



wwPDB EM Validation Summary Report ⓘ

Mar 14, 2026 – 06:52 AM UTC

PDB ID : 9J1W / pdb_00009j1w
EMDB ID : EMD-61080
Title : Endogenous dihydrolipoamide acetyltransferase (E2) core of pyruvate dehydrogenase complex from pig heart
Authors : Wang, C.; Zhang, X.; Chang, Y.J.
Deposited on : 2024-08-05
Resolution : 3.20 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 : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

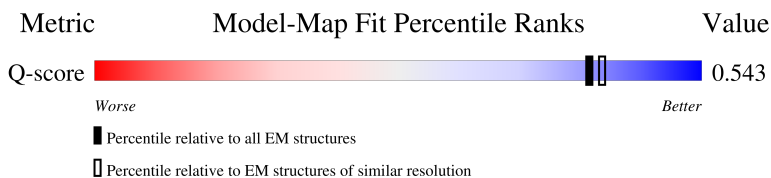
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.20 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.



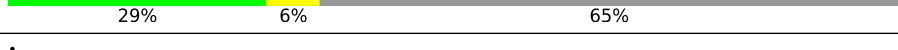
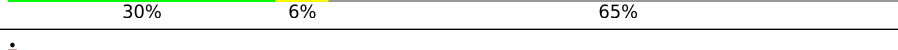
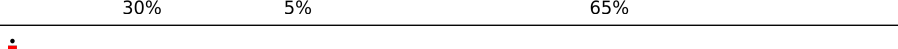
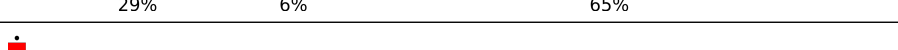
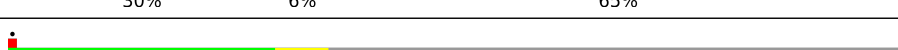
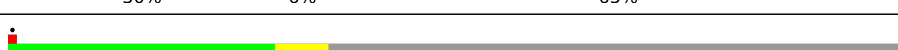
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15020 (2.70 - 3.70)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	647	 29% 6% 65%
1	AA	647	 30% 5% 65%
1	AB	647	 30% 6% 65%
1	B	647	 30% 6% 65%

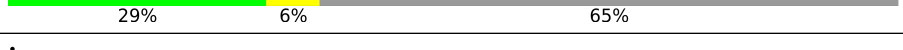
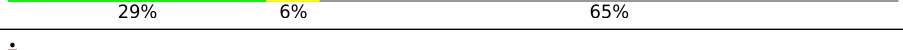
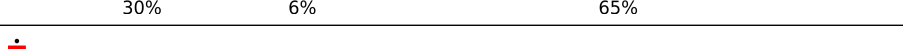
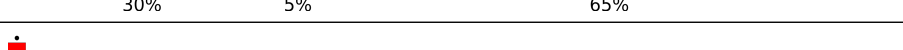
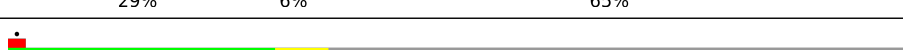
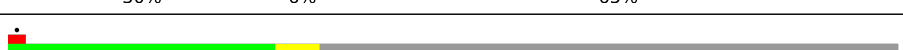
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Mol	Chain	Length	Quality of chain
1	BA	647	
1	BB	647	
1	C	647	
1	CA	647	
1	CB	647	
1	D	647	
1	DA	647	
1	DB	647	
1	E	647	
1	EA	647	
1	EB	647	
1	F	647	
1	FA	647	
1	FB	647	
1	G	647	
1	GA	647	
1	GB	647	
1	H	647	
1	HA	647	
1	HB	647	
1	I	647	
1	IA	647	
1	J	647	
1	JA	647	
1	K	647	

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Mol	Chain	Length	Quality of chain
1	KA	647	
1	L	647	
1	LA	647	
1	M	647	
1	MA	647	
1	N	647	
1	NA	647	
1	O	647	
1	OA	647	
1	P	647	
1	PA	647	
1	Q	647	
1	QA	647	
1	R	647	
1	RA	647	
1	S	647	
1	SA	647	
1	T	647	
1	TA	647	
1	UA	647	
1	V	647	
1	VA	647	
1	W	647	
1	WA	647	
1	X	647	

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Mol	Chain	Length	Quality of chain
1	XA	647	 29%6%65%
1	Y	647	 30%6%65%
1	YA	647	 30%6%65%
1	Z	647	 29%6%65%
1	ZA	647	 29%6%65%
1	a	647	 30%5%65%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 105420 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 Acetyltransferase component of pyruvate dehydrogenase complex.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	AA	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	AB	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	B	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	BA	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	BB	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	C	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	CA	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	CB	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	D	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	DA	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	DB	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	E	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	EA	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	EB	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	F	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	FA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	FB	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	G	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	GA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	GB	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	H	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	HA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	HB	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	I	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	IA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	J	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	JA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	K	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	KA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	L	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	LA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	M	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	MA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	N	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	NA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	O	228	Total 1757	C 1120	N 303	O 324	S 10	0	0

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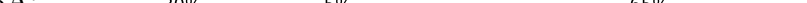
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1	OA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	P	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	PA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	Q	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	QA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	R	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	RA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	S	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	SA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	T	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	TA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	UA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	V	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	VA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	W	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	WA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	X	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	XA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	Y	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	YA	228	Total 1757	C 1120	N 303	O 324	S 10	0	0
1	Z	228	Total 1757	C 1120	N 303	O 324	S 10	0	0

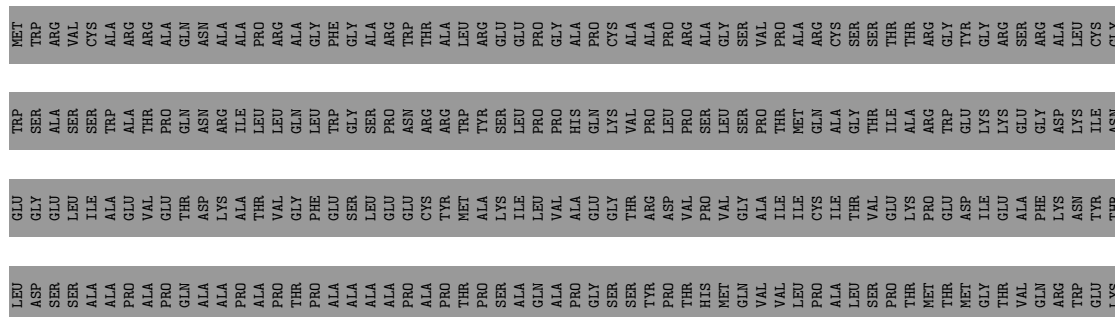
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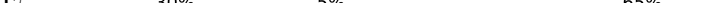
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Mol	Chain	Residues	Atoms					AltConf	Trace
1	ZA	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		
1	a	228	Total	C	N	O	S	0	0
			1757	1120	303	324	10		

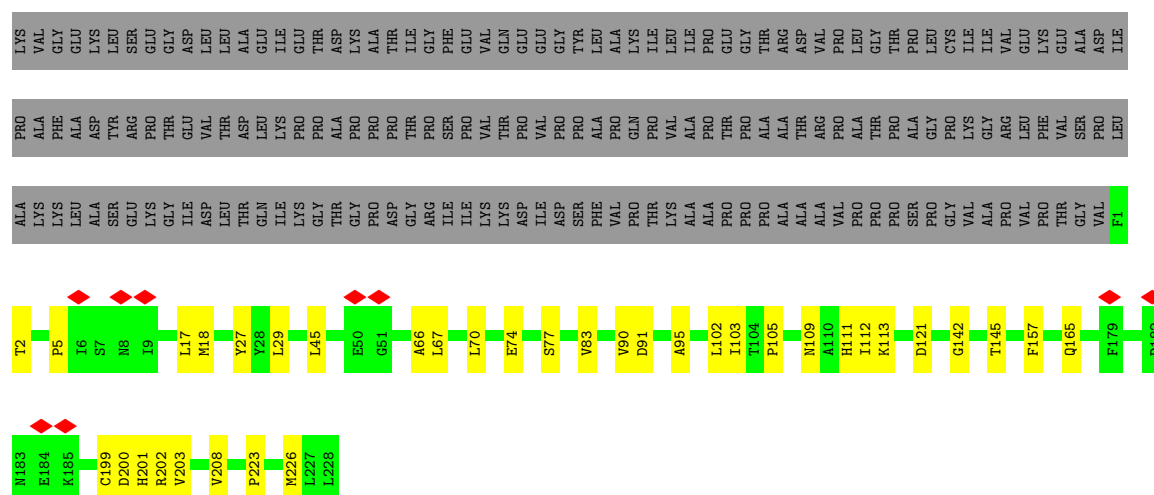
Frequency	Percentage
Daily	30%
Weekly	6%
Not at all	65%

- Chain BA: 



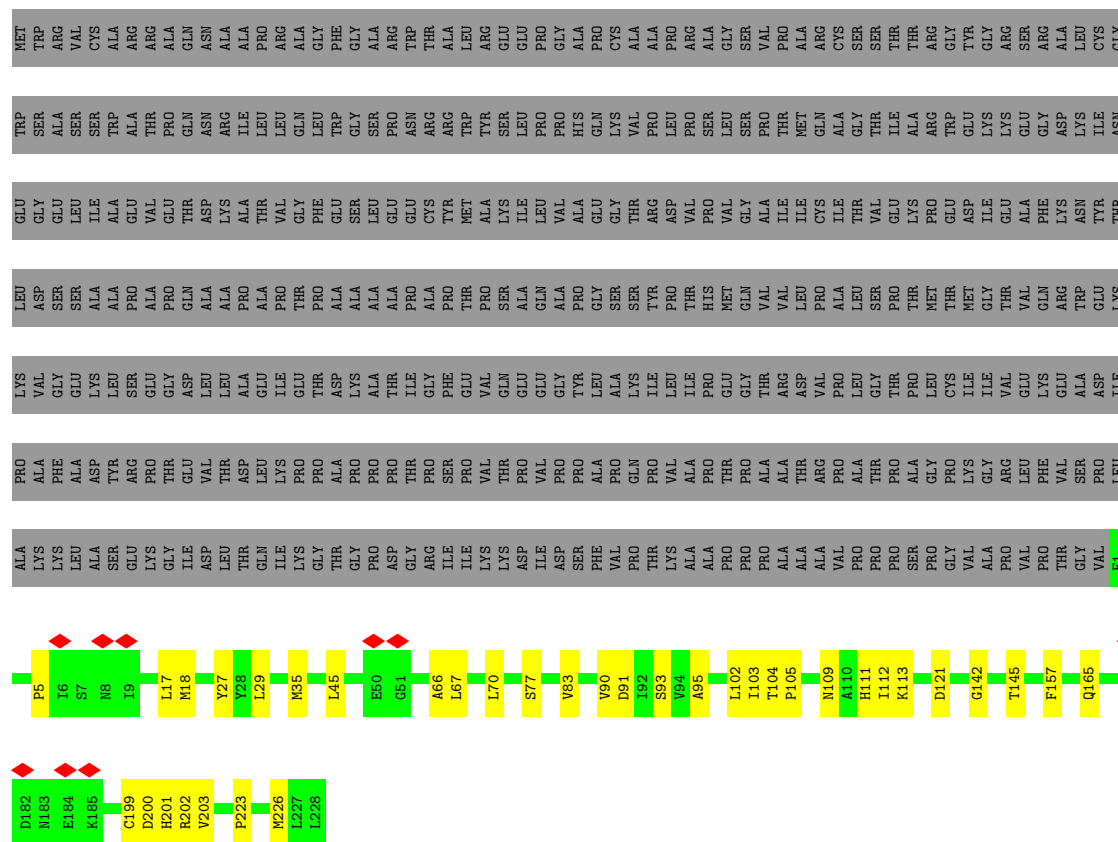
- Chain E:  30% 5% 65%





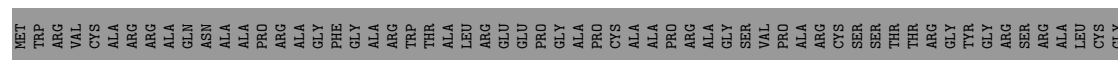
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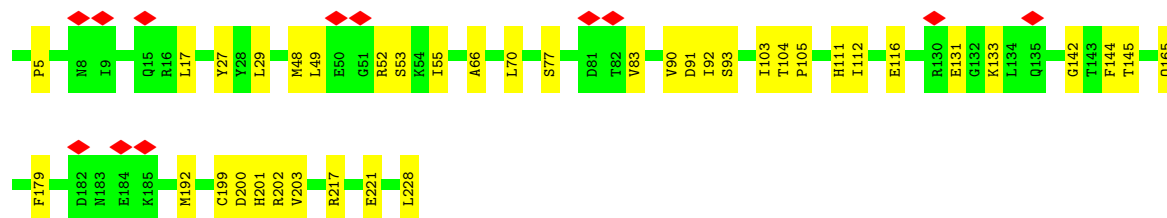
Chain FA: 30% 6% 65%



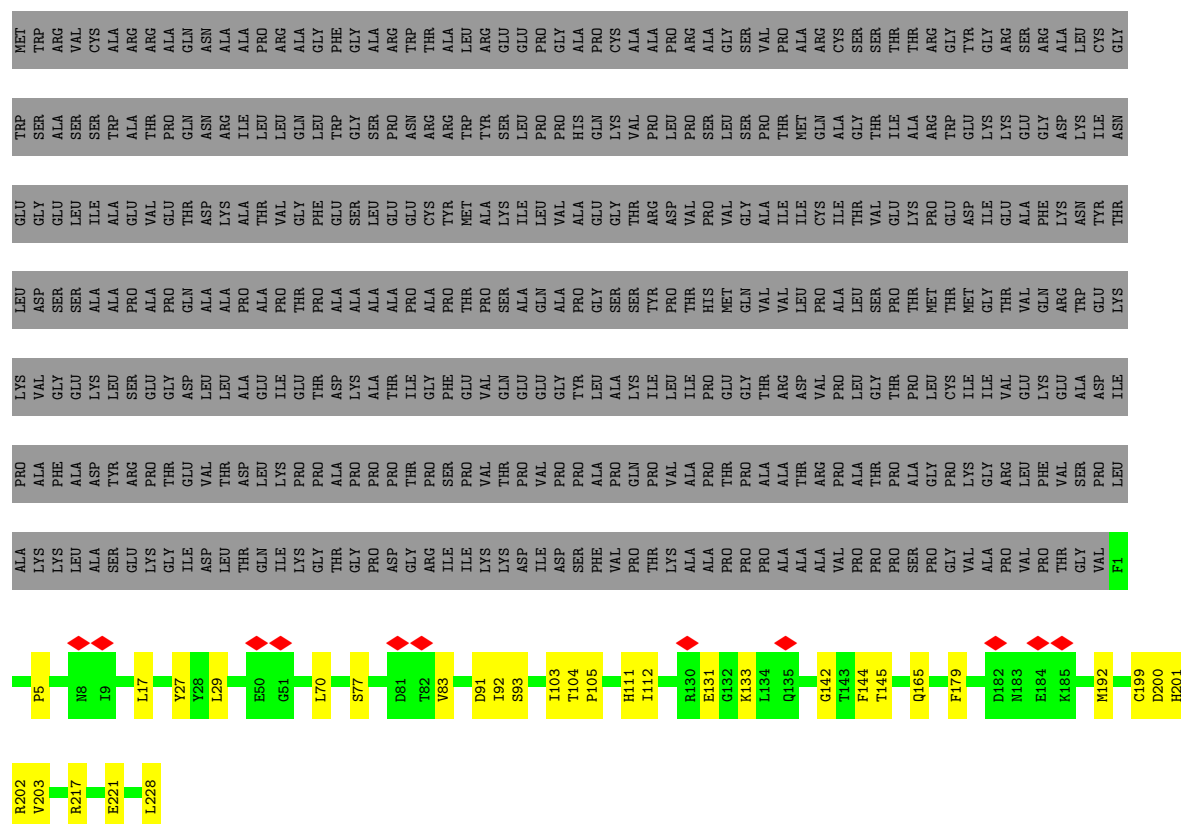
- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

Chain FB: 29% 6% 65%

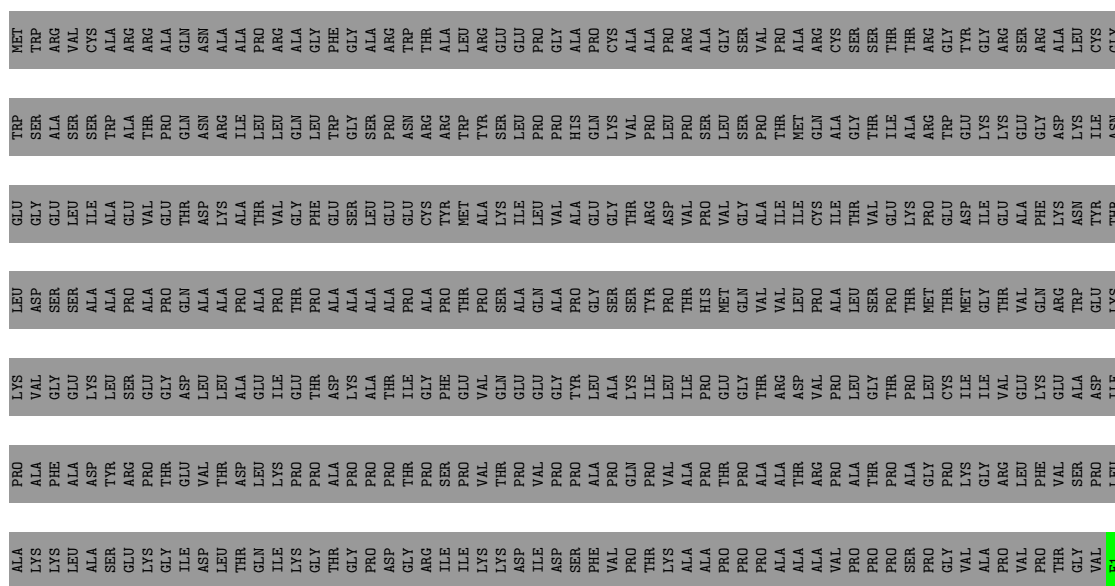




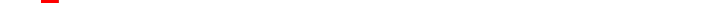
- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

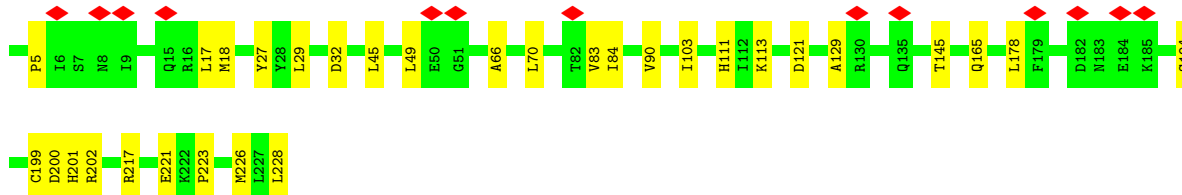


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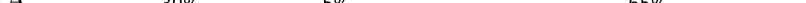


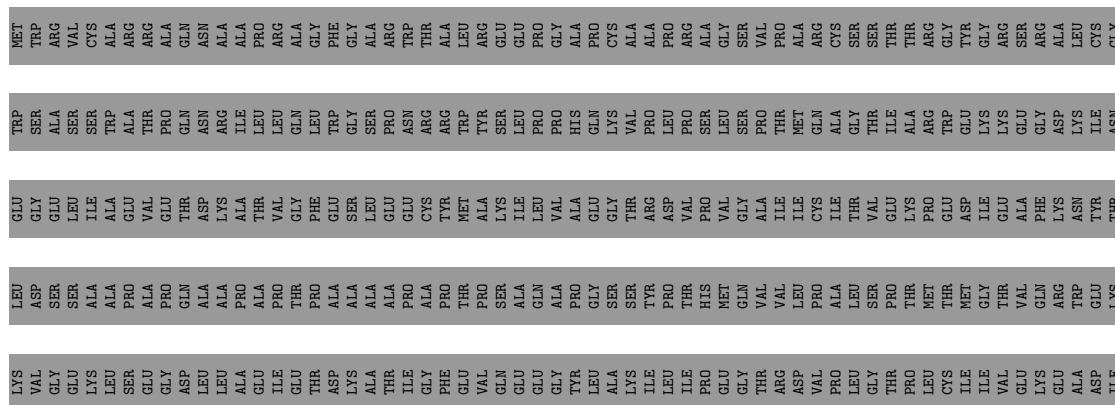
- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

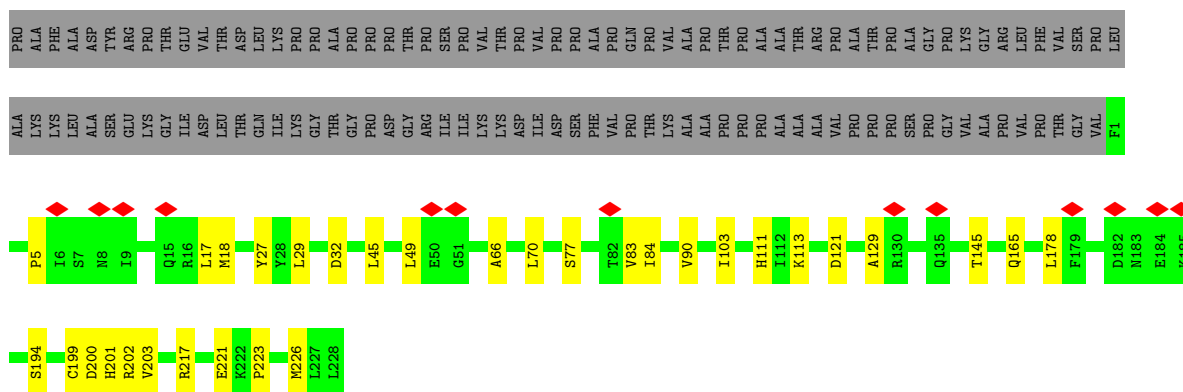
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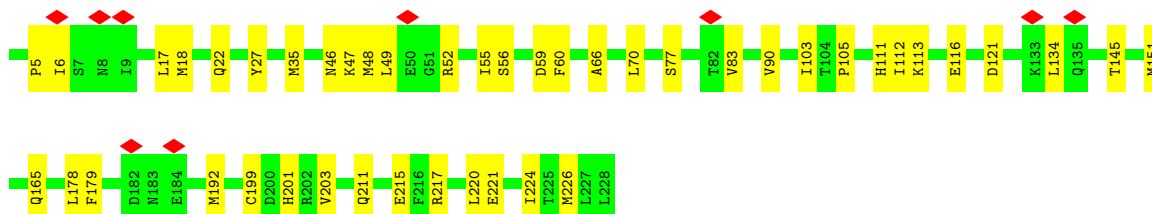
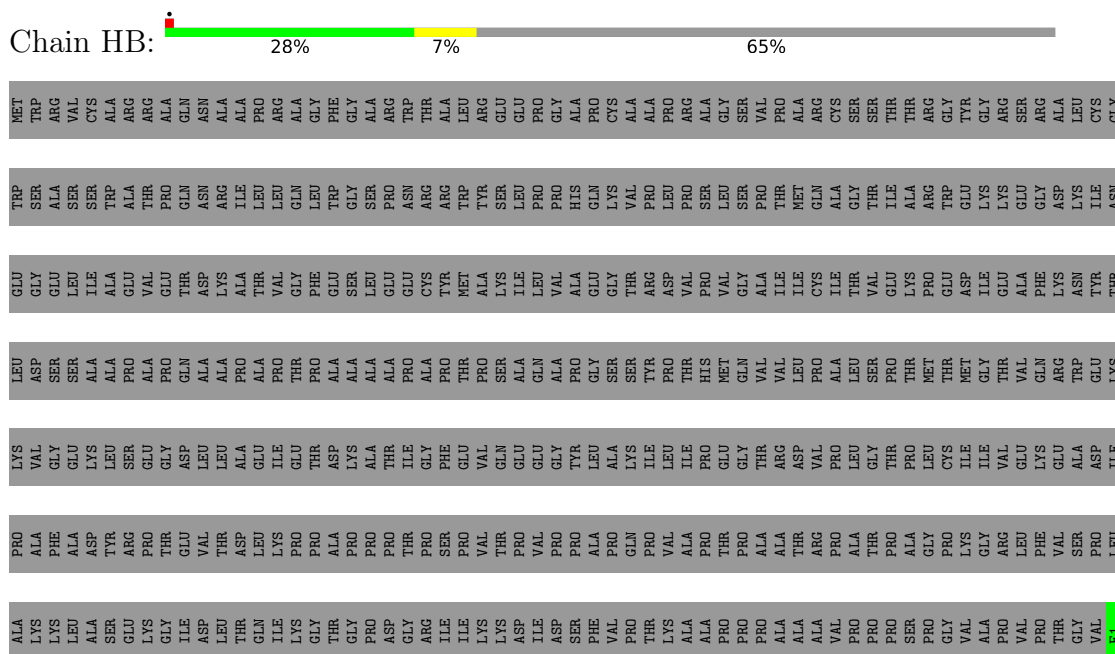
- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

Chain HA: 

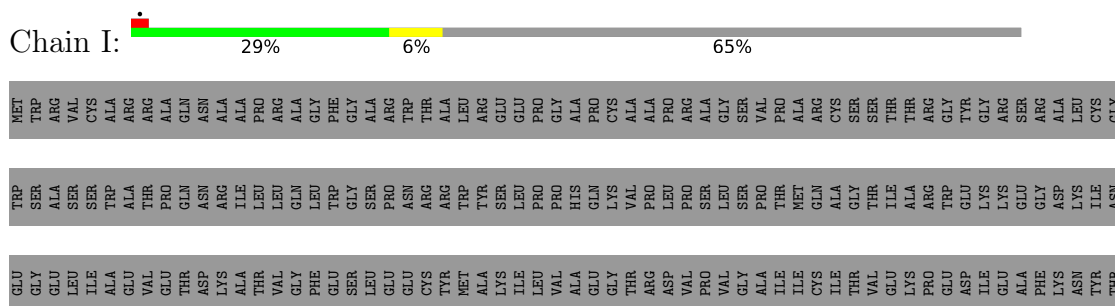


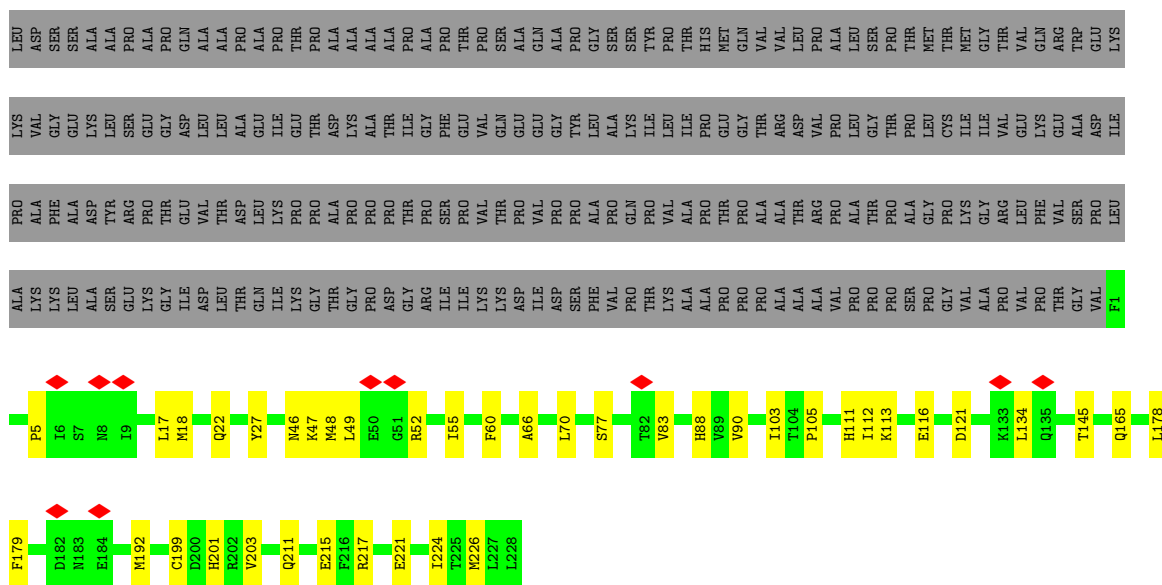


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex



- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex



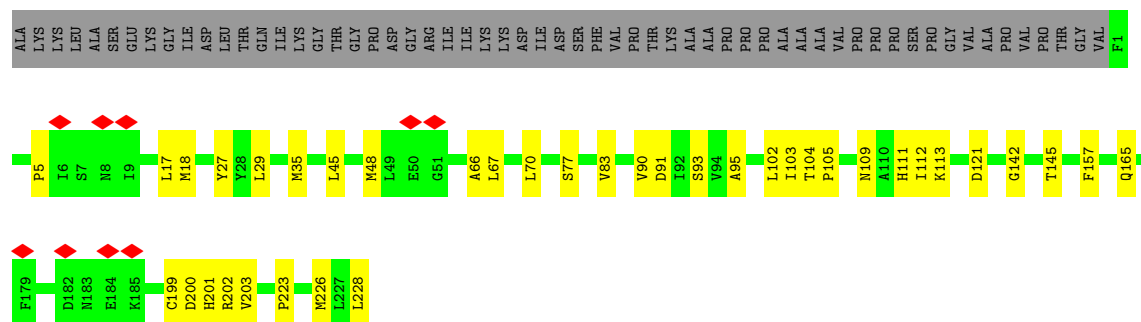


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

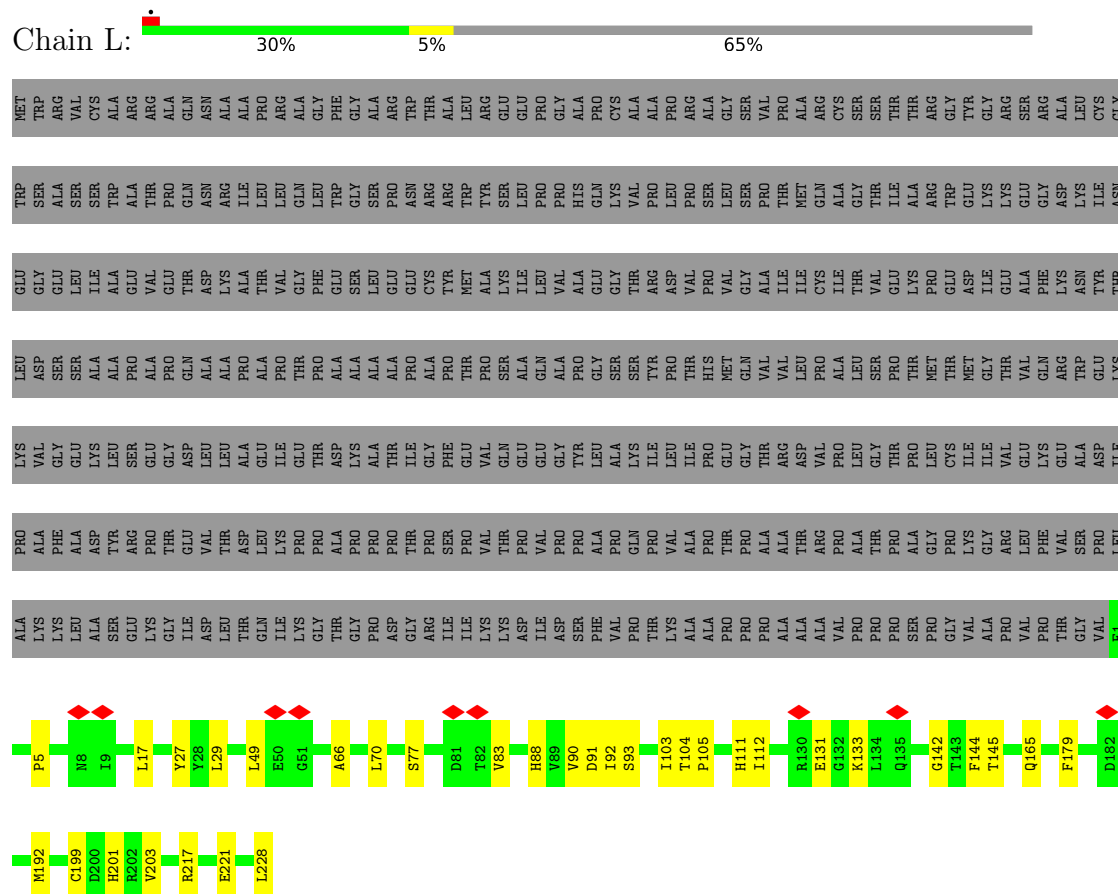


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

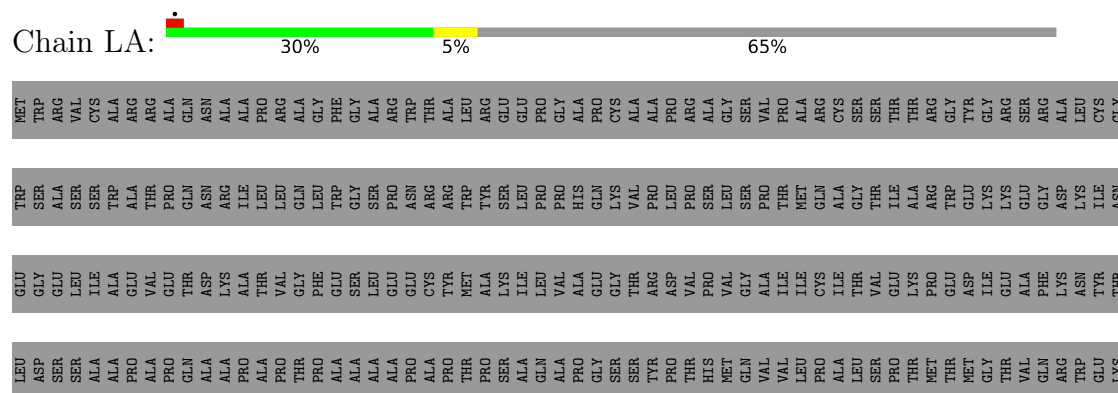


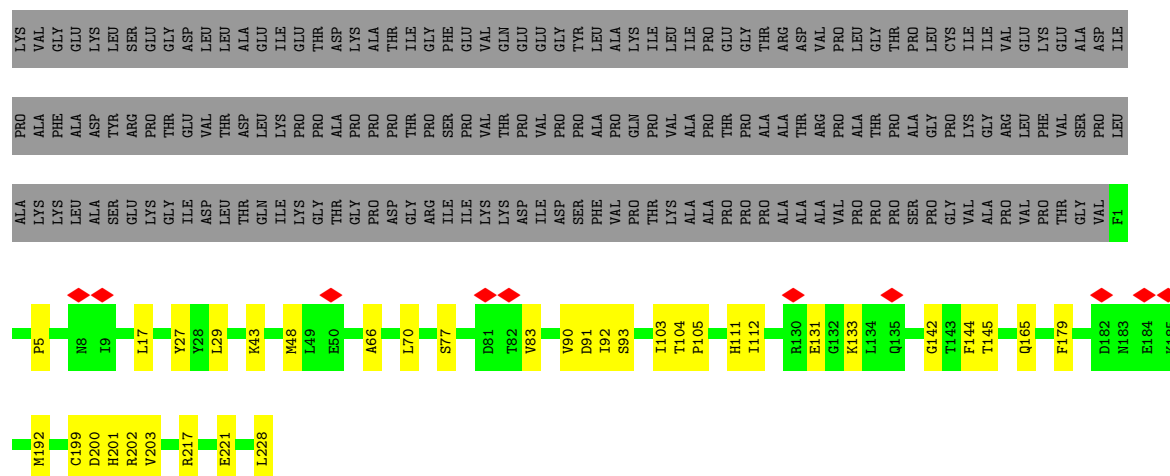


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

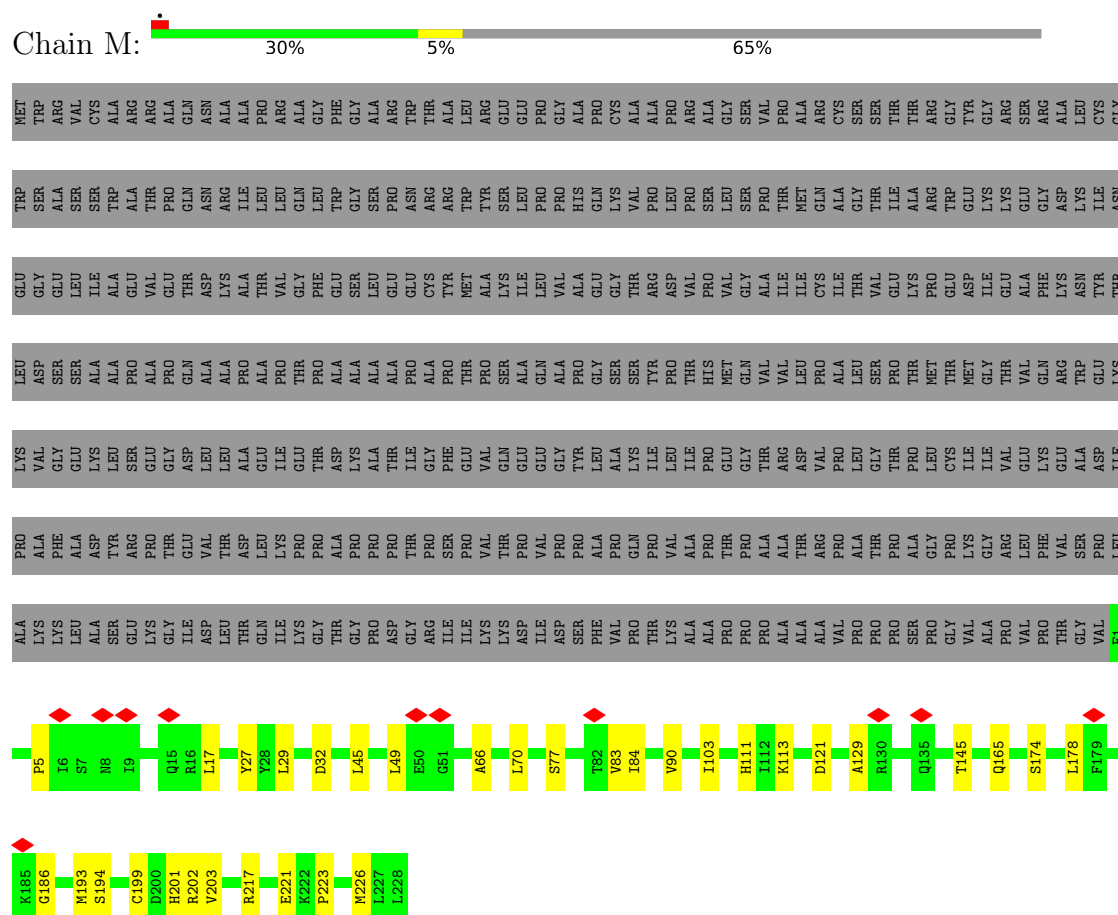


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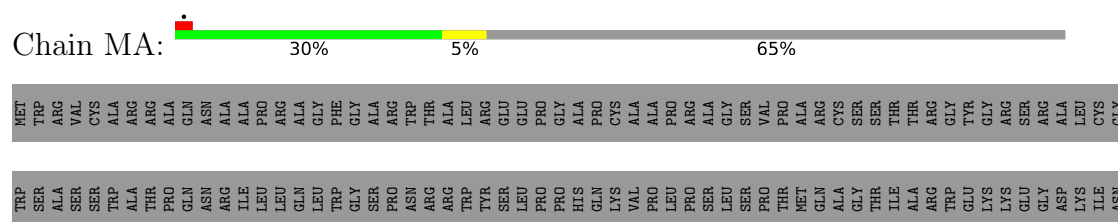


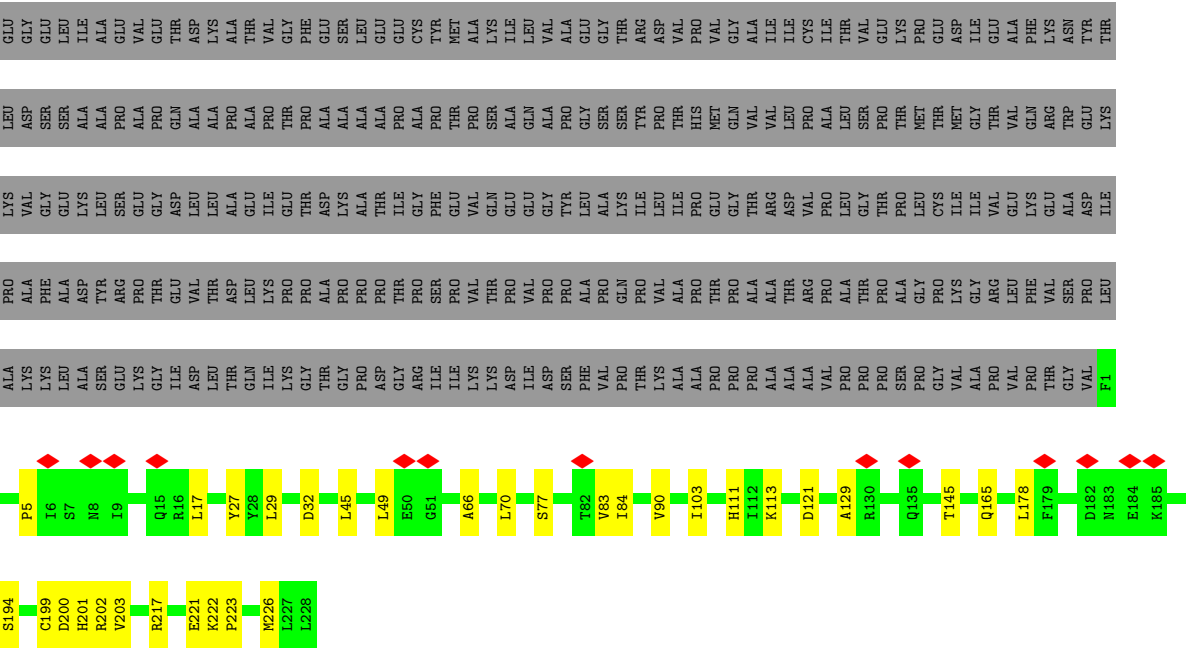


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

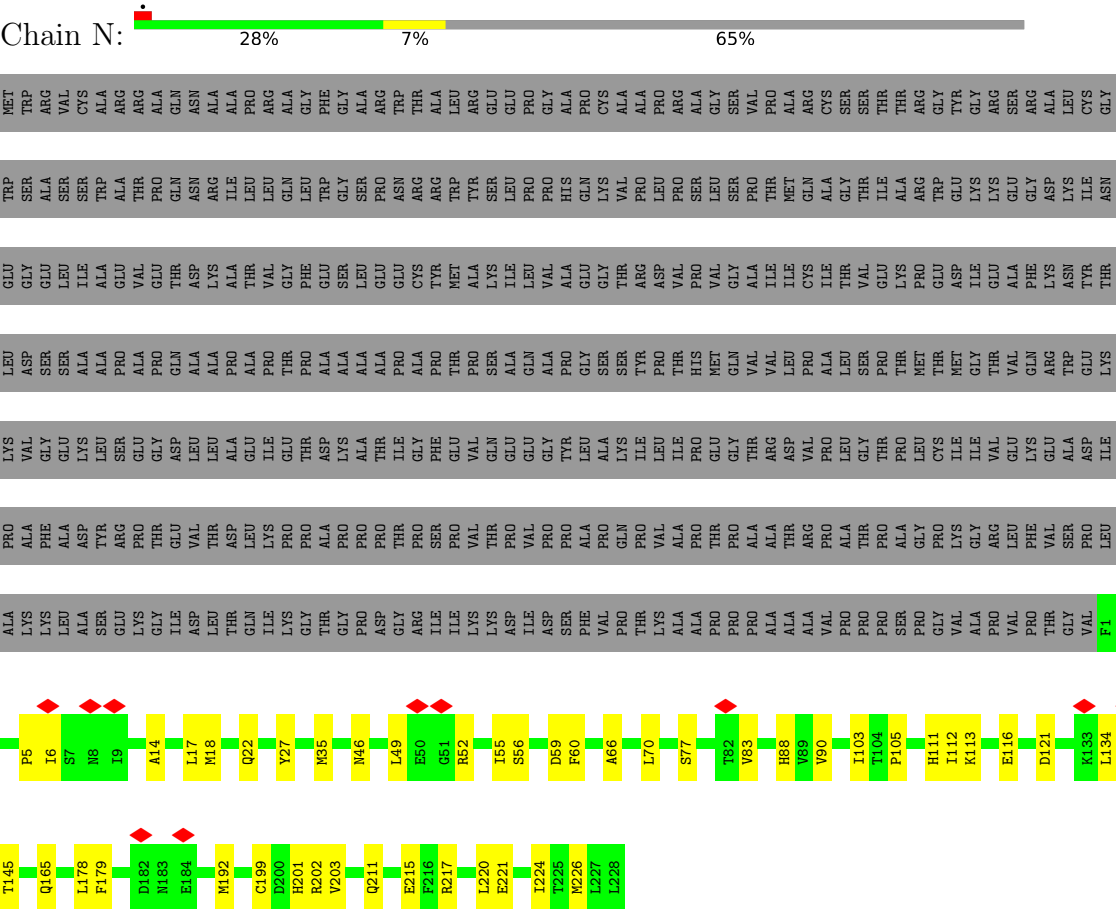


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex





• Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

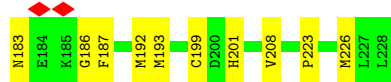
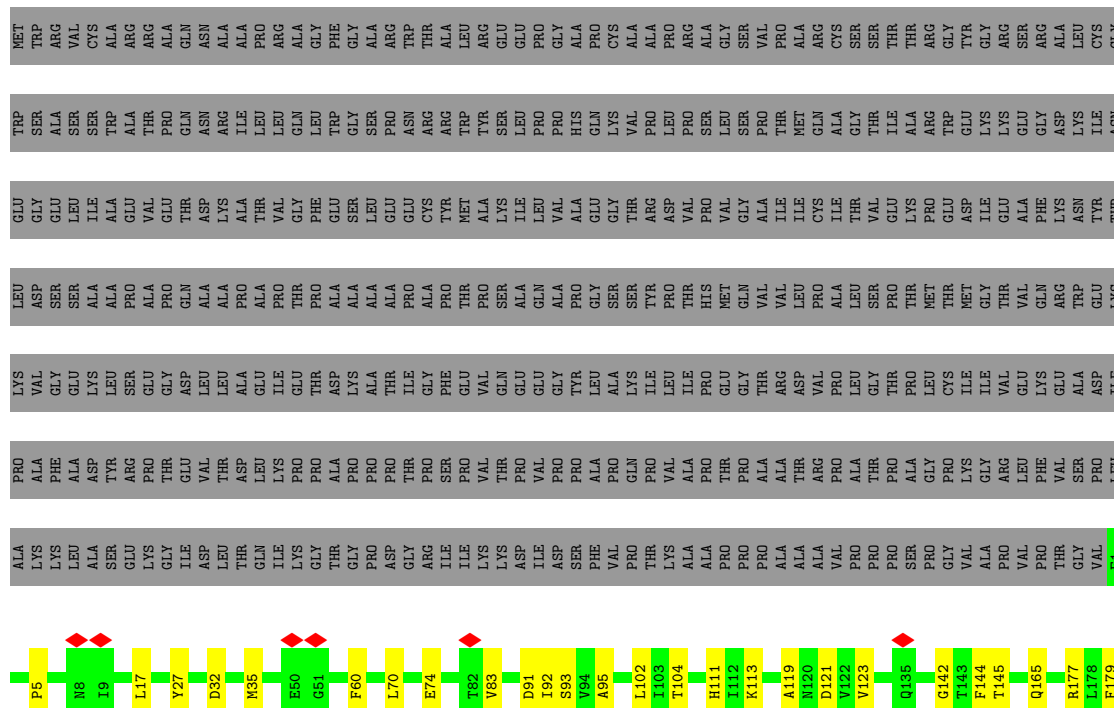


• Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

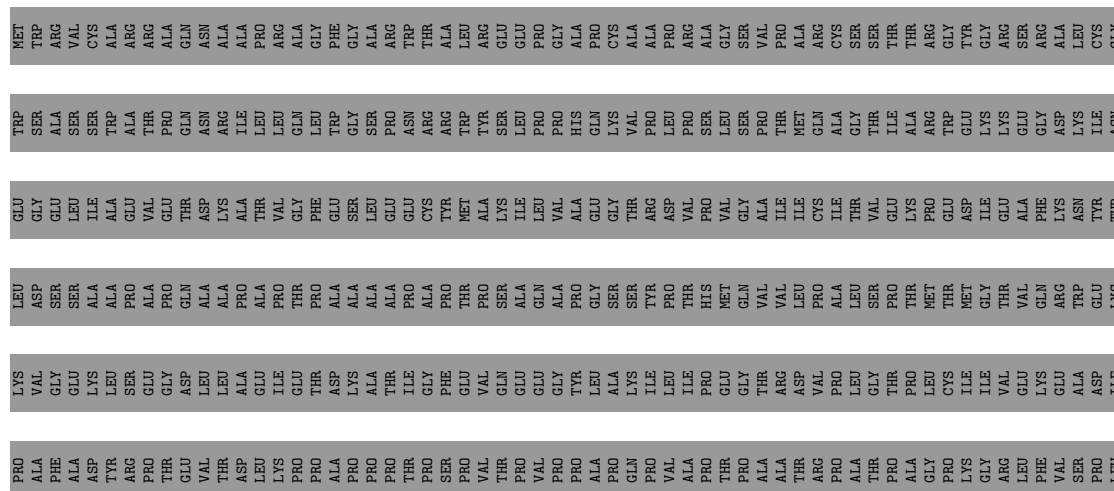


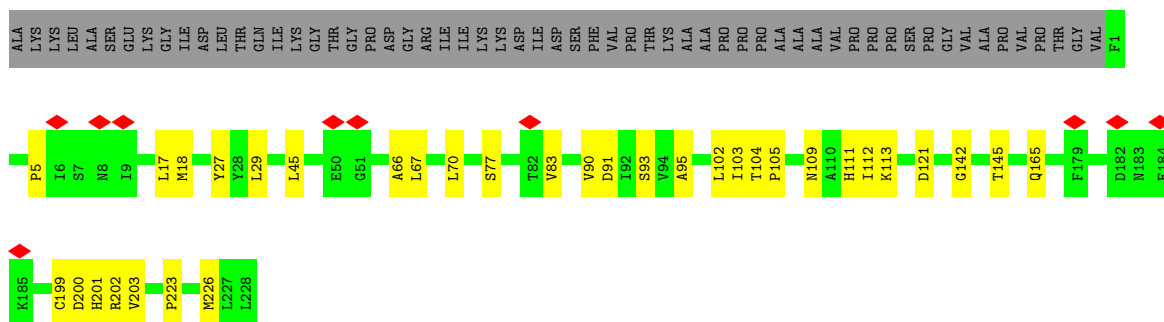


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

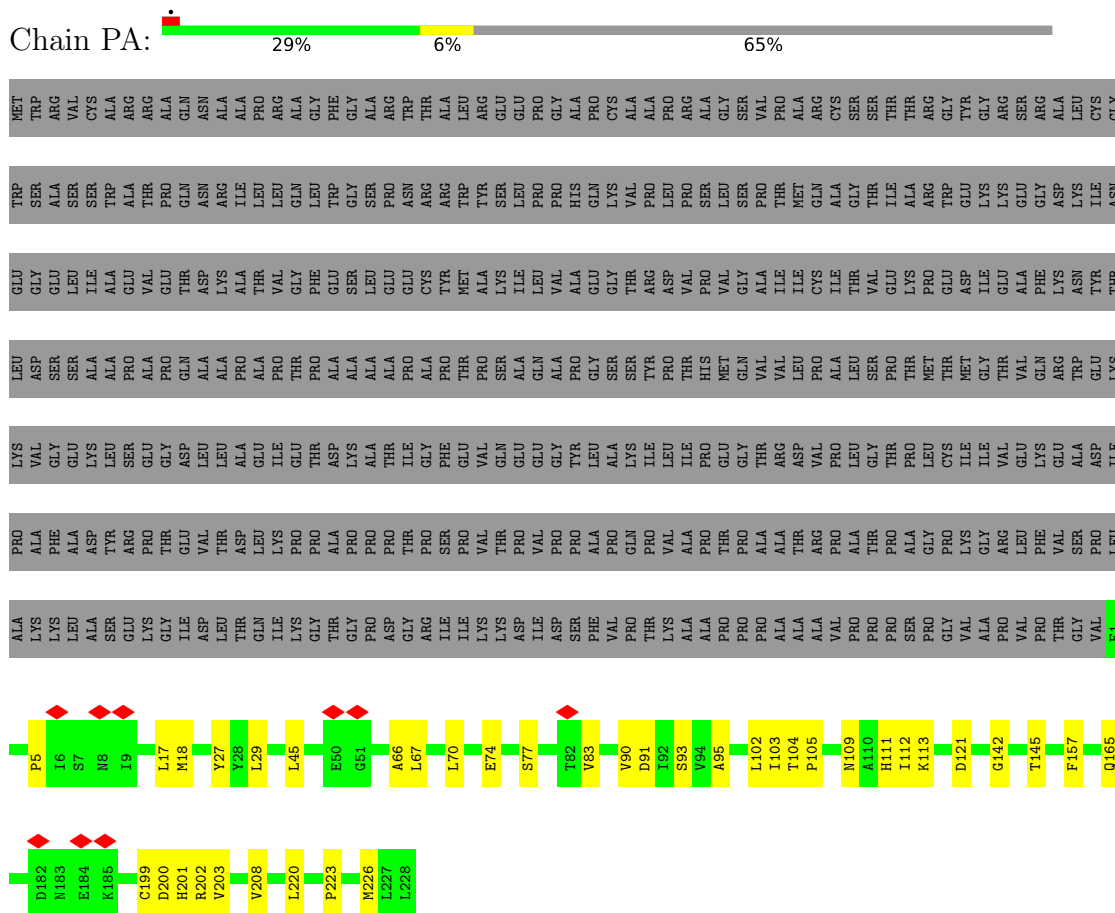


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

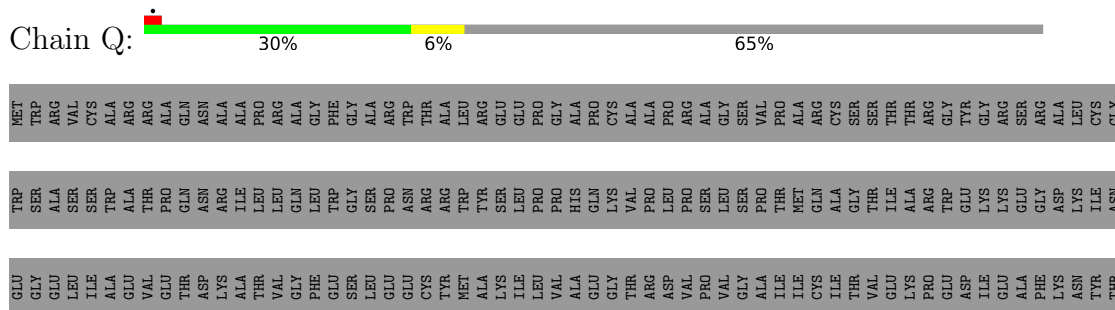


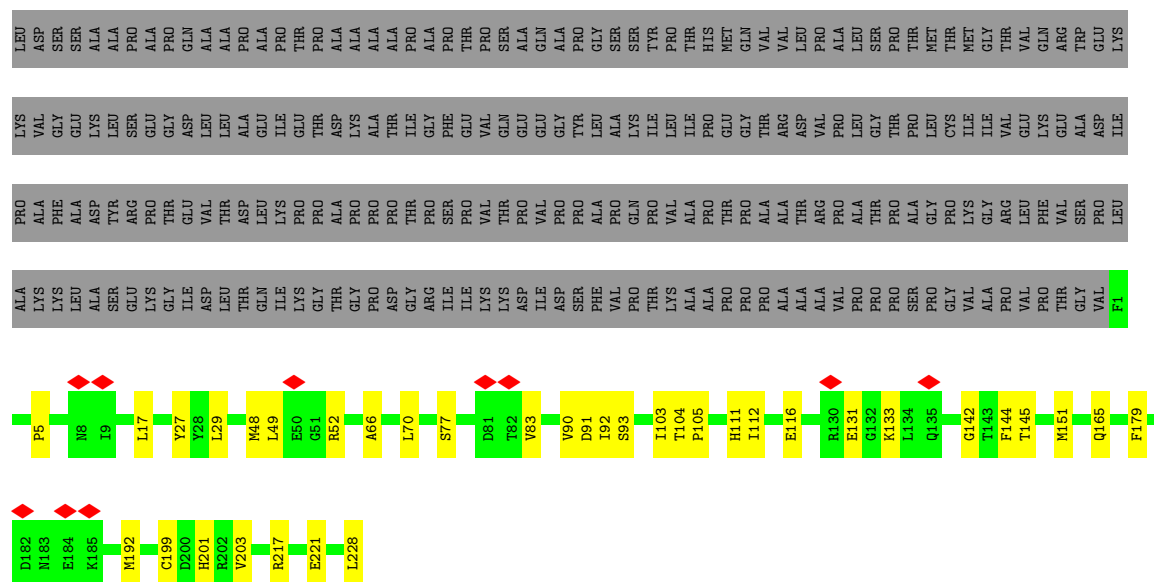


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

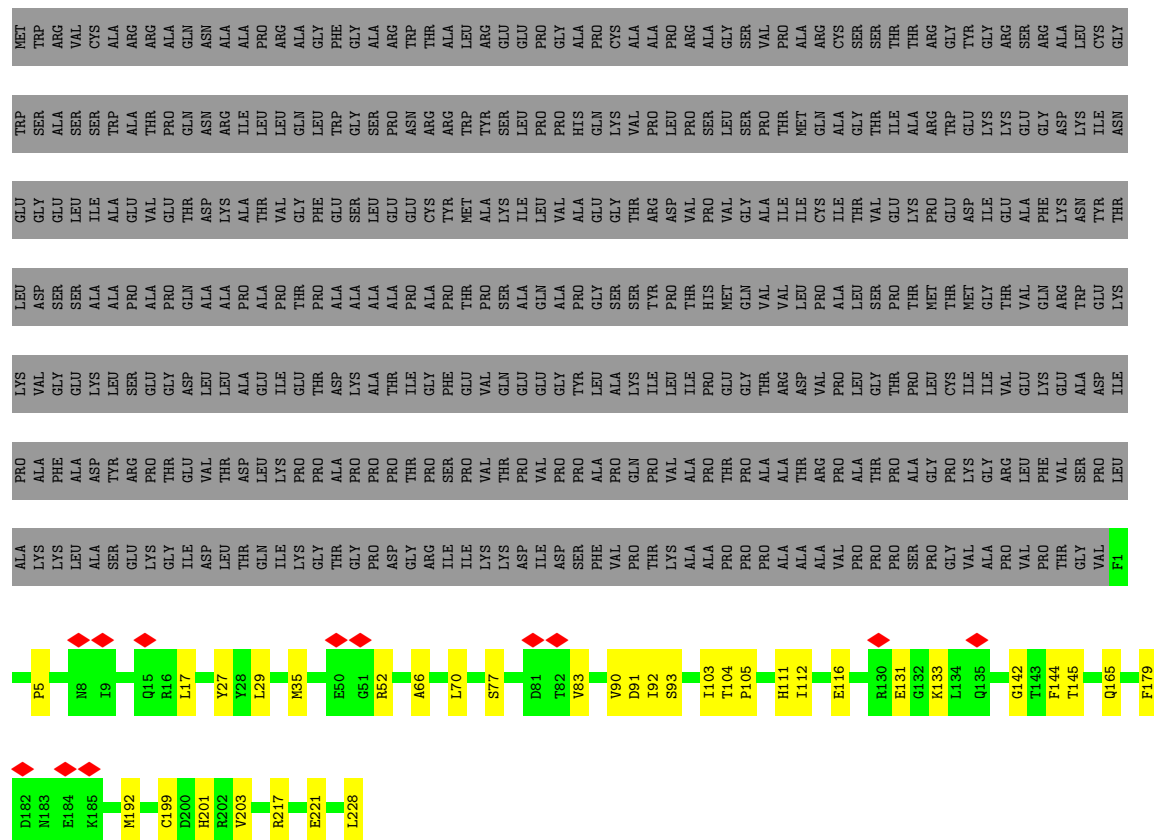


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex



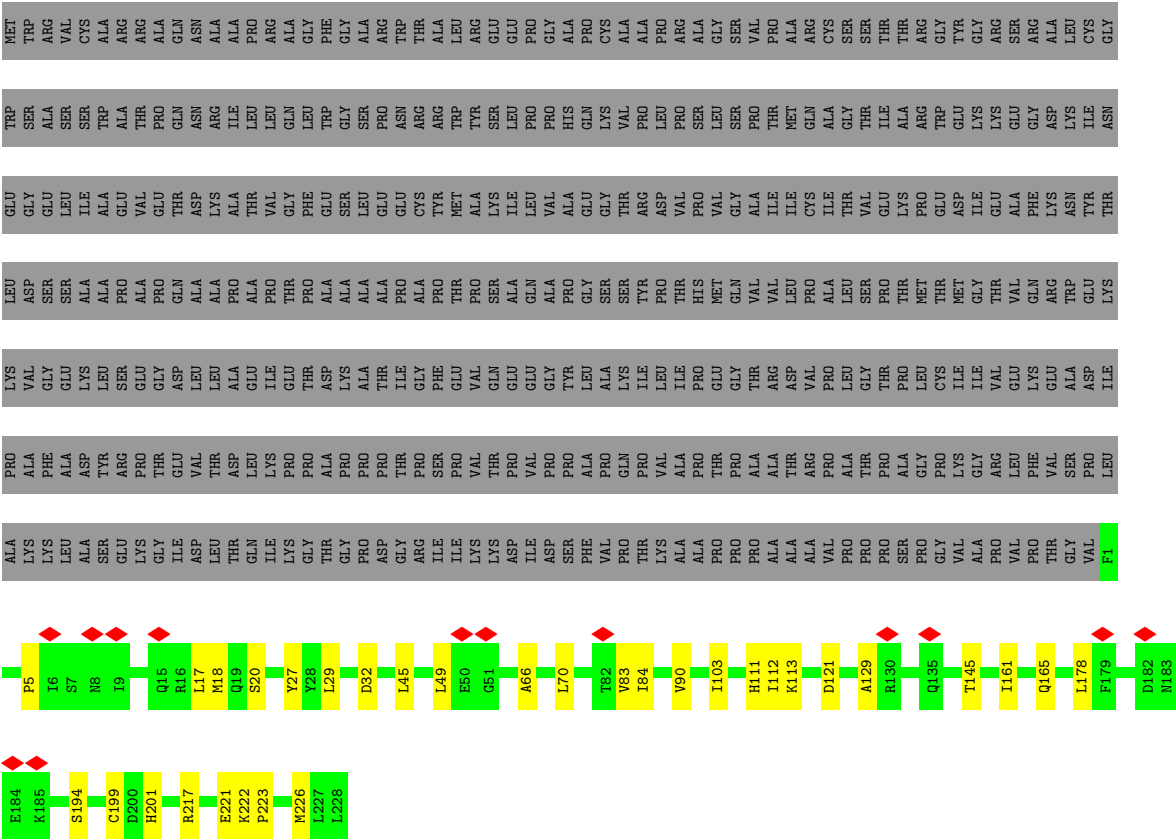


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

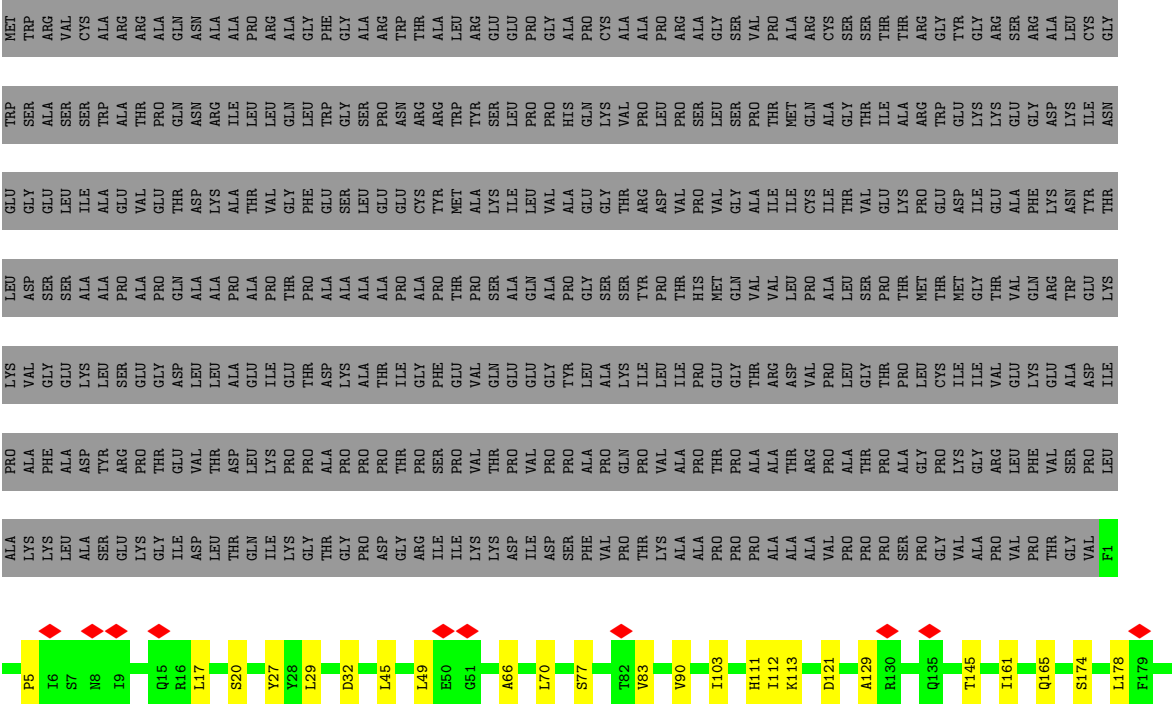


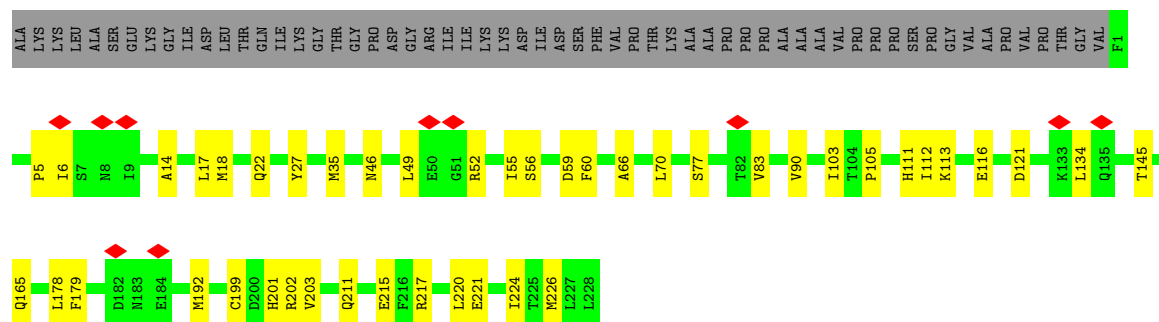
- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex



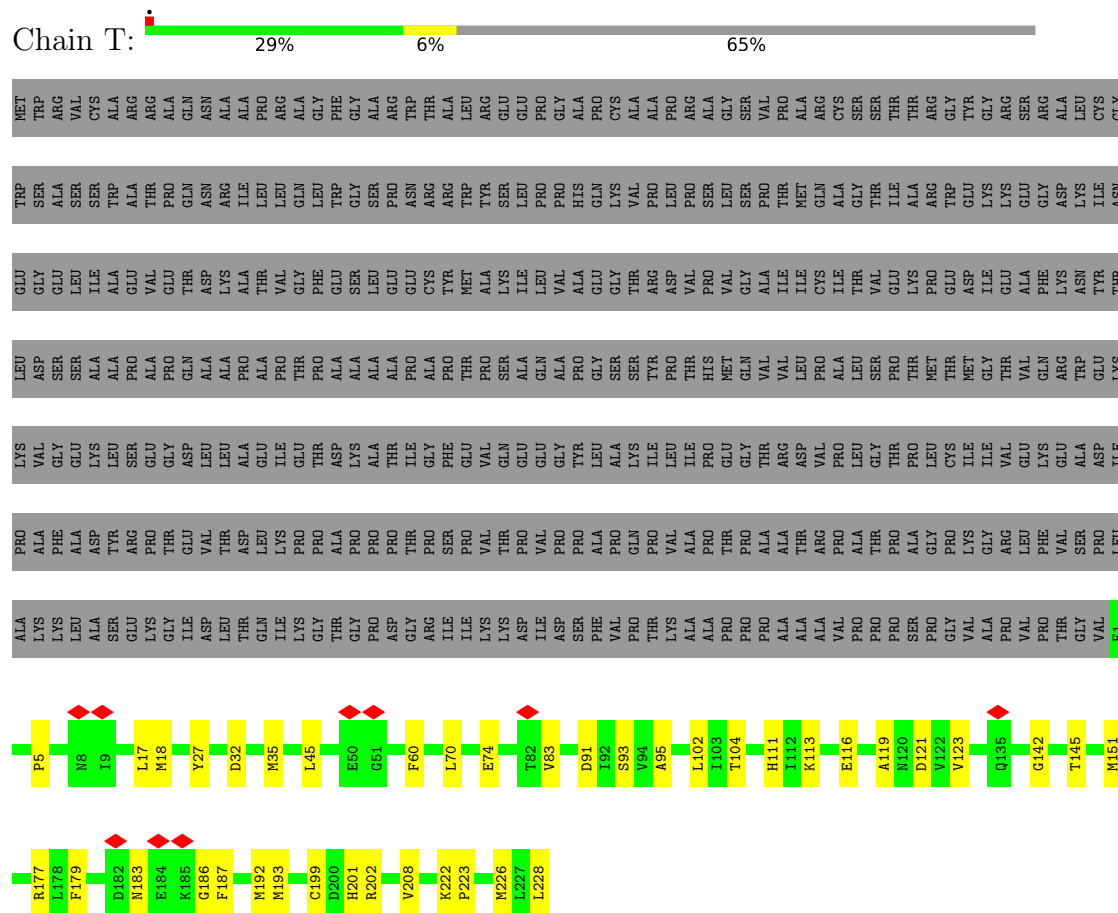


● Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

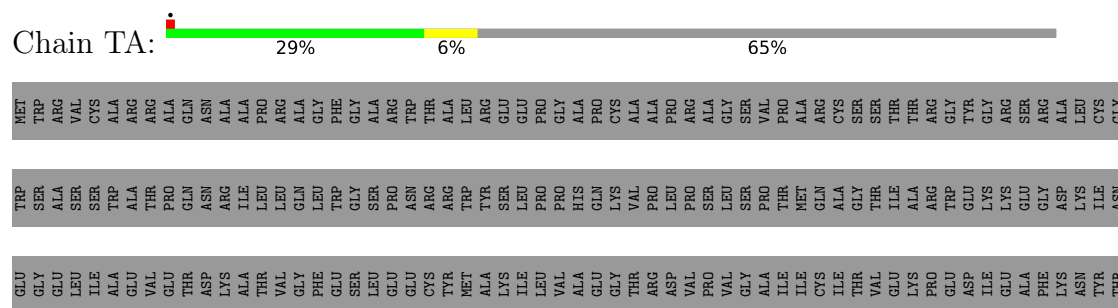


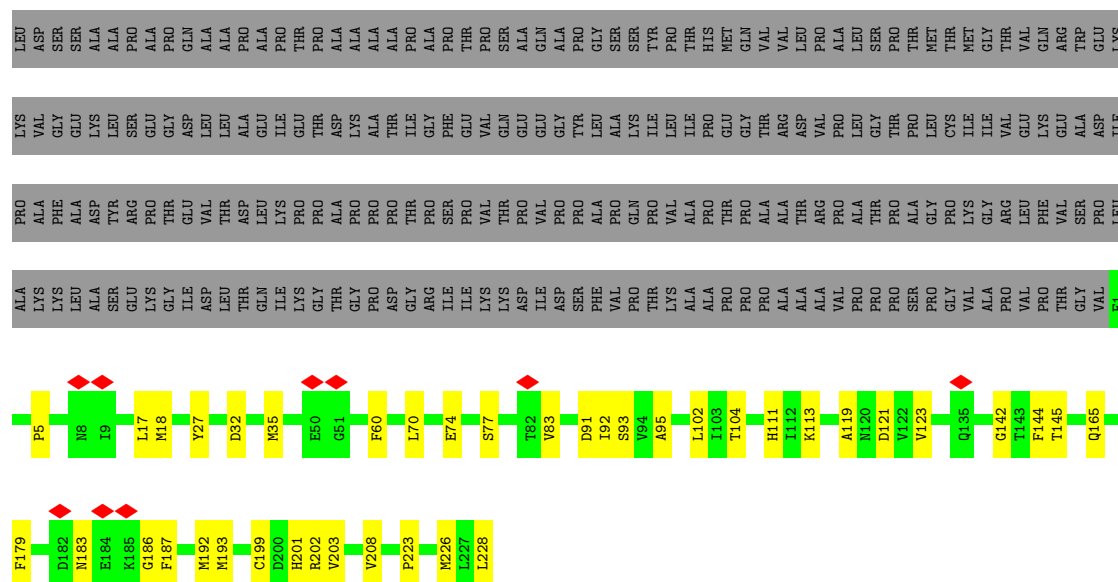


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

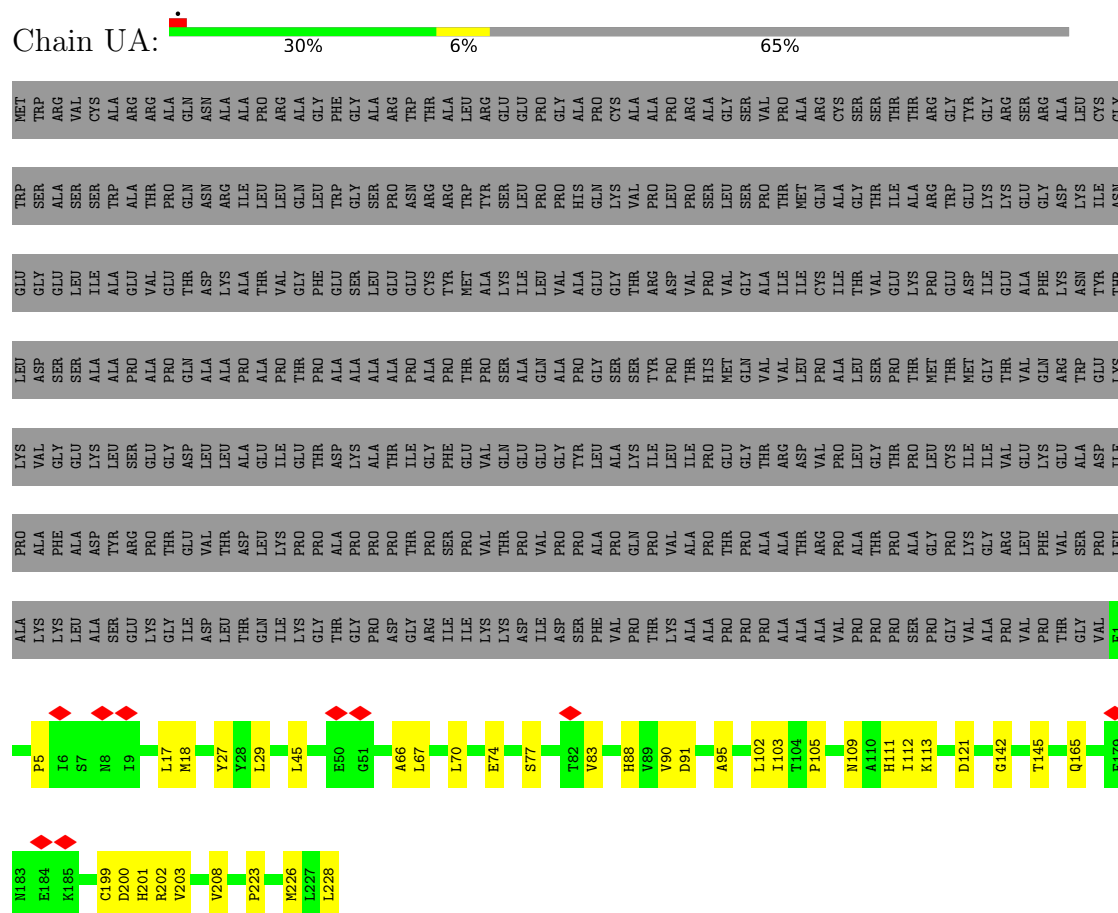


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex



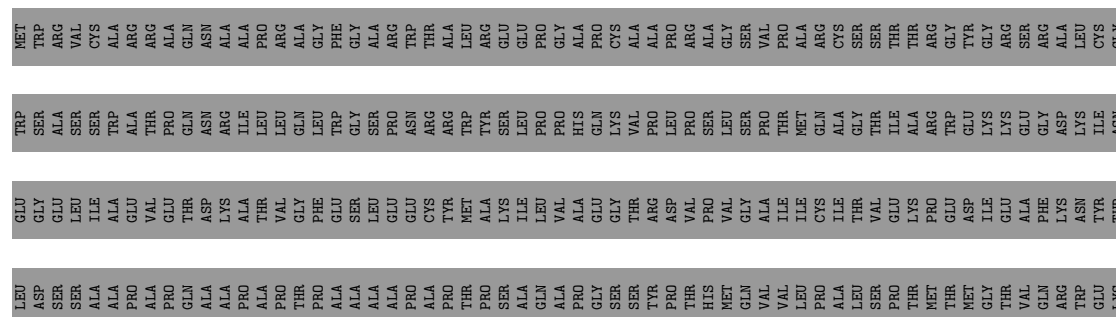


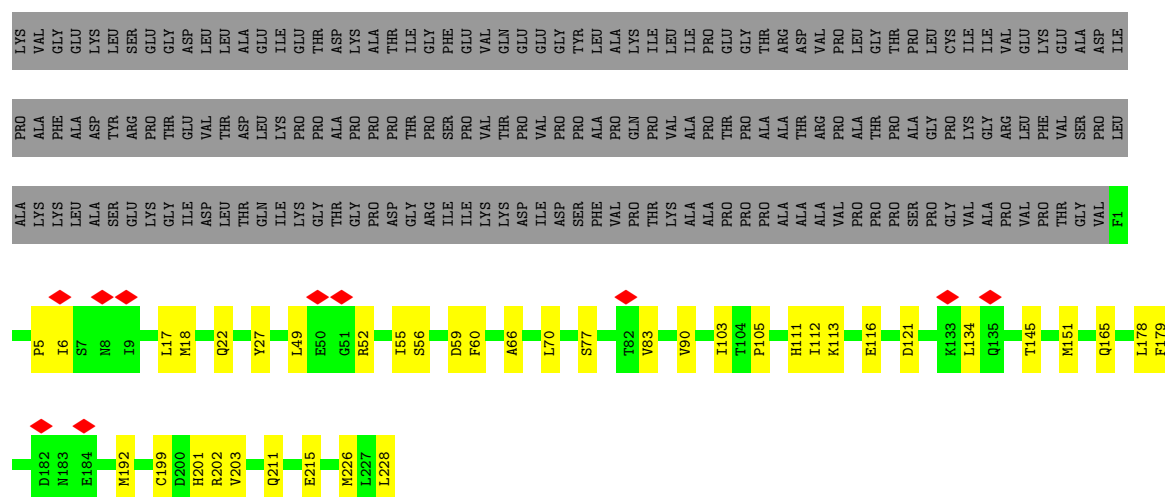
- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex



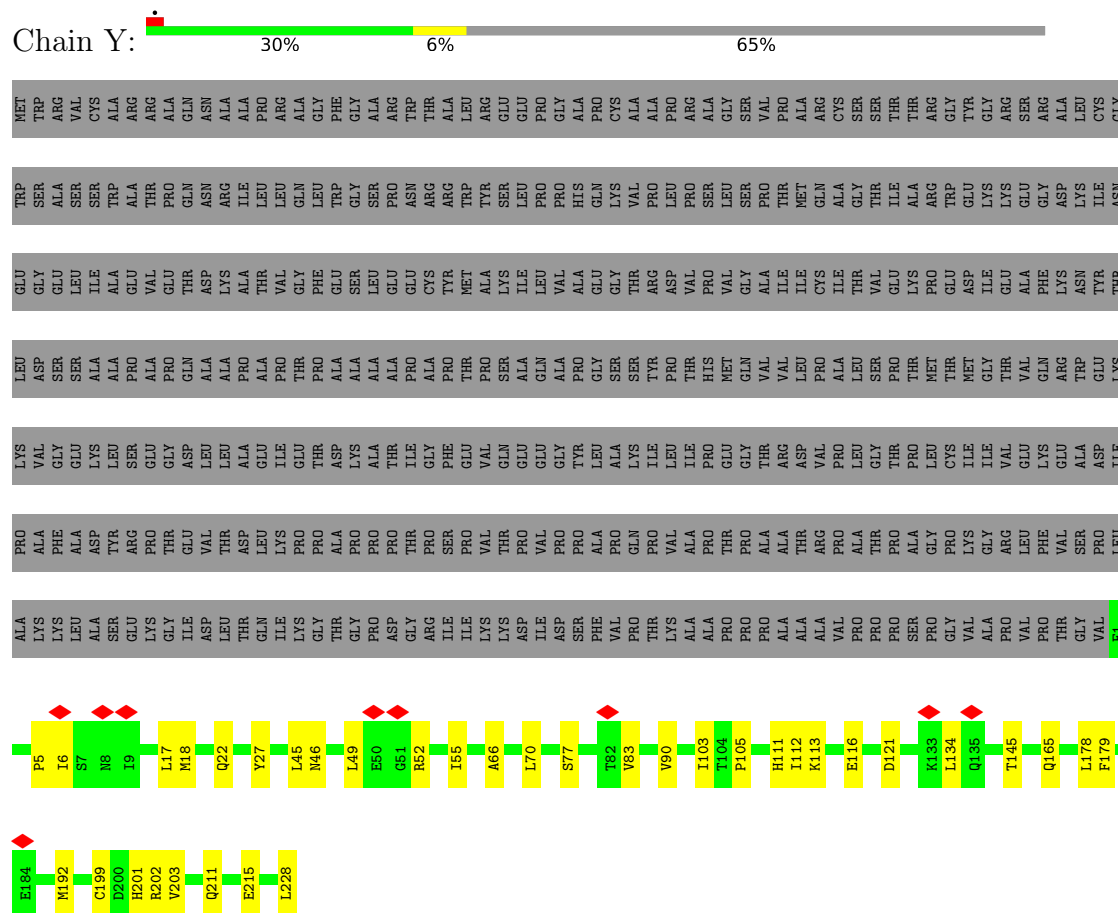
- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex



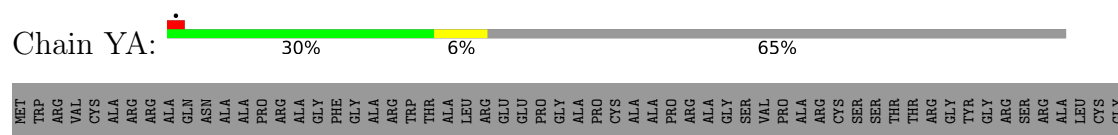




- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex

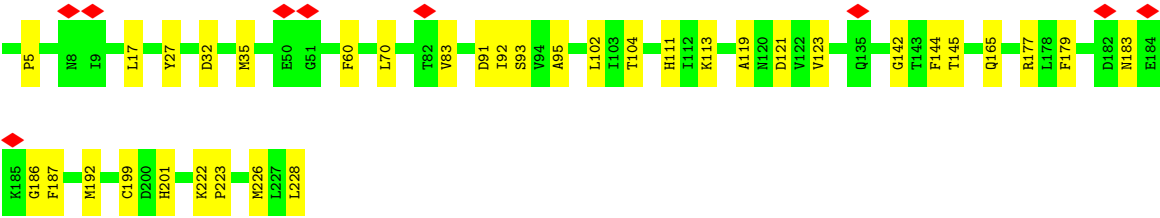


- Molecule 1: Acetyltransferase component of pyruvate dehydrogenase complex



- Chain ZA:





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	68914	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	52	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	2.424	Depositor
Minimum map value	-1.451	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.061	Depositor
Recommended contour level	0.34	Depositor
Map size ($\text{\AA}$)	614.4, 614.4, 614.4	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.2, 1.2, 1.2	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	A	0.12	0/1788	0.30	0/2422
1	AA	0.13	0/1788	0.31	0/2422
1	AB	0.13	0/1788	0.33	0/2422
1	B	0.13	0/1788	0.33	0/2422
1	BA	0.13	0/1788	0.31	0/2422
1	BB	0.13	0/1788	0.31	0/2422
1	C	0.13	0/1788	0.32	0/2422
1	CA	0.13	0/1788	0.33	0/2422
1	CB	0.13	0/1788	0.31	0/2422
1	D	0.13	0/1788	0.31	0/2422
1	DA	0.13	0/1788	0.32	0/2422
1	DB	0.13	0/1788	0.30	0/2422
1	E	0.13	0/1788	0.30	0/2422
1	EA	0.13	0/1788	0.30	0/2422
1	EB	0.13	0/1788	0.31	0/2422
1	F	0.13	0/1788	0.31	0/2422
1	FA	0.13	0/1788	0.31	0/2422
1	FB	0.13	0/1788	0.31	0/2422
1	G	0.13	0/1788	0.32	0/2422
1	GA	0.13	0/1788	0.32	0/2422
1	GB	0.14	0/1788	0.32	0/2422
1	H	0.13	0/1788	0.32	0/2422
1	HA	0.13	0/1788	0.32	0/2422
1	HB	0.13	0/1788	0.31	0/2422
1	I	0.13	0/1788	0.31	0/2422
1	IA	0.13	0/1788	0.31	0/2422
1	J	0.13	0/1788	0.30	0/2422
1	JA	0.13	0/1788	0.30	0/2422
1	K	0.13	0/1788	0.31	0/2422
1	KA	0.12	0/1788	0.30	0/2422
1	L	0.13	0/1788	0.31	0/2422
1	LA	0.13	0/1788	0.33	0/2422
1	M	0.14	0/1788	0.32	0/2422
1	MA	0.13	0/1788	0.31	0/2422

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	N	0.13	0/1788	0.32	0/2422
1	NA	0.13	0/1788	0.31	0/2422
1	O	0.13	0/1788	0.30	0/2422
1	OA	0.13	0/1788	0.30	0/2422
1	P	0.13	0/1788	0.31	0/2422
1	PA	0.13	0/1788	0.31	0/2422
1	Q	0.13	0/1788	0.32	0/2422
1	QA	0.13	0/1788	0.31	0/2422
1	R	0.13	0/1788	0.31	0/2422
1	RA	0.13	0/1788	0.32	0/2422
1	S	0.13	0/1788	0.31	0/2422
1	SA	0.13	0/1788	0.32	0/2422
1	T	0.13	0/1788	0.30	0/2422
1	TA	0.13	0/1788	0.30	0/2422
1	UA	0.13	0/1788	0.31	0/2422
1	V	0.13	0/1788	0.31	0/2422
1	VA	0.13	0/1788	0.32	0/2422
1	W	0.13	0/1788	0.33	0/2422
1	WA	0.13	0/1788	0.31	0/2422
1	X	0.13	0/1788	0.31	0/2422
1	XA	0.13	0/1788	0.33	0/2422
1	Y	0.13	0/1788	0.33	0/2422
1	YA	0.13	0/1788	0.30	0/2422
1	Z	0.13	0/1788	0.31	0/2422
1	ZA	0.12	0/1788	0.30	0/2422
1	a	0.13	0/1788	0.30	0/2422
All	All	0.13	0/107280	0.31	0/145320

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	A	1757	0	1808	45	0
1	AA	1757	0	1808	31	0
1	AB	1757	0	1808	26	0
1	B	1757	0	1808	25	0
1	BA	1757	0	1808	20	0
1	BB	1757	0	1808	25	0
1	C	1757	0	1808	23	0
1	CA	1757	0	1808	22	0
1	CB	1757	0	1808	27	0
1	D	1757	0	1808	28	0
1	DA	1757	0	1808	29	0
1	DB	1757	0	1808	24	0
1	E	1757	0	1808	21	0
1	EA	1757	0	1808	27	0
1	EB	1757	0	1808	37	0
1	F	1757	0	1808	38	0
1	FA	1757	0	1808	38	0
1	FB	1757	0	1808	25	0
1	G	1757	0	1808	20	0
1	GA	1757	0	1808	24	0
1	GB	1757	0	1808	24	0
1	H	1757	0	1808	24	0
1	HA	1757	0	1808	23	0
1	HB	1757	0	1808	30	0
1	I	1757	0	1808	27	0
1	IA	1757	0	1808	29	0
1	J	1757	0	1808	22	0
1	JA	1757	0	1808	26	0
1	K	1757	0	1808	25	0
1	KA	1757	0	1808	34	0
1	L	1757	0	1808	23	0
1	LA	1757	0	1808	24	0
1	M	1757	0	1808	24	0
1	MA	1757	0	1808	23	0
1	N	1757	0	1808	30	0
1	NA	1757	0	1808	27	0
1	O	1757	0	1808	26	0
1	OA	1757	0	1808	22	0
1	P	1757	0	1808	29	0
1	PA	1757	0	1808	32	0
1	Q	1757	0	1808	25	0
1	QA	1757	0	1808	22	0
1	R	1757	0	1808	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	RA	1757	0	1808	25	0
1	S	1757	0	1808	25	0
1	SA	1757	0	1808	28	0
1	T	1757	0	1808	30	0
1	TA	1757	0	1808	28	0
1	UA	1757	0	1808	31	0
1	V	1757	0	1808	31	0
1	VA	1757	0	1808	22	0
1	W	1757	0	1808	23	0
1	WA	1757	0	1808	23	0
1	X	1757	0	1808	22	0
1	XA	1757	0	1808	26	0
1	Y	1757	0	1808	25	0
1	YA	1757	0	1808	23	0
1	Z	1757	0	1808	26	0
1	ZA	1757	0	1808	41	0
1	a	1757	0	1808	23	0
All	All	105420	0	108480	1299	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 6.

The worst 5 of 1299 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:A:157:PHE:CB	1:ZA:29:LEU:HD23	1.57	1.34
1:AA:157:PHE:CB	1:F:29:LEU:HD23	1.57	1.34
1:F:157:PHE:CB	1:UA:29:LEU:HD23	1.57	1.33
1:FA:157:PHE:CB	1:P:29:LEU:HD23	1.57	1.33
1:EB:157:PHE:CB	1:V:29:LEU:HD23	1.57	1.33

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 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 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	A	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	AA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	AB	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	B	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	BA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	BB	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	C	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	CA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	CB	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	D	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	DA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	DB	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	E	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	EA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	EB	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	F	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	FA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	FB	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	G	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	GA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	GB	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	H	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	HA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	HB	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	I	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	IA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	J	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	JA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	K	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	KA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	L	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	LA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	M	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	MA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	N	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	NA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	O	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	OA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	P	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	PA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	Q	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	QA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	R	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	RA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	S	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	SA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	T	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	TA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	UA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	V	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	VA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	W	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	WA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	X	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	XA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	Y	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	YA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	Z	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	ZA	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
1	a	226/647 (35%)	222 (98%)	4 (2%)	0	100	100
All	All	13560/38820 (35%)	13320 (98%)	240 (2%)	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	A	195/526 (37%)	195 (100%)	0	100	100
1	AA	195/526 (37%)	195 (100%)	0	100	100
1	AB	195/526 (37%)	195 (100%)	0	100	100
1	B	195/526 (37%)	195 (100%)	0	100	100
1	BA	195/526 (37%)	195 (100%)	0	100	100
1	BB	195/526 (37%)	195 (100%)	0	100	100
1	C	195/526 (37%)	195 (100%)	0	100	100
1	CA	195/526 (37%)	195 (100%)	0	100	100
1	CB	195/526 (37%)	195 (100%)	0	100	100
1	D	195/526 (37%)	195 (100%)	0	100	100
1	DA	195/526 (37%)	195 (100%)	0	100	100
1	DB	195/526 (37%)	195 (100%)	0	100	100
1	E	195/526 (37%)	195 (100%)	0	100	100
1	EA	195/526 (37%)	195 (100%)	0	100	100
1	EB	195/526 (37%)	195 (100%)	0	100	100
1	F	195/526 (37%)	195 (100%)	0	100	100
1	FA	195/526 (37%)	195 (100%)	0	100	100
1	FB	195/526 (37%)	195 (100%)	0	100	100
1	G	195/526 (37%)	195 (100%)	0	100	100
1	GA	195/526 (37%)	195 (100%)	0	100	100
1	GB	195/526 (37%)	195 (100%)	0	100	100
1	H	195/526 (37%)	195 (100%)	0	100	100
1	HA	195/526 (37%)	195 (100%)	0	100	100
1	HB	195/526 (37%)	195 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	I	195/526 (37%)	195 (100%)	0	100	100
1	IA	195/526 (37%)	195 (100%)	0	100	100
1	J	195/526 (37%)	195 (100%)	0	100	100
1	JA	195/526 (37%)	195 (100%)	0	100	100
1	K	195/526 (37%)	195 (100%)	0	100	100
1	KA	195/526 (37%)	195 (100%)	0	100	100
1	L	195/526 (37%)	195 (100%)	0	100	100
1	LA	195/526 (37%)	195 (100%)	0	100	100
1	M	195/526 (37%)	195 (100%)	0	100	100
1	MA	195/526 (37%)	195 (100%)	0	100	100
1	N	195/526 (37%)	195 (100%)	0	100	100
1	NA	195/526 (37%)	195 (100%)	0	100	100
1	O	195/526 (37%)	195 (100%)	0	100	100
1	OA	195/526 (37%)	195 (100%)	0	100	100
1	P	195/526 (37%)	195 (100%)	0	100	100
1	PA	195/526 (37%)	195 (100%)	0	100	100
1	Q	195/526 (37%)	195 (100%)	0	100	100
1	QA	195/526 (37%)	195 (100%)	0	100	100
1	R	195/526 (37%)	195 (100%)	0	100	100
1	RA	195/526 (37%)	195 (100%)	0	100	100
1	S	195/526 (37%)	195 (100%)	0	100	100
1	SA	195/526 (37%)	195 (100%)	0	100	100
1	T	195/526 (37%)	195 (100%)	0	100	100
1	TA	195/526 (37%)	195 (100%)	0	100	100
1	UA	195/526 (37%)	195 (100%)	0	100	100
1	V	195/526 (37%)	195 (100%)	0	100	100
1	VA	195/526 (37%)	195 (100%)	0	100	100
1	W	195/526 (37%)	195 (100%)	0	100	100
1	WA	195/526 (37%)	195 (100%)	0	100	100
1	X	195/526 (37%)	195 (100%)	0	100	100
1	XA	195/526 (37%)	195 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	Y	195/526 (37%)	195 (100%)	0	100	100
1	YA	195/526 (37%)	195 (100%)	0	100	100
1	Z	195/526 (37%)	195 (100%)	0	100	100
1	ZA	195/526 (37%)	195 (100%)	0	100	100
1	a	195/526 (37%)	195 (100%)	0	100	100
All	All	11700/31560 (37%)	11700 (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 56 such sidechains are listed below:

Mol	Chain	Res	Type
1	L	137	HIS
1	a	76	ASN
1	OA	76	ASN
1	Z	76	ASN
1	W	137	HIS

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 ⓘ

There are no chain breaks in this entry.

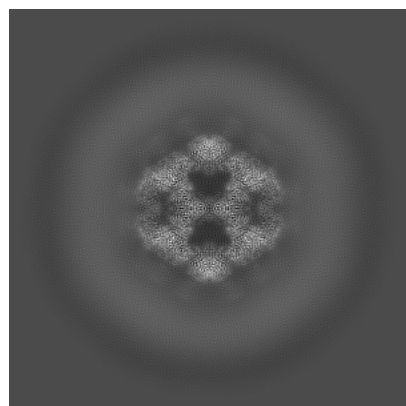
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-61080. These allow visual inspection of the internal detail of the map and identification of artifacts.

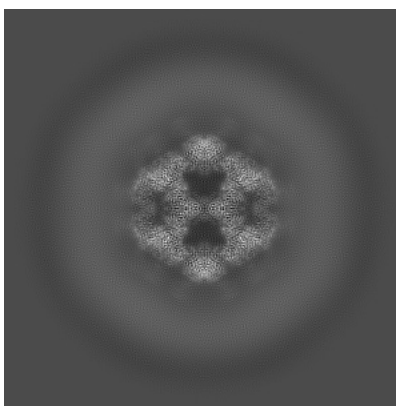
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

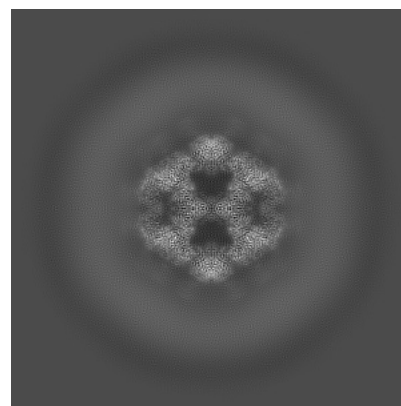
6.1.1 Primary map



X

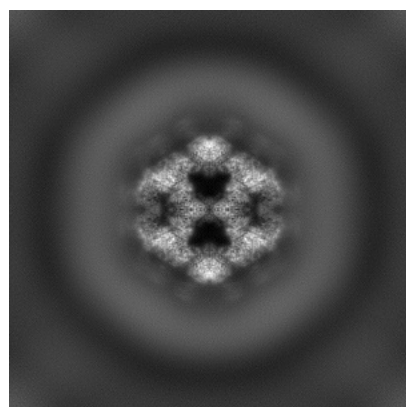


Y

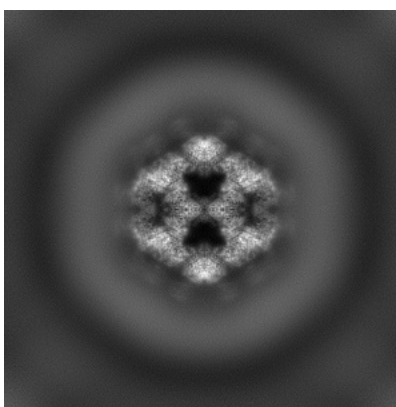


Z

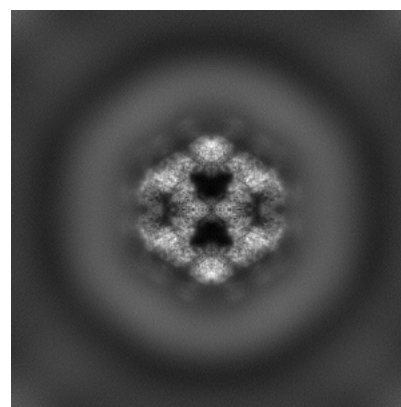
6.1.2 Raw map



X



Y

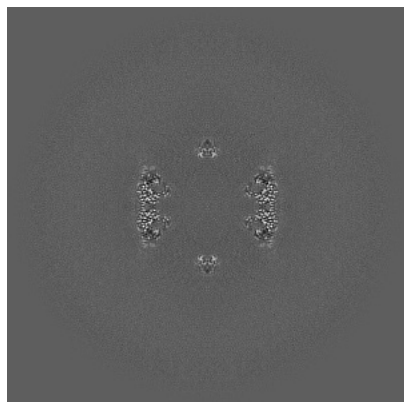


Z

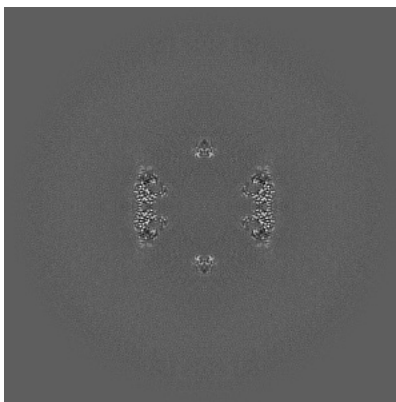
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

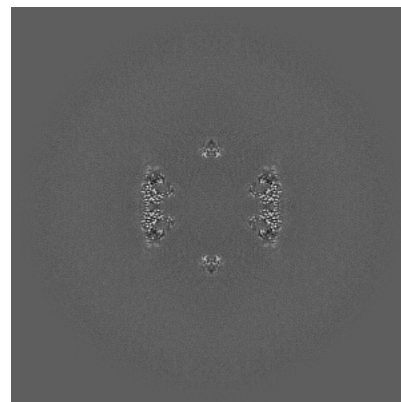
6.2.1 Primary map



X Index: 256

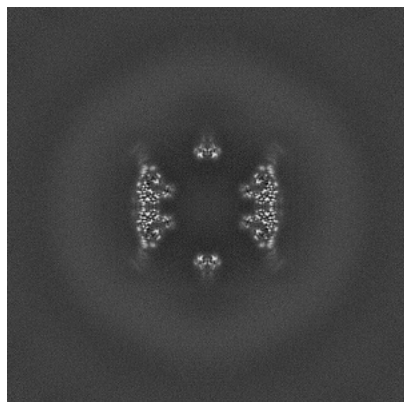


Y Index: 256

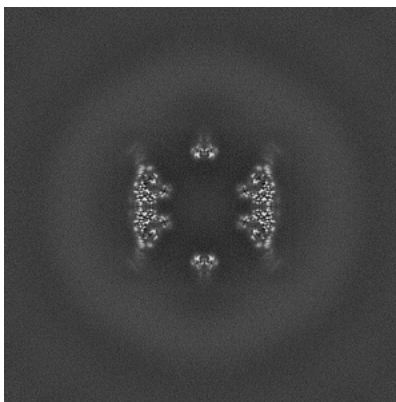


Z Index: 256

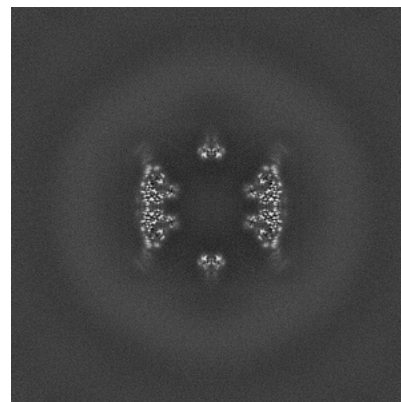
6.2.2 Raw map



X Index: 256



Y Index: 256

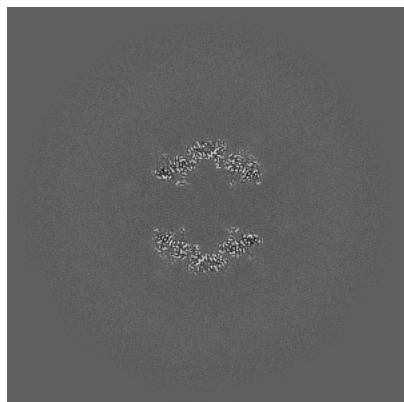


Z Index: 256

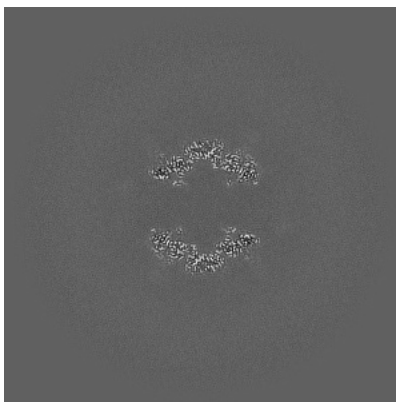
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

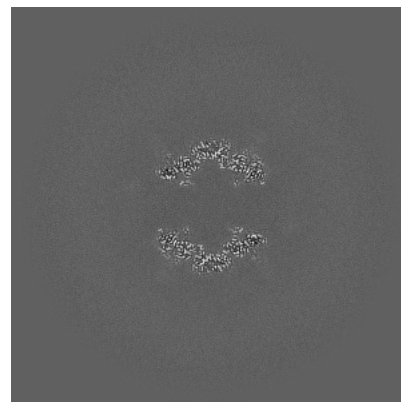
6.3.1 Primary map



X Index: 294

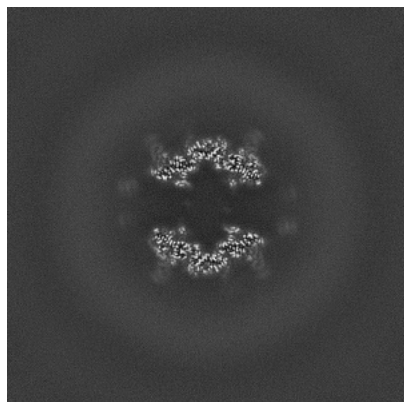


Y Index: 294

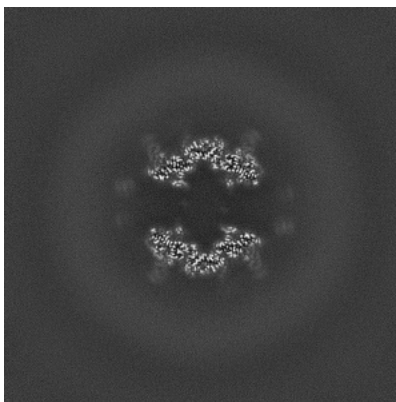


Z Index: 294

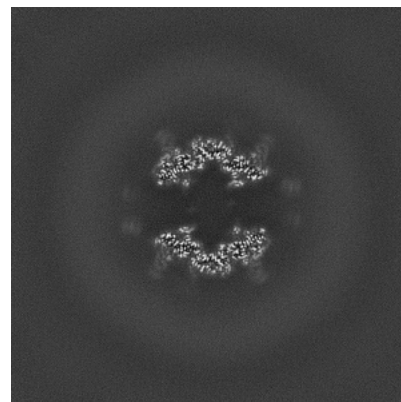
6.3.2 Raw map



X Index: 294



Y Index: 294

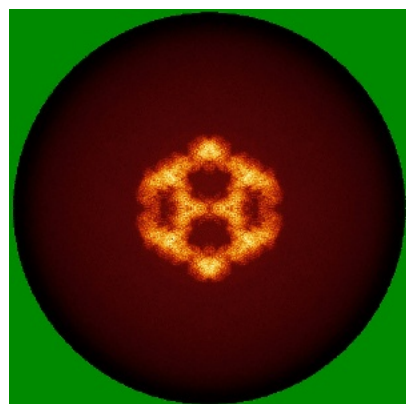


Z Index: 218

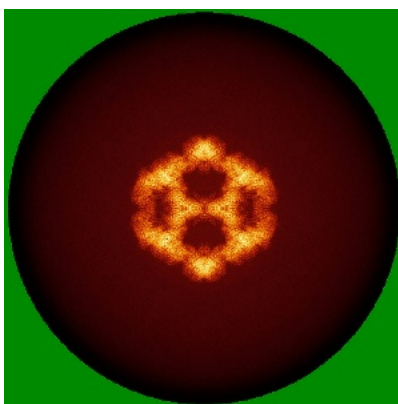
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

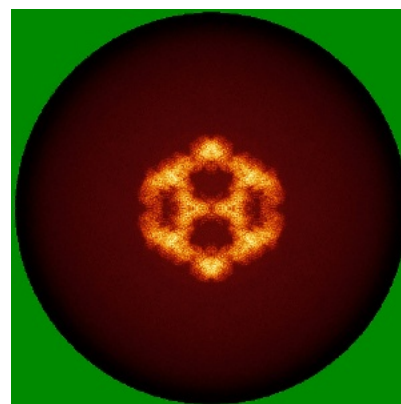
6.4.1 Primary map



X

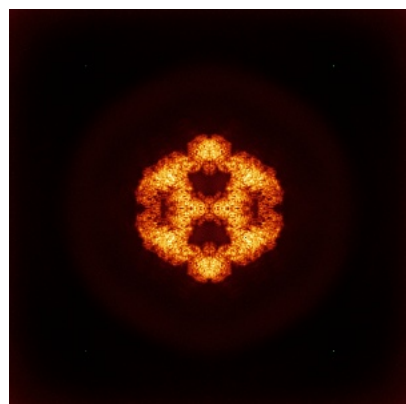


Y

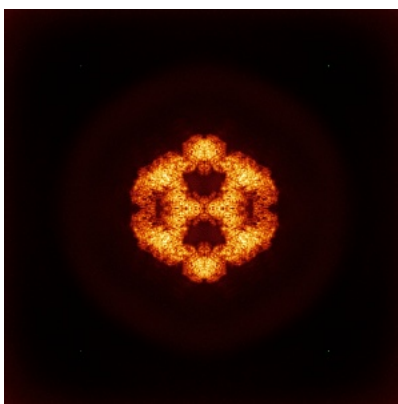


Z

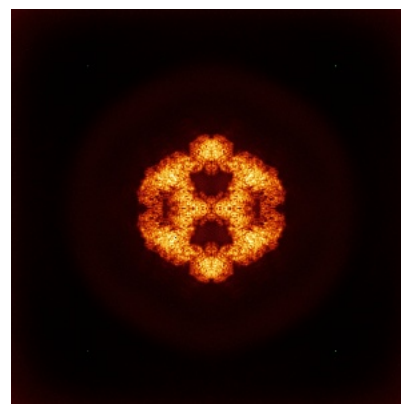
6.4.2 Raw map



X



Y

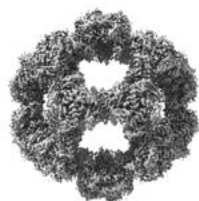


Z

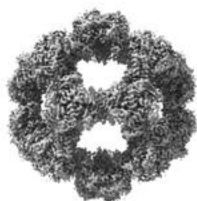
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

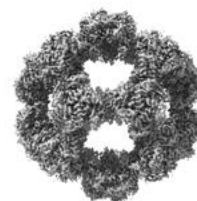
6.5.1 Primary map



X



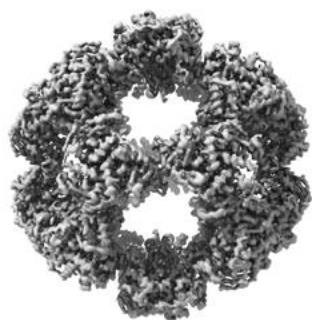
Y



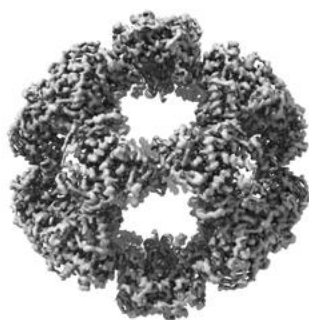
Z

The images above show the 3D surface view of the map at the recommended contour level 0.34. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

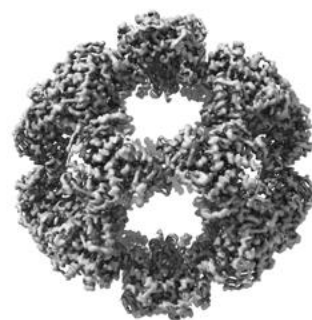
6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

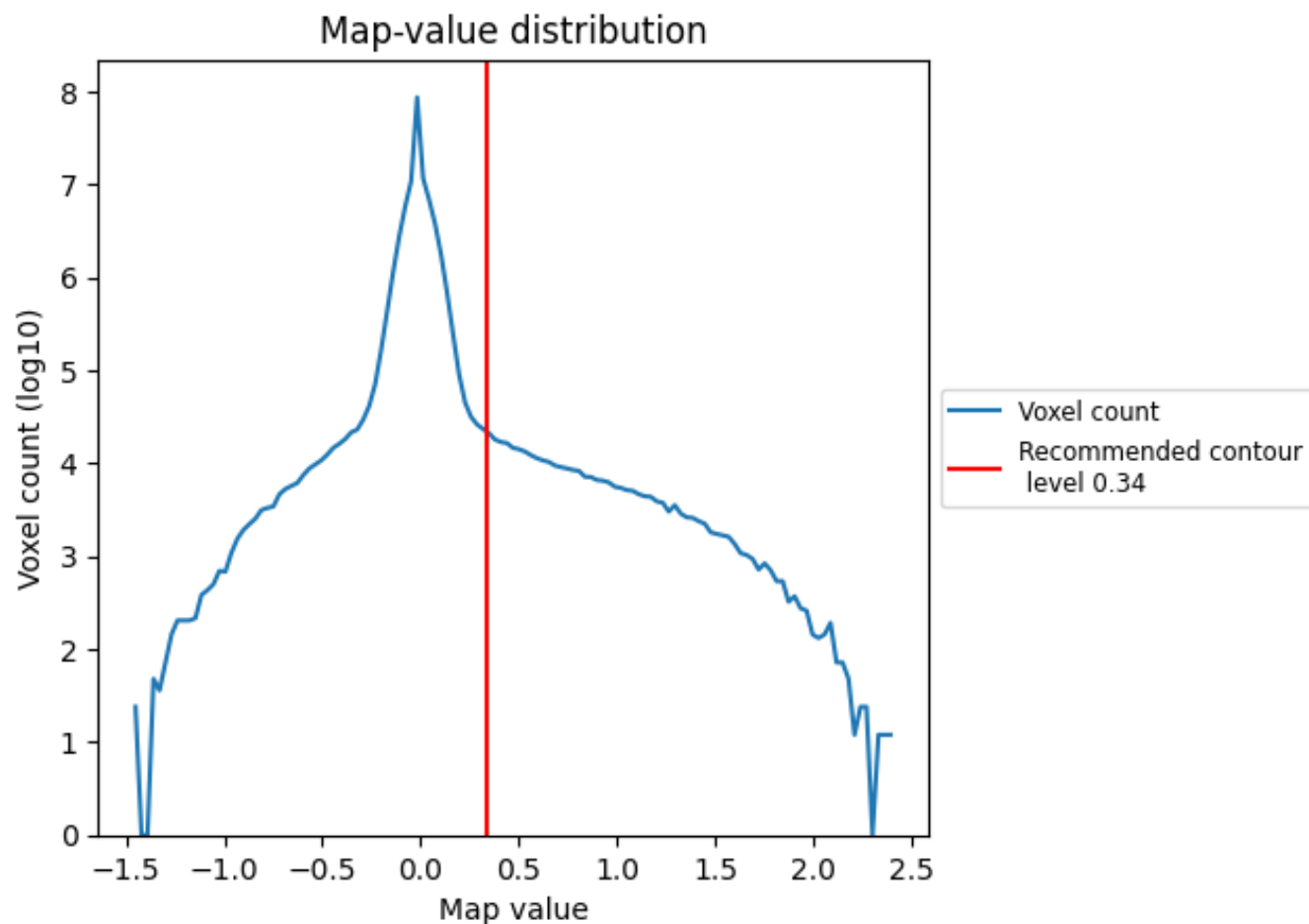
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

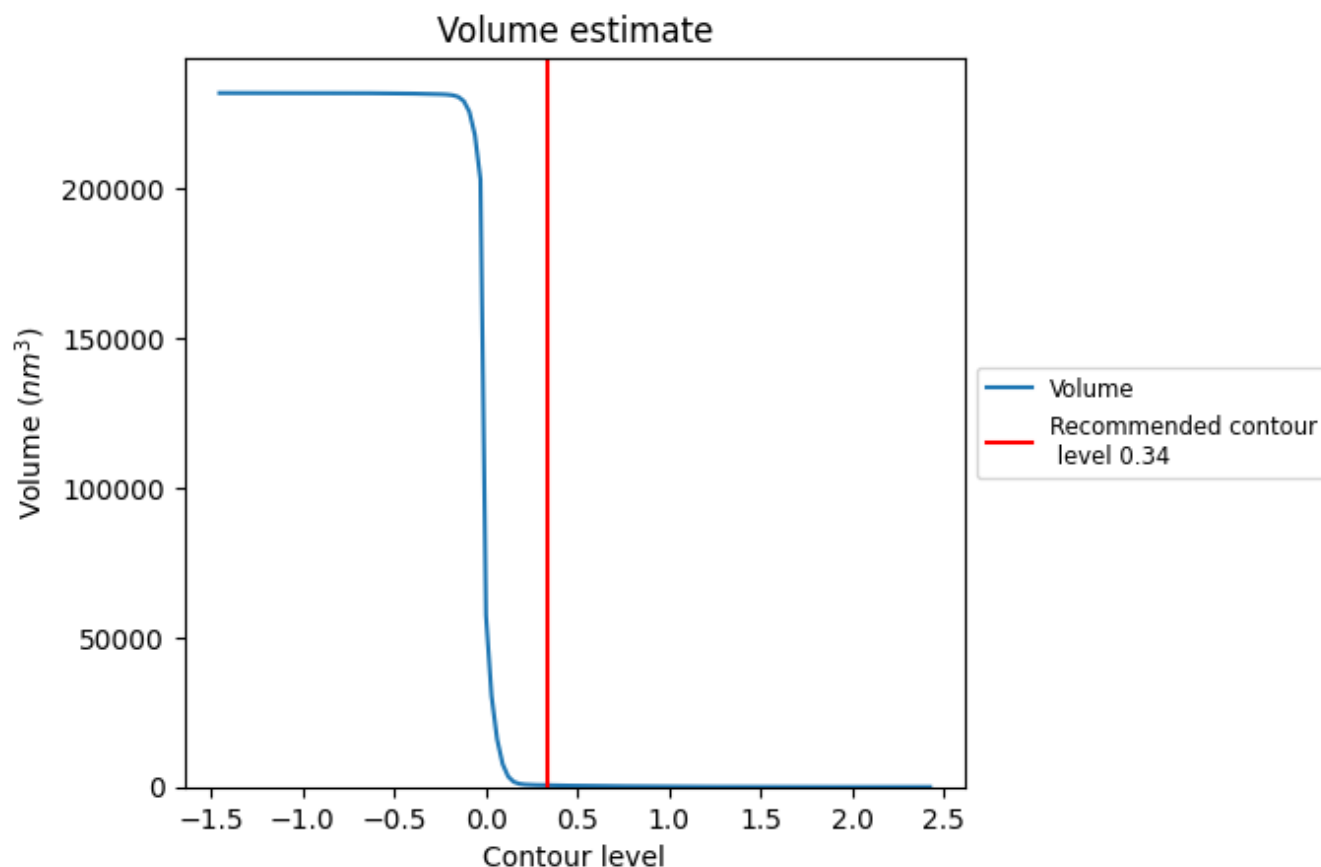
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

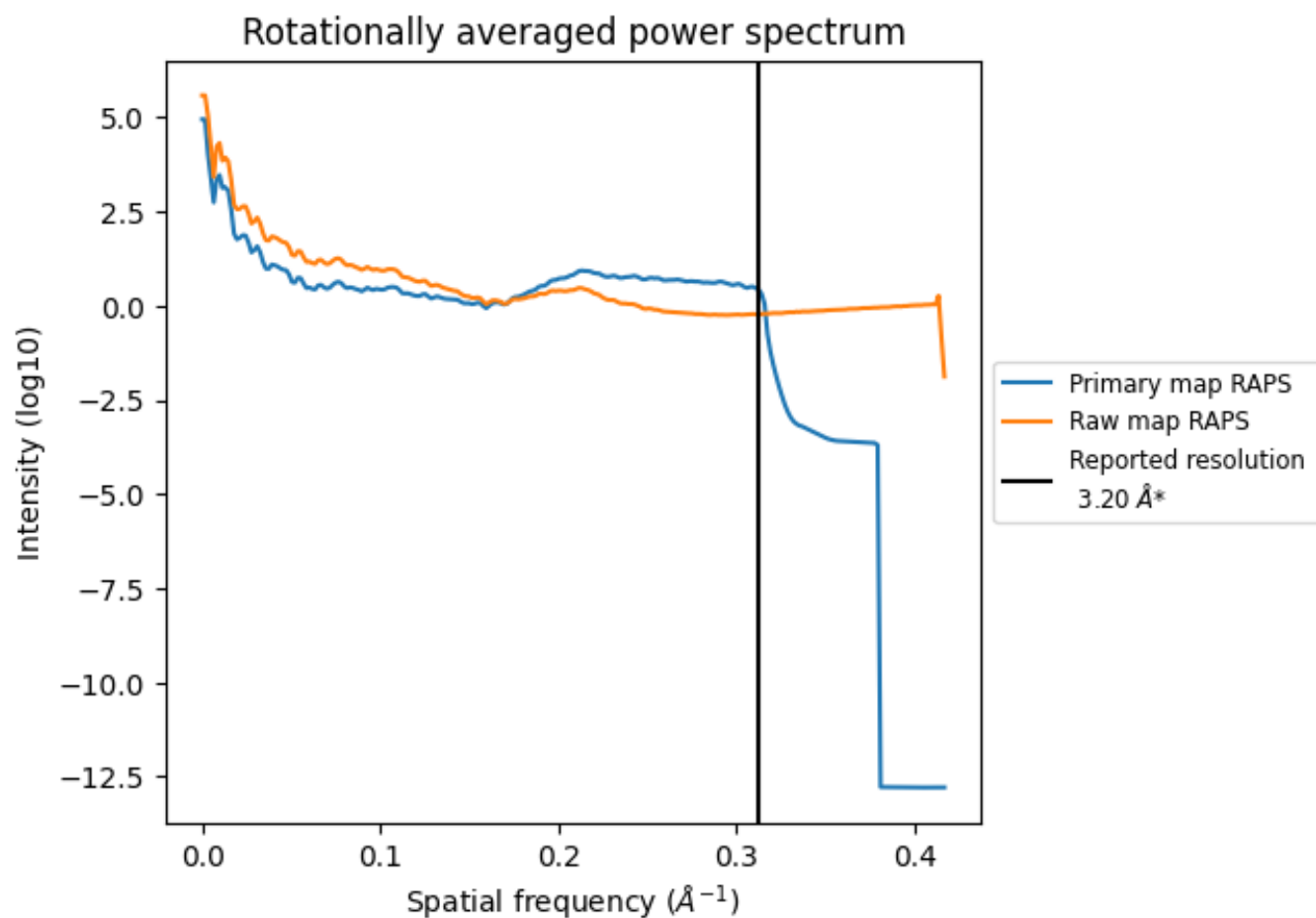
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 558 nm^3 ; this corresponds to an approximate mass of 504 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

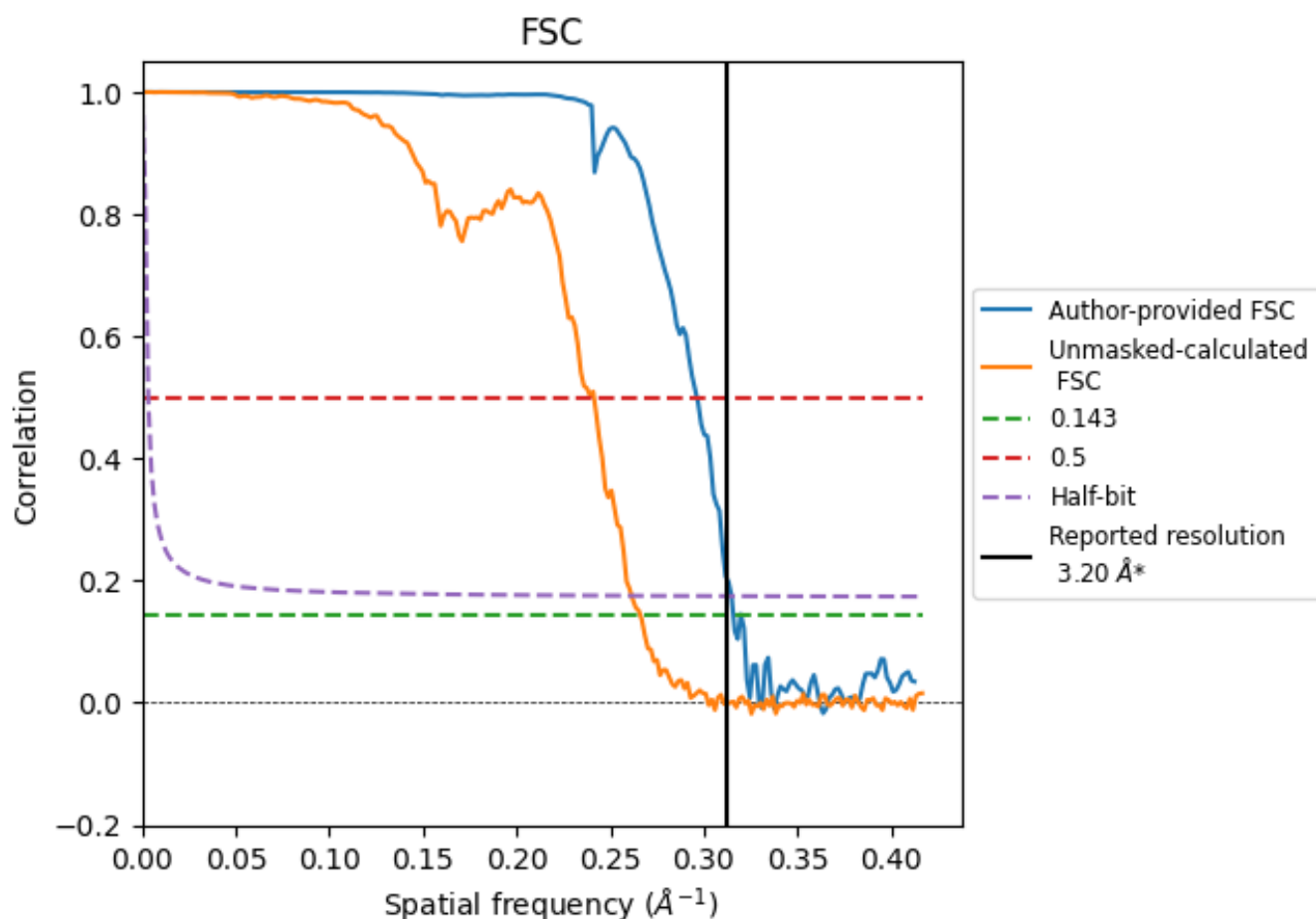


*Reported resolution corresponds to spatial frequency of 0.312 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.312 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

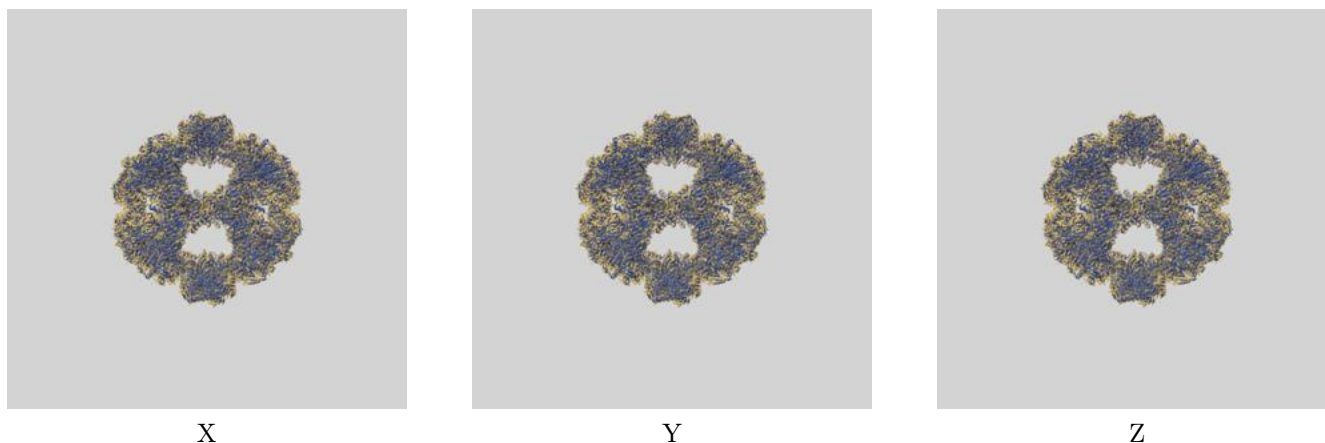
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	3.17	3.37	3.18
Unmasked-calculated*	3.76	4.14	3.82

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.76 differs from the reported value 3.2 by more than 10 %

9 Map-model fit [i](#)

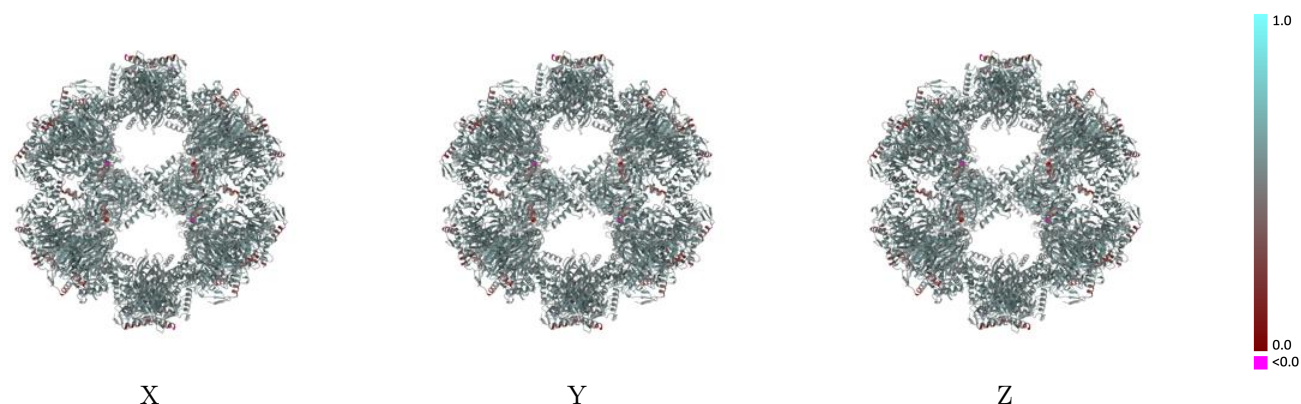
This section contains information regarding the fit between EMDB map EMD-61080 and PDB model 9J1W. Per-residue inclusion information can be found in section [3](#) on page [10](#).

9.1 Map-model overlay [i](#)



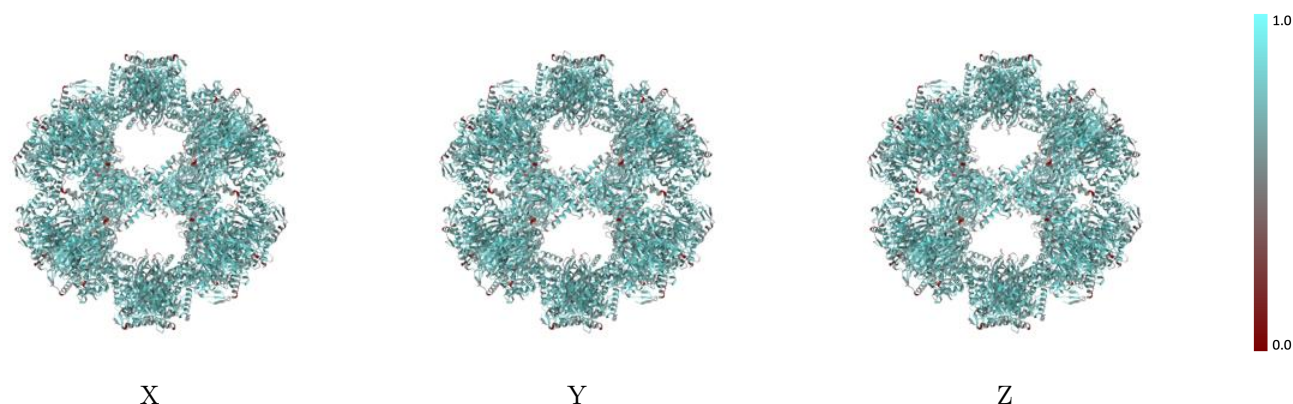
The images above show the 3D surface view of the map at the recommended contour level 0.34 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



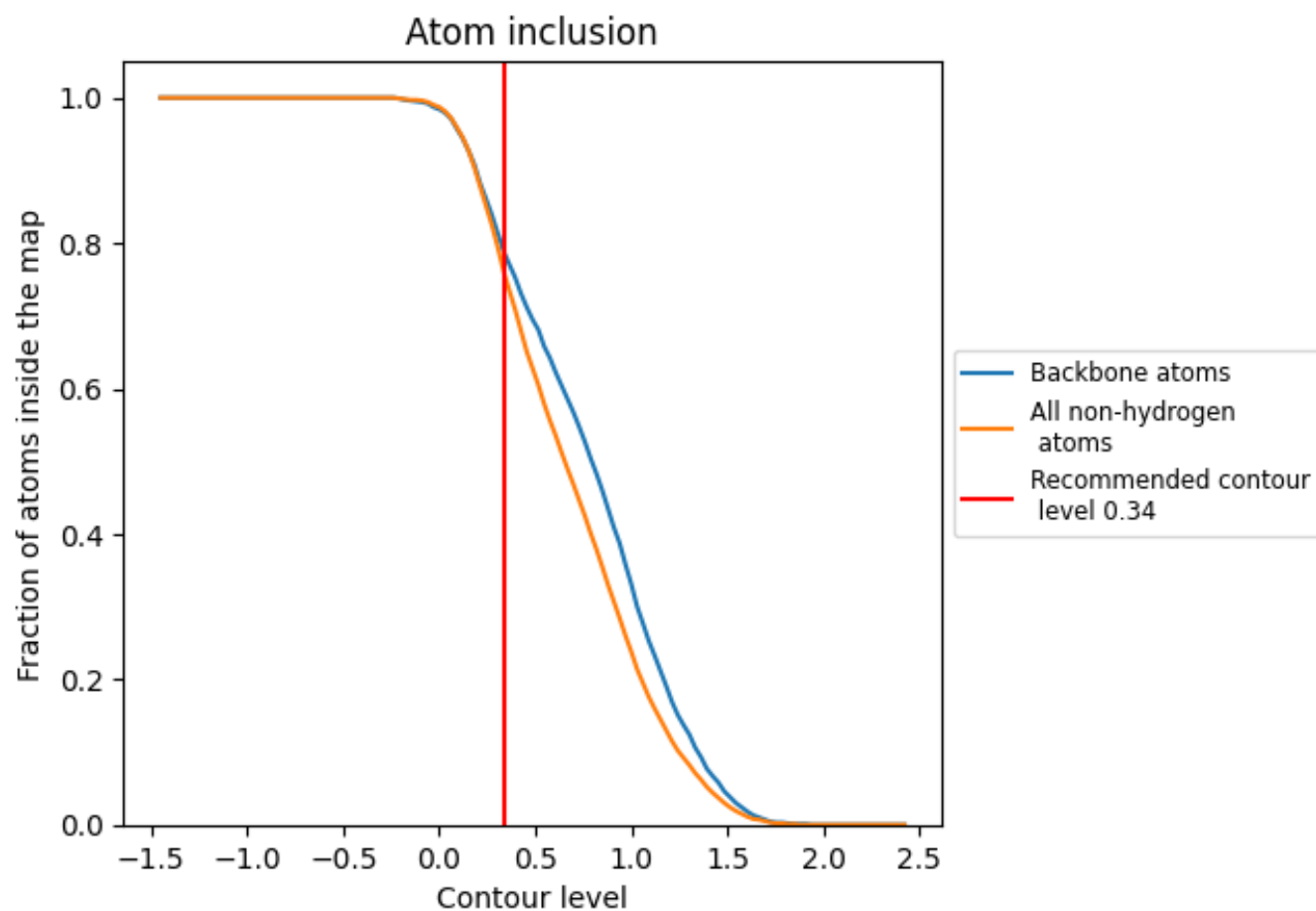
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.34).

9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 76% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.34) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7570	 0.5430
A	 0.7550	 0.5400
AA	 0.7590	 0.5420
AB	 0.7600	 0.5430
B	 0.7600	 0.5460
BA	 0.7590	 0.5450
BB	 0.7570	 0.5420
C	 0.7580	 0.5430
CA	 0.7590	 0.5410
CB	 0.7510	 0.5460
D	 0.7520	 0.5440
DA	 0.7540	 0.5440
DB	 0.7610	 0.5410
E	 0.7610	 0.5420
EA	 0.7610	 0.5410
EB	 0.7560	 0.5420
F	 0.7570	 0.5440
FA	 0.7570	 0.5420
FB	 0.7570	 0.5470
G	 0.7590	 0.5440
GA	 0.7570	 0.5460
GB	 0.7580	 0.5410
H	 0.7560	 0.5410
HA	 0.7570	 0.5420
HB	 0.7520	 0.5450
I	 0.7510	 0.5440
IA	 0.7500	 0.5430
J	 0.7600	 0.5420
JA	 0.7610	 0.5410
K	 0.7590	 0.5420
KA	 0.7570	 0.5420
L	 0.7610	 0.5450
LA	 0.7600	 0.5450
M	 0.7570	 0.5420
MA	 0.7580	 0.5410



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Chain	Atom inclusion	Q-score
N	 0.7530	 0.5440
NA	 0.7510	 0.5420
O	 0.7610	 0.5430
OA	 0.7610	 0.5400
P	 0.7590	 0.5420
PA	 0.7570	 0.5430
Q	 0.7600	 0.5450
QA	 0.7570	 0.5440
R	 0.7580	 0.5440
RA	 0.7570	 0.5410
S	 0.7510	 0.5450
SA	 0.7530	 0.5430
T	 0.7610	 0.5430
TA	 0.7610	 0.5410
UA	 0.7580	 0.5410
V	 0.7570	 0.5410
VA	 0.7600	 0.5450
W	 0.7600	 0.5440
WA	 0.7580	 0.5420
X	 0.7580	 0.5400
XA	 0.7520	 0.5470
Y	 0.7510	 0.5440
YA	 0.7610	 0.5400
Z	 0.7600	 0.5440
ZA	 0.7570	 0.5410
a	 0.7610	 0.5440