Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 08:38 PM EDT

PDB ID	:	3D3M
Title	:	The Crystal Structure of the C-terminal region of Death Associated Protein 5(DAP5)
Authors	:	Dym, O.; Israel Structural Proteomics Center (ISPC)
Deposited on	:	2008-05-12
Resolution	:	1.90 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 (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:

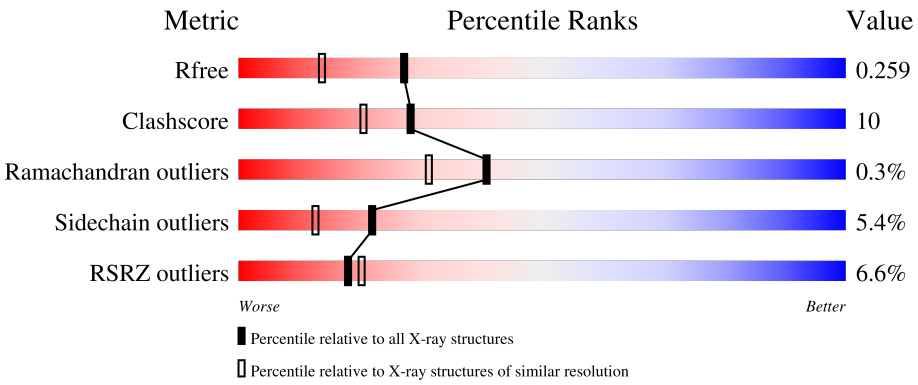
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	A	168	3% 77% 15% • •
1	B	168	10% 69% 20% • 7%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2714 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 Eukaryotic translation initiation factor 4 gamma 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	161	1337	881	212	239	5	0	0	0
1	B	157	1312	866	208	233	5	0	0	0

- Molecule 2 is water.

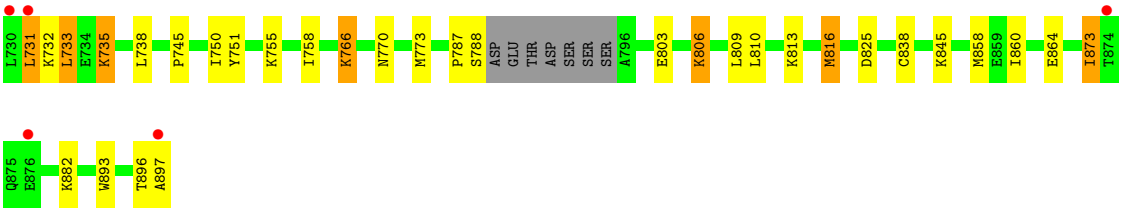
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	37	37	37	0	0
2	B	28	28	28	0	0

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.

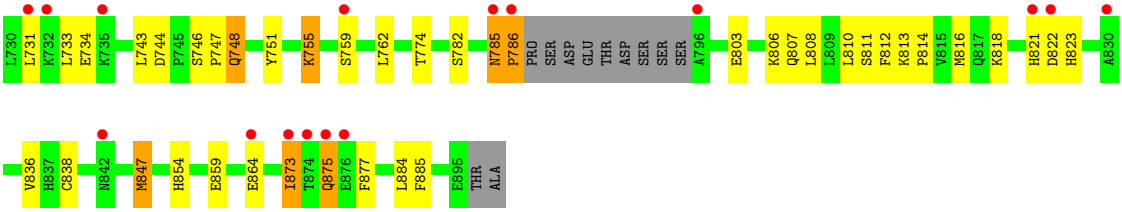
• Molecule 1: Eukaryotic translation initiation factor 4 gamma 2

Chain A:



• Molecule 1: Eukaryotic translation initiation factor 4 gamma 2

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	61.06Å 45.73Å 63.36Å 90.00° 103.81° 90.00°	Depositor
Resolution (Å)	25.00 – 1.90 24.98 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (25.00-1.90) 100.0 (24.98-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	4.90 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.4.0067	Depositor
R, R_{free}	0.211 , 0.259 0.211 , 0.259	Depositor DCC
R_{free} test set	1357 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	26.6	Xtriage
Anisotropy	0.333	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 48.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.022 for l,-k,h	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	2714	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.85% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	1.36	5/1373 (0.4%)	1.14	8/1857 (0.4%)
1	B	1.11	0/1347	0.97	5/1820 (0.3%)
All	All	1.24	5/2720 (0.2%)	1.06	13/3677 (0.4%)

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	1	0
1	B	0	1
All	All	1	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	806	LYS	CE-NZ	8.19	1.69	1.49
1	A	838	CYS	CB-SG	7.36	1.94	1.82
1	A	806	LYS	CD-CE	6.29	1.67	1.51
1	A	803	GLU	CG-CD	5.79	1.60	1.51
1	A	864	GLU	CG-CD	5.39	1.60	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	896	THR	C-N-CA	-11.52	92.91	121.70
1	A	766	LYS	CD-CE-NZ	9.30	133.10	111.70
1	A	816	MET	CG-SD-CE	7.39	112.02	100.20
1	A	738	LEU	CB-CG-CD2	-6.78	99.48	111.00
1	A	773	MET	CG-SD-CE	-6.66	89.54	100.20
1	A	806	LYS	CD-CE-NZ	6.64	126.96	111.70
1	B	785	ASN	CB-CA-C	-6.43	97.54	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	808	LEU	CB-CG-CD2	-5.84	101.08	111.00
1	A	896	THR	O-C-N	-5.44	113.99	122.70
1	B	744	ASP	CB-CG-OD1	5.39	123.16	118.30
1	A	897	ALA	N-CA-C	5.27	125.22	111.00
1	B	884	LEU	CB-CG-CD2	-5.23	102.11	111.00
1	B	847	MET	CG-SD-CE	5.06	108.30	100.20

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	897	ALA	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	785	ASN	Peptide

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	1337	0	1345	26	0
1	B	1312	0	1321	27	0
2	A	37	0	0	1	0
2	B	28	0	0	2	0
All	All	2714	0	2666	52	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 10.

All (52) 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:806:LYS:CE	1:A:806:LYS:NZ	1.69	1.50
1:A:858:MET:CE	1:A:860:ILE:CD1	2.31	1.09
1:A:766:LYS:NZ	1:A:825:ASP:OD1	1.86	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:858:MET:HE3	1:A:860:ILE:CD1	1.85	1.06
1:A:858:MET:CE	1:A:860:ILE:HD12	1.86	1.05
1:A:858:MET:HE3	1:A:860:ILE:HD11	1.38	0.99
1:A:858:MET:HE2	1:A:860:ILE:HD12	1.53	0.90
1:A:858:MET:CE	1:A:860:ILE:HD11	2.00	0.89
1:B:734:GLU:HG2	2:B:34:HOH:O	1.72	0.89
1:A:751:TYR:CE2	1:A:755:LYS:HD3	2.21	0.75
1:B:782:SER:O	1:B:786:PRO:HD2	1.87	0.75
1:A:893:TRP:C	1:A:893:TRP:CD1	2.61	0.74
1:B:746:SER:OG	1:B:748:GLN:OE1	2.13	0.64
1:B:873:ILE:O	1:B:875:GLN:O	2.15	0.63
1:B:747:PRO:HG3	1:B:811:SER:O	1.98	0.63
1:A:733:LEU:HD11	1:A:758:ILE:HD11	1.82	0.61
1:B:748:GLN:OE1	1:B:748:GLN:N	2.32	0.61
1:A:732:LYS:O	1:A:735:LYS:HB2	2.00	0.61
1:B:748:GLN:H	1:B:748:GLN:CD	2.03	0.60
1:B:751:TYR:CE2	1:B:755:LYS:HG3	2.39	0.57
1:B:734:GLU:OE2	1:B:774:THR:HG21	2.08	0.54
1:A:813:LYS:HA	1:A:816:MET:HE3	1.89	0.54
1:B:821:HIS:O	1:B:822:ASP:HB2	2.08	0.53
1:A:732:LYS:O	1:A:735:LYS:N	2.39	0.53
1:B:838:CYS:HA	1:B:847:MET:CE	2.39	0.52
1:A:770:ASN:ND2	2:A:108:HOH:O	2.44	0.51
1:A:751:TYR:CE2	1:A:755:LYS:CD	2.94	0.51
1:B:806:LYS:HE2	1:B:854:HIS:HE1	1.77	0.50
1:B:859:GLU:HA	2:B:40:HOH:O	2.11	0.50
1:B:838:CYS:SG	1:B:847:MET:CE	3.00	0.50
1:B:838:CYS:SG	1:B:847:MET:HE3	2.52	0.49
1:B:751:TYR:CE1	1:B:818:LYS:HG2	2.48	0.49
1:B:813:LYS:N	1:B:814:PRO:CD	2.75	0.49
1:B:821:HIS:O	1:B:823:HIS:HD2	1.97	0.47
1:B:821:HIS:O	1:B:823:HIS:CD2	2.67	0.46
1:A:893:TRP:CD1	1:B:885:PHE:CD1	3.04	0.46
1:B:864:GLU:H	1:B:864:GLU:CD	2.20	0.45
1:A:735:LYS:HD3	1:A:735:LYS:HA	1.34	0.44
1:A:809:LEU:HG	1:A:858:MET:HE1	2.00	0.44
1:B:813:LYS:N	1:B:814:PRO:HD2	2.31	0.44
1:A:882:LYS:HE2	1:A:882:LYS:HB3	1.81	0.43
1:B:759:SER:HB3	1:B:762:LEU:HD12	2.01	0.43
1:A:735:LYS:HZ2	1:A:735:LYS:HG2	1.69	0.43
1:B:803:GLU:OE2	1:B:807:GLN:NE2	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:812:PHE:C	1:B:816:MET:HE2	2.39	0.42
1:A:745:PRO:HA	1:A:750:ILE:HD11	2.02	0.41
1:A:755:LYS:HB2	1:A:755:LYS:NZ	2.34	0.41
1:B:836:VAL:CG2	1:B:877:PHE:HB3	2.50	0.41
1:A:731:LEU:HD22	1:A:735:LYS:NZ	2.35	0.41
1:A:873:ILE:O	1:A:873:ILE:HG13	2.09	0.41
1:A:733:LEU:HD22	1:A:733:LEU:HA	1.94	0.40
1:B:838:CYS:SG	1:B:847:MET:HE2	2.61	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	157/168 (94%)	153 (98%)	3 (2%)	1 (1%)	25	15
1	B	153/168 (91%)	151 (99%)	2 (1%)	0	100	100
All	All	310/336 (92%)	304 (98%)	5 (2%)	1 (0%)	41	31

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	787	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	150/157 (96%)	143 (95%)	7 (5%)	26	16
1	B	147/157 (94%)	138 (94%)	9 (6%)	18	9
All	All	297/314 (95%)	281 (95%)	16 (5%)	22	13

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	731	LEU
1	A	733	LEU
1	A	735	LYS
1	A	788	SER
1	A	810	LEU
1	A	845	LYS
1	A	873	ILE
1	B	731	LEU
1	B	733	LEU
1	B	743	LEU
1	B	748	GLN
1	B	755	LYS
1	B	786	PRO
1	B	810	LEU
1	B	873	ILE
1	B	875	GLN

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

Mol	Chain	Res	Type
1	A	823	HIS
1	B	823	HIS

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, 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	161/168 (95%)	0.06	5 (3%) 49 51	14, 25, 46, 55	0
1	B	157/168 (93%)	0.55	16 (10%) 6 8	18, 31, 52, 66	0
All	All	318/336 (94%)	0.30	21 (6%) 18 20	14, 28, 50, 66	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	874	THR	10.3
1	A	731	LEU	6.3
1	B	786	PRO	6.0
1	B	873	ILE	5.6
1	A	730	LEU	4.3
1	B	875	GLN	4.2
1	B	731	LEU	3.9
1	B	759	SER	3.3
1	A	876	GLU	3.3
1	B	732	LYS	3.1
1	B	822	ASP	2.9
1	B	821	HIS	2.8
1	B	796	ALA	2.7
1	B	785	ASN	2.7
1	B	876	GLU	2.5
1	B	830	ALA	2.5
1	B	842	ASN	2.3
1	A	874	THR	2.1
1	B	735	LYS	2.1
1	B	864	GLU	2.0
1	A	897	ALA	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 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.