

wwPDB EM Map/Model Validation Report i

Aug 23, 2016 – 03:12 PM EDT

PDB ID : 5T0V
EMDB ID: : EMD-8341
Title : Architecture of the Yeast Mitochondrial Iron-Sulfur Cluster Assembly Machinery: the Sub-Complex Formed by the Iron Donor, Yfh1, and the Scaffold, Isu1
Authors : Ranatunga, W.; Gakh, O.; Galeano, B.K.; Smith IV, D.Y.; Soderberg, C.A.; Al-Karadaghi, S.; Thompson, J.R.; Isaya, G.
Deposited on : 2016-08-16
Resolution : 17.50 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>

MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027939

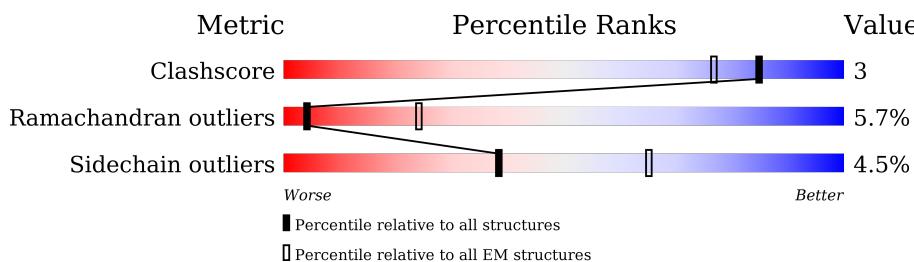
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

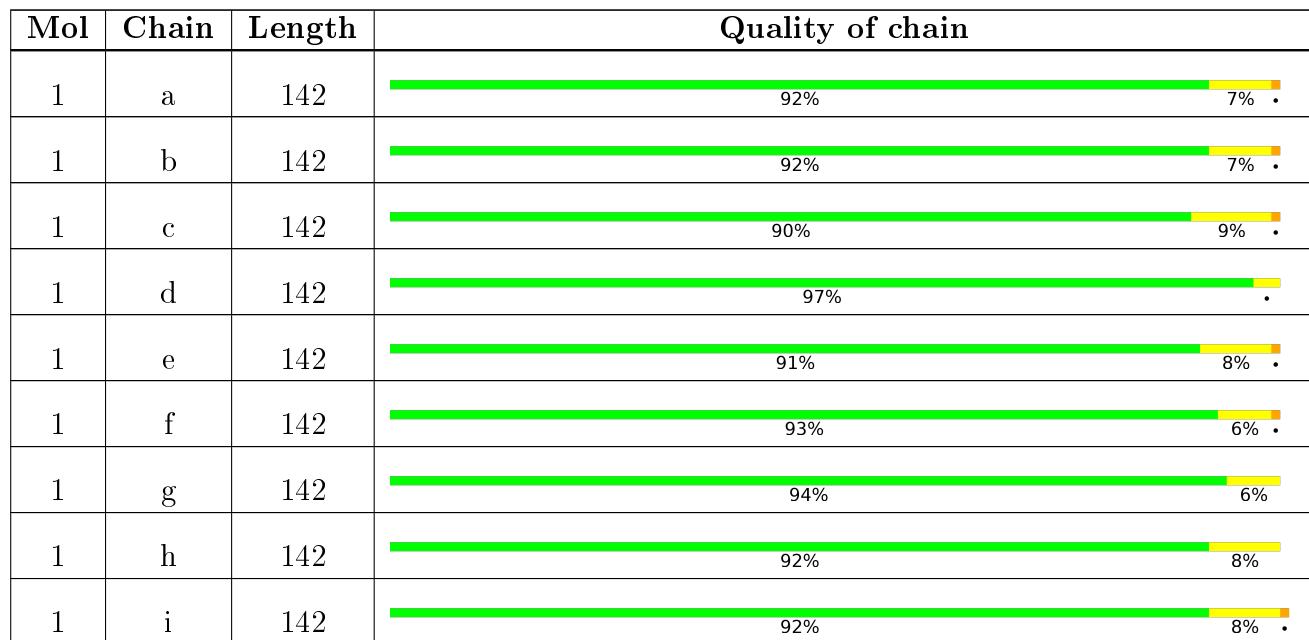
The reported resolution of this entry is 17.50 Å.

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	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=5%



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Mol	Chain	Length	Quality of chain		
1	j	142	92%	8%	.
1	k	142	90%	8%	.
1	l	142	89%	11%	
1	m	142	94%	6%	
1	n	142	93%	6%	.
1	o	142	94%	6%	
1	p	142	88%	12%	
1	q	142	91%	9%	
1	r	142	91%	8%	.
1	s	142	94%	6%	
1	t	142	91%	6%	.
1	u	142	94%	6%	
1	v	142	94%	6%	
1	w	142	94%	6%	
1	x	142	92%	7%	.
2	A	121	81%	15%	.
2	B	121	87%	12%	.
2	C	121	83%	14%	.
2	D	121	80%	18%	.
2	E	121	83%	13%	.
2	F	121	85%	14%	.
2	G	121	80%	17%	.
2	H	121	81%	16%	.
2	I	121	88%	12%	
2	J	121	81%	17%	..

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Mol	Chain	Length	Quality of chain		
2	K	121	79%	20%	.
2	L	121	84%	15%	.
2	M	121	87%	10%	.
2	N	121	76%	21%	.
2	O	121	88%	11%	.
2	P	121	83%	17%	
2	Q	121	85%	12%	.
2	R	121	88%	11%	.
2	S	121	85%	11%	..
2	T	121	86%	12%	.
2	U	121	77%	19%	.
2	V	121	84%	13%	.
2	W	121	79%	19%	.
2	X	121	90%	8%	.

2 Entry composition [\(i\)](#)

There are 2 unique types of molecules in this entry. The entry contains 48456 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 Iron sulfur cluster assembly protein 1, mitochondrial.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	a	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	b	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	c	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	d	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	e	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	f	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	g	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	h	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	i	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	j	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	k	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	l	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	m	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	n	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	o	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	p	142	Total	C	N	O	S	
			1072	672	186	205	9	0
1	q	142	Total	C	N	O	S	
			1072	672	186	205	9	0

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	r	142	Total	C	N	O	S		
			1072	672	186	205	9	0	0
1	s	142	Total	C	N	O	S		
			1072	672	186	205	9	0	0
1	t	142	Total	C	N	O	S		
			1072	672	186	205	9	0	0
1	u	142	Total	C	N	O	S		
			1072	672	186	205	9	0	0
1	v	142	Total	C	N	O	S		
			1072	672	186	205	9	0	0
1	w	142	Total	C	N	O	S		
			1072	672	186	205	9	0	0
1	x	142	Total	C	N	O	S		
			1072	672	186	205	9	0	0

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
a	24	GLY	-	expression tag	UNP Q03020
a	25	SER	-	expression tag	UNP Q03020
a	26	HIS	-	expression tag	UNP Q03020
a	27	MET	-	expression tag	UNP Q03020
b	24	GLY	-	expression tag	UNP Q03020
b	25	SER	-	expression tag	UNP Q03020
b	26	HIS	-	expression tag	UNP Q03020
b	27	MET	-	expression tag	UNP Q03020
c	24	GLY	-	expression tag	UNP Q03020
c	25	SER	-	expression tag	UNP Q03020
c	26	HIS	-	expression tag	UNP Q03020
c	27	MET	-	expression tag	UNP Q03020
d	24	GLY	-	expression tag	UNP Q03020
d	25	SER	-	expression tag	UNP Q03020
d	26	HIS	-	expression tag	UNP Q03020
d	27	MET	-	expression tag	UNP Q03020
e	24	GLY	-	expression tag	UNP Q03020
e	25	SER	-	expression tag	UNP Q03020
e	26	HIS	-	expression tag	UNP Q03020
e	27	MET	-	expression tag	UNP Q03020
f	24	GLY	-	expression tag	UNP Q03020
f	25	SER	-	expression tag	UNP Q03020
f	26	HIS	-	expression tag	UNP Q03020
f	27	MET	-	expression tag	UNP Q03020
g	24	GLY	-	expression tag	UNP Q03020

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Chain	Residue	Modelled	Actual	Comment	Reference
g	25	SER	-	expression tag	UNP Q03020
g	26	HIS	-	expression tag	UNP Q03020
g	27	MET	-	expression tag	UNP Q03020
h	24	GLY	-	expression tag	UNP Q03020
h	25	SER	-	expression tag	UNP Q03020
h	26	HIS	-	expression tag	UNP Q03020
h	27	MET	-	expression tag	UNP Q03020
i	24	GLY	-	expression tag	UNP Q03020
i	25	SER	-	expression tag	UNP Q03020
i	26	HIS	-	expression tag	UNP Q03020
i	27	MET	-	expression tag	UNP Q03020
j	24	GLY	-	expression tag	UNP Q03020
j	25	SER	-	expression tag	UNP Q03020
j	26	HIS	-	expression tag	UNP Q03020
j	27	MET	-	expression tag	UNP Q03020
k	24	GLY	-	expression tag	UNP Q03020
k	25	SER	-	expression tag	UNP Q03020
k	26	HIS	-	expression tag	UNP Q03020
k	27	MET	-	expression tag	UNP Q03020
l	24	GLY	-	expression tag	UNP Q03020
l	25	SER	-	expression tag	UNP Q03020
l	26	HIS	-	expression tag	UNP Q03020
l	27	MET	-	expression tag	UNP Q03020
m	24	GLY	-	expression tag	UNP Q03020
m	25	SER	-	expression tag	UNP Q03020
m	26	HIS	-	expression tag	UNP Q03020
m	27	MET	-	expression tag	UNP Q03020
n	24	GLY	-	expression tag	UNP Q03020
n	25	SER	-	expression tag	UNP Q03020
n	26	HIS	-	expression tag	UNP Q03020
n	27	MET	-	expression tag	UNP Q03020
o	24	GLY	-	expression tag	UNP Q03020
o	25	SER	-	expression tag	UNP Q03020
o	26	HIS	-	expression tag	UNP Q03020
o	27	MET	-	expression tag	UNP Q03020
p	24	GLY	-	expression tag	UNP Q03020
p	25	SER	-	expression tag	UNP Q03020
p	26	HIS	-	expression tag	UNP Q03020
p	27	MET	-	expression tag	UNP Q03020
q	24	GLY	-	expression tag	UNP Q03020
q	25	SER	-	expression tag	UNP Q03020
q	26	HIS	-	expression tag	UNP Q03020

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Chain	Residue	Modelled	Actual	Comment	Reference
q	27	MET	-	expression tag	UNP Q03020
r	24	GLY	-	expression tag	UNP Q03020
r	25	SER	-	expression tag	UNP Q03020
r	26	HIS	-	expression tag	UNP Q03020
r	27	MET	-	expression tag	UNP Q03020
s	24	GLY	-	expression tag	UNP Q03020
s	25	SER	-	expression tag	UNP Q03020
s	26	HIS	-	expression tag	UNP Q03020
s	27	MET	-	expression tag	UNP Q03020
t	24	GLY	-	expression tag	UNP Q03020
t	25	SER	-	expression tag	UNP Q03020
t	26	HIS	-	expression tag	UNP Q03020
t	27	MET	-	expression tag	UNP Q03020
u	24	GLY	-	expression tag	UNP Q03020
u	25	SER	-	expression tag	UNP Q03020
u	26	HIS	-	expression tag	UNP Q03020
u	27	MET	-	expression tag	UNP Q03020
v	24	GLY	-	expression tag	UNP Q03020
v	25	SER	-	expression tag	UNP Q03020
v	26	HIS	-	expression tag	UNP Q03020
v	27	MET	-	expression tag	UNP Q03020
w	24	GLY	-	expression tag	UNP Q03020
w	25	SER	-	expression tag	UNP Q03020
w	26	HIS	-	expression tag	UNP Q03020
w	27	MET	-	expression tag	UNP Q03020
x	24	GLY	-	expression tag	UNP Q03020
x	25	SER	-	expression tag	UNP Q03020
x	26	HIS	-	expression tag	UNP Q03020
x	27	MET	-	expression tag	UNP Q03020

- Molecule 2 is a protein called Frataxin homolog, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	121	Total	C	N	O	S	0	0
			947	597	153	195	2		
2	B	121	Total	C	N	O	S	0	0
			947	597	153	195	2		
2	C	121	Total	C	N	O	S	0	0
			947	597	153	195	2		
2	D	121	Total	C	N	O	S	0	0
			947	597	153	195	2		
2	E	121	Total	C	N	O	S	0	0
			947	597	153	195	2		

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	F	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	G	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	H	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	I	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	J	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	K	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	L	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	M	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	N	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	O	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	P	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	Q	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	R	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	S	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	T	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	U	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	V	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	W	121	Total	C	N	O	S		
			947	597	153	195	2	0	0
2	X	121	Total	C	N	O	S		
			947	597	153	195	2	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	73	ALA	TYR	conflict	UNP Q07540

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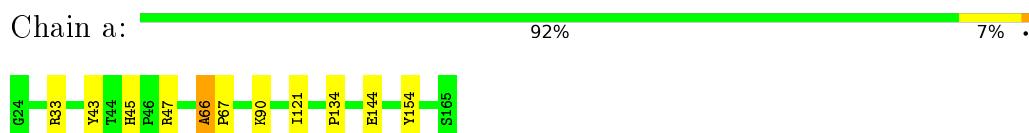
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Chain	Residue	Modelled	Actual	Comment	Reference
B	73	ALA	TYR	conflict	UNP Q07540
C	73	ALA	TYR	conflict	UNP Q07540
D	73	ALA	TYR	conflict	UNP Q07540
E	73	ALA	TYR	conflict	UNP Q07540
F	73	ALA	TYR	conflict	UNP Q07540
G	73	ALA	TYR	conflict	UNP Q07540
H	73	ALA	TYR	conflict	UNP Q07540
I	73	ALA	TYR	conflict	UNP Q07540
J	73	ALA	TYR	conflict	UNP Q07540
K	73	ALA	TYR	conflict	UNP Q07540
L	73	ALA	TYR	conflict	UNP Q07540
M	73	ALA	TYR	conflict	UNP Q07540
N	73	ALA	TYR	conflict	UNP Q07540
O	73	ALA	TYR	conflict	UNP Q07540
P	73	ALA	TYR	conflict	UNP Q07540
Q	73	ALA	TYR	conflict	UNP Q07540
R	73	ALA	TYR	conflict	UNP Q07540
S	73	ALA	TYR	conflict	UNP Q07540
T	73	ALA	TYR	conflict	UNP Q07540
U	73	ALA	TYR	conflict	UNP Q07540
V	73	ALA	TYR	conflict	UNP Q07540
W	73	ALA	TYR	conflict	UNP Q07540
X	73	ALA	TYR	conflict	UNP Q07540

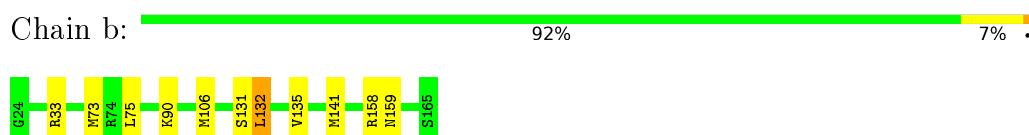
3 Residue-property plots [\(i\)](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

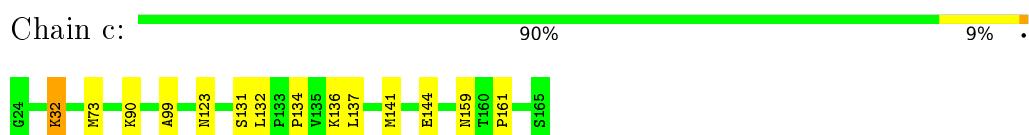
- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial



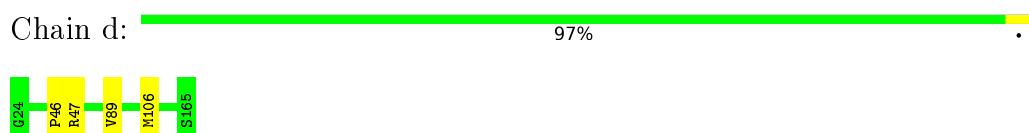
- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial





- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain g: 94% 6%



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain h: 92% 8%



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain i: 92% 8%



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain j: 92% 8%



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain k: 90% 8%



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain l: 89% 11%



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain m: 94% 6%



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain n: 93% 6%

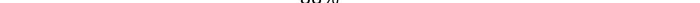


- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain o: 94% 6%

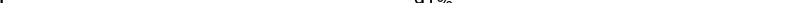


- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain p:  88% 12%

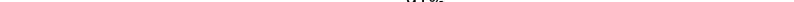


- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain q:  91% 9%

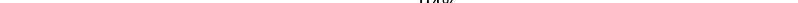


- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain r:  91% 8% ;



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chains:  94% 6%

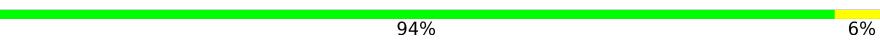


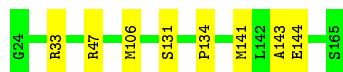
- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain t: 91% 6% :

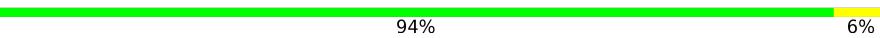


- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain u: 

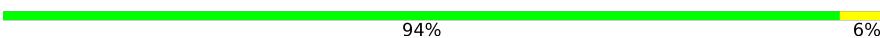


- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain v: 



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain w: 



- Molecule 1: Iron sulfur cluster assembly protein 1, mitochondrial

Chain x: 



- Molecule 2: Frataxin homolog, mitochondrial

Chain A: 



- Molecule 2: Frataxin homolog, mitochondrial

Chain B: 



- Molecule 2: Frataxin homolog, mitochondrial

Chain C: 

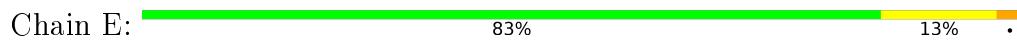


- Molecule 2: Frataxin homolog, mitochondrial

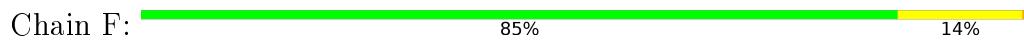
Chain D: 



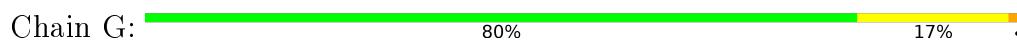
- Molecule 2: Frataxin homolog, mitochondrial



- Molecule 2: Frataxin homolog, mitochondrial



- Molecule 2: Frataxin homolog, mitochondrial



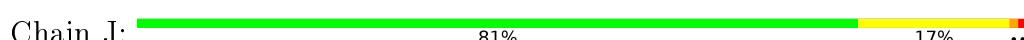
- Molecule 2: Frataxin homolog, mitochondrial



- Molecule 2: Frataxin homolog, mitochondrial



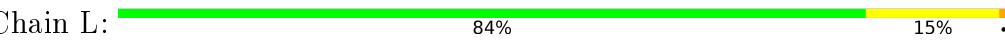
- Molecule 2: Frataxin homolog, mitochondrial

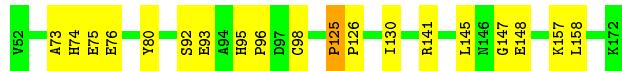


- Molecule 2: Frataxin homolog, mitochondrial

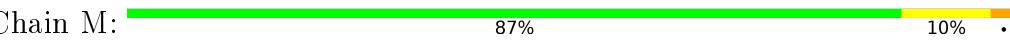


- Molecule 2: Frataxin homolog, mitochondrial

Chain L:  84% 15%



- Molecule 2: Frataxin homolog, mitochondrial

Chain M:  87% 10%

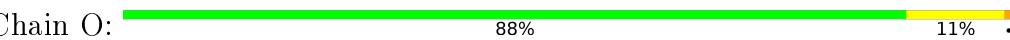


- Molecule 2: Frataxin homolog, mitochondrial

Chain N:  76% 21%

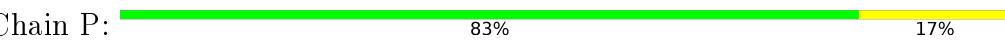


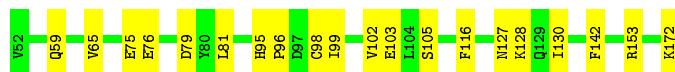
- Molecule 2: Frataxin homolog, mitochondrial

Chain O:  88% 11%

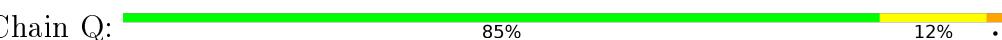


- Molecule 2: Frataxin homolog, mitochondrial

Chain P:  83% 17%



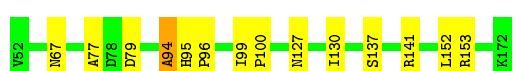
- Molecule 2: Frataxin homolog, mitochondrial

Chain Q:  85% 12%



- Molecule 2: Frataxin homolog, mitochondrial

Chain R:  88% 11%



- Molecule 2: Frataxin homolog, mitochondrial

Chain S:  85% 11% ..



- Molecule 2: Frataxin homolog, mitochondrial

Chain T:  86% 12% ..



- Molecule 2: Frataxin homolog, mitochondrial

Chain U:  77% 19% ..



- Molecule 2: Frataxin homolog, mitochondrial

Chain V:  84% 13% ..



- Molecule 2: Frataxin homolog, mitochondrial

Chain W:  79% 19% ..



- Molecule 2: Frataxin homolog, mitochondrial

Chain X:  90% 8% ..



4 Experimental information (i)

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of particles used	4153	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	Not provided	Depositor
Microscope	FEI TECNAI F30	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	Not provided	Depositor
Minimum defocus (nm)	210	Depositor
Maximum defocus (nm)	2800	Depositor
Magnification	115000	Depositor
Image detector	Not provided	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 > 2$	RMSZ	# $ Z > 2$
1	a	1.04	0/1089	1.23	3/1466 (0.2%)
1	b	1.04	0/1089	1.28	6/1466 (0.4%)
1	c	1.02	0/1089	1.24	2/1466 (0.1%)
1	d	1.05	0/1089	1.18	0/1466
1	e	1.03	0/1089	1.29	5/1466 (0.3%)
1	f	1.06	0/1089	1.24	1/1466 (0.1%)
1	g	1.04	0/1089	1.18	1/1466 (0.1%)
1	h	1.00	0/1089	1.27	4/1466 (0.3%)
1	i	1.03	0/1089	1.19	1/1466 (0.1%)
1	j	1.02	0/1089	1.26	1/1466 (0.1%)
1	k	1.04	0/1089	1.27	6/1466 (0.4%)
1	l	1.04	0/1089	1.32	5/1466 (0.3%)
1	m	1.01	0/1089	1.17	0/1466
1	n	0.98	0/1089	1.22	3/1466 (0.2%)
1	o	0.98	0/1089	1.18	2/1466 (0.1%)
1	p	1.04	0/1089	1.27	5/1466 (0.3%)
1	q	1.04	0/1089	1.21	2/1466 (0.1%)
1	r	1.01	0/1089	1.24	1/1466 (0.1%)
1	s	1.05	0/1089	1.22	2/1466 (0.1%)
1	t	1.07	0/1089	1.24	2/1466 (0.1%)
1	u	1.05	0/1089	1.18	0/1466
1	v	1.04	0/1089	1.22	3/1466 (0.2%)
1	w	1.05	0/1089	1.22	0/1466
1	x	1.06	0/1089	1.25	3/1466 (0.2%)
2	A	0.99	0/967	1.21	2/1319 (0.2%)
2	B	1.02	0/967	1.24	0/1319
2	C	1.03	0/967	1.28	2/1319 (0.2%)
2	D	1.02	0/967	1.28	5/1319 (0.4%)
2	E	1.01	0/967	1.27	4/1319 (0.3%)
2	F	1.06	0/967	1.30	3/1319 (0.2%)
2	G	1.02	0/967	1.25	1/1319 (0.1%)
2	H	1.01	0/967	1.25	4/1319 (0.3%)
2	I	1.02	0/967	1.23	1/1319 (0.1%)
2	J	1.05	0/967	1.27	2/1319 (0.2%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >2	RMSZ	# Z >2
2	K	0.98	0/967	1.27	3/1319 (0.2%)
2	L	1.01	0/967	1.23	1/1319 (0.1%)
2	M	1.02	0/967	1.26	3/1319 (0.2%)
2	N	1.00	0/967	1.32	7/1319 (0.5%)
2	O	1.01	0/967	1.19	1/1319 (0.1%)
2	P	1.03	0/967	1.31	4/1319 (0.3%)
2	Q	1.01	0/967	1.29	5/1319 (0.4%)
2	R	1.02	0/967	1.26	3/1319 (0.2%)
2	S	1.01	0/967	1.29	3/1319 (0.2%)
2	T	1.05	0/967	1.20	3/1319 (0.2%)
2	U	1.02	0/967	1.29	2/1319 (0.2%)
2	V	1.04	0/967	1.29	2/1319 (0.2%)
2	W	1.05	0/967	1.26	4/1319 (0.3%)
2	X	0.99	0/967	1.23	2/1319 (0.2%)
All	All	1.03	0/49344	1.25	125/66840 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	a	0	1
1	n	0	1
1	r	0	2
1	t	0	1
2	I	0	1
2	M	0	1
All	All	0	7

There are no bond length outliers.

The worst 5 of 125 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	119	TYR	CB-CG-CD2	-10.14	114.91	121.00
2	C	119	TYR	CB-CG-CD1	10.11	127.07	121.00
1	l	105	TYR	CB-CG-CD2	-9.20	115.48	121.00
1	x	68	ALA	N-CA-CB	9.11	122.86	110.10
1	p	154	TYR	CB-CG-CD2	-8.95	115.63	121.00

There are no chirality outliers.

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	I	119	TYR	Sidechain
2	M	119	TYR	Sidechain
1	a	33	ARG	Sidechain
1	n	105	TYR	Sidechain
1	r	33	ARG	Sidechain

5.2 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 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	1072	0	1106	0	0
1	b	1072	0	1108	0	0
1	c	1072	0	1108	0	0
1	d	1072	0	1108	0	0
1	e	1072	0	1108	0	0
1	f	1072	0	1108	0	0
1	g	1072	0	1108	0	0
1	h	1072	0	1108	0	0
1	i	1072	0	1106	0	0
1	j	1072	0	1108	0	0
1	k	1072	0	1106	0	0
1	l	1072	0	1108	0	0
1	m	1072	0	1108	0	0
1	n	1072	0	1106	0	0
1	o	1072	0	1108	0	0
1	p	1072	0	1108	0	0
1	q	1072	0	1108	0	0
1	r	1072	0	1105	0	0
1	s	1072	0	1106	0	0
1	t	1072	0	1108	0	0
1	u	1072	0	1108	0	0
1	v	1072	0	1108	0	0
1	w	1072	0	1106	0	0
1	x	1072	0	1108	0	0
2	A	947	0	921	2	0
2	B	947	0	921	2	0
2	C	947	0	921	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	947	0	921	1	0
2	E	947	0	921	3	0
2	F	947	0	921	1	0
2	G	947	0	921	2	0
2	H	947	0	921	3	0
2	I	947	0	921	1	0
2	J	947	0	921	3	0
2	K	947	0	921	2	0
2	L	947	0	920	19	0
2	M	947	0	921	5	0
2	N	947	0	920	26	0
2	O	947	0	920	17	0
2	P	947	0	921	2	0
2	Q	947	0	921	1	0
2	R	947	0	921	2	0
2	S	947	0	921	6	0
2	T	947	0	921	2	0
2	U	947	0	921	6	0
2	V	947	0	921	3	0
2	W	947	0	920	13	0
2	X	947	0	921	1	0
All	All	48456	0	48677	118	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 3.

The worst 5 of 118 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)
2:O:113:ILE:CG2	2:O:116:PHE:CD2	2.12	1.31
2:L:76:GLU:O	2:L:80:TYR:CD1	1.86	1.29
2:O:113:ILE:HG22	2:O:116:PHE:CD2	1.68	1.27
2:N:75:GLU:CD	2:N:80:TYR:CD2	2.17	1.18
2:O:113:ILE:HG21	2:O:116:PHE:CD2	1.82	1.15

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	140/142 (99%)	124 (89%)	11 (8%)	5 (4%)	4 38
1	b	140/142 (99%)	120 (86%)	17 (12%)	3 (2%)	9 50
1	c	140/142 (99%)	119 (85%)	15 (11%)	6 (4%)	3 34
1	d	140/142 (99%)	123 (88%)	16 (11%)	1 (1%)	26 71
1	e	140/142 (99%)	117 (84%)	18 (13%)	5 (4%)	4 38
1	f	140/142 (99%)	118 (84%)	18 (13%)	4 (3%)	6 43
1	g	140/142 (99%)	121 (86%)	13 (9%)	6 (4%)	3 34
1	h	140/142 (99%)	116 (83%)	19 (14%)	5 (4%)	4 38
1	i	140/142 (99%)	122 (87%)	12 (9%)	6 (4%)	3 34
1	j	140/142 (99%)	123 (88%)	12 (9%)	5 (4%)	4 38
1	k	140/142 (99%)	119 (85%)	14 (10%)	7 (5%)	3 31
1	l	140/142 (99%)	119 (85%)	13 (9%)	8 (6%)	2 28
1	m	140/142 (99%)	126 (90%)	10 (7%)	4 (3%)	6 43
1	n	140/142 (99%)	127 (91%)	9 (6%)	4 (3%)	6 43
1	o	140/142 (99%)	121 (86%)	16 (11%)	3 (2%)	9 50
1	p	140/142 (99%)	123 (88%)	10 (7%)	7 (5%)	3 31
1	q	140/142 (99%)	121 (86%)	12 (9%)	7 (5%)	3 31
1	r	140/142 (99%)	120 (86%)	15 (11%)	5 (4%)	4 38
1	s	140/142 (99%)	124 (89%)	13 (9%)	3 (2%)	9 50
1	t	140/142 (99%)	118 (84%)	16 (11%)	6 (4%)	3 34
1	u	140/142 (99%)	127 (91%)	10 (7%)	3 (2%)	9 50
1	v	140/142 (99%)	127 (91%)	8 (6%)	5 (4%)	4 38
1	w	140/142 (99%)	127 (91%)	11 (8%)	2 (1%)	14 58
1	x	140/142 (99%)	116 (83%)	19 (14%)	5 (4%)	4 38
2	A	119/121 (98%)	89 (75%)	15 (13%)	15 (13%)	0 8

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	B	119/121 (98%)	89 (75%)	21 (18%)	9 (8%)	1 20
2	C	119/121 (98%)	96 (81%)	12 (10%)	11 (9%)	1 17
2	D	119/121 (98%)	83 (70%)	21 (18%)	15 (13%)	0 8
2	E	119/121 (98%)	95 (80%)	14 (12%)	10 (8%)	1 18
2	F	119/121 (98%)	88 (74%)	23 (19%)	8 (7%)	1 24
2	G	119/121 (98%)	87 (73%)	15 (13%)	17 (14%)	0 6
2	H	119/121 (98%)	94 (79%)	13 (11%)	12 (10%)	1 14
2	I	119/121 (98%)	87 (73%)	24 (20%)	8 (7%)	1 24
2	J	119/121 (98%)	88 (74%)	18 (15%)	13 (11%)	0 11
2	K	119/121 (98%)	93 (78%)	13 (11%)	13 (11%)	0 11
2	L	119/121 (98%)	94 (79%)	16 (13%)	9 (8%)	1 20
2	M	119/121 (98%)	99 (83%)	14 (12%)	6 (5%)	3 31
2	N	119/121 (98%)	98 (82%)	14 (12%)	7 (6%)	2 27
2	O	119/121 (98%)	97 (82%)	17 (14%)	5 (4%)	3 34
2	P	119/121 (98%)	90 (76%)	20 (17%)	9 (8%)	1 20
2	Q	119/121 (98%)	98 (82%)	15 (13%)	6 (5%)	3 31
2	R	119/121 (98%)	93 (78%)	21 (18%)	5 (4%)	3 34
2	S	119/121 (98%)	98 (82%)	9 (8%)	12 (10%)	1 14
2	T	119/121 (98%)	98 (82%)	16 (13%)	5 (4%)	3 34
2	U	119/121 (98%)	85 (71%)	18 (15%)	16 (13%)	0 7
2	V	119/121 (98%)	90 (76%)	19 (16%)	10 (8%)	1 18
2	W	119/121 (98%)	89 (75%)	18 (15%)	12 (10%)	1 14
2	X	119/121 (98%)	95 (80%)	16 (13%)	8 (7%)	1 24
All	All	6216/6312 (98%)	5131 (82%)	729 (12%)	356 (6%)	4 28

5 of 356 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	a	47	ARG
1	a	66	ALA
2	A	59	GLN
2	A	93	GLU
2	A	94	ALA

5.3.2 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 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	121/121 (100%)	118 (98%)	3 (2%)	55 81
1	b	121/121 (100%)	117 (97%)	4 (3%)	45 76
1	c	121/121 (100%)	114 (94%)	7 (6%)	25 61
1	d	121/121 (100%)	118 (98%)	3 (2%)	55 81
1	e	121/121 (100%)	117 (97%)	4 (3%)	45 76
1	f	121/121 (100%)	115 (95%)	6 (5%)	30 66
1	g	121/121 (100%)	119 (98%)	2 (2%)	68 87
1	h	121/121 (100%)	117 (97%)	4 (3%)	45 76
1	i	121/121 (100%)	115 (95%)	6 (5%)	30 66
1	j	121/121 (100%)	115 (95%)	6 (5%)	30 66
1	k	121/121 (100%)	118 (98%)	3 (2%)	55 81
1	l	121/121 (100%)	117 (97%)	4 (3%)	45 76
1	m	121/121 (100%)	117 (97%)	4 (3%)	45 76
1	n	121/121 (100%)	118 (98%)	3 (2%)	55 81
1	o	121/121 (100%)	117 (97%)	4 (3%)	45 76
1	p	121/121 (100%)	115 (95%)	6 (5%)	30 66
1	q	121/121 (100%)	117 (97%)	4 (3%)	45 76
1	r	121/121 (100%)	115 (95%)	6 (5%)	30 66
1	s	121/121 (100%)	117 (97%)	4 (3%)	45 76
1	t	121/121 (100%)	113 (93%)	8 (7%)	21 57
1	u	121/121 (100%)	116 (96%)	5 (4%)	37 71
1	v	121/121 (100%)	120 (99%)	1 (1%)	86 94
1	w	121/121 (100%)	114 (94%)	7 (6%)	25 61
1	x	121/121 (100%)	115 (95%)	6 (5%)	30 66
2	A	109/109 (100%)	101 (93%)	8 (7%)	17 54
2	B	109/109 (100%)	104 (95%)	5 (5%)	33 68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	C	109/109 (100%)	104 (95%)	5 (5%)	33 68
2	D	109/109 (100%)	103 (94%)	6 (6%)	27 63
2	E	109/109 (100%)	104 (95%)	5 (5%)	33 68
2	F	109/109 (100%)	102 (94%)	7 (6%)	22 58
2	G	109/109 (100%)	103 (94%)	6 (6%)	27 63
2	H	109/109 (100%)	104 (95%)	5 (5%)	33 68
2	I	109/109 (100%)	106 (97%)	3 (3%)	51 78
2	J	109/109 (100%)	102 (94%)	7 (6%)	22 58
2	K	109/109 (100%)	100 (92%)	9 (8%)	14 49
2	L	109/109 (100%)	106 (97%)	3 (3%)	51 78
2	M	109/109 (100%)	105 (96%)	4 (4%)	41 73
2	N	109/109 (100%)	98 (90%)	11 (10%)	9 38
2	O	109/109 (100%)	104 (95%)	5 (5%)	33 68
2	P	109/109 (100%)	104 (95%)	5 (5%)	33 68
2	Q	109/109 (100%)	99 (91%)	10 (9%)	11 43
2	R	109/109 (100%)	106 (97%)	3 (3%)	51 78
2	S	109/109 (100%)	106 (97%)	3 (3%)	51 78
2	T	109/109 (100%)	102 (94%)	7 (6%)	22 58
2	U	109/109 (100%)	101 (93%)	8 (7%)	17 54
2	V	109/109 (100%)	104 (95%)	5 (5%)	33 68
2	W	109/109 (100%)	104 (95%)	5 (5%)	33 68
2	X	109/109 (100%)	106 (97%)	3 (3%)	51 78
All	All	5520/5520 (100%)	5272 (96%)	248 (4%)	38 69

5 of 248 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	K	156	THR
2	N	93	GLU
1	w	52	LEU
1	l	58	ASN
2	M	64	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 60 such sidechains are listed below:

Mol	Chain	Res	Type
1	1	45	HIS
2	M	59	GLN
2	V	95	HIS
2	L	74	HIS
2	M	95	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 carbohydrates 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.