Full wwPDB X-ray Structure Validation Report (i)

Jan 4, 2021 – 01:04 pm GMT

PDB ID	:	6XZM
Title	:	Arabidopsis UV-B photoreceptor UVR8 mutant D96N D107N W285A
Authors	:	Lau, K.; Hothorn, M.
Deposited on	:	2020-02-04
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 following versions of software and data (see references (i)) were used in the production of this report:

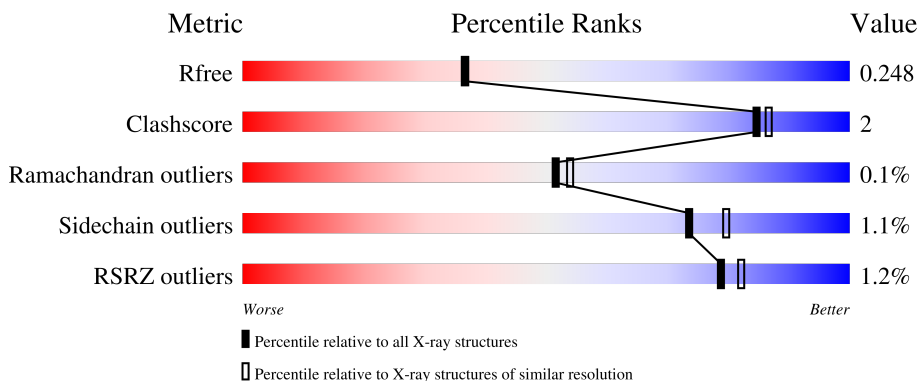|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.16
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.16

1 Overall quality at a glance (i)

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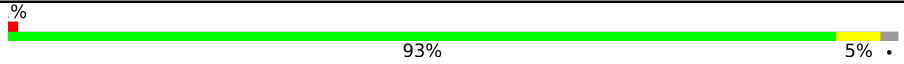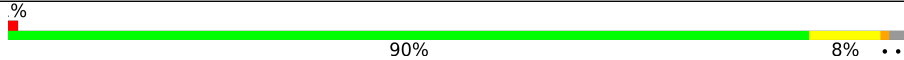*X-RAY DIFFRACTION*

The reported resolution of this entry is 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	Percentile Ranks	Value
Rfree		0.248
Clashscore		2
Ramachandran outliers		0.1%
Sidechain outliers		1.1%
RSRZ outliers		1.2%

Worse ← → *Better*

■ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
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	A	373	93% 5% ·
1	B	373	90% 8% ··

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 10791 atoms, of which 5170 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 Ultraviolet-B receptor UVR8.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	367	Total 5293	C 1711	H 2547	N 492	O 529	S 14	0	0	0
1	B	367	Total 5331	C 1715	H 2579	N 492	O 530	S 15	0	2	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	GLY	-	expression tag	UNP Q9FN03
A	10	ALA	-	expression tag	UNP Q9FN03
A	11	MET	-	expression tag	UNP Q9FN03
A	96	ASN	ASP	engineered mutation	UNP Q9FN03
A	107	ASN	ASP	engineered mutation	UNP Q9FN03
A	285	ALA	TRP	engineered mutation	UNP Q9FN03
B	9	GLY	-	expression tag	UNP Q9FN03
B	10	ALA	-	expression tag	UNP Q9FN03
B	11	MET	-	expression tag	UNP Q9FN03
B	96	ASN	ASP	engineered mutation	UNP Q9FN03
B	107	ASN	ASP	engineered mutation	UNP Q9FN03
B	285	ALA	TRP	engineered mutation	UNP Q9FN03

- Molecule 2 is NITRATE ION (three-letter code: NO3) (formula: NO_3).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 4	N 1	O 3	0	0
2	A	1	Total 4	N 1	O 3	0	0
2	B	1	Total 4	N 1	O 3	0	0
2	B	1	Total 4	N 1	O 3	0	0

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).
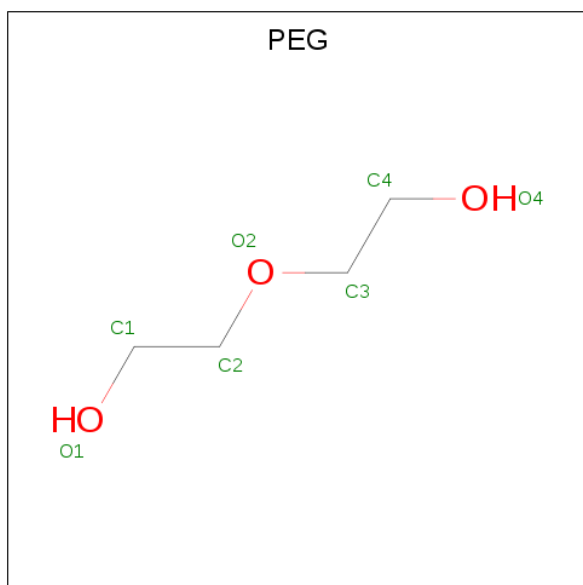
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 17	C 4	H 10	O 3	0	0
3	B	1	Total 17	C 4	H 10	O 3	0	0
3	B	1	Total 17	C 4	H 10	O 3	0	0

- Molecule 4 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).
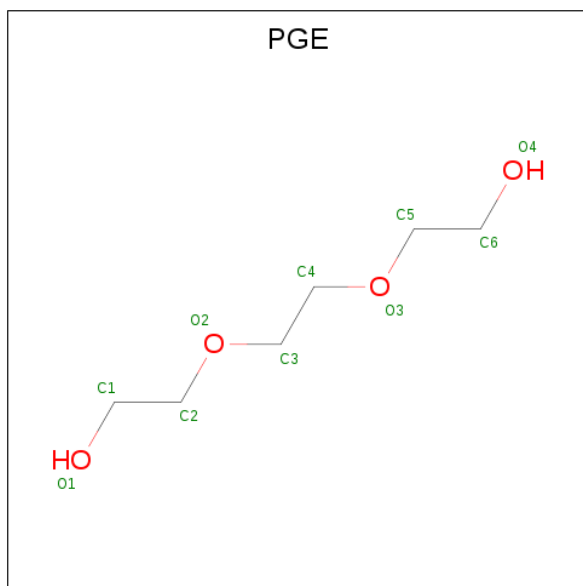
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total 24	C 6	H 14	O 4	0	0

- Molecule 5 is water.

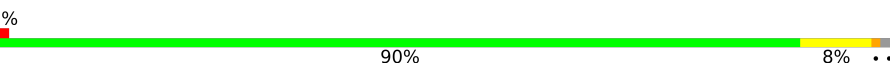| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	39	Total 39	O 39	0	0
5	B	36	Total 37	O 37	0	1

3 Residue-property plots ⓘ

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.
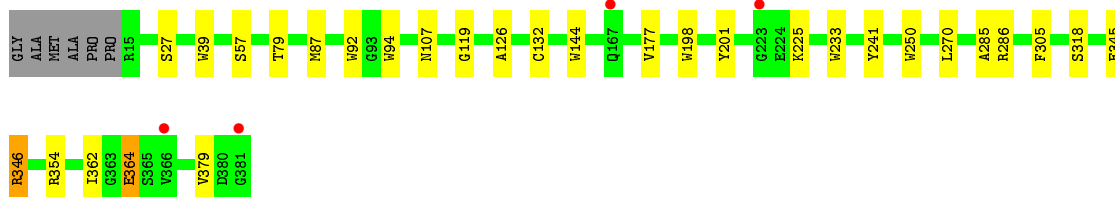
• Molecule 1: Ultraviolet-B receptor UVR8

• Molecule 1: Ultraviolet-B receptor UVR8

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	98.27Å　98.27Å　138.84Å 90.00°　90.00°　120.00°	Depositor
Resolution (Å)	49.13　–　2.10 49.13　–　2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.13-2.10) 88.5 (49.13-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.93 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.197　,　0.248 0.197　,　0.248	Depositor DCC
R_{free} test set	2200 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å^2)	36.8	Xtriage
Anisotropy	0.510	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 38.2	EDS
L-test for twinning[2]	$< |L| > = 0.50, < L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	10791	wwPDB-VP
Average B, all atoms (Å^2)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% of the height of the origin peak. No significant pseudotranslation is detected.*

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[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)

Bond lengths and bond angles in the following residue types are not validated in this section: CME, PEG, PGE, NO3

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	0.54	0/2799	0.68	0/3797
1	B	0.54	0/2811	0.72	3/3813 (0.1%)
All	All	0.54	0/5610	0.70	3/7610 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	346	ARG	NE-CZ-NH2	5.26	122.93	120.30
1	B	270	LEU	CA-CB-CG	5.22	127.31	115.30
1	B	270	LEU	CB-CG-CD2	-5.15	102.25	111.00

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	2746	2547	2602	11	0
1	B	2752	2579	2611	16	0
2	A	8	0	0	1	0
2	B	8	0	0	0	0
3	A	7	10	10	2	1
3	B	14	20	20	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	10	14	14	2	0
5	A	39	0	0	0	0
5	B	37	0	0	0	0
All	All	5621	5170	5257	25	1

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 (25) 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:B:198:TRP:CZ3	4:B:405:PGE:O1	2.55	0.58
1:B:305:PHE:CD1	1:B:354:ARG:HD3	2.39	0.57
1:B:286:ARG:NH2	4:B:405:PGE:H6	2.22	0.55
1:A:89:VAL:O	1:A:113:PRO:HA	2.12	0.49
1:A:304:LYS:HE2	1:B:379:VAL:O	2.13	0.48
1:A:285:ALA:HA	2:A:402:NO3:O3	2.13	0.48
1:A:330:VAL:O	3:A:403:PEG:H22	2.13	0.48
1:B:225:LYS:HB2	1:B:241:TYR:HD2	1.78	0.48
1:A:185:ALA:HB1	1:A:229:VAL:HG21	1.97	0.46
1:B:87:MET:HB3	1:B:119:GLY:HA2	1.97	0.46
1:A:205:GLY:O	1:A:218:VAL:HG13	2.17	0.45
1:B:126:ALA:HB3	1:B:177:VAL:HG23	1.99	0.44
1:B:362:ILE:HG13	1:B:364:GLU:HG2	1.99	0.44
1:B:94:TRP:HZ3	1:B:107:ASN:HD21	1.66	0.44
1:B:345:GLU:HG3	1:B:346:ARG:HG2	2.00	0.43
1:A:331:VAL:HG12	3:A:403:PEG:H31	2.01	0.43
1:A:34:ASP:OD1	1:A:68:GLN:HA	2.18	0.43
1:B:132:CYS:HB2	1:B:144:TRP:CE2	2.55	0.42
1:B:198:TRP:CH2	1:B:201:TYR:HE1	2.37	0.42
1:B:27:SER:HB2	1:B:39:TRP:CE2	2.54	0.42
1:B:79:THR:HB	1:B:92:TRP:CE2	2.54	0.42
1:A:304:LYS:CE	1:B:379:VAL:O	2.68	0.42
1:A:345:GLU:C	1:A:347:ASN:H	2.24	0.41
1:A:132:CYS:HB2	1:A:144:TRP:CE2	2.56	0.41
1:B:233:TRP:CD2	1:B:285:ALA:HB1	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:403:PEG:O4	3:A:403:PEG:O4[6_554]	1.99	0.21

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	364/373 (98%)	354 (97%)	10 (3%)	0	100	100
1	B	366/373 (98%)	352 (96%)	13 (4%)	1 (0%)	41	41
All	All	730/746 (98%)	706 (97%)	23 (3%)	1 (0%)	51	54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	364	GLU

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	281/291 (97%)	278 (99%)	3 (1%)	73	79
1	B	283/291 (97%)	280 (99%)	3 (1%)	73	79
All	All	564/582 (97%)	558 (99%)	6 (1%)	73	79

All (6) residues with a non-rotameric sidechain are listed below:

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Mol	Chain	Res	Type
1	A	248	TYR
1	A	251	SER
1	A	327	ASP
1	B	57	SER
1	B	250	TRP
1	B	318	SER

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

Mol	Chain	Res	Type
1	B	102	HIS
1	B	107	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)

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

8 ligands are modelled in this entry.

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 (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(Å²)	Q<0.9
1	A	366/373 (98%)	0.19	5 (1%)	75	78	35, 46, 63, 77	0
1	B	366/373 (98%)	0.11	4 (1%)	80	84	37, 47, 64, 80	0
All	All	732/746 (98%)	0.15	9 (1%)	79	82	35, 47, 65, 80	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	223	GLY	4.3
1	A	381	GLY	3.4
1	B	381	GLY	2.9
1	A	28	VAL	2.3
1	A	250	TRP	2.2
1	A	285	ALA	2.2
1	B	366	VAL	2.1
1	B	167	GLN	2.0
1	A	172	ILE	2.0

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(Å²)	Q<0.9
1	CME	B	317	10/11	0.91	0.18	39,56,109,109	0
1	CME	A	317	10/11	0.97	0.12	38,51,71,71	0

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, 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	PEG	B	403	7/7	0.81	0.17	60,72,82,84	0
3	PEG	A	403	7/7	0.87	0.33	49,59,71,79	0
4	PGE	B	405	10/10	0.88	0.20	46,61,73,80	0
2	NO3	A	402	4/4	0.90	0.16	40,47,49,53	0
3	PEG	B	404	7/7	0.92	0.12	50,61,73,78	0
2	NO3	B	402	4/4	0.95	0.22	48,51,53,59	0
2	NO3	B	401	4/4	0.95	0.16	43,47,52,52	0
2	NO3	A	401	4/4	0.98	0.13	37,38,42,46	0

6.5 Other polymers (i)

There are no such residues in this entry.