



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 04:00 PM EDT

PDB ID : 6QDK / pdb_00006qdk
Title : Molecular features of the UNC-45 chaperone critical for binding and folding muscle myosin
Authors : Meinhart, A.; Clausen, T.; Hellerschmied, D.
Deposited on : 2019-01-02
Resolution : 3.40 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

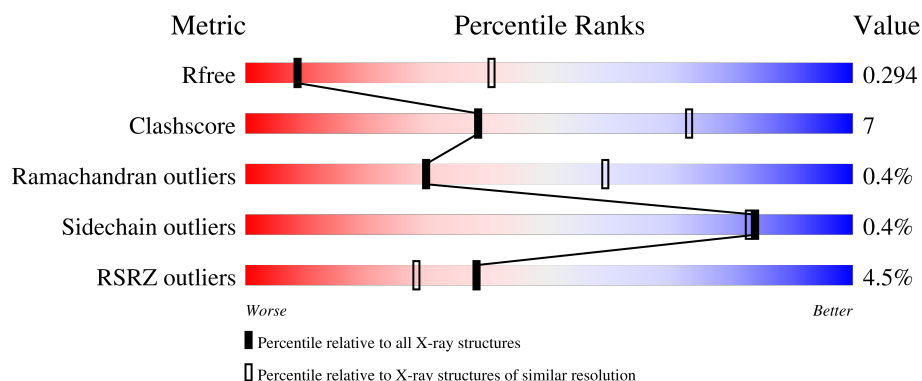
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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}	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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 $\leq 5\%$. 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	964	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6848 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 UNC-45,UNC-45.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	882	Total	C	N	O	S	Se	0	0	0
			6848	4318	1183	1303	19	25			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	initiating methionine	UNP G5EG62
A	427	GLU	GLY	engineered mutation	UNP G5EG62



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	86.91Å 86.91Å 718.54Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.96 – 3.40 29.96 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.96-3.40) 99.5 (29.96-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 3.39Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???), CNS	Depositor
R, R_{free}	0.287 , 0.319 0.289 , 0.294	Depositor DCC
R_{free} test set	1217 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	98.9	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 19.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6848	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% 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	0.37	0/6835	0.66	1/9181 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	890	MSE	N-CA-C	6.29	124.19	110.80

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	6848	0	6883	94	0
All	All	6848	0	6883	94	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 7.

All (94) 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:853:CYS:SG	1:A:889:ILE:HG23	1.59	1.41
1:A:847:THR:HG21	1:A:889:ILE:CG1	1.52	1.39

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:847:THR:CG2	1:A:889:ILE:CG1	2.31	1.08
1:A:847:THR:CG2	1:A:889:ILE:HG12	1.89	1.02
1:A:101:PHE:CZ	1:A:105:LYS:HD3	1.96	1.01
1:A:853:CYS:SG	1:A:889:ILE:CG2	2.52	0.98
1:A:847:THR:CG2	1:A:889:ILE:HG13	1.91	0.97
1:A:847:THR:HG21	1:A:889:ILE:HG13	0.95	0.93
1:A:381:VAL:CG2	1:A:382:PHE:N	2.33	0.91
1:A:847:THR:HG23	1:A:889:ILE:HG12	1.60	0.83
1:A:856:ILE:HG21	1:A:889:ILE:HD13	1.61	0.83
1:A:101:PHE:CE2	1:A:105:LYS:HD3	2.14	0.82
1:A:381:VAL:HG23	1:A:382:PHE:N	1.94	0.81
1:A:381:VAL:HG22	1:A:382:PHE:H	1.45	0.81
1:A:879:GLN:O	1:A:883:LEU:HG	1.82	0.79
1:A:856:ILE:HD11	1:A:896:LEU:HD22	1.66	0.78
1:A:376:LEU:HD12	1:A:376:LEU:O	1.84	0.78
1:A:856:ILE:HD13	1:A:889:ILE:HG21	1.65	0.78
1:A:381:VAL:CG2	1:A:382:PHE:H	1.99	0.74
1:A:767:ARG:HH12	1:A:804:PHE:H	1.37	0.72
1:A:101:PHE:O	1:A:105:LYS:HG3	1.90	0.71
1:A:856:ILE:CD1	1:A:896:LEU:HD22	2.27	0.64
1:A:844:ALA:HA	1:A:885:GLY:HA3	1.83	0.60
1:A:841:ALA:HB2	1:A:881:ARG:HD3	1.84	0.59
1:A:879:GLN:HG2	1:A:883:LEU:HD11	1.84	0.59
1:A:524:ALA:N	1:A:527:SER:HG	2.01	0.59
1:A:57:ARG:NH1	1:A:69:ASP:OD2	2.36	0.58
1:A:141:LYS:HG3	1:A:145:MSE:HE2	1.85	0.58
1:A:101:PHE:CE2	1:A:105:LYS:NZ	2.74	0.56
1:A:380:MSE:O	1:A:382:PHE:N	2.38	0.55
1:A:374:GLN:HG2	1:A:378:GLU:OE1	2.07	0.55
1:A:91:ARG:NH1	1:A:103:ASP:OD2	2.41	0.54
1:A:853:CYS:CB	1:A:889:ILE:HG23	2.37	0.54
1:A:303:ASP:O	1:A:311:ARG:NH2	2.41	0.54
1:A:101:PHE:CE2	1:A:105:LYS:CD	2.88	0.53
1:A:311:ARG:NH1	1:A:360:PRO:O	2.42	0.53
1:A:101:PHE:CE2	1:A:105:LYS:CE	2.92	0.52
1:A:12:ARG:HG2	1:A:49:LEU:HD21	1.91	0.52
1:A:779:ILE:HG23	1:A:799:LEU:HD11	1.92	0.51
1:A:700:SER:OG	1:A:701:GLY:N	2.44	0.50
1:A:156:ASP:HB3	1:A:159:GLN:HB2	1.94	0.50
1:A:397:PHE:HE2	1:A:438:ASN:HD22	1.60	0.50
1:A:856:ILE:HD11	1:A:896:LEU:CD2	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:LEU:HD23	1:A:163:ALA:HB2	1.94	0.49
1:A:543:SER:H	1:A:546:ILE:HD12	1.76	0.49
1:A:498:VAL:HG21	1:A:546:ILE:HG23	1.95	0.49
1:A:745:GLU:O	1:A:749:ASN:ND2	2.45	0.49
1:A:730:TYR:HB3	1:A:769:ARG:HG2	1.95	0.48
1:A:145:MSE:HE1	1:A:170:LEU:HD13	1.95	0.48
1:A:709:HIS:CE1	1:A:713:LYS:HE3	2.49	0.48
1:A:640:ARG:HD2	1:A:679:LEU:HD11	1.96	0.47
1:A:856:ILE:CD1	1:A:896:LEU:CD2	2.92	0.47
1:A:64:GLU:OE2	1:A:218:ASN:ND2	2.48	0.47
1:A:32:THR:OG1	1:A:53:ARG:NH2	2.48	0.46
1:A:263:VAL:HG12	1:A:267:MSE:HE2	1.96	0.46
1:A:844:ALA:HA	1:A:885:GLY:CA	2.44	0.46
1:A:886:ILE:HD11	1:A:896:LEU:HD21	1.97	0.46
1:A:51:ARG:HD2	1:A:82:LYS:HB2	1.98	0.46
1:A:51:ARG:NH1	1:A:52:ASN:OD1	2.49	0.46
1:A:362:SER:OG	1:A:363:ALA:N	2.49	0.45
1:A:429:VAL:HG12	1:A:431:ILE:H	1.82	0.45
1:A:891:HIS:O	1:A:891:HIS:CG	2.70	0.44
1:A:87:ARG:HH21	1:A:91:ARG:HH22	1.64	0.44
1:A:774:LYS:HB3	1:A:777:PRO:HG2	2.00	0.44
1:A:101:PHE:CD2	1:A:105:LYS:HE2	2.53	0.44
1:A:590:TYR:O	1:A:594:THR:OG1	2.26	0.44
1:A:390:LYS:HG2	1:A:431:ILE:HD12	2.00	0.44
1:A:680:ARG:HD2	1:A:714:LEU:HA	1.99	0.44
1:A:16:ASN:OD1	1:A:31:TYR:OH	2.34	0.43
1:A:405:THR:OG1	1:A:407:ASP:OD1	2.34	0.43
1:A:271:ASP:H	1:A:276:MSE:HB3	1.83	0.43
1:A:651:CYS:O	1:A:655:SER:OG	2.35	0.43
1:A:767:ARG:NH1	1:A:804:PHE:H	2.10	0.43
1:A:767:ARG:HA	1:A:770:ILE:HG12	2.00	0.43
1:A:663:LEU:HD23	1:A:666:ILE:HD12	2.01	0.42
1:A:149:ALA:HB2	1:A:163:ALA:HB3	2.00	0.42
1:A:408:ASP:HA	1:A:411:HIS:CD2	2.54	0.42
1:A:42:ASP:HB3	1:A:45:LEU:HB3	2.01	0.42
1:A:857:MSE:HB3	1:A:863:TRP:HZ3	1.85	0.42
1:A:823:TRP:HD1	1:A:839:SER:HA	1.85	0.42
1:A:301:LEU:O	1:A:311:ARG:NE	2.48	0.41
1:A:87:ARG:HH21	1:A:91:ARG:NH2	2.18	0.41
1:A:709:HIS:CD2	1:A:748:ALA:HA	2.54	0.41
1:A:468:VAL:HG23	1:A:506:LYS:HD2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:632:GLU:HA	1:A:635:VAL:HG22	2.03	0.41
1:A:670:LEU:HA	1:A:673:PHE:HD2	1.85	0.41
1:A:853:CYS:SG	1:A:889:ILE:HG12	2.60	0.41
1:A:821:LYS:HE3	1:A:825:LEU:HD12	2.02	0.41
1:A:85:PHE:O	1:A:88:SER:OG	2.32	0.41
1:A:376:LEU:CD1	1:A:380:MSE:HE2	2.50	0.41
1:A:465:VAL:HG21	1:A:502:VAL:HG22	2.03	0.41
1:A:847:THR:O	1:A:888:ASN:HB3	2.21	0.41
1:A:736:LEU:HD23	1:A:736:LEU:HA	1.91	0.40
1:A:574:LEU:HD23	1:A:577:LEU:HD12	2.03	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	857/964 (89%)	809 (94%)	45 (5%)	3 (0%)	30 59

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	381	VAL
1	A	890	MSE
1	A	864	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	729/745 (98%)	726 (100%)	3 (0%)	84	83

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	378	GLU
1	A	381	VAL
1	A	468	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	GLN
1	A	102	GLN
1	A	128	ASN
1	A	196	ASN
1	A	265	ASN
1	A	302	GLN
1	A	411	HIS
1	A	888	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 oligosaccharides 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	912:VAL	C	918:UNK	N	10.06

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	840/964 (87%)	0.43	38 (4%) 38 28	74, 93, 123, 133	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	796	ALA	4.8
1	A	828	ALA	4.6
1	A	866	VAL	4.4
1	A	843	PHE	3.9
1	A	775	ALA	3.7
1	A	822	LEU	3.5
1	A	101	PHE	3.4
1	A	860	ILE	3.3
1	A	778	LYS	3.1
1	A	870	ILE	3.0
1	A	446	LEU	2.9
1	A	856	ILE	2.9
1	A	826	TYR	2.7
1	A	740	LEU	2.7
1	A	426	GLN	2.6
1	A	889	ILE	2.6
1	A	105	LYS	2.6
1	A	776	ILE	2.6
1	A	888	ASN	2.6
1	A	886	ILE	2.5
1	A	771	LEU	2.5
1	A	835	LEU	2.5
1	A	779	ILE	2.4
1	A	675	GLU	2.4
1	A	737	CYS	2.4
1	A	786	THR	2.4
1	A	418	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	799	LEU	2.3
1	A	823	TRP	2.3
1	A	827	SER	2.3
1	A	774	LYS	2.2
1	A	800	LEU	2.2
1	A	732	VAL	2.1
1	A	803	LEU	2.1
1	A	838	ALA	2.1
1	A	472	GLU	2.1
1	A	504	LEU	2.0
1	A	121	LEU	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 oligosaccharides 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.