



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 06:06 PM UTC

PDB ID : 3NHE / pdb_00003nhe
Title : High Resolution Structure (1.26Å) of USP2a in Complex with Ubiquitin
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Deposited on : 2010-06-14
Resolution : 1.26 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

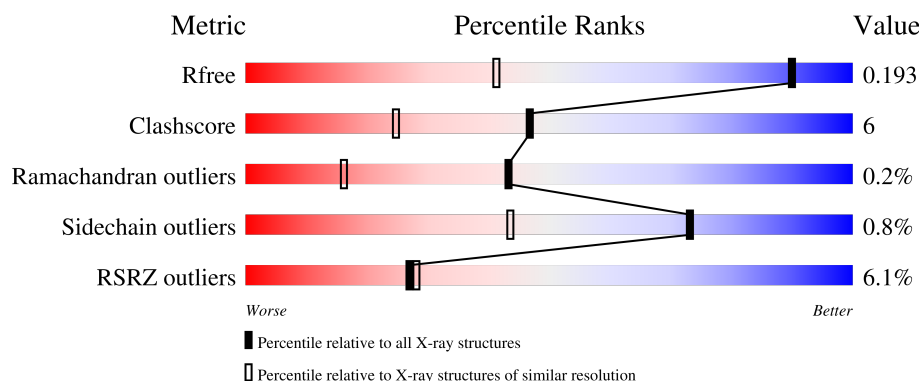
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.26 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	1693 (1.28-1.24)
Clashscore	190562	1730 (1.28-1.24)
Ramachandran outliers	187476	1695 (1.28-1.24)
Sidechain outliers	187428	1694 (1.28-1.24)
RSRZ outliers	180081	1693 (1.28-1.24)

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	348	7% 87% 9% ..
2	B	76	% 93% 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	CME	A	276	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3950 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 Ubiquitin carboxyl-terminal hydrolase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	5	17	0
			2820	1772	504	521	23			

- Molecule 2 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	76	Total	C	N	O	S	11	3	0
			621	389	110	121	1			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

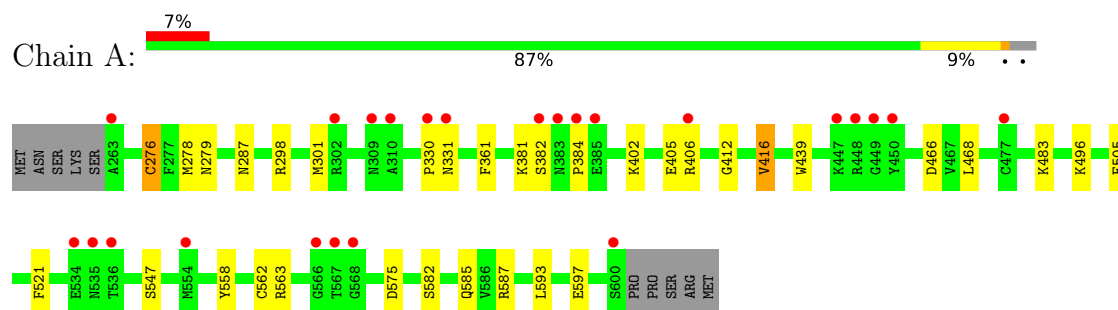
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	402	Total	O	0	0
			402	402		
4	B	106	Total	O	0	0
			106	106		

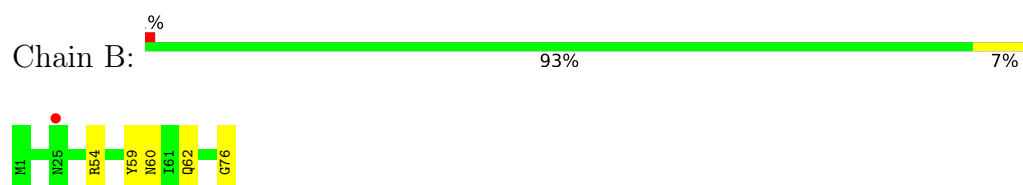
3 Residue-property plots [i](#)

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: Ubiquitin carboxyl-terminal hydrolase 2



- Molecule 2: Ubiquitin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.77 Å 53.96 Å 74.72 Å 90.00° 107.71° 90.00°	Depositor
Resolution (Å)	19.18 – 1.26 19.18 – 1.26	Depositor EDS
% Data completeness (in resolution range)	98.1 (19.18-1.26) 98.1 (19.18-1.26)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.73 (at 1.26 Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.169 , 0.189 0.171 , 0.193	Depositor DCC
R_{free} test set	5146 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	11.9	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 48.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3950	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.92% 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: CME, 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	0.37	0/2915	0.66	1/3934 (0.0%)
2	B	0.39	0/633	0.66	0/849
All	All	0.38	0/3548	0.66	1/4783 (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	384	PRO	N-CA-CB	8.18	110.59	103.31

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	2820	0	2791	37	0
2	B	621	0	652	4	0
3	A	1	0	0	0	0
4	A	402	0	0	2	0
4	B	106	0	0	1	0
All	All	3950	0	3443	39	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 6.

All (39) 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:287:ASN:HB2	1:A:562[B]:CYS:SG	2.00	1.02
1:A:276:CME:HE3	1:A:575:ASP:HB3	1.50	0.91
1:A:287:ASN:CB	1:A:562[B]:CYS:SG	2.59	0.90
1:A:276:CME:CE	1:A:575:ASP:HB3	2.10	0.80
1:A:276:CME:HZ3	1:A:575:ASP:OD1	1.82	0.80
1:A:298:ARG:HB3	1:A:301:MET:HE2	1.67	0.77
1:A:412:GLY:HA2	1:A:416[A]:VAL:HG23	1.73	0.70
1:A:582[B]:SER:H	1:A:585:GLN:HE21	1.42	0.68
1:A:582[A]:SER:H	1:A:585:GLN:HE21	1.43	0.67
1:A:416[B]:VAL:HG12	1:A:439:TRP:CE2	2.37	0.59
1:A:521:PHE:O	1:A:587[B]:ARG:HD3	2.04	0.58
1:A:276:CME:SG	1:A:558:TYR:O	2.63	0.57
1:A:276:CME:HB3	1:A:276:CME:CZ	2.34	0.57
1:A:496:LYS:HE2	4:A:732:HOH:O	2.05	0.56
1:A:412:GLY:O	1:A:416[B]:VAL:HG22	2.06	0.56
1:A:582[B]:SER:H	1:A:585:GLN:NE2	2.06	0.53
2:B:62:GLN:HB3	4:B:508:HOH:O	2.10	0.52
1:A:483:LYS:NZ	4:A:657:HOH:O	2.42	0.52
1:A:582[A]:SER:H	1:A:585:GLN:NE2	2.07	0.51
1:A:563[B]:ARG:NH2	1:A:597:GLU:OE2	2.45	0.50
1:A:402:LYS:O	1:A:405[A]:GLU:HG2	2.12	0.49
1:A:278[B]:MET:HG3	1:A:361:PHE:CZ	2.48	0.48
1:A:276:CME:SD	1:A:279:ASN:HB2	2.55	0.47
1:A:416[B]:VAL:CG1	1:A:439:TRP:CE2	2.97	0.47
1:A:412:GLY:CA	1:A:416[A]:VAL:HG23	2.45	0.44
1:A:406:ARG:HH22	2:B:60[B]:ASN:HD22	1.64	0.44
1:A:412:GLY:HA2	1:A:416[A]:VAL:CG2	2.43	0.44
1:A:287:ASN:HB3	1:A:562[B]:CYS:SG	2.54	0.44
1:A:330:PRO:O	1:A:331:ASN:HB2	2.18	0.44
1:A:276:CME:CZ	1:A:575:ASP:OD1	2.61	0.43
1:A:276:CME:HB2	2:B:76:GLY:O	2.19	0.42
1:A:416[A]:VAL:HG22	1:A:439:TRP:CE2	2.54	0.42
2:B:54:ARG:HD2	2:B:59:TYR:OH	2.20	0.42
1:A:278[A]:MET:HG3	1:A:361:PHE:CZ	2.55	0.41
1:A:412:GLY:HA2	1:A:416[B]:VAL:HG13	2.02	0.41
1:A:547:SER:HB3	1:A:593:LEU:HB2	2.03	0.41
1:A:416[B]:VAL:HG12	1:A:439:TRP:CD2	2.55	0.41
1:A:276:CME:CA	1:A:276:CME:HE2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:466:ASP:O	1:A:468:LEU:CD1	2.69	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	352/348 (101%)	344 (98%)	7 (2%)	1 (0%)	36	14
2	B	77/76 (101%)	75 (97%)	2 (3%)	0	100	100
All	All	429/424 (101%)	419 (98%)	9 (2%)	1 (0%)	43	16

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	382	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	319/315 (101%)	315 (99%)	4 (1%)	61	24
2	B	71/68 (104%)	71 (100%)	0	100	100
All	All	390/383 (102%)	386 (99%)	4 (1%)	73	34

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	381	LYS
1	A	416[A]	VAL
1	A	416[B]	VAL
1	A	505	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	306	HIS
1	A	309	ASN
1	A	398	GLN
1	A	537	ASN
1	A	548	ASN
1	A	585	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	CME	A	276	1	8,9,10	0.79	0	6,9,11	1.41	1 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CME	A	276	1	-	3/5/8/10	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	276	CME	CA-CB-SG	-2.26	105.19	114.45

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	276	CME	CE-SD-SG-CB
1	A	276	CME	SD-CE-CZ-OH
1	A	276	CME	N-CA-CB-SG

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	276	CME	9	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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, 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	337/348 (96%)	0.35	24 (7%) 22 23	5, 12, 24, 33	17 (5%)
2	B	75/76 (98%)	-0.04	1 (1%) 75 75	7, 11, 16, 22	2 (2%)
All	All	412/424 (97%)	0.28	25 (6%) 27 28	5, 12, 23, 33	19 (4%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	567	THR	5.5
1	A	566	GLY	5.4
1	A	477	CYS	4.8
1	A	450	TYR	4.2
1	A	384	PRO	3.9
1	A	449	GLY	3.6
1	A	309	ASN	3.5
1	A	535	ASN	3.4
1	A	568	GLY	3.4
1	A	534	GLU	3.4
1	A	447	LYS	3.3
1	A	554[A]	MET	3.3
1	A	382	SER	2.8
1	A	406	ARG	2.7
1	A	302	ARG	2.6
1	A	310	ALA	2.6
1	A	600	SER	2.6
1	A	331	ASN	2.6
1	A	330	PRO	2.5
1	A	536	THR	2.4
1	A	385	GLU	2.4
2	B	25	ASN	2.4
1	A	383	ASN	2.4
1	A	263	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	448	ARG	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	CME	A	276	10/11	0.60	0.23	10,14,25,26	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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(Å ²)	Q<0.9
3	ZN	A	1	1/1	1.00	0.03	11,11,11,11	0

6.5 Other polymers [i](#)

There are no such residues in this entry.