

wwPDB X-ray Structure Validation Summary Report (i)

Apr 24, 2024 – 04:08 PM EDT

PDB ID	:	8VFN
Title	:	Crystal Structure of WT D-Dopachrome Tautomerase (D-DT) at 310K
Authors	:	Parkins, A.; Pilien, A.; Wolff, A.; Thompson, M.C.; Pantouris, G.
Deposited on		
Resolution	:	1.29 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:

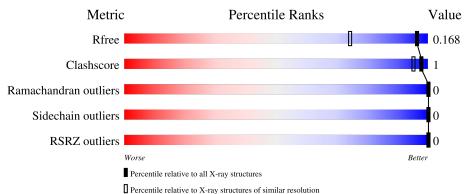
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 1.29 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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% 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	117	100%
1	В	117	97% .
1	С	117	100%



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2894 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	1 A	117	Total	С	Ν	0	S	0	0	0
		117	885	564	154	163	4			
1	1 B	117	Total	С	Ν	0	S	0	0	0
			885	564	154	163	4			
1 C	117	Total	С	Ν	0	S	0	1	0	
		890	567	154	165	4				

• Molecule 1 is a protein called D-dopachrome decarboxylase.

• Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	86	Total O 86 86	0	0
2	В	79	Total O 79 79	0	0
2	С	69	Total O 69 69	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 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: D-dopachrome decarboxylase

Chain A:

in A: 100%

There are no outlier residues recorded for this chain.

• Molecule 1: D-dopachrome decarboxylase

Chain B:



• Molecule 1: D-dopachrome decarboxylase

Chain C:

100%

97%

There are no outlier residues recorded for this chain.



4 Data and refinement statistics (i)

Property	Value	Source
Space group	Р 3	Depositor
Cell constants	84.29Å 84.29Å 41.06Å	Dereciter
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	73.00 - 1.29	Depositor
Resolution (A)	73.00 - 1.29	EDS
% Data completeness	97.9 (73.00-1.29)	Depositor
(in resolution range)	97.9 (73.00-1.29)	EDS
R _{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.43 (at 1.29 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
D D	0.159 , 0.167	Depositor
R, R_{free}	0.160 , 0.168	DCC
R_{free} test set	4085 reflections $(5.08%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	12.7	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35 , 43.3	EDS
L-test for twinning ²	$< L > = 0.51, < L^2 > = 0.34$	Xtriage
	0.015 for -h,-k,l	
Estimated twinning fraction	0.023 for h,-h-k,-l	Xtriage
	0.012 for -k,-h,-l	
F_o, F_c correlation	0.97	EDS
Total number of atoms	2894	wwPDB-VP
Average B, all atoms $(Å^2)$	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 19.64% 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)

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	Bo	nd lengths	Bond angles		
	Ullalli	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.73	0/902	0.83	0/1221	
1	В	0.76	1/902~(0.1%)	0.82	0/1221	
1	С	0.70	0/910	0.80	0/1232	
All	All	0.73	1/2714~(0.0%)	0.82	0/3674	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	В	22	LEU	C-N	-5.49	1.21	1.34

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	885	0	906	0	0
1	В	885	0	906	4	0
1	С	890	0	910	0	0
2	А	86	0	0	0	0
2	В	79	0	0	0	0
2	С	69	0	0	0	0
All	All	2894	0	2722	4	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 (4) 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:B:82:GLU:HG3	1:B:86:LYS:NZ	1.72	1.03
1:B:82:GLU:HG3	1:B:86:LYS:HZ1	1.48	0.75
1:B:82:GLU:HG3	1:B:86:LYS:HZ2	1.61	0.61
1:B:82:GLU:O	1:B:86:LYS:NZ	2.30	0.55

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	115/117~(98%)	112 (97%)	3~(3%)	0	100	100	
1	В	115/117~(98%)	112 (97%)	3~(3%)	0	100	100	
1	С	116/117~(99%)	113 (97%)	3(3%)	0	100	100	
All	All	346/351~(99%)	337~(97%)	9~(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	А	95/95~(100%)	95~(100%)	0	100 100	
1	В	95/95~(100%)	95 (100%)	0	100 100	
1	С	96/95~(101%)	96 (100%)	0	100 100	
All	All	286/285~(100%)	286 (100%)	0	100 100	

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	А	92	GLN
1	В	92	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)

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 (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	$\langle RSRZ \rangle$	#	₽RSR	Z>2	$OWAB(Å^2)$	Q < 0.9
1	А	117/117~(100%)	-0.49	0	100	100	9, 14, 26, 35	0
1	В	117/117~(100%)	-0.43	0	100	100	9, 14, 26, 32	0
1	С	$117/117 \ (100\%)$	-0.41	0	100	100	11, 17, 28, 38	0
All	All	351/351~(100%)	-0.44	0	100	100	9, 15, 27, 38	0

There are no RSRZ outliers to report.

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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.

