



wwPDB EM Validation Summary Report ⓘ

May 24, 2026 – 03:10 AM JST

PDB ID : 23AT / pdb_000023at
EMDB ID : EMD-68806
Title : Structure of Arabidopsis SNX1 (Class II, 6-fold)
Authors : Li, Y.B.; Tao, R.; Zhang, H.; Wen, X.K.; Leung, S.K.P.; Lau, W.C.Y.; Jiang, L.W.; Cui, Y.
Deposited on : 2026-01-30
Resolution : 8.70 Å(reported)

This is a wwPDB EM Validation Summary 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/EMValidationReportHelp>
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:

EMDB validation analysis : **FAILED**
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : **NOT EXECUTED**
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

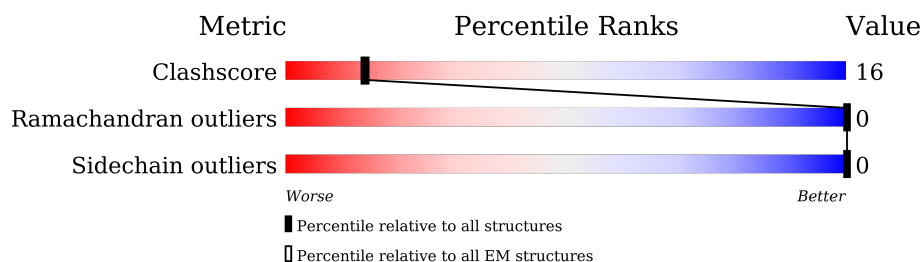
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 8.70 Å.

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)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	0	402	61% 33% 6%
1	1	402	61% 33% 6%
1	2	402	61% 33% 6%
1	3	402	60% 34% 6%
1	4	402	61% 33% 6%
1	5	402	61% 34% 6%
1	6	402	59% 35% 6%
1	7	402	62% 32% 6%
1	8	402	60% 34% 6%

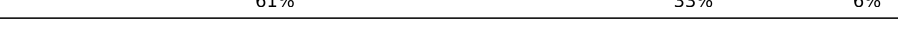
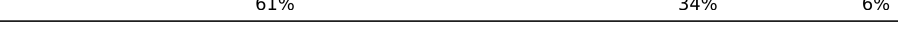
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Mol	Chain	Length	Quality of chain
1	9	402	
1	A	402	
1	AA	402	
1	AB	402	
1	B	402	
1	C	402	
1	D	402	
1	E	402	
1	F	402	
1	G	402	
1	H	402	
1	I	402	
1	J	402	
1	K	402	
1	L	402	
1	M	402	
1	N	402	
1	O	402	
1	P	402	
1	Q	402	
1	R	402	
1	S	402	
1	T	402	
1	U	402	
1	V	402	

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Mol	Chain	Length	Quality of chain
1	W	402	
1	X	402	
1	Y	402	
1	Z	402	
1	a	402	
1	b	402	
1	c	402	
1	d	402	
1	e	402	
1	f	402	
1	g	402	
1	h	402	
1	i	402	
1	j	402	
1	k	402	
1	l	402	
1	m	402	
1	n	402	
1	o	402	
1	p	402	
1	q	402	
1	r	402	
1	s	402	
1	t	402	
1	u	402	

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Mol	Chain	Length	Quality of chain
1	v	402	 61% 33% 6%
1	w	402	 61% 34% 6%
1	x	402	 61% 34% 6%
1	y	402	 60% 34% 6%
1	z	402	 61% 33% 6%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 198528 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Sorting nexin 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	0	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	1	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	2	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	3	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	4	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	5	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	6	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	7	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	8	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	9	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	A	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	AA	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	AB	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	B	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	C	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	D	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	E	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	F	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	G	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	H	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	I	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	J	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	K	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	L	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	M	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	N	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	O	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	P	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	Q	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	R	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	S	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	T	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	U	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	V	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	W	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	X	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	Y	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	Z	379	Total 3102	C 1973	N 536	O 583	S 10	0	0

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	a	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	b	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	c	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	d	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	e	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	f	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	g	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	h	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	i	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	j	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	k	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	l	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	m	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	n	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	o	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	p	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	q	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	r	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	s	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	t	379	Total 3102	C 1973	N 536	O 583	S 10	0	0
1	u	379	Total 3102	C 1973	N 536	O 583	S 10	0	0

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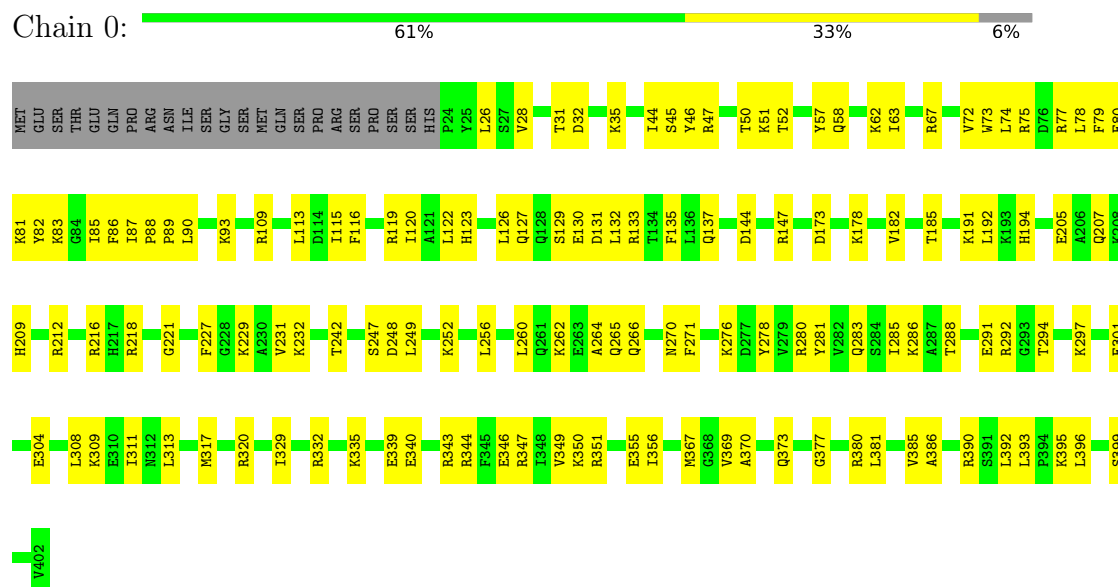
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Mol	Chain	Residues	Atoms					AltConf	Trace
1	v	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	w	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	x	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	y	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		
1	z	379	Total	C	N	O	S	0	0
			3102	1973	536	583	10		

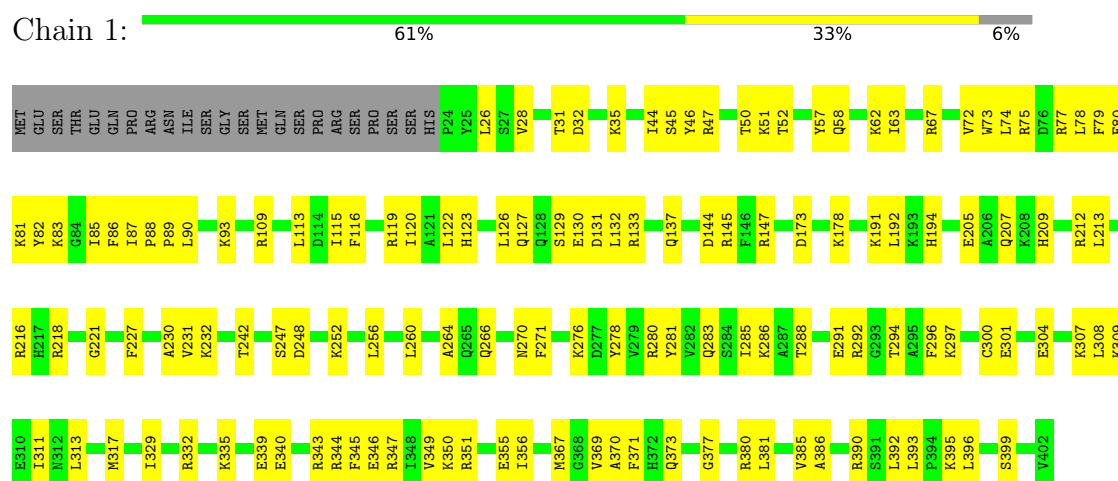
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-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 outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Sorting nexin 1



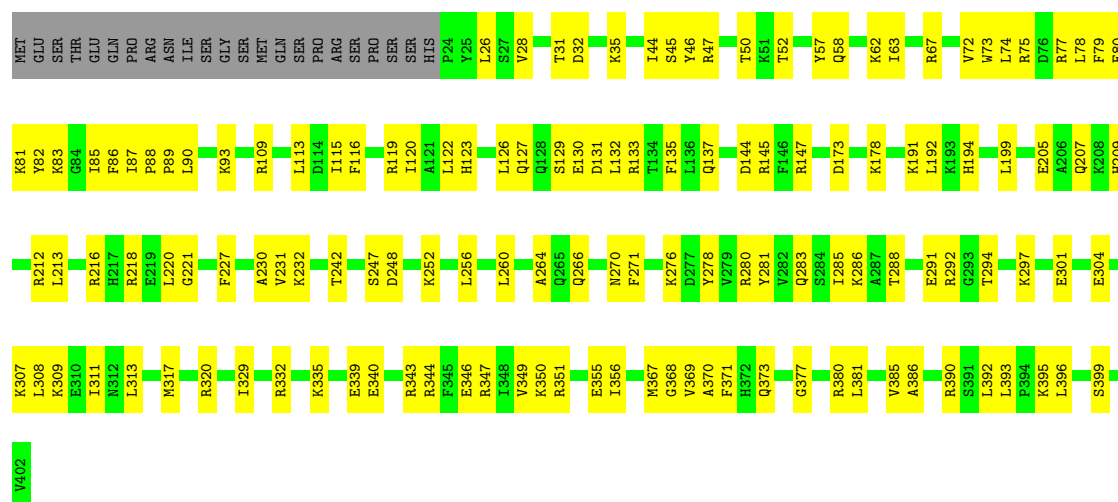
- Molecule 1: Sorting nexin 1



- Molecule 1: Sorting nexin 1

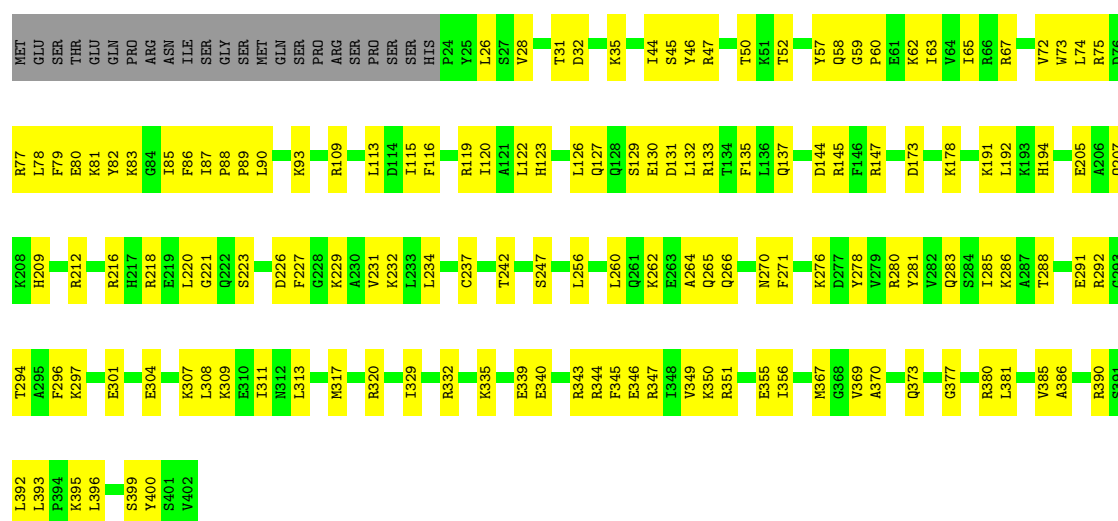


Chain 5:  61% 34% 6%



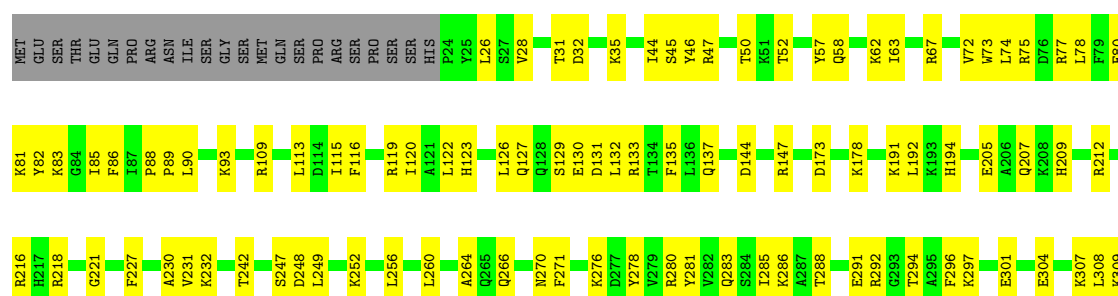
• Molecule 1: Sorting nexin 1

Chain 6:  59% 35% 6%



• Molecule 1: Sorting nexin 1

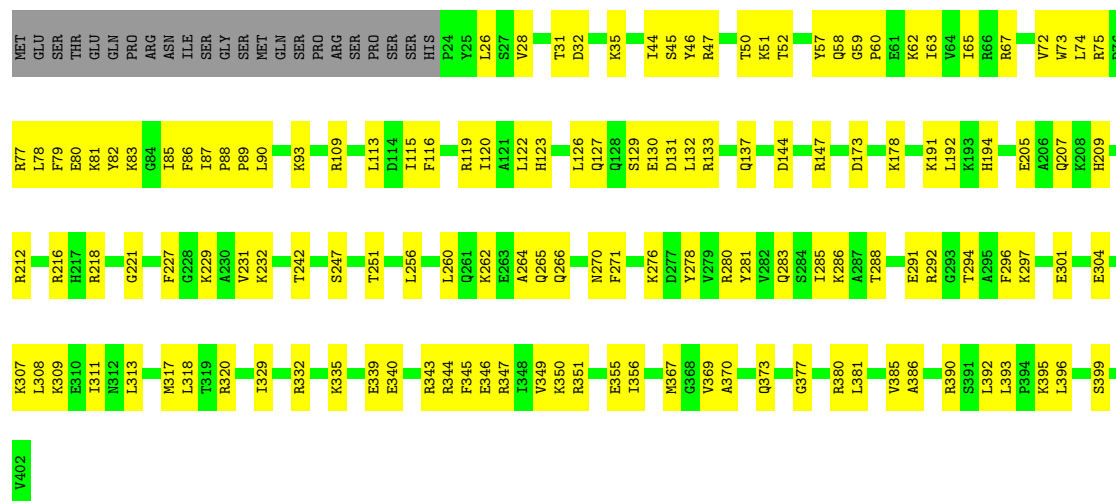
Chain 7:  62% 32% 6%





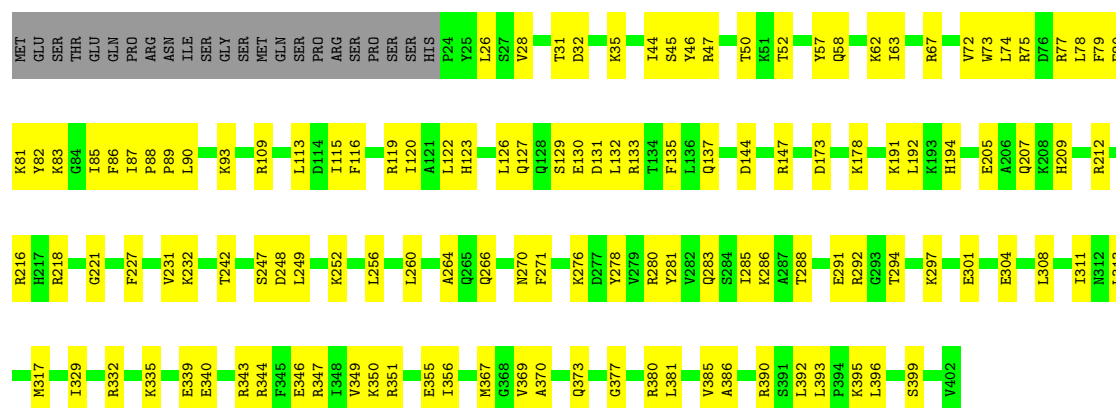
• Molecule 1: Sorting nexin 1

Chain 8: 60% 34% 6%



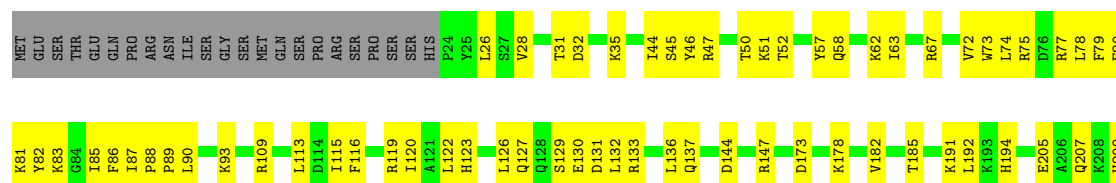
• Molecule 1: Sorting nexin 1

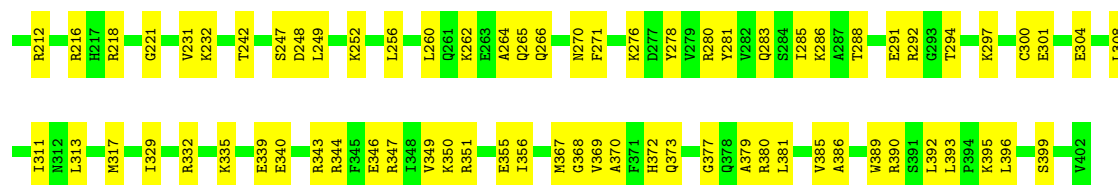
Chain 9: 63% 31% 6%



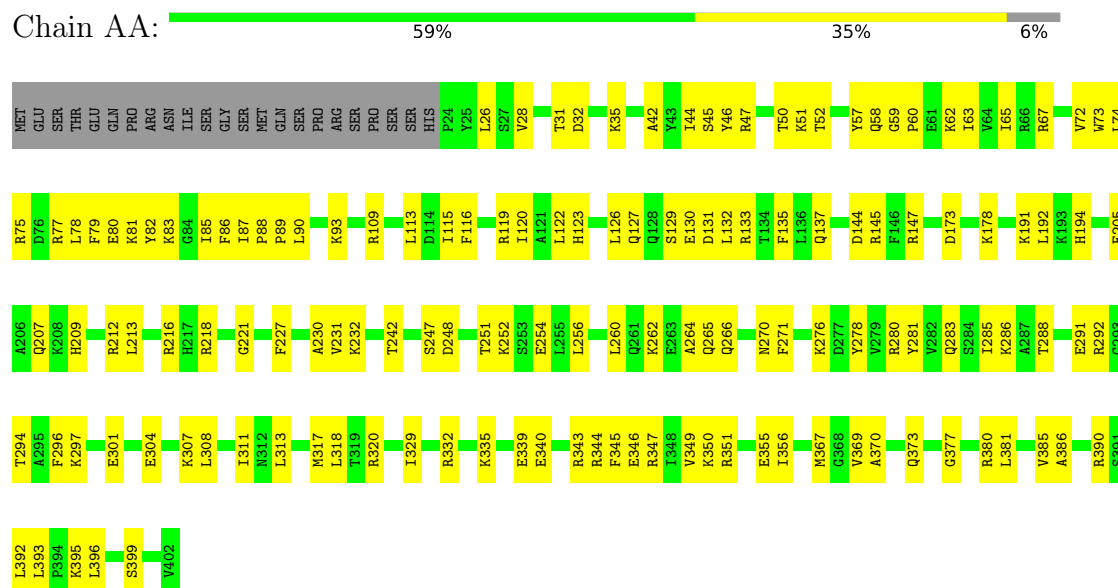
• Molecule 1: Sorting nexin 1

Chain A: 61% 34% 6%

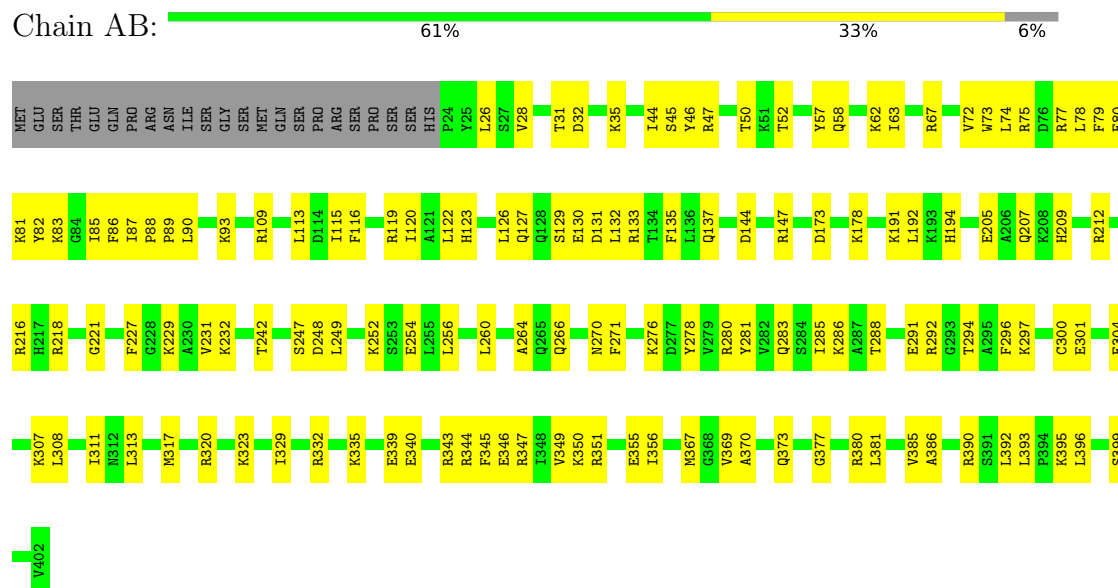




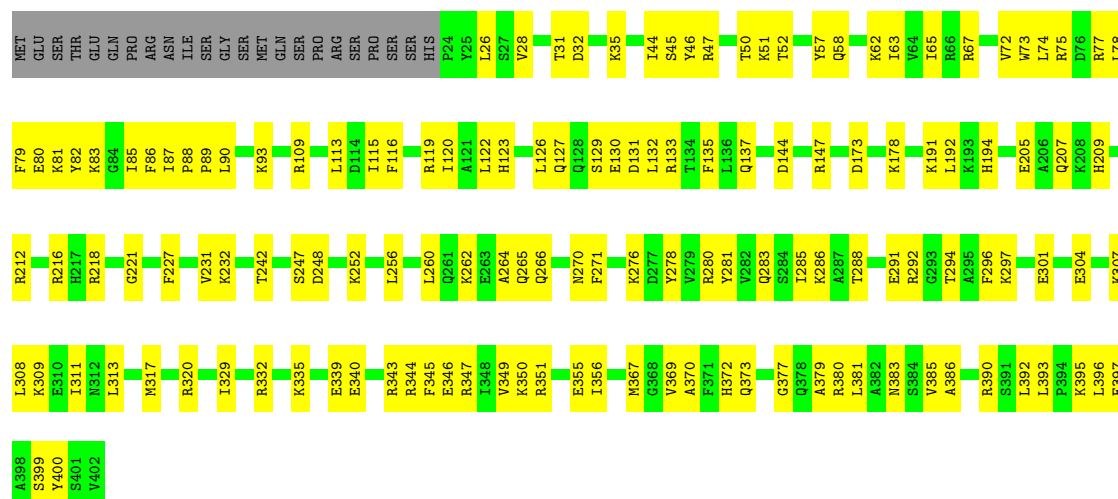
- Molecule 1: Sorting nexin 1



- Molecule 1: Sorting nexin 1

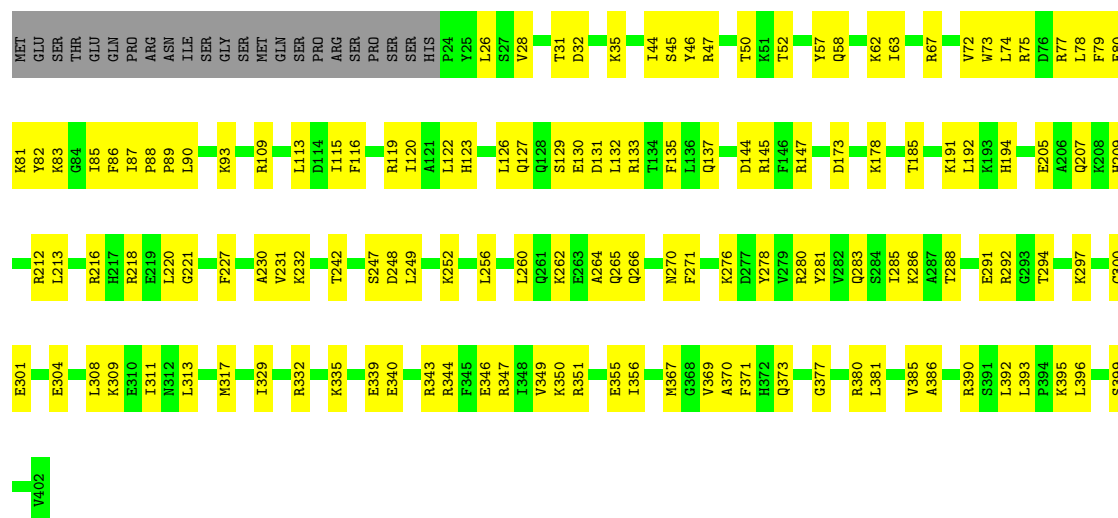


- Molecule 1: Sorting nexin 1



• Molecule 1: Sorting nexin 1

Chain C: 60% 34% 6%



• Molecule 1: Sorting nexin 1

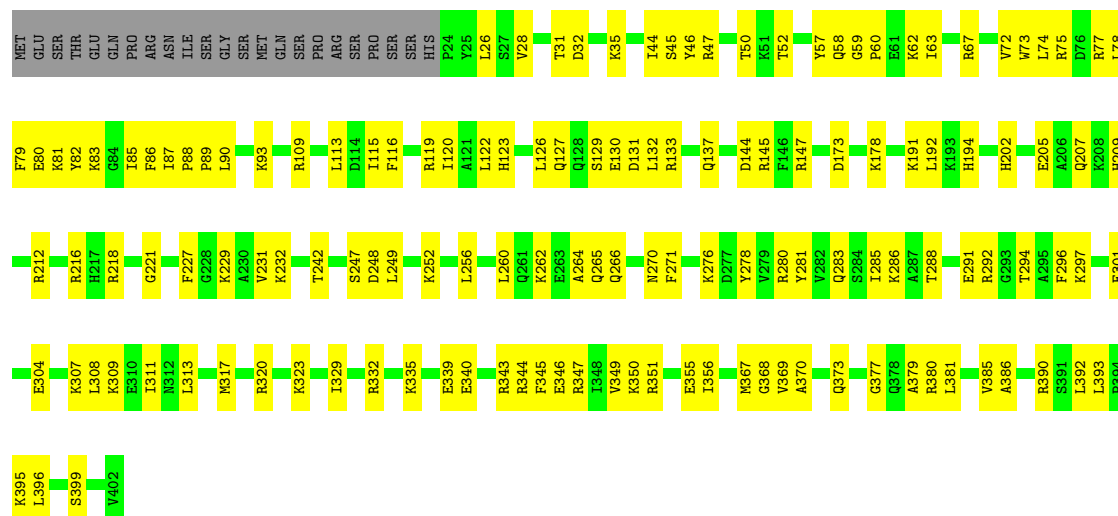
Chain D: 62% 32% 6%





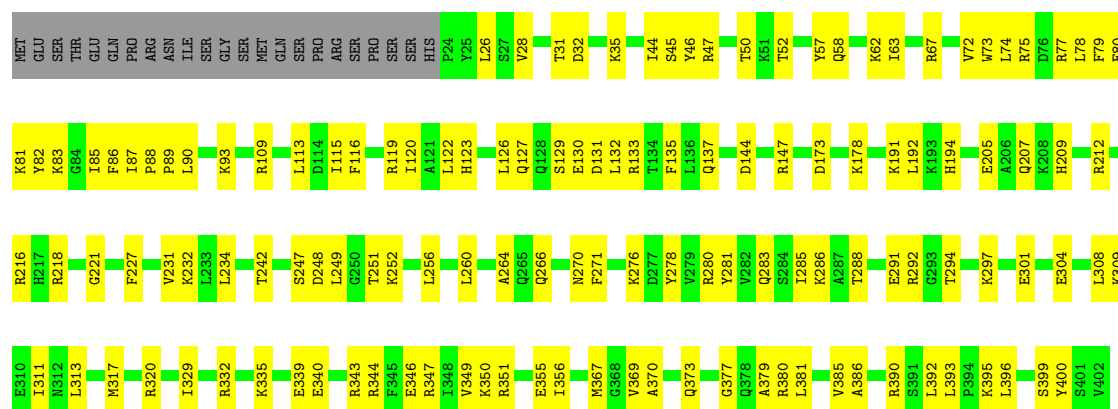
• Molecule 1: Sorting nexin 1

Chain E: 59% 35% 6%



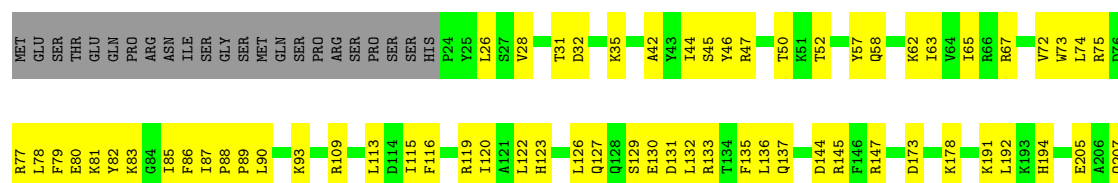
• Molecule 1: Sorting nexin 1

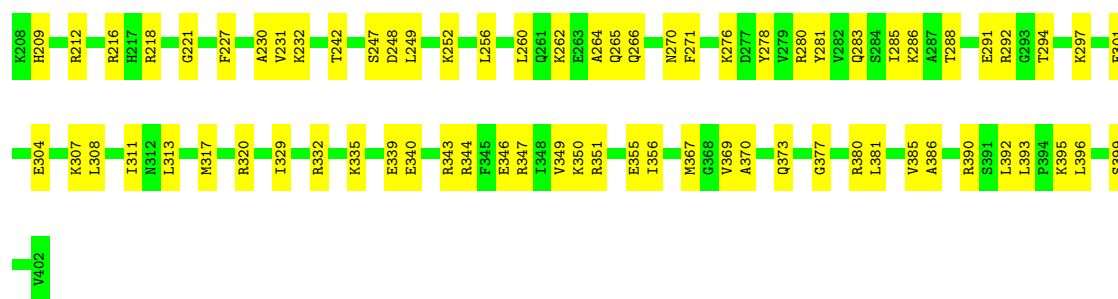
Chain F: 61% 33% 6%



• Molecule 1: Sorting nexin 1

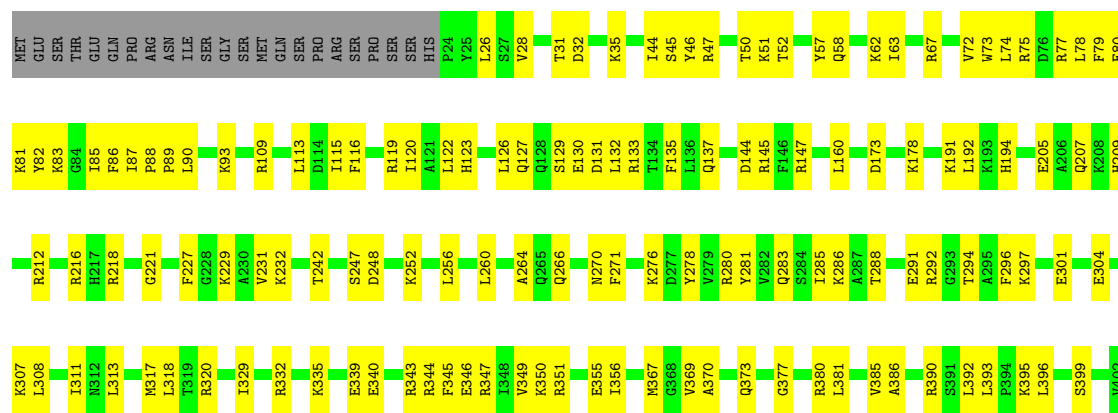
Chain G: 61% 34% 6%





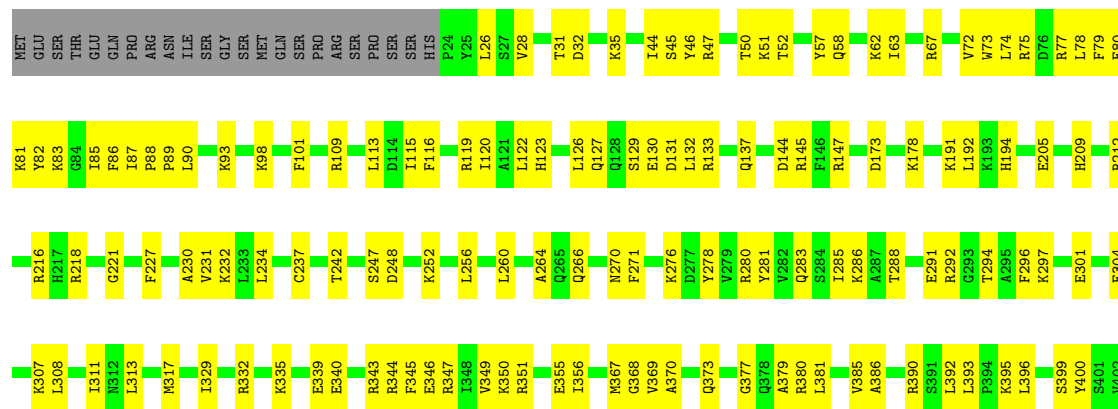
- Molecule 1: Sorting nexin 1

Chain H:



- Molecule 1: Sorting nexin 1

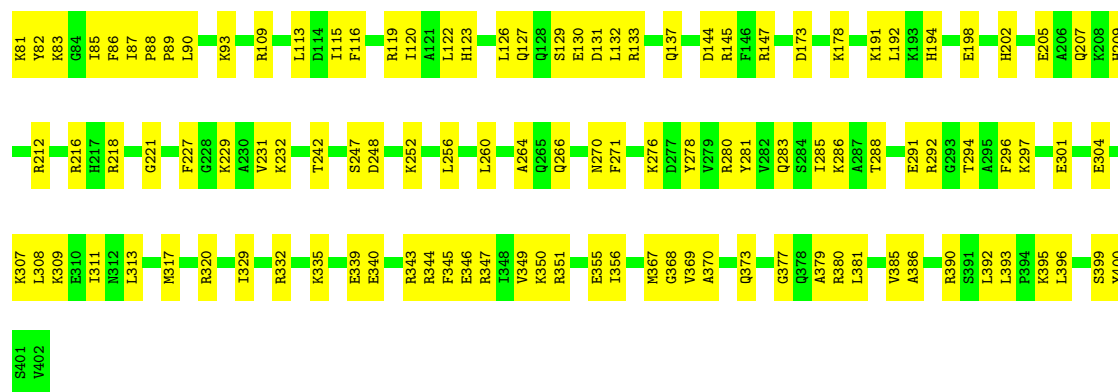
Chain I:



- Molecule 1: Sorting nexin 1

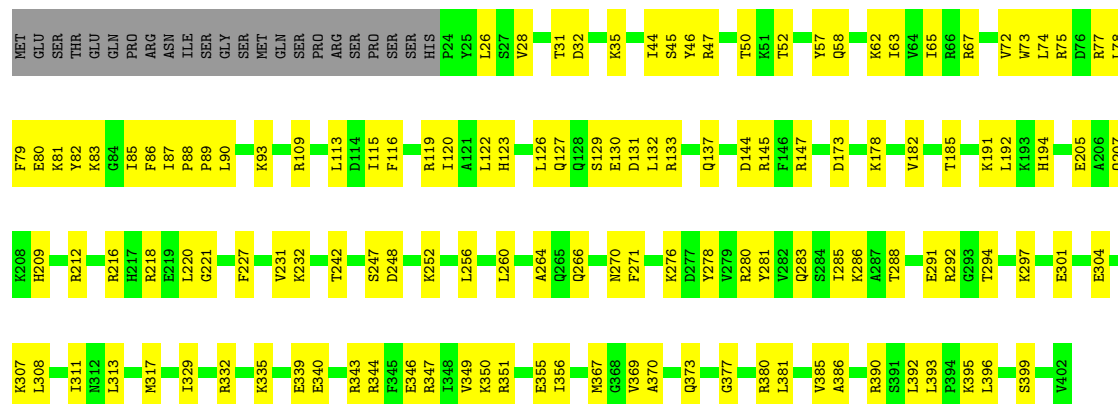
Chain J:





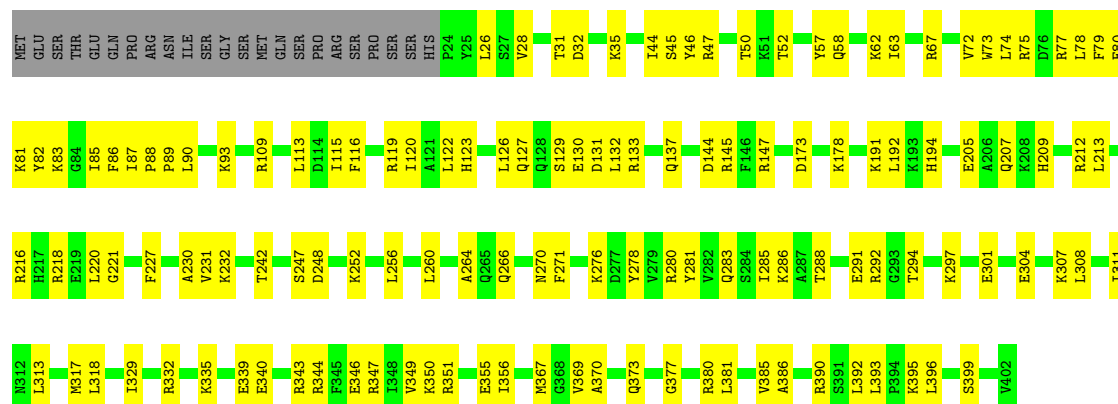
• Molecule 1: Sorting nexin 1

Chain K: 62% 32% 6%



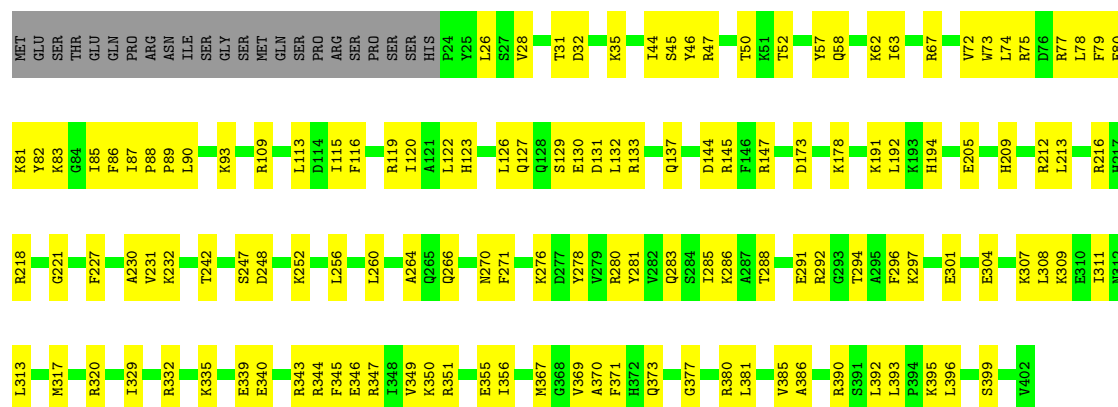
• Molecule 1: Sorting nexin 1

Chain L: 62% 32% 6%



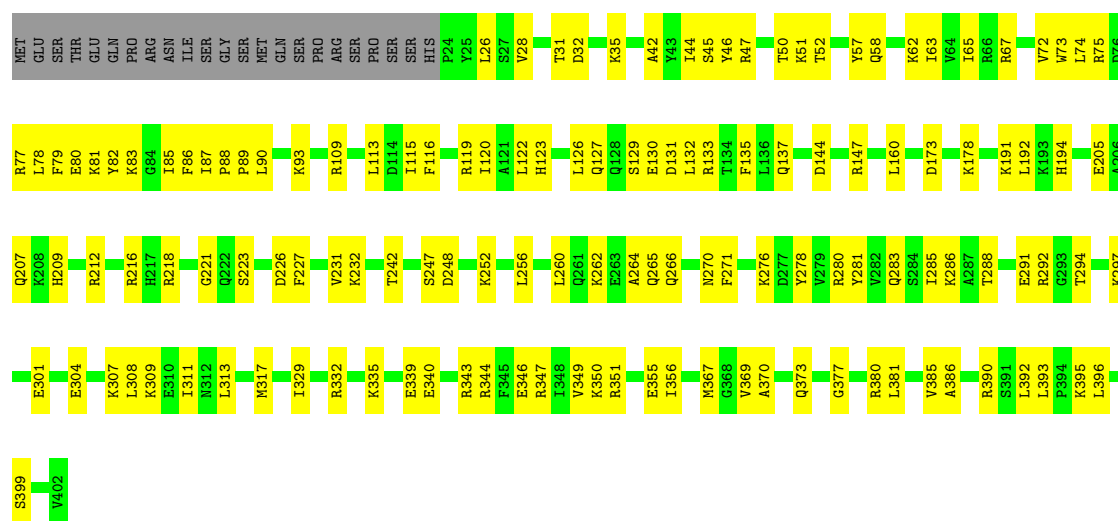
• Molecule 1: Sorting nexin 1

Chain M: 61% 33% 6%



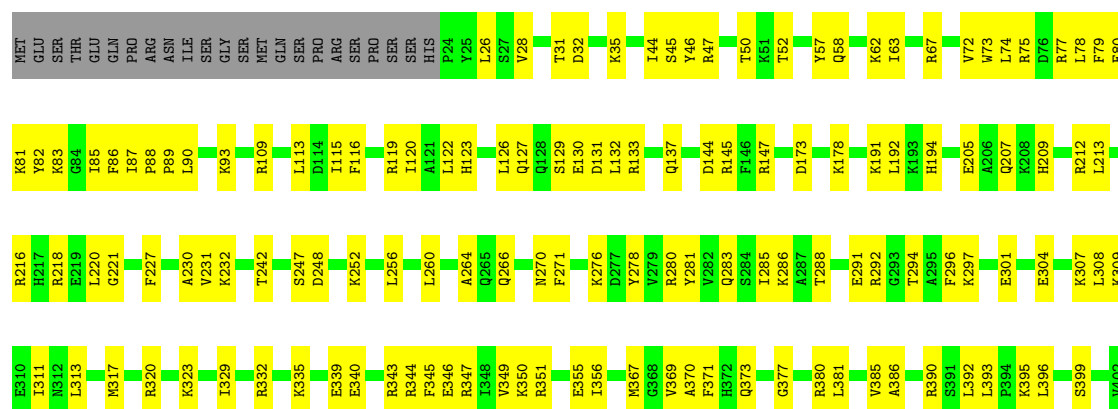
• Molecule 1: Sorting nexin 1

Chain N: 61% 34% 6%



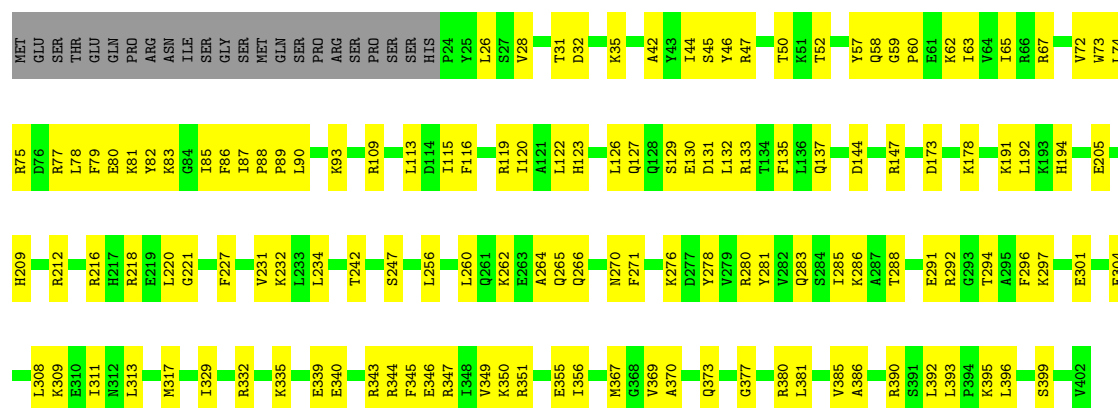
• Molecule 1: Sorting nexin 1

Chain O: 61% 34% 6%



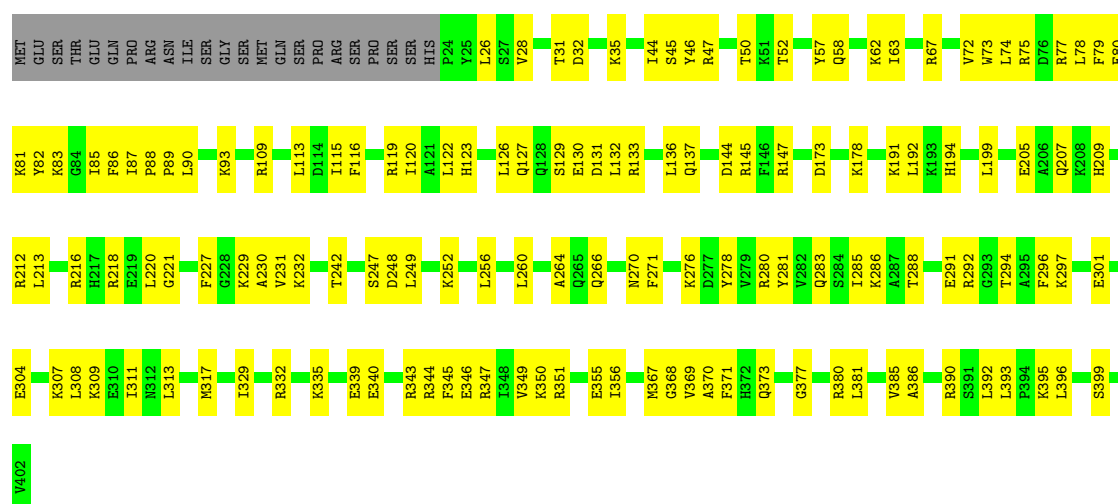
• Molecule 1: Sorting nexin 1

Chain P: 



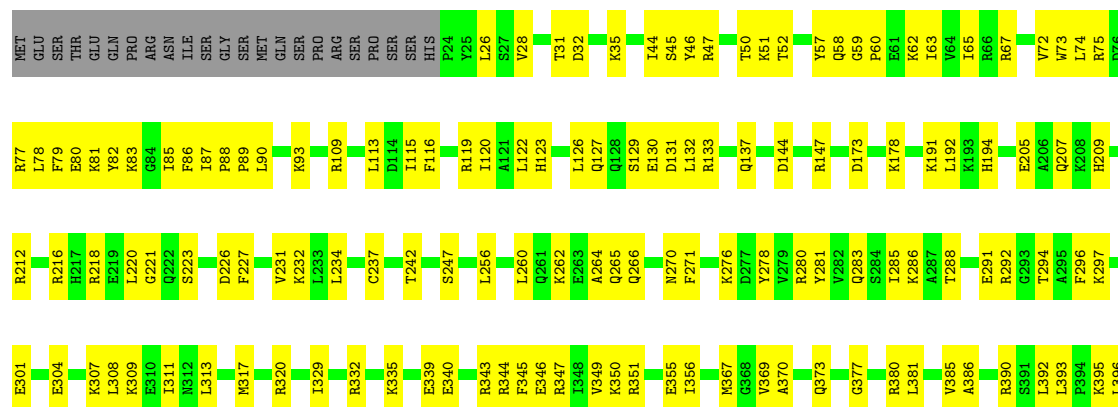
• Molecule 1: Sorting nexin 1

Chain Q: 



• Molecule 1: Sorting nexin 1

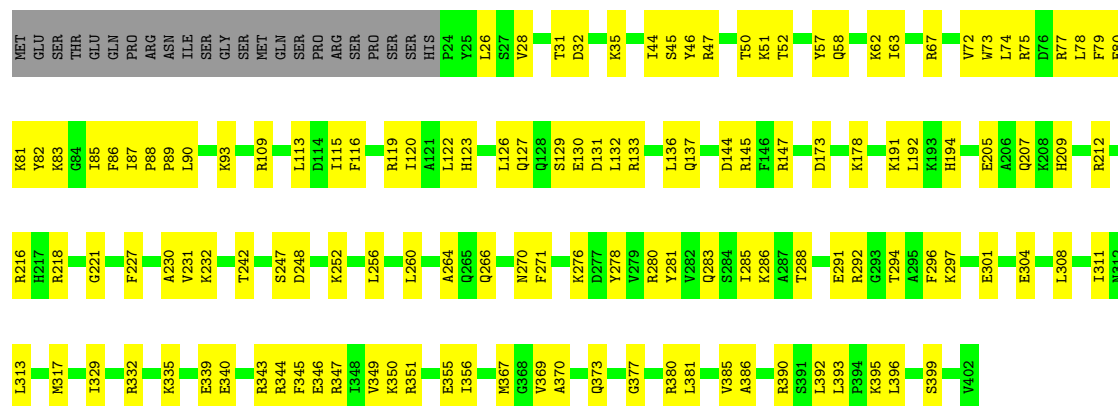
Chain R: 

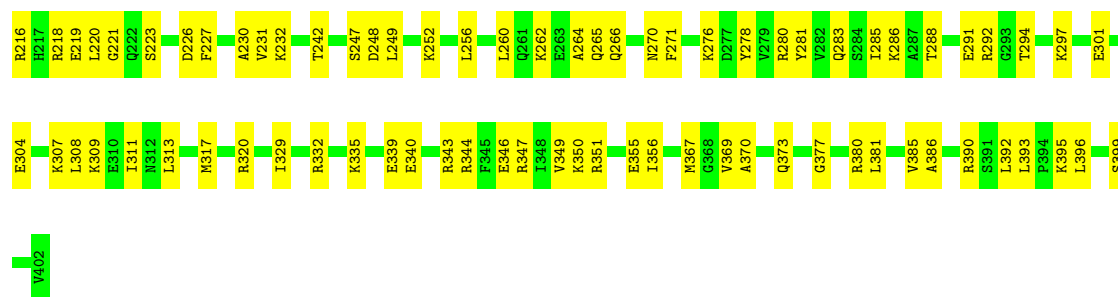




• Molecule 1: Sorting nexin 1

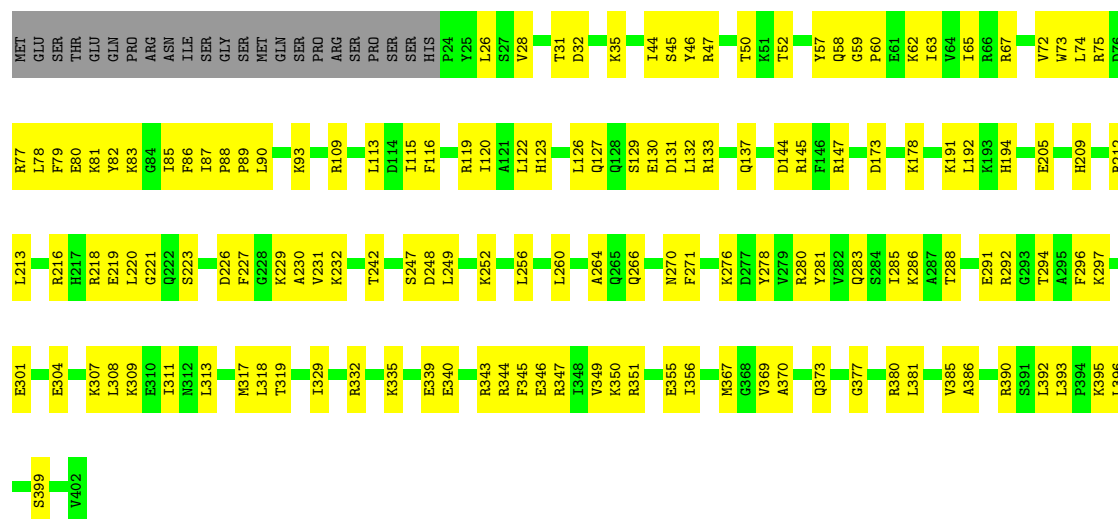
Chain S: 62% 32% 6%





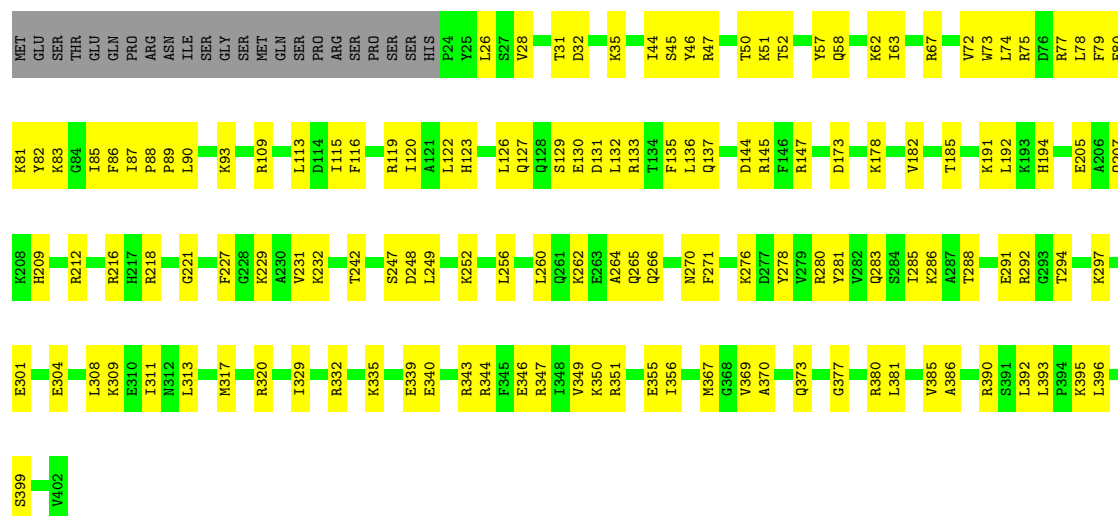
- Molecule 1: Sorting nexin 1

Chain V: 59% 35% 6%



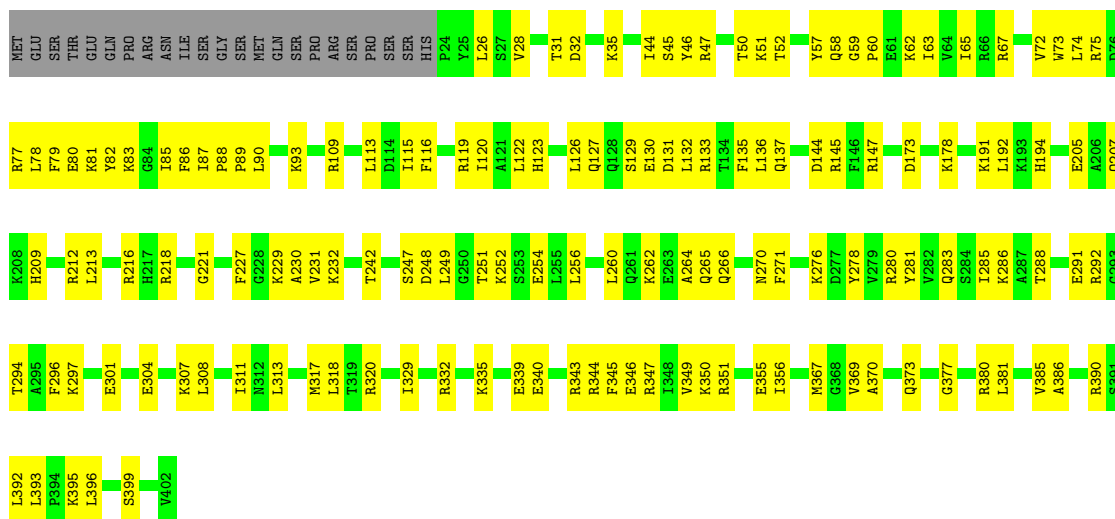
- Molecule 1: Sorting nexin 1

Chain W: 60% 34% 6%



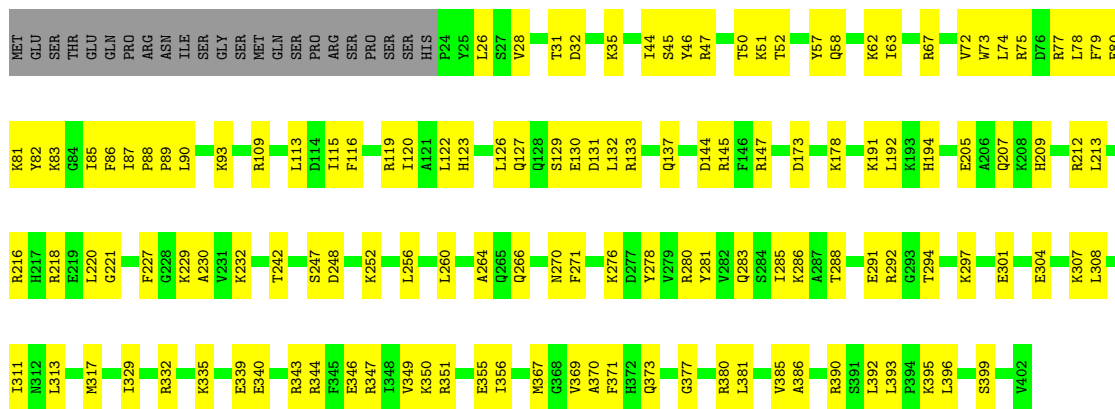
- Molecule 1: Sorting nexin 1

Chain X: 58% 36% 6%

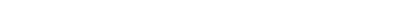


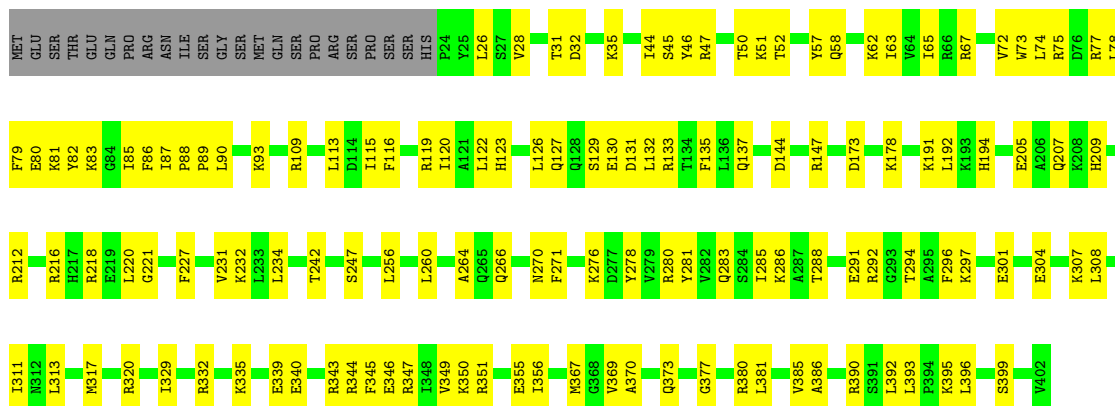
- Molecule 1: Sorting nexin 1

Chain Y:  62% 33% 6%



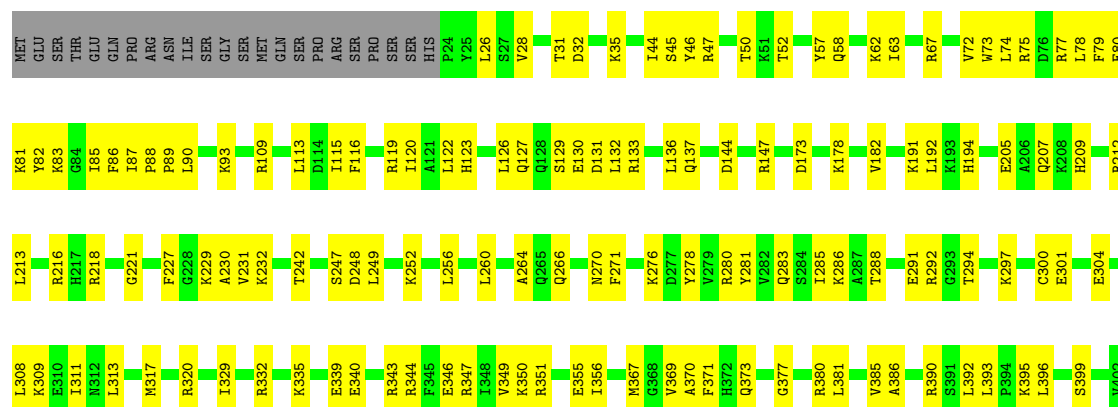
- Molecule 1: Sorting nexin 1

Chain Z:  62% 33% 6%



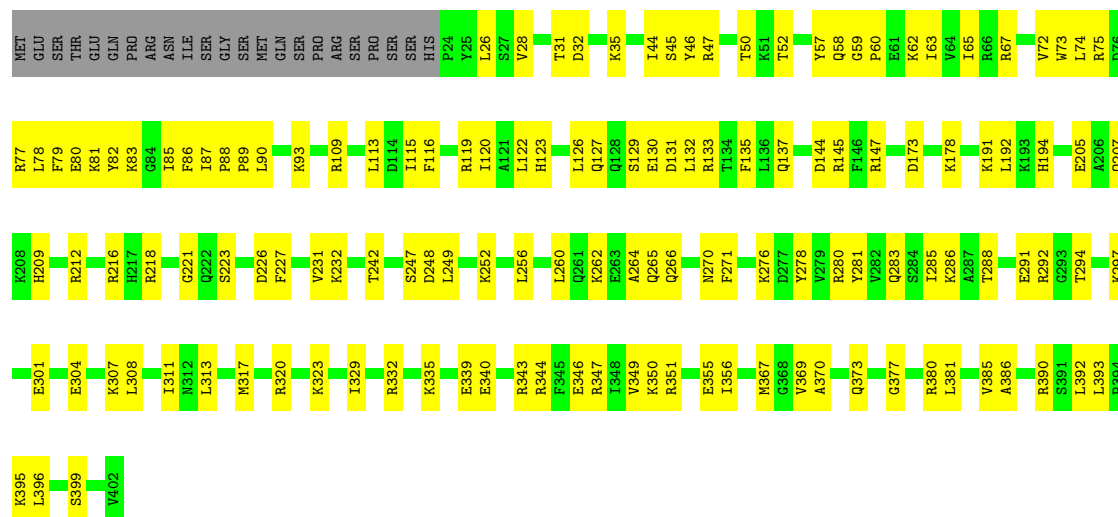
- Molecule 1: Sorting nexin 1

Chain a:  61% 33% 6%



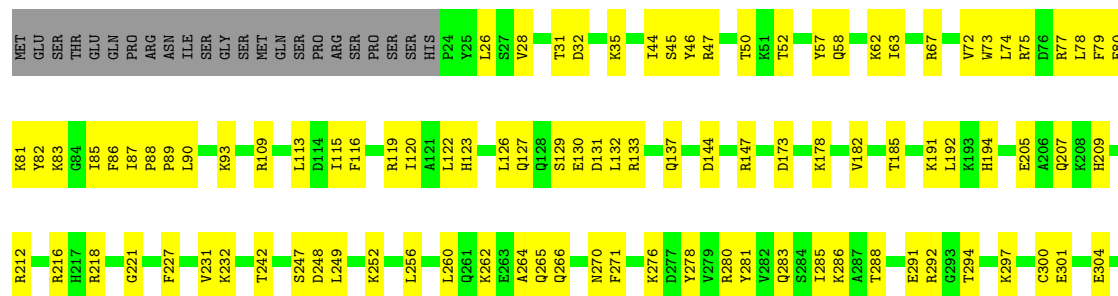
- Molecule 1: Sorting nexin 1

Chain b:  60% 34% 6%



- Molecule 1: Sorting nexin 1

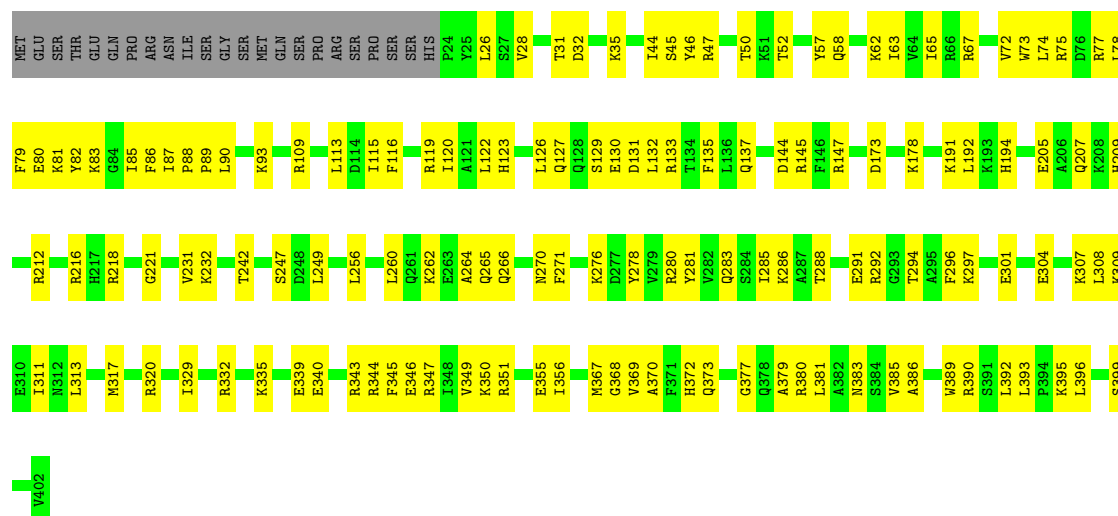
Chain c:  61% 34% 6%





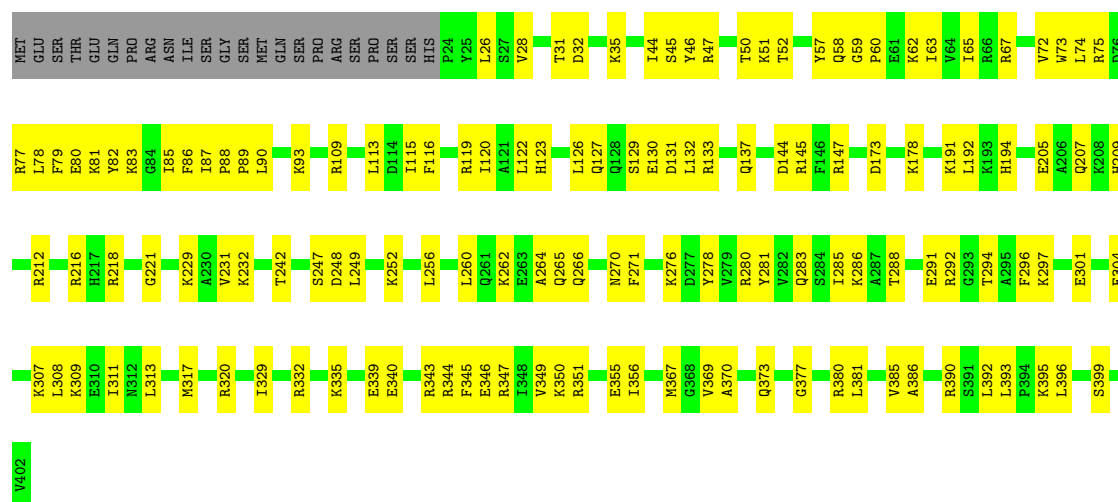
- Molecule 1: Sorting nexin 1

Chain d: 60% 34% 6%



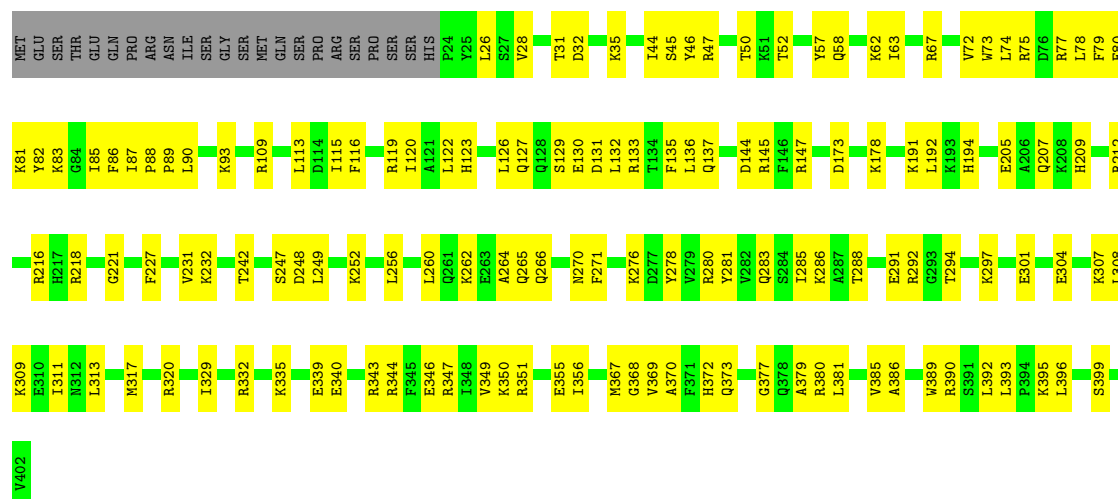
- Molecule 1: Sorting nexin 1

Chain e: 60% 34% 6%



- Molecule 1: Sorting nexin 1

Chain f: 60% 34% 6%



- Molecule 1: Sorting nexin 1

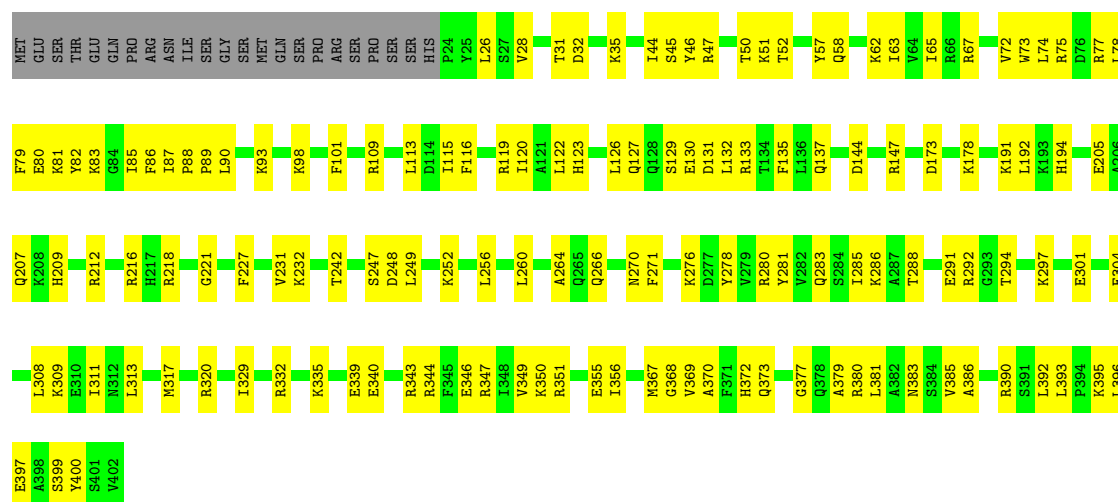


- Molecule 1: Sorting nexin 1



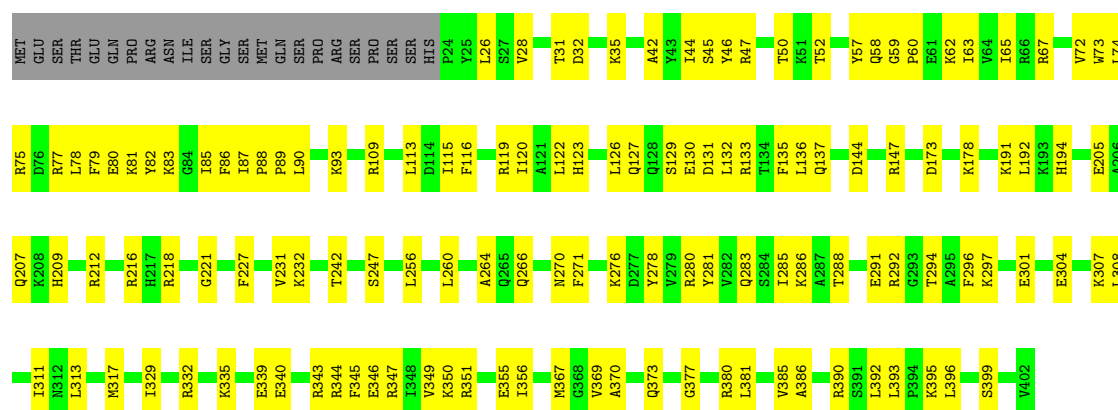
- Molecule 1: Sorting nexin 1

Chain i: 



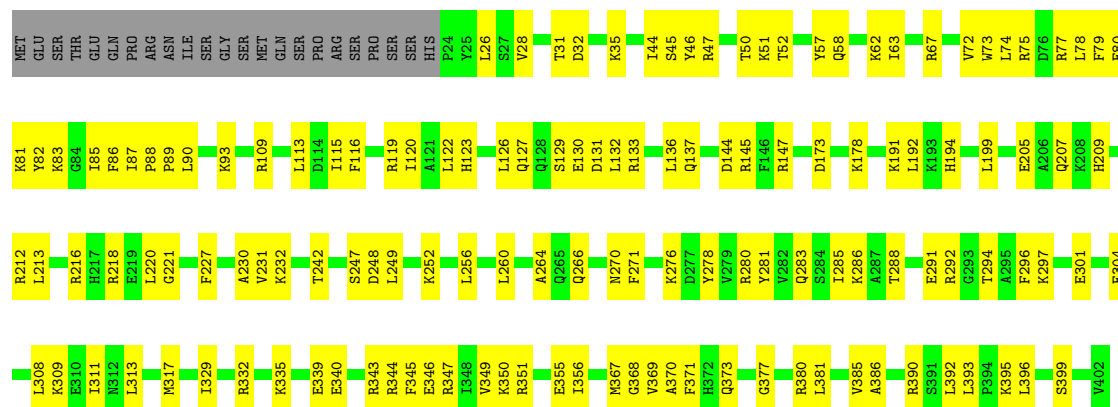
• Molecule 1: Sorting nexin 1

Chain j: 



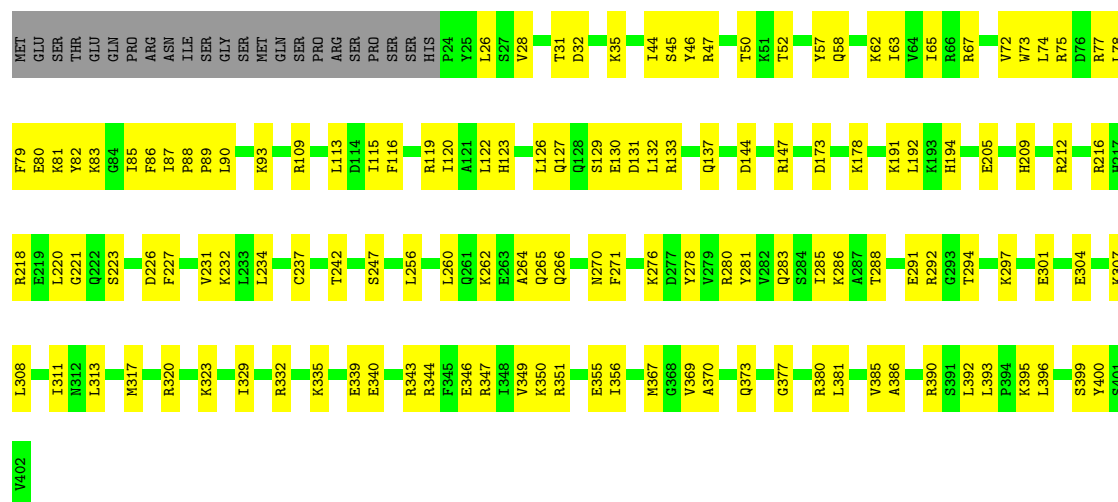
• Molecule 1: Sorting nexin 1

Chain k: 



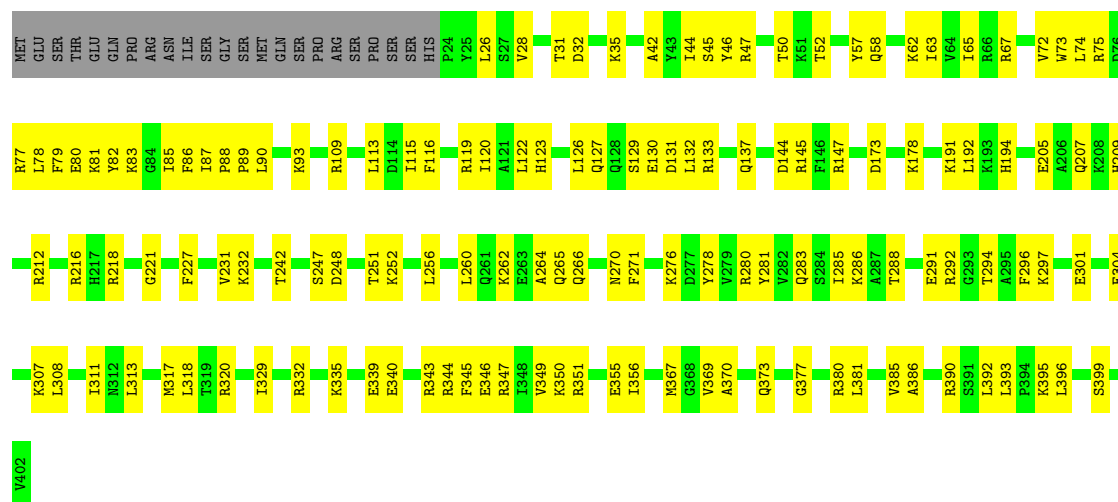
- Molecule 1: Sorting nexin 1

Chain l: 



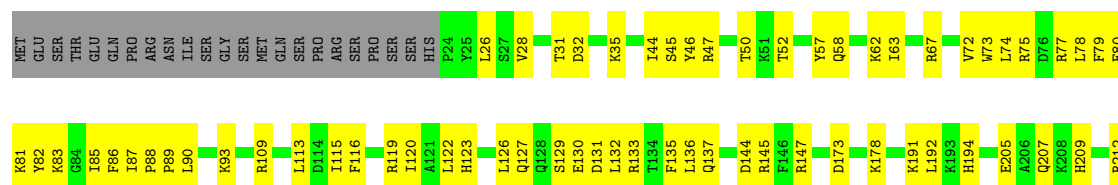
- Molecule 1: Sorting nexin 1

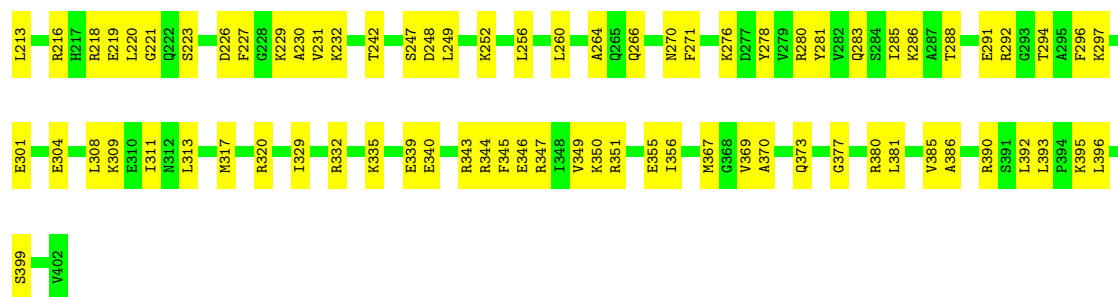
Chain m: 



- Molecule 1: Sorting nexin 1

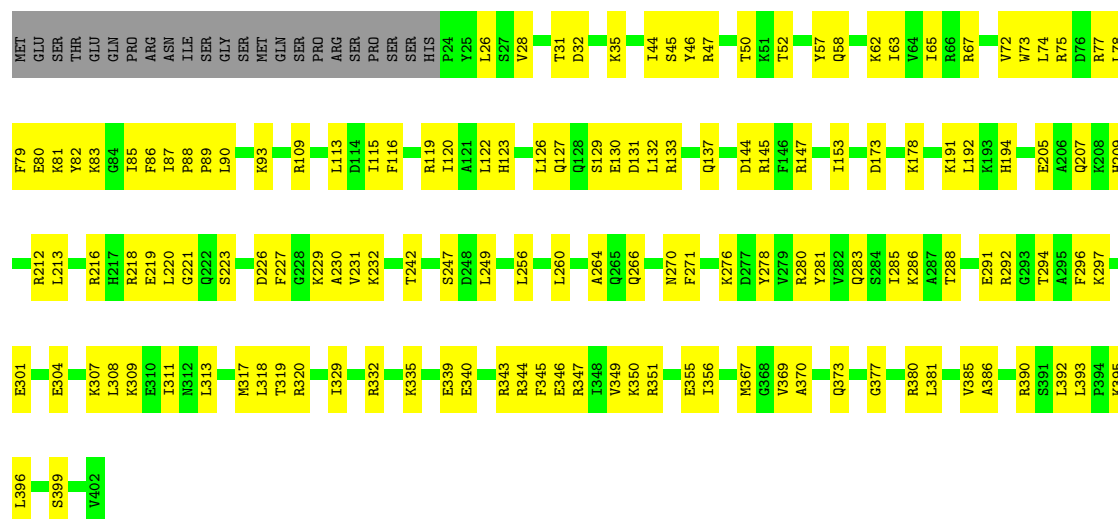
Chain n: 





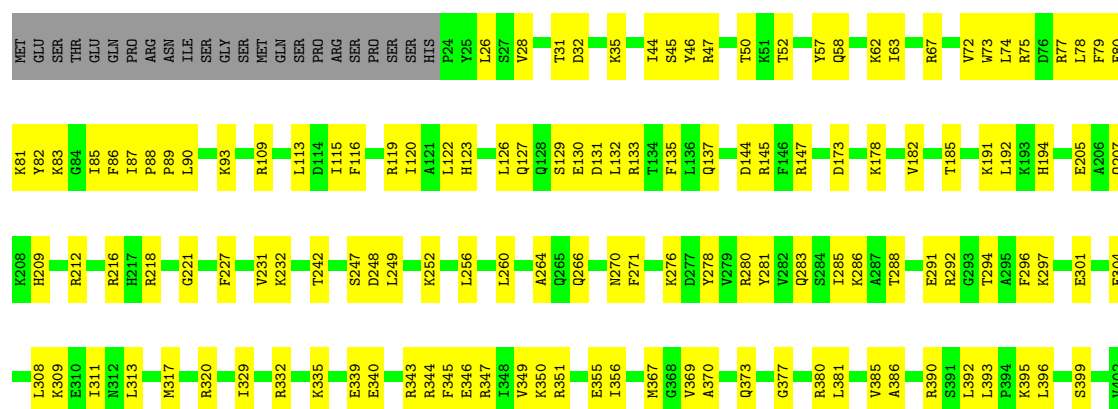
- Molecule 1: Sorting nexin 1

Chain o: 59% 35% 6%



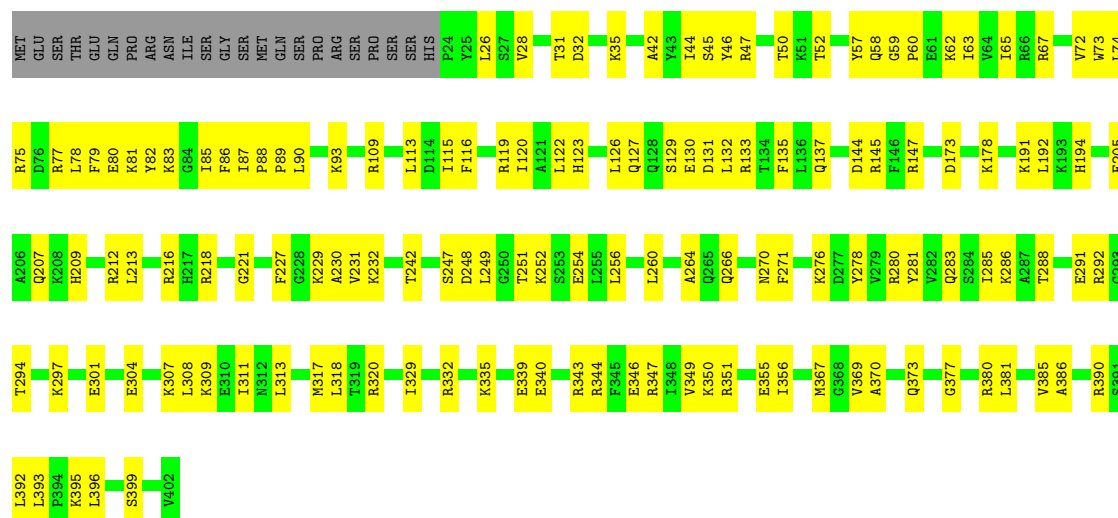
- Molecule 1: Sorting nexin 1

Chain p: 61% 33% 6%

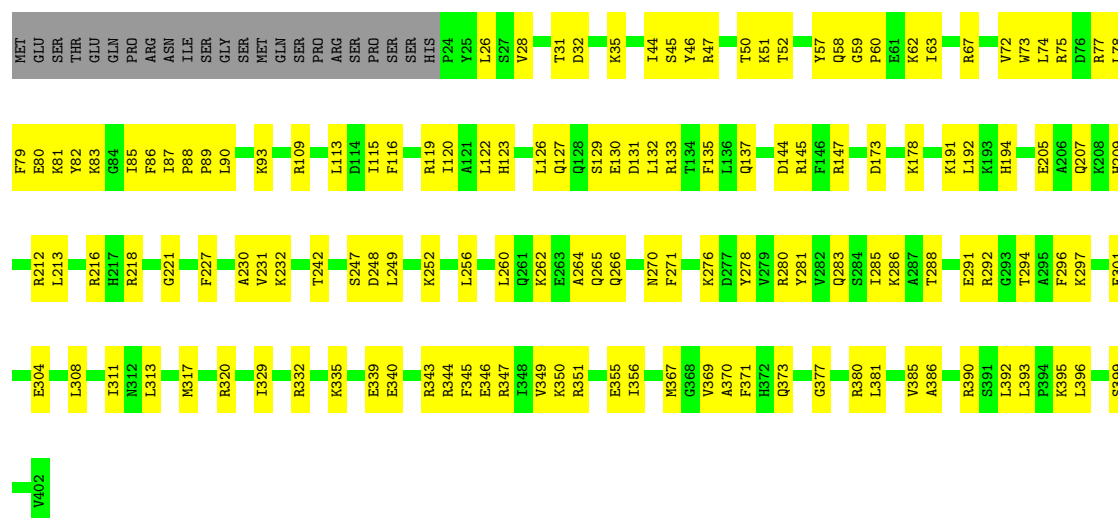


- Molecule 1: Sorting nexin 1

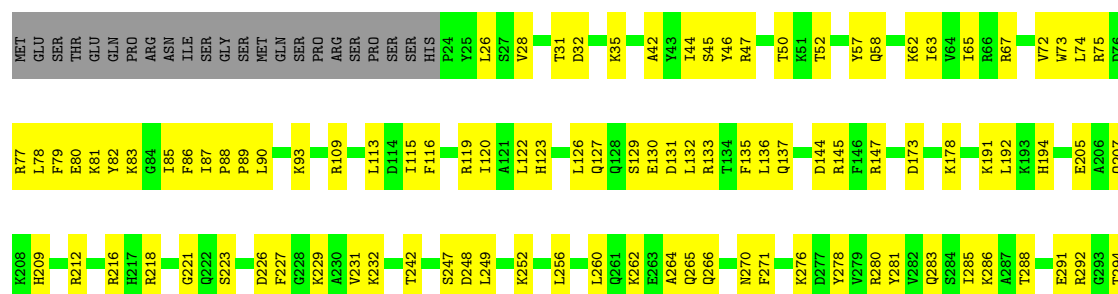
Chain q: 59% 35% 6%

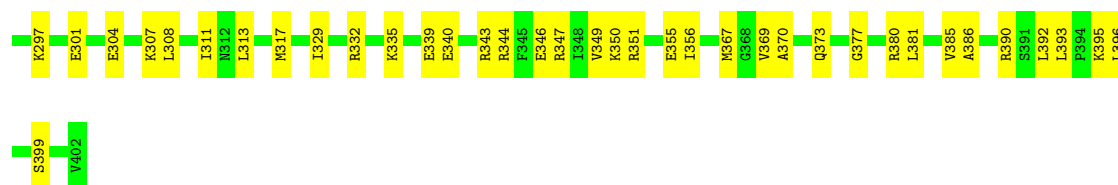


• Molecule 1: Sorting nexin 1



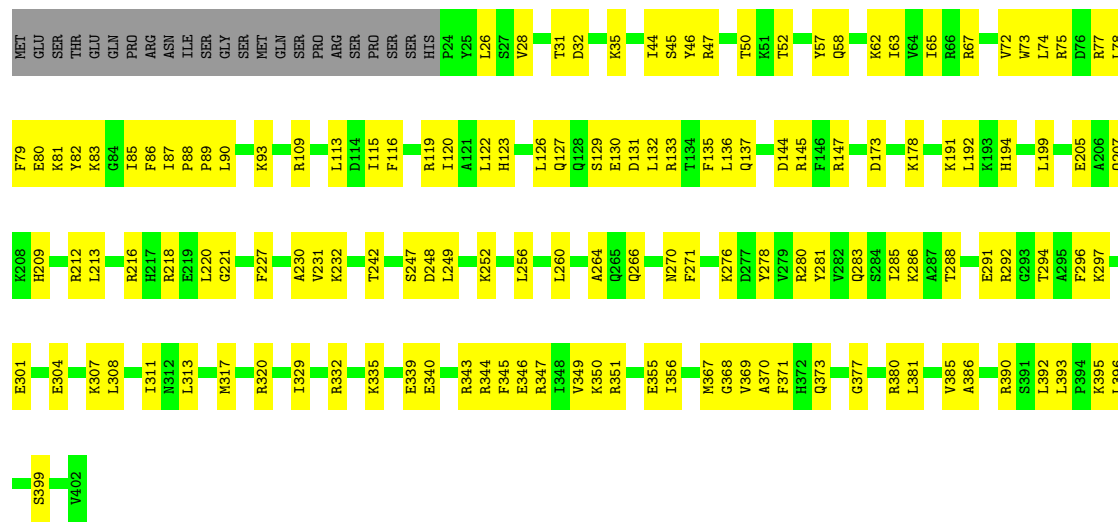
• Molecule 1: Sorting nexin 1





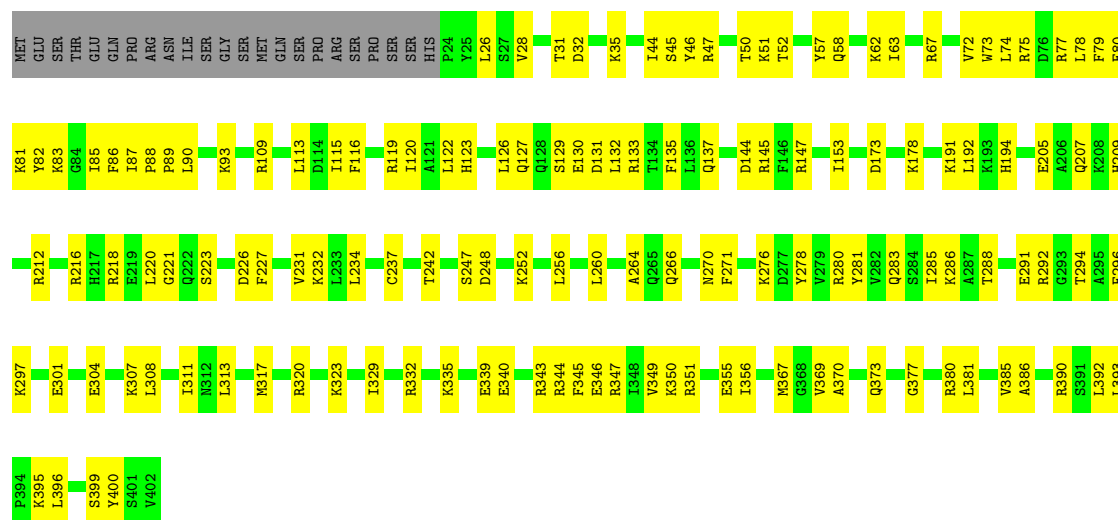
- Molecule 1: Sorting nexin 1

Chain t: 60% 35% 6%



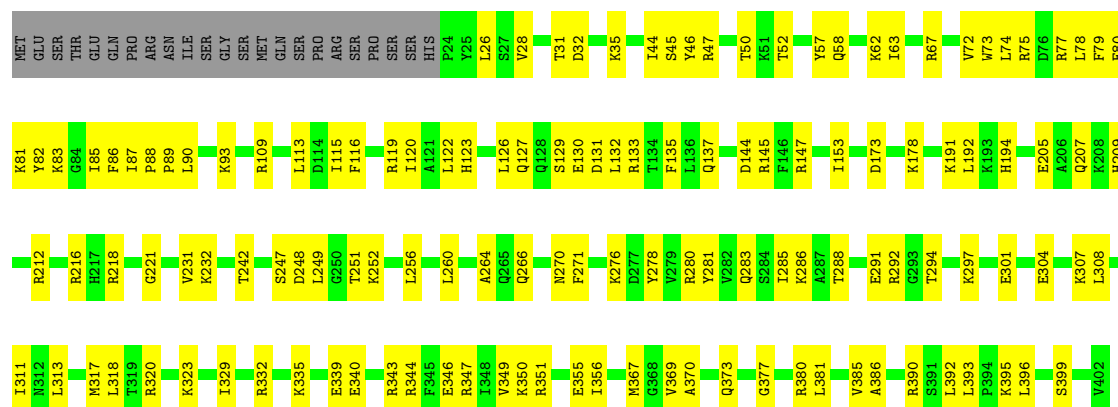
- Molecule 1: Sorting nexin 1

Chain u: 60% 35% 6%

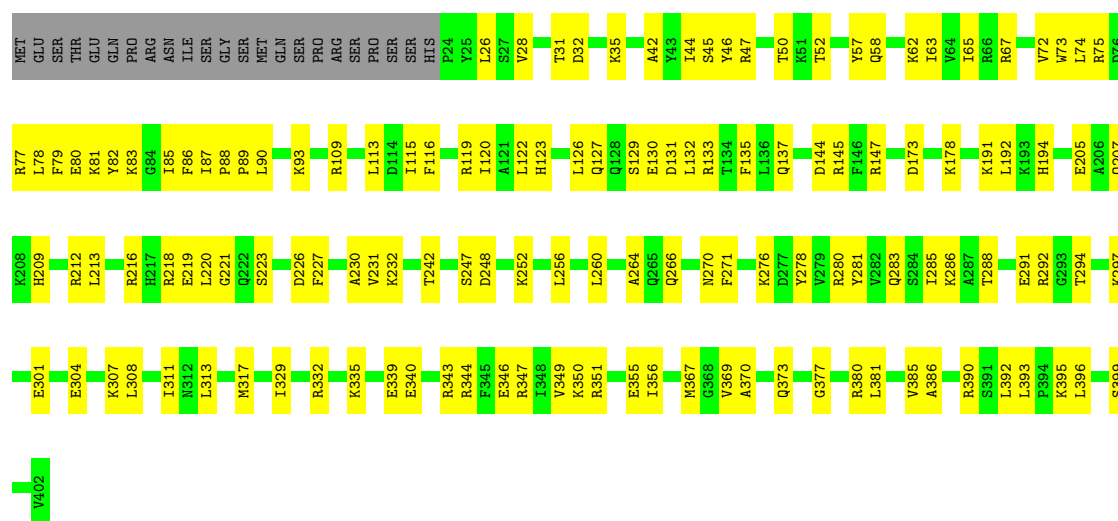


- Molecule 1: Sorting nexin 1

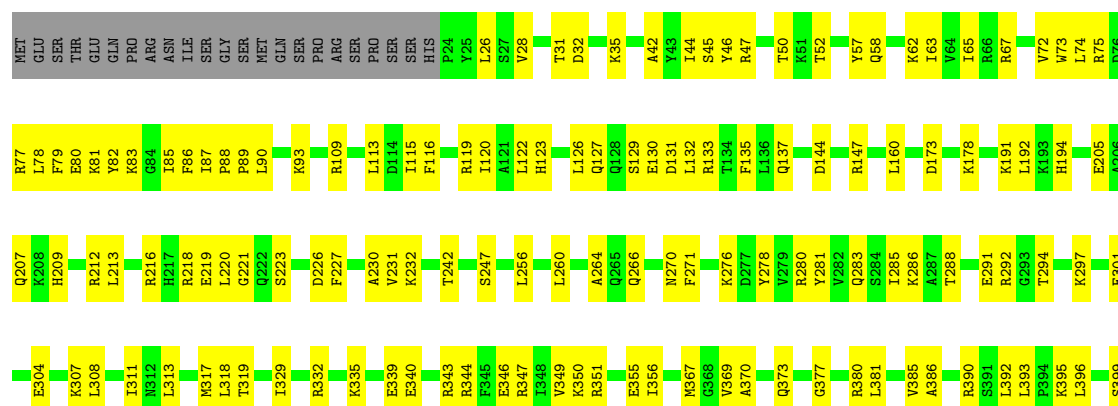
Chain v: 61% 33% 6%



• Molecule 1: Sorting nexin 1



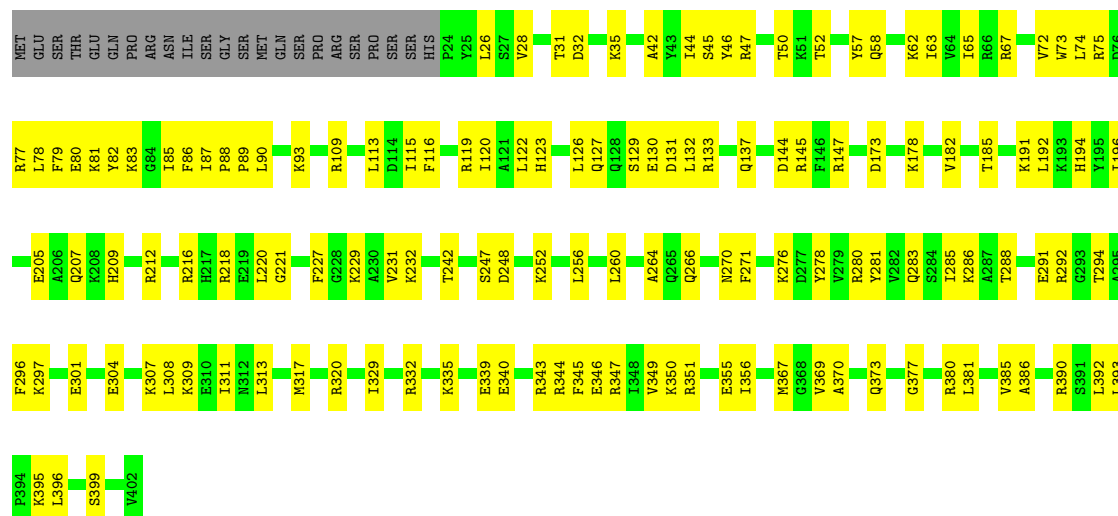
• Molecule 1: Sorting nexin 1





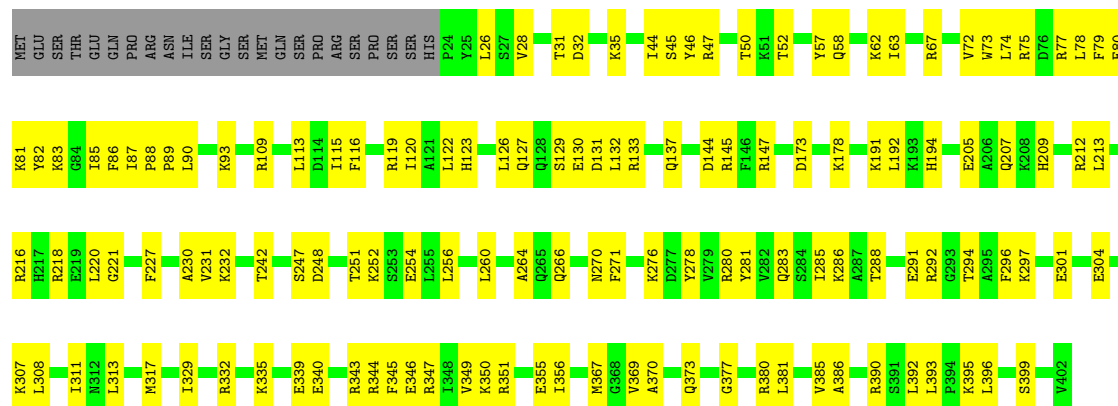
• Molecule 1: Sorting nexin 1

Chain y:  60% 34% 6%



• Molecule 1: Sorting nexin 1

Chain z:  61% 33% 6%



4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=-60.3°, rise=9.55 Å, axial sym=C1	Depositor
Number of segments used	20728	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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 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	0	0.11	0/3159	0.29	0/4242
1	1	0.11	0/3159	0.29	0/4242
1	2	0.11	0/3159	0.29	0/4242
1	3	0.11	0/3159	0.29	0/4242
1	4	0.11	0/3159	0.29	0/4242
1	5	0.11	0/3159	0.29	0/4242
1	6	0.11	0/3159	0.29	0/4242
1	7	0.11	0/3159	0.29	0/4242
1	8	0.11	0/3159	0.29	0/4242
1	9	0.11	0/3159	0.29	0/4242
1	A	0.11	0/3159	0.29	0/4242
1	AA	0.11	0/3159	0.29	0/4242
1	AB	0.11	0/3159	0.29	0/4242
1	B	0.11	0/3159	0.29	0/4242
1	C	0.11	0/3159	0.29	0/4242
1	D	0.11	0/3159	0.29	0/4242
1	E	0.11	0/3159	0.29	0/4242
1	F	0.11	0/3159	0.29	0/4242
1	G	0.11	0/3159	0.29	0/4242
1	H	0.11	0/3159	0.29	0/4242
1	I	0.11	0/3159	0.29	0/4242
1	J	0.11	0/3159	0.29	0/4242
1	K	0.11	0/3159	0.29	0/4242
1	L	0.11	0/3159	0.29	0/4242
1	M	0.11	0/3159	0.29	0/4242
1	N	0.11	0/3159	0.29	0/4242
1	O	0.11	0/3159	0.29	0/4242
1	P	0.11	0/3159	0.29	0/4242
1	Q	0.11	0/3159	0.29	0/4242
1	R	0.11	0/3159	0.29	0/4242
1	S	0.11	0/3159	0.29	0/4242
1	T	0.11	0/3159	0.29	0/4242
1	U	0.11	0/3159	0.29	0/4242
1	V	0.11	0/3159	0.29	0/4242

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	W	0.11	0/3159	0.29	0/4242
1	X	0.11	0/3159	0.29	0/4242
1	Y	0.11	0/3159	0.29	0/4242
1	Z	0.11	0/3159	0.29	0/4242
1	a	0.11	0/3159	0.29	0/4242
1	b	0.11	0/3159	0.29	0/4242
1	c	0.11	0/3159	0.29	0/4242
1	d	0.11	0/3159	0.29	0/4242
1	e	0.11	0/3159	0.29	0/4242
1	f	0.11	0/3159	0.29	0/4242
1	g	0.11	0/3159	0.29	0/4242
1	h	0.11	0/3159	0.29	0/4242
1	i	0.11	0/3159	0.29	0/4242
1	j	0.11	0/3159	0.29	0/4242
1	k	0.11	0/3159	0.29	0/4242
1	l	0.11	0/3159	0.29	0/4242
1	m	0.11	0/3159	0.29	0/4242
1	n	0.11	0/3159	0.29	0/4242
1	o	0.11	0/3159	0.29	0/4242
1	p	0.11	0/3159	0.29	0/4242
1	q	0.11	0/3159	0.29	0/4242
1	r	0.11	0/3159	0.29	0/4242
1	s	0.11	0/3159	0.29	0/4242
1	t	0.11	0/3159	0.29	0/4242
1	u	0.11	0/3159	0.29	0/4242
1	v	0.11	0/3159	0.29	0/4242
1	w	0.11	0/3159	0.29	0/4242
1	x	0.11	0/3159	0.29	0/4242
1	y	0.11	0/3159	0.29	0/4242
1	z	0.11	0/3159	0.29	0/4242
All	All	0.11	0/202176	0.29	0/271488

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	0	3102	0	3129	101	0
1	1	3102	0	3129	98	0
1	2	3102	0	3129	100	0
1	3	3102	0	3129	106	0
1	4	3102	0	3129	98	0
1	5	3102	0	3129	106	0
1	6	3102	0	3129	109	0
1	7	3102	0	3129	91	0
1	8	3102	0	3129	100	0
1	9	3102	0	3129	89	0
1	A	3102	0	3129	106	0
1	AA	3102	0	3129	108	0
1	AB	3102	0	3129	98	0
1	B	3102	0	3129	107	0
1	C	3102	0	3129	99	0
1	D	3102	0	3129	96	0
1	E	3102	0	3129	102	0
1	F	3102	0	3129	98	0
1	G	3102	0	3129	96	0
1	H	3102	0	3129	100	0
1	I	3102	0	3129	109	0
1	J	3102	0	3129	109	0
1	K	3102	0	3129	95	0
1	L	3102	0	3129	93	0
1	M	3102	0	3129	96	0
1	N	3102	0	3129	101	0
1	O	3102	0	3129	97	0
1	P	3102	0	3129	97	0
1	Q	3102	0	3129	106	0
1	R	3102	0	3129	108	0
1	S	3102	0	3129	90	0
1	T	3102	0	3129	99	0
1	U	3102	0	3129	114	0
1	V	3102	0	3129	115	0
1	W	3102	0	3129	103	0
1	X	3102	0	3129	105	0
1	Y	3102	0	3129	93	0
1	Z	3102	0	3129	97	0
1	a	3102	0	3129	101	0
1	b	3102	0	3129	105	0
1	c	3102	0	3129	113	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	d	3102	0	3129	108	0
1	e	3102	0	3129	96	0
1	f	3102	0	3129	104	0
1	g	3102	0	3129	100	0
1	h	3102	0	3129	111	0
1	i	3102	0	3129	118	0
1	j	3102	0	3129	91	0
1	k	3102	0	3129	107	0
1	l	3102	0	3129	106	0
1	m	3102	0	3129	98	0
1	n	3102	0	3129	118	0
1	o	3102	0	3129	121	0
1	p	3102	0	3129	100	0
1	q	3102	0	3129	101	0
1	r	3102	0	3129	102	0
1	s	3102	0	3129	98	0
1	t	3102	0	3129	109	0
1	u	3102	0	3129	107	0
1	v	3102	0	3129	96	0
1	w	3102	0	3129	111	0
1	x	3102	0	3129	118	0
1	y	3102	0	3129	103	0
1	z	3102	0	3129	95	0
All	All	198528	0	200256	6191	0

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 16.

The worst 5 of 6191 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:h:379:ALA:HB1	1:i:390:ARG:HH11	0.94	1.10
1:3:300:CYS:HB3	1:x:318:LEU:HD21	1.31	1.04
1:E:379:ALA:HB1	1:F:390:ARG:HH11	1.17	1.02
1:f:379:ALA:HB1	1:g:390:ARG:HH11	1.20	1.02
1:A:379:ALA:HB1	1:B:390:ARG:HH11	1.23	1.02

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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 EM 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	0	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	1	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	2	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	3	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	4	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	5	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	6	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	7	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	8	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	9	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	A	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	AA	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	AB	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	B	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	C	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	D	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	E	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	F	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	G	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	H	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	I	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	J	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	K	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	L	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	M	377/402 (94%)	372 (99%)	5 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	N	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	O	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	P	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	Q	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	R	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	S	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	T	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	U	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	V	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	W	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	X	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	Y	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	Z	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	a	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	b	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	c	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	d	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	e	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	f	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	g	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	h	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	i	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	j	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	k	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	l	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	m	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	n	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	o	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	p	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	q	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	r	377/402 (94%)	372 (99%)	5 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	s	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	t	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	u	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	v	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	w	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	x	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	y	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
1	z	377/402 (94%)	372 (99%)	5 (1%)	0	100	100
All	All	24128/25728 (94%)	23808 (99%)	320 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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 EM 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	0	335/357 (94%)	335 (100%)	0	100	100
1	1	335/357 (94%)	335 (100%)	0	100	100
1	2	335/357 (94%)	335 (100%)	0	100	100
1	3	335/357 (94%)	335 (100%)	0	100	100
1	4	335/357 (94%)	335 (100%)	0	100	100
1	5	335/357 (94%)	335 (100%)	0	100	100
1	6	335/357 (94%)	335 (100%)	0	100	100
1	7	335/357 (94%)	335 (100%)	0	100	100
1	8	335/357 (94%)	335 (100%)	0	100	100
1	9	335/357 (94%)	335 (100%)	0	100	100
1	A	335/357 (94%)	335 (100%)	0	100	100
1	AA	335/357 (94%)	335 (100%)	0	100	100
1	AB	335/357 (94%)	335 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	335/357 (94%)	335 (100%)	0	100	100
1	C	335/357 (94%)	335 (100%)	0	100	100
1	D	335/357 (94%)	335 (100%)	0	100	100
1	E	335/357 (94%)	335 (100%)	0	100	100
1	F	335/357 (94%)	335 (100%)	0	100	100
1	G	335/357 (94%)	335 (100%)	0	100	100
1	H	335/357 (94%)	335 (100%)	0	100	100
1	I	335/357 (94%)	335 (100%)	0	100	100
1	J	335/357 (94%)	335 (100%)	0	100	100
1	K	335/357 (94%)	335 (100%)	0	100	100
1	L	335/357 (94%)	335 (100%)	0	100	100
1	M	335/357 (94%)	335 (100%)	0	100	100
1	N	335/357 (94%)	335 (100%)	0	100	100
1	O	335/357 (94%)	335 (100%)	0	100	100
1	P	335/357 (94%)	335 (100%)	0	100	100
1	Q	335/357 (94%)	335 (100%)	0	100	100
1	R	335/357 (94%)	335 (100%)	0	100	100
1	S	335/357 (94%)	335 (100%)	0	100	100
1	T	335/357 (94%)	335 (100%)	0	100	100
1	U	335/357 (94%)	335 (100%)	0	100	100
1	V	335/357 (94%)	335 (100%)	0	100	100
1	W	335/357 (94%)	335 (100%)	0	100	100
1	X	335/357 (94%)	335 (100%)	0	100	100
1	Y	335/357 (94%)	335 (100%)	0	100	100
1	Z	335/357 (94%)	335 (100%)	0	100	100
1	a	335/357 (94%)	335 (100%)	0	100	100
1	b	335/357 (94%)	335 (100%)	0	100	100
1	c	335/357 (94%)	335 (100%)	0	100	100
1	d	335/357 (94%)	335 (100%)	0	100	100
1	e	335/357 (94%)	335 (100%)	0	100	100
1	f	335/357 (94%)	335 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	g	335/357 (94%)	335 (100%)	0	100	100
1	h	335/357 (94%)	335 (100%)	0	100	100
1	i	335/357 (94%)	335 (100%)	0	100	100
1	j	335/357 (94%)	335 (100%)	0	100	100
1	k	335/357 (94%)	335 (100%)	0	100	100
1	l	335/357 (94%)	335 (100%)	0	100	100
1	m	335/357 (94%)	335 (100%)	0	100	100
1	n	335/357 (94%)	335 (100%)	0	100	100
1	o	335/357 (94%)	335 (100%)	0	100	100
1	p	335/357 (94%)	335 (100%)	0	100	100
1	q	335/357 (94%)	335 (100%)	0	100	100
1	r	335/357 (94%)	335 (100%)	0	100	100
1	s	335/357 (94%)	335 (100%)	0	100	100
1	t	335/357 (94%)	335 (100%)	0	100	100
1	u	335/357 (94%)	335 (100%)	0	100	100
1	v	335/357 (94%)	335 (100%)	0	100	100
1	w	335/357 (94%)	335 (100%)	0	100	100
1	x	335/357 (94%)	335 (100%)	0	100	100
1	y	335/357 (94%)	335 (100%)	0	100	100
1	z	335/357 (94%)	335 (100%)	0	100	100
All	All	21440/22848 (94%)	21440 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 351 such sidechains are listed below:

Mol	Chain	Res	Type
1	e	162	GLN
1	o	261	GLN
1	f	207	GLN
1	j	378	GLN
1	r	201	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.