



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:53 PM UTC

PDB ID : 1SQH / pdb_00001sqh
Title : X-RAY STRUCTURE OF DROSOPHILA MALONOGASTER PROTEIN Q9VR51 NORTHEAST STRUCTURAL GENOMICS CONSORTIUM TARGET FR87.
Authors : Kuzin, A.P.; Chen, Y.; Forouhar, F.; Acton, T.B.; Xiao, R.; Cooper, B.; Ho, C.K.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2004-03-18
Resolution : 2.00 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)

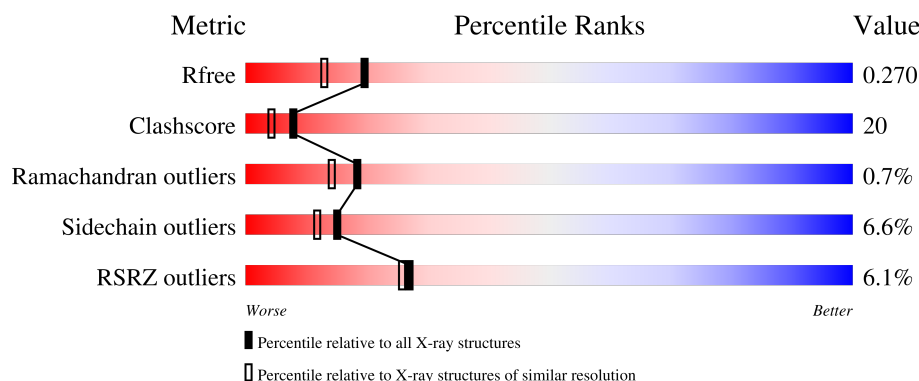
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

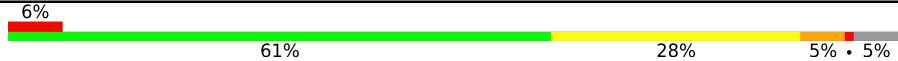
The reported resolution of this entry is 2.00 Å.

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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	312	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2557 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 hypothetical protein CG14615-PA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	Se	0	0	0
			2405	1530	431	435	5	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	155	MSE	MET	modified residue	UNP Q9VR51
A	186	MSE	MET	modified residue	UNP Q9VR51
A	236	MSE	MET	modified residue	UNP Q9VR51
A	256	MSE	MET	modified residue	UNP Q9VR51
A	305	LEU	-	expression tag	UNP Q9VR51
A	306	GLU	-	expression tag	UNP Q9VR51
A	307	HIS	-	expression tag	UNP Q9VR51
A	308	HIS	-	expression tag	UNP Q9VR51
A	309	HIS	-	expression tag	UNP Q9VR51
A	310	HIS	-	expression tag	UNP Q9VR51
A	311	HIS	-	expression tag	UNP Q9VR51
A	312	HIS	-	expression tag	UNP Q9VR51

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	152	Total	O	0	0
			152	152		

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: hypothetical protein CG14615-PA



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	82.81Å 82.81Å 279.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.84 – 2.00 19.84 – 2.00	Depositor EDS
% Data completeness (in resolution range)	69.5 (19.84-2.00) 84.4 (19.84-2.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 1.99Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.262 0.237 , 0.270	Depositor DCC
R_{free} test set	2144 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	25.3	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2557	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% 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

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	0/2455	0.95	12/3317 (0.4%)

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	A	0	1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	228	GLN	N-CA-C	-7.77	96.90	109.96
1	A	99	GLN	CA-C-N	7.34	129.02	119.84
1	A	99	GLN	C-N-CA	7.34	129.02	119.84
1	A	89	ILE	N-CA-C	-6.60	105.31	111.45
1	A	108	GLU	N-CA-C	-6.54	104.15	111.28
1	A	28	LYS	N-CA-C	6.04	117.54	111.07
1	A	174	PHE	N-CA-C	5.85	119.10	109.85
1	A	132	CYS	N-CA-C	5.82	118.12	111.02
1	A	59	LEU	N-CA-C	5.19	117.61	111.33
1	A	156	VAL	N-CA-C	5.11	116.07	108.46
1	A	18	VAL	N-CA-C	-5.06	105.56	110.42
1	A	201	TYR	N-CA-C	-5.02	105.72	111.14

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	278	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	2405	0	2386	95	0
2	A	152	0	0	6	0
All	All	2557	0	2386	95	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 20.

All (95) 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:245:ARG:NH2	1:A:278:ARG:NE	1.99	1.10
1:A:170:LEU:HD22	1:A:176:ILE:HG23	1.44	0.97
1:A:176:ILE:H	1:A:176:ILE:HD13	1.42	0.84
1:A:146:ALA:HB3	1:A:300:VAL:HG21	1.61	0.82
1:A:245:ARG:HH22	1:A:278:ARG:NE	1.76	0.79
1:A:223:ILE:HD12	1:A:240:LEU:HD23	1.67	0.75
1:A:245:ARG:NH2	1:A:278:ARG:CZ	2.50	0.74
1:A:245:ARG:HH22	1:A:278:ARG:CD	2.01	0.73
1:A:256:MSE:HE2	1:A:260:ILE:HG13	1.73	0.71
1:A:99:GLN:HE22	1:A:105:GLU:H	1.39	0.70
1:A:153:PHE:HB2	1:A:155:MSE:HE1	1.72	0.70
1:A:75:ASN:HB3	1:A:78:THR:HG23	1.75	0.69
1:A:165:ALA:O	1:A:258:ARG:HD3	1.94	0.68
1:A:210:LYS:H	1:A:228:GLN:HE21	1.43	0.65
1:A:245:ARG:NH2	1:A:278:ARG:HE	1.92	0.65
1:A:176:ILE:HG22	1:A:214:ILE:HD13	1.79	0.64
1:A:245:ARG:HD2	1:A:245:ARG:N	2.11	0.63
1:A:99:GLN:HB3	1:A:103:ALA:HA	1.81	0.63
1:A:22:LEU:HA	1:A:39:LEU:HD23	1.79	0.63
1:A:24:LEU:HD11	1:A:77:ARG:HA	1.82	0.61
1:A:58:ASP:O	1:A:62:ILE:HG23	2.01	0.61
1:A:44:LYS:HG2	1:A:48:GLN:HE21	1.64	0.61
1:A:121:GLY:HA2	1:A:299:LEU:O	2.01	0.60
1:A:170:LEU:CD2	1:A:176:ILE:HG23	2.26	0.59
1:A:180:ARG:HG3	1:A:182:GLU:HG2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:GLU:HB3	1:A:120:GLN:HB2	1.85	0.58
1:A:14:SER:O	1:A:18:VAL:HG23	2.05	0.57
1:A:134:ARG:HG3	2:A:384:HOH:O	2.06	0.56
1:A:245:ARG:CZ	1:A:278:ARG:HE	2.19	0.55
1:A:245:ARG:CZ	1:A:278:ARG:NE	2.68	0.55
1:A:21:LEU:HG	1:A:39:LEU:HD21	1.88	0.55
1:A:119:THR:HG23	1:A:302:ASN:CG	2.31	0.55
1:A:146:ALA:O	1:A:300:VAL:HG22	2.06	0.55
1:A:146:ALA:HB3	1:A:300:VAL:CG2	2.34	0.55
1:A:264:GLU:HG3	1:A:266:ILE:HG13	1.89	0.54
1:A:219:THR:HG23	1:A:221:GLU:H	1.73	0.53
1:A:226:ILE:HD13	1:A:227:PHE:N	2.23	0.53
1:A:153:PHE:CB	1:A:155:MSE:HE1	2.39	0.52
1:A:44:LYS:HG2	1:A:48:GLN:NE2	2.25	0.52
1:A:212:LEU:HB2	1:A:226:ILE:HG22	1.92	0.52
1:A:223:ILE:HA	1:A:240:LEU:HD22	1.92	0.51
1:A:223:ILE:HD11	1:A:242:LYS:HE2	1.92	0.51
1:A:84:SER:OG	1:A:94:SER:HB3	2.10	0.51
1:A:40:TYR:CE2	1:A:266:ILE:HG21	2.46	0.50
1:A:106:LEU:O	1:A:109:CYS:HB2	2.11	0.50
1:A:132:CYS:O	1:A:135:VAL:HG12	2.11	0.50
1:A:184:ALA:HB3	2:A:422:HOH:O	2.10	0.50
1:A:56:ARG:O	1:A:57:ASN:HB2	2.10	0.50
1:A:185:ALA:HB3	1:A:186:MSE:HE2	1.92	0.50
1:A:245:ARG:HH22	1:A:278:ARG:HD2	1.74	0.50
1:A:38:LEU:C	1:A:38:LEU:HD12	2.37	0.49
1:A:165:ALA:HB3	1:A:258:ARG:HG2	1.94	0.49
1:A:176:ILE:H	1:A:176:ILE:CD1	2.21	0.49
1:A:179:LEU:HD12	1:A:211:SER:HB2	1.95	0.49
1:A:210:LYS:HD3	1:A:259:GLU:HG2	1.95	0.49
1:A:176:ILE:HD13	1:A:176:ILE:N	2.19	0.48
1:A:31:ILE:HG23	1:A:267:THR:OG1	2.13	0.48
1:A:176:ILE:HG22	1:A:214:ILE:CD1	2.44	0.48
1:A:38:LEU:HD12	1:A:38:LEU:O	2.13	0.48
1:A:256:MSE:HE3	1:A:256:MSE:HA	1.95	0.48
1:A:210:LYS:H	1:A:228:GLN:NE2	2.07	0.48
1:A:302:ASN:HD22	1:A:302:ASN:HA	1.54	0.47
1:A:24:LEU:CD1	1:A:77:ARG:HA	2.43	0.46
1:A:26:LYS:HE3	1:A:31:ILE:CD1	2.46	0.46
1:A:210:LYS:HE3	2:A:327:HOH:O	2.16	0.46
1:A:119:THR:O	1:A:119:THR:CG2	2.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:LEU:HD11	1:A:64:LEU:HB3	1.97	0.45
1:A:218:ASP:CG	1:A:219:THR:H	2.25	0.45
1:A:38:LEU:HD11	1:A:66:LYS:HE2	1.99	0.45
1:A:256:MSE:HE3	1:A:259:GLU:HB3	1.99	0.44
1:A:273:VAL:HG12	1:A:275:THR:H	1.82	0.44
1:A:242:LYS:HG2	2:A:341:HOH:O	2.17	0.44
1:A:187:VAL:HG23	1:A:225:TRP:NE1	2.33	0.44
1:A:99:GLN:HA	1:A:100:PRO:HD3	1.83	0.44
1:A:146:ALA:HA	2:A:378:HOH:O	2.16	0.43
1:A:88:ASP:HB3	2:A:338:HOH:O	2.19	0.43
1:A:124:LEU:HB2	1:A:297:ILE:HB	2.00	0.43
1:A:184:ALA:O	1:A:187:VAL:HG13	2.18	0.43
1:A:107:TRP:CE2	1:A:134:ARG:HD3	2.54	0.43
1:A:155:MSE:CE	1:A:272:ILE:HG12	2.49	0.42
1:A:87:ARG:HH22	1:A:117:GLU:CD	2.27	0.42
1:A:301:PRO:O	1:A:302:ASN:C	2.61	0.42
1:A:20:GLU:OE2	1:A:77:ARG:NH1	2.52	0.42
1:A:96:PHE:CD2	1:A:97:SER:N	2.88	0.41
1:A:132:CYS:O	1:A:136:LYS:HG3	2.20	0.41
1:A:23:ASP:O	1:A:27:VAL:HG23	2.20	0.41
1:A:199:LEU:HD22	1:A:203:GLN:NE2	2.35	0.41
1:A:176:ILE:CD1	1:A:176:ILE:N	2.83	0.41
1:A:42:GLN:OE1	1:A:66:LYS:HG3	2.20	0.41
1:A:22:LEU:HD13	1:A:39:LEU:HB3	2.03	0.41
1:A:119:THR:O	1:A:119:THR:HG22	2.21	0.41
1:A:132:CYS:SG	1:A:295:GLU:HB2	2.61	0.41
1:A:245:ARG:NH2	1:A:278:ARG:CD	2.67	0.41
1:A:155:MSE:HE2	1:A:155:MSE:HB2	1.80	0.40
1:A:223:ILE:HA	1:A:240:LEU:CD2	2.51	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	295/312 (95%)	278 (94%)	15 (5%)	2 (1%)	18 14

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	302	ASN
1	A	218	ASP

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	256/267 (96%)	239 (93%)	17 (7%)	15 12

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ILE
1	A	38	LEU
1	A	53	GLN
1	A	87	ARG
1	A	124	LEU
1	A	155	MSE
1	A	169	ASP
1	A	176	ILE
1	A	187	VAL
1	A	199	LEU
1	A	200	THR
1	A	226	ILE
1	A	240	LEU
1	A	252	LEU
1	A	266	ILE
1	A	278	ARG
1	A	302	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	GLN
1	A	99	GLN
1	A	120	GLN
1	A	133	ASN
1	A	148	GLN
1	A	203	GLN
1	A	228	GLN
1	A	238	GLN
1	A	294	ASN
1	A	302	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, 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	293/312 (93%)	0.41	18 (6%)	27 26	16, 30, 49, 64	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	193	ASN	4.3
1	A	59	LEU	3.9
1	A	101	ASP	3.8
1	A	176	ILE	3.7
1	A	302	ASN	3.0
1	A	216	ARG	2.8
1	A	194	LYS	2.8
1	A	191	TRP	2.8
1	A	303	SER	2.5
1	A	192	PRO	2.5
1	A	54	ILE	2.4
1	A	223	ILE	2.4
1	A	55	PRO	2.3
1	A	60	ASN	2.2
1	A	195	GLY	2.1
1	A	57	ASN	2.1
1	A	31	ILE	2.1
1	A	277	TRP	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.