

wwPDB X-ray Structure Validation Summary Report (i)

Aug 7, 2023 – 01:55 PM EDT

PDB ID	:	1NI6
Title	:	Comparisions of the Heme-Free and-Bound Crystal Structures of Human Heme
		Oxygenase-1
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Deposited on		
Resolution	:	2.10 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 https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

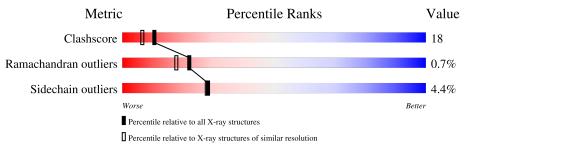
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.10 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain					
1	А	224	68%	24% • •				
1	В	224	61%	33% • •				
1	С	224	62%	35% •				
1	D	224	56%	38% • 5%				
2	Е	2	100%					



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7529 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.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	٨	215	Total	С	Ν	0	S	0	0	0
1	А	215	1754	1125	300	324	5	0		0
1	В	214	Total	С	Ν	Ο	S	0	0	0
1	I D	214	1745	1121	299	320	5			
1	С	223	Total	С	Ν	Ο	S	0	0	0
1	U	223	1818	1164	312	335	7	0	0	0
1	Л	213	Total	С	Ν	Ο	S	0	0	0
		213	1738	1116	298	319	5		U	0

• Molecule 1 is a protein called Heme oxygenase 1.

• Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-1)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace	
2	Е	2	Total 23	C 12	0 11	0	0	0

• Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	5	ZeroOcc	AltConf
3	С	1	Total C	Cl 1	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	156	Total O 156 156	0	0
4	В	110	Total O 110 110	0	0

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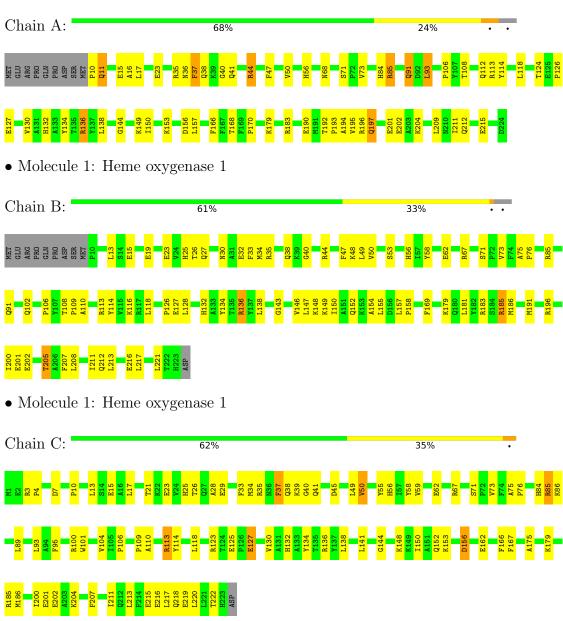
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	С	131	Total O 131 131	0	0
4	D	53	$\begin{array}{cc} \text{Total} & \text{O} \\ 53 & 53 \end{array}$	0	0



Residue-property plots (i) 3

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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: Heme oxygenase 1

• Molecule 1: Heme oxygenase 1





• Molecule 2: alpha-D-glucopyranose-(1-1)-alpha-D-glucopyranose

Chain E:

100%







4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	76.21Å 55.51Å 108.00Å	Depositor
a, b, c, α , β , γ	90.00° 98.88° 90.00°	Depositor
Resolution (Å)	50.00 - 2.10	Depositor
% Data completeness	93.9 (50.00-2.10)	Depositor
(in resolution range)	30.3 (00.00 2.10)	Depositor
R_{merge}	0.07	Depositor
R _{sym}	0.05	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.217 , 0.267	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7529	wwPDB-VP
Average B, all atoms $(Å^2)$	37.0	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: GLC, CL

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	
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.36	0/1795	0.57	0/2425
1	В	0.35	0/1786	0.56	0/2414
1	С	0.33	0/1861	0.56	0/2516
1	D	0.31	0/1778	0.52	0/2403
All	All	0.34	0/7220	0.55	0/9758

There are no bond length outliers.

There are no bond angle outliers.

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	А	1754	0	1742	45	0
1	В	1745	0	1738	66	0
1	С	1818	0	1808	73	0
1	D	1738	0	1730	79	0
2	Е	23	0	21	0	0
3	С	1	0	0	0	0
4	А	156	0	0	12	0
4	В	110	0	0	7	0
4	С	131	0	0	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	53	0	0	5	0
All	All	7529	0	7039	261	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 18.

The worst 5 of 261 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:C:113:ARG:HB3	1:C:113:ARG:HH11	1.12	1.09
1:B:44:ARG:NH1	1:B:48:LYS:HD2	1.86	0.91
1:A:192:THR:HG22	1:A:194:ALA:H	1.36	0.89
1:A:192:THR:HB	1:A:195:VAL:HG23	1.58	0.85
1:A:130:VAL:HG23	4:A:355:HOH:O	1.78	0.83

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	А	213/224~(95%)	207~(97%)	5(2%)	1 (0%)	29 26
1	В	212/224~(95%)	204 (96%)	7 (3%)	1 (0%)	29 26
1	С	221/224~(99%)	213~(96%)	7 (3%)	1 (0%)	29 26
1	D	211/224 (94%)	196 (93%)	12 (6%)	3 (1%)	11 6
All	All	857/896~(96%)	820 (96%)	31 (4%)	6 (1%)	22 18

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	11	GLN
	<i>a</i>	7	

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Mol	Chain	Res	Type
1	В	154	ALA
1	D	159	SER
1	D	154	ALA
1	D	156	ASP

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric Outliers		Percentiles		
1	А	185/194~(95%)	174 (94%)	11 (6%)	19	17	
1	В	184/194~(95%)	178 (97%)	6 (3%)	38	40	
1	С	193/194~(100%)	182 (94%)	11 (6%)	20	18	
1	D	183/194~(94%)	178 (97%)	5(3%)	44	48	
All	All	745/776~(96%)	712~(96%)	33 (4%)	28	28	

 $5~{\rm of}~33$ residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	D	32	GLU
1	D	93	LEU
1	D	157	LEU
1	В	116	LYS
1	В	114	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 46 such sidechains are listed below:

Mol	Chain	Res	Type
1	С	38	GLN
1	С	171	ASN
1	С	41	GLN
1	С	84	HIS
1	D	11	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)

2 monosaccharides 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).

Mal	Turne	Chain	Dec	Link			nd lengths		Bond angles		
	Type	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2	
2	GLC	Е	1	2	11,11,12	1.69	1 (9%)	$15,\!15,\!17$	1.02	1 (6%)	
2	GLC	Е	2	2	12,12,12	1.81	3 (25%)	$17,\!17,\!17$	1.19	1 (5%)	

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	GLC	Е	1	2	-	2/2/19/22	0/1/1/1
2	GLC	Е	2	2	-	0/2/22/22	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Ε	2	GLC	C4-C5	3.49	1.60	1.53
2	Е	1	GLC	O5-C1	3.35	1.49	1.43
2	Е	2	GLC	O5-C1	2.68	1.49	1.42
2	Е	2	GLC	C3-C2	2.29	1.58	1.52

All (2) bond angle outliers are listed below:



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Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
2	Е	1	GLC	C2-C3-C4	-3.02	105.66	110.89
2	Е	2	GLC	C4-C3-C2	-2.70	106.11	110.82

There are no chirality outliers.

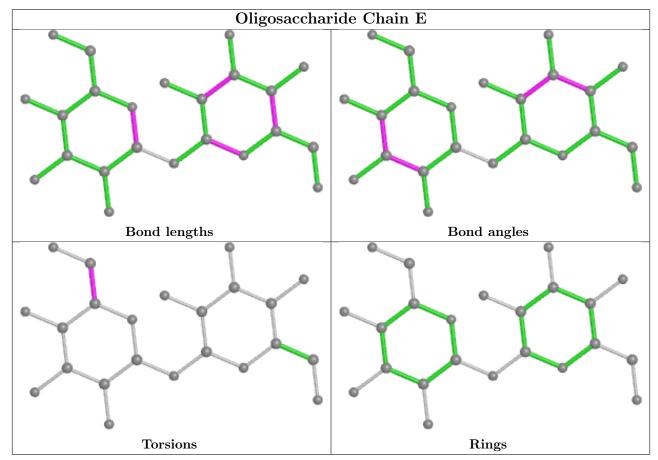
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	Е	1	GLC	C4-C5-C6-O6
2	Е	1	GLC	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.



There are no bond length outliers. There are no bond angle outliers. There are no chirality outliers. There are no torsion outliers. There are no ring outliers. No monomer is involved in short contacts.

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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

