



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 02:29 AM GMT

PDB ID : 2H88
Title : Avian Mitochondrial Respiratory Complex II at 1.8 Angstrom Resolution
Authors : Huang, L.S.; Shen, J.T.; Wang, A.C.; Berry, E.A.
Deposited on : 2006-06-06
Resolution : 1.74 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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

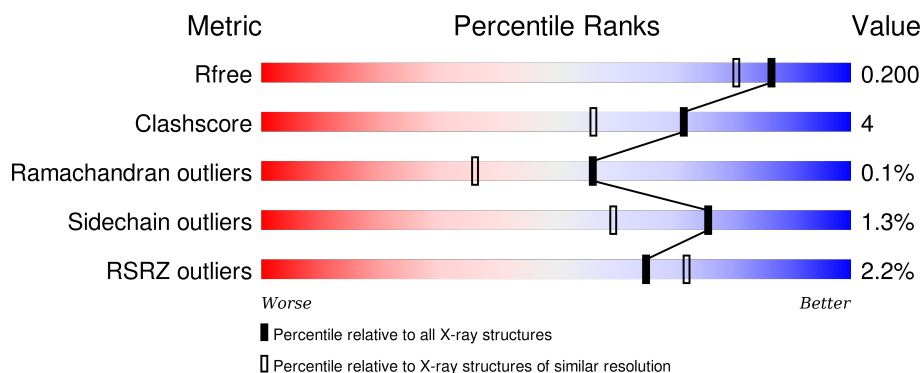
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.74 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))
R_{free}	91344	2417 (1.76-1.72)
Clashscore	102246	2570 (1.76-1.72)
Ramachandran outliers	100387	2544 (1.76-1.72)
Sidechain outliers	100360	2544 (1.76-1.72)
RSRZ outliers	91569	2420 (1.76-1.72)

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.

Mol	Chain	Length	Quality of chain
1	A	621	<div> <div>90%</div> <div>8%</div> <div>.</div> </div>
1	N	621	<div> <div>91%</div> <div>7%</div> <div>.</div> </div>
2	B	252	<div> <div>89%</div> <div>6%</div> <div>5%</div> </div>
2	O	252	<div> <div>90%</div> <div>.</div> <div>5%</div> </div>
3	C	140	<div> <div>81%</div> <div>18%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
3	P	140	
4	D	103	
4	Q	103	

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	GOL	B	1009	-	-	-	X
13	GOL	O	1009	-	-	-	X
14	BHG	C	141	X	-	-	-
14	BHG	P	205	X	-	-	-
8	TEO	A	1002	X	-	-	-
8	TEO	N	1002	X	-	-	-
9	UNL	B	1007	-	-	-	X
9	UNL	C	145	-	-	-	X
9	UNL	D	107	-	-	-	X
9	UNL	D	116	-	-	-	X
9	UNL	N	1007	-	-	X	-
9	UNL	O	1006	-	-	X	-
9	UNL	P	208	-	-	X	-
9	UNL	P	209	-	-	X	-
9	UNL	P	220	-	-	-	X
9	UNL	Q	212	-	-	-	X

2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 19414 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 SUCCINATE DEHYDROGENASE FLAVOPROTEIN SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	612	Total	C	N	O	S	0	0	0
			4726	2956	843	898	29			
1	N	612	Total	C	N	O	S	0	0	0
			4726	2956	843	898	29			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	501	ARG	CYS	CONFLICT	UNP Q9YHT1
A	556	LEU	PHE	CONFLICT	UNP Q9YHT1
A	560	GLU	ASP	CONFLICT	UNP Q9YHT1
N	501	ARG	CYS	CONFLICT	UNP Q9YHT1
N	556	LEU	PHE	CONFLICT	UNP Q9YHT1
N	560	GLU	ASP	CONFLICT	UNP Q9YHT1

- Molecule 2 is a protein called Succinate dehydrogenase Ip subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	239	Total	C	N	O	S	0	0	0
			1916	1213	325	356	22			
2	O	239	Total	C	N	O	S	0	0	0
			1916	1213	325	356	22			

- Molecule 3 is a protein called SUCCINATE DEHYDROGENASE CYTOCHROME B, LARGE SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	139	Total	C	N	O	S	0	0	0
			1077	706	178	189	4			
3	P	139	Total	C	N	O	S	0	0	0
			1077	706	178	189	4			

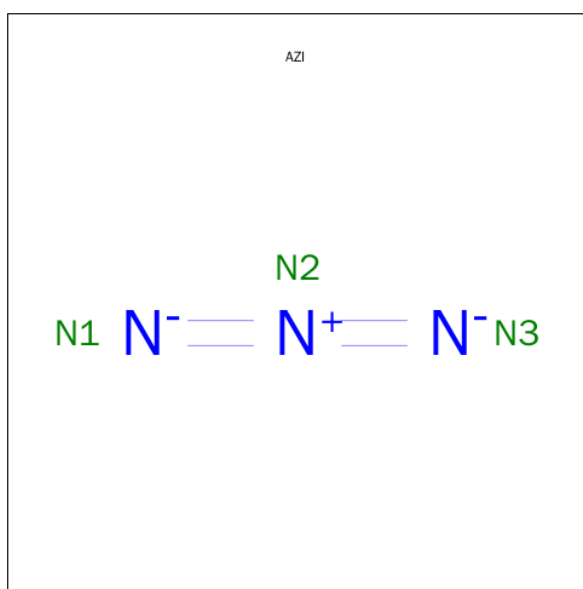
- Molecule 4 is a protein called Succinate dehydrogenase cytochrome B, small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	101	Total	C	N	O	S	0	0	0
			766	505	121	137	3			
4	Q	101	Total	C	N	O	S	0	0	0
			766	505	121	137	3			

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	O	1	Total K 1 1	0	0
5	B	1	Total K 1 1	0	0
5	A	1	Total K 1 1	0	0
5	N	1	Total K 1 1	0	0

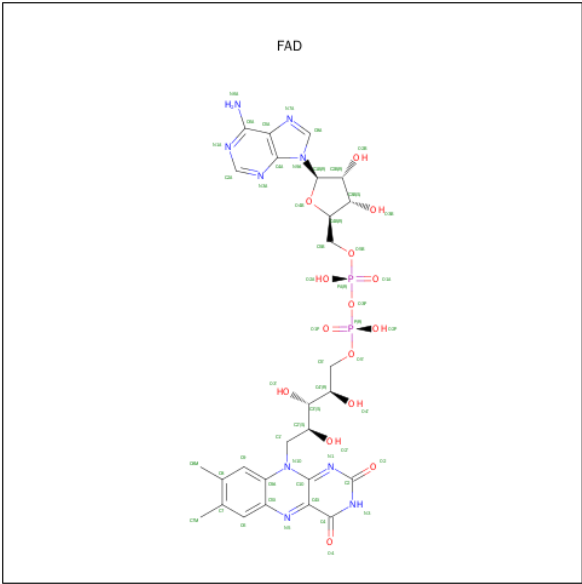
- Molecule 6 is AZIDE ION (three-letter code: AZI) (formula: N₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total N 3 3	0	0
6	N	1	Total N 3 3	0	0

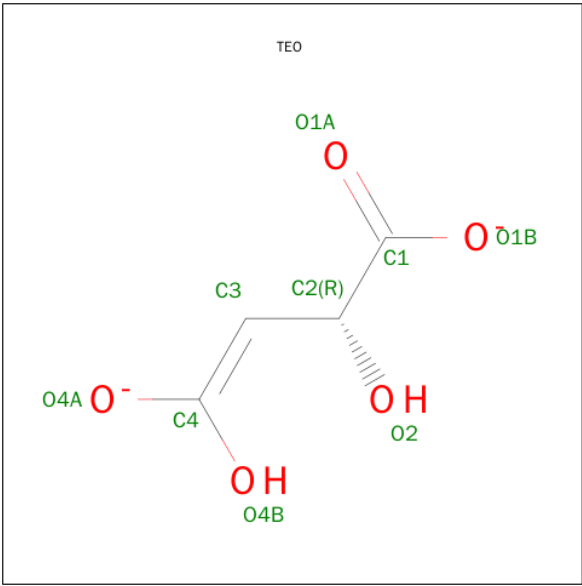
- Molecule 7 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:

C₂₇H₃₃N₉O₁₅P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
7	N	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 8 is MALATE LIKE INTERMEDIATE (three-letter code: TEO) (formula: C₄H₄O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			9	4	5		

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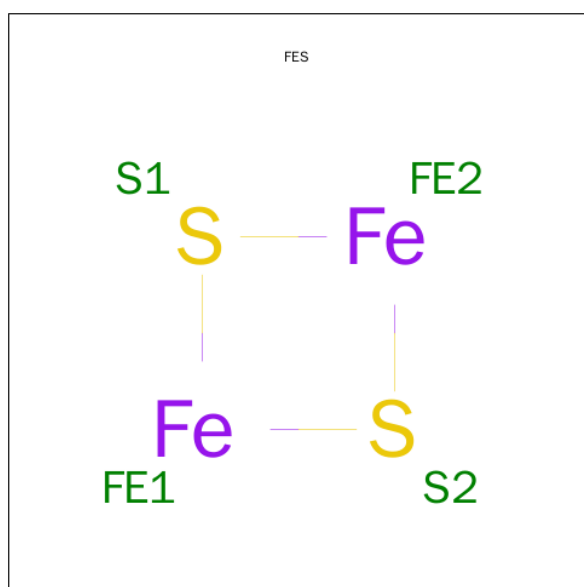
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	N	1	Total	C	O	0	0
			9	4	5		

- Molecule 9 is UNKNOWN LIGAND (three-letter code: UNL) (formula:).

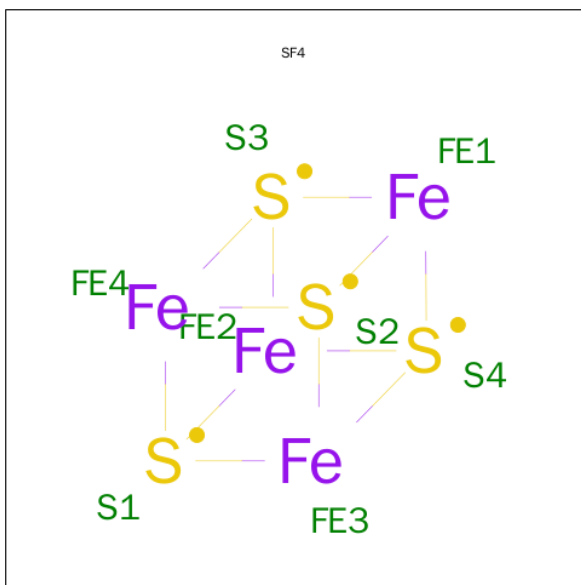
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	P	14	Total	C	O	0	0
			22	6	16		
9	Q	6	Total	C	O	0	0
			16	12	4		
9	D	11	Total	C	O	0	0
			32	21	11		
9	B	4	Total	O		0	0
			4	4			
9	C	9	Total	C	O	0	0
			16	6	10		
9	A	6	Total	O		0	0
			7	7			
9	N	10	Total	O		0	0
			11	11			
9	O	4	Total	O		0	0
			4	4			

- Molecule 10 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe₂S₂).



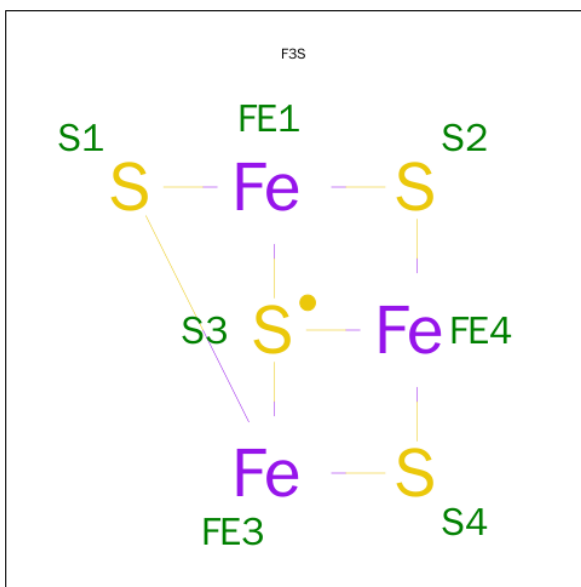
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	Fe	S	0	0
			4	2	2		
10	O	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 11 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



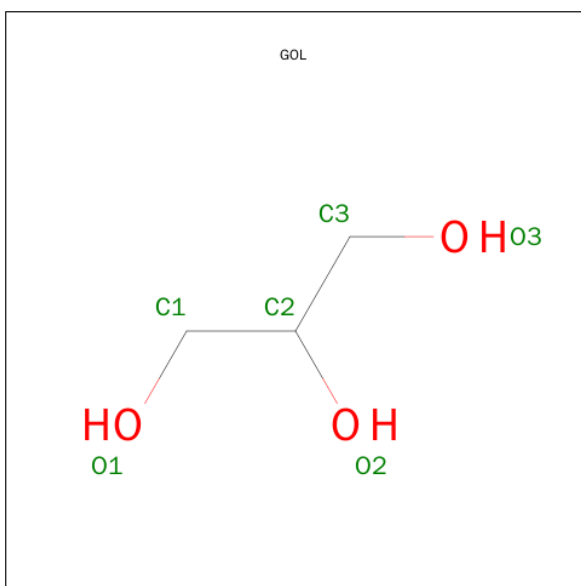
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	B	1	Total	Fe	S	0	0
			8	4	4		
11	O	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 12 is FE3-S4 CLUSTER (three-letter code: F3S) (formula: Fe_3S_4).



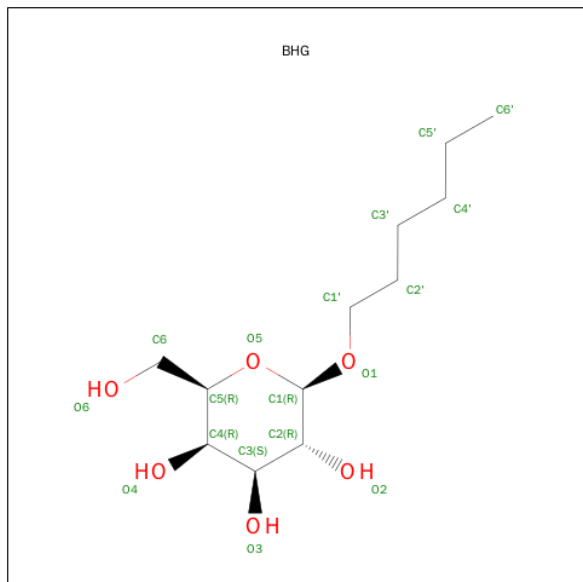
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	B	1	Total	Fe	S	0	0
			7	3	4		
12	O	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 13 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



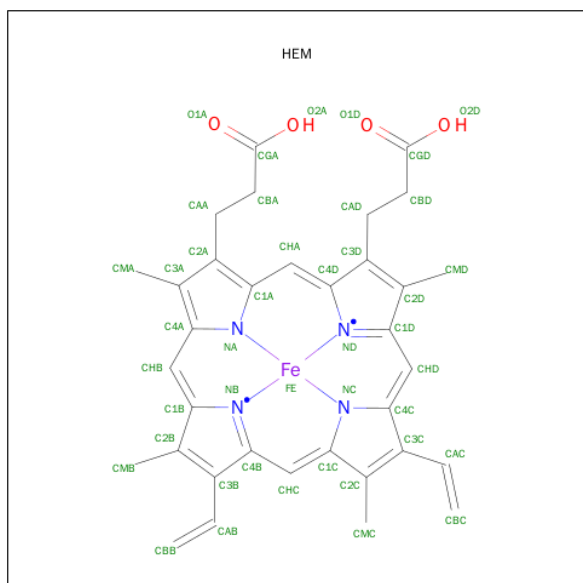
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
13	B	1	Total	C	O	0	0
			6	3	3		
13	O	1	Total	C	O	0	0
			6	3	3		

- Molecule 14 is SUGAR (2-HEXYLOXY-6-HYDROXYMETHYL-TETRAHYDRO-PYRAN-3,4,5-TRIOL) (three-letter code: BHG) (formula: $C_{12}H_{24}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
14	C	1	Total	C	O	0	0
			18	12	6		
14	P	1	Total	C	O	0	0
			18	12	6		

- Molecule 15 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
15	C	1	Total 41	C 32	Fe 1	N 4	O 4	0	0
15	P	1	Total 41	C 32	Fe 1	N 4	O 4	0	0

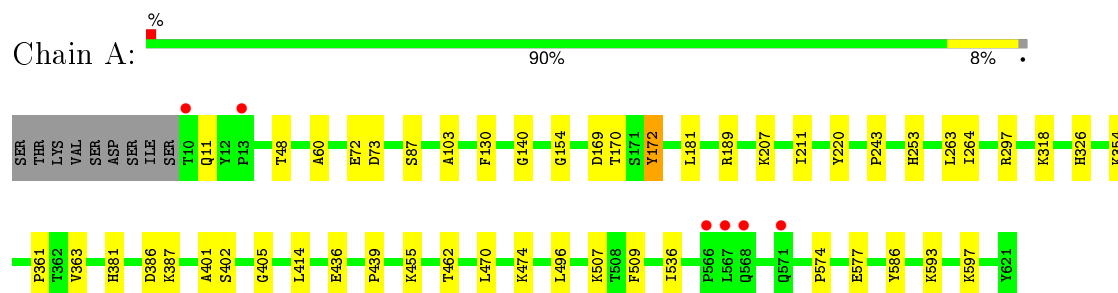
- Molecule 16 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	A	579	Total 579	O 579	0	0
16	B	296	Total 296	O 296	0	0
16	C	104	Total 104	O 104	0	0
16	D	51	Total 51	O 51	0	0
16	N	569	Total 569	O 569	0	0
16	O	281	Total 281	O 281	0	0
16	P	99	Total 99	O 99	0	0
16	Q	51	Total 51	O 51	0	0

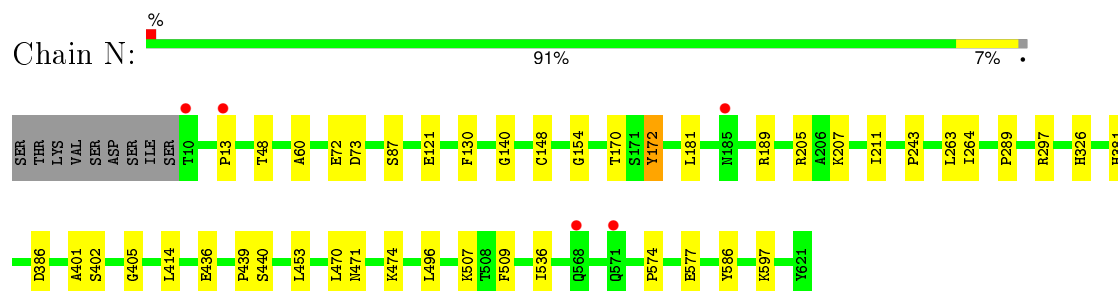
3 Residue-property plots [i](#)

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 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.

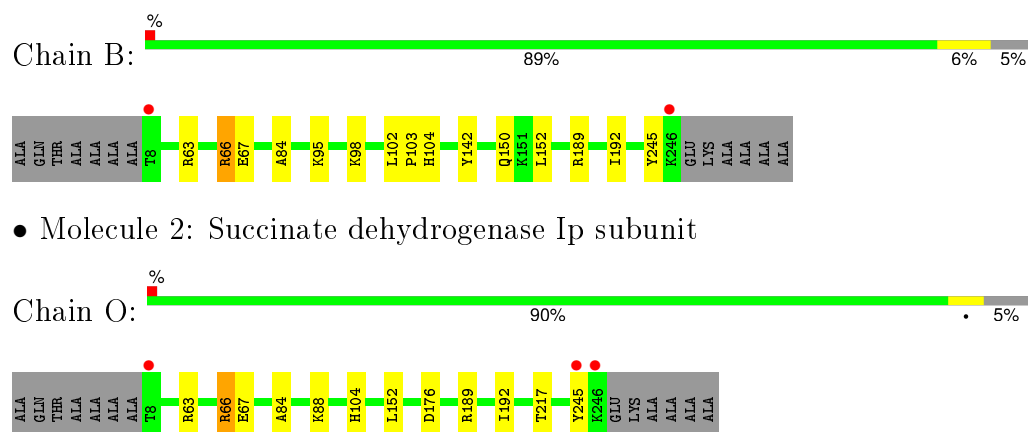
- Molecule 1: SUCCINATE DEHYDROGENASE FLAVOPROTEIN SUBUNIT



- Molecule 1: SUCCINATE DEHYDROGENASE FLAVOPROTEIN SUBUNIT

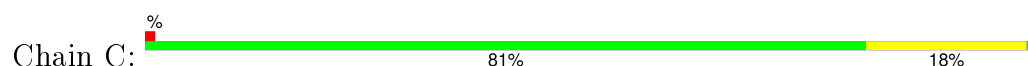


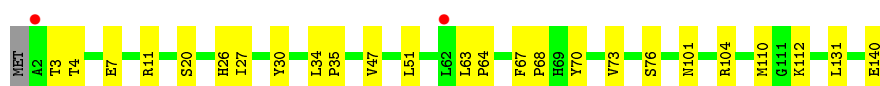
- Molecule 2: Succinate dehydrogenase Ip subunit



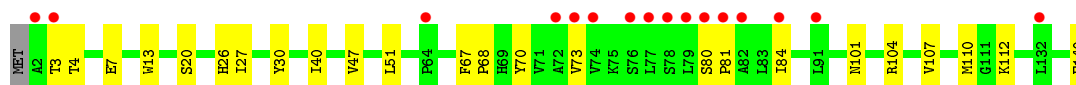
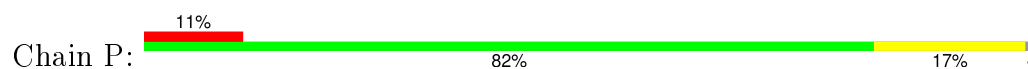
- Molecule 2: Succinate dehydrogenase Ip subunit

- Molecule 3: SUCCINATE DEHYDROGENASE CYTOCHROME B, LARGE SUBUNIT

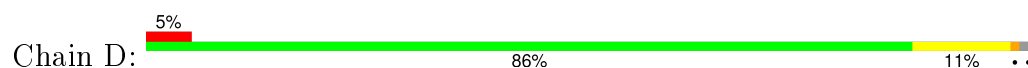




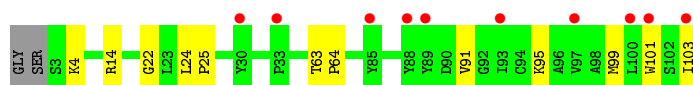
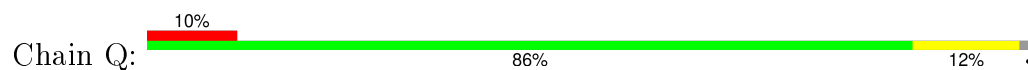
- Molecule 3: SUCCINATE DEHYDROGENASE CYTOCHROME B, LARGE SUBUNIT



- Molecule 4: Succinate dehydrogenase cytochrome B, small subunit



- Molecule 4: Succinate dehydrogenase cytochrome B, small subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	119.97Å 199.39Å 68.06Å 90.00° 90.35° 90.00°	Depositor
Resolution (Å)	45.14 – 1.74 45.14 – 1.70	Depositor EDS
% Data completeness (in resolution range)	89.2 (45.14-1.74) 86.9 (45.14-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 1.70Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.178 , 0.206 0.175 , 0.200	Depositor DCC
R_{free} test set	14646 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.319	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.0	EDS
Estimated twinning fraction	0.227 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 304101 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19414	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.84% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: AZI, GOL, TEO, SF4, BHG, F3S, FES, HEM, UNL, K, FAD

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	0.99	1/4827 (0.0%)	0.86	2/6535 (0.0%)
1	N	1.01	1/4827 (0.0%)	0.87	1/6535 (0.0%)
2	B	0.99	2/1958 (0.1%)	0.83	0/2640
2	O	1.00	1/1958 (0.1%)	0.84	0/2640
3	C	0.73	0/1106	0.68	0/1503
3	P	0.73	0/1106	0.68	0/1503
4	D	0.60	0/789	0.64	0/1082
4	Q	0.54	0/789	0.62	0/1082
All	All	0.94	5/17360 (0.0%)	0.82	3/23520 (0.0%)

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
3	C	0	1
3	P	0	1
All	All	0	4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	O	84	ALA	CA-CB	6.40	1.65	1.52
2	B	142	TYR	CD1-CE1	6.09	1.48	1.39
1	A	103	ALA	CA-CB	5.94	1.65	1.52
2	B	84	ALA	CA-CB	5.23	1.63	1.52
1	N	148	CYS	CB-SG	5.16	1.91	1.82

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	140	GLY	N-CA-C	6.00	128.10	113.10
1	N	140	GLY	N-CA-C	5.91	127.87	113.10
1	A	318	LYS	CD-CE-NZ	5.25	123.78	111.70

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	172	TYR	Sidechain
3	C	30	TYR	Sidechain
1	N	172	TYR	Sidechain
3	P	30	TYR	Sidechain

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	4726	0	4609	33	0
1	N	4726	0	4609	34	0
2	B	1916	0	1913	10	0
2	O	1916	0	1913	8	0
3	C	1077	0	1112	22	0
3	P	1077	0	1112	20	0
4	D	766	0	761	10	0
4	Q	766	0	761	10	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	N	1	0	0	0	0
5	O	1	0	0	0	0
6	A	3	0	0	0	0
6	N	3	0	0	0	0
7	A	53	0	30	3	0
7	N	53	0	30	3	0
8	A	9	0	2	3	0
8	N	9	0	2	1	0
9	A	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	B	4	0	0	1	0
9	C	16	0	0	0	0
9	D	32	0	0	1	0
9	N	11	0	0	3	0
9	O	4	0	0	2	0
9	P	22	0	0	4	0
9	Q	16	0	0	0	0
10	B	4	0	0	0	0
10	O	4	0	0	0	0
11	B	8	0	0	0	0
11	O	8	0	0	0	0
12	B	7	0	0	0	0
12	O	7	0	0	0	0
13	B	6	0	8	1	0
13	O	6	0	8	1	0
14	C	18	0	24	0	0
14	P	18	0	24	0	0
15	C	41	0	24	0	0
15	P	41	0	24	0	0
16	A	579	0	0	8	0
16	B	296	0	0	3	0
16	C	104	0	0	3	0
16	D	51	0	0	1	0
16	N	569	0	0	7	0
16	O	281	0	0	0	0
16	P	99	0	0	0	0
16	Q	51	0	0	0	0
All	All	19414	0	16966	146	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 4.

The worst 5 of 146 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)
9:O:1005:UNL:O1	9:O:1006:UNL:O1	1.69	1.09
1:N:13:PRO:O	9:N:1011:UNL:O1	1.75	1.04
9:B:1005:UNL:O1	9:B:1006:UNL:O1	1.84	0.96
4:Q:91:VAL:HG11	4:Q:99:MET:HE1	1.56	0.85
3:P:101:ASN:HD21	3:P:104:ARG:HH11	1.27	0.82

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	610/621 (98%)	591 (97%)	19 (3%)	0	100	100
1	N	610/621 (98%)	591 (97%)	19 (3%)	0	100	100
2	B	237/252 (94%)	231 (98%)	5 (2%)	1 (0%)	39	20
2	O	237/252 (94%)	232 (98%)	4 (2%)	1 (0%)	39	20
3	C	137/140 (98%)	136 (99%)	1 (1%)	0	100	100
3	P	137/140 (98%)	136 (99%)	1 (1%)	0	100	100
4	D	99/103 (96%)	98 (99%)	1 (1%)	0	100	100
4	Q	99/103 (96%)	98 (99%)	1 (1%)	0	100	100
All	All	2166/2232 (97%)	2113 (98%)	51 (2%)	2 (0%)	56	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	66	ARG
2	O	66	ARG

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 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	497/506 (98%)	493 (99%)	4 (1%)	86	77

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	N	497/506 (98%)	493 (99%)	4 (1%)	86	77
2	B	214/219 (98%)	210 (98%)	4 (2%)	65	42
2	O	214/219 (98%)	209 (98%)	5 (2%)	58	33
3	C	118/119 (99%)	117 (99%)	1 (1%)	86	77
3	P	118/119 (99%)	117 (99%)	1 (1%)	86	77
4	D	78/79 (99%)	76 (97%)	2 (3%)	54	28
4	Q	78/79 (99%)	76 (97%)	2 (3%)	54	28
All	All	1814/1846 (98%)	1791 (99%)	23 (1%)	76	60

5 of 23 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	D	101	TRP
1	N	72	GLU
4	Q	14	ARG
1	N	48	THR
1	N	73	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	101	ASN
1	N	326	HIS
3	P	26	HIS
3	C	66	GLN
2	O	137	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 ⓘ

Of 86 ligands modelled in this entry, 64 are unknown and 4 are monoatomic - leaving 18 for Mogul analysis.

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$
7	FAD	A	1001	1	48,58,58	1.88	16 (33%)	54,89,89	2.90	13 (24%)
8	TEO	A	1002	-	0,8,8	0.00	-	1,10,10	12.37	1 (100%)
6	AZI	A	623	-	0,2,2	0.00	-	0,1,1	0.00	-
10	FES	B	1002	2	0,4,4	0.00	-	0,4,4	0.00	-
11	SF4	B	1003	2	0,12,12	0.00	-	0,24,24	0.00	-
12	F3S	B	1004	2	0,9,9	0.00	-	0,15,15	0.00	-
13	GOL	B	1009	-	5,5,5	1.12	0	5,5,5	0.58	0
14	BHG	C	141	-	18,18,18	1.73	4 (22%)	23,23,23	0.78	1 (4%)
15	HEM	C	142	3,4	29,48,50	2.20	9 (31%)	24,80,82	2.77	10 (41%)
7	FAD	N	1001	1	48,58,58	2.07	15 (31%)	54,89,89	2.96	13 (24%)
8	TEO	N	1002	-	0,8,8	0.00	-	1,10,10	13.17	1 (100%)
6	AZI	N	623	-	0,2,2	0.00	-	0,1,1	0.00	-
10	FES	O	1002	2	0,4,4	0.00	-	0,4,4	0.00	-
11	SF4	O	1003	2	0,12,12	0.00	-	0,24,24	0.00	-
12	F3S	O	1004	2	0,9,9	0.00	-	0,15,15	0.00	-
13	GOL	O	1009	-	5,5,5	1.15	0	5,5,5	0.73	0
15	HEM	P	201	3,4	29,48,50	2.24	9 (31%)	24,80,82	2.76	10 (41%)
14	BHG	P	205	-	18,18,18	1.34	2 (11%)	23,23,23	0.66	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
7	FAD	A	1001	1	-	0/30/50/50	0/6/6/6
8	TEO	A	1002	-	1/1/3/4	0/1/8/8	0/0/0/0
6	AZI	A	623	-	-	0/0/0/0	0/0/0/0
10	FES	B	1002	2	-	0/0/4/4	0/1/1/1
11	SF4	B	1003	2	-	0/0/48/48	0/6/5/5
12	F3S	B	1004	2	-	0/0/24/24	0/0/3/3
13	GOL	B	1009	-	-	0/4/4/4	0/0/0/0
14	BHG	C	141	-	1/1/5/5	0/9/29/29	0/1/1/1
15	HEM	C	142	3,4	-	0/6/50/54	0/0/8/8
7	FAD	N	1001	1	-	0/30/50/50	0/6/6/6
8	TEO	N	1002	-	1/1/3/4	0/1/8/8	0/0/0/0
6	AZI	N	623	-	-	0/0/0/0	0/0/0/0
10	FES	O	1002	2	-	0/0/4/4	0/1/1/1
11	SF4	O	1003	2	-	0/0/48/48	0/6/5/5
12	F3S	O	1004	2	-	0/0/24/24	0/0/3/3
13	GOL	O	1009	-	-	0/4/4/4	0/0/0/0
15	HEM	P	201	3,4	-	0/6/50/54	0/0/8/8
14	BHG	P	205	-	1/1/5/5	0/9/29/29	0/1/1/1

The worst 5 of 55 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	P	201	HEM	C2D-C3D	-4.98	1.39	1.54
15	C	142	HEM	CAB-C3B	-4.75	1.42	1.53
15	C	142	HEM	C2D-C3D	-4.71	1.40	1.54
15	P	201	HEM	CAB-C3B	-4.68	1.42	1.53
7	A	1001	FAD	PA-O2A	-3.63	1.39	1.54

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	N	1001	FAD	N3A-C2A-N1A	-13.41	118.62	128.89
7	A	1001	FAD	N3A-C2A-N1A	-13.13	118.84	128.89
7	N	1001	FAD	C4X-C10-N10	-7.15	116.31	120.52
7	A	1001	FAD	C4X-C10-N10	-6.85	116.48	120.52
7	N	1001	FAD	C4-C4X-C10	-5.80	116.23	119.94

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
14	C	141	BHG	C4
14	P	205	BHG	C4
8	A	1002	TEO	C2

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Mol	Chain	Res	Type	Atom
8	N	1002	TEO	C2

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1001	FAD	3	0
8	A	1002	TEO	3	0
13	B	1009	GOL	1	0
7	N	1001	FAD	3	0
8	N	1002	TEO	1	0
13	O	1009	GOL	1	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 ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	612/621 (98%)	-0.38	6 (0%) 84 89	14, 23, 45, 74	0
1	N	612/621 (98%)	-0.38	5 (0%) 87 91	12, 22, 43, 69	0
2	B	239/252 (94%)	-0.46	2 (0%) 87 91	15, 24, 41, 70	0
2	O	239/252 (94%)	-0.43	3 (1%) 79 85	14, 23, 37, 69	0
3	C	139/140 (99%)	-0.15	2 (1%) 78 84	21, 35, 50, 61	0
3	P	139/140 (99%)	0.34	16 (11%) 6 8	21, 37, 81, 92	0
4	D	101/103 (98%)	0.14	5 (4%) 32 37	26, 42, 58, 66	0
4	Q	101/103 (98%)	0.51	10 (9%) 9 11	27, 47, 69, 79	0
All	All	2182/2232 (97%)	-0.27	49 (2%) 65 72	12, 25, 54, 92	0

The worst 5 of 49 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	P	79	LEU	6.6
3	P	78	SER	4.6
1	A	10	THR	4.6
3	P	64	PRO	4.3
4	Q	88	TYR	4.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
9	UNL	D	107	6/-	0.92	0.29	9.44	66,68,69,70	0
9	UNL	Q	212	7/-	0.63	0.51	9.19	74,75,76,76	0
9	UNL	B	1007	1/-	0.79	0.19	7.74	61,61,61,61	0
9	UNL	P	220	2/-	0.77	0.28	5.96	81,81,81,81	0
13	GOL	O	1009	6/6	0.94	0.10	5.36	35,40,41,43	0
9	UNL	C	145	4/-	0.92	0.20	4.56	100,100,100,100	0
9	UNL	D	116	10/-	0.79	0.22	4.16	65,66,69,69	0
13	GOL	B	1009	6/6	0.94	0.09	3.44	33,36,38,38	0
9	UNL	D	114	8/-	0.93	0.16	1.93	61,63,64,65	0
9	UNL	P	209	2/-	0.82	0.15	1.80	49,49,49,49	0
9	UNL	P	229	1/-	0.89	0.11	1.75	54,54,54,54	0
12	F3S	B	1004	7/7	0.99	0.08	1.53	18,20,21,22	0
14	BHG	P	205	18/18	0.95	0.12	1.35	36,42,59,61	0
14	BHG	C	141	18/18	0.86	0.16	1.32	48,70,75,75	0
9	UNL	Q	219	1/-	0.74	0.12	1.11	60,60,60,60	0
11	SF4	O	1003	8/8	1.00	0.09	0.98	16,17,18,18	0
9	UNL	O	1006	1/-	0.82	0.12	0.97	54,54,54,54	0
6	AZI	A	623	3/3	0.90	0.11	0.95	62,62,63,64	0
12	F3S	O	1004	7/7	0.99	0.09	0.92	19,20,21,22	0
15	HEM	P	201	41/43	0.97	0.11	0.83	28,33,50,53	0
11	SF4	B	1003	8/8	1.00	0.08	0.81	17,19,19,19	0
9	UNL	Q	213	5/-	0.86	0.19	0.72	74,74,74,74	0
9	UNL	N	1011	1/-	0.89	0.13	0.67	37,37,37,37	0
9	UNL	B	1006	1/-	0.76	0.10	0.38	60,60,60,60	0
10	FES	O	1002	4/4	1.00	0.09	0.32	13,14,14,15	0
15	HEM	C	142	41/43	0.98	0.09	0.27	27,32,49,52	0
7	FAD	N	1001	53/53	0.98	0.08	0.19	11,14,18,21	0
7	FAD	A	1001	53/53	0.98	0.07	0.15	12,14,18,22	0
10	FES	B	1002	4/4	1.00	0.08	0.12	14,16,16,16	0
5	K	A	622	1/1	0.98	0.07	-0.28	24,24,24,24	0
5	K	N	622	1/1	0.98	0.07	-0.29	25,25,25,25	0
8	TEO	N	1002	9/9	0.97	0.06	-0.37	17,18,19,20	1
5	K	B	253	1/1	0.99	0.06	-0.38	38,38,38,38	0
5	K	O	253	1/1	0.99	0.06	-0.38	39,39,39,39	0
8	TEO	A	1002	9/9	0.99	0.06	-0.97	18,20,23,24	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
9	UNL	C	237	1/-	0.84	0.14	-	66,66,66,66	0
9	UNL	C	147	1/-	0.85	0.14	-	59,59,59,59	0
9	UNL	A	1005	1/-	0.78	0.13	-	63,63,63,63	0
9	UNL	Q	228	1/-	0.88	0.11	-	64,64,64,64	0
9	UNL	C	144	1/-	0.84	0.21	-	62,62,62,62	0
9	UNL	P	215	4/-	0.65	0.17	-	74,74,75,76	0
9	UNL	N	1009	1/-	0.85	0.17	-	54,54,54,54	0
9	UNL	A	1004	1/-	0.78	0.23	-	61,61,61,61	0
9	UNL	P	216	1/-	0.87	0.19	-	53,53,53,53	0
9	UNL	A	1008	1/-	0.84	0.25	-	64,64,64,64	0
9	UNL	N	1004	1/-	0.90	0.14	-	60,60,60,60	0
9	UNL	N	1007	1/-	0.95	0.18	-	44,44,44,44	0
9	UNL	N	1010	1/-	0.79	0.11	-	60,60,60,60	0
9	UNL	P	242	1/-	0.82	0.17	-	57,57,57,57	0
9	UNL	D	249	1/-	0.65	0.31	-	72,72,72,72	0
9	UNL	N	1006	1/-	0.90	0.14	-	54,54,54,54	0
6	AZI	N	623	3/3	0.82	0.09	-	68,68,70,71	0
9	UNL	C	238	1/-	0.77	0.10	-	59,59,59,59	0
9	UNL	D	108	1/-	0.46	0.14	-	63,63,63,63	0
9	UNL	D	250	1/-	0.75	0.12	-	73,73,73,73	0
9	UNL	Q	218	1/-	0.95	0.11	-	53,53,53,53	0
9	UNL	Q	256	1/-	0.66	0.20	-	84,84,84,84	0
9	UNL	P	217	1/-	-0.03	0.43	-	113,113,113,113	0
9	UNL	N	1005	1/-	0.77	0.25	-	76,76,76,76	0
9	UNL	N	1012	1/-	0.92	0.12	-	53,53,53,53	0
9	UNL	D	252	1/-	0.49	0.12	-	73,73,73,73	0
9	UNL	C	149	1/-	0.88	0.17	-	81,81,81,81	0
9	UNL	N	1003	2/-	0.32	0.39	-	114,114,114,114	0
9	UNL	P	231	1/-	0.67	0.21	-	74,74,74,74	0
9	UNL	P	240	1/-	0.85	0.12	-	56,56,56,56	0
9	UNL	O	1008	1/-	0.50	0.20	-	63,63,63,63	0
9	UNL	B	1008	1/-	0.65	0.20	-	69,69,69,69	0
9	UNL	P	230	1/-	0.81	0.18	-	69,69,69,69	0
9	UNL	P	241	1/-	0.81	0.23	-	59,59,59,59	0
9	UNL	P	214	4/-	0.60	0.76	-	94,94,94,95	0
9	UNL	A	1006	1/-	0.74	0.35	-	63,63,63,63	0
9	UNL	A	1003	1/-	0.88	0.25	-	69,69,69,69	0
9	UNL	N	1008	1/-	0.77	0.15	-	65,65,65,65	0
9	UNL	A	1007	2/-	0.62	0.33	-	65,65,65,66	0
9	UNL	O	1007	1/-	0.68	0.16	-	59,59,59,59	0
9	UNL	D	119	1/-	0.97	0.18	-	63,63,63,63	0
9	UNL	O	1005	1/-	0.98	0.10	-	35,35,35,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
9	UNL	C	146	4/-	0.76	0.46	-	108,109,109,109	0
9	UNL	C	148	1/-	0.96	0.14	-	56,56,56,56	0
9	UNL	D	254	1/-	0.67	0.10	-	67,67,67,67	0
9	UNL	B	1005	1/-	0.97	0.07	-	39,39,39,39	0
9	UNL	P	232	1/-	0.91	0.29	-	61,61,61,61	0
9	UNL	D	251	1/-	0.86	0.10	-	73,73,73,73	0
9	UNL	D	253	1/-	0.81	0.15	-	72,72,72,72	0
9	UNL	P	208	1/-	0.94	0.09	-	44,44,44,44	0
9	UNL	C	143	2/-	0.54	0.17	-	65,65,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.