



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:45 PM UTC

PDB ID : 2ROM / pdb_00002rom
Title : CRYSTAL STRUCTURE OF NITRIC REDUCTASE FROM DENITRIFYING FUNGUS FUSARIUM OXYSPORUM COMPLEX WITH CARBON MONOXIDE
Authors : Park, S.-Y.; Nakagawa, A.
Deposited on : 1997-03-24
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

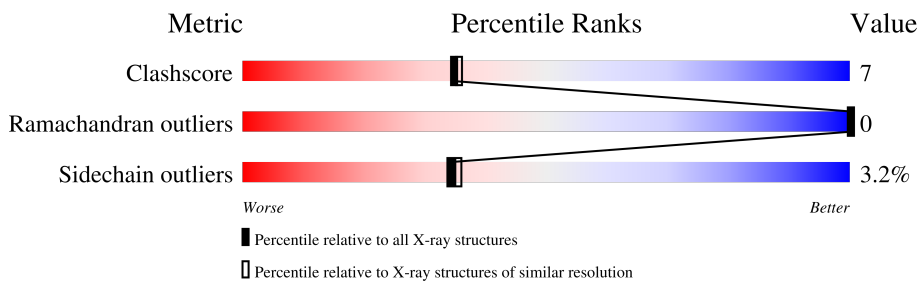
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	403	 79% 19% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3283 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 CYTOCHROME P450.

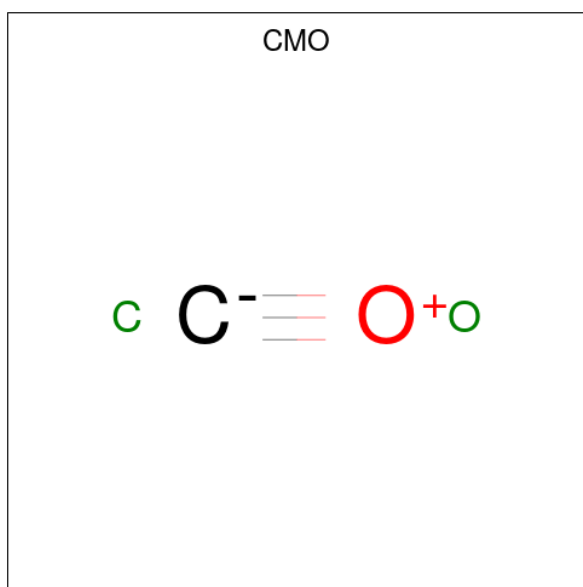
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	399	3099	1971	529	586	13	0	0	0

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: 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 CARBON MONOXIDE (CCD ID: CMO) (formula: CO).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	139	Total	O	0	0
			139	139		

4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.14Å 82.82Å 87.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00	Depositor
% Data completeness (in resolution range)	81.6 (10.00-2.00)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.199 , 0.269	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	3283	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

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: HEM, CMO

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	0.45	0/3166	0.98	13/4303 (0.3%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	303	LYS	N-CA-C	9.28	125.25	113.55
1	A	134	ALA	N-CA-C	7.00	118.56	111.07
1	A	29	ASN	CA-C-N	6.17	126.36	119.32
1	A	29	ASN	C-N-CA	6.17	126.36	119.32
1	A	371	GLN	N-CA-C	-5.87	104.83	112.23
1	A	239	ALA	N-CA-C	5.86	118.16	111.02
1	A	217	GLU	N-CA-C	5.76	119.48	112.23
1	A	164	LEU	N-CA-C	5.39	116.84	111.07
1	A	337	TRP	N-CA-C	5.30	116.50	109.93
1	A	341	ASP	N-CA-C	5.23	117.51	110.29
1	A	181	ALA	N-CA-C	-5.09	105.86	111.71
1	A	268	ASN	CA-C-N	5.09	124.75	119.56
1	A	268	ASN	C-N-CA	5.09	124.75	119.56

There are no chirality outliers.

There are no planarity outliers.

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	3099	0	3115	47	0
2	A	43	0	30	1	0
3	A	2	0	0	0	0
4	A	139	0	0	3	0
All	All	3283	0	3145	47	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 7.

All (47) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:GLN:HE22	1:A:334:ASN:HD21	1.21	0.84
1:A:264:GLN:HE22	1:A:334:ASN:ND2	1.87	0.73
1:A:264:GLN:NE2	1:A:334:ASN:HD21	1.89	0.69
1:A:171:ASN:ND2	1:A:174:ARG:HE	1.93	0.67
1:A:377:LYS:HE2	1:A:402:ILE:HD11	1.79	0.64
1:A:141:VAL:HA	1:A:145:ALA:HB3	1.78	0.64
1:A:129:LYS:HD2	1:A:372:LYS:HZ3	1.62	0.64
1:A:377:LYS:CE	1:A:402:ILE:HD11	2.28	0.63
1:A:180:THR:HG22	1:A:182:ARG:H	1.64	0.63
1:A:45:VAL:HG22	1:A:313:ALA:HA	1.81	0.62
1:A:126:GLU:O	1:A:130:GLN:HG2	2.00	0.62
1:A:220:LYS:HB2	1:A:221:PRO:HD3	1.81	0.62
1:A:174:ARG:HD2	1:A:188:ASN:OD1	1.99	0.62
1:A:58:GLU:H	1:A:58:GLU:CD	2.08	0.61
1:A:129:LYS:HD2	1:A:372:LYS:NZ	2.16	0.61
1:A:133:CYS:SG	1:A:402:ILE:HA	2.47	0.55
1:A:62:LYS:HE3	4:A:516:HOH:O	2.07	0.55
1:A:148:VAL:HB	1:A:149:PRO:HD3	1.87	0.55
1:A:288:LEU:HD11	4:A:577:HOH:O	2.08	0.54
1:A:146:LEU:HB3	1:A:147:PRO:HD3	1.90	0.54
1:A:392:ARG:HB3	1:A:392:ARG:NH1	2.26	0.50
1:A:86:PHE:HB3	1:A:97:GLN:HB3	1.92	0.50
1:A:59:LYS:HE3	1:A:296:GLU:OE2	2.13	0.48
1:A:354:ALA:HB2	2:A:501:HEM:HBC2	1.96	0.47
1:A:58:GLU:CD	1:A:58:GLU:N	2.73	0.47
1:A:326:GLU:O	1:A:335:ARG:NH2	2.48	0.47
1:A:7:SER:HA	1:A:35:LYS:O	2.16	0.45
1:A:377:LYS:HE3	1:A:402:ILE:HD11	1.98	0.45
1:A:255:LEU:HB3	1:A:262:LEU:HD13	1.99	0.45
1:A:180:THR:HG22	1:A:181:ALA:N	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:PRO:HG3	4:A:600:HOH:O	2.17	0.44
1:A:322:GLU:H	1:A:322:GLU:CD	2.26	0.43
1:A:85:THR:OG1	1:A:87:VAL:HG12	2.18	0.43
1:A:207:LYS:H	1:A:212:SER:CB	2.31	0.43
1:A:78:GLN:O	1:A:81:LYS:HG3	2.19	0.43
1:A:159:VAL:HA	1:A:160:PRO:HD3	1.89	0.43
1:A:241:ASN:O	1:A:245:VAL:HG23	2.19	0.42
1:A:128:MET:HE3	1:A:373:PHE:CE2	2.55	0.42
1:A:202:ARG:HD2	1:A:212:SER:OG	2.18	0.42
1:A:389:PRO:HG2	1:A:392:ARG:CD	2.49	0.42
1:A:226:LYS:HB2	1:A:226:LYS:HE3	1.78	0.42
1:A:111:LYS:HE3	1:A:111:LYS:HB2	1.89	0.41
1:A:51:VAL:HG13	1:A:313:ALA:HB1	2.02	0.41
1:A:380:VAL:HG22	1:A:381:PRO:HD2	2.03	0.41
1:A:303:LYS:O	1:A:304:LEU:HB3	2.21	0.41
1:A:64:ARG:HH22	1:A:79:ALA:HB3	1.85	0.40
1:A:70:PRO:HG2	1:A:291:LYS:HB3	2.02	0.40

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	397/403 (98%)	388 (98%)	9 (2%)	0	100 100

There are no Ramachandran outliers to report.

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	340/342 (99%)	329 (97%)	11 (3%)	34 35

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	LYS
1	A	72	LEU
1	A	81	LYS
1	A	146	LEU
1	A	167	LEU
1	A	182	ARG
1	A	217	GLU
1	A	237	LEU
1	A	250	LEU
1	A	394	VAL
1	A	399	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	114	GLN
1	A	171	ASN
1	A	201	GLN
1	A	241	ASN
1	A	274	GLN
1	A	316	GLN
1	A	332	ASN
1	A	334	ASN
1	A	371	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
3	CMO	A	502	2	0,1,1	-	-	-		
2	HEM	A	501	3,1	50,50,50	1.17	4 (8%)	67,82,82	1.24	7 (10%)

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	3,1	-	5/14/54/54	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	CAB-C3B	-3.40	1.38	1.47
2	A	501	HEM	CAC-C3C	-3.08	1.39	1.47
2	A	501	HEM	CMB-C2B	2.23	1.55	1.50
2	A	501	HEM	C4A-C3A	2.05	1.47	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	C4A-C3A-C2A	-3.65	102.64	106.82
2	A	501	HEM	C3B-C4B-NB	2.83	111.50	109.47
2	A	501	HEM	CMA-C3A-C4A	2.81	129.70	125.42
2	A	501	HEM	C2B-C1B-NB	2.79	113.04	109.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	C3B-C2B-C1B	-2.44	104.58	106.41
2	A	501	HEM	C1A-C2A-C3A	2.21	110.28	106.87
2	A	501	HEM	CAA-C2A-C3A	-2.13	122.29	127.07

There are no chirality outliers.

All (5) torsion outliers are listed below:

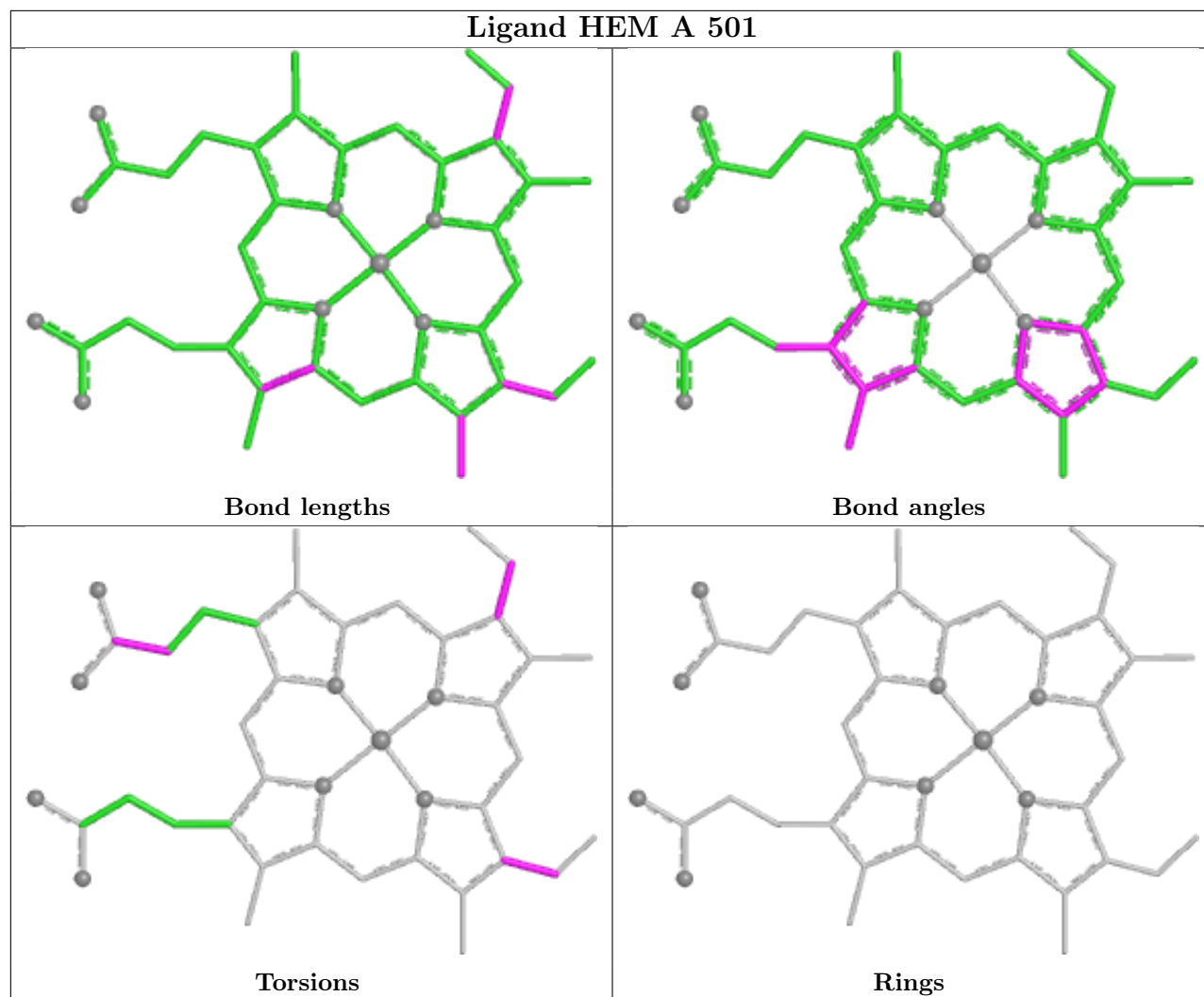
Mol	Chain	Res	Type	Atoms
2	A	501	HEM	C2C-C3C-CAC-CBC
2	A	501	HEM	C4C-C3C-CAC-CBC
2	A	501	HEM	C2B-C3B-CAB-CBB
2	A	501	HEM	C4B-C3B-CAB-CBB
2	A	501	HEM	CAD-CBD-CGD-O2D

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	HEM	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.