



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 4REG / pdb_00004reg
Title : Crystal Structure Analysis of PF0642
Authors : He, J.H.; Zhou, H.; Liu, X.P.
Deposited on : 2014-09-23
Resolution : 2.49 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

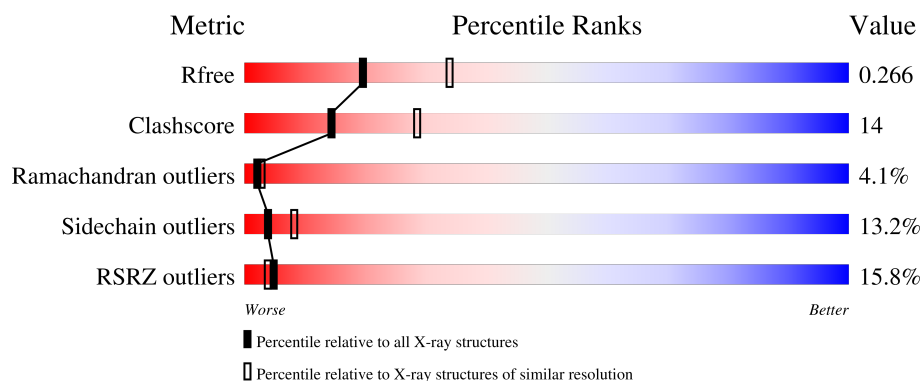
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	340	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4429 atoms, of which 2196 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	279	Total	C	H	N	O	S	0	0	0
			4385	1406	2196	374	405	4			

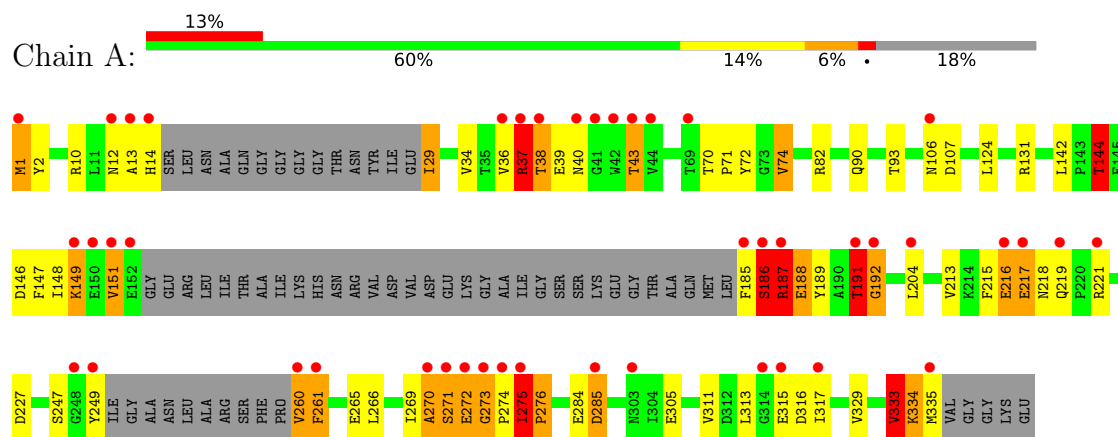
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	44	Total	O	0	0
			44	44		

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	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	79.26Å 79.26Å 136.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.37 – 2.49 45.37 – 2.49	Depositor EDS
% Data completeness (in resolution range)	99.4 (45.37-2.49) 99.6 (45.37-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.211 , 0.265 0.220 , 0.266	Depositor DCC
R_{free} test set	878 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 35.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4429	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% 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	1.06	5/2230 (0.2%)	1.21	23/3021 (0.8%)

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	2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	151	VAL	CA-CB	6.67	1.63	1.54
1	A	269	ILE	C-O	-5.80	1.16	1.23
1	A	261	PHE	CA-C	-5.55	1.45	1.52
1	A	144	THR	CB-CG2	-5.28	1.35	1.52
1	A	333	VAL	C-O	-5.01	1.18	1.24

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	285	ASP	N-CA-C	9.73	124.20	111.75
1	A	37	ARG	CA-C-N	8.53	137.05	121.70
1	A	37	ARG	C-N-CA	8.53	137.05	121.70
1	A	284	GLU	CA-C-N	7.52	131.37	120.38
1	A	284	GLU	C-N-CA	7.52	131.37	120.38
1	A	273	GLY	N-CA-C	-7.50	97.04	112.34
1	A	269	ILE	N-CA-C	7.02	119.06	108.80
1	A	271	SER	N-CA-C	-6.15	99.21	107.88
1	A	106	ASN	N-CA-C	5.95	120.27	112.89
1	A	221	ARG	CA-C-N	-5.84	113.74	119.76
1	A	221	ARG	C-N-CA	-5.84	113.74	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	37	ARG	CA-C-O	-5.79	114.34	121.06
1	A	192	GLY	N-CA-C	5.71	119.25	112.33
1	A	273	GLY	CA-C-N	-5.60	112.84	119.84
1	A	273	GLY	C-N-CA	-5.60	112.84	119.84
1	A	191	THR	N-CA-C	-5.51	99.06	110.80
1	A	275	ILE	CA-C-N	-5.36	113.14	119.84
1	A	275	ILE	C-N-CA	-5.36	113.14	119.84
1	A	215	PHE	N-CA-C	5.22	117.82	110.14
1	A	37	ARG	O-C-N	-5.16	117.29	123.27
1	A	247	SER	N-CA-C	5.05	116.86	111.36
1	A	142	LEU	CA-C-N	-5.00	114.61	119.76
1	A	142	LEU	C-N-CA	-5.00	114.61	119.76

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	260	VAL	Peptide
1	A	275	ILE	Peptide

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	2189	2196	2217	62	0
2	A	44	0	0	1	0
All	All	2233	2196	2217	62	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 14.

All (62) 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:271:SER:OG	1:A:272:GLU:OE2	1.85	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:HIS:HB2	1:A:260:VAL:CG2	2.04	0.87
1:A:14:HIS:HB2	1:A:260:VAL:HG23	1.61	0.82
1:A:273:GLY:O	1:A:275:ILE:N	2.18	0.76
1:A:186:SER:O	1:A:188:GLU:N	2.20	0.73
1:A:275:ILE:HG23	1:A:276:PRO:N	2.04	0.72
1:A:275:ILE:CG2	1:A:276:PRO:N	2.55	0.70
1:A:10:ARG:HH21	1:A:192:GLY:HA2	1.57	0.68
1:A:186:SER:O	1:A:189:TYR:N	2.26	0.66
1:A:333:VAL:C	1:A:335:MET:H	2.05	0.65
1:A:334:LYS:CG	1:A:335:MET:O	2.46	0.63
1:A:148:ILE:O	1:A:149:LYS:HB2	1.98	0.63
1:A:275:ILE:HG23	1:A:276:PRO:CD	2.30	0.61
1:A:2:TYR:HB2	1:A:271:SER:HA	1.83	0.61
1:A:12:ASN:HA	1:A:191:THR:HG22	1.84	0.60
1:A:334:LYS:HG2	1:A:335:MET:O	2.02	0.60
1:A:216:GLU:HG3	1:A:217:GLU:HA	1.83	0.60
1:A:37:ARG:HA	1:A:38:THR:OG1	2.03	0.58
1:A:14:HIS:HB2	1:A:260:VAL:HG21	1.85	0.56
1:A:13:ALA:C	1:A:14:HIS:CD2	2.84	0.55
1:A:275:ILE:HG22	1:A:276:PRO:O	2.06	0.55
1:A:334:LYS:HG3	1:A:335:MET:O	2.06	0.55
1:A:270:ALA:HB2	1:A:305:GLU:HB2	1.88	0.55
1:A:275:ILE:HG23	1:A:276:PRO:HD2	1.88	0.55
1:A:144:THR:HG22	1:A:147:PHE:H	1.72	0.55
1:A:10:ARG:NH2	1:A:192:GLY:HA2	2.22	0.54
1:A:1:MET:N	2:A:409:HOH:O	2.32	0.53
1:A:216:GLU:HA	1:A:218:ASN:N	2.25	0.52
1:A:29:ILE:HD12	1:A:260:VAL:HG22	1.91	0.52
1:A:271:SER:O	1:A:275:ILE:HD11	2.10	0.52
1:A:271:SER:CB	1:A:272:GLU:HA	2.40	0.52
1:A:186:SER:C	1:A:187:ARG:HG3	2.36	0.51
1:A:38:THR:OG1	1:A:43:THR:O	2.29	0.50
1:A:333:VAL:O	1:A:335:MET:N	2.44	0.50
1:A:271:SER:HB2	1:A:272:GLU:HA	1.94	0.50
1:A:273:GLY:C	1:A:275:ILE:N	2.70	0.49
1:A:329:VAL:O	1:A:333:VAL:HG13	2.13	0.49
1:A:265:GLU:HB2	1:A:311:VAL:HG13	1.94	0.49
1:A:185:PHE:C	1:A:187:ARG:HG3	2.37	0.49
1:A:273:GLY:H	1:A:275:ILE:HD13	1.77	0.49
1:A:216:GLU:CG	1:A:217:GLU:HA	2.43	0.49
1:A:273:GLY:O	1:A:275:ILE:HD13	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:275:ILE:CG2	1:A:276:PRO:O	2.63	0.47
1:A:273:GLY:C	1:A:275:ILE:H	2.23	0.46
1:A:151:VAL:O	1:A:151:VAL:HG23	2.16	0.45
1:A:13:ALA:O	1:A:14:HIS:CG	2.70	0.45
1:A:13:ALA:O	1:A:14:HIS:CD2	2.70	0.45
1:A:333:VAL:C	1:A:335:MET:N	2.69	0.43
1:A:13:ALA:C	1:A:14:HIS:CG	2.97	0.43
1:A:1:MET:HB2	1:A:272:GLU:HG3	1.99	0.43
1:A:70:THR:HB	1:A:71:PRO:HD2	2.00	0.43
1:A:216:GLU:HA	1:A:218:ASN:H	1.83	0.43
1:A:227:ASP:OD1	1:A:227:ASP:C	2.60	0.43
1:A:144:THR:CG2	1:A:146:ASP:CG	2.93	0.42
1:A:2:TYR:N	1:A:2:TYR:CD1	2.88	0.42
1:A:72:TYR:O	1:A:74:VAL:N	2.53	0.42
1:A:265:GLU:HB2	1:A:311:VAL:CG1	2.51	0.41
1:A:39:GLU:O	1:A:149:LYS:NZ	2.54	0.41
1:A:187:ARG:HH11	1:A:187:ARG:HD3	1.74	0.41
1:A:313:LEU:HD12	1:A:313:LEU:HA	1.94	0.41
1:A:147:PHE:C	1:A:148:ILE:O	2.60	0.40
1:A:266:LEU:C	1:A:266:LEU:HD12	2.45	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	271/340 (80%)	243 (90%)	17 (6%)	11 (4%)	2 3

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	THR

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Mol	Chain	Res	Type
1	A	187	ARG
1	A	334	LYS
1	A	40	ASN
1	A	186	SER
1	A	216	GLU
1	A	261	PHE
1	A	275	ILE
1	A	270	ALA
1	A	274	PRO
1	A	276	PRO

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	234/279 (84%)	203 (87%)	31 (13%)	4 8

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	29	ILE
1	A	34	VAL
1	A	36	VAL
1	A	37	ARG
1	A	43	THR
1	A	74	VAL
1	A	82	ARG
1	A	90	GLN
1	A	93	THR
1	A	107	ASP
1	A	124	LEU
1	A	131	ARG
1	A	144	THR
1	A	149	LYS
1	A	186	SER

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Mol	Chain	Res	Type
1	A	187	ARG
1	A	188	GLU
1	A	191	THR
1	A	204	LEU
1	A	213	VAL
1	A	217	GLU
1	A	219	GLN
1	A	249	TYR
1	A	272	GLU
1	A	275	ILE
1	A	285	ASP
1	A	315	GLU
1	A	316	ASP
1	A	317	ILE
1	A	333	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	HIS

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

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 ⓘ

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	279/340 (82%)	0.78	44 (15%) 5 4	22, 39, 84, 126	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	249	TYR	12.4
1	A	42	TRP	8.4
1	A	274	PRO	6.4
1	A	275	ILE	6.3
1	A	260	VAL	6.0
1	A	273	GLY	6.0
1	A	14	HIS	5.6
1	A	261	PHE	5.3
1	A	272	GLU	5.3
1	A	270	ALA	4.9
1	A	335	MET	4.9
1	A	185	PHE	4.8
1	A	315	GLU	4.7
1	A	152	GLU	4.6
1	A	151	VAL	4.6
1	A	216	GLU	4.5
1	A	44	VAL	4.1
1	A	43	THR	4.1
1	A	40	ASN	3.8
1	A	285	ASP	3.7
1	A	13	ALA	3.7
1	A	186	SER	3.6
1	A	38	THR	3.5
1	A	271	SER	3.4
1	A	217	GLU	3.3
1	A	149	LYS	3.0
1	A	36	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	1	MET	2.9
1	A	303	ASN	2.8
1	A	37	ARG	2.8
1	A	314	GLY	2.6
1	A	192	GLY	2.6
1	A	204	LEU	2.6
1	A	150	GLU	2.6
1	A	248	GLY	2.5
1	A	219	GLN	2.4
1	A	221	ARG	2.3
1	A	69	THR	2.3
1	A	12	ASN	2.2
1	A	187	ARG	2.3
1	A	191	THR	2.2
1	A	106	ASN	2.2
1	A	317	ILE	2.1
1	A	41	GLY	2.1

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.