

Full wwPDB X-ray Structure Validation Report (i)

Jul 31, 2023 – 01:26 AM EDT

PDB ID : 1BN5

Title : HUMAN METHIONINE AMINOPEPTIDASE 2

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Deposited on : 1998-07-31

Resolution : 1.80 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 types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) 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)

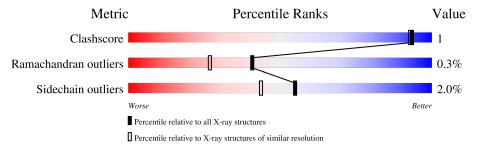
Validation Pipeline (wwPDB-VP) : 2.34

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 1.80 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	Similar resolution		
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$		
Clashscore	141614	6793 (1.80-1.80)		
Ramachandran outliers	138981	6697 (1.80-1.80)		
Sidechain outliers	138945	6696 (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 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	A	478	67%	6% •	26%				



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3031 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 METHIONINE AMINOPEPTIDASE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	A	355	Total 2780	C 1750	N 475	O 529	S 26	0	3	0

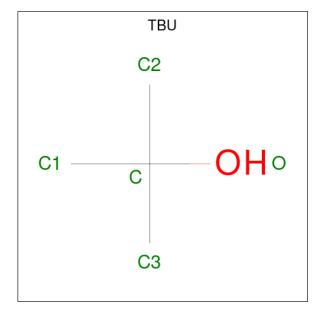
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference	
Α	347	ILE	VAL	conflict	UNP P50579	

• Molecule 2 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Co 2 2	0	0

• Molecule 3 is TERTIARY-BUTYL ALCOHOL (three-letter code: TBU) (formula: C₄H₁₀O).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 5 4 1	0	0
3	A	1	Total C O 5 4 1	0	0

\bullet Molecule 4 is water.

M	lol	Chain	Residues	Atoms	ZeroOcc	AltConf
	4	A	239	Total O 239 239	0	0

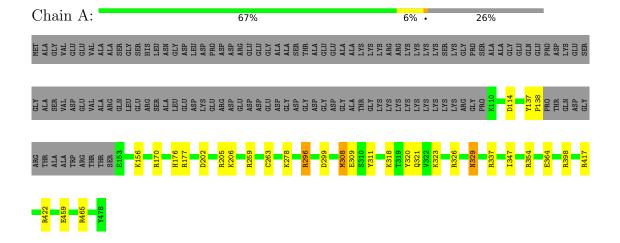


3 Residue-property plots (i)

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: METHIONINE AMINOPEPTIDASE





4 Data and refinement statistics (i)

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

Property	Value	Source	
Space group	C 2 2 21	Depositor	
Cell constants	90.49Å 99.38Å 101.59Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	25.00 - 1.80	Depositor	
% Data completeness	98.4 (25.00-1.80)	Depositor	
(in resolution range)	30.4 (29.00 1.00)		
R_{merge}	0.07	Depositor	
R_{sym}	0.07	Depositor	
Refinement program	REFMAC	Depositor	
R, R_{free}	0.183 , 0.228	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	3031	wwPDB-VP	
Average B, all atoms (Å ²)	21.0	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: TBU, CO

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	Boı	nd lengths	Во	ond angles
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.61	$1/2850 \ (0.0\%)$	1.26	$21/3855 \ (0.5\%)$

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	A	138	PRO	N-CD	5.38	1.55	1.47

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^{o})$	$\operatorname{Ideal}(^{o})$
1	A	417	ARG	NE-CZ-NH2	-20.14	110.23	120.30
1	A	422	ARG	NE-CZ-NH2	-19.58	110.51	120.30
1	A	205	ARG	NE-CZ-NH1	-11.76	114.42	120.30
1	A	465	ARG	NE-CZ-NH1	9.77	125.18	120.30
1	A	296	ARG	CD-NE-CZ	9.49	136.89	123.60
1	A	422	ARG	NE-CZ-NH1	8.67	124.64	120.30
1	A	259	ARG	NE-CZ-NH1	7.63	124.11	120.30
1	A	308	MET	CG-SD-CE	7.42	112.08	100.20
1	A	354	ARG	CD-NE-CZ	6.77	133.08	123.60
1	A	417	ARG	NH1-CZ-NH2	6.69	126.76	119.40
1	A	177	ARG	NE-CZ-NH2	-6.54	117.03	120.30
1	A	170	ARG	NE-CZ-NH1	-6.42	117.09	120.30
1	A	465	ARG	NE-CZ-NH2	-6.15	117.22	120.30
1	A	422	ARG	CD-NE-CZ	5.82	131.75	123.60
1	A	398	ARG	CD-NE-CZ	5.80	131.72	123.60
1	A	329	ASN	CA-CB-CG	5.51	125.52	113.40
1	A	417	ARG	NE-CZ-NH1	5.37	122.98	120.30
1	A	337	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	A	354	ARG	NE-CZ-NH1	5.31	122.96	120.30
1	A	137	TYR	C-N-CD	5.24	139.41	128.40

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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
1	A	205	ARG	NH1-CZ-NH2	5.18	125.09	119.40

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	2780	0	2722	8	0
2	A	2	0	0	0	0
3	A	10	0	20	0	0
4	A	239	0	0	1	0
All	All	3031	0	2742	8	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 1.

All (8) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{array}{c} { m Clash} \\ { m overlap} \ ({ m \AA}) \end{array}$
1:A:326:ARG:HB2	1:A:347:ILE:HD13	1.74	0.70
1:A:296:ARG:HD3	1:A:299:ASP:OD2	2.13	0.49
1:A:318:LYS:HE2	1:A:320:TYR:CE2	2.48	0.48
1:A:364:GLU:HB3	1:A:459:GLU:HG3	1.98	0.45
1:A:309:GLU:O	1:A:321:GLN:NE2	2.52	0.43
1:A:176:HIS:CE1	1:A:263:CYS:HB3	2.54	0.42
1:A:347:ILE:HD12	4:A:599:HOH:O	2.20	0.40
1:A:278:LYS:NZ	1:A:311:TYR:OH	2.55	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	354/478 (74%)	346 (98%)	7 (2%)	1 (0%)	41 27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	ASP

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	alysed Rotameric C		Percentiles	
1	A	306/399 (77%)	300 (98%)	6 (2%)	55 44	

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	156	LYS
1	A	202	ASP
1	A	206	LYS
1	A	308	MET
1	A	323	LYS
1	A	329	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:



Mol	Chain	Res	Type
1	A	160	GLN
1	A	329	ASN

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 monosaccharides 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 lengths (or angles).

Mal	Mal True Chair Dag		Dog Link		Bond lengths			Bond angles		
Mol	Type	Chain	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
3	TBU	A	689	-	4,4,4	0.81	0	6,6,6	1.87	2 (33%)
3	TBU	A	690	-	4,4,4	0.86	0	6,6,6	1.91	1 (16%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
3	A	690	TBU	O-C-C2	3.84	118.12	107.90
3	A	689	TBU	O-C-C2	3.52	117.26	107.90
3	A	689	TBU	O-C-C1	-2.27	101.87	107.90



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.

