



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:36 AM UTC

PDB ID : 5H11 / pdb_00005h11
Title : Crystal structure of Zebrafish Sec10
Authors : Chen, J.; Yamagata, A.; Fukai, S.
Deposited on : 2016-10-07
Resolution : 2.73 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

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
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

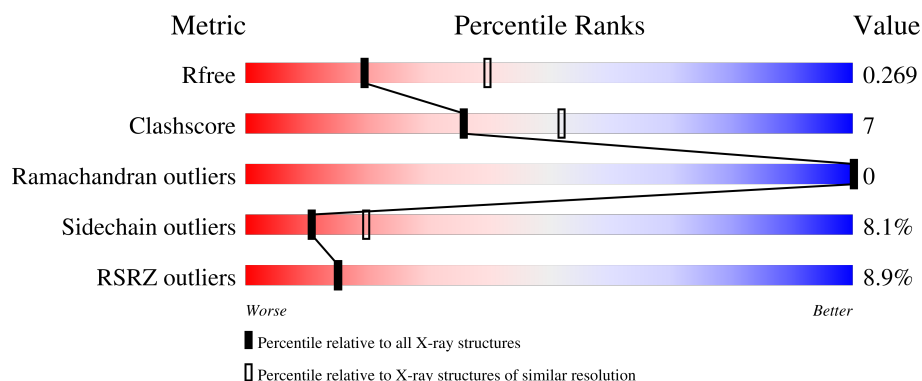
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.73 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	509	9% 76% 18% • •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BR	A	801	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BR	A	804	-	-	X	-
2	BR	A	806	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4003 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	494	Total	C	N	O	S	0	0	0
			3991	2519	700	745	27			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP B0BLY0
A	-3	PRO	-	expression tag	UNP B0BLY0
A	-2	LEU	-	expression tag	UNP B0BLY0
A	-1	GLY	-	expression tag	UNP B0BLY0
A	0	SER	-	expression tag	UNP B0BLY0
A	?	-	GLN	deletion	UNP B0BLY0
A	?	-	GLU	deletion	UNP B0BLY0
A	?	-	LEU	deletion	UNP B0BLY0
A	?	-	LYS	deletion	UNP B0BLY0
A	?	-	GLU	deletion	UNP B0BLY0
A	?	-	ARG	deletion	UNP B0BLY0
A	?	-	ILE	deletion	UNP B0BLY0
A	?	-	ARG	deletion	UNP B0BLY0
A	?	-	GLN	deletion	UNP B0BLY0
A	?	-	ARG	deletion	UNP B0BLY0

- Molecule 2 is BROMIDE ION (CCD ID: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	10	Total	Br	0	0
			10	10		

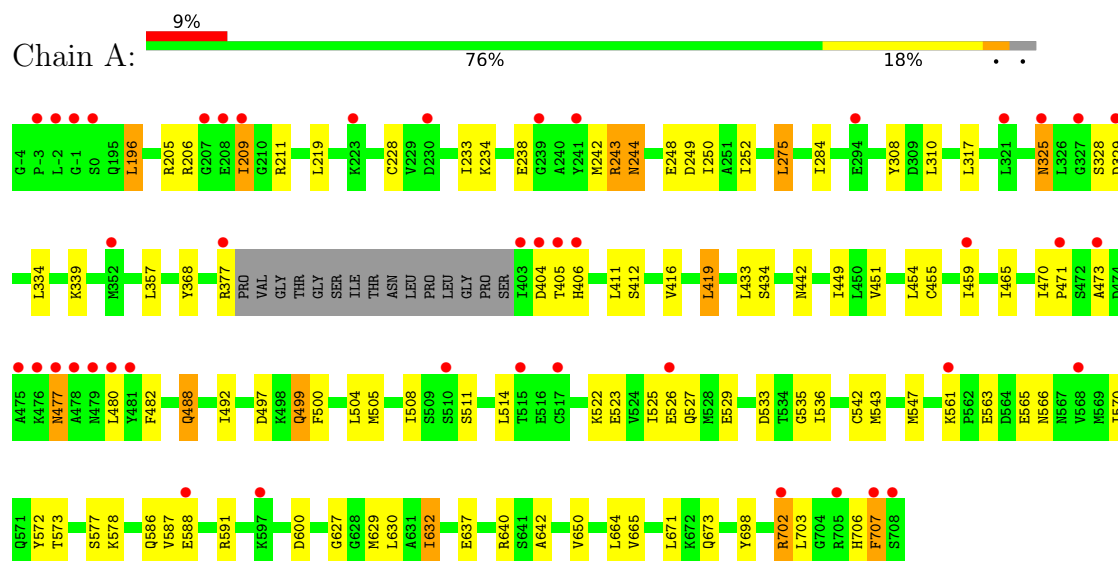
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	O	0	0
			2	2		

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: Uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	147.69Å 162.65Å 45.57Å 90.00° 95.55° 90.00°	Depositor
Resolution (Å)	46.92 – 2.73 46.92 – 2.73	Depositor EDS
% Data completeness (in resolution range)	95.2 (46.92-2.73) 95.2 (46.92-2.73)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.236 , 0.267 0.240 , 0.269	Depositor DCC
R_{free} test set	1384 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.902	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 23.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4003	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.73% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BR

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	0.28	0/4058	0.71	1/5470 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	196	LEU	N-CA-C	-6.03	105.90	113.20

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	3991	0	4001	58	0
2	A	10	0	0	10	0
3	A	2	0	0	0	0
All	All	4003	0	4001	58	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 (58) 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:A:205:ARG:NH1	2:A:808:BR:BR	2.48	1.01
1:A:328:SER:HB3	2:A:801:BR:BR	2.17	1.00
1:A:499:GLN:HB2	2:A:804:BR:BR	2.18	0.98
1:A:499:GLN:CB	2:A:804:BR:BR	2.84	0.80
1:A:328:SER:CB	2:A:801:BR:BR	2.85	0.79
1:A:416:VAL:HG11	1:A:492:ILE:HD12	1.64	0.77
1:A:632:ILE:HD11	1:A:665:VAL:HG21	1.70	0.73
1:A:637:GLU:OE1	1:A:640:ARG:NH1	2.28	0.66
1:A:473:ALA:HA	1:A:542:CYS:HA	1.78	0.64
1:A:209:ILE:HB	1:A:211:ARG:H	1.64	0.61
1:A:698:TYR:HA	1:A:703:LEU:HD12	1.82	0.61
1:A:572:TYR:HE1	1:A:577:SER:HB3	1.67	0.59
1:A:233:ILE:HG23	1:A:275:LEU:HD23	1.85	0.59
1:A:325:ASN:OD1	1:A:325:ASN:N	2.28	0.55
1:A:523:GLU:O	1:A:527:GLN:HG2	2.05	0.55
1:A:434:SER:OG	1:A:442:ASN:ND2	2.31	0.55
1:A:219:LEU:HD13	1:A:228:CYS:HB2	1.89	0.54
1:A:404:ASP:HB2	1:A:406:HIS:CD2	2.43	0.54
1:A:627:GLY:HA2	1:A:630:LEU:HD12	1.90	0.53
1:A:511:SER:HB3	1:A:514:LEU:HB2	1.92	0.52
1:A:416:VAL:HG21	1:A:492:ILE:HG23	1.90	0.52
1:A:308:TYR:HA	1:A:433:LEU:HD11	1.92	0.52
1:A:471:PRO:HG2	1:A:477:ASN:HB3	1.92	0.52
1:A:499:GLN:HB3	2:A:804:BR:BR	2.65	0.51
1:A:328:SER:HB2	2:A:801:BR:BR	2.66	0.51
1:A:449:ILE:CD1	2:A:806:BR:BR	3.14	0.51
1:A:488:GLN:O	1:A:492:ILE:HG12	2.11	0.51
1:A:642:ALA:HB1	1:A:650:VAL:HG12	1.93	0.50
1:A:455:CYS:HA	1:A:459:ILE:HG22	1.94	0.49
1:A:702:ARG:HA	1:A:702:ARG:HE	1.78	0.49
1:A:505:MET:HA	1:A:508:ILE:HG22	1.95	0.48
1:A:250:ILE:HD13	1:A:284:ILE:HD11	1.95	0.47
1:A:543:MET:O	1:A:547:MET:HG3	2.15	0.47
1:A:522:LYS:O	1:A:526:GLU:HG2	2.14	0.47
1:A:629:MET:SD	1:A:629:MET:N	2.87	0.46
1:A:449:ILE:HD13	2:A:806:BR:BR	2.71	0.46
1:A:368:TYR:CE1	1:A:412:SER:HB2	2.50	0.46
1:A:248:GLU:O	1:A:252:ILE:HG12	2.17	0.45
1:A:434:SER:HG	1:A:442:ASN:HD22	1.62	0.44
1:A:243:ARG:NH1	1:A:244:ASN:OD1	2.51	0.43
1:A:234:LYS:O	1:A:238:GLU:HG2	2.19	0.43
1:A:543:MET:HE3	1:A:586:GLN:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:482:PHE:CE2	1:A:535:GLY:HA3	2.54	0.43
1:A:533:ASP:HA	1:A:536:ILE:HG22	2.01	0.43
1:A:377:ARG:HH11	1:A:377:ARG:HA	1.84	0.43
1:A:419:LEU:HD23	1:A:454:LEU:HB2	2.01	0.42
1:A:206:ARG:HD2	1:A:206:ARG:HA	1.81	0.42
1:A:587:VAL:O	1:A:591:ARG:HG3	2.20	0.42
1:A:497:ASP:HA	1:A:500:PHE:HB3	2.02	0.41
1:A:572:TYR:CE1	1:A:577:SER:HB3	2.51	0.41
1:A:525:ILE:O	1:A:529:GLU:HB2	2.20	0.41
1:A:563:GLU:H	1:A:563:GLU:CD	2.28	0.41
1:A:411:LEU:O	1:A:488:GLN:NE2	2.53	0.41
1:A:339:LYS:HA	1:A:339:LYS:HD2	1.80	0.41
1:A:470:ILE:HA	1:A:471:PRO:HD3	1.77	0.41
1:A:449:ILE:HD12	2:A:806:BR:BR	2.76	0.40
1:A:451:VAL:O	1:A:455:CYS:HB2	2.20	0.40
1:A:707:PHE:HD1	1:A:707:PHE:HA	1.63	0.40

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	490/509 (96%)	464 (95%)	26 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	444/456 (97%)	408 (92%)	36 (8%)	11	20

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	196	LEU
1	A	209	ILE
1	A	242	MET
1	A	243	ARG
1	A	244	ASN
1	A	249	ASP
1	A	275	LEU
1	A	310	LEU
1	A	317	LEU
1	A	325	ASN
1	A	329	ASP
1	A	334	LEU
1	A	357	LEU
1	A	405	THR
1	A	419	LEU
1	A	465	ILE
1	A	477	ASN
1	A	480	LEU
1	A	488	GLN
1	A	499	GLN
1	A	504	LEU
1	A	561	LYS
1	A	565	GLU
1	A	566	ASN
1	A	570	ILE
1	A	573	THR
1	A	578	LYS
1	A	588	GLU
1	A	600	ASP
1	A	632	ILE
1	A	664	LEU
1	A	671	LEU
1	A	673	GLN
1	A	702	ARG
1	A	706	HIS

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Mol	Chain	Res	Type
1	A	707	PHE

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

Mol	Chain	Res	Type
1	A	221	HIS
1	A	406	HIS
1	A	458	HIS
1	A	501	ASN
1	A	546	GLN
1	A	706	HIS

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 ⓘ

Of 10 ligands modelled in this entry, 10 are monoatomic - leaving 0 for Mogul analysis.

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

6.1 Protein, DNA and RNA chains

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	494/509 (97%)	0.63	44 (8%) 15 15	34, 52, 92, 134	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	403	ILE	8.2
1	A	707	PHE	6.0
1	A	480	LEU	4.9
1	A	-3	PRO	3.9
1	A	230	ASP	3.8
1	A	477	ASN	3.8
1	A	510	SER	3.8
1	A	325	ASN	3.6
1	A	-2	LEU	3.4
1	A	478	ALA	3.4
1	A	404	ASP	3.4
1	A	479	ASN	3.4
1	A	405	THR	3.3
1	A	561	LYS	3.3
1	A	475	ALA	3.2
1	A	471	PRO	3.2
1	A	208	GLU	3.1
1	A	209	ILE	3.1
1	A	481	TYR	3.1
1	A	517	CYS	3.0
1	A	207	GLY	3.0
1	A	473	ALA	2.9
1	A	702	ARG	2.8
1	A	476	LYS	2.7
1	A	352	MET	2.6
1	A	-1	GLY	2.6
1	A	459	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	241	TYR	2.6
1	A	223	LYS	2.6
1	A	294	GLU	2.4
1	A	327	GLY	2.4
1	A	708	SER	2.3
1	A	705	ARG	2.2
1	A	321	LEU	2.2
1	A	329	ASP	2.2
1	A	377	ARG	2.2
1	A	597	LYS	2.2
1	A	406	HIS	2.1
1	A	588	GLU	2.1
1	A	239	GLY	2.1
1	A	0	SER	2.1
1	A	515	THR	2.1
1	A	526	GLU	2.0
1	A	568	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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	BR	A	808	1/1	0.64	0.29	209,209,209,209	0
2	BR	A	809	1/1	0.69	0.30	168,168,168,168	0
2	BR	A	810	1/1	0.77	0.32	183,183,183,183	0
2	BR	A	807	1/1	0.81	0.37	172,172,172,172	0
2	BR	A	801	1/1	0.87	0.18	108,108,108,108	0
2	BR	A	806	1/1	0.89	0.14	147,147,147,147	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BR	A	802	1/1	0.90	0.25	151,151,151,151	0
2	BR	A	803	1/1	0.91	0.19	137,137,137,137	0
2	BR	A	804	1/1	0.94	0.12	105,105,105,105	0
2	BR	A	805	1/1	0.96	0.11	100,100,100,100	0

6.5 Other polymers [i](#)

There are no such residues in this entry.