



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:09 PM UTC

PDB ID : 7XV7 / pdb_00007xv7
Title : Crystal structure of ZYG11B bound to ORF10 peptide
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Deposited on : 2022-05-21
Resolution : 2.60 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

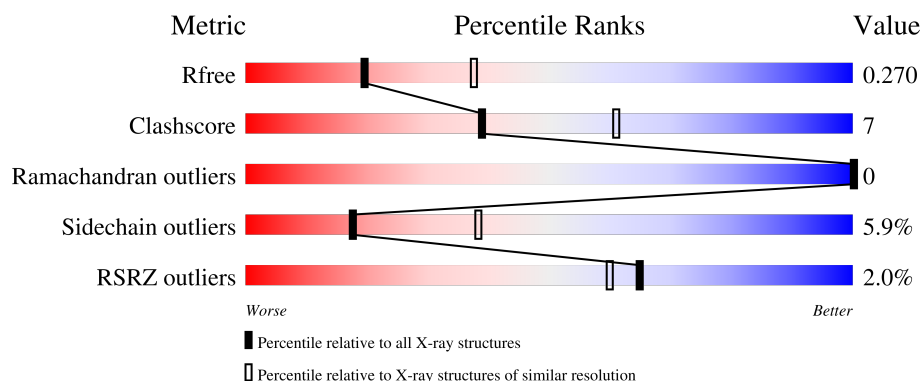
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	250	 77% 18% ..
1	B	250	 80% 17% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3988 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 Protein zyg-11 homolog B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	246	Total	C	N	O	S	0	1	0
			1981	1266	335	370	10			
1	A	244	Total	C	N	O	S	0	1	0
			1963	1255	333	365	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	479	GLY	-	expression tag	UNP Q9C0D3
B	480	TYR	-	expression tag	UNP Q9C0D3
B	481	ILE	-	expression tag	UNP Q9C0D3
B	482	ASN	-	expression tag	UNP Q9C0D3
B	483	VAL	-	expression tag	UNP Q9C0D3
B	484	GLY	-	expression tag	UNP Q9C0D3
A	479	GLY	-	expression tag	UNP Q9C0D3
A	480	TYR	-	expression tag	UNP Q9C0D3
A	481	ILE	-	expression tag	UNP Q9C0D3
A	482	ASN	-	expression tag	UNP Q9C0D3
A	483	VAL	-	expression tag	UNP Q9C0D3
A	484	GLY	-	expression tag	UNP Q9C0D3

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

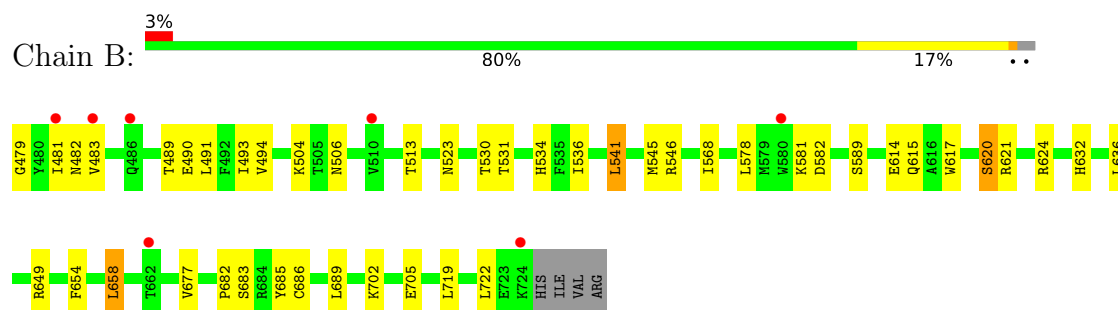
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	13	Total	O	0	0
			13	13		
3	A	16	Total	O	0	0
			16	16		

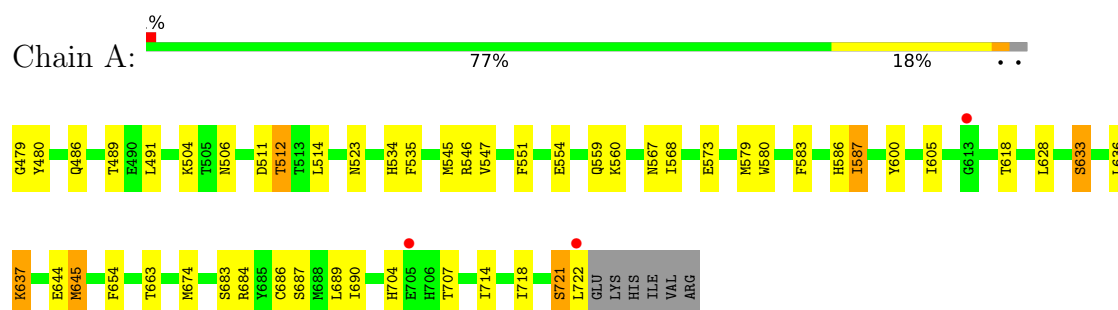
3 Residue-property plots [i](#)

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: Protein zyg-11 homolog B



- Molecule 1: Protein zyg-11 homolog B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.50Å 96.85Å 121.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.21 – 2.60 31.21 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.5 (31.21-2.60) 98.4 (31.21-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.55 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.203 , 0.269 (Not available) , 0.270	Depositor DCC
R_{free} test set	1579 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	58.1	Xtriage
Anisotropy	0.500	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.4	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	0.95	EDS
Total number of atoms	3988	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.55% 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: SO4

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	0.42	0/2010	0.59	0/2734
1	B	0.38	0/2028	0.57	0/2757
All	All	0.40	0/4038	0.58	0/5491

There are no bond length outliers.

There are no bond angle outliers.

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	1963	0	1933	32	0
1	B	1981	0	1950	28	0
2	A	10	0	0	0	0
2	B	5	0	0	0	0
3	A	16	0	0	1	0
3	B	13	0	0	0	0
All	All	3988	0	3883	53	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 (53) 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:489:THR:HG21	1:A:523:ASN:HD22	1.50	0.76
1:A:506:ASN:OD1	1:A:546:ARG:NH2	2.18	0.69
1:A:551:PHE:HB3	1:A:554:GLU:HG3	1.78	0.64
1:A:545:MET:HE1	1:A:586:HIS:HB2	1.79	0.63
1:B:632:HIS:CE1	1:B:636:LEU:HD11	2.35	0.62
1:B:490:GLU:HA	1:B:493:ILE:HD12	1.82	0.62
1:A:633:SER:O	1:A:637:LYS:HD2	2.04	0.57
1:A:479:GLY:N	3:A:901:HOH:O	2.39	0.55
1:A:504:LYS:HG3	1:A:514:LEU:HB2	1.89	0.54
1:B:523:ASN:OD1	1:A:480:TYR:HB2	2.07	0.54
1:A:545:MET:HE1	1:A:586:HIS:CB	2.37	0.54
1:B:649:ARG:HH22	1:A:486:GLN:HG3	1.74	0.53
1:B:541:LEU:HD21	1:B:578:LEU:HD23	1.91	0.52
1:A:511:ASP:OD1	1:A:512:THR:N	2.44	0.51
1:A:545:MET:HG3	1:A:580:TRP:CZ3	2.45	0.51
1:B:617:TRP:NE1	1:B:624:ARG:HB2	2.26	0.50
1:B:614:GLU:H	1:B:614:GLU:CD	2.20	0.50
1:B:491:LEU:HD13	1:B:534:HIS:ND1	2.27	0.49
1:A:579:MET:HE2	1:A:618:THR:H	1.77	0.49
1:A:535:PHE:CD1	1:A:568:ILE:HD11	2.48	0.49
1:A:600:TYR:CG	1:A:645:MET:HE1	2.48	0.48
1:A:491:LEU:HB3	1:A:534:HIS:CD2	2.48	0.48
1:A:686:CYS:O	1:A:690:ILE:HG13	2.14	0.48
1:B:649:ARG:HH21	1:A:486:GLN:HE21	1.60	0.48
1:A:714:ILE:O	1:A:718:ILE:HG13	2.13	0.48
1:A:559:GLN:HG3	1:A:560:LYS:N	2.27	0.48
1:A:628:LEU:HD11	1:A:663:THR:HG22	1.96	0.47
1:B:654:PHE:O	1:B:658:LEU:HD12	2.15	0.47
1:B:617:TRP:NE1	1:B:620:SER:O	2.40	0.47
1:A:721:SER:O	1:A:722:LEU:HD23	2.14	0.46
1:A:704:HIS:O	1:A:707:THR:OG1	2.28	0.45
1:B:479:GLY:N	1:A:567:ASN:OD1	2.50	0.45
1:B:682:PRO:O	1:B:686:CYS:HB2	2.17	0.45
1:B:677:VAL:HB	1:B:685:TYR:CD2	2.53	0.44
1:B:504:LYS:HE3	1:B:513:THR:HB	1.99	0.44
1:B:689:LEU:HD23	1:B:722:LEU:HD11	2.00	0.43
1:B:491:LEU:HB3	1:B:534:HIS:CD2	2.54	0.43
1:B:719:LEU:HD23	1:B:719:LEU:HA	1.85	0.43
1:B:483:VAL:HG11	1:A:684:ARG:CZ	2.48	0.43
1:B:581:LYS:HE2	1:B:581:LYS:HB3	1.80	0.43
1:A:545:MET:SD	1:A:583:PHE:HA	2.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:649:ARG:NH2	1:A:486:GLN:HG3	2.34	0.43
1:A:587:ILE:HD12	1:A:605:ILE:HG21	2.01	0.42
1:A:689:LEU:HD12	1:A:689:LEU:HA	1.79	0.42
1:B:481:ILE:HG23	1:B:482:ASN:H	1.83	0.42
1:B:536:ILE:HD11	1:B:568:ILE:HG21	2.02	0.42
1:B:481:ILE:HG23	1:B:482:ASN:N	2.35	0.42
1:B:545:MET:HE3	1:B:545:MET:HB2	1.76	0.42
1:B:494:VAL:HG21	1:B:531:THR:HG23	2.01	0.41
1:B:506:ASN:OD1	1:B:546:ARG:NH2	2.53	0.41
1:A:636:LEU:HD23	1:A:636:LEU:HA	1.81	0.40
1:A:573:GLU:CD	1:A:573:GLU:H	2.29	0.40
1:A:654:PHE:CB	1:A:674:MET:HE3	2.52	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	243/250 (97%)	229 (94%)	14 (6%)	0	100	100
1	B	245/250 (98%)	242 (99%)	3 (1%)	0	100	100
All	All	488/500 (98%)	471 (96%)	17 (4%)	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	A	195/226 (86%)	184 (94%)	11 (6%)	19	40
1	B	181/226 (80%)	170 (94%)	11 (6%)	17	37
All	All	376/452 (83%)	354 (94%)	22 (6%)	18	38

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	530	THR
1	B	541	LEU
1	B	582	ASP
1	B	589	SER
1	B	615	GLN
1	B	620	SER
1	B	621	ARG
1	B	658	LEU
1	B	683	SER
1	B	702	LYS
1	B	705	GLU
1	A	489	THR
1	A	512	THR
1	A	547	VAL
1	A	587	ILE
1	A	633	SER
1	A	637	LYS
1	A	644	GLU
1	A	645	MET
1	A	683	SER
1	A	687	SER
1	A	721	SER

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

Mol	Chain	Res	Type
1	B	538	ASN
1	B	592	HIS
1	B	632	HIS
1	B	710	HIS
1	A	486	GLN
1	A	608	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 ligands are 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	$\# Z > 2$
2	SO4	B	801	-	4,4,4	0.29	0	6,6,6	0.08	0
2	SO4	A	801	-	4,4,4	0.25	0	6,6,6	0.28	0
2	SO4	A	802	-	4,4,4	0.21	0	6,6,6	0.33	0

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 ⓘ

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	244/250 (97%)	0.13	3 (1%) 76 73	29, 67, 99, 140	1 (0%)
1	B	246/250 (98%)	0.24	7 (2%) 55 49	34, 68, 107, 153	1 (0%)
All	All	490/500 (98%)	0.19	10 (2%) 65 60	29, 67, 105, 153	2 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	483	VAL	3.6
1	B	580	TRP	3.5
1	B	486	GLN	2.7
1	B	662	THR	2.6
1	B	510	VAL	2.5
1	A	722	LEU	2.5
1	A	613	GLY	2.4
1	A	705	GLU	2.2
1	B	481	ILE	2.2
1	B	724	LYS	2.1

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](#)

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
2	SO4	B	801	5/5	0.71	0.14	146,146,146,147	0
2	SO4	A	802	5/5	0.85	0.14	97,99,102,103	0
2	SO4	A	801	5/5	0.96	0.09	71,71,75,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.