



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 9, 2026 – 10:02 am BST

PDB ID : 9SBF / pdb\_00009sbf  
Title : Drosophila Elav, eSH3 fragment (RRM3) - hexagonal crystal form  
Authors : Futterer, K.; Soller, M.  
Deposited on : 2025-08-08  
Resolution : 3.73 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
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

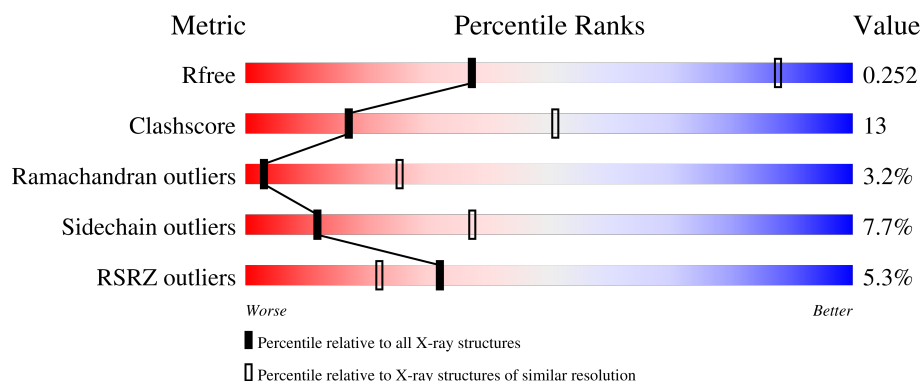
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.73 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	1234 (3.88-3.60)
Clashscore	190562	1274 (3.88-3.60)
Ramachandran outliers	187476	1226 (3.88-3.60)
Sidechain outliers	187428	1221 (3.88-3.60)
RSRZ outliers	180081	1233 (3.88-3.60)

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  $\geq 3$ , 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	109	<div> <div>0%</div> <div> <div>46%</div> <div>21%</div> <div>..</div> <div>31%</div> </div> </div>
1	B	109	<div> <div>2%</div> <div> <div>44%</div> <div>20%</div> <div>.</div> <div>33%</div> </div> </div>
1	C	109	<div> <div>6%</div> <div> <div>50%</div> <div>17%</div> <div>.</div> <div>29%</div> </div> </div>
1	D	109	<div> <div>6%</div> <div> <div>53%</div> <div>14%</div> <div>.</div> <div>31%</div> </div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4056 atoms, of which 1965 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.

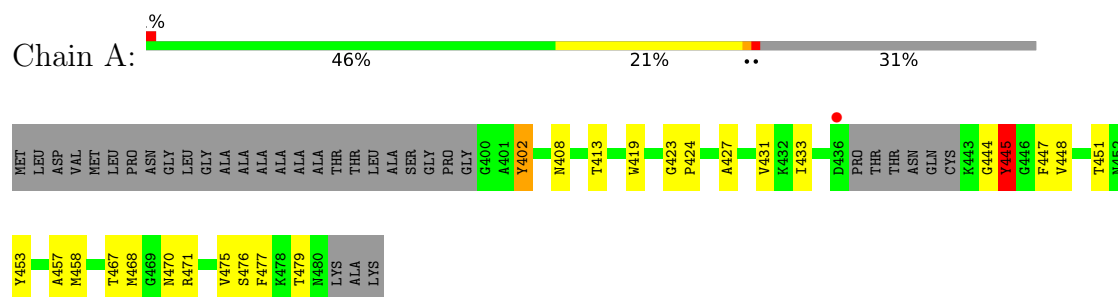
- Molecule 1 is a protein called Protein elav.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	75	Total	C	H	N	O	S	32	0	0
			1052	357	510	83	99	3			
1	B	73	Total	C	H	N	O	S	29	0	0
			1024	345	500	83	93	3			
1	C	77	Total	C	H	N	O	S	47	0	0
			1011	336	489	87	95	4			
1	D	75	Total	C	H	N	O	S	45	0	0
			969	324	466	80	95	4			

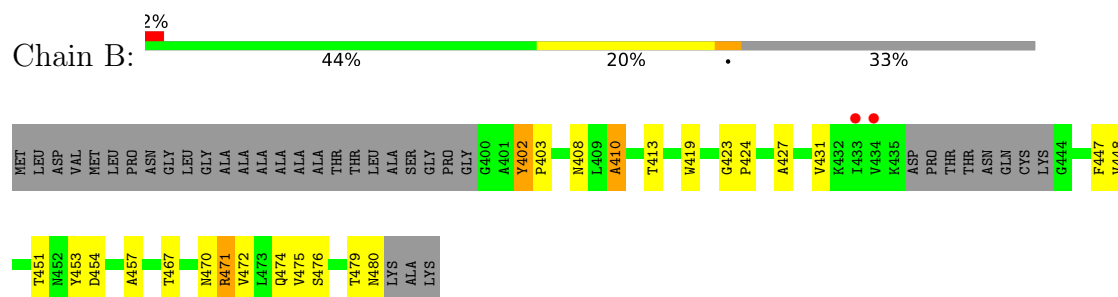
### 3 Residue-property plots [i](#)

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.

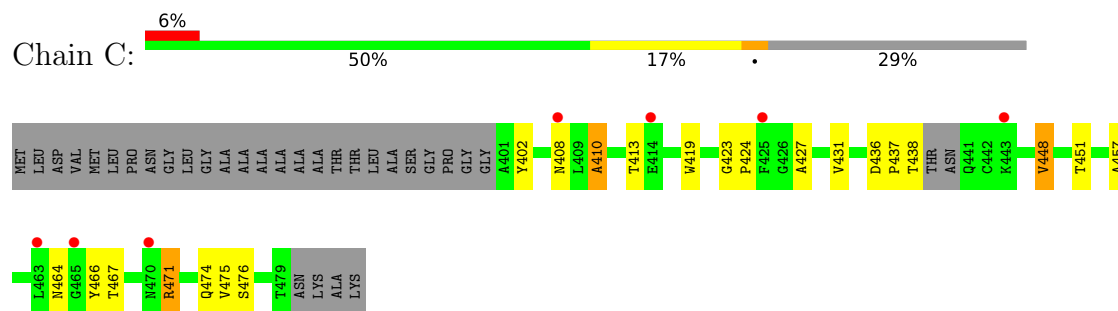
- Molecule 1: Protein elav



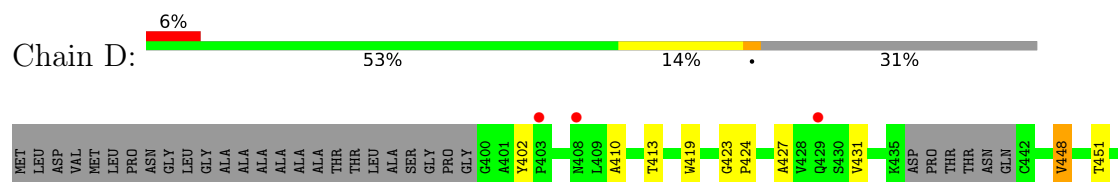
- Molecule 1: Protein elav

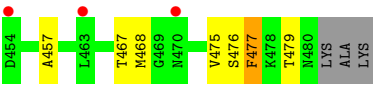


- Molecule 1: Protein elav



- Molecule 1: Protein elav





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.10Å 92.10Å 307.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	79.76 – 3.73 79.76 – 3.73	Depositor EDS
% Data completeness (in resolution range)	99.8 (79.76-3.73) 99.8 (79.76-3.73)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.75 (at 3.78Å)	Xtriage
Refinement program	REFMAC 5.8.0430 (refmacat 0.4.100)	Depositor
R, $R_{free}$	0.234 , 0.256 0.235 , 0.252	Depositor DCC
$R_{free}$ test set	413 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	133.4	Xtriage
Anisotropy	0.475	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.43 , 205.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	4056	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	138.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.10% 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	0.60	0/555	1.12	1/756 (0.1%)
1	B	0.58	0/536	1.16	2/729 (0.3%)
1	C	0.57	0/532	1.13	1/724 (0.1%)
1	D	0.56	0/512	1.06	1/699 (0.1%)
All	All	0.58	0/2135	1.12	5/2908 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
All	All	0	2

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	402	TYR	CB-CA-C	8.99	120.15	110.17
1	A	402	TYR	CB-CA-C	8.71	119.84	110.17
1	C	402	TYR	CB-CA-C	8.39	119.48	110.17
1	D	402	TYR	CB-CA-C	8.16	119.23	110.17
1	B	454	ASP	CA-CB-CG	5.23	117.83	112.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	471	ARG	Sidechain
1	C	471	ARG	Sidechain

## 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	542	510	488	18	0
1	B	524	500	479	19	0
1	C	522	489	451	11	0
1	D	503	466	430	12	0
All	All	2091	1965	1848	50	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 13.

All (50) 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:C:408:ASN:O	1:C:471:ARG:HD3	1.88	0.72
1:B:431:VAL:HG22	1:B:448:VAL:HG22	1.80	0.63
1:C:437:PRO:O	1:C:438:THR:C	2.42	0.63
1:C:431:VAL:HG22	1:C:448:VAL:HG22	1.83	0.61
1:A:453:TYR:HE1	1:B:453:TYR:HE1	1.49	0.61
1:D:413:THR:HG21	1:D:468:MET:HE3	1.84	0.59
1:A:431:VAL:HG22	1:A:448:VAL:HG22	1.84	0.59
1:C:423:GLY:N	1:C:424:PRO:CD	2.67	0.58
1:D:431:VAL:HG22	1:D:448:VAL:HG22	1.84	0.58
1:D:423:GLY:N	1:D:424:PRO:CD	2.66	0.58
1:B:423:GLY:N	1:B:424:PRO:CD	2.67	0.57
1:A:444:GLY:O	1:A:445:TYR:HB3	2.04	0.57
1:A:423:GLY:N	1:A:424:PRO:CD	2.67	0.56
1:C:427:ALA:HB3	1:C:451:THR:HB	1.87	0.56
1:A:427:ALA:HB3	1:A:451:THR:HB	1.88	0.56
1:B:427:ALA:HB3	1:B:451:THR:HB	1.88	0.54
1:A:413:THR:HG21	1:A:468:MET:HE3	1.88	0.54
1:D:427:ALA:HB3	1:D:451:THR:HB	1.90	0.53
1:A:453:TYR:HE1	1:B:453:TYR:CE1	2.27	0.52
1:A:445:TYR:C	1:A:445:TYR:CD1	2.88	0.52
1:D:477:PHE:O	1:D:479:THR:HG23	2.09	0.52
1:A:477:PHE:O	1:A:479:THR:HG23	2.12	0.50
1:A:475:VAL:O	1:A:476:SER:HB3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:475:VAL:O	1:D:476:SER:HB3	2.13	0.49
1:A:453:TYR:CE1	1:B:453:TYR:HE1	2.31	0.48
1:C:475:VAL:O	1:C:476:SER:HB3	2.13	0.48
1:B:408:ASN:O	1:B:471:ARG:HD3	2.15	0.47
1:B:410:ALA:HB3	1:B:413:THR:OG1	2.15	0.47
1:C:410:ALA:HB3	1:C:413:THR:OG1	2.15	0.47
1:B:475:VAL:O	1:B:476:SER:HB3	2.13	0.46
1:B:419:TRP:CD2	1:D:419:TRP:HB3	2.51	0.46
1:B:408:ASN:ND2	1:B:472:VAL:O	2.48	0.45
1:B:419:TRP:HB3	1:D:419:TRP:CD2	2.53	0.44
1:A:419:TRP:CD2	1:C:419:TRP:HB3	2.53	0.43
1:D:477:PHE:CD1	1:D:477:PHE:N	2.86	0.43
1:A:419:TRP:HB3	1:C:419:TRP:CD2	2.54	0.43
1:A:477:PHE:HB3	1:B:402:TYR:CZ	2.53	0.43
1:C:423:GLY:N	1:C:424:PRO:HD3	2.34	0.42
1:B:479:THR:OG1	1:B:480:ASN:N	2.53	0.42
1:A:402:TYR:CZ	1:B:403:PRO:HG2	2.55	0.42
1:A:408:ASN:O	1:A:471:ARG:HD3	2.19	0.42
1:B:423:GLY:N	1:B:424:PRO:HD3	2.35	0.42
1:D:423:GLY:N	1:D:424:PRO:HD3	2.34	0.42
1:B:419:TRP:CE3	1:D:419:TRP:HB3	2.55	0.41
1:A:470:ASN:O	1:A:471:ARG:HG3	2.21	0.41
1:D:477:PHE:N	1:D:477:PHE:HD1	2.19	0.41
1:B:470:ASN:O	1:B:471:ARG:HG3	2.21	0.41
1:C:464:ASN:OD1	1:C:464:ASN:O	2.39	0.41
1:A:447:PHE:CD1	1:A:447:PHE:C	2.99	0.41
1:B:447:PHE:CD1	1:B:447:PHE:C	2.99	0.41

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	71/109 (65%)	63 (89%)	6 (8%)	2 (3%)	4	27
1	B	69/109 (63%)	62 (90%)	5 (7%)	2 (3%)	3	26
1	C	73/109 (67%)	65 (89%)	5 (7%)	3 (4%)	2	20
1	D	71/109 (65%)	65 (92%)	4 (6%)	2 (3%)	4	27
All	All	284/436 (65%)	255 (90%)	20 (7%)	9 (3%)	3	25

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	445	TYR
1	C	457	ALA
1	D	457	ALA
1	A	457	ALA
1	B	457	ALA
1	C	436	ASP
1	D	410	ALA
1	B	410	ALA
1	C	410	ALA

### 5.3.2 Protein sidechains ⓘ

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	46/81 (57%)	42 (91%)	4 (9%)	9	34
1	B	45/81 (56%)	43 (96%)	2 (4%)	25	50
1	C	39/81 (48%)	35 (90%)	4 (10%)	7	29
1	D	38/81 (47%)	35 (92%)	3 (8%)	11	37
All	All	168/324 (52%)	155 (92%)	13 (8%)	12	37

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	433	ILE
1	A	445	TYR

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Mol	Chain	Res	Type
1	A	458	MET
1	A	467	THR
1	B	467	THR
1	B	474	GLN
1	C	448	VAL
1	C	466	TYR
1	C	467	THR
1	C	474	GLN
1	D	448	VAL
1	D	467	THR
1	D	477	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

### 5.7 Other polymers ⓘ

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, 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	75/109 (68%)	-0.12	1 (1%) 75 50	87, 121, 164, 185	0
1	B	73/109 (66%)	0.05	2 (2%) 56 35	79, 117, 161, 172	0
1	C	77/109 (70%)	0.60	7 (9%) 15 13	108, 151, 204, 261	0
1	D	75/109 (68%)	0.34	6 (8%) 18 15	97, 146, 187, 208	0
All	All	300/436 (68%)	0.22	16 (5%) 32 22	79, 134, 186, 261	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	436	ASP	5.4
1	C	470	ASN	5.3
1	D	470	ASN	3.8
1	D	403	PRO	3.8
1	B	433	ILE	3.6
1	D	408	ASN	3.3
1	C	463	LEU	3.0
1	C	414	GLU	2.9
1	C	465	GLY	2.9
1	B	434	VAL	2.5
1	D	454	ASP	2.5
1	C	408	ASN	2.5
1	C	443	LYS	2.3
1	D	463	LEU	2.3
1	C	425	PHE	2.3
1	D	429	GLN	2.2

### 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.