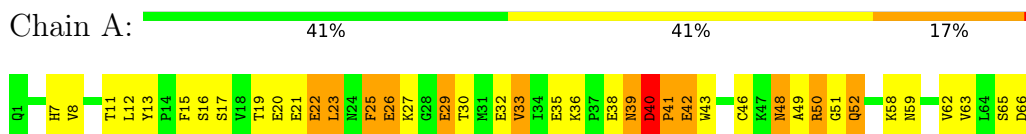


- Molecule 2: PINCH protein

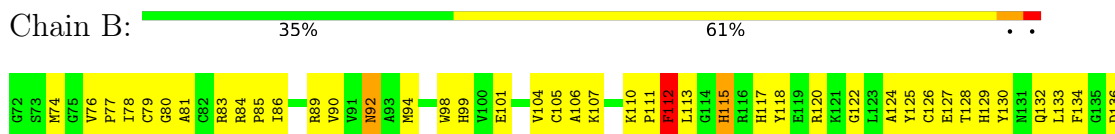


4.2.15 Score per residue for model 15

- Molecule 1: Cytoplasmic protein NCK2



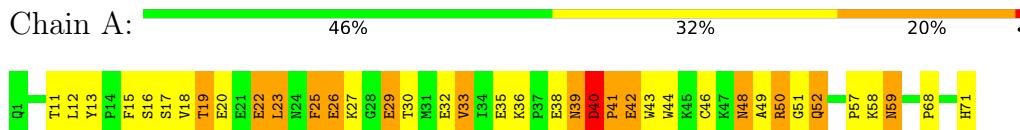
- Molecule 2: PINCH protein



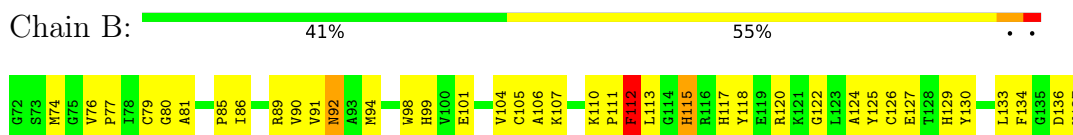
Y137

4.2.16 Score per residue for model 16

- Molecule 1: Cytoplasmic protein NCK2

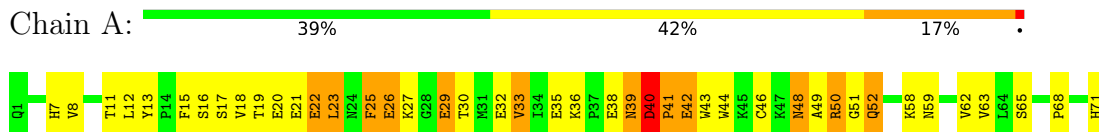


- Molecule 2: PINCH protein

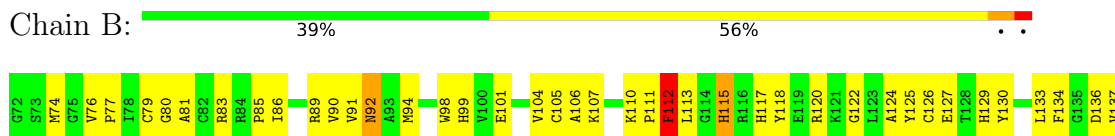


4.2.17 Score per residue for model 17

- Molecule 1: Cytoplasmic protein NCK2

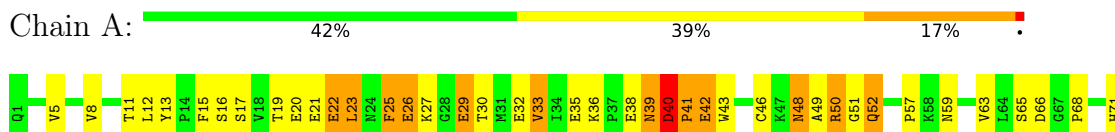


- Molecule 2: PINCH protein



4.2.18 Score per residue for model 18

- Molecule 1: Cytoplasmic protein NCK2



- Molecule 2: PINCH protein



G72	S73	M74	G75	V76	P77	L78	C79	G80	A81	C82	P85	I86	R89	Y90	Y91	M92	A93	Y94	W98	H99	V100	E101	V104	C105	A106	K107	K110	P111	F112	L113	G114	H115	R116	H117	Y118	E119	R120	K121	G122	L123	A124	Y125	C126	E127	T128	H129	Y130	M131	Q132	L133	F134	G135	D136	V137
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5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing, molecular dynamics, rigid body/torsion angle dynamics*.

Of the 50 calculated structures, 18 were deposited, based on the following criterion: *structures with the most representative side chains orientations on the complex interface*.

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

Software name	Classification	Version
X-PLOR	structure solution	3.1
Xplor-NIH	structure solution	
Xplor-NIH	refinement	

No chemical shift data was provided.

6 Model quality i

6.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

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	1.15±0.00	0±0/580 (0.0± 0.0%)	0.94±0.00	0±0/789 (0.0± 0.0%)
2	B	2.80±0.00	3±0/540 (0.6± 0.0%)	2.51±0.00	5±0/725 (0.7± 0.0%)
All	All	2.11	54/20160 (0.3%)	1.87	90/27252 (0.3%)

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
2	B	0.0±0.0	1.0±0.0
All	All	0	18

All unique bond 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(Å)	Ideal(Å)	Models	
								Worst	Total
2	B	137	VAL	C-O	-39.80	0.47	1.23	9	18
2	B	112	PHE	CB-CG	36.08	2.12	1.51	10	18
2	B	112	PHE	CA-CB	27.75	2.15	1.53	5	18

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
2	B	112	PHE	CB-CG-CD2	-40.16	92.69	120.80	7	18
2	B	112	PHE	CB-CG-CD1	35.48	145.63	120.80	12	18
2	B	137	VAL	CA-C-O	-23.48	70.79	120.10	6	18
2	B	112	PHE	CA-CB-CG	-17.86	71.04	113.90	5	18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	112	PHE	N-CA-CB	13.57	135.02	110.60	3	18

There are no chirality outliers.

All unique planar outliers are listed below.

Mol	Chain	Res	Type	Group	Models (Total)
2	B	112	PHE	Sidechain	18

6.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	566	555	555	47±3
2	B	525	502	501	52±2
All	All	19674	19026	19008	1769

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 46.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:112:PHE:CB	2:B:112:PHE:CG	1.33	2.12	16	18
2:B:112:PHE:CB	2:B:112:PHE:CA	1.25	2.15	2	18
2:B:112:PHE:CG	2:B:112:PHE:CA	0.97	2.48	7	18
1:A:19:THR:HG22	1:A:20:GLU:OE1	0.91	1.65	6	2
2:B:112:PHE:CB	2:B:112:PHE:CD2	0.86	2.59	7	18
2:B:112:PHE:HB3	2:B:115:HIS:HB2	0.79	1.55	1	18
2:B:112:PHE:HB2	2:B:115:HIS:O	0.77	1.80	5	18
1:A:48:ASN:N	1:A:52:GLN:O	0.77	2.18	2	18
2:B:79:CYS:SG	2:B:81:ALA:HB3	0.76	2.21	7	18
1:A:19:THR:HG22	1:A:20:GLU:H	0.73	1.42	13	11
1:A:18:VAL:O	1:A:18:VAL:HG23	0.72	1.84	5	11
2:B:104:VAL:HG12	2:B:111:PRO:CA	0.71	2.16	15	18
2:B:118:TYR:CD1	2:B:118:TYR:N	0.70	2.59	15	18
1:A:48:ASN:ND2	1:A:52:GLN:H	0.70	1.85	1	18

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:22:GLU:N	1:A:22:GLU:OE1	0.67	2.28	1	18
1:A:38:GLU:O	1:A:40:ASP:N	0.67	2.28	10	18
1:A:48:ASN:ND2	1:A:51:GLY:N	0.67	2.43	4	18
1:A:59:ASN:OD1	1:A:59:ASN:N	0.66	2.28	16	1
1:A:48:ASN:ND2	1:A:48:ASN:O	0.65	2.29	1	18
1:A:41:PRO:O	1:A:42:GLU:HB2	0.64	1.92	3	12
1:A:48:ASN:ND2	1:A:52:GLN:N	0.64	2.46	9	18
2:B:94:MET:SD	2:B:122:GLY:O	0.64	2.56	11	18
2:B:112:PHE:CG	2:B:112:PHE:N	0.61	2.69	13	18
2:B:89:ARG:HH11	2:B:89:ARG:CG	0.60	2.10	16	17
1:A:59:ASN:OD1	2:B:83:ARG:NH2	0.59	2.35	15	2
1:A:40:ASP:N	1:A:41:PRO:CD	0.59	2.66	13	18
2:B:89:ARG:CG	2:B:89:ARG:NH1	0.59	2.66	12	18
1:A:48:ASN:HD22	1:A:48:ASN:C	0.59	2.01	1	18
1:A:48:ASN:O	1:A:51:GLY:N	0.58	2.35	4	18
2:B:89:ARG:CG	2:B:89:ARG:HH11	0.58	2.10	10	1
1:A:41:PRO:O	1:A:42:GLU:CB	0.58	2.52	6	18
2:B:112:PHE:CB	2:B:115:HIS:O	0.58	2.52	12	18
2:B:104:VAL:HG12	2:B:111:PRO:HA	0.57	1.77	3	18
2:B:117:HIS:C	2:B:118:TYR:CD1	0.57	2.78	6	18
1:A:42:GLU:OE1	1:A:58:LYS:NZ	0.57	2.38	10	2
2:B:112:PHE:HB3	2:B:115:HIS:C	0.56	2.21	4	18
1:A:62:VAL:HG11	2:B:78:ILE:HD13	0.56	1.77	6	2
2:B:112:PHE:CB	2:B:115:HIS:C	0.56	2.74	10	18
1:A:20:GLU:OE1	1:A:20:GLU:N	0.56	2.38	6	2
1:A:43:TRP:N	1:A:43:TRP:CD1	0.56	2.72	17	18
1:A:41:PRO:O	1:A:42:GLU:HB3	0.55	2.01	12	6
1:A:66:ASP:OD2	2:B:84:ARG:NH1	0.55	2.40	1	1
2:B:79:CYS:SG	2:B:81:ALA:CB	0.55	2.95	5	18
2:B:133:LEU:O	2:B:134:PHE:CD2	0.54	2.61	16	18
1:A:8:VAL:N	1:A:65:SER:OG	0.54	2.40	10	8
2:B:118:TYR:O	2:B:125:TYR:N	0.54	2.41	8	18
2:B:92:ASN:HD22	2:B:92:ASN:N	0.54	2.02	13	18
2:B:118:TYR:CZ	2:B:127:GLU:HB2	0.53	2.39	7	18
1:A:18:VAL:O	1:A:18:VAL:CG2	0.53	2.56	5	5
2:B:106:ALA:N	2:B:124:ALA:O	0.53	2.41	13	18
2:B:112:PHE:HB3	2:B:115:HIS:CB	0.53	2.33	8	18
1:A:40:ASP:N	1:A:41:PRO:HD2	0.53	2.19	2	18
1:A:26:GLU:CD	1:A:26:GLU:H	0.53	2.07	12	10
2:B:76:VAL:HG11	2:B:85:PRO:HB3	0.53	1.81	5	18
1:A:26:GLU:H	1:A:26:GLU:CD	0.52	2.07	6	8

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:65:SER:O	1:A:66:ASP:OD1	0.52	2.28	9	3
2:B:80:GLY:C	2:B:98:TRP:CZ3	0.52	2.83	5	18
2:B:111:PRO:HB2	2:B:113:LEU:CD1	0.52	2.35	16	18
1:A:35:GLU:CA	1:A:36:LYS:HZ3	0.52	2.17	15	16
1:A:41:PRO:O	1:A:42:GLU:HG2	0.51	2.06	1	2
2:B:112:PHE:O	2:B:113:LEU:C	0.51	2.49	18	18
1:A:41:PRO:O	1:A:42:GLU:CG	0.51	2.59	14	8
1:A:48:ASN:HD22	1:A:51:GLY:N	0.51	2.04	1	18
2:B:125:TYR:CE1	2:B:133:LEU:HD22	0.51	2.41	1	18
1:A:46:CYS:O	1:A:46:CYS:SG	0.50	2.70	2	18
2:B:86:ILE:HG21	2:B:90:VAL:HG22	0.50	1.83	10	18
2:B:84:ARG:NH2	2:B:99:HIS:CE1	0.50	2.79	8	1
1:A:11:THR:HG22	1:A:13:TYR:H	0.50	1.67	14	18
1:A:44:TRP:CZ2	1:A:58:LYS:HD2	0.50	2.42	3	7
1:A:66:ASP:O	1:A:66:ASP:OD1	0.49	2.30	8	7
1:A:33:VAL:O	1:A:36:LYS:NZ	0.49	2.45	2	9
1:A:62:VAL:CG1	1:A:63:VAL:N	0.49	2.75	13	1
2:B:78:ILE:HG23	2:B:84:ARG:C	0.49	2.28	2	7
1:A:5:VAL:CG1	1:A:65:SER:OG	0.49	2.60	12	6
1:A:23:LEU:C	1:A:23:LEU:CD1	0.48	2.81	12	18
1:A:35:GLU:CA	1:A:36:LYS:NZ	0.48	2.77	12	18
1:A:12:LEU:O	1:A:13:TYR:CD1	0.48	2.67	4	18
1:A:62:VAL:HG22	1:A:63:VAL:N	0.48	2.24	9	10
1:A:26:GLU:CD	1:A:26:GLU:N	0.48	2.68	8	18
1:A:49:ALA:C	1:A:51:GLY:N	0.48	2.68	3	18
1:A:19:THR:HG22	1:A:20:GLU:N	0.48	2.20	13	2
1:A:25:PHE:CD1	1:A:25:PHE:N	0.48	2.82	18	18
2:B:77:PRO:O	2:B:86:ILE:N	0.47	2.46	2	18
2:B:99:HIS:C	2:B:101:GLU:N	0.47	2.68	3	18
1:A:35:GLU:HA	1:A:36:LYS:HZ3	0.47	1.70	15	10
1:A:57:PRO:CB	1:A:59:ASN:HD21	0.47	2.23	18	1
1:A:21:GLU:C	1:A:22:GLU:OE1	0.47	2.53	13	13
1:A:48:ASN:C	1:A:50:ARG:N	0.47	2.68	4	18
1:A:29:GLU:OE2	1:A:30:THR:O	0.47	2.33	5	18
2:B:129:HIS:O	2:B:133:LEU:CD1	0.47	2.63	12	18
1:A:19:THR:OG1	1:A:22:GLU:CG	0.47	2.63	3	2
1:A:33:VAL:HG12	1:A:36:LYS:NZ	0.47	2.25	16	18
2:B:128:THR:O	2:B:132:GLN:N	0.47	2.37	18	11
1:A:41:PRO:O	1:A:42:GLU:HG3	0.46	2.10	14	3
2:B:112:PHE:N	2:B:112:PHE:CD1	0.46	2.84	11	18
2:B:105:CYS:HB2	2:B:126:CYS:SG	0.46	2.51	4	18

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:125:TYR:CE1	2:B:133:LEU:CD2	0.46	2.99	13	18
1:A:48:ASN:ND2	1:A:50:ARG:C	0.46	2.70	1	18
1:A:39:ASN:O	1:A:40:ASP:CB	0.46	2.64	7	18
1:A:59:ASN:ND2	1:A:59:ASN:H	0.46	2.09	18	1
1:A:15:PHE:CG	1:A:16:SER:N	0.45	2.85	1	18
1:A:62:VAL:HG12	1:A:63:VAL:N	0.45	2.27	13	1
2:B:129:HIS:O	2:B:133:LEU:N	0.45	2.47	1	13
1:A:5:VAL:HG12	1:A:65:SER:OG	0.45	2.12	12	4
1:A:7:HIS:CA	1:A:65:SER:OG	0.45	2.65	17	8
1:A:59:ASN:OD1	1:A:59:ASN:C	0.45	2.55	10	1
1:A:36:LYS:HB3	1:A:44:TRP:NE1	0.45	2.26	9	5
1:A:62:VAL:CG2	1:A:63:VAL:N	0.45	2.80	15	5
2:B:115:HIS:CD2	2:B:115:HIS:N	0.44	2.86	1	18
1:A:33:VAL:HG12	1:A:36:LYS:HZ2	0.44	1.72	15	5
2:B:84:ARG:CZ	2:B:84:ARG:CB	0.44	2.95	8	1
1:A:49:ALA:C	1:A:51:GLY:H	0.44	2.16	9	18
1:A:64:LEU:O	2:B:84:ARG:NH1	0.44	2.51	14	2
2:B:77:PRO:CG	2:B:90:VAL:HG21	0.43	2.43	1	18
2:B:118:TYR:CZ	2:B:127:GLU:CB	0.43	3.02	7	18
2:B:118:TYR:HB3	2:B:130:TYR:CD2	0.43	2.49	11	18
2:B:105:CYS:SG	2:B:107:LYS:N	0.43	2.92	12	18
1:A:48:ASN:ND2	1:A:48:ASN:C	0.43	2.72	12	14
2:B:89:ARG:HH11	2:B:89:ARG:HG2	0.43	1.74	3	13
1:A:57:PRO:CB	1:A:59:ASN:ND2	0.43	2.82	18	1
1:A:12:LEU:C	1:A:13:TYR:CG	0.43	2.93	13	18
1:A:35:GLU:N	1:A:36:LYS:NZ	0.42	2.67	3	18
1:A:64:LEU:O	2:B:84:ARG:NH2	0.42	2.52	7	1
2:B:130:TYR:OH	2:B:136:ASP:OD1	0.42	2.37	14	15
2:B:76:VAL:HG11	2:B:85:PRO:CB	0.42	2.45	5	15
1:A:11:THR:HG22	1:A:13:TYR:N	0.42	2.30	1	14
2:B:104:VAL:HG12	2:B:111:PRO:CB	0.41	2.45	12	17
1:A:63:VAL:CG1	2:B:82:CYS:O	0.41	2.68	18	1
1:A:22:GLU:OE1	1:A:22:GLU:CA	0.41	2.66	13	4
1:A:57:PRO:C	1:A:59:ASN:H	0.41	2.19	1	13
2:B:105:CYS:SG	2:B:125:TYR:HA	0.41	2.56	15	18
1:A:44:TRP:CE2	1:A:58:LYS:HD2	0.41	2.50	9	2
2:B:79:CYS:HA	2:B:86:ILE:HD11	0.41	1.92	10	12
2:B:90:VAL:HG12	2:B:91:VAL:N	0.41	2.31	9	12
1:A:57:PRO:HB3	1:A:59:ASN:HD21	0.41	1.75	18	1
2:B:130:TYR:O	2:B:134:PHE:N	0.41	2.53	1	1
1:A:34:ILE:C	1:A:36:LYS:HZ3	0.41	2.19	12	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:43:TRP:HB2	1:A:55:LEU:HD12	0.41	1.93	12	3
1:A:26:GLU:N	1:A:26:GLU:OE1	0.41	2.43	6	1
1:A:18:VAL:HG23	1:A:18:VAL:O	0.41	2.15	9	1
1:A:62:VAL:CG2	2:B:78:ILE:HD13	0.41	2.46	13	1
1:A:42:GLU:O	1:A:42:GLU:HG2	0.41	2.15	17	1
2:B:125:TYR:CZ	2:B:133:LEU:HD22	0.41	2.51	5	10
2:B:108:CYS:O	2:B:109:GLU:CB	0.41	2.69	14	6
1:A:50:ARG:NH1	1:A:50:ARG:CG	0.41	2.84	1	2
1:A:8:VAL:CG1	1:A:30:THR:CG2	0.41	2.99	7	6
2:B:129:HIS:O	2:B:133:LEU:HD12	0.40	2.16	4	10
2:B:79:CYS:CA	2:B:86:ILE:HD11	0.40	2.47	11	5
1:A:17:SER:OG	1:A:22:GLU:CG	0.40	2.69	3	1
2:B:99:HIS:O	2:B:101:GLU:N	0.40	2.55	7	2
2:B:129:HIS:O	2:B:130:TYR:C	0.40	2.60	16	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	69/71 (97%)	50±0 (73±1%)	13±0 (18±1%)	6±0 (9±0%)	1	12
2	B	64/66 (97%)	49±0 (77±0%)	14±0 (22±0%)	1±0 (2±0%)	13	57
All	All	2394/2466 (97%)	1789 (75%)	479 (20%)	126 (5%)	3	23

All 7 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	33	VAL	18
1	A	39	ASN	18
1	A	40	ASP	18
1	A	41	PRO	18
1	A	42	GLU	18
1	A	68	PRO	18
2	B	74	MET	18

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	64/64 (100%)	50±1 (79±1%)	14±1 (21±1%)	3	32
2	B	54/54 (100%)	49±0 (91±0%)	5±0 (9±0%)	12	58
All	All	2124/2124 (100%)	1790 (84%)	334 (16%)	5	42

All 23 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	17	SER	18
1	A	23	LEU	18
1	A	25	PHE	18
1	A	26	GLU	18
1	A	27	LYS	18
1	A	29	GLU	18
1	A	32	GLU	18
1	A	40	ASP	18
1	A	48	ASN	18
1	A	50	ARG	18
1	A	52	GLN	18
1	A	71	HIS	18
2	B	92	ASN	18
2	B	110	LYS	18
2	B	112	PHE	18
2	B	115	HIS	18
2	B	120	ARG	18
1	A	22	GLU	16
1	A	19	THR	8
2	B	84	ARG	1
1	A	58	LYS	1
1	A	60	TYR	1
1	A	59	ASN	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 monosaccharides in this entry.

6.6 Ligand geometry [i](#)

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

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided