



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 06:28 AM BST

PDB ID : 6L5J
Title : Crystal structure of human rootletin 1108-1317
Authors : Kim, J.; Choi, H.J.
Deposited on : 2019-10-23
Resolution : 2.77 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

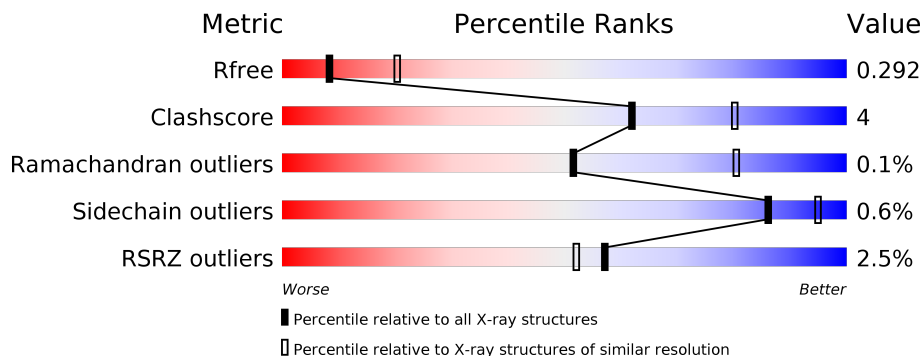
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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}	130704	4107 (2.80-2.76)
Clashscore	141614	4575 (2.80-2.76)
Ramachandran outliers	138981	4487 (2.80-2.76)
Sidechain outliers	138945	4489 (2.80-2.76)
RSRZ outliers	127900	4027 (2.80-2.76)

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 on 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	215	
1	B	215	
1	C	215	
1	D	215	

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 6665 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 Rootletin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	208	1678	994	344	334	1	5	0	0	0
1	B	208	1679	994	344	336	1	4	0	0	0
1	C	208	1679	994	344	336	1	4	0	0	0
1	D	201	1629	964	337	323	1	4	0	0	0

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1104	GLY	-	expression tag	UNP Q5TZA2
A	1105	ALA	-	expression tag	UNP Q5TZA2
A	1106	MSE	-	expression tag	UNP Q5TZA2
A	1107	GLU	-	expression tag	UNP Q5TZA2
A	1153	MSE	LEU	engineered mutation	UNP Q5TZA2
A	1166	MSE	LEU	engineered mutation	UNP Q5TZA2
A	1318	TRP	-	expression tag	UNP Q5TZA2
B	1104	GLY	-	expression tag	UNP Q5TZA2
B	1105	ALA	-	expression tag	UNP Q5TZA2
B	1106	MSE	-	expression tag	UNP Q5TZA2
B	1107	GLU	-	expression tag	UNP Q5TZA2
B	1153	MSE	LEU	engineered mutation	UNP Q5TZA2
B	1166	MSE	LEU	engineered mutation	UNP Q5TZA2
B	1318	TRP	-	expression tag	UNP Q5TZA2
C	1104	GLY	-	expression tag	UNP Q5TZA2
C	1105	ALA	-	expression tag	UNP Q5TZA2
C	1106	MSE	-	expression tag	UNP Q5TZA2
C	1107	GLU	-	expression tag	UNP Q5TZA2
C	1153	MSE	LEU	engineered mutation	UNP Q5TZA2
C	1166	MSE	LEU	engineered mutation	UNP Q5TZA2
C	1318	TRP	-	expression tag	UNP Q5TZA2

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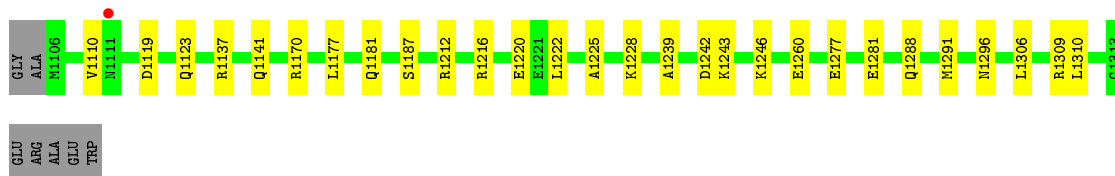
Chain	Residue	Modelled	Actual	Comment	Reference
D	1104	GLY	-	expression tag	UNP Q5TZA2
D	1105	ALA	-	expression tag	UNP Q5TZA2
D	1106	MSE	-	expression tag	UNP Q5TZA2
D	1107	GLU	-	expression tag	UNP Q5TZA2
D	1153	MSE	LEU	engineered mutation	UNP Q5TZA2
D	1166	MSE	LEU	engineered mutation	UNP Q5TZA2
D	1318	TRP	-	expression tag	UNP Q5TZA2

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.

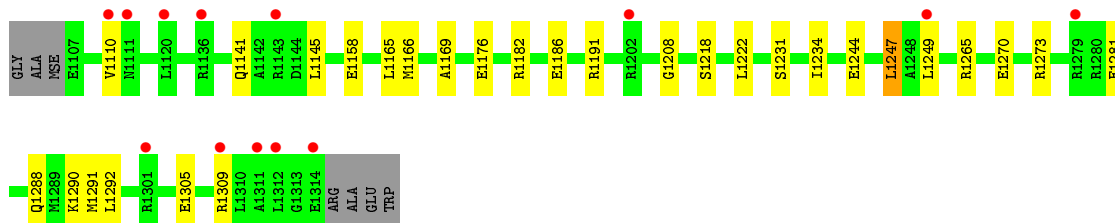
- Molecule 1: Rootletin

Chain A: 




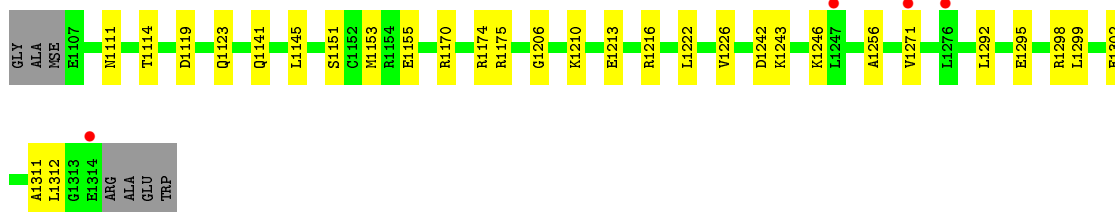
- Molecule 1: Rootletin

Chain B: 




- Molecule 1: Rootletin

Chain C: 



- Molecule 1: Rootletin

Chain D: 

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.39Å 68.00Å 79.77Å 92.62° 108.50° 113.84°	Depositor
Resolution (Å)	47.47 – 2.77 47.47 – 2.77	Depositor EDS
% Data completeness (in resolution range)	92.7 (47.47-2.77) 94.4 (47.47-2.77)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.77Å)	Xtrriage
Refinement program	PHENIX 1.17-3549, BUSTER 2.13.0	Depositor
R, R_{free}	0.283 , 0.290 0.285 , 0.292	Depositor DCC
R_{free} test set	1254 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	49.7	Xtrriage
Anisotropy	1.011	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 35.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.36$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6665	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.77% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.38	0/1674	0.61	0/2214
1	B	0.41	0/1675	0.61	1/2216 (0.0%)
1	C	0.42	0/1675	0.66	0/2216
1	D	0.43	0/1625	0.62	0/2148
All	All	0.41	0/6649	0.63	1/8794 (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	B	1290	LYS	CD-CE-NZ	-6.56	96.61	111.70

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	1678	0	1710	22	0
1	B	1679	0	1707	23	0
1	C	1679	0	1708	18	0
1	D	1629	0	1661	12	0
All	All	6665	0	6786	59	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 (59) 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:D:1157:GLU:HA	1:D:1160:ARG:HG2	1.75	0.67
1:C:1243:LYS:HE2	1:D:1244:GLU:HB3	1.78	0.66
1:C:1111:ASN:HA	1:C:1114:THR:HG22	1.79	0.64
1:A:1239:ALA:O	1:A:1243:LYS:HG2	2.02	0.60
1:C:1153:MSE:HE3	1:D:1152:CYS:SG	2.42	0.60
1:A:1170:ARG:HE	1:B:1166:MSE:SE	2.34	0.60
1:B:1305:GLU:O	1:B:1309:ARG:HG2	2.02	0.59
1:C:1170:ARG:HH11	1:C:1174:ARG:HH21	1.49	0.59
1:B:1182:ARG:NH1	1:B:1186:GLU:OE2	2.36	0.59
1:D:1166:MSE:HA	1:D:1169:ALA:HB3	1.85	0.58
1:A:1177:LEU:O	1:A:1181:GLN:HG3	2.07	0.55
1:C:1119:ASP:O	1:C:1123:GLN:HG2	2.07	0.55
1:C:1206:GLY:O	1:C:1210:LYS:HG2	2.08	0.54
1:C:1222:LEU:O	1:C:1226:VAL:HG23	2.08	0.54
1:B:1244:GLU:HA	1:B:1247:LEU:HD23	1.90	0.53
1:A:1110:VAL:HG22	1:B:1110:VAL:HG22	1.91	0.52
1:D:1192:GLU:HG2	1:D:1195:ARG:HH21	1.75	0.52
1:A:1137:ARG:O	1:A:1141:GLN:HG3	2.10	0.52
1:A:1225:ALA:HA	1:A:1228:LYS:HE3	1.91	0.52
1:A:1119:ASP:O	1:A:1123:GLN:HG3	2.10	0.51
1:B:1281:GLU:OE2	1:C:1210:LYS:NZ	2.36	0.51
1:A:1260:GLU:OE2	1:B:1265:ARG:NH2	2.45	0.50
1:D:1144:ASP:O	1:D:1148:GLN:HG3	2.11	0.50
1:B:1288:GLN:HA	1:B:1291:MSE:HE2	1.95	0.48
1:C:1295:GLU:O	1:C:1299:LEU:HG	2.13	0.48
1:B:1176:GLU:HG2	1:C:1311:ALA:HB1	1.96	0.48
1:D:1192:GLU:HG2	1:D:1195:ARG:NH2	2.28	0.48
1:A:1177:LEU:HD11	1:B:1176:GLU:HG3	1.94	0.48
1:C:1271:VAL:HG12	1:D:1271:VAL:HG12	1.96	0.48
1:A:1242:ASP:O	1:A:1246:LYS:HG2	2.13	0.47
1:B:1141:GLN:O	1:B:1145:LEU:HD13	2.14	0.47
1:A:1306:LEU:HD12	1:A:1309:ARG:NH1	2.30	0.47
1:B:1231:SER:HA	1:B:1234:ILE:HG22	1.97	0.46
1:A:1277:GLU:O	1:A:1281:GLU:HG3	2.16	0.46
1:A:1177:LEU:HD23	1:C:1312:LEU:HB3	1.97	0.46
1:C:1151:SER:O	1:C:1155:GLU:HG3	2.17	0.45
1:B:1270:GLU:O	1:B:1273:ARG:HB3	2.17	0.45
1:C:1298:ARG:HH12	1:C:1302:GLU:HB3	1.82	0.45
1:D:1161:THR:O	1:D:1164:ARG:HG2	2.16	0.44
1:A:1177:LEU:CD1	1:B:1176:GLU:HG3	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1187:SER:HA	1:B:1191:ARG:NH2	2.33	0.44
1:A:1296:ASN:ND2	1:B:1292:LEU:HD11	2.32	0.43
1:B:1165:LEU:O	1:B:1169:ALA:N	2.37	0.43
1:B:1288:GLN:HA	1:B:1291:MSE:CE	2.48	0.43
1:A:1310:LEU:O	1:A:1310:LEU:HG	2.18	0.43
1:C:1292:LEU:HB3	1:D:1292:LEU:HD22	1.99	0.43
1:A:1288:GLN:HA	1:A:1291:MSE:HE2	2.00	0.43
1:D:1216:ARG:O	1:D:1220:GLU:HG3	2.18	0.42
1:A:1212:ARG:HD3	1:B:1208:GLY:O	2.19	0.42
1:C:1213:GLU:OE2	1:C:1216:ARG:NH1	2.53	0.42
1:A:1306:LEU:HD12	1:A:1309:ARG:HH11	1.84	0.42
1:A:1216:ARG:NH1	1:A:1220:GLU:OE2	2.53	0.42
1:C:1141:GLN:O	1:C:1145:LEU:HD13	2.19	0.42
1:B:1218:SER:O	1:B:1222:LEU:HG	2.20	0.42
1:B:1244:GLU:CA	1:B:1247:LEU:HD23	2.50	0.41
1:B:1222:LEU:HD23	1:B:1222:LEU:HA	1.84	0.41
1:D:1165:LEU:O	1:D:1169:ALA:N	2.45	0.41
1:A:1222:LEU:HB3	1:B:1222:LEU:HB3	2.03	0.41
1:C:1242:ASP:OD1	1:C:1246:LYS:HE3	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	206/215 (96%)	203 (98%)	3 (2%)	0	100	100
1	B	206/215 (96%)	205 (100%)	1 (0%)	0	100	100
1	C	206/215 (96%)	204 (99%)	1 (0%)	1 (0%)	29	58
1	D	199/215 (93%)	199 (100%)	0	0	100	100
All	All	817/860 (95%)	811 (99%)	5 (1%)	1 (0%)	51	80

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1256	ALA

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	172/171 (101%)	172 (100%)	0	100	100
1	B	172/171 (101%)	169 (98%)	3 (2%)	60	85
1	C	172/171 (101%)	171 (99%)	1 (1%)	86	95
1	D	166/171 (97%)	166 (100%)	0	100	100
All	All	682/684 (100%)	678 (99%)	4 (1%)	86	95

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1158	GLU
1	B	1247	LEU
1	B	1249	LEU
1	C	1175	ARG

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

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 monosaccharides 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

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	203/215 (94%)	0.14	1 (0%) 91 90	46, 80, 105, 128	0
1	B	204/215 (94%)	0.29	13 (6%) 19 14	53, 87, 118, 139	0
1	C	204/215 (94%)	0.16	4 (1%) 65 61	44, 69, 104, 111	0
1	D	197/215 (91%)	0.26	2 (1%) 82 80	48, 71, 108, 137	0
All	All	808/860 (93%)	0.22	20 (2%) 57 52	44, 77, 109, 139	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1314	GLU	4.7
1	B	1314	GLU	4.6
1	B	1312	LEU	3.9
1	D	1111	ASN	3.6
1	B	1202	ARG	3.6
1	D	1311	ALA	3.4
1	A	1111	ASN	2.8
1	B	1136	ARG	2.7
1	B	1249	LEU	2.5
1	B	1279	ARG	2.5
1	C	1276	LEU	2.4
1	B	1143	ARG	2.3
1	B	1301	ARG	2.3
1	B	1309	ARG	2.3
1	B	1311	ALA	2.2
1	B	1111	ASN	2.1
1	B	1110	VAL	2.0
1	B	1120	LEU	2.0
1	C	1271	VAL	2.0
1	C	1247	LEU	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 monosaccharides 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.