



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2018 – 05:25 pm GMT

PDB ID : 5E3C
Title : Structure of human DPP3 in complex with hemorphin like opioid peptide IVYPW
Authors : Kumar, P.; Reithofer, V.; Reisinger, M.; Pavkov-Keller, T.; Wallner, S.; Machereux, P.; Gruber, K.
Deposited on : 2015-10-02
Resolution : 2.77 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)	:	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

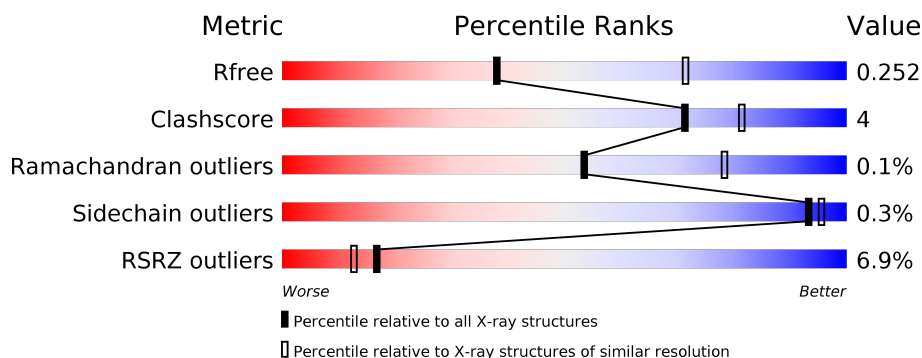
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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	4013 (2.80-2.72)
Clashscore	122126	1029 (2.78-2.74)
Ramachandran outliers	120053	1013 (2.78-2.74)
Sidechain outliers	120020	1013 (2.78-2.74)
RSRZ outliers	108989	3920 (2.80-2.72)

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. 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	726	<div> <div>7%</div> <div> <div></div> <div>90%</div> <div>10%</div> </div> </div>
2	B	5	<div> <div>20%</div> <div> <div></div> <div>60%</div> <div>40%</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5835 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 Dipeptidyl peptidase 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	724	Total	C	N	O	S	0	1	0
			5753	3667	974	1100	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	SER	CYS	engineered mutation	UNP Q9NY33
A	207	CYS	GLU	engineered mutation	UNP Q9NY33
A	451	ALA	GLU	engineered mutation	UNP Q9NY33
A	491	CYS	SER	engineered mutation	UNP Q9NY33
A	519	SER	CYS	engineered mutation	UNP Q9NY33
A	654	SER	CYS	engineered mutation	UNP Q9NY33

- Molecule 2 is a protein called IVYPW.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	5	Total	C	N	O	0	0	0
			49	36	6	7			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	K 1	0	0

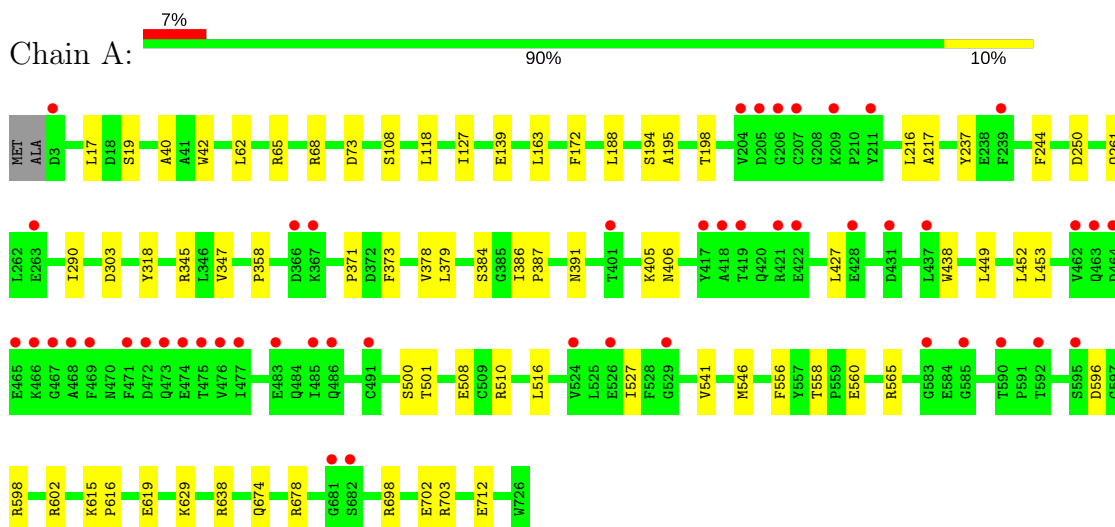
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	29	Total 29	O 29	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 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: Dipeptidyl peptidase 3



• Molecule 2: IVYPW



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.67Å 106.39Å 65.12Å 90.00° 93.44° 90.00°	Depositor
Resolution (Å)	45.56 – 2.77 45.56 – 2.76	Depositor EDS
% Data completeness (in resolution range)	97.8 (45.56-2.77) 97.9 (45.56-2.76)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	0.21	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.22 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.204 , 0.253 0.207 , 0.252	Depositor DCC
R_{free} test set	1033 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	57.5	Xtriage
Anisotropy	0.128	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 33.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5835	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: K, MG, ZN

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	0.24	0/5888	0.36	0/7976
2	B	0.21	0/52	0.40	0/71
All	All	0.24	0/5940	0.36	0/8047

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 [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	5753	0	5630	41	0
2	B	49	0	48	9	0
3	A	1	0	0	0	0
4	A	2	0	0	0	0
5	A	1	0	0	0	0
6	A	29	0	0	0	0
All	All	5835	0	5678	44	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 (44) 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)
2:B:1:ILE:HG22	2:B:2:VAL:H	1.19	1.00
2:B:1:ILE:HG22	2:B:2:VAL:N	1.75	1.00
1:A:516:LEU:HD11	1:A:546:MET:HG2	1.71	0.71
1:A:68[A]:ARG:NH1	1:A:712:GLU:OE2	2.25	0.69
2:B:1:ILE:CG2	2:B:2:VAL:N	2.50	0.65
1:A:387:PRO:HB3	2:B:2:VAL:HG13	1.81	0.63
1:A:391:ASN:O	2:B:1:ILE:N	2.35	0.59
1:A:438:TRP:HB2	1:A:541:VAL:HG21	1.83	0.59
1:A:508:GLU:OE1	2:B:1:ILE:HG21	2.03	0.58
1:A:65:ARG:NH2	1:A:139:GLU:OE2	2.37	0.57
1:A:73:ASP:OD2	1:A:698:ARG:NH2	2.31	0.56
1:A:318:TYR:OH	2:B:1:ILE:CG2	2.55	0.55
1:A:358:PRO:HB3	1:A:619:GLU:HG3	1.90	0.53
1:A:198:THR:HG22	1:A:216:LEU:HD23	1.91	0.53
1:A:558:THR:HG22	1:A:560:GLU:HG2	1.90	0.53
1:A:19:SER:HB3	1:A:379:LEU:HD23	1.91	0.53
1:A:678:ARG:NH2	1:A:702:GLU:O	2.42	0.52
1:A:427:LEU:O	1:A:602:ARG:NH1	2.45	0.49
1:A:449:LEU:HD22	1:A:453:LEU:HD12	1.94	0.49
1:A:345:ARG:HB2	1:A:527:ILE:HG23	1.95	0.49
1:A:108:SER:HB3	1:A:384:SER:HB2	1.95	0.48
1:A:194:SER:HB3	1:A:290:ILE:HG13	1.95	0.47
1:A:501:THR:HG1	1:A:638:ARG:HH22	1.60	0.47
1:A:596:ASP:OD2	1:A:598:ARG:NE	2.46	0.46
1:A:217:ALA:HA	1:A:250:ASP:HB3	1.97	0.45
1:A:188:LEU:HD13	1:A:195:ALA:HB2	1.99	0.45
1:A:347:VAL:HG13	1:A:371:PRO:HD2	1.99	0.45
1:A:371:PRO:HG2	1:A:452:LEU:HD11	1.98	0.44
1:A:386:ILE:HD11	1:A:674:GLN:NE2	2.33	0.44
1:A:62:LEU:HD21	1:A:127:ILE:HD13	2.00	0.44
1:A:42:TRP:CD2	1:A:703:ARG:HD3	2.52	0.43
1:A:345:ARG:NH1	1:A:527:ILE:O	2.51	0.43
1:A:391:ASN:H	2:B:1:ILE:H2	1.68	0.42
1:A:391:ASN:ND2	1:A:406:ASN:OD1	2.53	0.41
1:A:237:TYR:N	1:A:244:PHE:O	2.49	0.41
1:A:556:PHE:HA	1:A:565:ARG:NH2	2.36	0.41
1:A:303:ASP:O	1:A:405:LYS:NZ	2.44	0.41
1:A:510:ARG:HH21	1:A:629:LYS:HE2	1.85	0.41
1:A:40:ALA:HB2	1:A:261:GLN:HB2	2.02	0.41
1:A:17:LEU:HD21	1:A:378:VAL:HB	2.02	0.41
1:A:615:LYS:HB3	1:A:616:PRO:HD3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:LEU:HD21	1:A:172:PHE:CD1	2.56	0.40
1:A:65:ARG:HA	1:A:68[B]:ARG:NH2	2.37	0.40
1:A:387:PRO:CB	2:B:2:VAL:HG13	2.51	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	723/726 (100%)	701 (97%)	21 (3%)	1 (0%)	53	78
2	B	3/5 (60%)	3 (100%)	0	0	100	100
All	All	726/731 (99%)	704 (97%)	21 (3%)	1 (0%)	53	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	500	SER

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	609/609 (100%)	607 (100%)	2 (0%)	93	95
2	B	5/5 (100%)	5 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	614/614 (100%)	612 (100%)	2 (0%)	93	95

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	118	LEU
1	A	373	PHE

Some 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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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 ⓘ

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, 95th 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 ²)	Q<0.9
1	A	724/726 (99%)	0.31	49 (6%) 17 13	28, 53, 87, 112	0
2	B	5/5 (100%)	1.56	1 (20%) 1 0	65, 67, 70, 73	0
All	All	729/731 (99%)	0.32	50 (6%) 17 13	28, 53, 87, 112	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	207	CYS	6.7
1	A	465	GLU	4.6
1	A	467	GLY	4.4
1	A	206	GLY	4.1
1	A	595	SER	4.1
1	A	475	THR	4.0
1	A	681	GLY	4.0
1	A	417	TYR	3.9
1	A	491	CYS	3.7
1	A	418	ALA	3.7
1	A	474	GLU	3.7
1	A	366	ASP	3.5
1	A	466	LYS	3.4
1	A	472	ASP	3.4
1	A	485	ILE	3.3
1	A	468	ALA	3.2
1	A	401	THR	3.1
1	A	529	GLY	3.1
1	A	469	PHE	3.0
1	A	477	ILE	3.0
1	A	471	PHE	3.0
1	A	585	GLY	2.9
1	A	211	TYR	2.9
1	A	464	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	583	GLY	2.8
1	A	462	VAL	2.7
1	A	419	THR	2.7
1	A	422	GLU	2.6
1	A	437	LEU	2.6
1	A	209	LYS	2.5
1	A	524	VAL	2.5
1	A	421	ARG	2.5
1	A	486	GLN	2.4
1	A	590	THR	2.4
1	A	3	ASP	2.4
2	B	4	PRO	2.4
1	A	205	ASP	2.4
1	A	592	THR	2.3
1	A	431	ASP	2.3
1	A	473	GLN	2.3
1	A	476	VAL	2.2
1	A	263	GLU	2.2
1	A	682	SER	2.1
1	A	367	LYS	2.1
1	A	239	PHE	2.1
1	A	526	GLU	2.1
1	A	483	GLU	2.1
1	A	428	GLU	2.1
1	A	463	GLN	2.0
1	A	204	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, 95th 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(\AA^2)	Q<0.9
3	ZN	A	801	1/1	0.90	0.20	70,70,70,70	1
4	MG	A	804	1/1	0.93	0.13	38,38,38,38	0
4	MG	A	802	1/1	0.95	0.11	46,46,46,46	0
5	K	A	803	1/1	0.97	0.10	57,57,57,57	0

6.5 Other polymers [i](#)

There are no such residues in this entry.