



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 12:03 AM GMT

PDB ID : 1CMJ
Title : CRYSTAL STRUCTURES OF FERRIC-NO COMPLEXES OF FUNGAL NITRIC OXIDE REDUCTASE AND THEIR SER286 MUTANTS AT CRYOGENIC TEMPERATURE
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Deposited on : 1999-05-07
Resolution : 1.70 Å(reported)

This is a full wwPDB validation 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/ValidationPDFNotes.html>

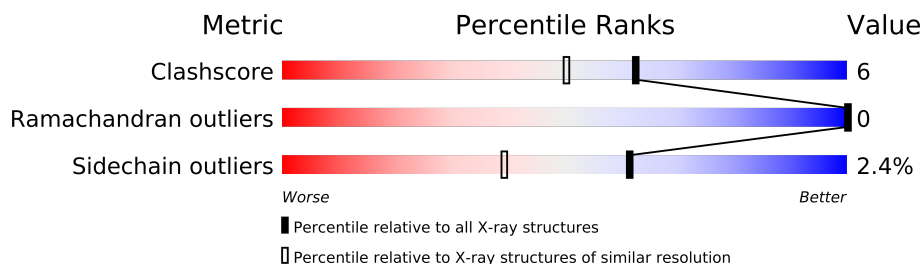
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	21963
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.70 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	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	402	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 3482 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CYTOCHROME P450.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	399	3100	1972	529	586	13	0	0	0

There is a discrepancy between the modelled and reference sequences:

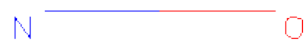
Chain	Residue	Modelled	Actual	Comment	Reference
A	286	THR	SER	ENGINEERED	UNP P23295

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 3 is NITRIC OXIDE (three-letter code: NO) (formula: NO).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	N	O	0	0
			2	1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	337	Total	O	0	0
			337	337		

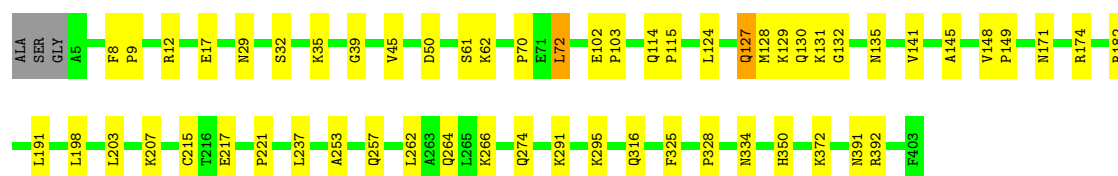
3 Residue-property plots

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: CYTOCHROME P450

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.69Å 81.13Å 85.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.70	Depositor
% Data completeness (in resolution range)	97.6 (50.00-1.70)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.197 , 0.250	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3482	wwPDB-VP
Average B, all atoms (Å ²)	16.0	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: HEM, NO

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	0.32	0/3167	0.55	0/4305

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3100	0	3117	35	0
2	A	43	0	30	0	0
3	A	2	0	0	0	0
4	A	337	0	0	5	0
All	All	3482	0	3147	35	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (35) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:264:GLN:HE22	1:A:334:ASN:HD21	1.32	0.75
1:A:130:GLN:C	1:A:132:GLY:H	1.91	0.72
1:A:12:ARG:HB3	1:A:17:GLU:O	1.91	0.70
1:A:141:VAL:HA	1:A:145:ALA:HB3	1.78	0.64
1:A:130:GLN:C	1:A:132:GLY:N	2.56	0.59
1:A:8:PHE:CG	1:A:9:PRO:HA	2.39	0.58
1:A:253:ALA:O	1:A:257:GLN:HG2	2.07	0.55
1:A:70:PRO:HG2	1:A:291:LYS:HB3	1.89	0.54
1:A:114:GLN:HB3	1:A:115:PRO:HD3	1.94	0.49
1:A:132:GLY:HA3	4:A:793:HOH:O	2.13	0.49
1:A:128:MET:HA	4:A:611:HOH:O	2.14	0.48
1:A:102:GLU:N	1:A:103:PRO:HD2	2.28	0.48
1:A:62:LYS:HD3	1:A:350:HIS:CE1	2.50	0.47
1:A:61:SER:HB2	1:A:295:LYS:HD2	1.96	0.47
1:A:127:GLN:HG2	4:A:765:HOH:O	2.15	0.47
1:A:274:GLN:HG3	4:A:614:HOH:O	2.15	0.46
1:A:171:ASN:O	1:A:174:ARG:HB2	2.17	0.45
1:A:129:LYS:HD2	1:A:372:LYS:HE3	1.99	0.45
1:A:29:ASN:ND2	1:A:32:SER:HB3	2.32	0.44
1:A:102:GLU:N	1:A:103:PRO:CD	2.80	0.44
1:A:262:LEU:HG	1:A:266:LYS:HE3	1.99	0.44
1:A:391:ASN:OD1	1:A:392:ARG:HG2	2.18	0.44
1:A:217:GLU:O	1:A:221:PRO:HG2	2.19	0.43
1:A:325:PHE:O	1:A:328:PRO:HD3	2.18	0.43
1:A:35:LYS:HE3	1:A:39:GLY:HA2	2.01	0.42
1:A:131:LYS:HG3	1:A:131:LYS:O	2.19	0.42
1:A:130:GLN:O	1:A:132:GLY:N	2.52	0.42
1:A:316:GLN:HG2	4:A:770:HOH:O	2.19	0.41
1:A:203:LEU:HD23	1:A:215:CYS:HB3	2.02	0.41
1:A:198:LEU:HA	1:A:198:LEU:HD23	1.87	0.41
1:A:45:VAL:CG1	1:A:50:ASP:HB3	2.51	0.41
1:A:8:PHE:CD1	1:A:9:PRO:HA	2.56	0.41
1:A:9:PRO:HD3	1:A:72:LEU:HG	2.02	0.41
1:A:148:VAL:HB	1:A:149:PRO:CD	2.50	0.40
1:A:148:VAL:HB	1:A:149:PRO:HD3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	397/402 (99%)	385 (97%)	12 (3%)	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 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	340/341 (100%)	332 (98%)	8 (2%)	61	39

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	LEU
1	A	124	LEU
1	A	127	GLN
1	A	135	ASN
1	A	182	ARG
1	A	191	LEU
1	A	207	LYS
1	A	237	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	118	GLN
1	A	130	GLN
1	A	170	GLN

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Mol	Chain	Res	Type
1	A	171	ASN
1	A	231	GLN
1	A	264	GLN
1	A	371	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

2 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	$\# Z > 2$
2	HEM	A	501	1,3	49,50,50	2.89	12 (24%)	46,82,82	1.39	4 (8%)
3	NO	A	502	2	0,1,1	0.00	-	0,0,0	0.00	-

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
2	HEM	A	501	1,3	-	0/14/114/114	0/0/8/8
3	NO	A	502	2	-	0/0/0/0	0/0/0/0

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C2D-C1D	12.01	1.47	1.44
2	A	501	HEM	C2B-C1B	9.98	1.47	1.44
2	A	501	HEM	C3D-C4D	6.03	1.46	1.44
2	A	501	HEM	C3D-C2D	-4.11	1.36	1.43
2	A	501	HEM	C4A-C3A	4.10	1.45	1.40
2	A	501	HEM	C3C-C2C	-3.39	1.37	1.43
2	A	501	HEM	CHA-C4D	3.27	1.40	1.35
2	A	501	HEM	C3B-C2B	-3.10	1.38	1.43
2	A	501	HEM	C4A-NA	2.49	1.41	1.36
2	A	501	HEM	CHB-C1B	2.42	1.39	1.35
2	A	501	HEM	CMB-C2B	2.34	1.54	1.47
2	A	501	HEM	CHD-C4C	2.21	1.40	1.36

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	C3B-C4B-NB	-4.02	111.12	114.00
2	A	501	HEM	C4A-CHB-C1B	-3.89	122.36	127.47
2	A	501	HEM	C1A-CHA-C4D	-3.67	122.64	127.47
2	A	501	HEM	CBA-CAA-C2A	-2.79	107.78	112.69

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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.