Full wwPDB X-ray Structure Validation Report ⓘ

May 12, 2020 – 11:13 pm BST

PDB ID	:	1PNG
Title	:	CRYSTAL STRUCTURE OF PEPTIDE-N(4)-(N-ACETYL-BETA-D-GLUCOSAMINYL) ASPARAGINE AMIDASE AT 2.2 ANGSTROMS RESOLUTION
Authors	:	Van Roey, P.; Kuhn, P.
Deposited on	:	1994-06-02
Resolution	:	2.20 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 ⓘ symbol.

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

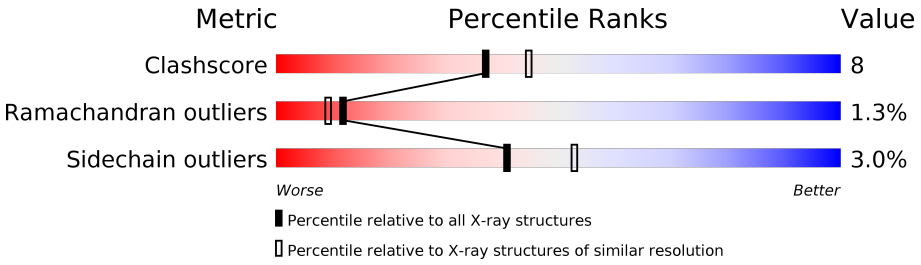
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.20 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	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	314	81% 16%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2655 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 PEPTIDE-N(4)-(N-ACETYL-BETA-D-GLUCOSAMINYL) ASPARAGINE AMIDASE F.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	310	2433	1550	408	467	8	0	0	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	222	222	222	0	0

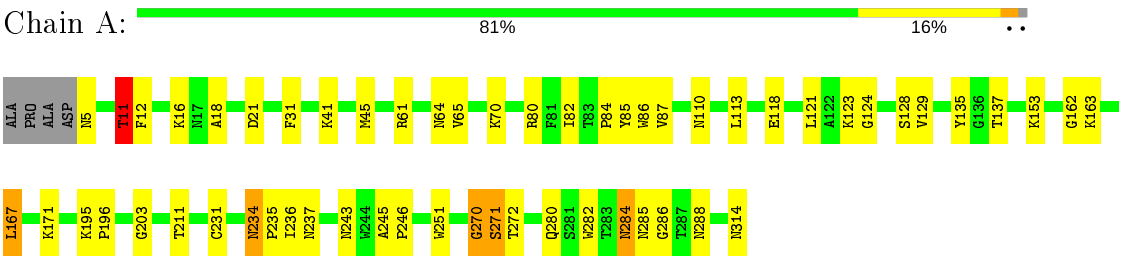
3 Residue-property plots (i)

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. 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: PEPTIDE-N(4)-(N-ACETYL-BETA-D-GLUCOSAMINYL)ASPARAGINE AMIDASE F

Chain A:



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 a, b, c, α, β, γ	87.16Å 125.10Å 79.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.10 – 2.20	Depositor
% Data completeness (in resolution range)	93.5 (49.10-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	unknown	Depositor
R, R_{free}	0.183 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2655	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	0.58	0/2500	0.84	2/3407 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	203	GLY	O-C-N	-6.53	112.26	122.70
1	A	11	THR	N-CA-C	5.28	125.24	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	11	THR	Mainchain

5.2 Too-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 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	2433	0	2342	37	0
2	A	222	0	0	4	0
All	All	2655	0	2342	37	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 8.

All (37) 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:A:45:MET:HE1	1:A:65:VAL:HG11	1.63	0.80
1:A:163:LYS:HE2	2:A:389:HOH:O	1.85	0.75
1:A:41:LYS:HE3	1:A:137:THR:H	1.51	0.74
1:A:21:ASP:OD1	1:A:21:ASP:O	2.12	0.67
1:A:70:LYS:HE3	2:A:342:HOH:O	1.95	0.65
1:A:195:LYS:HG3	1:A:288:ASN:HB2	1.79	0.65
1:A:21:ASP:OD1	1:A:21:ASP:C	2.36	0.64
1:A:284:ASN:ND2	1:A:286:GLY:H	1.96	0.64
1:A:41:LYS:HE3	1:A:137:THR:N	2.14	0.61
1:A:234:ASN:HD22	1:A:235:PRO:N	2.01	0.58
1:A:270:GLY:O	1:A:271:SER:HB3	2.04	0.58
1:A:195:LYS:HG3	1:A:288:ASN:CB	2.35	0.57
1:A:45:MET:CE	1:A:65:VAL:HG11	2.34	0.56
1:A:211:THR:H	1:A:280:GLN:HE21	1.54	0.55
1:A:171:LYS:NZ	1:A:314:ASN:HD21	2.11	0.48
1:A:12:PHE:O	1:A:128:SER:HA	2.12	0.48
1:A:16:LYS:HE3	1:A:123:LYS:O	2.14	0.47
1:A:195:LYS:HB3	1:A:196:PRO:HA	1.95	0.47
1:A:70:LYS:HE2	1:A:110:ASN:O	2.14	0.47
1:A:234:ASN:ND2	1:A:236:ILE:H	2.12	0.47
1:A:64:ASN:HB3	1:A:80:ARG:HB3	1.97	0.46
1:A:11:THR:O	1:A:31:PHE:CZ	2.71	0.44
1:A:5:ASN:ND2	1:A:135:TYR:CD2	2.85	0.44
1:A:61:ARG:HD2	1:A:118:GLU:O	2.17	0.44
1:A:284:ASN:HD22	1:A:285:ASN:N	2.15	0.44
1:A:171:LYS:HZ3	1:A:314:ASN:HD21	1.65	0.44
1:A:211:THR:H	1:A:280:GLN:NE2	2.14	0.44
1:A:153:LYS:HG2	2:A:424:HOH:O	2.17	0.44
1:A:231:CYS:HB2	1:A:245:ALA:O	2.17	0.44
1:A:162:GLY:HA2	1:A:282:TRP:O	2.18	0.44
1:A:45:MET:HE3	1:A:129:VAL:HG11	2.00	0.43
1:A:167:LEU:HA	2:A:394:HOH:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:ILE:HG13	1:A:251:TRP:CE3	2.53	0.42
1:A:243:ASN:HD21	1:A:246:PRO:HG2	1.85	0.42
1:A:16:LYS:HE2	1:A:18:ALA:HB2	2.01	0.41
1:A:85:TYR:O	1:A:87:VAL:N	2.52	0.41
1:A:234:ASN:C	1:A:234:ASN:HD22	2.23	0.41

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	308/314 (98%)	294 (96%)	10 (3%)	4 (1%)	12 9

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	271	SER
1	A	86	TRP
1	A	124	GLY
1	A	270	GLY

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	A	263/265 (99%)	255 (97%)	8 (3%)	41 53

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	PRO
1	A	113	LEU
1	A	121	LEU
1	A	167	LEU
1	A	234	ASN
1	A	237	ASN
1	A	272	THR
1	A	284	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	8	ASN
1	A	17	ASN
1	A	110	ASN
1	A	234	ASN
1	A	237	ASN
1	A	243	ASN
1	A	280	GLN
1	A	284	ASN
1	A	314	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 carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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 is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.