



Full wwPDB NMR Structure Validation Report ⓘ

Apr 24, 2023 – 10:20 am BST

PDB ID : 7ZEZ
BMRB ID : 34728
Title : Trimolecular complex Cyp33-RRMdelta alpha : MLL1-PHD3 : H3K4me3
Authors : Blatter, M.; Allain, F.; Meylan, C.
Deposited on : 2022-03-31

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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
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) : 2.32.2

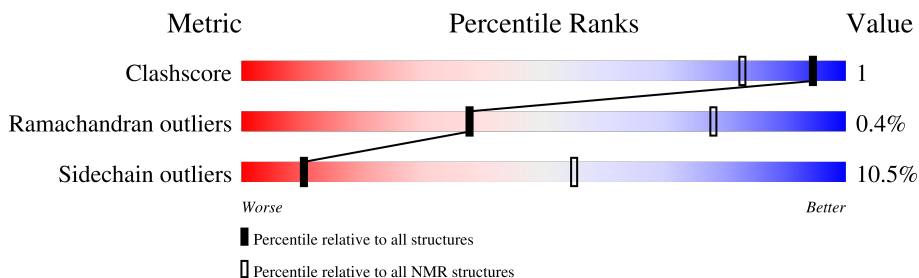
1 Overall quality at a glance

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	93	
2	B	64	
3	D	13	

2 Ensemble composition and analysis i

This entry contains 20 models. Model 1 is the overall representative, medoid model (most similar to other models).

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

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:3-A:87, B:1569-B:1627, D:1-D:3, D:5-D:6 (149)	0.44	1

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 2 clusters and 1 single-model cluster was found.

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

3 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2620 atoms, of which 1284 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Isoform 3 of Peptidyl-prolyl cis-trans isomerase E.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
1	A	93	1443	462	716	123	139	3	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ALA	-	expression tag	UNP Q9UNP9
A	-2	GLY	-	expression tag	UNP Q9UNP9
A	-1	HIS	-	expression tag	UNP Q9UNP9

- Molecule 2 is a protein called MLL cleavage product N320.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
2	B	64	970	310	460	83	106	11	0

- Molecule 3 is a protein called Histone H3.

Mol	Chain	Residues	Atoms					Trace
			Total	C	H	N	O	
3	D	13	205	57	108	22	18	0

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

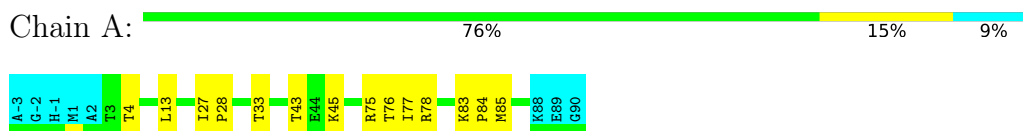
Mol	Chain	Residues	Atoms	
			Total	Zn
4	B	2	2	2

4 Residue-property plots

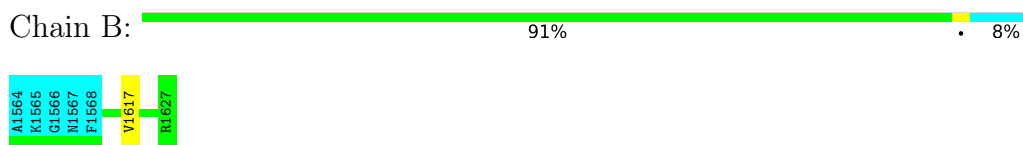
4.1 Average score per residue in the NMR ensemble

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.

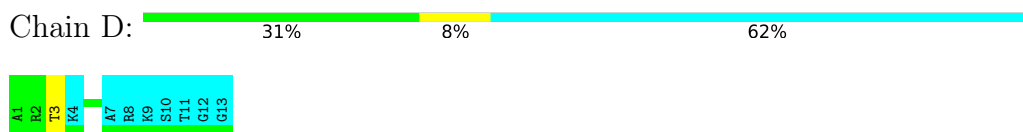
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320



- Molecule 3: Histone H3

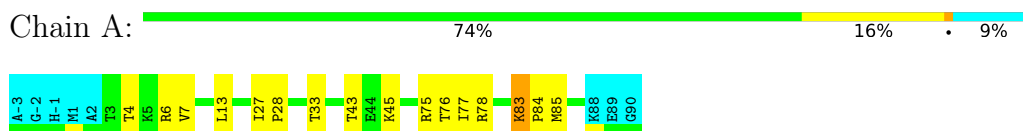


4.2 Scores per residue for each member of the ensemble

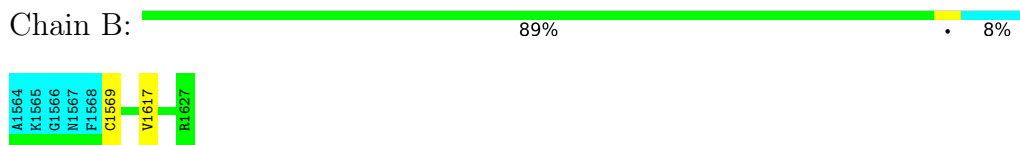
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1 (medoid)

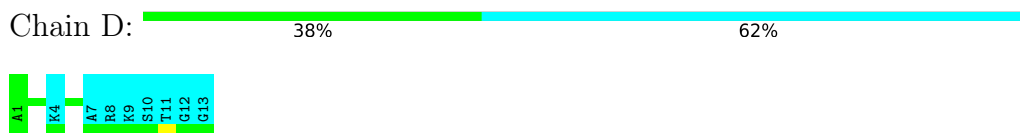
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

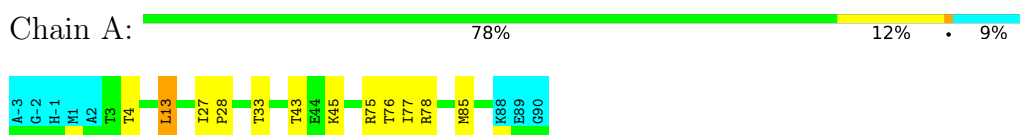


- Molecule 3: Histone H3

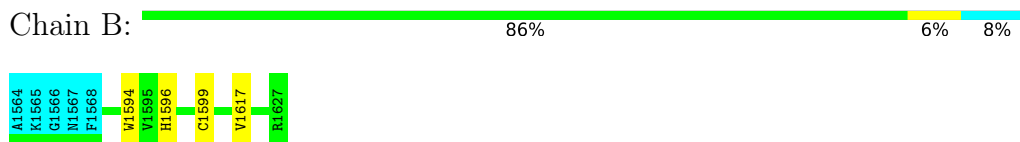


4.2.2 Score per residue for model 2

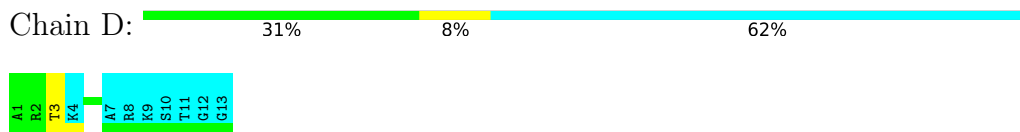
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

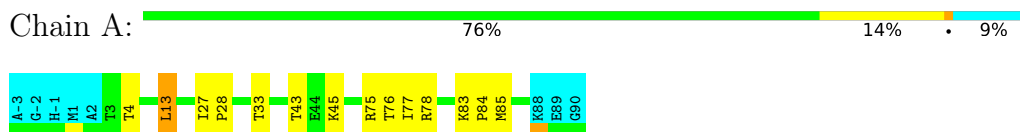


- Molecule 3: Histone H3




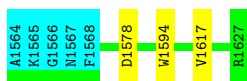
4.2.3 Score per residue for model 3

- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



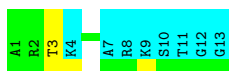
- Molecule 2: MLL cleavage product N320

Chain B:  88% 5% 8%




- Molecule 3: Histone H3

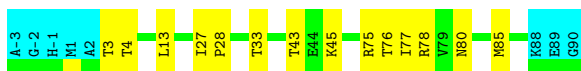
Chain D:  31% 8% 62%




4.2.4 Score per residue for model 4

- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E

Chain A:  76% 15% 9%



- Molecule 2: MLL cleavage product N320

Chain B:  89% 8%




- Molecule 3: Histone H3

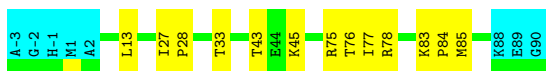
Chain D:  23% 15% 62%




4.2.5 Score per residue for model 5

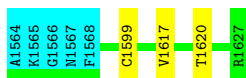
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E

Chain A:  77% 14% 9%

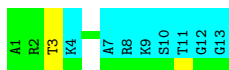


- Molecule 2: MLL cleavage product N320

Chain B:  88% 5% 8%

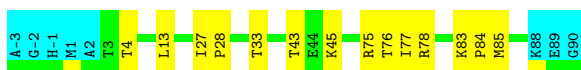
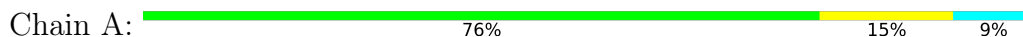


- Molecule 3: Histone H3

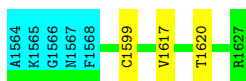
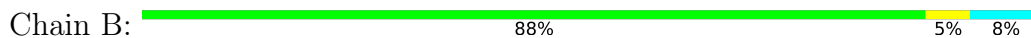


4.2.6 Score per residue for model 6

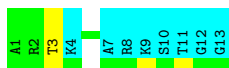
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320



- Molecule 3: Histone H3



4.2.7 Score per residue for model 7

- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E

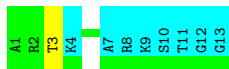


- Molecule 2: MLL cleavage product N320



- Molecule 3: Histone H3

Chain D: 




4.2.8 Score per residue for model 8

- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E

Chain A: 



- Molecule 2: MLL cleavage product N320

Chain B: 




- Molecule 3: Histone H3

Chain D: 




4.2.9 Score per residue for model 9

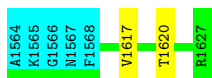
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E

Chain A: 

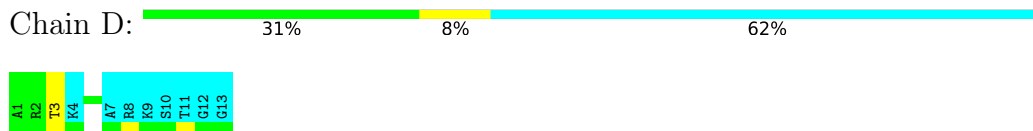


- Molecule 2: MLL cleavage product N320

Chain B: 

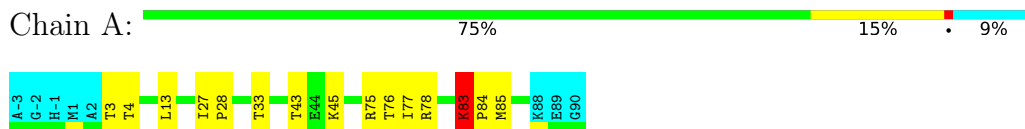


- Molecule 3: Histone H3

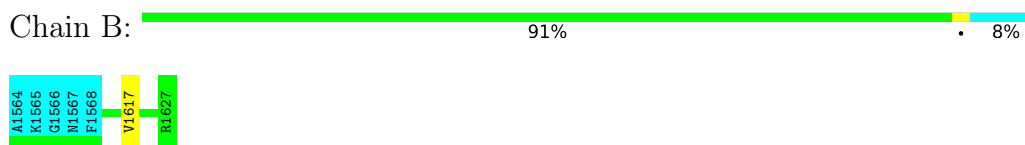


4.2.10 Score per residue for model 10

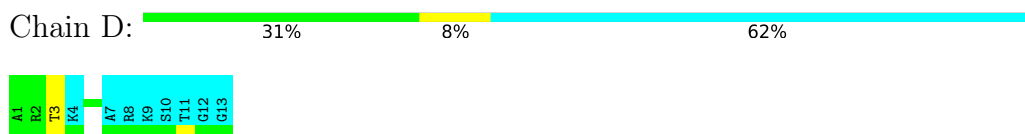
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

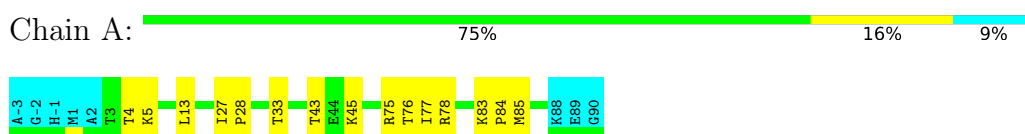


- Molecule 3: Histone H3

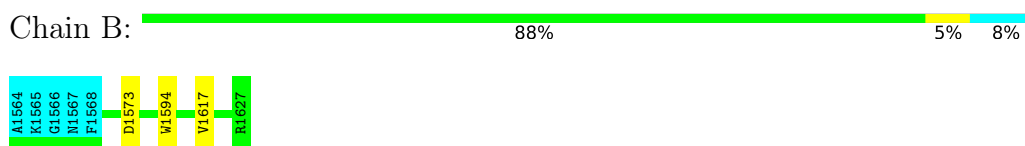


4.2.11 Score per residue for model 11

- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E

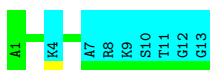


- Molecule 2: MLL cleavage product N320



- Molecule 3: Histone H3



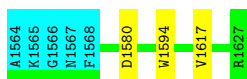
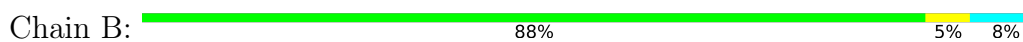


4.2.12 Score per residue for model 12

- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

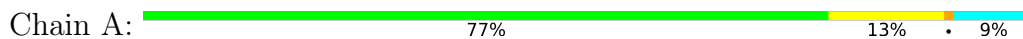


- Molecule 3: Histone H3

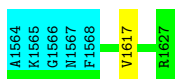
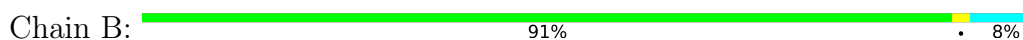


4.2.13 Score per residue for model 13

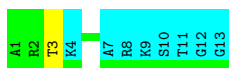
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

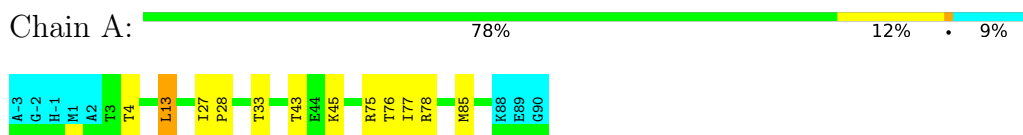


- Molecule 3: Histone H3

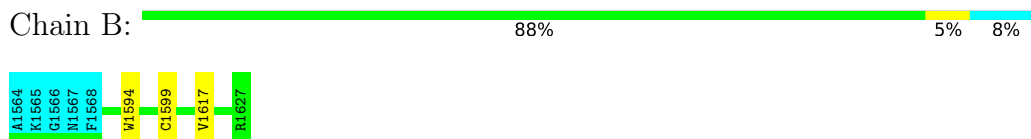


4.2.14 Score per residue for model 14

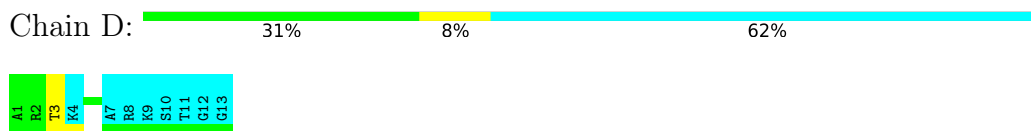
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

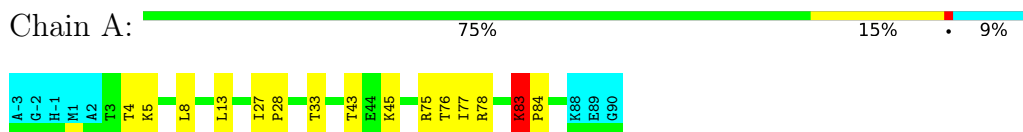


- Molecule 3: Histone H3

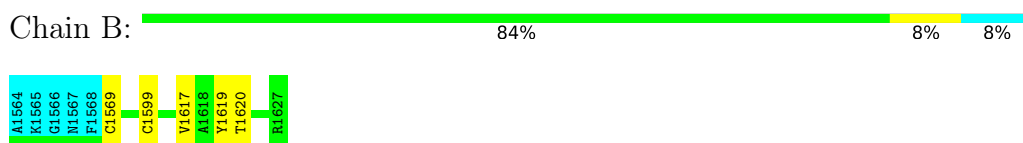


4.2.15 Score per residue for model 15

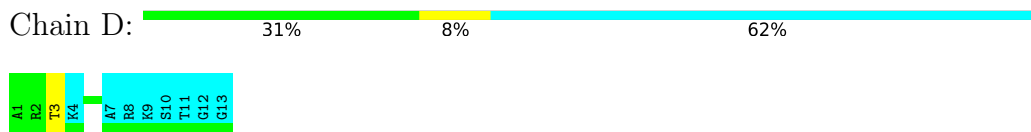
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

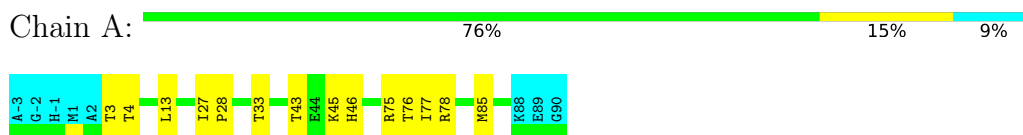


- Molecule 3: Histone H3

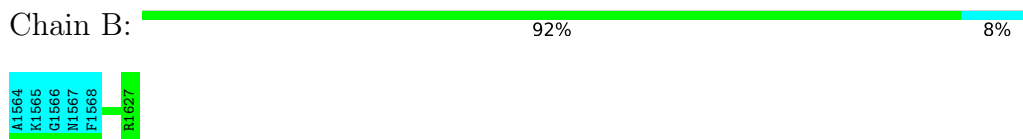


4.2.16 Score per residue for model 16

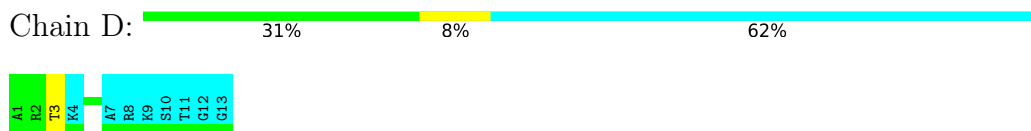
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

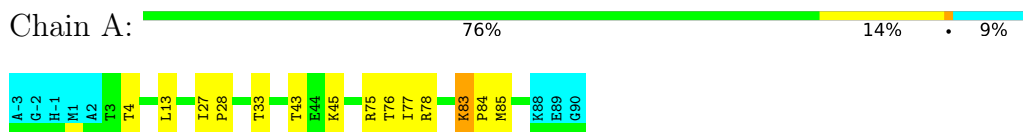


- Molecule 3: Histone H3

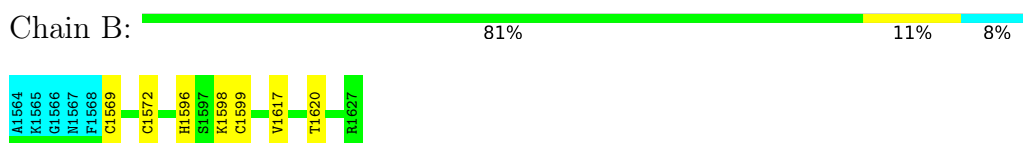


4.2.17 Score per residue for model 17

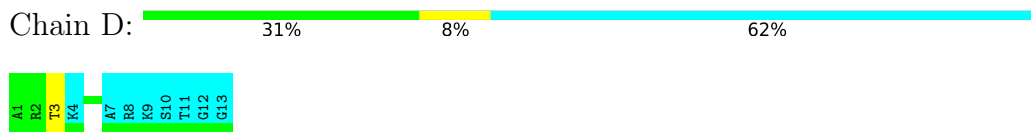
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

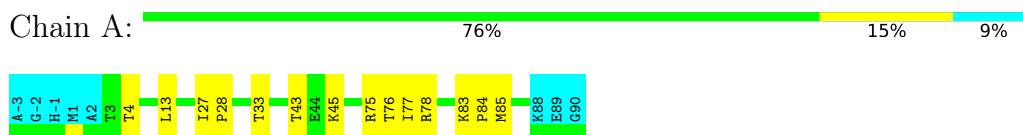


- Molecule 3: Histone H3

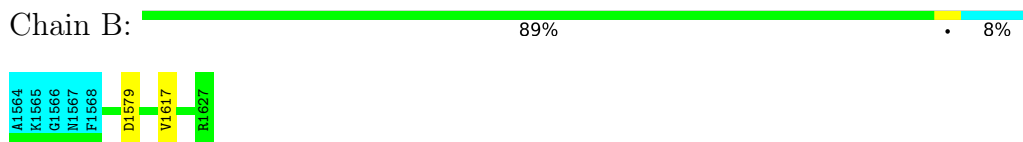


4.2.18 Score per residue for model 18

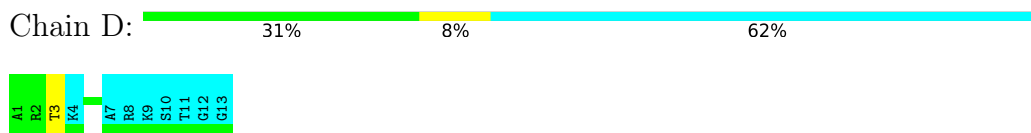
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320

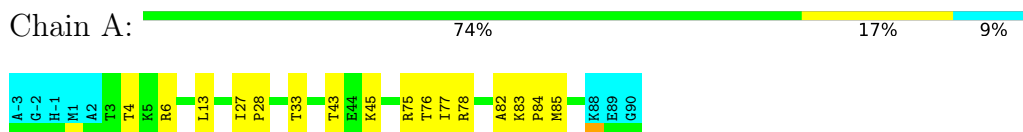


- Molecule 3: Histone H3

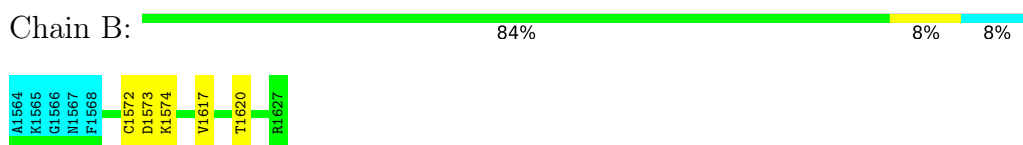


4.2.19 Score per residue for model 19

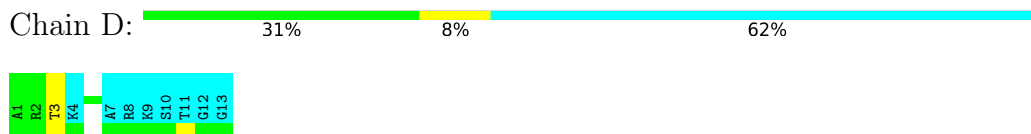
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E



- Molecule 2: MLL cleavage product N320




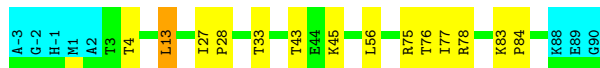
- Molecule 3: Histone H3




4.2.20 Score per residue for model 20

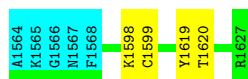
- Molecule 1: Isoform 3 of Peptidyl-prolyl cis-trans isomerase E

Chain A:  76% 14% 9%



- Molecule 2: MLL cleavage product N320

Chain B:  86% 6% 8%



- Molecule 3: Histone H3

Chain D:  23% 15% 62%



5 Refinement protocol and experimental data overview

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

Of the 250 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
Amber	refinement	
CYANA	structure calculation	3

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: M3L, ZN

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	673	663	663	2±1
2	B	473	424	424	0±0
3	D	39	40	42	0±0
All	All	23740	22540	22580	49

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	
				Worst	Total
1:A:83:LYS:H	1:A:84:PRO:HD2	0.48	1.67	10	4
2:B:1572:CYS:SG	2:B:1574:LYS:HE2	0.48	2.47	19	1
1:A:83:LYS:N	1:A:84:PRO:CD	0.48	2.77	5	12
1:A:27:ILE:N	1:A:28:PRO:CD	0.45	2.80	17	20
1:A:8:LEU:HD12	1:A:8:LEU:N	0.43	2.28	7	1
1:A:83:LYS:H	1:A:84:PRO:CD	0.43	2.26	17	1
2:B:1596:HIS:HB2	2:B:1599:CYS:HB2	0.42	1.91	2	1
1:A:27:ILE:N	1:A:28:PRO:HD2	0.42	2.30	11	5
2:B:1619:TYR:CG	2:B:1620:THR:N	0.42	2.88	20	2
2:B:1596:HIS:HB2	2:B:1599:CYS:SG	0.41	2.55	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:83:LYS:N	1:A:84:PRO:HD2	0.41	2.31	18	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	85/93 (91%)	80±1 (94±1%)	5±1 (5±1%)	0±0 (1±1%)	29	74
2	B	58/64 (91%)	55±2 (94±3%)	3±1 (6±3%)	0±0 (0±1%)	44	80
3	D	4/13 (31%)	4±1 (94±13%)	0±1 (6±13%)	0±0 (0±0%)	100	100
All	All	2940/3400 (86%)	2763 (94%)	164 (6%)	13 (0%)	38	78

All 6 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	13	LEU	5
1	A	83	LYS	4
2	B	1578	ASP	1
2	B	1580	ASP	1
2	B	1579	ASP	1
1	A	82	ALA	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	70/74 (95%)	59±1 (84±2%)	11±1 (16±2%)	5	42
2	B	57/60 (95%)	55±1 (97±2%)	2±1 (3±2%)	42	88

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	D	4/8 (50%)	3±0 (75±11%)	1±0 (25±11%)	2 25
All	All	2620/2840 (92%)	2344 (89%)	276 (11%)	10 55

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)
1	A	13	LEU	20
1	A	33	THR	20
1	A	43	THR	20
1	A	45	LYS	20
1	A	75	ARG	20
1	A	76	THR	20
1	A	77	ILE	20
1	A	78	ARG	20
1	A	4	THR	18
2	B	1617	VAL	18
3	D	3	THR	17
1	A	85	MET	15
1	A	6	ARG	6
1	A	3	THR	5
2	B	1599	CYS	5
2	B	1620	THR	5
1	A	7	VAL	3
2	B	1569	CYS	3
1	A	80	ASN	3
3	D	6	THR	3
1	A	8	LEU	2
1	A	56	LEU	2
1	A	83	LYS	2
1	A	5	LYS	2
2	B	1573	ASP	2
2	B	1598	LYS	2
1	A	40	ASP	1
1	A	46	HIS	1
2	B	1572	CYS	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](#)

1 non-standard protein/DNA/RNA residue is 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
3	M3L	D	4	3	10,11,12	0.70±0.01	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
3	M3L	D	4	3	9,14,16	0.38±0.05	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
3	M3L	D	4	3	-	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 [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