



Full wwPDB X-ray Structure Validation Report

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PDB ID : 1NTF
Title : Crystal Structure of Cimex Nitrophorin
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Deposited on : 2003-01-29
Resolution : 1.80 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>

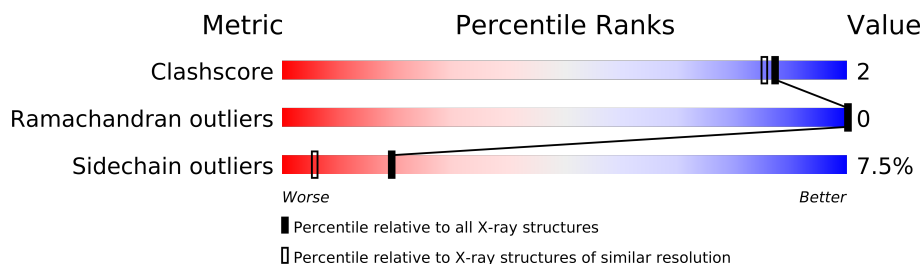
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.80 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	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (1.80-1.80)

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	282	

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 2433 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 salivary nitrophorin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	280	Total	C	N	O	S	0	0	0
			2227	1407	373	442	5			

- 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		

- Molecule 3 is water.

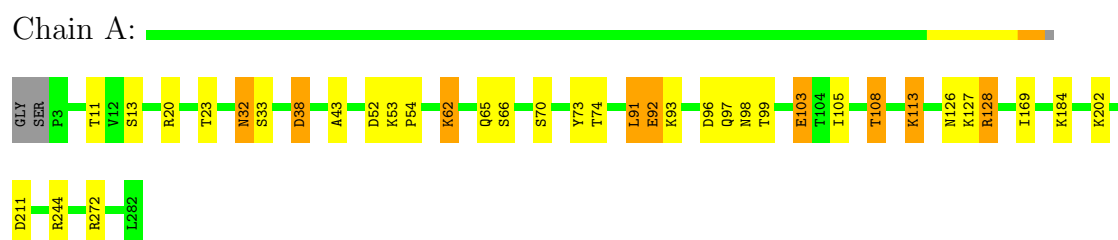
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	163	Total	O	0	0
			163	163		

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: salivary nitrophorin



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.20Å 42.10Å 65.85Å 90.00° 95.88° 90.00°	Depositor
Resolution (Å)	21.00 – 1.80	Depositor
% Data completeness (in resolution range)	99.7 (21.00-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
Refinement program	REFMAC 5.1.22	Depositor
R, R_{free}	0.151 , 0.198	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2433	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: HEM

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	1.06	1/2276 (0.0%)	1.03	6/3088 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	184	LYS	CD-CE	6.10	1.66	1.51

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	272	ARG	NE-CZ-NH2	7.20	123.90	120.30
1	A	96	ASP	CB-CG-OD1	6.42	124.08	118.30
1	A	38	ASP	CB-CG-OD2	5.68	123.41	118.30
1	A	244	ARG	NE-CZ-NH2	-5.53	117.54	120.30
1	A	52	ASP	CB-CG-OD2	5.42	123.18	118.30
1	A	211	ASP	CB-CG-OD2	5.30	123.07	118.30

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	2227	0	2188	11	0
2	A	43	0	30	0	0
3	A	163	0	0	0	0
All	All	2433	0	2218	11	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 2.

All (11) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:74:THR:OG1	1:A:92:GLU:HG2	1.82	0.78
1:A:32:ASN:HD22	1:A:32:ASN:H	1.46	0.64
1:A:38:ASP:HA	1:A:91:LEU:HD22	1.84	0.58
1:A:73:TYR:CE2	1:A:91:LEU:HD13	2.42	0.55
1:A:13:SER:HA	1:A:43:ALA:O	2.17	0.45
1:A:126:ASN:C	1:A:127:LYS:HD3	2.37	0.45
1:A:103:GLU:HG3	1:A:128:ARG:NH2	2.32	0.44
1:A:108:THR:HG23	1:A:113:LYS:HA	1.98	0.44
1:A:62:LYS:O	1:A:66:SER:OG	2.30	0.42
1:A:53:LYS:N	1:A:54:PRO:HD2	2.35	0.42
1:A:11:THR:HB	1:A:169:ILE:HG21	2.02	0.41

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	278/282 (99%)	271 (98%)	7 (2%)	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	252/253 (100%)	233 (92%)	19 (8%)	19 6

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ARG
1	A	23	THR
1	A	32	ASN
1	A	33	SER
1	A	62	LYS
1	A	65	GLN
1	A	70	SER
1	A	91	LEU
1	A	92	GLU
1	A	93	LYS
1	A	97	GLN
1	A	98	ASN
1	A	99	THR
1	A	103	GLU
1	A	105	ILE
1	A	108	THR
1	A	113	LYS
1	A	128	ARG
1	A	202	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	63	ASN
1	A	145	ASN

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 ⓘ

1 ligand is 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	300	1,3	49,50,50	2.54	16 (32%)	46,82,82	2.60	14 (30%)

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	300	1,3	-	0/14/114/114	0/0/8/8

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	300	HEM	C2B-C1B	9.29	1.46	1.44
2	A	300	HEM	C3C-C2C	-5.12	1.34	1.43
2	A	300	HEM	C3D-C2D	5.08	1.52	1.43
2	A	300	HEM	C3C-CAC	4.49	1.54	1.40
2	A	300	HEM	C3B-C2B	-4.03	1.36	1.43
2	A	300	HEM	FE-NB	3.97	2.12	1.97
2	A	300	HEM	FE-ND	3.92	2.12	1.97
2	A	300	HEM	C3B-CAB	3.83	1.52	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	300	HEM	FE-NC	3.78	2.12	1.97
2	A	300	HEM	FE-NA	3.72	2.08	1.92
2	A	300	HEM	C4A-C3A	3.52	1.44	1.40
2	A	300	HEM	CMC-C2C	2.68	1.55	1.47
2	A	300	HEM	C1C-NC	-2.44	1.34	1.38
2	A	300	HEM	CMB-C2B	2.37	1.54	1.47
2	A	300	HEM	CHA-C4D	2.14	1.38	1.35
2	A	300	HEM	CMD-C2D	2.04	1.53	1.47

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	300	HEM	C3B-C4B-NB	-9.59	107.14	114.00
2	A	300	HEM	C4D-ND-C1D	7.29	112.63	105.16
2	A	300	HEM	C1B-NB-C4B	5.04	110.32	105.16
2	A	300	HEM	CHC-C4B-NB	3.96	127.88	124.58
2	A	300	HEM	CMA-C3A-C4A	-3.93	122.57	128.62
2	A	300	HEM	CBA-CAA-C2A	-3.81	105.98	112.69
2	A	300	HEM	C2D-C1D-ND	-3.49	108.81	112.93
2	A	300	HEM	CMA-C3A-C2A	3.05	130.70	124.94
2	A	300	HEM	CAD-C3D-C4D	2.42	128.89	124.53
2	A	300	HEM	CBD-CAD-C3D	-2.41	109.10	114.37
2	A	300	HEM	C4C-NC-C1C	2.33	107.96	105.53
2	A	300	HEM	O1D-CGD-CBD	-2.29	115.17	123.03
2	A	300	HEM	O1A-CGA-CBA	-2.20	115.47	123.03
2	A	300	HEM	CHD-C4C-NC	2.14	126.59	124.73

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.