

Full wwPDB X-ray Structure Validation Report (i)

May 17, 2020 – 08:49 pm BST

PDB ID	:	4KNV
Title	:	The crystal structure of APO HUMAN HDHD4 FROM SE-MAD
Authors	:	Klei, H.E.
Deposited on		
Resolution	:	1.99 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

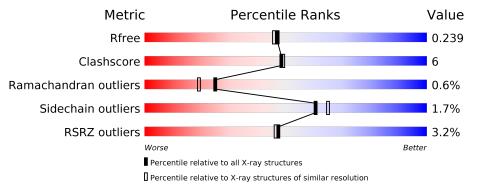
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 on 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% 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	А	241	3% 87%	11% •
1	В	241	3% 82%	14% • •



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4012 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	А	241	Total 1896	C 1197		O 361		${ m Se} 6$	0	2	0
1	В	234	Total 1836	C 1158		O 348		Se 6	0	1	0

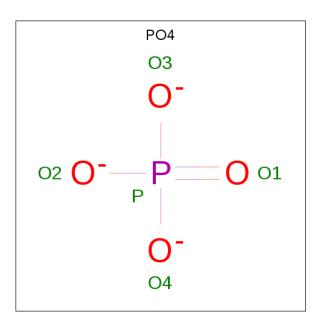
• Molecule 1 is a protein called N-acylneuraminate-9-phosphatase.

Chain	Residue	Modelled	Actual	Comment	Reference
А	6	MSE	-	EXPRESSION TAG	UNP Q8TBE9
A	243	LEU	-	EXPRESSION TAG	UNP Q8TBE9
А	244	VAL	-	EXPRESSION TAG	UNP Q8TBE9
А	245	PRO	-	EXPRESSION TAG	UNP Q8TBE9
А	246	ARG	-	EXPRESSION TAG	UNP Q8TBE9
В	6	MSE	-	EXPRESSION TAG	UNP Q8TBE9
В	243	LEU	-	EXPRESSION TAG	UNP Q8TBE9
В	244	VAL	-	EXPRESSION TAG	UNP Q8TBE9
В	245	PRO	-	EXPRESSION TAG	UNP Q8TBE9
В	246	ARG	-	EXPRESSION TAG	UNP Q8TBE9

There are 10 discrepancies between the modelled and reference sequences:

• Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0
2	В	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0

• Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	В	1	Total Mg 1 1	0	0
3	А	1	Total Mg 1 1	0	0

• Molecule 4 is water.

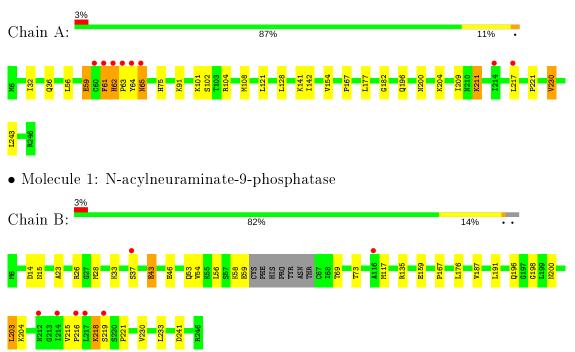
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	134	Total O 134 134	0	0
4	В	134	Total O 134 134	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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 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.

• Molecule 1: N-acylneuraminate-9-phosphatase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants	46.34Å 53.44 Å 63.96 Å	Depositor
a, b, c, α , β , γ	65.50° 74.99° 85.17°	Depositor
Resolution (Å)	33.36 - 1.99	Depositor
Resolution (A)	33.36 - 1.99	EDS
% Data completeness	91.0 (33.36-1.99)	Depositor
(in resolution range)	91.5(33.36-1.99)	EDS
R _{merge}	0.07	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	4.43 (at 2.00Å)	Xtriage
Refinement program	PHENIX ?	Depositor
D D	0.200 , 0.246	Depositor
R, R_{free}	0.194 , 0.239	DCC
R_{free} test set	991 reflections (2.95%)	wwPDB-VP
Wilson B-factor (Å ²)	17.4	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.36 , 43.6	EDS
L-test for twinning ²	$ \langle L \rangle = 0.50, \langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4012	wwPDB-VP
Average B, all atoms $(Å^2)$	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 10.16% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: PO4, MG $\,$

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	0.39	0/1934	0.54	0/2609	
1	В	0.39	0/1864	0.52	0/2509	
All	All	0.39	0/3798	0.53	0/5118	

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	1896	0	1904	22	0
1	В	1836	0	1859	24	0
2	А	5	0	0	0	0
2	В	5	0	0	0	0
3	А	1	0	0	0	0
3	В	1	0	0	0	0
4	А	134	0	0	1	0
4	В	134	0	0	2	0
All	All	4012	0	3763	45	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 6.



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:218:LYS:HE2	1:B:219:SER:H	1.35	0.88
1:B:218:LYS:HE2	1:B:219:SER:N	1.88	0.88
1:A:61:PHE:O	1:A:62:HIS:HB2	1.93	0.69
1:B:53:GLN:NE2	4:B:456:HOH:O	2.27	0.67
1:B:117:MSE:HE2	1:B:230:VAL:O	1.97	0.65
1:B:187:VAL:HG21	1:B:233:LEU:HD11	1.82	0.61
1:A:128:LEU:HB2	1:A:177:LEU:HD11	1.82	0.61
1:A:167:PRO:HG3	1:A:200:ASN:HB3	1.82	0.61
1:B:216:PRO:HG2	4:B:430:HOH:O	2.01	0.60
1:A:64:TYR:O	1:A:65:ASN:HB2	2.01	0.59
1:A:56:LEU:O	1:A:59:GLU:HG3	2.01	0.59
1:B:28:MSE:HE3	1:B:53:GLN:HG2	1.85	0.58
1:B:167:PRO:HG3	1:B:200:ASN:HB3	1.85	0.58
1:B:23:ALA:HA	1:B:26[B]:ARG:HH11	1.69	0.57
1:A:209:ILE:HD11	1:A:230:VAL:HG23	1.87	0.55
1:A:102:SER:HB2	1:B:176:LEU:HD23	1.89	0.54
1:A:91:LYS:HD3	4:A:525:HOH:O	2.07	0.53
1:A:61:PHE:O	1:A:62:HIS:CB	2.57	0.52
1:A:182:GLY:O	1:A:204:LYS:HE2	2.10	0.52
1:B:56:LEU:O	1:B:59:GLU:HB3	2.09	0.51
1:B:196:GLN:OE1	1:B:221:PRO:HG3	2.11	0.51
1:B:69:THR:O	1:B:73:THR:HG23	2.11	0.51
1:A:32:ILE:O	1:A:36:GLN:HG3	2.11	0.50
1:B:117:MSE:HE1	1:B:230:VAL:HG22	1.94	0.50
1:B:198:GLY:HA2	1:B:203:LEU:HD22	1.94	0.50
1:B:135:ARG:HG3	1:B:159:GLU:HG2	1.95	0.49
1:B:54:VAL:CG1	1:B:58:LYS:HE3	2.44	0.47
1:A:62:HIS:HB2	1:A:63:PRO:HD3	1.96	0.47
1:B:191:LEU:HD11	1:B:216:PRO:HG3	1.96	0.46
1:A:108:MSE:SE	1:A:141:LYS:HG2	2.65	0.46
1:A:121:LEU:CD2	1:A:243:LEU:HD11	2.45	0.45
1:B:218:LYS:O	1:B:219:SER:CB	2.64	0.45
1:B:204:LYS:HE3	1:B:241:ASP:O	2.17	0.45
1:A:59:GLU:OE1	1:A:75:HIS:HE1	2.02	0.43
1:A:61:PHE:HD1	1:A:62:HIS:H	1.65	0.43
1:B:54:VAL:HG12	1:B:58:LYS:HE3	2.00	0.43
1:A:182:GLY:O	1:A:204:LYS:HB2	2.19	0.42
1:A:101:LYS:O	1:A:104:ARG:HG2	2.20	0.42
1:A:142:ILE:HD13	1:A:154:VAL:HG11	2.02	0.42
1:B:14:ASP:O	1:B:15:ASN:HB2	2.18	0.42

All (45) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:211:LYS:HD3	1:A:211:LYS:HA	1.48	0.42
1:A:217:LEU:H	1:A:217:LEU:HG	1.65	0.41
1:B:33:LYS:O	1:B:37:SER:HB3	2.21	0.41
1:A:196:GLN:OE1	1:A:221:PRO:HG3	2.21	0.41
1:B:43:GLU:O	1:B:46:GLU:HG2	2.21	0.40

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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	241/241~(100%)	230~(95%)	8(3%)	3~(1%)	13 7
1	В	231/241~(96%)	223~(96%)	8 (4%)	0	100 100
All	All	472/482~(98%)	453~(96%)	16 (3%)	3 (1%)	25 19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	61	PHE
1	А	62	HIS
1	А	65	ASN

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	А	211/204~(103%)	208~(99%)	3 (1%)	67 72
1	В	203/204~(100%)	199~(98%)	4 (2%)	55 58
All	All	414/408~(102%)	407~(98%)	7 (2%)	60 65

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	59	GLU
1	А	211	LYS
1	А	230	VAL
1	В	43	GLU
1	В	203	LEU
1	В	215	VAL
1	В	218	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	75	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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 length (or angles).

ſ	Mal	Iol Type Chain Res	Chain I	Dog	Link	B	ond leng	gths	В	ond ang	gles
	WIOI			Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2		
	2	PO4	А	301	3	4,4,4	1.14	0	6,6,6	0.63	0
	2	PO4	В	301	3	4,4,4	1.11	0	6,6,6	0.38	0

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)

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, 95^{th} 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	А	235/241~(97%)	-0.03	8 (3%) 45 44	9, 17, 32, 85	0
1	В	$228/241 \ (94\%)$	-0.01	7 (3%) 49 48	9, 18, 33, 50	0
All	All	463/482~(96%)	-0.02	15 (3%) 47 46	9, 18, 33, 85	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	64	TYR	12.9
1	А	61	PHE	8.7
1	А	65	ASN	7.2
1	В	217	LEU	6.9
1	А	60	CYS	5.4
1	А	63	PRO	5.4
1	А	62	HIS	4.4
1	В	219	SER	4.4
1	А	217	LEU	4.0
1	В	216	PRO	3.3
1	В	214	ILE	2.3
1	А	214	ILE	2.1
1	В	116	ALA	2.1
1	В	212	ASN	2.1
1	В	37	SER	2.0

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. The B-factors column lists the minimum, median, 95^{th} 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	\mathbf{RSR}	$\mathbf{B} ext{-factors}(\mathbf{A}^2)$	$Q{<}0.9$
3	MG	В	302	1/1	0.97	0.04	$16,\!16,\!16,\!16$	0
2	PO4	А	301	5/5	0.99	0.08	8,10,11,14	0
2	PO4	В	301	5/5	0.99	0.08	$10,\!10,\!12,\!13$	0
3	MG	А	302	1/1	0.99	0.06	$13,\!13,\!13,\!13$	0

6.5 Other polymers (i)

There are no such residues in this entry.

