



Full wwPDB NMR Structure Validation Report ⓘ

Jun 7, 2020 – 03:29 am BST

PDB ID : 6MNL
Title : NMR solution structures of second bromodomain of BRD4 with FOXO3a peptide
Authors : Zeng, L.; Zhou, M.-M.
Deposited on : 2018-10-02

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:

Cyrange : Kirchner and Güntert (2011)
NmrClust : Kelley et al. (1996)
MolProbity : 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 : 2.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

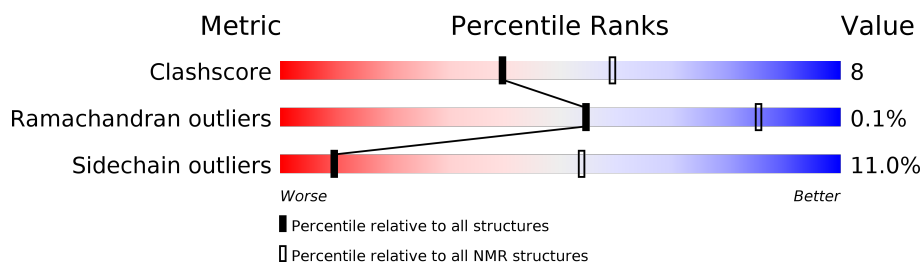
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 84%.

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	16	100%
2	B	128	58% 25% 17%

2 Ensemble composition and analysis

This entry contains 20 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: *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:349-B:454 (106)	0.21	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 4 clusters and 3 single-model clusters were found.

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

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 2304 atoms, of which 1144 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called FOXO3a peptide.

Mol	Chain	Residues	Atoms					Trace
1	A	16	Total	C	H	N	O	0
			249	72	126	28	23	

- Molecule 2 is a protein called Bromodomain-containing protein 4.

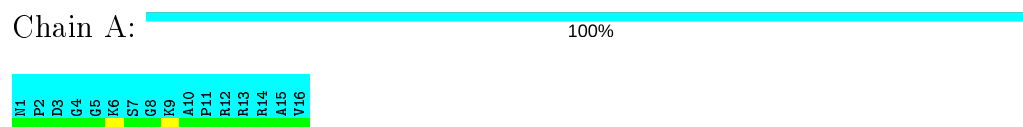
Mol	Chain	Residues	Atoms						Trace
2	B	128	Total	C	H	N	O	S	0
			2055	658	1018	175	193	11	

4 Residue-property plots

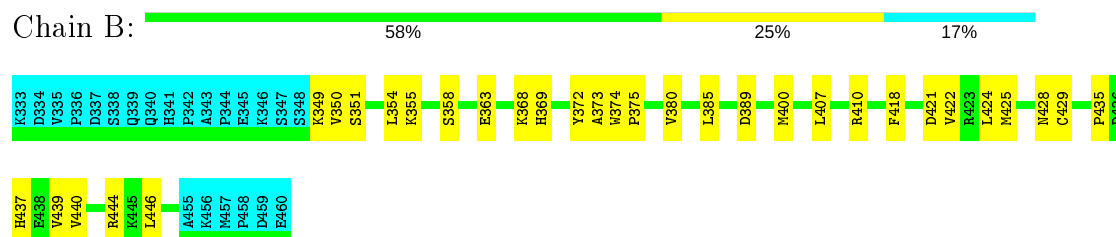
4.1 Average score per residue in the NMR ensemble

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: FOXO3a peptide



- Molecule 2: Bromodomain-containing protein 4

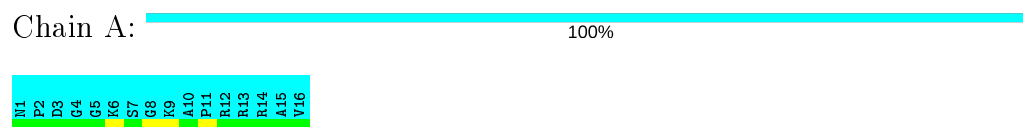


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

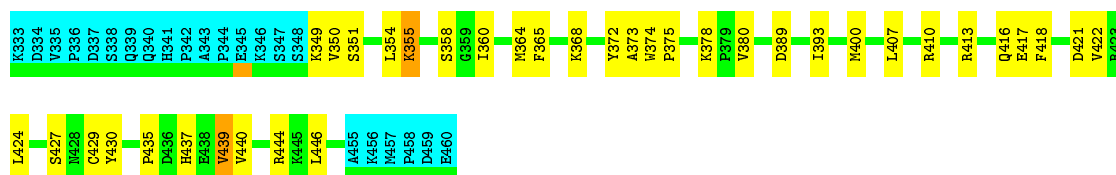
4.2.1 Score per residue for model 1

- Molecule 1: FOXO3a peptide



- Molecule 2: Bromodomain-containing protein 4

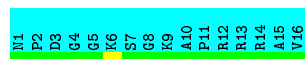




4.2.2 Score per residue for model 2

- Molecule 1: FOXO3a peptide

Chain A: 100%



- Molecule 2: Bromodomain-containing protein 4

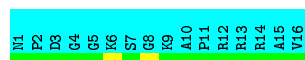
Chain B: 59% 23% 17%



4.2.3 Score per residue for model 3

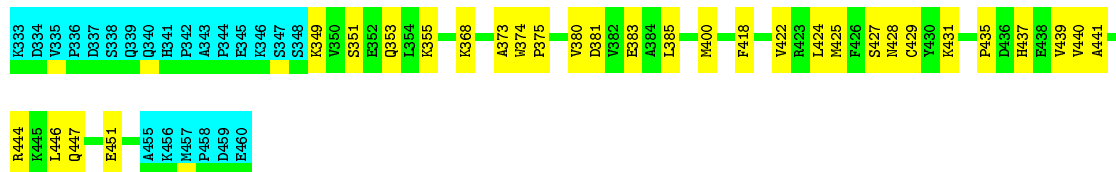
- Molecule 1: FOXO3a peptide

Chain A: 100%



- Molecule 2: Bromodomain-containing protein 4

Chain B: 59% 23% 17%



4.2.4 Score per residue for model 4

- Molecule 1: FOXO3a peptide

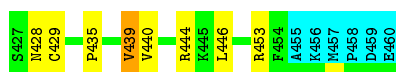
Chain A:  100%



- Molecule 2: Bromodomain-containing protein 4

Chain B:  55%  26%  17%

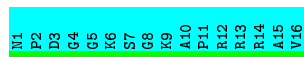




4.2.7 Score per residue for model 7

- Molecule 1: FOXO3a peptide

Chain A: 100%



- Molecule 2: Bromodomain-containing protein 4

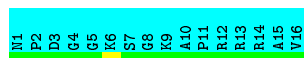
Chain B: 55% 27% 17%



4.2.8 Score per residue for model 8

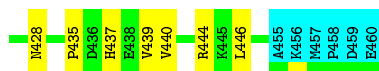
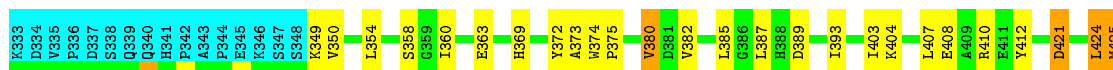
- Molecule 1: FOXO3a peptide

Chain A: 100%



- Molecule 2: Bromodomain-containing protein 4

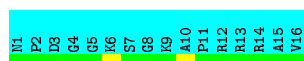
Chain B: 57% 23% 17%



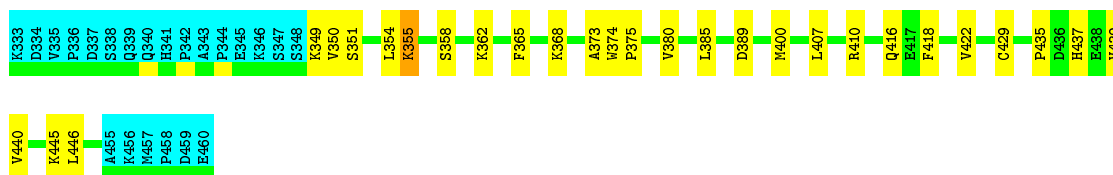
4.2.9 Score per residue for model 9

- Molecule 1: FOXO3a peptide

Chain A: 100%

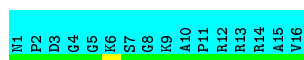


- Molecule 2: Bromodomain-containing protein 4

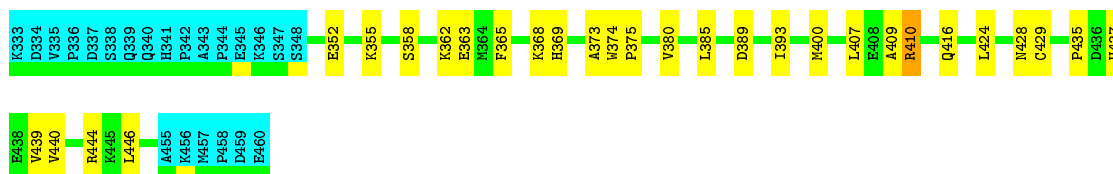


4.2.10 Score per residue for model 10 (medoid)

- Molecule 1: FOXO3a peptide



- Molecule 2: Bromodomain-containing protein 4

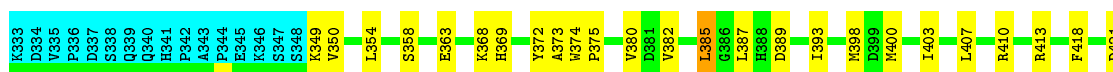


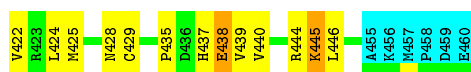
4.2.11 Score per residue for model 11

- Molecule 1: FOXO3a peptide



- Molecule 2: Bromodomain-containing protein 4

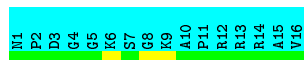




4.2.12 Score per residue for model 12

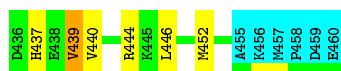
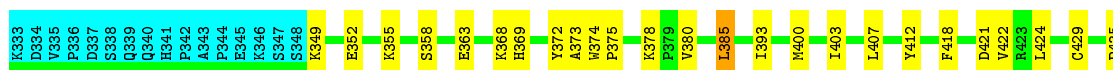
- Molecule 1: FOXO3a peptide

Chain A:  100%



- Molecule 2: Bromodomain-containing protein 4

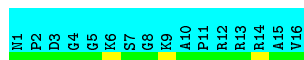
Chain B:  59% 23% 17%



4.2.13 Score per residue for model 13

- Molecule 1: FOXO3a peptide

Chain A:  100%



- Molecule 2: Bromodomain-containing protein 4

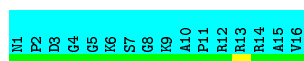
Chain B:  55% 25% 17%



4.2.14 Score per residue for model 14

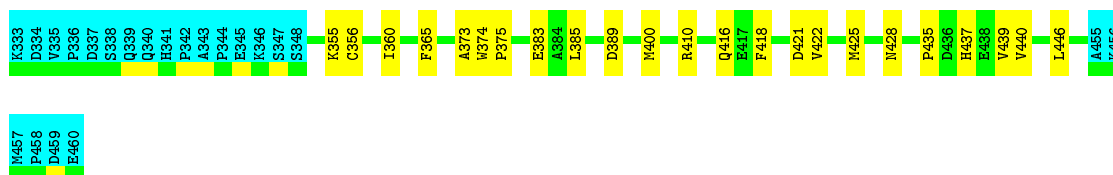
- Molecule 1: FOXO3a peptide

Chain A:  100%



- Molecule 2: Bromodomain-containing protein 4

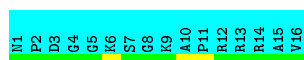
Chain B: 65% 18% 17%



4.2.15 Score per residue for model 15

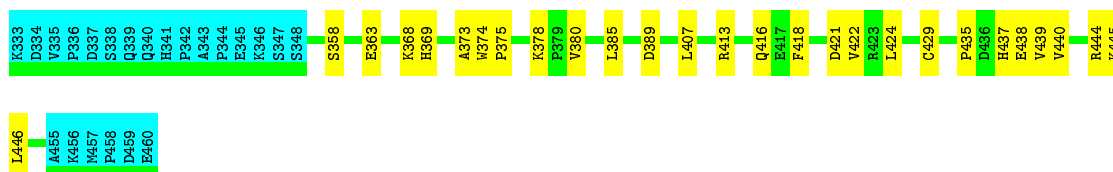
- Molecule 1: FOXO3a peptide

Chain A: 100%



- Molecule 2: Bromodomain-containing protein 4

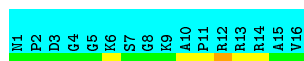
Chain B: 62% 21% 17%



4.2.16 Score per residue for model 16

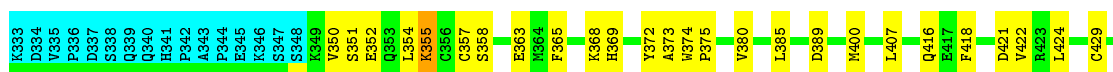
- Molecule 1: FOXO3a peptide

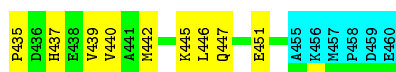
Chain A: 100%



- Molecule 2: Bromodomain-containing protein 4

Chain B: 55% 27% 17%

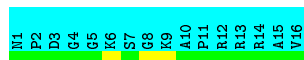




4.2.17 Score per residue for model 17

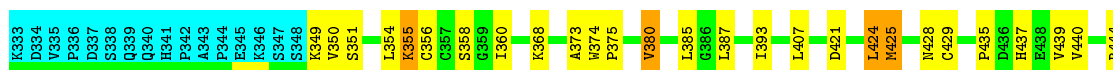
- Molecule 1: FOXO3a peptide

Chain A: 100%



- Molecule 2: Bromodomain-containing protein 4

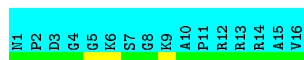
Chain B: 61% 19% 17%



4.2.18 Score per residue for model 18

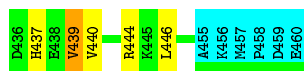
- Molecule 1: FOXO3a peptide

Chain A: 100%



- Molecule 2: Bromodomain-containing protein 4

Chain B: 57% 23% 17%



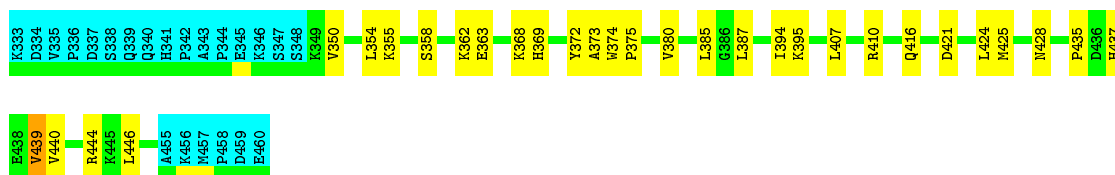
4.2.19 Score per residue for model 19

- Molecule 1: FOXO3a peptide

Chain A: 100%



- Molecule 2: Bromodomain-containing protein 4

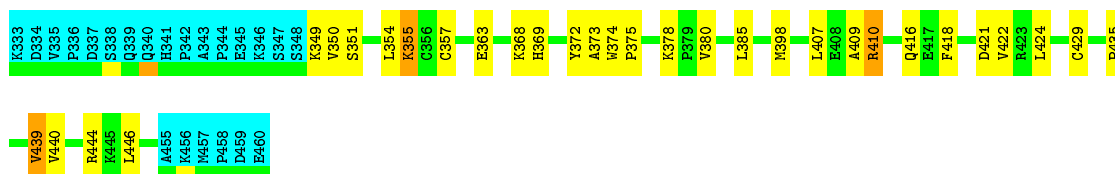


4.2.20 Score per residue for model 20

- Molecule 1: FOXO3a peptide



- Molecule 2: Bromodomain-containing protein 4



5 Refinement protocol and experimental data overview

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

Of the 200 calculated structures, 20 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
ARIA	refinement	2.3
CNS	refinement	
CNS	structure calculation	

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)	input_cs.cif
Number of chemical shift lists	1
Total number of shifts	1511
Number of shifts mapped to atoms	1511
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	84%

No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality ⓘ

6.1 Standard geometry ⓘ

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

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 ⓘ

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	868	858	854	13±3
All	All	17360	17160	17080	266

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 8.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:440:VAL:O	2:B:444:ARG:HG2	0.64	1.93	17	12
2:B:373:ALA:HB2	2:B:446:LEU:HD13	0.63	1.69	6	1
2:B:357:CYS:SG	2:B:407:LEU:HD13	0.61	2.34	20	3
2:B:373:ALA:HB2	2:B:446:LEU:HD21	0.60	1.73	19	19
2:B:438:GLU:HG3	2:B:442:MET:SD	0.57	2.40	5	1
2:B:435:PRO:HA	2:B:440:VAL:HG21	0.56	1.77	19	20
2:B:439:VAL:HA	2:B:442:MET:SD	0.56	2.41	5	1
2:B:349:LYS:O	2:B:352:GLU:HG2	0.55	2.02	12	2
2:B:394:ILE:HG23	2:B:428:ASN:OD1	0.55	2.03	19	2
2:B:360:ILE:HG22	2:B:364:MET:SD	0.54	2.42	1	2
2:B:447:GLN:O	2:B:451:GLU:HG2	0.54	2.02	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:351:SER:O	2:B:355:LYS:HE3	0.54	2.03	17	7
2:B:425:MET:HA	2:B:428:ASN:OD1	0.53	2.03	14	6
2:B:409:ALA:C	2:B:410:ARG:HG2	0.53	2.24	2	3
2:B:350:VAL:O	2:B:354:LEU:HG	0.53	2.03	19	14
2:B:447:GLN:O	2:B:451:GLU:HB2	0.51	2.05	13	3
2:B:358:SER:O	2:B:362:LYS:HD3	0.51	2.05	19	5
2:B:374:TRP:N	2:B:375:PRO:HD2	0.50	2.22	12	20
2:B:358:SER:HA	2:B:407:LEU:HD21	0.49	1.85	6	13
2:B:418:PHE:O	2:B:422:VAL:HG23	0.49	2.07	13	15
2:B:437:HIS:O	2:B:440:VAL:HG22	0.48	2.08	19	18
2:B:363:GLU:O	2:B:369:HIS:HD2	0.48	1.91	18	13
2:B:351:SER:O	2:B:355:LYS:HD3	0.48	2.09	18	3
2:B:358:SER:CA	2:B:407:LEU:HD21	0.47	2.39	8	9
2:B:444:ARG:NE	2:B:444:ARG:HA	0.47	2.24	12	2
2:B:374:TRP:CD2	2:B:375:PRO:HD3	0.47	2.44	7	9
2:B:433:ASN:HD22	2:B:433:ASN:N	0.46	2.07	13	1
2:B:365:PHE:CE1	2:B:400:MET:HB3	0.46	2.46	4	9
2:B:421:ASP:O	2:B:424:LEU:HB2	0.46	2.11	8	3
2:B:356:CYS:O	2:B:360:ILE:HG13	0.46	2.10	6	3
2:B:382:VAL:HG13	2:B:387:LEU:HB2	0.46	1.87	7	4
2:B:360:ILE:HG12	2:B:453:ARG:HE	0.46	1.70	6	1
2:B:357:CYS:SG	2:B:407:LEU:CD1	0.45	3.05	16	3
2:B:413:ARG:NE	2:B:413:ARG:HA	0.45	2.26	6	1
2:B:403:ILE:HG13	2:B:412:TYR:OH	0.45	2.12	12	2
2:B:425:MET:HA	2:B:428:ASN:ND2	0.43	2.28	8	2
2:B:400:MET:HA	2:B:403:ILE:HG22	0.43	1.90	4	5
2:B:426:PHE:N	2:B:426:PHE:CD2	0.43	2.86	6	1
2:B:445:LYS:O	2:B:448:ASP:HB3	0.43	2.14	7	1
2:B:413:ARG:HA	2:B:413:ARG:NE	0.43	2.29	11	3
2:B:422:VAL:HA	2:B:425:MET:HE2	0.42	1.90	11	1
2:B:445:LYS:HG3	2:B:446:LEU:N	0.42	2.30	16	2
2:B:427:SER:O	2:B:431:LYS:HG2	0.42	2.14	3	1
2:B:393:ILE:N	2:B:393:ILE:HD12	0.42	2.29	12	3
2:B:393:ILE:HD12	2:B:393:ILE:N	0.42	2.29	1	3
2:B:444:ARG:HA	2:B:444:ARG:NE	0.41	2.31	20	1
2:B:441:ALA:O	2:B:444:ARG:HB2	0.41	2.16	3	1
2:B:389:ASP:O	2:B:393:ILE:HD13	0.41	2.16	1	1
2:B:394:ILE:HG22	2:B:397:PRO:HG3	0.41	1.92	2	1
2:B:368:LYS:HD3	2:B:369:HIS:CE1	0.41	2.50	18	1
2:B:429:CYS:HA	2:B:433:ASN:HD22	0.41	1.75	4	1
2:B:413:ARG:HB3	2:B:417:GLU:OE1	0.41	2.16	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:404:LYS:O	2:B:408:GLU:HG3	0.40	2.16	8	1
2:B:420:ALA:O	2:B:424:LEU:HD23	0.40	2.17	2	1
2:B:360:ILE:O	2:B:363:GLU:HB3	0.40	2.16	8	1
2:B:427:SER:HA	2:B:430:TYR:CD2	0.40	2.51	1	1
2:B:351:SER:O	2:B:355:LYS:HE2	0.40	2.16	7	1
2:B:360:ILE:O	2:B:364:MET:HG3	0.40	2.16	1	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	106/128 (83%)	102±1 (96±1%)	4±1 (4±1%)	0±0 (0±0%)	54	85
All	All	2120/2880 (74%)	2042 (96%)	75 (4%)	3 (0%)	54	85

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)
2	B	438	GLU	2
2	B	380	VAL	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	0	-	-	-	
2	B	93/113 (82%)	83±2 (89±2%)	10±2 (11±2%)	10	54
All	All	1860/2440 (76%)	1656 (89%)	204 (11%)	10	54

All 29 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	439	VAL	20
2	B	385	LEU	19
2	B	380	VAL	17
2	B	368	LYS	15
2	B	355	LYS	14
2	B	421	ASP	14
2	B	424	LEU	14
2	B	410	ARG	13
2	B	389	ASP	12
2	B	372	TYR	11
2	B	349	LYS	10
2	B	416	GLN	10
2	B	378	LYS	6
2	B	398	MET	4
2	B	425	MET	3
2	B	383	GLU	3
2	B	445	LYS	3
2	B	400	MET	2
2	B	444	ARG	2
2	B	428	ASN	2
2	B	390	TYR	2
2	B	442	MET	1
2	B	433	ASN	1
2	B	396	HIS	1
2	B	452	MET	1
2	B	395	LYS	1
2	B	413	ARG	1
2	B	381	ASP	1
2	B	429	CYS	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
1	ALY	A	6	1	10,11,12	0.63±0.02	0±0 (0±0%)
1	ALY	A	9	1	10,11,12	0.61±0.02	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.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
1	ALY	A	6	1	7,12,14	0.99±0.05	0±0 (0±0%)
1	ALY	A	9	1	7,12,14	0.97±0.03	0±0 (0±0%)

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	ALY	A	6	1	-	0±0,9,10,12	-
1	ALY	A	9	1	-	0±0,9,10,12	-

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates ⓘ

There are no carbohydrates 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

The completeness of assignment taking into account all chemical shift lists is 84% for the well-defined parts and 79% for the entire structure.

7.1 Chemical shift list 1

File name: input_cs.cif

Chemical shift list name: *assigned_chem_shift_list_0*

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	1511
Number of shifts mapped to atoms	1511
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	17

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	133	-0.36 ± 0.11	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	127	0.28 ± 0.11	None needed (< 0.5 ppm)
$^{13}\text{C}'$	8	—	None (insufficient data)
^{15}N	115	0.53 ± 0.39	None needed (imprecise)

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 84%, i.e. 1151 atoms were assigned a chemical shift out of a possible 1367. 14 out of 14 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	414/520 (80%)	203/207 (98%)	112/212 (53%)	99/101 (98%)
Sidechain	609/701 (87%)	384/417 (92%)	223/253 (88%)	2/31 (6%)

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	Total	¹H	¹³C	¹⁵N
Aromatic	128/146 (88%)	68/76 (89%)	59/61 (97%)	1/9 (11%)
Overall	1151/1367 (84%)	655/700 (94%)	394/526 (75%)	102/141 (72%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 79%, i.e. 1411 atoms were assigned a chemical shift out of a possible 1790. 16 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	515/688 (75%)	259/273 (95%)	141/284 (50%)	115/131 (88%)
Sidechain	766/948 (81%)	495/567 (87%)	269/335 (80%)	2/46 (4%)
Aromatic	130/154 (84%)	69/80 (86%)	60/63 (95%)	1/11 (9%)
Overall	1411/1790 (79%)	823/920 (89%)	470/682 (69%)	118/188 (63%)

7.1.4 Statistically unusual chemical shifts ⓘ

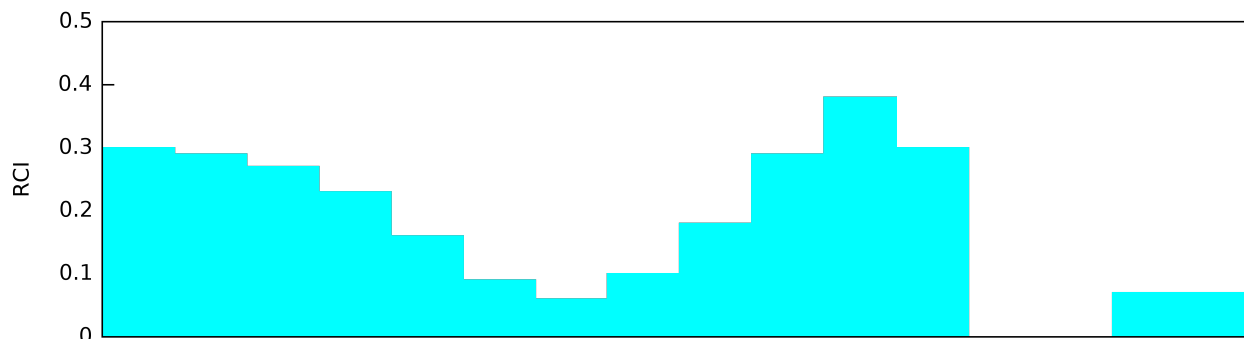
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
2	B	347	SER	C	8.45	183.48 – 165.88	-94.4
2	B	386	GLY	C	7.90	183.33 – 164.53	-88.3
2	B	346	LYS	C	8.51	186.58 – 166.78	-84.9
2	B	385	LEU	C	7.48	187.06 – 166.96	-84.3
2	B	447	GLN	C	11.57	186.20 – 166.50	-83.6
2	B	388	HIS	C	8.30	185.27 – 165.27	-83.5
2	B	437	HIS	C	9.41	185.27 – 165.27	-82.9
2	B	366	ALA	C	9.11	188.57 – 166.97	-78.1
2	B	375	PRO	HG2	-0.72	3.48 – 0.38	-8.5
2	B	447	GLN	HB2	0.07	3.30 – 0.80	-7.9
2	B	375	PRO	HB2	-0.39	3.82 – 0.32	-7.0
2	B	375	PRO	HD2	1.57	5.45 – 1.85	-5.8
2	B	422	VAL	HG11	-0.67	2.13 – -0.47	-5.8
2	B	422	VAL	HG13	-0.67	2.13 – -0.47	-5.8
2	B	422	VAL	HG12	-0.67	2.13 – -0.47	-5.8
2	B	357	CYS	HB2	0.38	5.20 – 0.70	-5.7
2	B	439	VAL	H	11.72	11.69 – 4.89	5.0

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

The images below report *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:



Random coil index (RCI) for chain B:

