



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 19, 2023 – 12:12 PM JST

PDB ID : 8I0A
Title : Structure of alpha-L-Arabinofuranosidase from Trametes hirsuta
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Deposited on : 2023-01-10
Resolution : 2.01 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

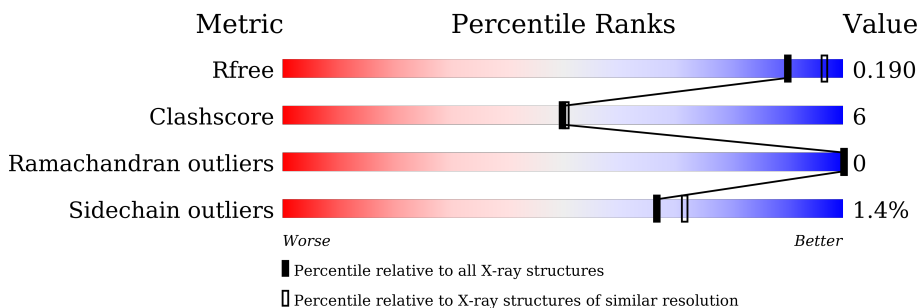
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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))
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)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	534	 87% 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8492 atoms, of which 4026 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 alpha-L-arabinofuranosidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	524	8208	2670	4026	734	768	10	0	0	0

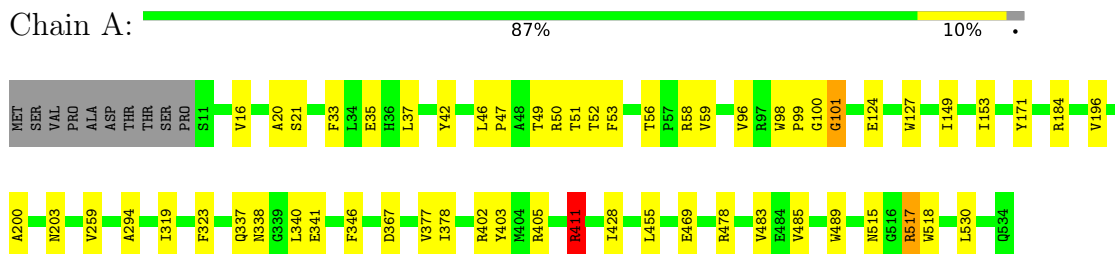
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	284	Total	O	0	0
			284	284		

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.

- Molecule 1: alpha-L-arabinofuranosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	134.94Å 134.94Å 175.56Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	55.44 – 2.01 116.86 – 1.21	Depositor EDS
% Data completeness (in resolution range)	100.0 (55.44-2.01) 48.0 (116.86-1.21)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.37 (at 1.21Å)	Xtrriage
Refinement program	PHENIX 1.20.1	Depositor
R, R_{free}	0.190 , 0.194 0.191 , 0.190	Depositor DCC
R_{free} test set	1735 reflections (0.80%)	wwPDB-VP
Wilson B-factor (Å ²)	5.8	Xtrriage
Anisotropy	0.035	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8492	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.69	3/4303 (0.1%)	0.75	1/5871 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	411	ARG	CZ-NH1	6.70	1.41	1.33
1	A	411	ARG	CZ-NH2	6.37	1.41	1.33
1	A	101	GLY	N-CA	-6.22	1.36	1.46

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	101	GLY	N-CA-C	-5.72	98.80	113.10

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	4182	4026	4047	48	0
2	A	284	0	0	3	0
All	All	4466	4026	4047	48	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.

All (48) 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:37:LEU:HD21	1:A:340:LEU:CD1	1.75	1.15
1:A:37:LEU:HD21	1:A:340:LEU:HD12	1.35	1.03
1:A:294:ALA:HB2	1:A:428:ILE:HD11	1.47	0.93
1:A:37:LEU:CD2	1:A:340:LEU:HD12	2.01	0.90
1:A:37:LEU:HD21	1:A:340:LEU:HD11	1.58	0.85
1:A:483:VAL:HG23	1:A:530:LEU:HD23	1.62	0.82
1:A:377:VAL:HG12	1:A:378:ILE:HG12	1.63	0.80
1:A:58:ARG:HG3	2:A:710:HOH:O	1.92	0.69
1:A:49:THR:HG22	1:A:51:THR:H	1.57	0.68
1:A:52:THR:HG22	1:A:53:PHE:O	1.94	0.68
1:A:515:ASN:OD1	1:A:517:ARG:HG2	1.94	0.68
1:A:405:ARG:HG2	2:A:692:HOH:O	1.98	0.63
1:A:483:VAL:CG2	1:A:530:LEU:HD23	2.30	0.61
1:A:37:LEU:CD2	1:A:340:LEU:CD1	2.61	0.60
1:A:101:GLY:HA2	1:A:203:ASN:OD1	2.01	0.59
1:A:294:ALA:HB2	1:A:428:ILE:CD1	2.30	0.55
1:A:485:VAL:HG21	1:A:518:TRP:CH2	2.43	0.54
1:A:259:VAL:CG2	1:A:319:ILE:HG13	2.39	0.53
1:A:377:VAL:HG12	1:A:378:ILE:CG1	2.36	0.51
1:A:37:LEU:HD13	1:A:127:TRP:CG	2.46	0.51
1:A:101:GLY:HA2	1:A:203:ASN:CG	2.31	0.50
1:A:340:LEU:HB2	1:A:377:VAL:HG13	1.94	0.49
1:A:171:TYR:O	1:A:184:ARG:HD3	2.12	0.49
1:A:20:ALA:HB3	1:A:411:ARG:HG3	1.96	0.48
1:A:49:THR:HG22	1:A:51:THR:N	2.26	0.48
1:A:478:ARG:HG3	1:A:478:ARG:HH11	1.78	0.48
1:A:49:THR:CG2	1:A:50:ARG:N	2.77	0.47
1:A:469:GLU:OE2	1:A:517:ARG:NH1	2.48	0.46
1:A:402:ARG:HD3	1:A:403:TYR:CZ	2.51	0.46
1:A:346:PHE:CZ	1:A:489:TRP:HA	2.50	0.46
1:A:485:VAL:HG21	1:A:518:TRP:HH2	1.80	0.46
1:A:259:VAL:HG22	1:A:319:ILE:HG13	1.97	0.46
1:A:33:PHE:HE1	1:A:100:GLY:HA2	1.81	0.45
1:A:42:TYR:O	1:A:52:THR:HG23	2.16	0.45
1:A:101:GLY:C	1:A:203:ASN:ND2	2.70	0.45
1:A:153:ILE:HG13	1:A:196:VAL:HG11	1.99	0.44
1:A:56:THR:CG2	1:A:59:VAL:H	2.31	0.43
1:A:98:TRP:CE2	1:A:99:PRO:HB3	2.54	0.43
1:A:20:ALA:CB	1:A:411:ARG:HG3	2.49	0.43
1:A:377:VAL:O	1:A:378:ILE:C	2.57	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:VAL:HG23	1:A:149:ILE:HG21	2.00	0.42
1:A:35:GLU:HA	1:A:99:PRO:O	2.19	0.42
1:A:46:LEU:HB3	1:A:47:PRO:HD2	2.02	0.42
1:A:16:VAL:HG11	1:A:455:LEU:HD21	2.01	0.42
1:A:98:TRP:CG	1:A:99:PRO:HA	2.55	0.41
1:A:337:GLN:OE1	1:A:337:GLN:N	2.51	0.41
1:A:338:ASN:CG	1:A:341:GLU:HG2	2.42	0.40
1:A:200:ALA:HB1	2:A:693:HOH:O	2.22	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	522/534 (98%)	506 (97%)	16 (3%)	0	100 100

There are no Ramachandran outliers to report.

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	441/450 (98%)	435 (99%)	6 (1%)	67 72

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

Mol	Chain	Res	Type
1	A	21	SER
1	A	124	GLU
1	A	323	PHE
1	A	367	ASP
1	A	411	ARG
1	A	517	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

There are no ligands in this entry.

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

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.