

Full wwPDB X-ray Structure Validation Report (i)

Feb 26, 2024 – 06:20 PM JST

PDB ID	:	8XPN
Title	:	The Crystal Structure of USP8 from Biortus.
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Deposited on	:	2024-01-04
Resolution	:	2.10 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:

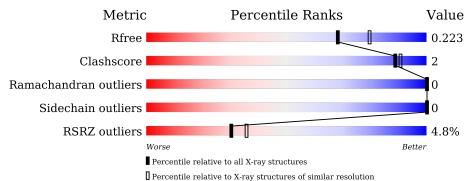
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\,DIFFRACTION$

The reported resolution of this entry is 2.10 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	$\begin{array}{l} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R _{free}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	А	377	87%		• 10%
1	В	377	90%		• 8%
1	С	377	8%	7%	13%



8XPN

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 8768 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Δ	341	Total	С	Ν	0	\mathbf{S}	0	1	0
	1 A	541	2749	1745	478	510	16	0		
1	D	347	Total	С	Ν	0	S	0	1	0
	I B	347	2827	1795	499	518	15			
1	1 C	327	Total	С	Ν	0	S	0	1	0
			2639	1684	456	485	14	0		U

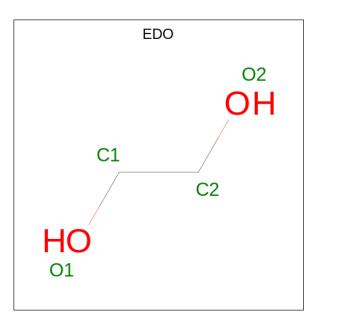
• Molecule 1 is a protein called Ubiquitin carboxyl-terminal hydrolase 8.

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	Total Zn 1 1	0	0
2	В	1	Total Zn 1 1	0	0
2	С	1	Total Zn 1 1	0	0

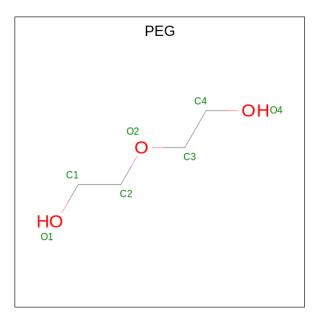
• Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
3	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0

• Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	С	1	Total 7	$\begin{array}{c} \mathrm{C} \\ 4 \end{array}$	O 3	0	0

• Molecule 5 is water.

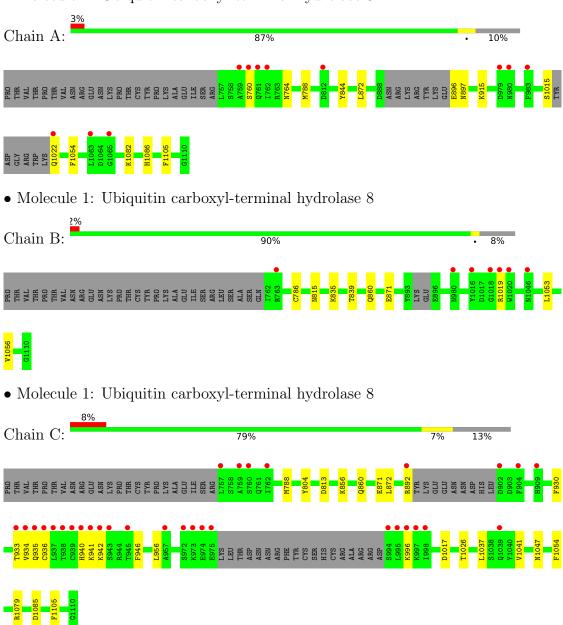


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	152	Total O 152 152	0	0
5	В	237	Total O 237 237	0	0
5	С	146	Total O 146 146	0	0



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: Ubiquitin carboxyl-terminal hydrolase 8



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	57.83Å 109.34Å 180.99Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.78 - 2.10	Depositor
Resolution (A)	48.73 - 2.10	EDS
% Data completeness	99.4 (48.78-2.10)	Depositor
(in resolution range)	99.4 (48.73-2.10)	EDS
R _{merge}	0.15	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$3.06 (at 2.10 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
D D.	0.176 , 0.216	Depositor
R, R_{free}	0.185 , 0.223	DCC
R_{free} test set	3285 reflections $(4.87%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	29.0	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 46.4	EDS
L-test for twinning ²	$ \langle L \rangle = 0.49, \langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8768	wwPDB-VP
Average B, all atoms $(Å^2)$	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.85% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: EDO, ZN, PEG

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		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.33	0/2812	0.63	0/3788	
1	В	0.36	0/2895	0.67	0/3899	
1	С	0.34	0/2701	0.66	1/3637~(0.0%)	
All	All	0.34	0/8408	0.65	1/11324~(0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	С	1085	ASP	CB-CA-C	5.84	122.07	110.40

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	А	2749	0	2695	9	0
1	В	2827	0	2772	6	0
1	С	2639	0	2606	17	0
2	А	1	0	0	0	0
2	В	1	0	0	0	0
2	С	1	0	0	0	0
3	В	8	0	12	1	0



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	С	7	0	10	1	0
5	А	152	0	0	3	0
5	В	237	0	0	1	0
5	С	146	0	0	4	0
All	All	8768	0	8095	31	0

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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 2.

All (31) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash
1:A:760:SER:O	1:A:764:ASN:HB2	1.99	overlap (Å) 0.62
1:A:788:MET:HA	1:A:872:LEU:HD22	1.85	0.58
1:C:860:GLN:NE2	5:C:1302:HOH:O	2.35	0.57
1:C:935:GLN:HA	1:C:941:LYS:O	2.06	0.55
1:C:860:GLN:NE2	1:C:871:GLU:OE1	2.40	0.52
1:C:892:ARG:NH1	5:C:1306:HOH:O	2.42	0.52
1:C:933:THR:O	1:C:996:LYS:HA	2.11	0.51
1:C:1079:ARG:HG2	5:C:1323:HOH:O	2.10	0.51
1:A:896:GLU:OE1	1:B:1019:ARG:NH2	2.44	0.51
1:B:860:GLN:NE2	1:B:871:GLU:OE2	2.43	0.50
1:C:956:LEU:HD13	1:C:1026:THR:HB	1.94	0.49
1:C:936:CYS:O	1:C:940:HIS:HA	2.13	0.49
1:A:1015:SER:HB3	1:A:1022:GLN:HB2	1.95	0.48
1:B:786:CYS:SG	5:B:1424:HOH:O	2.55	0.48
1:B:835:LYS:O	1:B:839:THR:HG23	2.13	0.47
1:C:788:MET:HA	1:C:872:LEU:CD2	2.44	0.47
1:C:1047:ASN:O	5:C:1301:HOH:O	2.20	0.47
1:C:934:VAL:O	1:C:942:LYS:HA	2.14	0.47
1:C:1037:LEU:O	1:C:1041:VAL:HG23	2.14	0.47
1:C:1017:ASP:OD1	1:C:1017:ASP:C	2.54	0.46
1:B:1053:LEU:HD21	1:B:1056:VAL:CG2	2.46	0.46
1:B:815:ASN:HD21	3:B:1201:EDO:H11	1.80	0.46
1:C:930:PHE:O	1:C:946:PHE:HA	2.16	0.45
1:A:1054:PHE:CZ	1:A:1105:PHE:HB3	2.52	0.45
1:A:1082:LYS:NZ	5:A:1306:HOH:O	2.45	0.44
1:C:856:LYS:HA	4:C:1201:PEG:H41	2.01	0.42
1:A:897:ASN:HA	5:A:1395:HOH:O	2.19	0.42
1:C:1054:PHE:CZ	1:C:1105:PHE:HB3	2.55	0.41
1:A:844:TYR:CG	1:A:1086:HIS:HB2	2.56	0.40



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:915:LYS:NZ	5:A:1312:HOH:O	2.53	0.40
1:C:804:TYR:OH	1:C:813:ASP:OD2	2.28	0.40

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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	Perce	\mathbf{n} tiles
1	А	336/377~(89%)	323~(96%)	13~(4%)	0	100	100
1	В	344/377~(91%)	337~(98%)	7~(2%)	0	100	100
1	С	322/377~(85%)	313 (97%)	9(3%)	0	100	100
All	All	1002/1131~(89%)	973~(97%)	29 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	А	303/336~(90%)	303~(100%)	0	100 100
1	В	309/336~(92%)	309 (100%)	0	100 100
1	С	289/336~(86%)	289 (100%)	0	100 100
All	All	901/1008 (89%)	901 (100%)	0	100 100



There are no protein residues with a non-rotameric sidechain to report.

Sometimes 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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 for Mogul analysis.

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).

Mal	Mol Type	Chain	Res	Link	B	Bond lengths			Bond angles		
	туре	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2	
4	PEG	С	1201	-	$6,\!6,\!6$	0.29	0	$5,\!5,\!5$	0.21	0	
3	EDO	В	1202	-	$3,\!3,\!3$	0.22	0	2,2,2	0.25	0	
3	EDO	В	1201	-	$3,\!3,\!3$	0.20	0	2,2,2	0.29	0	

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
4	PEG	С	1201	-	-	3/4/4/4	-
3	EDO	В	1202	-	-	1/1/1/1	-



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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	В	1201	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	С	1201	PEG	O1-C1-C2-O2
4	С	1201	PEG	O2-C3-C4-O4
3	В	1202	EDO	O1-C1-C2-O2
3	В	1201	EDO	O1-C1-C2-O2
4	С	1201	PEG	C4-C3-O2-C2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	С	1201	PEG	1	0
3	В	1201	EDO	1	0

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, 95^{th} 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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	$\mathbf{Q}{<}0.9$
1	А	341/377~(90%)	-0.00	11 (3%) 47 54	20, 37, 77, 93	0
1	В	347/377~(92%)	-0.27	7 (2%) 65 69	18, 28, 54, 96	0
1	С	327/377~(86%)	0.18	31 (9%) 8 10	18, 34, 76, 101	0
All	All	1015/1131 (89%)	-0.03	49 (4%) 30 36	18, 33, 73, 101	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	759 ALA		8.3
1	С	938	THR	4.6
1	А	980	ASN	4.6
1	В	1020	TRP	4.4
1	В	1046	ASN	4.3
1	В	980 ASN		4.3
1	В	1016	TYR	4.0
1	С	994 SER		3.9
1	С	937	LEU	3.9
1	С	940	HIS	3.8
1	С	759	ALA	3.8
1	С	945	THR	3.7
1	С	936	CYS	3.7
1	С	941	LYS	3.7
1	С	974	GLU	3.7
1	С	757	LEU	3.6
1	С	995	LEU	3.6
1	А	762	ILE	3.6
1	С	957	ALA	3.5
1	С	942	2 LYS 3.	
1	В	1019	ARG	3.0
1	А	1063	LEU	3.0
1	С	933	THR	2.9



Mol	Chain	$\frac{1 \text{ previous page.}}{\text{Res} \text{Type}}$		RSRZ
1	В	1018	GLY	2.9
1	А	760	SER	2.9
1	А	979	ASP	2.8
1	С	997	LYS	2.8
1	С	934	VAL	2.7
1	В	763	ARG	2.6
1	С	998	ILE	2.6
1	С	762	ILE	2.5
1	А	1022	GLN	2.5
1	С	902	ASP	2.5
1	С	943	SER	2.5
1	С	1039	GLN	2.5
1	А	761	GLN	2.5
1	С	939	CYS	2.4
1	С	996	LYS	2.4
1	С	760	SER	2.4
1	С	972	SER	2.4
1	А	983	PHE	2.3
1	С	904	PHE	2.2
1	С	935	GLN	2.2
1	А	1065	GLY	2.2
1	С	975	GLU	2.1
1	А	812		
1	С	909	HIS	2.1
1	С	892	ARG	2.1
1	С	973	LYS	2.0

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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)

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^{th} 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	$\operatorname{B-factors}(\operatorname{\AA}^2)$	Q<0.9
2	ZN	С	1202	1/1	0.76	0.09	99,99,99,99	0
3	EDO	В	1201	4/4	0.79	0.20	49,49,51,52	0
4	PEG	С	1201	7/7	0.84	0.14	38,48,51,52	0
3	EDO	В	1202	4/4	0.90	0.10	41,46,51,55	0
2	ZN	А	1201	1/1	0.90	0.06	56, 56, 56, 56	0
2	ZN	В	1203	1/1	1.00	0.10	22,22,22,22	0

6.5 Other polymers (i)

There are no such residues in this entry.

