



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 12:58 am GMT

PDB ID : 3CSK  
Title : Structure of DPP III from *Saccharomyces cerevisiae*  
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Deposited on : 2008-04-10  
Resolution : 1.95 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](mailto: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.

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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 : trunk30967  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk30967

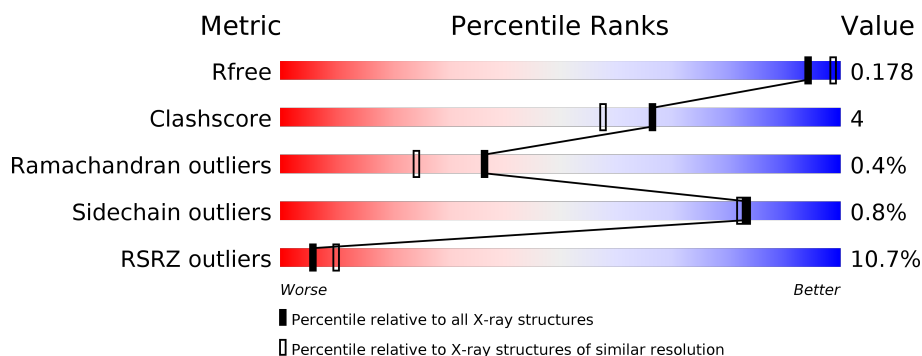
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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))
$R_{free}$	111664	2220 (1.96-1.96)
Clashscore	122126	2333 (1.96-1.96)
Ramachandran outliers	120053	2314 (1.96-1.96)
Sidechain outliers	120020	2314 (1.96-1.96)
RSRZ outliers	108989	2174 (1.96-1.96)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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  $\leq 5\%$ . 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	A	711	<div> <div>11%</div> <div>90%</div> <div>10%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6153 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 Probable dipeptidyl-peptidase 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	710	Total	C	N	O	S	0	3	0
			5700	3654	942	1087	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	130	SER	CYS	ENGINEERED	UNP Q08225
A	458	ASP	GLY	CONFLICT	UNP Q08225

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		

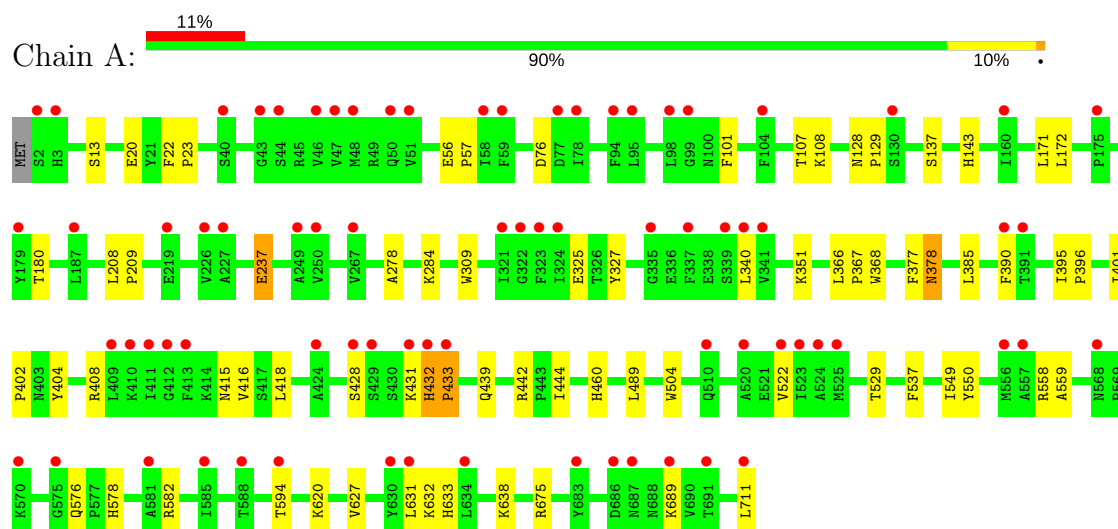
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	450	Total	O	0	0
			450	450		

### 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: Probable dipeptidyl-peptidase 3



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.62Å 110.12Å 67.91Å 90.00° 113.45° 90.00°	Depositor
Resolution (Å)	30.00 – 1.95 30.64 – 1.95	Depositor EDS
% Data completeness (in resolution range)	95.5 (30.00-1.95) 96.1 (30.64-1.95)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.97 (at 1.95Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.192 , 0.228 0.188 , 0.178	Depositor DCC
$R_{free}$ test set	2856 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.6	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 50.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6153	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 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	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.22	0/5842	0.37	0/7899

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

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	5700	0	5606	49	0
2	A	1	0	0	0	0
3	A	2	0	0	0	0
4	A	450	0	0	2	0
All	All	6153	0	5606	49	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 4.

All (49) 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:432:HIS:HB3	1:A:433:PRO:HD2	1.61	0.83
1:A:432:HIS:HB3	1:A:433:PRO:CD	2.16	0.76
1:A:444:ILE:HD13	1:A:550:TYR:HD2	1.53	0.72
1:A:431:LYS:O	1:A:432:HIS:HB2	1.91	0.69
1:A:378:ASN:HD22	1:A:378:ASN:N	1.92	0.66
1:A:489:LEU:HD13	1:A:633:HIS:CD2	2.33	0.63
1:A:401:ILE:HB	1:A:402:PRO:HA	1.80	0.61
1:A:431:LYS:O	1:A:432:HIS:CB	2.51	0.59
1:A:432:HIS:CB	1:A:433:PRO:CD	2.81	0.59
1:A:576:GLN:HB3	1:A:578[B]:HIS:CE1	2.39	0.58
1:A:620:LYS:HE2	1:A:620:LYS:HA	1.88	0.54
1:A:378:ASN:ND2	1:A:378:ASN:N	2.56	0.53
1:A:504:TRP:CZ3	1:A:638:LYS:HG3	2.44	0.52
1:A:377:PHE:C	1:A:378:ASN:HD22	2.12	0.51
1:A:367:PRO:O	1:A:632:LYS:HG2	2.10	0.51
1:A:439:GLN:HG2	1:A:442:ARG:NH2	2.27	0.49
1:A:340:LEU:HD13	1:A:390:PHE:HD1	1.78	0.49
1:A:522:VAL:HG13	1:A:627:VAL:HG13	1.96	0.48
1:A:396:PRO:O	1:A:418:LEU:HD12	2.14	0.48
1:A:278:ALA:HB3	1:A:284:LYS:HG3	1.95	0.47
1:A:172:LEU:HA	1:A:180:THR:O	2.15	0.46
1:A:489:LEU:HD11	1:A:633:HIS:HA	1.96	0.46
1:A:432:HIS:ND1	1:A:433:PRO:HD3	2.29	0.46
1:A:620:LYS:HG2	4:A:789:HOH:O	2.16	0.44
1:A:504:TRP:CE3	1:A:638:LYS:HG3	2.52	0.44
1:A:366:LEU:HD13	1:A:368:TRP:CH2	2.53	0.44
1:A:428:SER:CB	1:A:558:ARG:HH22	2.30	0.44
1:A:107:THR:HG22	1:A:171:LEU:CD2	2.47	0.44
1:A:711:LEU:O	1:A:711:LEU:HD23	2.17	0.43
1:A:101:PHE:CZ	1:A:108:LYS:HB2	2.53	0.42
1:A:431:LYS:O	1:A:432:HIS:CD2	2.72	0.42
1:A:237:GLU:H	1:A:237:GLU:CD	2.22	0.42
1:A:404:TYR:O	1:A:408:ARG:HG3	2.19	0.42
1:A:522:VAL:CG1	1:A:631:LEU:HB2	2.50	0.42
1:A:309:TRP:CH2	1:A:401:ILE:HG13	2.54	0.42
1:A:559:ALA:HB1	1:A:582:ARG:HD3	2.01	0.41
1:A:594:THR:HG23	4:A:1069:HOH:O	2.20	0.41
1:A:22:PHE:N	1:A:23:PRO:CD	2.84	0.41
1:A:13:SER:OG	1:A:675:ARG:HD3	2.20	0.41
1:A:529:THR:HG22	1:A:549:ILE:HD13	2.02	0.41
1:A:128:ASN:HA	1:A:129:PRO:HD3	1.94	0.41
1:A:137:SER:CB	1:A:143:HIS:HE1	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:LYS:HD2	1:A:537:PHE:HA	2.02	0.41
1:A:325:GLU:HB3	1:A:327:TYR:CE1	2.56	0.41
1:A:56:GLU:HB3	1:A:57:PRO:HD3	2.03	0.41
1:A:208:LEU:HA	1:A:209:PRO:HD3	1.69	0.41
1:A:460:HIS:CD2	1:A:460:HIS:C	2.94	0.40
1:A:385:LEU:HD12	1:A:416:VAL:HG11	2.04	0.40
1:A:395:ILE:HA	1:A:396:PRO:HD2	1.92	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	711/711 (100%)	691 (97%)	17 (2%)	3 (0%)	36 24

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	432	HIS
1	A	76	ASP
1	A	433	PRO

### 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	633/631 (100%)	628 (99%)	5 (1%)	83	82

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	GLU
1	A	237	GLU
1	A	378	ASN
1	A	415	ASN
1	A	689	LYS

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

Mol	Chain	Res	Type
1	A	378	ASN
1	A	688	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 [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	710/711 (99%)	0.50	76 (10%) 6 10	24, 45, 74, 116	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	78	ILE	6.8
1	A	321	ILE	5.8
1	A	432	HIS	5.6
1	A	46	VAL	5.1
1	A	409	LEU	5.0
1	A	337	PHE	4.9
1	A	585	ILE	4.5
1	A	570	LYS	4.5
1	A	323	PHE	4.3
1	A	47	VAL	4.1
1	A	431	LYS	4.1
1	A	683	TYR	4.0
1	A	523	ILE	3.9
1	A	98	LEU	3.8
1	A	412	GLY	3.7
1	A	428	SER	3.7
1	A	391	THR	3.6
1	A	219	GLU	3.6
1	A	687	ASN	3.5
1	A	44	SER	3.5
1	A	522	VAL	3.4
1	A	631	LEU	3.4
1	A	94	PHE	3.4
1	A	95	LEU	3.3
1	A	413	PHE	3.2
1	A	48	MET	3.2
1	A	2	SER	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	556	MET	3.1
1	A	59	PHE	3.0
1	A	594	THR	2.9
1	A	411	ILE	2.9
1	A	410	LYS	2.9
1	A	227	ALA	2.8
1	A	711	LEU	2.8
1	A	322	GLY	2.8
1	A	390	PHE	2.7
1	A	510	GLN	2.7
1	A	160	ILE	2.7
1	A	689	LYS	2.7
1	A	525	MET	2.7
1	A	581	ALA	2.6
1	A	99	GLY	2.6
1	A	339	SER	2.6
1	A	335	GLY	2.6
1	A	634	LEU	2.6
1	A	51	VAL	2.6
1	A	179	TYR	2.6
1	A	43	GLY	2.6
1	A	433	PRO	2.6
1	A	557	ALA	2.5
1	A	691	THR	2.5
1	A	340	LEU	2.5
1	A	130	SER	2.5
1	A	187	LEU	2.4
1	A	630	TYR	2.4
1	A	249	ALA	2.4
1	A	524	ALA	2.4
1	A	267	VAL	2.4
1	A	424	ALA	2.4
1	A	226	VAL	2.3
1	A	520	ALA	2.3
1	A	40	SER	2.3
1	A	324	ILE	2.3
1	A	686	ASP	2.3
1	A	568	ASN	2.3
1	A	341	VAL	2.2
1	A	58	ILE	2.2
1	A	50	GLN	2.2
1	A	175	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	3	HIS	2.1
1	A	77	ASP	2.1
1	A	588	THR	2.1
1	A	104	PHE	2.1
1	A	429	SER	2.1
1	A	575	GLY	2.1
1	A	250	VAL	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<sup>th</sup> 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	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	MG	A	714	1/1	0.94	0.10	66,66,66,66	0
3	MG	A	713	1/1	0.98	0.05	49,49,49,49	0
2	ZN	A	712	1/1	1.00	0.10	32,32,32,32	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.