



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 3NLD
Title : Structure of endothelial nitric oxide synthase heme domain complexed with 6- $\{[(3'S,4'S)-3'-[2''-(3'''-fluorophenethylamino)ethoxy]pyrrolidin-4'-yl\}methyl\}$ -4-methylpyridin-2-amine
Authors : Ji, H.; Delker, S.L.; Li, H.; Martasek, P.; Roman, L.; Poulos, T.L.; Silverman, R.B.
Deposited on : 2010-06-21
Resolution : 2.29 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

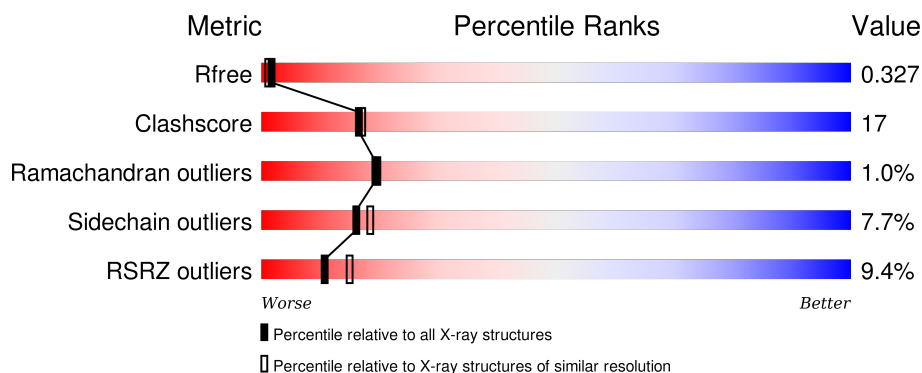
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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	444	
1	B	444	

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
4	ACT	A	850	-	-	X	X
4	ACT	B	850	-	-	X	-

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 6750 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 Nitric oxide synthase, endothelial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	403	Total	C	N	O	S	0	0	0
			3202	2036	563	587	16			
1	B	403	Total	C	N	O	S	0	0	0
			3209	2040	566	587	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ARG	CYS	SEE REMARK 999	UNP P29473
B	100	ARG	CYS	SEE REMARK 999	UNP P29473

- 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
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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

- Molecule 3 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula: $C_9H_{15}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	9	5	3		
3	B	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- [illegible]

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total 27	C 21	F 1	N 4	O 1	0	0
6	B	1	Total 27	C 21	F 1	N 4	O 1	0	0

-
- Chemical structure of As₂O₃ (arsenic trioxide). The central atom is As (Arsenic, pink). It is bonded to three oxygen atoms (red) and one arsenic atom (pink). The bonds are: a single bond to C1 (left), a double bond to O (right), and a single bond to O H (bottom). The arsenic atom is also bonded to C2 (top) via a single bond. The labels are: C1 (green), C2 (green), AS (green), O (red), O H (red), and O1 (green).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	As	C	0	0
			3	1	2		
7	B	1	Total	As	C	0	0
			3	1	2		

- Molecule 8 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Zn	0	0
			1	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	70	Total	O	0	0
			70	70		
9	B	68	Total	O	0	0
			68	68		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.40Å 106.89Å 157.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.92 – 2.29 36.92 – 2.29	Depositor EDS
% Data completeness (in resolution range)	98.8 (36.92-2.29) 98.8 (36.92-2.29)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0089	Depositor
R, R_{free}	0.222 , 0.292 0.270 , 0.327	Depositor DCC
R_{free} test set	2201 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.927	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 39.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 44143 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6750	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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: GOL, ZN, H4B, JSS, ACT, HEM, CAD

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	0.89	3/3291 (0.1%)	0.87	4/4483 (0.1%)
1	B	1.08	9/3298 (0.3%)	0.98	7/4491 (0.2%)
All	All	0.99	12/6589 (0.2%)	0.93	11/8974 (0.1%)

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	253	TYR	CE2-CZ	11.37	1.53	1.38
1	B	253	TYR	CG-CD2	8.30	1.50	1.39
1	B	383	VAL	CB-CG1	7.09	1.67	1.52
1	B	385	MET	CG-SD	6.66	1.98	1.81
1	B	283	PRO	C-O	6.21	1.35	1.23

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	374	ARG	NE-CZ-NH1	7.92	124.26	120.30
1	A	367	ARG	NE-CZ-NH1	-7.11	116.75	120.30
1	B	367	ARG	NE-CZ-NH2	-6.35	117.13	120.30
1	B	374	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	B	395	LEU	CA-CB-CG	-5.70	102.18	115.30

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	3202	0	3107	88	0
1	B	3209	0	3117	127	0
2	A	43	0	30	3	0
2	B	43	0	30	3	0
3	A	17	0	15	0	0
3	B	17	0	15	1	0
4	A	4	0	3	4	0
4	B	4	0	3	6	0
5	A	6	0	8	1	0
5	B	6	0	8	0	0
6	A	27	0	29	2	0
6	B	27	0	29	1	0
7	A	3	0	0	1	0
7	B	3	0	0	1	0
8	A	1	0	0	0	0
9	A	70	0	0	11	0
9	B	68	0	0	5	0
All	All	6750	0	6394	216	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 17.

The worst 5 of 216 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:B:384:CYS:SG	7:B:950:CAD:AS	2.48	1.32
1:A:384:CYS:SG	7:A:950:CAD:AS	2.66	1.13
1:B:368:ASN:HD21	4:B:850:ACT:H2	1.10	1.11
1:A:236:ARG:HD3	1:A:242:ASP:OD1	1.68	0.93
1:B:368:ASN:ND2	4:B:850:ACT:H2	1.82	0.92

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	399/444 (90%)	363 (91%)	31 (8%)	5 (1%)	15	14
1	B	399/444 (90%)	352 (88%)	44 (11%)	3 (1%)	24	26
All	All	798/888 (90%)	715 (90%)	75 (9%)	8 (1%)	19	20

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	455	SER
1	A	240	ARG
1	B	103	GLY
1	B	108	PRO
1	A	103	GLY

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	342/377 (91%)	317 (93%)	25 (7%)	17	20
1	B	343/377 (91%)	315 (92%)	28 (8%)	14	16
All	All	685/754 (91%)	632 (92%)	53 (8%)	16	18

5 of 53 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	458	LEU
1	B	105	LEU
1	B	398	ASP
1	A	469	TYR
1	B	99	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
1	B	91	GLN
1	B	178	GLN
1	B	340	ASN
1	A	468	ASN
1	B	368	ASN

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 13 ligands modelled in this entry, 1 is monoatomic - leaving 12 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
2	HEM	A	500	1	30,50,50	2.60	9 (30%)	24,82,82	2.64	12 (50%)
3	H4B	A	600	-	13,18,18	1.12	1 (7%)	11,26,26	2.42	6 (54%)
6	JSS	A	800	-	26,29,29	0.78	1 (3%)	27,38,38	1.53	3 (11%)
4	ACT	A	850	-	1,3,3	0.83	0	0,3,3	0.00	-
5	GOL	A	880	-	5,5,5	0.56	0	5,5,5	0.48	0
7	CAD	A	950	-	0,2,4	0.00	-	0,1,6	0.00	-
2	HEM	B	500	1	30,50,50	2.43	10 (33%)	24,82,82	2.62	12 (50%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	H4B	B	600	-	13,18,18	1.69	4 (30%)	11,26,26	3.11	9 (81%)
6	JSS	B	800	-	26,29,29	0.92	1 (3%)	27,38,38	1.83	5 (18%)
4	ACT	B	850	-	1,3,3	0.07	0	0,3,3	0.00	-
5	GOL	B	880	-	5,5,5	0.56	0	5,5,5	0.49	0
7	CAD	B	950	-	0,2,4	0.00	-	0,1,6	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	500	1	-	0/10/54/54	0/0/8/8
3	H4B	A	600	-	-	0/8/17/17	0/2/2/2
6	JSS	A	800	-	-	0/13/23/23	0/3/3/3
4	ACT	A	850	-	-	0/0/0/0	0/0/0/0
5	GOL	A	880	-	-	0/4/4/4	0/0/0/0
7	CAD	A	950	-	-	0/0/0/0	0/0/0/0
2	HEM	B	500	1	-	0/10/54/54	0/0/8/8
3	H4B	B	600	-	-	0/8/17/17	0/2/2/2
6	JSS	B	800	-	-	0/13/23/23	0/3/3/3
4	ACT	B	850	-	-	0/0/0/0	0/0/0/0
5	GOL	B	880	-	-	0/4/4/4	0/0/0/0
7	CAD	B	950	-	-	0/0/0/0	0/0/0/0

The worst 5 of 26 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	HEM	C3B-C4B	-8.58	1.44	1.51
2	A	500	HEM	C3B-C4B	-8.12	1.44	1.51
2	B	500	HEM	C3D-C4D	-5.39	1.44	1.51
2	A	500	HEM	C3D-C4D	-5.18	1.44	1.51
3	B	600	H4B	C4A-C8A	-4.05	1.33	1.41

The worst 5 of 47 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	600	H4B	N3-C2-N1	-5.73	116.15	125.53
2	B	500	HEM	CBA-CAA-C2A	-5.30	103.04	112.53
2	A	500	HEM	C3B-CAB-CBB	-4.91	116.92	124.46
2	B	500	HEM	CAA-C2A-C1A	-3.91	122.76	127.01
3	A	600	H4B	N3-C2-N1	-3.80	119.31	125.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	HEM	3	0
6	A	800	JSS	2	0
4	A	850	ACT	4	0
5	A	880	GOL	1	0
7	A	950	CAD	1	0
2	B	500	HEM	3	0
3	B	600	H4B	1	0
6	B	800	JSS	1	0
4	B	850	ACT	6	0
7	B	950	CAD	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	403/444 (90%)	0.49	43 (10%) 8 11	37, 58, 84, 99	0
1	B	403/444 (90%)	0.54	33 (8%) 14 19	43, 61, 83, 107	0
All	All	806/888 (90%)	0.52	76 (9%) 11 15	37, 60, 84, 107	0

The worst 5 of 76 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	144	GLY	5.0
1	A	67	GLY	4.3
1	A	200	ALA	4.1
1	B	140	ILE	4.0
1	A	337	ALA	4.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
4	ACT	A	850	4/4	0.87	0.35	3.03	66,67,70,71	0
6	JSS	A	800	27/27	0.89	0.28	1.50	40,57,92,94	0
4	ACT	B	850	4/4	0.95	0.23	1.10	63,64,67,67	0
6	JSS	B	800	27/27	0.90	0.21	0.73	35,48,82,86	0
5	GOL	B	880	6/6	0.93	0.18	0.69	55,55,57,58	0
3	H4B	A	600	17/17	0.96	0.21	0.47	43,49,54,54	0
2	HEM	A	500	43/43	0.97	0.20	0.33	41,48,60,68	0
3	H4B	B	600	17/17	0.94	0.17	0.16	39,43,50,54	0
2	HEM	B	500	43/43	0.97	0.17	0.12	41,48,60,61	0
8	ZN	A	900	1/1	0.99	0.13	-0.19	60,60,60,60	0
7	CAD	A	950	3/5	0.99	0.12	-1.20	87,87,88,88	0
7	CAD	B	950	3/5	0.98	0.10	-1.84	85,85,85,87	0
5	GOL	A	880	6/6	0.90	0.40	-	75,76,78,79	0

6.5 Other polymers [i](#)

There are no such residues in this entry.