



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 01:00 AM UTC

PDB ID : 2L1L / pdb_00002111
Title : NMR Solution Structure of the Phi0 PKI NES Peptide in Complex with CRM1-RanGTP
Authors : Madl, T.; Sattler, M.
Deposited on : 2010-07-29

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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

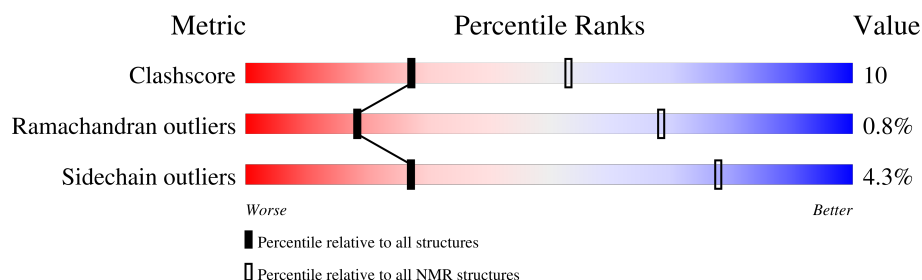
1 Overall quality at a glance

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	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

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	27	100%
2	B	127	87% 13%

2 Ensemble composition and analysis ⓘ

This entry contains 10 models. Model 2 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	B:504-B:630 (127)	0.00	2

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: No clusters in NmrClust output

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2475 atoms, of which 1247 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called cAMP-dependent protein kinase inhibitor alpha.

Mol	Chain	Residues	Atoms						Trace
1	A	27	Total	C	H	N	O	S	0
			380	114	189	32	44	1	

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP P61925
A	7	LEU	SER	conflict	UNP P61925
A	27	CYS	-	expression tag	UNP P61925

- Molecule 2 is a protein called Exportin-1.

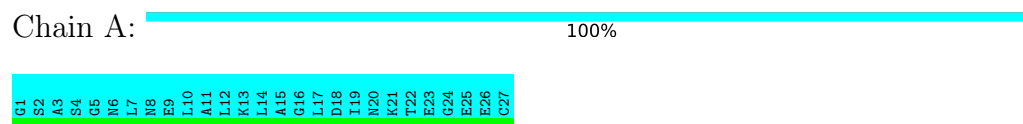
Mol	Chain	Residues	Atoms						Trace
2	B	127	Total	C	H	N	O	S	0
			2095	663	1058	181	184	9	

4 Residue-property plots [i](#)

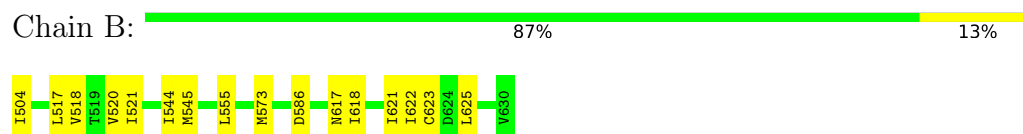
4.1 Average score per residue in the NMR ensemble

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: cAMP-dependent protein kinase inhibitor alpha



- Molecule 2: Exportin-1

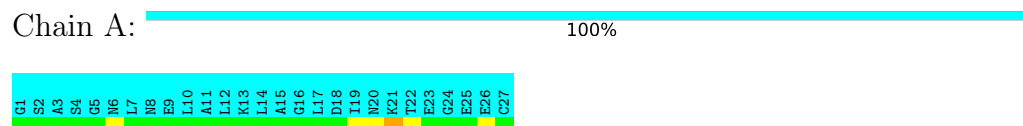


4.2 Scores per residue for each member of the ensemble

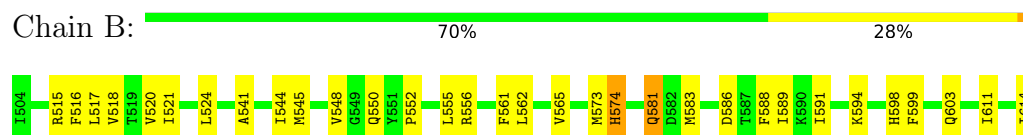
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: cAMP-dependent protein kinase inhibitor alpha



- Molecule 2: Exportin-1

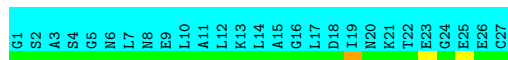




4.2.2 Score per residue for model 2 (medoid)

- Molecule 1: cAMP-dependent protein kinase inhibitor alpha

Chain A: 100%



- Molecule 2: Exportin-1

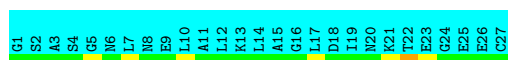
Chain B: 76% 23%



4.2.3 Score per residue for model 3

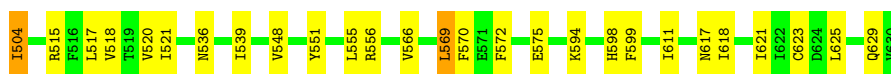
- Molecule 1: cAMP-dependent protein kinase inhibitor alpha

Chain A: 100%



- Molecule 2: Exportin-1

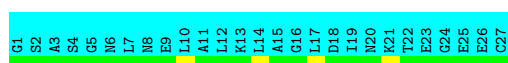
Chain B: 79% 20%



4.2.4 Score per residue for model 4

- Molecule 1: cAMP-dependent protein kinase inhibitor alpha

Chain A: 100%



- Molecule 2: Exportin-1

Chain B: 72% 25%



4.2.5 Score per residue for model 5

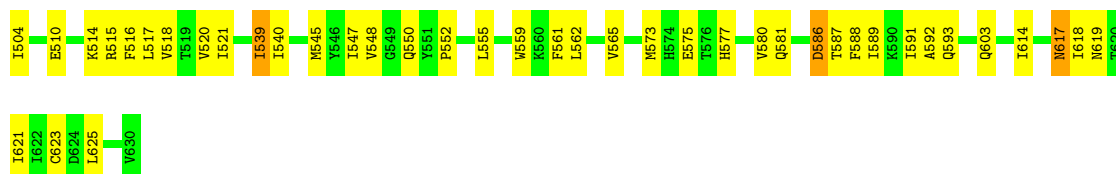
- Molecule 1: cAMP-dependent protein kinase inhibitor alpha

Chain A: 100%



- Molecule 2: Exportin-1

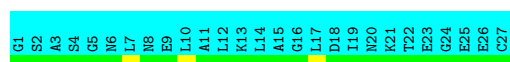
Chain B: 68% 30% .



4.2.6 Score per residue for model 6

- Molecule 1: cAMP-dependent protein kinase inhibitor alpha

Chain A: 100%



- Molecule 2: Exportin-1

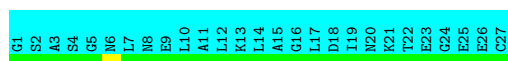
Chain B: 74% 24% .



4.2.7 Score per residue for model 7

- Molecule 1: cAMP-dependent protein kinase inhibitor alpha

Chain A: 100%



• Molecule 2: Exportin-1

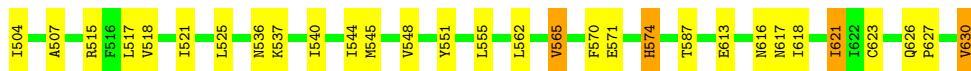
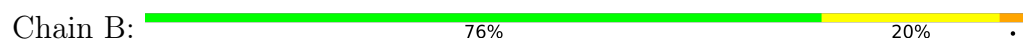


4.2.8 Score per residue for model 8

• Molecule 1: cAMP-dependent protein kinase inhibitor alpha



• Molecule 2: Exportin-1

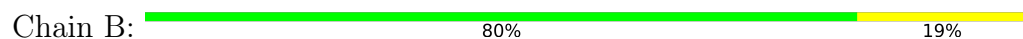


4.2.9 Score per residue for model 9

• Molecule 1: cAMP-dependent protein kinase inhibitor alpha



• Molecule 2: Exportin-1



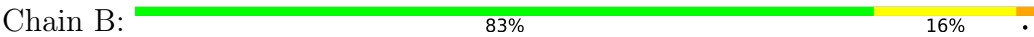
4.2.10 Score per residue for model 10

• Molecule 1: cAMP-dependent protein kinase inhibitor alpha





● Molecule 2: Exportin-1



5 Refinement protocol and experimental data overview

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

Of the 1000 calculated structures, 10 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
Sparky	structure solution	
ARIA	geometry optimization	1.2
CNS	geometry optimization	
CNS	refinement	

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [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	0	0	0	0±0
2	B	1037	1058	1050	20±5
All	All	10370	10580	10500	204

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 10.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:569:LEU:HD11	2:B:584:ALA:O	0.75	1.82	6	1
2:B:562:LEU:O	2:B:565:VAL:HG22	0.75	1.80	5	1
2:B:504:ILE:HG13	2:B:507:ALA:HB3	0.74	1.60	9	2
2:B:573:MET:HB3	2:B:625:LEU:HD11	0.69	1.65	10	5
2:B:577:HIS:O	2:B:580:VAL:HG12	0.68	1.88	7	2
2:B:504:ILE:HG12	2:B:507:ALA:HB3	0.68	1.66	8	1
2:B:627:PRO:O	2:B:630:VAL:HG22	0.67	1.89	6	2
2:B:626:GLN:O	2:B:630:VAL:HG13	0.62	1.95	6	2
2:B:622:ILE:HA	2:B:625:LEU:HD13	0.62	1.70	10	5
2:B:573:MET:HB3	2:B:625:LEU:HD21	0.61	1.72	7	2
2:B:545:MET:HB3	2:B:587:THR:HG21	0.61	1.73	7	5
2:B:516:PHE:O	2:B:520:VAL:HG13	0.60	1.97	5	2
2:B:577:HIS:O	2:B:580:VAL:HG22	0.59	1.97	5	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:517:LEU:HA	2:B:520:VAL:HG22	0.58	1.74	1	2
2:B:575:GLU:HB3	2:B:580:VAL:HG21	0.58	1.75	5	1
2:B:520:VAL:HG23	2:B:521:ILE:HG13	0.57	1.75	1	1
2:B:617:ASN:HD22	2:B:617:ASN:C	0.57	2.08	1	2
2:B:569:LEU:HD21	2:B:584:ALA:O	0.57	2.00	6	1
2:B:586:ASP:O	2:B:589:ILE:HG22	0.56	2.01	4	4
2:B:569:LEU:HD13	2:B:569:LEU:O	0.54	2.02	6	1
2:B:548:VAL:HG13	2:B:555:LEU:HD11	0.53	1.79	8	1
2:B:517:LEU:HD21	2:B:551:TYR:CG	0.53	2.39	8	1
2:B:536:ASN:O	2:B:539:ILE:HG22	0.52	2.05	6	2
2:B:626:GLN:O	2:B:630:VAL:HG23	0.52	2.05	7	1
2:B:583:MET:HE3	2:B:583:MET:HA	0.52	1.81	1	1
2:B:520:VAL:HG23	2:B:521:ILE:HG23	0.52	1.80	5	1
2:B:622:ILE:HA	2:B:625:LEU:HD23	0.52	1.81	7	1
2:B:534:LYS:HD3	2:B:534:LYS:H	0.51	1.63	6	1
2:B:599:PHE:CD1	2:B:599:PHE:C	0.51	2.88	9	1
2:B:617:ASN:ND2	2:B:617:ASN:C	0.51	2.68	2	1
2:B:581:GLN:NE2	2:B:581:GLN:H	0.51	2.03	1	1
2:B:517:LEU:O	2:B:521:ILE:HG22	0.51	2.06	3	3
2:B:521:ILE:HG22	2:B:547:ILE:CD1	0.51	2.36	7	1
2:B:566:VAL:HA	2:B:569:LEU:HD23	0.50	1.83	3	1
2:B:520:VAL:HG13	2:B:521:ILE:HG23	0.50	1.83	6	2
2:B:556:ARG:HG3	2:B:598:HIS:NE2	0.50	2.21	7	4
2:B:575:GLU:CG	2:B:580:VAL:HG11	0.50	2.37	7	2
2:B:588:PHE:O	2:B:591:ILE:HG22	0.50	2.06	1	1
2:B:539:ILE:CG2	2:B:540:ILE:N	0.49	2.74	5	1
2:B:525:LEU:HD22	2:B:544:ILE:HD13	0.49	1.85	6	1
2:B:545:MET:SD	2:B:583:MET:HB3	0.48	2.49	4	1
2:B:558:HIS:ND1	2:B:561:PHE:HB2	0.48	2.23	4	1
2:B:525:LEU:HD12	2:B:544:ILE:HD13	0.48	1.85	8	1
2:B:599:PHE:C	2:B:599:PHE:HD1	0.48	2.16	9	1
2:B:562:LEU:HD12	2:B:591:ILE:HD11	0.47	1.86	1	1
2:B:580:VAL:HG13	2:B:581:GLN:N	0.47	2.24	7	2
2:B:517:LEU:HD12	2:B:518:VAL:N	0.47	2.25	2	1
2:B:517:LEU:HA	2:B:520:VAL:HG12	0.47	1.85	6	3
2:B:517:LEU:O	2:B:517:LEU:HD23	0.47	2.09	5	1
2:B:542:SER:O	2:B:545:MET:HG2	0.47	2.10	4	1
2:B:524:LEU:HD23	2:B:544:ILE:HB	0.47	1.86	1	1
2:B:516:PHE:O	2:B:520:VAL:HG23	0.47	2.10	9	3
2:B:563:LYS:HG3	2:B:603:GLN:HE22	0.47	1.69	2	1
2:B:551:TYR:O	2:B:555:LEU:HD23	0.46	2.11	3	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:544:ILE:O	2:B:547:ILE:CG1	0.46	2.64	7	1
2:B:548:VAL:HG11	2:B:565:VAL:HG21	0.46	1.85	8	1
2:B:534:LYS:H	2:B:534:LYS:HD2	0.46	1.69	9	1
2:B:599:PHE:O	2:B:611:ILE:HG23	0.46	2.11	3	2
2:B:515:ARG:O	2:B:518:VAL:HG12	0.46	2.11	4	2
2:B:515:ARG:O	2:B:518:VAL:HG22	0.46	2.11	5	3
2:B:559:TRP:HB2	2:B:603:GLN:NE2	0.46	2.26	5	2
2:B:544:ILE:HG13	2:B:545:MET:N	0.46	2.26	2	2
2:B:545:MET:SD	2:B:583:MET:C	0.46	2.99	4	1
2:B:544:ILE:O	2:B:547:ILE:HG12	0.46	2.11	7	1
2:B:622:ILE:CG2	2:B:630:VAL:HG22	0.46	2.42	7	1
2:B:548:VAL:HG23	2:B:555:LEU:HD21	0.45	1.88	3	1
2:B:571:GLU:O	2:B:574:HIS:ND1	0.45	2.50	8	1
2:B:563:LYS:O	2:B:566:VAL:HG12	0.45	2.12	2	2
2:B:562:LEU:O	2:B:565:VAL:HG12	0.45	2.11	8	1
2:B:580:VAL:HG23	2:B:581:GLN:N	0.45	2.27	5	1
2:B:618:ILE:N	2:B:618:ILE:HD13	0.45	2.27	6	1
2:B:572:PHE:HA	2:B:575:GLU:HG2	0.45	1.87	3	2
2:B:561:PHE:O	2:B:565:VAL:HG13	0.44	2.12	5	1
2:B:618:ILE:O	2:B:621:ILE:HG12	0.44	2.13	1	5
2:B:504:ILE:CG1	2:B:507:ALA:HB3	0.44	2.40	8	1
2:B:613:GLU:HA	2:B:616:ASN:ND2	0.44	2.27	8	1
2:B:614:ILE:O	2:B:617:ASN:ND2	0.44	2.50	2	3
2:B:628:GLN:NE2	2:B:628:GLN:H	0.44	2.10	2	1
2:B:545:MET:HE3	2:B:583:MET:HG2	0.44	1.90	1	1
2:B:555:LEU:HD12	2:B:562:LEU:HA	0.44	1.88	2	1
2:B:559:TRP:HB2	2:B:603:GLN:HE21	0.44	1.71	6	1
2:B:618:ILE:O	2:B:621:ILE:CG2	0.44	2.66	3	2
2:B:575:GLU:HG3	2:B:580:VAL:HG11	0.44	1.88	7	1
2:B:569:LEU:HD13	2:B:569:LEU:C	0.43	2.38	6	1
2:B:514:LYS:O	2:B:517:LEU:HG	0.43	2.13	2	1
2:B:591:ILE:HG13	2:B:592:ALA:N	0.43	2.29	5	2
2:B:618:ILE:HG12	2:B:619:ASN:N	0.43	2.28	6	1
2:B:617:ASN:O	2:B:618:ILE:C	0.43	2.61	1	10
2:B:574:HIS:ND1	2:B:574:HIS:C	0.43	2.76	1	2
2:B:537:LYS:HA	2:B:540:ILE:HG22	0.43	1.89	4	2
2:B:514:LYS:O	2:B:518:VAL:HG13	0.43	2.14	5	1
2:B:618:ILE:O	2:B:621:ILE:HG22	0.43	2.14	8	3
2:B:596:ARG:HA	2:B:599:PHE:CE2	0.43	2.48	9	1
2:B:548:VAL:HG22	2:B:555:LEU:HD11	0.42	1.91	1	3
2:B:504:ILE:N	2:B:504:ILE:HD13	0.42	2.29	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:536:ASN:O	2:B:539:ILE:CG1	0.42	2.67	2	1
2:B:625:LEU:HD22	2:B:629:GLN:HE21	0.42	1.74	3	1
2:B:599:PHE:O	2:B:611:ILE:HG12	0.42	2.14	9	1
2:B:550:GLN:C	2:B:552:PRO:HD3	0.42	2.39	4	3
2:B:541:ALA:O	2:B:544:ILE:HG12	0.42	2.13	2	2
2:B:536:ASN:O	2:B:539:ILE:CG2	0.42	2.68	3	1
2:B:570:PHE:HA	2:B:621:ILE:HD11	0.42	1.90	8	1
2:B:586:ASP:O	2:B:589:ILE:HG12	0.42	2.14	10	1
2:B:569:LEU:C	2:B:569:LEU:HD12	0.42	2.39	3	1
2:B:617:ASN:HB2	2:B:620:THR:HG22	0.42	1.91	4	1
2:B:541:ALA:O	2:B:544:ILE:HG22	0.42	2.14	4	1
2:B:618:ILE:HG13	2:B:619:ASN:N	0.42	2.30	4	2
2:B:518:VAL:O	2:B:522:LYS:HB2	0.42	2.14	9	1
2:B:618:ILE:O	2:B:621:ILE:CG1	0.41	2.68	2	2
2:B:628:GLN:HE21	2:B:628:GLN:HA	0.41	1.74	4	1
2:B:547:ILE:O	2:B:550:GLN:HG2	0.41	2.15	5	1
2:B:575:GLU:CB	2:B:580:VAL:HG21	0.41	2.45	5	1
2:B:521:ILE:O	2:B:525:LEU:HD13	0.41	2.14	8	1
2:B:622:ILE:HG21	2:B:630:VAL:HG13	0.41	1.91	7	1
2:B:541:ALA:O	2:B:544:ILE:CG2	0.41	2.68	1	1
2:B:552:PRO:HG2	2:B:594:LYS:HG2	0.41	1.91	4	1
2:B:621:ILE:O	2:B:621:ILE:HD13	0.41	2.16	8	1
2:B:594:LYS:N	2:B:594:LYS:HD2	0.41	2.31	1	1
2:B:588:PHE:HA	2:B:591:ILE:HG12	0.41	1.92	5	2
2:B:618:ILE:HG23	2:B:619:ASN:N	0.41	2.31	5	1
2:B:615:LEU:HD13	2:B:618:ILE:HD11	0.41	1.93	9	1
2:B:626:GLN:HB3	2:B:627:PRO:HD2	0.41	1.93	7	1
2:B:617:ASN:O	2:B:620:THR:HG22	0.41	2.15	10	1
2:B:556:ARG:HG3	2:B:598:HIS:HE2	0.40	1.76	1	1
2:B:569:LEU:HD21	2:B:584:ALA:HA	0.40	1.92	6	1
2:B:569:LEU:CD2	2:B:584:ALA:HA	0.40	2.47	6	1
2:B:575:GLU:HB3	2:B:580:VAL:HG11	0.40	1.92	9	1
2:B:561:PHE:O	2:B:565:VAL:HG23	0.40	2.16	1	1
2:B:568:LYS:HA	2:B:568:LYS:HE2	0.40	1.94	2	1
2:B:536:ASN:O	2:B:540:ILE:HD13	0.40	2.16	8	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	0	-	-	-	-	
2	B	125/127 (98%)	118±4 (94±3%)	5±0 (4±0%)	1±0 (1±0%)	18	68
All	All	1237/1540 (80%)	1178 (95%)	49 (4%)	10 (1%)	18	68

All 1 unique Ramachandran outliers are listed below.

Mol	Chain	Res	Type	Models (Total)
2	B	623	CYS	10

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	0	-	-	-	
2	B	115/115 (100%)	110±1 (96±1%)	5±1 (4±1%)	27	78
All	All	1150/1350 (85%)	1101 (96%)	49 (4%)	27	78

All 34 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)
2	B	504	ILE	4
2	B	574	HIS	3
2	B	617	ASN	3
2	B	570	PHE	3
2	B	510	GLU	3
2	B	568	LYS	2
2	B	594	LYS	2
2	B	545	MET	2

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Mol	Chain	Res	Type	Models (Total)
2	B	630	VAL	2
2	B	581	GLN	1
2	B	603	GLN	1
2	B	569	LEU	1
2	B	531	LYS	1
2	B	628	GLN	1
2	B	539	ILE	1
2	B	586	ASP	1
2	B	593	GLN	1
2	B	560	LYS	1
2	B	590	LYS	1
2	B	618	ILE	1
2	B	511	GLU	1
2	B	529	GLU	1
2	B	559	TRP	1
2	B	596	ARG	1
2	B	537	LYS	1
2	B	565	VAL	1
2	B	621	ILE	1
2	B	525	LEU	1
2	B	534	LYS	1
2	B	550	GLN	1
2	B	598	HIS	1
2	B	599	PHE	1
2	B	517	LEU	1
2	B	622	ILE	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.6 Ligand geometry

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided