



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 10:15 PM GMT

PDB ID : 1SQP
Title : Crystal Structure Analysis of Bovine Bcl with Myxothiazol
Authors : Esser, L.; Quinn, B.; Li, Y.F.; Zhang, M.; Elberry, M.; Yu, L.; Yu, C.A.; Xia, D.
Deposited on : 2004-03-19
Resolution : 2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
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) : trunk26865

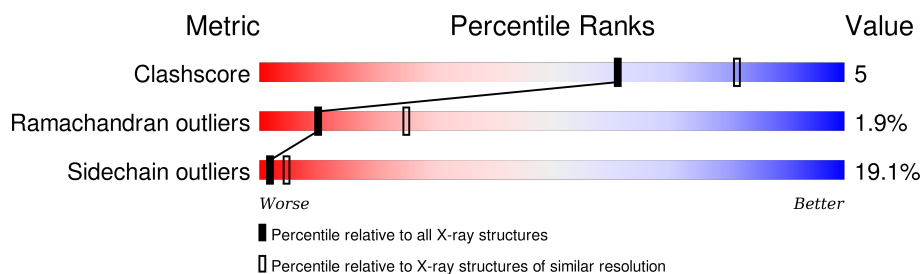
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	480	
2	B	453	
3	C	379	
4	D	241	
5	E	196	
6	F	110	
7	G	81	

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Mol	Chain	Length	Quality of chain
8	H	78	
9	I	78	
10	J	62	
11	K	56	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
13	CDL	A	447	X	-	-	-
13	CDL	D	242	X	-	-	-
13	CDL	G	82	X	-	-	-

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 17274 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 Ubiquinol-cytochrome-c reductase complex core protein I, mitochondrial precursor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	446	Total	C	N	O	S	0	0	0
			3458	2161	609	668	20			

- Molecule 2 is a protein called Ubiquinol-cytochrome-c reductase complex core protein 2, mitochondrial precursor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	423	Total	C	N	O	S	0	0	0
			3172	1993	562	610	7			

- Molecule 3 is a protein called Cytochrome b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	378	Total	C	N	O	S	0	0	0
			3003	2013	471	501	18			

- Molecule 4 is a protein called Cytochrome c1, heme protein, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	241	Total	C	N	O	S	0	0	0
			1919	1225	330	349	15			

- Molecule 5 is a protein called Ubiquinol-cytochrome c reductase iron-sulfur subunit, mitochondrial precursor (EC 1.10.2.2) (Rieske iron-sulfur protein) (RISP) [Contains: Ubiquinol-cytochrome c reductase 8 kDa protein (Complex III subunit IX)].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	196	Total	C	N	O	S	0	0	0
			1519	957	263	291	8			

- Molecule 6 is a protein called sub6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	105	Total	C	N	O	S	0	0	0
			911	576	165	168	2			

- Molecule 7 is a protein called Ubiquinol-cytochrome c reductase complex ubiquinone-binding protein QP-C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	75	Total	C	N	O	S	0	0	0
			628	410	118	99	1			

- Molecule 8 is a protein called Ubiquinol-cytochrome c reductase complex 11 kDa protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	67	Total	C	N	O	S	0	0	0
			548	332	99	112	5			

- Molecule 9 is a protein called Ubiquinol-cytochrome c reductase iron-sulfur subunit, mitochondrial precursor (EC 1.10.2.2) (Rieske iron-sulfur protein) (RISP) [Contains: Ubiquinol-cytochrome c reductase 8 kDa protein (Complex III subunit IX)].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	57	Total	C	N	O	S	0	0	0
			406	253	77	74	2			

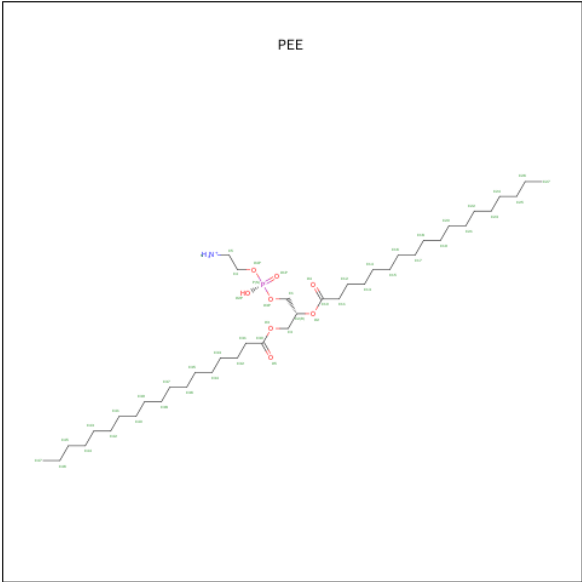
- Molecule 10 is a protein called Ubiquinol-cytochrome c reductase complex 7.2 kDa protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
10	J	61	Total	C	N	O	0	0	0
			502	329	87	86			

- Molecule 11 is a protein called Ubiquinol-cytochrome c reductase complex 6.4 kDa protein.

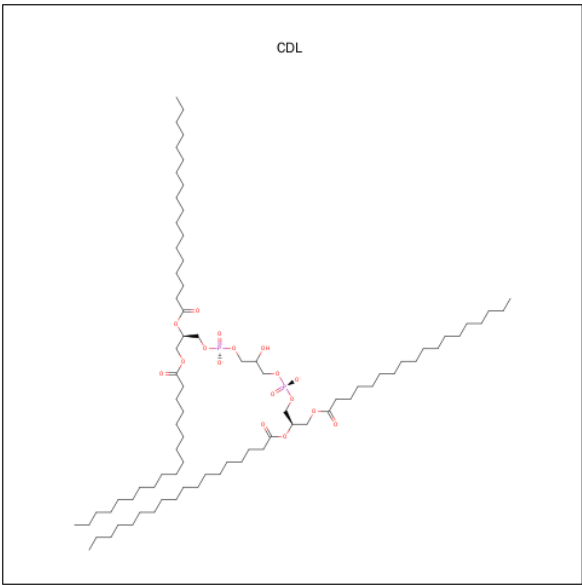
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	53	Total	C	N	O	S	0	0	0
			436	292	78	65	1			

- Molecule 12 is 1,2-DIOLEOYL-SN-GLYCERO-3-PHOSPHOETHANOLAMINE (three-letter code: PEE) (formula: C₄₁H₈₃NO₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
12	E	1	Total	C	N	O	P	0	0
			49	39	1	8	1		
12	A	1	Total	C	N	O	P	0	0
			49	39	1	8	1		
12	C	1	Total	C	N	O	P	0	0
			49	39	1	8	1		

- Molecule 13 is CARDIOLIPIN (three-letter code: CDL) (formula: C₈₁H₁₅₆O₁₇P₂).



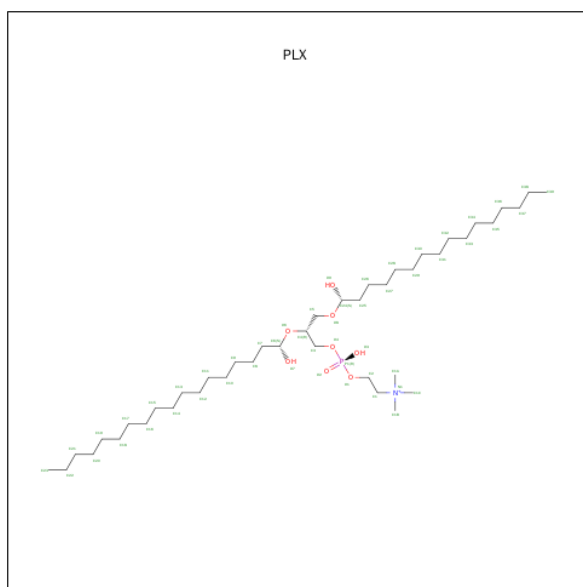
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
13	A	1	Total	C	O	P	0	0
			64	45	17	2		

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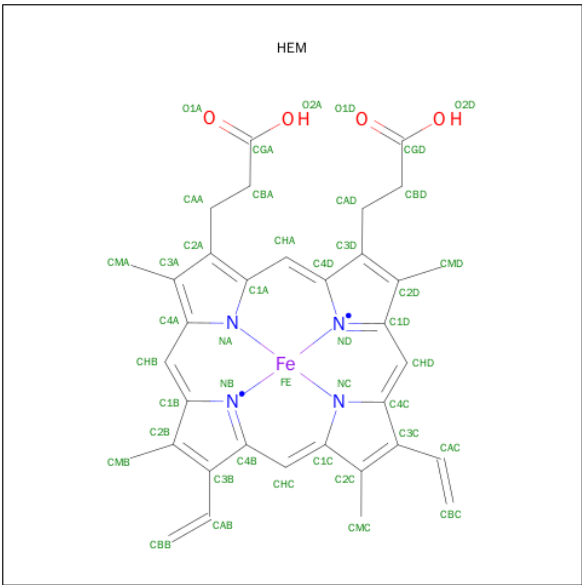
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
13	D	1	Total	C	O	P	0	0
			64	45	17	2		
13	G	1	Total	C	O	P	0	0
			64	45	17	2		

- Molecule 14 is (9R,11S)-9-({[(1S)-1-HYDROXYHEXADECYL]OXY}METHYL)-2,2-DIMETHYL-5,7,10-TRIOXA-2LAMBDA 5 -AZA-6LAMBDA 5 -PHOSPHAOCTACOSANE-6,6,11-TRIOL (three-letter code: PLX) (formula: C₄₂H₈₉NO₈P).



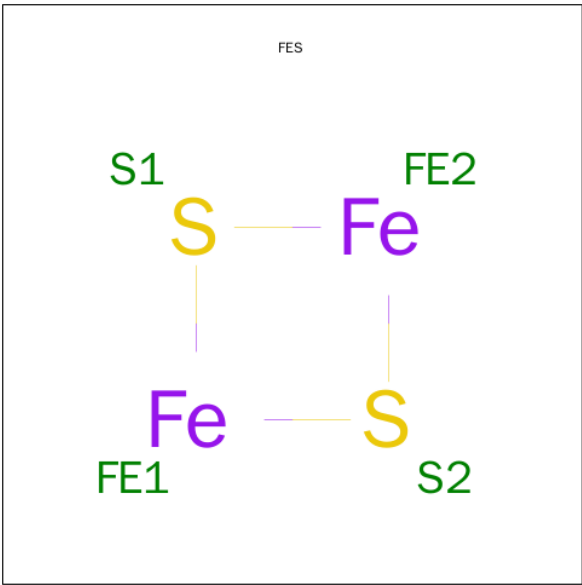
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
14	J	1	Total	C	N	O	P	0	0
			52	42	1	8	1		

- Molecule 15 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C₃₄H₃₂FeN₄O₄).



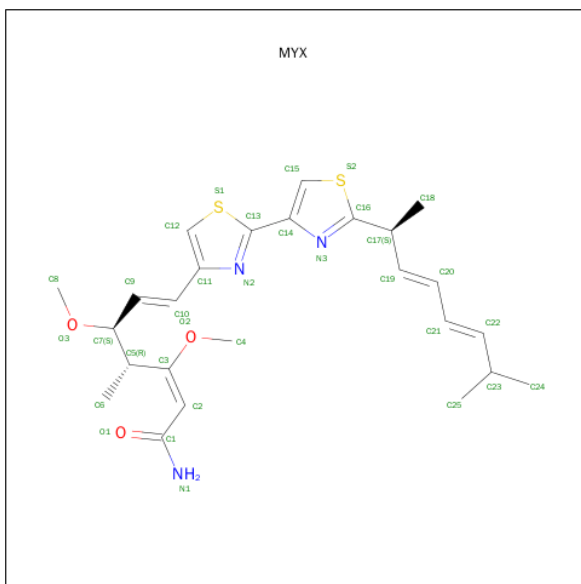
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
15	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
15	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
15	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 16 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe₂S₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
16	E	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 17 is (2Z,6E)-7-{2'-[(2E,4E)-1,6-DIMETHYLHEPTA-2,4-DIENYL]-2,4'-BI-1,3-THIAZOL-4-YL}-3,5-DIMETHOXY-4-METHYLHEPTA-2,6-DIENAMIDE (three-letter code: MYX) (formula: C₂₅H₃₃N₃O₃S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
17	C	1	Total	C	N	O	S	0	0
			33	25	3	3	2		

- Molecule 18 is water.

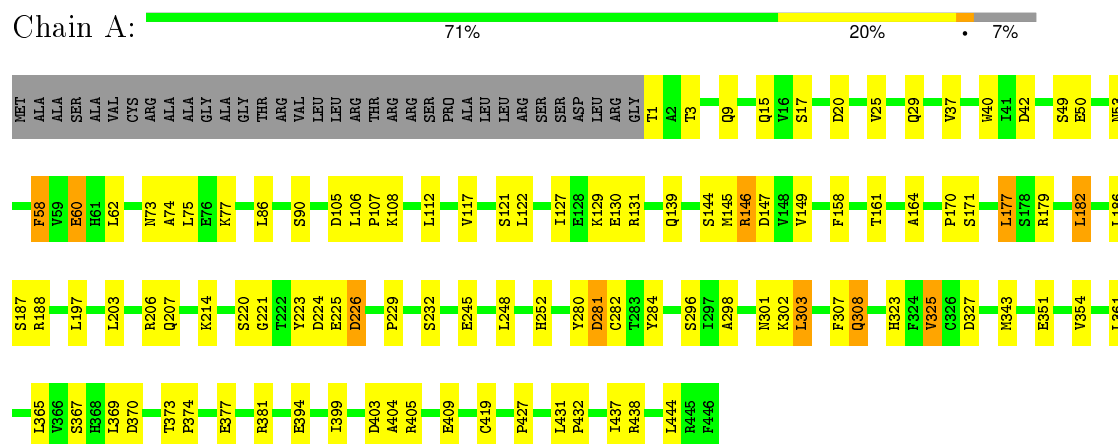
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	A	51	Total	O	0	0
			51	51		
18	B	84	Total	O	0	0
			84	84		
18	C	36	Total	O	0	0
			36	36		
18	D	11	Total	O	0	0
			11	11		
18	E	2	Total	O	0	0
			2	2		
18	F	15	Total	O	0	0
			15	15		
18	G	11	Total	O	0	0
			11	11		
18	I	2	Total	O	0	0
			2	2		
18	K	3	Total	O	0	0
			3	3		

3 Residue-property plots

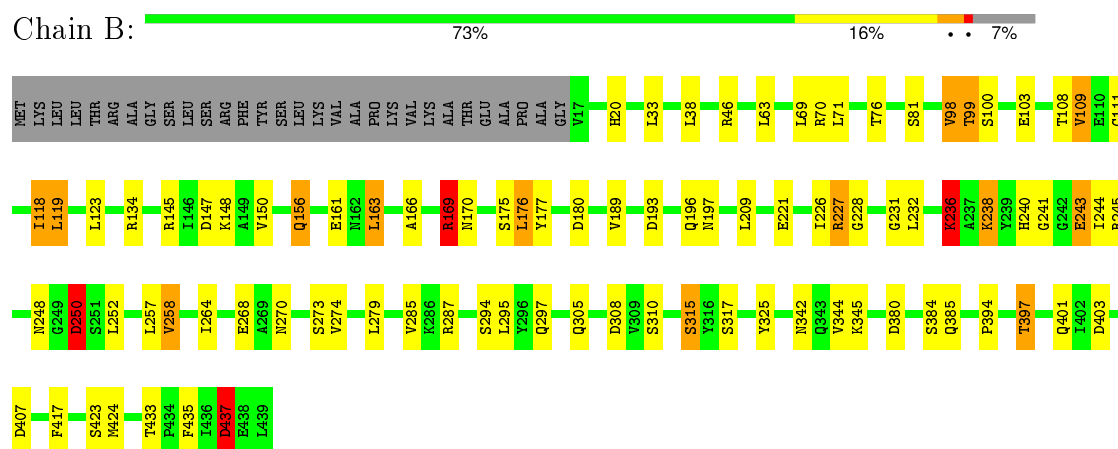
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

Note EDS was not executed.

- Molecule 1: Ubiquinol-cytochrome-c reductase complex core protein I, mitochondrial precursor



- Molecule 2: Ubiquinol-cytochrome-c reductase complex core protein 2, mitochondrial precursor

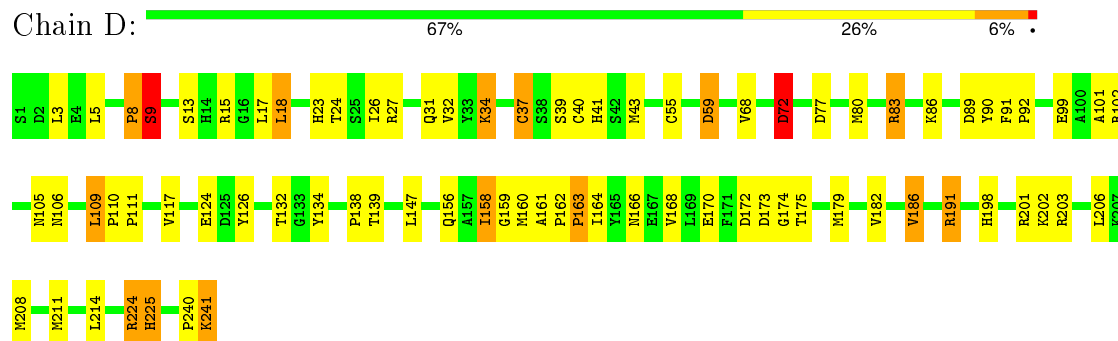


- Molecule 3: Cytochrome b

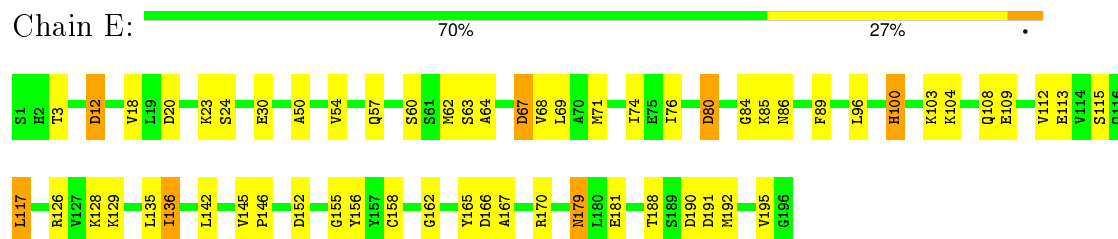




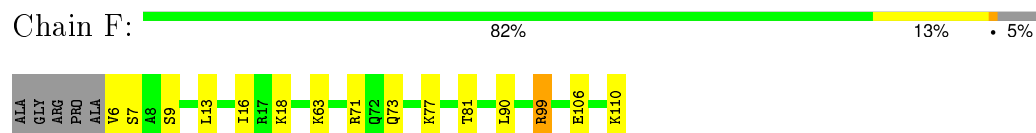
- Molecule 4: Cytochrome c1, heme protein, mitochondrial



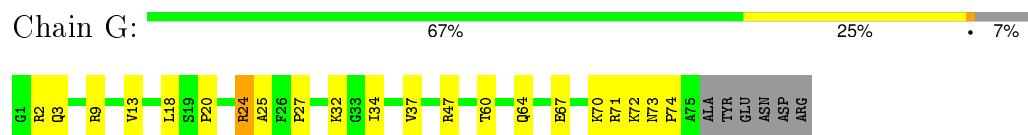
- Molecule 5: Ubiquinol-cytochrome c reductase iron-sulfur subunit, mitochondrial precursor (EC 1.10.2.2) (Rieske iron-sulfur protein) (RISP) [Contains: Ubiquinol-cytochrome c reductase 8 kDa protein (Complex III subunit IX)]



- Molecule 6: sub6



- Molecule 7: Ubiquinol-cytochrome c reductase complex ubiquinone-binding protein QP-C

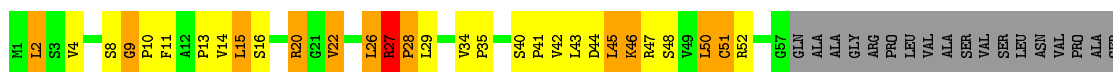
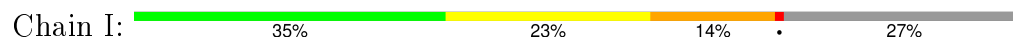


- Molecule 8: Ubiquinol-cytochrome c reductase complex 11 kDa protein





- Molecule 9: Ubiquinol-cytochrome c reductase iron-sulfur subunit, mitochondrial precursor (EC 1.10.2.2) (Rieske iron-sulfur protein) (RISP) [Contains: Ubiquinol-cytochrome c reductase 8 kDa protein (Complex III subunit IX)]



- Molecule 10: Ubiquinol-cytochrome c reductase complex 7.2 kDa protein



- Molecule 11: Ubiquinol-cytochrome c reductase complex 6.4 kDa protein



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	153.70Å 153.70Å 596.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.70	Depositor
% Data completeness (in resolution range)	94.1 (40.00-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.263 , 0.314	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	17274	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MYX, CDL, PLX, FES, HEM, PEE

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	0.90	0/3531	0.84	5/4792 (0.1%)
2	B	0.99	0/3232	0.88	8/4386 (0.2%)
3	C	1.01	0/3100	0.82	3/4242 (0.1%)
4	D	0.97	0/1978	0.86	5/2684 (0.2%)
5	E	1.01	0/1553	0.82	6/2100 (0.3%)
6	F	0.99	0/930	0.88	0/1246
7	G	1.15	0/649	0.78	0/878
8	H	0.90	0/553	0.88	0/741
9	I	1.31	0/411	1.15	2/558 (0.4%)
10	J	1.06	0/515	0.85	0/696
11	K	1.13	0/452	0.82	0/618
All	All	0.99	0/16904	0.86	29/22941 (0.1%)

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
2	B	0	1
6	F	0	1
All	All	0	3

There are no bond length outliers.

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	172	ASP	CB-CG-OD2	6.79	124.41	118.30
1	A	327	ASP	CB-CG-OD2	6.56	124.20	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	308	ASP	CB-CG-OD2	6.40	124.06	118.30
2	B	147	ASP	CB-CG-OD2	6.28	123.95	118.30
3	C	320	LEU	CA-CB-CG	5.78	128.59	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	303	LEU	Peptide
2	B	169	ARG	Mainchain
6	F	99	ARG	Mainchain

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	3458	0	3356	32	0
2	B	3172	0	3152	44	0
3	C	3003	0	3065	33	0
4	D	1919	0	1869	35	0
5	E	1519	0	1504	12	0
6	F	911	0	902	3	0
7	G	628	0	636	8	0
8	H	548	0	530	3	0
9	I	406	0	437	27	0
10	J	502	0	505	3	0
11	K	436	0	445	11	0
12	A	49	0	75	1	0
12	C	49	0	75	0	0
12	E	49	0	75	3	0
13	A	64	0	72	1	0
13	D	64	0	72	0	0
13	G	64	0	72	2	0
14	J	52	0	88	4	0
15	C	86	0	60	4	0
15	D	43	0	30	11	0
16	E	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
17	C	33	0	33	1	0
18	A	51	0	0	0	0
18	B	84	0	0	2	0
18	C	36	0	0	0	0
18	D	11	0	0	0	0
18	E	2	0	0	1	0
18	F	15	0	0	0	0
18	G	11	0	0	1	0
18	I	2	0	0	0	0
18	K	3	0	0	0	0
All	All	17274	0	17053	186	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 5.

The worst 5 of 186 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)
4:D:40:CYS:SG	15:D:243:HEM:HAC	1.33	1.68
3:C:11:MET:CE	3:C:11:MET:SD	2.02	1.48
11:K:38:TRP:HE3	11:K:41:ILE:CD1	1.25	1.48
11:K:38:TRP:CE3	11:K:41:ILE:CD1	2.04	1.41
4:D:37:CYS:SG	15:D:243:HEM:HAB	1.62	1.35

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 X-ray entries followed by that with respect to entries of similar resolution.

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	444/480 (92%)	418 (94%)	20 (4%)	6 (1%)	14 35
2	B	421/453 (93%)	390 (93%)	27 (6%)	4 (1%)	19 45

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	376/379 (99%)	353 (94%)	19 (5%)	4 (1%)	17	42
4	D	239/241 (99%)	211 (88%)	21 (9%)	7 (3%)	6	14
5	E	194/196 (99%)	162 (84%)	29 (15%)	3 (2%)	13	32
6	F	103/110 (94%)	98 (95%)	5 (5%)	0	100	100
7	G	73/81 (90%)	66 (90%)	5 (7%)	2 (3%)	6	16
8	H	65/78 (83%)	58 (89%)	5 (8%)	2 (3%)	5	12
9	I	55/78 (70%)	31 (56%)	16 (29%)	8 (14%)	0	0
10	J	59/62 (95%)	50 (85%)	7 (12%)	2 (3%)	5	10
11	K	51/56 (91%)	43 (84%)	7 (14%)	1 (2%)	9	24
All	All	2080/2214 (94%)	1880 (90%)	161 (8%)	39 (2%)	10	25

5 of 39 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	250	ASP
3	C	154	PRO
4	D	8	PRO
7	G	73	ASN
9	I	29	LEU

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 X-ray entries followed by that with respect to entries of similar resolution.

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	370/394 (94%)	315 (85%)	55 (15%)	4	9
2	B	332/355 (94%)	288 (87%)	44 (13%)	5	11
3	C	326/327 (100%)	262 (80%)	64 (20%)	1	4
4	D	206/206 (100%)	158 (77%)	48 (23%)	1	2
5	E	168/168 (100%)	130 (77%)	38 (23%)	1	3
6	F	96/99 (97%)	85 (88%)	11 (12%)	7	16
7	G	66/71 (93%)	55 (83%)	11 (17%)	3	7

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
8	H	64/74 (86%)	42 (66%)	22 (34%)	0	0
9	I	44/60 (73%)	32 (73%)	12 (27%)	0	1
10	J	51/52 (98%)	32 (63%)	19 (37%)	0	0
11	K	42/45 (93%)	29 (69%)	13 (31%)	0	0
All	All	1765/1851 (95%)	1428 (81%)	337 (19%)	2	4

5 of 337 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	296	PHE
4	D	106	ASN
10	J	25	VAL
3	C	334	THR
4	D	27	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	342	ASN
3	C	114	ASN
6	F	73	GLN
2	B	277	HIS
7	G	3	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

12 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
13	CDL	A	447	-	63,63,99	1.66	8 (12%)	65,75,111	1.59	8 (12%)
12	PEE	A	448	-	48,48,50	1.37	4 (8%)	49,53,55	1.41	6 (12%)
12	PEE	C	380	-	48,48,50	1.30	4 (8%)	49,53,55	1.35	7 (14%)
15	HEM	C	381	3	30,50,50	2.56	10 (33%)	24,82,82	3.27	10 (41%)
15	HEM	C	382	3	30,50,50	2.65	11 (36%)	24,82,82	3.25	11 (45%)
17	MYX	C	383	-	30,34,34	2.57	6 (20%)	20,45,45	1.52	3 (15%)
13	CDL	D	242	-	63,63,99	1.68	9 (14%)	65,75,111	1.57	7 (10%)
15	HEM	D	243	4	30,50,50	2.75	11 (36%)	24,82,82	2.94	12 (50%)
12	PEE	E	197	-	48,48,50	1.32	4 (8%)	49,53,55	1.28	4 (8%)
16	FES	E	198	5	0,4,4	0.00	-	0,4,4	0.00	-
13	CDL	G	82	-	63,63,99	1.73	8 (12%)	65,75,111	1.69	10 (15%)
14	PLX	J	63	-	51,51,51	0.80	3 (5%)	49,59,59	0.65	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
13	CDL	A	447	-	1/1/9/9	0/74/74/110	0/0/0/0
12	PEE	A	448	-	-	0/52/52/54	0/0/0/0
12	PEE	C	380	-	-	0/52/52/54	0/0/0/0
15	HEM	C	381	3	-	0/10/54/54	0/0/8/8
15	HEM	C	382	3	-	0/10/54/54	0/0/8/8
17	MYX	C	383	-	-	0/24/36/36	0/2/2/2
13	CDL	D	242	-	1/1/9/9	0/74/74/110	0/0/0/0
15	HEM	D	243	4	-	0/10/54/54	0/0/8/8
12	PEE	E	197	-	-	0/52/52/54	0/0/0/0
16	FES	E	198	5	-	0/0/4/4	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
13	CDL	G	82	-	1/1/9/9	0/74/74/110	0/0/0/0
14	PLX	J	63	-	-	0/53/55/55	0/0/0/0

The worst 5 of 78 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	D	243	HEM	C3D-C4D	-7.51	1.41	1.51
15	C	381	HEM	C3D-C4D	-6.77	1.42	1.51
15	C	382	HEM	C3D-C4D	-6.70	1.43	1.51
15	D	243	HEM	C2D-C3D	-6.49	1.35	1.54
15	C	382	HEM	C2D-C3D	-6.02	1.36	1.54

The worst 5 of 78 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	C	381	HEM	C3B-CAB-CBB	-8.24	111.82	124.46
15	C	382	HEM	C3C-CAC-CBC	-7.40	113.10	124.46
15	D	243	HEM	C3C-CAC-CBC	-7.05	113.64	124.46
15	C	382	HEM	C3B-CAB-CBB	-6.95	113.79	124.46
13	G	82	CDL	OB8-CB7-OB9	-6.29	107.25	123.49

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
13	D	242	CDL	CA4
13	G	82	CDL	CA4
13	A	447	CDL	CA4

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	A	447	CDL	1	0
12	A	448	PEE	1	0
15	C	381	HEM	3	0
15	C	382	HEM	1	0
17	C	383	MYX	1	0
15	D	243	HEM	11	0
12	E	197	PEE	3	0
13	G	82	CDL	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
14	J	63	PLX	4	0

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.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.