



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 22, 2023 – 11:46 AM EDT

PDB ID : 2Z46
Title : Crystal Structure of Native-ORF134
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Deposited on : 2007-06-12
Resolution : 2.97 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

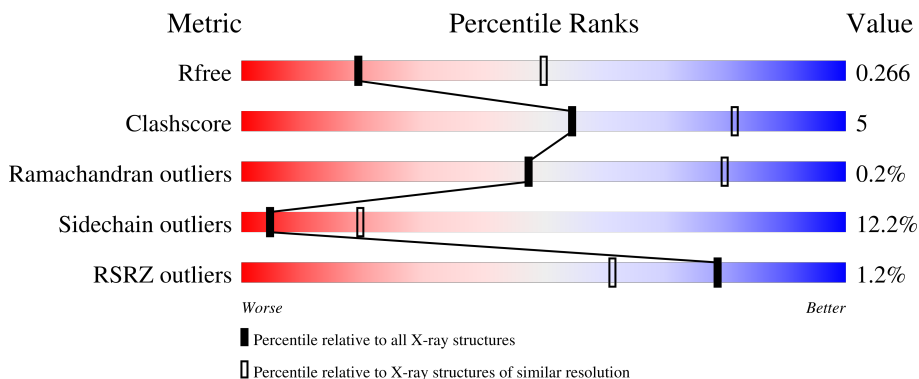
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.97 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	2754 (3.00-2.96)
Clashscore	141614	3103 (3.00-2.96)
Ramachandran outliers	138981	2993 (3.00-2.96)
Sidechain outliers	138945	2996 (3.00-2.96)
RSRZ outliers	127900	2644 (3.00-2.96)

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	134	
1	B	134	
1	C	134	
1	D	134	
1	E	134	

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Mol	Chain	Length	Quality of chain
1	F	134	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a small red segment at the beginning labeled '2%', a large green segment labeled '64%', a yellow segment labeled '16%', and a grey segment at the end labeled '17%'. There are two small black dots on the grey segment.</p>

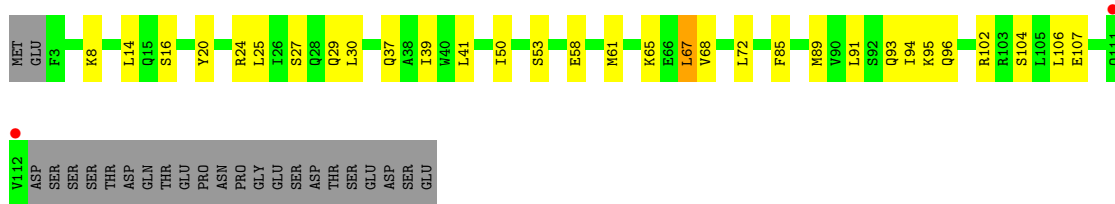
2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5347 atoms, of which 0 are hydrogens and 0 are deuteriums.

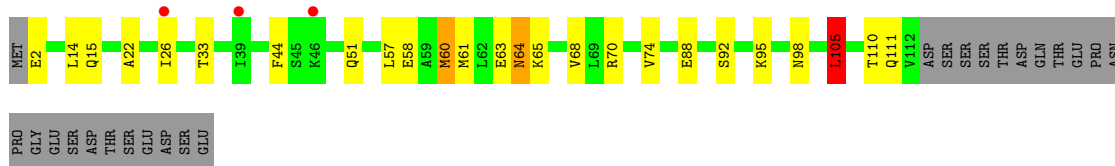
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called ORF134.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	109	886	562	154	167	3	0	0	0
1	B	109	886	562	154	167	3	0	0	0
1	C	109	886	562	155	166	3	0	0	0
1	D	110	894	567	155	168	4	0	0	0
1	E	110	893	567	156	167	3	0	0	0
1	F	111	902	572	157	170	3	0	0	0



● Molecule 1: ORF134



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	93.79Å 93.79Å 410.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.97 49.70 – 2.97	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.00-2.97) 99.7 (49.70-2.97)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 2.96Å)	Xtrriage
Refinement program	REFMAC 5.3.0034	Depositor
R, R_{free}	0.248 , 0.277 0.236 , 0.266	Depositor DCC
R_{free} test set	1955 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	74.3	Xtrriage
Anisotropy	0.078	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 25.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5347	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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	0.48	0/898	0.61	0/1210
1	B	0.50	0/898	0.62	0/1210
1	C	0.45	0/898	0.61	0/1210
1	D	0.45	0/906	0.63	1/1220 (0.1%)
1	E	0.45	0/905	0.60	0/1220
1	F	0.43	0/914	0.64	1/1232 (0.1%)
All	All	0.46	0/5419	0.62	2/7302 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	105	LEU	CA-CB-CG	7.00	131.39	115.30
1	D	105	LEU	CA-CB-CG	5.36	127.63	115.30

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	886	0	912	10	0
1	B	886	0	912	15	0
1	C	886	0	914	8	0
1	D	894	0	924	10	0
1	E	893	0	923	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	902	0	929	13	0
All	All	5347	0	5514	56	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 5.

All (56) 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:E:89:MET:O	1:E:93:GLN:HG2	1.90	0.70
1:E:104:SER:HA	1:E:107:GLU:HB3	1.76	0.66
1:B:28:GLN:O	1:B:31:SER:HB2	1.96	0.64
1:A:101:HIS:HE1	1:B:54:ASP:OD2	1.82	0.63
1:A:90:VAL:HG21	1:B:14:LEU:HD23	1.80	0.63
1:C:108:ARG:HA	1:C:111:GLN:HB3	1.80	0.63
1:A:80:GLU:HG3	1:B:91:LEU:HD11	1.84	0.59
1:C:44:PHE:CZ	1:C:60:MET:HB3	2.37	0.59
1:D:26:ILE:HG21	1:D:67:LEU:HD11	1.86	0.58
1:E:50:ILE:O	1:E:50:ILE:HG13	2.06	0.55
1:B:69:LEU:HB3	1:D:43:GLU:HG2	1.88	0.55
1:B:24:ARG:HG3	1:B:24:ARG:NH1	2.22	0.55
1:A:104:SER:O	1:A:108:ARG:HB2	2.09	0.52
1:F:44:PHE:CZ	1:F:60:MET:HA	2.45	0.51
1:B:24:ARG:HG3	1:B:24:ARG:HH11	1.76	0.51
1:F:64:ASN:C	1:F:64:ASN:HD22	2.14	0.51
1:F:61:MET:HB3	1:F:68:VAL:HG21	1.94	0.50
1:E:61:MET:HG3	1:F:105:LEU:HD22	1.94	0.49
1:D:43:GLU:OE2	1:D:47:ARG:NE	2.45	0.49
1:B:68:VAL:O	1:B:72:LEU:HG	2.13	0.48
1:B:70:ARG:HH11	1:B:70:ARG:HG3	1.79	0.48
1:E:16:SER:HA	1:E:53:SER:HB3	1.93	0.48
1:E:58:GLU:HA	1:E:61:MET:HE3	1.96	0.48
1:E:20:TYR:HB2	1:E:50:ILE:HG12	1.94	0.48
1:C:88:GLU:O	1:C:92:SER:HB3	2.14	0.47
1:B:53:SER:O	1:B:57:LEU:HG	2.15	0.46
1:A:65:LYS:H	1:A:65:LYS:HD3	1.80	0.46
1:F:22:ALA:HB1	1:F:74:VAL:HG23	1.97	0.46
1:D:1:MET:HG2	1:D:3:PHE:HD2	1.81	0.45
1:F:88:GLU:O	1:F:92:SER:HB2	2.17	0.45
1:E:61:MET:HG3	1:F:105:LEU:CD2	2.47	0.44
1:F:64:ASN:HD22	1:F:65:LYS:N	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:LEU:HD12	1:D:61:MET:HG2	2.00	0.43
1:A:14:LEU:HD12	1:A:14:LEU:HA	1.86	0.43
1:B:14:LEU:HD12	1:B:14:LEU:HA	1.83	0.43
1:C:83:LEU:HD13	1:D:87:PRO:HA	2.00	0.42
1:E:14:LEU:HD13	1:F:14:LEU:HD13	2.01	0.42
1:C:15:GLN:NE2	1:D:93:GLN:HB3	2.34	0.42
1:D:98:ASN:O	1:D:102:ARG:HG3	2.20	0.42
1:C:36:GLY:H	1:C:39:ILE:HG23	1.83	0.42
1:A:14:LEU:HD11	1:B:14:LEU:HD13	2.01	0.42
1:E:67:LEU:HA	1:E:67:LEU:HD12	1.67	0.42
1:A:105:LEU:HB2	1:B:61:MET:SD	2.60	0.42
1:D:50:ILE:HG13	1:D:50:ILE:O	2.19	0.41
1:E:85:PHE:O	1:E:89:MET:HB2	2.20	0.41
1:A:68:VAL:O	1:A:72:LEU:HG	2.20	0.41
1:A:73:THR:HG22	1:A:77:ASN:HD21	1.86	0.41
1:F:22:ALA:O	1:F:26:ILE:HG12	2.20	0.41
1:B:67:LEU:O	1:B:71:ILE:HG13	2.21	0.40
1:B:89:MET:HG2	1:B:93:GLN:NE2	2.35	0.40
1:C:14:LEU:HD23	1:D:90:VAL:HG21	2.02	0.40
1:E:68:VAL:O	1:E:72:LEU:HG	2.21	0.40
1:E:93:GLN:HB3	1:F:15:GLN:OE1	2.21	0.40
1:E:65:LYS:HB3	1:F:105:LEU:HD21	2.04	0.40
1:E:30:LEU:HD13	1:E:37:GLN:HE21	1.87	0.40
1:E:72:LEU:O	1:F:98:ASN:ND2	2.53	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	107/134 (80%)	100 (94%)	7 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	107/134 (80%)	103 (96%)	4 (4%)	0	100	100
1	C	107/134 (80%)	104 (97%)	3 (3%)	0	100	100
1	D	108/134 (81%)	104 (96%)	4 (4%)	0	100	100
1	E	108/134 (81%)	102 (94%)	5 (5%)	1 (1%)	17	53
1	F	109/134 (81%)	100 (92%)	9 (8%)	0	100	100
All	All	646/804 (80%)	613 (95%)	32 (5%)	1 (0%)	47	80

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	29	GLN

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	99/123 (80%)	84 (85%)	15 (15%)	3	12
1	B	99/123 (80%)	87 (88%)	12 (12%)	5	20
1	C	99/123 (80%)	87 (88%)	12 (12%)	5	20
1	D	100/123 (81%)	92 (92%)	8 (8%)	12	38
1	E	100/123 (81%)	87 (87%)	13 (13%)	4	17
1	F	101/123 (82%)	88 (87%)	13 (13%)	4	17
All	All	598/738 (81%)	525 (88%)	73 (12%)	5	20

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	13	THR
1	A	14	LEU
1	A	20	TYR
1	A	25	LEU

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Mol	Chain	Res	Type
1	A	50	ILE
1	A	52	GLU
1	A	53	SER
1	A	65	LYS
1	A	68	VAL
1	A	78	LEU
1	A	102	ARG
1	A	103	ARG
1	A	107	GLU
1	A	108	ARG
1	B	4	LYS
1	B	8	LYS
1	B	16	SER
1	B	20	TYR
1	B	24	ARG
1	B	25	LEU
1	B	31	SER
1	B	43	GLU
1	B	67	LEU
1	B	105	LEU
1	B	106	LEU
1	B	107	GLU
1	C	5	LYS
1	C	25	LEU
1	C	29	GLN
1	C	37	GLN
1	C	39	ILE
1	C	46	LYS
1	C	50	ILE
1	C	61	MET
1	C	62	LEU
1	C	68	VAL
1	C	92	SER
1	C	106	LEU
1	D	2	GLU
1	D	29	GLN
1	D	43	GLU
1	D	68	VAL
1	D	102	ARG
1	D	105	LEU
1	D	106	LEU
1	D	108	ARG

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Mol	Chain	Res	Type
1	E	8	LYS
1	E	24	ARG
1	E	25	LEU
1	E	27	SER
1	E	39	ILE
1	E	41	LEU
1	E	67	LEU
1	E	91	LEU
1	E	94	ILE
1	E	95	LYS
1	E	96	GLN
1	E	102	ARG
1	E	106	LEU
1	F	2	GLU
1	F	33	THR
1	F	51	GLN
1	F	57	LEU
1	F	58	GLU
1	F	60	MET
1	F	63	GLU
1	F	64	ASN
1	F	70	ARG
1	F	95	LYS
1	F	105	LEU
1	F	110	THR
1	F	111	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	77	ASN
1	A	101	HIS
1	B	28	GLN
1	B	37	GLN
1	B	48	HIS
1	B	93	GLN
1	B	101	HIS
1	C	15	GLN
1	C	101	HIS
1	D	34	ASN
1	D	37	GLN

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Mol	Chain	Res	Type
1	D	51	GLN
1	D	93	GLN
1	D	101	HIS
1	E	15	GLN
1	E	34	ASN
1	E	37	GLN
1	E	51	GLN
1	F	37	GLN
1	F	64	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 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 [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	109/134 (81%)	0.21	0 100 100	49, 62, 79, 87	0
1	B	109/134 (81%)	0.21	1 (0%) 84 69	48, 61, 86, 93	0
1	C	109/134 (81%)	0.06	0 100 100	48, 68, 89, 94	0
1	D	110/134 (82%)	0.23	2 (1%) 68 48	42, 54, 92, 102	0
1	E	110/134 (82%)	0.21	2 (1%) 68 48	47, 71, 103, 112	0
1	F	111/134 (82%)	0.44	3 (2%) 54 35	65, 91, 136, 137	0
All	All	658/804 (81%)	0.23	8 (1%) 79 61	42, 66, 110, 137	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	110	THR	4.7
1	E	112	VAL	4.7
1	D	1	MET	3.3
1	E	111	GLN	2.7
1	F	26	ILE	2.5
1	F	39	ILE	2.3
1	F	46	LYS	2.2
1	B	110	THR	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 monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.