



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:09 PM UTC

PDB ID : 7SD2 / pdb\_00007sd2  
Title : Murine Fab that recognizes Hev b 8 (profilin for Hevea brasiliensis)  
Authors : Rodriguez-Romero, A.; Garcia-Ramirez, B.  
Deposited on : 2021-09-29  
Resolution : 3.75 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Xtrriage (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

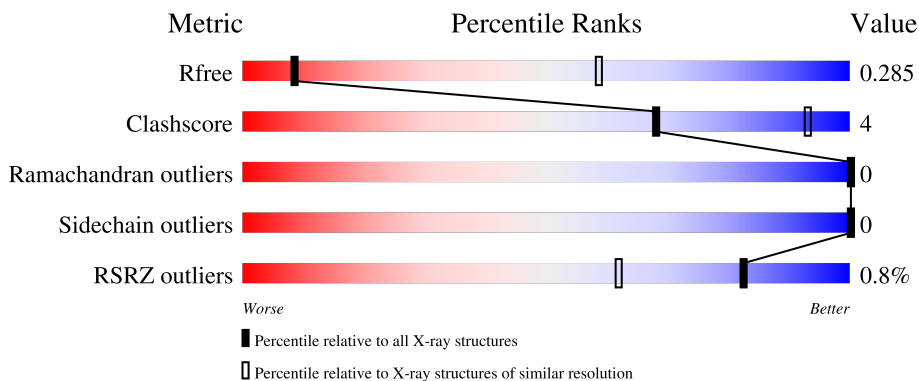
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.75 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	1029 (3.90-3.62)
Clashscore	190562	1061 (3.90-3.62)
Ramachandran outliers	187476	1014 (3.90-3.62)
Sidechain outliers	187428	1009 (3.90-3.62)
RSRZ outliers	180081	1028 (3.90-3.62)

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  $\geq 3$ , 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	209	 90% 6%
1	C	209	 88% 6% 6%
1	H	209	 88% 10%
2	B	214	 95%
2	D	214	 89% 8%

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Mol	Chain	Length	Quality of chain
2	L	214	 91% 7%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8148 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 Heavy Chain Antibody IgE/Fab anti-profilin Hev b 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	H	204	1406	887	249	262	8	0	0	0
1	A	201	1345	853	238	247	7	0	0	0
1	C	197	1164	728	211	219	6	0	0	0

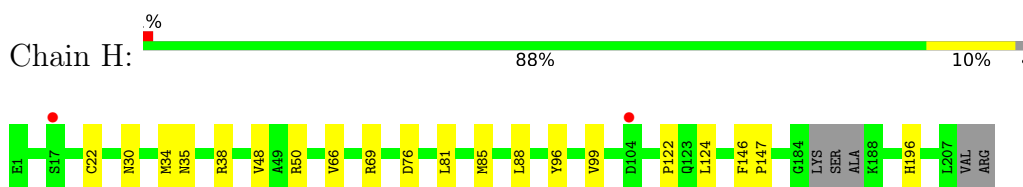
- Molecule 2 is a protein called Light Chain Antibody IgE/Fab anti-profilin Hev b 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	210	1434	904	247	277	6	0	0	0
2	B	210	1406	886	242	272	6	0	0	0
2	D	209	1393	882	243	263	5	0	0	0

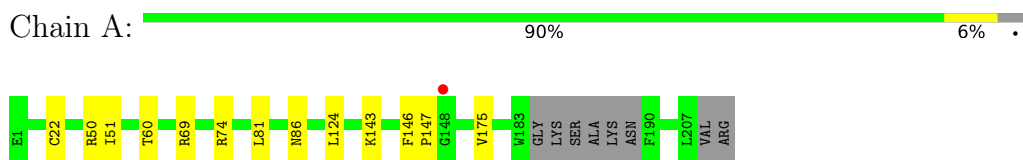
### 3 Residue-property plots [i](#)

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.

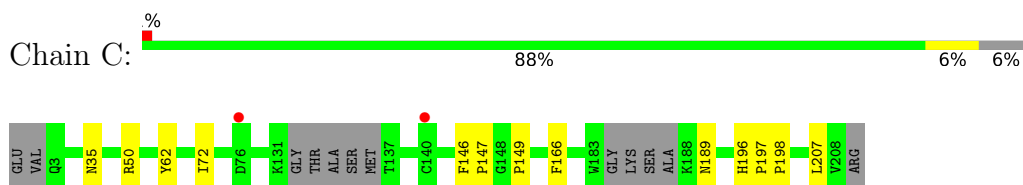
- Molecule 1: Heavy Chain Antibody IgE/Fab anti-profilin Hev b 8



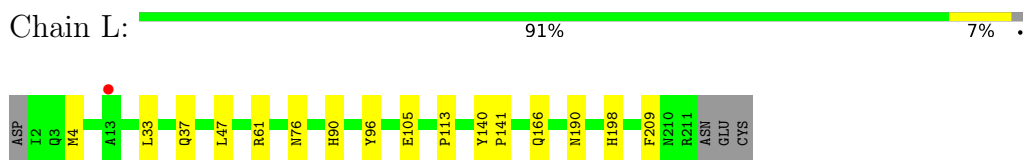
- Molecule 1: Heavy Chain Antibody IgE/Fab anti-profilin Hev b 8



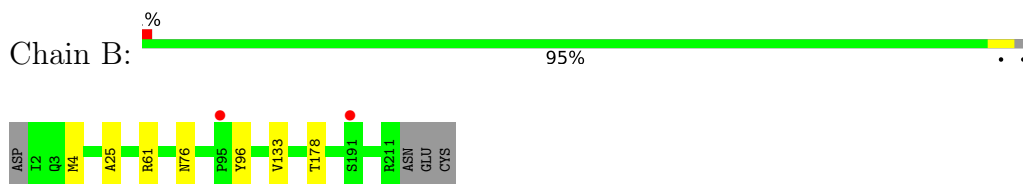
- Molecule 1: Heavy Chain Antibody IgE/Fab anti-profilin Hev b 8



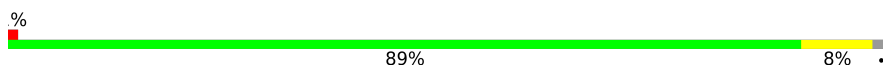
- Molecule 2: Light Chain Antibody IgE/Fab anti-profilin Hev b 8



- Molecule 2: Light Chain Antibody IgE/Fab anti-profilin Hev b 8



- Molecule 2: Light Chain Antibody IgE/Fab anti-profilin Hev b 8

Chain D:  %



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.00Å 103.84Å 180.20Å 90.00° 94.26° 90.00°	Depositor
Resolution (Å)	44.96 – 3.75 44.96 – 3.75	Depositor EDS
% Data completeness (in resolution range)	99.2 (44.96-3.75) 99.1 (44.96-3.75)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.76 (at 3.77Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, $R_{free}$	0.246 , 0.284 0.247 , 0.285	Depositor DCC
$R_{free}$ test set	805 reflections (4.65%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	99.8	Xtrriage
Anisotropy	0.287	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 152.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	8148	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	102.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

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	0.13	0/1379	0.33	0/1904
1	C	0.13	0/1190	0.31	0/1652
1	H	0.13	0/1442	0.32	0/1984
2	B	0.12	0/1445	0.29	0/1992
2	D	0.12	0/1432	0.30	0/1974
2	L	0.11	0/1473	0.29	0/2028
All	All	0.12	0/8361	0.30	0/11534

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

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	1345	0	1067	8	0
1	C	1164	0	726	8	0
1	H	1406	0	1140	13	0
2	B	1406	0	1083	4	0
2	D	1393	0	1061	11	0
2	L	1434	0	1146	10	0
All	All	8148	0	6223	51	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 4.

All (51) 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)
2:L:105:GLU:HA	2:L:166:GLN:HE22	1.53	0.73
2:D:200:THR:HG22	2:D:202:THR:H	1.54	0.70
1:H:85:MET:HE2	1:H:88:LEU:HD21	1.79	0.65
1:A:51:ILE:HG22	1:A:60:THR:HG22	1.81	0.63
2:L:4:MET:HE1	2:L:33:LEU:HD12	1.81	0.62
2:L:37:GLN:HB2	2:L:47:LEU:HD11	1.82	0.61
1:A:50:ARG:HD2	2:B:96:TYR:HE2	1.66	0.60
1:H:50:ARG:HD2	2:L:96:TYR:HE2	1.67	0.59
1:H:34:MET:HB3	1:H:81:LEU:HD22	1.87	0.56
2:B:133:VAL:HG13	2:B:178:THR:HG22	1.87	0.56
1:A:22:CYS:HB3	1:A:81:LEU:HB3	1.87	0.55
2:D:61:ARG:HB2	2:D:76:ASN:O	2.08	0.53
2:D:150:ILE:HG21	2:D:155:ARG:HE	1.74	0.53
1:C:35:ASN:HA	1:C:50:ARG:HA	1.92	0.52
2:D:22:THR:HA	2:D:72:SER:HA	1.92	0.51
1:A:124:LEU:HD23	1:A:124:LEU:H	1.76	0.51
2:D:59:PRO:HB2	2:D:61:ARG:HG2	1.94	0.49
1:A:143:LYS:HA	1:A:175:VAL:HG12	1.95	0.48
1:C:62:TYR:HE1	1:C:72:ILE:H	1.60	0.48
2:B:61:ARG:HB2	2:B:76:ASN:O	2.14	0.48
1:H:30:ASN:HB3	1:H:76:ASP:HB3	1.96	0.47
1:A:69:ARG:HB3	1:A:86:ASN:O	2.14	0.47
2:L:4:MET:HE3	2:L:90:HIS:HB3	1.96	0.47
1:C:149:PRO:HD2	1:C:197:PRO:HG3	1.96	0.46
2:L:113:PRO:HD3	2:L:198:HIS:NE2	2.30	0.46
1:C:166:PHE:CD2	2:D:176:SER:HB3	2.51	0.46
1:H:35:ASN:HD21	1:H:50:ARG:HH11	1.64	0.45
2:B:4:MET:HE2	2:B:25:ALA:HB2	1.99	0.45
2:D:113:PRO:HA	2:D:138:ASN:O	2.16	0.45
2:L:140:TYR:CG	2:L:141:PRO:HA	2.52	0.45
1:C:189:ASN:HA	1:C:207:LEU:HA	1.99	0.45
1:H:38:ARG:HB3	1:H:96:TYR:CD1	2.53	0.44
1:H:124:LEU:H	1:H:124:LEU:HD23	1.83	0.44
1:H:22:CYS:HB3	1:H:81:LEU:HB3	1.99	0.43
1:C:196:HIS:ND1	1:C:198:PRO:HD2	2.33	0.43
1:H:122:PRO:HD3	1:H:196:HIS:CD2	2.54	0.43
1:C:146:PHE:HA	1:C:147:PRO:HA	1.84	0.43
1:H:48:VAL:HG23	1:H:66:VAL:HG21	2.01	0.43
2:D:89:GLN:HG3	2:D:97:THR:O	2.19	0.42
2:L:61:ARG:HB2	2:L:76:ASN:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:134:CYS:HB2	2:D:148:TRP:CZ2	2.54	0.42
1:C:196:HIS:ND1	1:C:197:PRO:HD2	2.35	0.42
1:H:66:VAL:HG12	1:H:69:ARG:NH2	2.35	0.41
2:D:54:LEU:HD23	2:D:54:LEU:HA	1.92	0.41
2:D:113:PRO:HB3	2:D:139:PHE:HB3	2.03	0.41
1:A:146:PHE:HA	1:A:147:PRO:HA	1.88	0.41
2:L:105:GLU:HA	2:L:166:GLN:NE2	2.30	0.41
1:A:51:ILE:HD12	1:A:74:ARG:HD2	2.03	0.40
1:H:35:ASN:HB2	1:H:99:VAL:CG2	2.52	0.40
1:H:146:PHE:HA	1:H:147:PRO:HA	1.89	0.40
2:L:190:ASN:O	2:L:209:PHE:HA	2.22	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	197/209 (94%)	193 (98%)	4 (2%)	0	100	100
1	C	191/209 (91%)	183 (96%)	8 (4%)	0	100	100
1	H	200/209 (96%)	190 (95%)	10 (5%)	0	100	100
2	B	208/214 (97%)	204 (98%)	4 (2%)	0	100	100
2	D	207/214 (97%)	203 (98%)	4 (2%)	0	100	100
2	L	208/214 (97%)	202 (97%)	6 (3%)	0	100	100
All	All	1211/1269 (95%)	1175 (97%)	36 (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	A	100/183 (55%)	100 (100%)	0	100	100
1	C	51/183 (28%)	51 (100%)	0	100	100
1	H	113/183 (62%)	113 (100%)	0	100	100
2	B	109/188 (58%)	109 (100%)	0	100	100
2	D	101/188 (54%)	101 (100%)	0	100	100
2	L	118/188 (63%)	118 (100%)	0	100	100
All	All	592/1113 (53%)	592 (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. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	L	189	HIS
1	A	84	GLN
1	A	86	ASN
2	B	161	ASN
2	D	138	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](#)

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](#)

There are no ligands in this entry.

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	201/209 (96%)	0.25	1 (0%) 87 70	75, 108, 149, 179	0
1	C	197/209 (94%)	0.21	2 (1%) 79 57	94, 144, 177, 204	0
1	H	204/209 (97%)	-0.06	2 (0%) 79 57	48, 80, 137, 181	0
2	B	210/214 (98%)	0.09	2 (0%) 79 57	59, 94, 129, 150	0
2	D	209/214 (97%)	0.10	2 (0%) 79 57	74, 108, 142, 169	0
2	L	210/214 (98%)	0.03	1 (0%) 87 70	46, 77, 120, 146	0
All	All	1231/1269 (97%)	0.10	10 (0%) 82 62	46, 101, 157, 204	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	76	ASP	3.1
1	C	140	CYS	2.9
1	A	148	GLY	2.9
2	B	95	PRO	2.6
2	D	4	MET	2.5
2	L	13	ALA	2.4
1	H	17	SER	2.4
1	H	104	ASP	2.1
2	B	191	SER	2.1
2	D	171	SER	2.0

### 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](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.