

Full wwPDB X-ray Structure Validation Report (i)

Dec 28, 2024 – 04:01 PM EST

PDB ID : 8VSQ

Title : Crystal structure of Esub1 Authors : Muna, T.H.; Prehna, G.

Deposited on : 2024-01-24

Resolution : 2.90 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 : 2022.3.0, CSD as543be (2022)

Xtriage (Phenix) : 1.21 EDS : 3.0

Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)

CCP4 : 9.0.004 (Gargrove)

Density-Fitness : 1.0.11

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

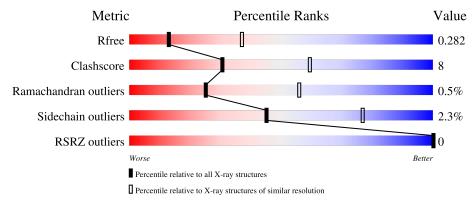
Validation Pipeline (wwPDB-VP) : 2.40

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 2.90 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	Similar resolution
Metric	$(\# ext{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	164625	2335 (2.90-2.90)
Clashscore	180529	2564 (2.90-2.90)
Ramachandran outliers	177936	2514 (2.90-2.90)
Sidechain outliers	177891	2516 (2.90-2.90)
RSRZ outliers	164620	2337 (2.90-2.90)

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	206	76%	18%	• 5%
1	В	206	76%	18%	• 5%



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3179 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 Esub1.

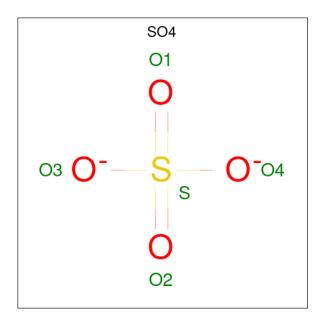
\mathbf{Mol}	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	196	Total	С	N	О	S	0	0	0
1	Λ	190	1562	998	273	287	4	0	U	U
1	B	196	Total	С	N	О	S	0	0	0
1	Ъ	190	1562	998	273	287	4		0 0	U

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	199	LEU	-	expression tag	UNP A0A0H2UUU0
A	200	GLU	-	expression tag	UNP A0A0H2UUU0
A	201	HIS	-	expression tag	UNP A0A0H2UUU0
A	202	HIS	-	expression tag	UNP A0A0H2UUU0
A	203	HIS	-	expression tag	UNP A0A0H2UUU0
A	204	HIS	-	expression tag	UNP A0A0H2UUU0
A	205	HIS	-	expression tag	UNP A0A0H2UUU0
A	206	HIS	-	expression tag	UNP A0A0H2UUU0
В	199	LEU	-	expression tag	UNP A0A0H2UUU0
В	200	GLU	-	expression tag	UNP A0A0H2UUU0
В	201	HIS	-	expression tag	UNP A0A0H2UUU0
В	202	HIS	-	expression tag	UNP A0A0H2UUU0
В	203	HIS	-	expression tag	UNP A0A0H2UUU0
В	204	HIS	-	expression tag	UNP A0A0H2UUU0
В	205	HIS	-	expression tag	UNP A0A0H2UUU0
В	206	HIS	-	expression tag	UNP A0A0H2UUU0

• Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).





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

• Molecule 3 is water.

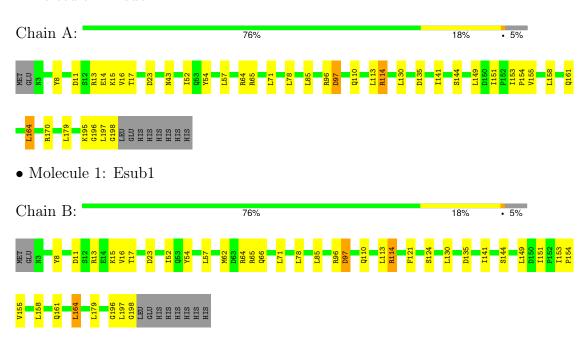
\mathbf{Mol}	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
3	A	24	Total O 24 24	0	0
3	В	26	Total O 26 26	0	0



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: Esub1





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants	82.36Å 43.87Å 85.84Å	Donogitor
a, b, c, α , β , γ	90.00° 118.67° 90.00°	Depositor
Resolution (Å)	43.91 - 2.90	Depositor
Resolution (A)	43.91 - 2.90	EDS
% Data completeness	98.5 (43.91-2.90)	Depositor
(in resolution range)	98.5 (43.91-2.90)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.15 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
P.P.	0.222 , 0.281	Depositor
R, R_{free}	0.224 , 0.282	DCC
R_{free} test set	583 reflections $(4.96%)$	wwPDB-VP
Wilson B-factor (Å ²)	64.7	Xtriage
Anisotropy	1.097	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.33, 54.2	EDS
L-test for twinning ²	$< L > = 0.49, < L^2> = 0.32$	Xtriage
	0.002 for l,k,-h-l	
	0.002 for -h-l,k,h	
Estimated twinning fraction	0.478 for h,-k,-h-l	Xtriage
	0.009 for l,-k,h	
	0.012 for -h-l,-k,l	
F_o, F_c correlation	0.98	EDS
Total number of atoms	3179	wwPDB-VP
Average B, all atoms (\mathring{A}^2)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.91% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



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: 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		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.77	0/1588	1.00	2/2145 (0.1%)	
1	В	0.77	0/1588	1.01	0/2145	
All	All	0.77	0/3176	1.00	2/4290 (0.0%)	

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
1	A	170	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	A	170	ARG	NE-CZ-NH2	-5.17	117.71	120.30

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	1562	0	1606	28	1
1	В	1562	0	1606	27	0
2	A	5	0	0	0	0
3	A	24	0	0	3	0
3	В	26	0	0	0	0
All	All	3179	0	3212	52	1



The all-atom clash score is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clash score for this structure is 8.

All (52) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${\rm distance}({\rm \AA})$	overlap (Å)
1:B:64:ARG:HG2	1:B:65:ARG:H	1.28	0.93
1:A:64:ARG:HG2	1:A:65:ARG:H	1.46	0.81
1:B:13:ARG:O	1:B:15:LYS:HG3	1.90	0.70
1:A:13:ARG:O	1:A:15:LYS:HG3	1.93	0.68
1:A:64:ARG:NH1	3:A:401:HOH:O	1.99	0.64
1:B:64:ARG:HG2	1:B:65:ARG:N	2.09	0.63
1:A:8:TYR:OH	1:A:154:PRO:HD3	1.98	0.62
1:B:8:TYR:OH	1:B:154:PRO:HD3	1.98	0.62
1:A:17:THR:HG22	1:A:149:LEU:O	1.99	0.62
1:A:164:LEU:HD11	1:B:196:GLY:HA3	1.83	0.61
1:B:17:THR:HG22	1:B:149:LEU:O	2.00	0.60
1:B:16:VAL:HG13	1:B:155:VAL:HG21	1.87	0.55
1:A:114:ARG:HH11	1:A:158:LEU:HD22	1.71	0.55
1:A:179:LEU:HD22	1:B:179:LEU:HD22	1.88	0.55
1:A:16:VAL:HG13	1:A:155:VAL:HG21	1.88	0.55
1:A:196:GLY:HA3	1:B:164:LEU:HD11	1.90	0.53
1:B:62:MET:HG3	1:B:66:GLN:OE1	2.10	0.51
1:A:11:ASP:OD1	1:A:13:ARG:HG3	2.10	0.51
1:A:64:ARG:HG2	1:A:65:ARG:N	2.21	0.51
1:B:11:ASP:OD1	1:B:13:ARG:HG3	2.10	0.51
1:B:114:ARG:HH11	1:B:158:LEU:HD22	1.76	0.51
1:A:135:ASP:OD1	1:A:141:ILE:HA	2.11	0.51
1:B:135:ASP:OD1	1:B:141:ILE:HA	2.12	0.50
1:A:114:ARG:HB2	1:A:161:GLN:OE1	2.11	0.50
1:B:114:ARG:HB2	1:B:161:GLN:OE1	2.11	0.50
1:A:130:LEU:HD13	1:A:151:ILE:HG23	1.92	0.50
1:B:23:ASP:OD2	1:B:110:GLN:NE2	2.41	0.49
1:A:52:ILE:CG2	1:A:54:TYR:CE1	2.97	0.48
1:B:130:LEU:HD13	1:B:151:ILE:HG23	1.95	0.48
1:A:23:ASP:OD2	1:A:110:GLN:NE2	2.42	0.48
1:B:52:ILE:CG2	1:B:54:TYR:CE1	2.98	0.47
1:B:71:LEU:HD21	1:B:97:ASP:HB3	1.97	0.47
1:A:52:ILE:HG21	1:A:54:TYR:CE1	2.50	0.46
1:B:54:TYR:HA	1:B:57:LEU:HG	1.97	0.46
1:B:78:LEU:N	1:B:78:LEU:HD23	2.30	0.45
1:B:62:MET:CG	1:B:66:GLN:OE1	2.65	0.45
1:B:52:ILE:HG21	1:B:54:TYR:CE1	2.52	0.45

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Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${f distance} ({f A})$	overlap (Å)
1:A:71:LEU:HD21	1:A:97:ASP:HB3	1.99	0.45
1:A:78:LEU:N	1:A:78:LEU:HD23	2.31	0.45
1:A:54:TYR:HA	1:A:57:LEU:HG	1.98	0.44
1:B:121:PHE:O	1:B:124:SER:OG	2.25	0.44
1:A:195:LYS:HE2	3:A:407:HOH:O	2.17	0.43
1:A:197:LEU:O	1:A:198:GLY:C	2.56	0.43
1:B:197:LEU:O	1:B:198:GLY:C	2.56	0.43
1:A:113:LEU:HD13	1:A:153:ILE:HG21	2.01	0.42
1:A:54:TYR:O	1:A:96:ARG:NH2	2.53	0.42
1:B:8:TYR:O	1:B:11:ASP:HB2	2.19	0.42
1:B:54:TYR:O	1:B:96:ARG:NH2	2.53	0.42
1:A:8:TYR:O	1:A:11:ASP:HB2	2.20	0.41
1:A:43:ASN:ND2	3:A:402:HOH:O	2.54	0.41
1:A:64:ARG:CG	1:A:65:ARG:H	2.25	0.41
1:B:113:LEU:HD13	1:B:153:ILE:HG21	2.03	0.41

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	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} ({\rm \AA}) \end{array}$	Clash overlap (Å)	
1:A:14:GLU:OE1	1:A:17:THR:OG1[2_554]	2.04	0.16	

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	Analysed Favoured Allowed		Outliers	Perce	entiles
1	A	194/206 (94%)	182 (94%)	11 (6%)	1 (0%)	25	56
1	В	194/206 (94%)	182 (94%)	11 (6%)	1 (0%)	25	56
All	All	388/412 (94%)	364 (94%)	22 (6%)	2 (0%)	25	56

All (2) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	144	SER
1	В	144	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	Analysed Rotameric Outlier		Percentiles
1	A	172/183 (94%)	168 (98%)	4 (2%)	45 77
1	В	172/183~(94%)	168 (98%)	4 (2%)	45 77
All	All	344/366 (94%)	336 (98%)	8 (2%)	45 77

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	85	LEU
1	A	97	ASP
1	A	114	ARG
1	A	164	LEU
1	В	85	LEU
1	В	97	ASP
1	В	114	ARG
1	В	164	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	В	61	GLN

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)

1 ligand 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	В	ond leng	$_{ m gths}$	В	ond ang	gles
WIOI I	Type		n nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	301	-	4,4,4	0.33	0	6,6,6	0.13	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 (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	$\langle { m RSRZ} \rangle$	# RSRZ > 2		$OWAB(Å^2)$	Q<0.9
1	A	196/206 (95%)	-1.64	0 100	100	53, 82, 115, 144	0
1	В	196/206 (95%)	-1.67	0 100	100	54, 81, 117, 157	0
All	All	392/412 (95%)	-1.65	0 100	100	53, 81, 117, 157	0

There are no RSRZ outliers to report.

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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	SO4	A	301	5/5	0.99	0.04	104,105,114,122	0

6.5 Other polymers (i)

There are no such residues in this entry.

