



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 04:48 PM GMT

PDB ID : 1GCM
Title : GCN4 LEUCINE ZIPPER CORE MUTANT P-LI
Authors : Harbury, P.B.; Kim, P.S.; Alber, T.
Deposited on : 1995-04-25
Resolution : 1.80 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

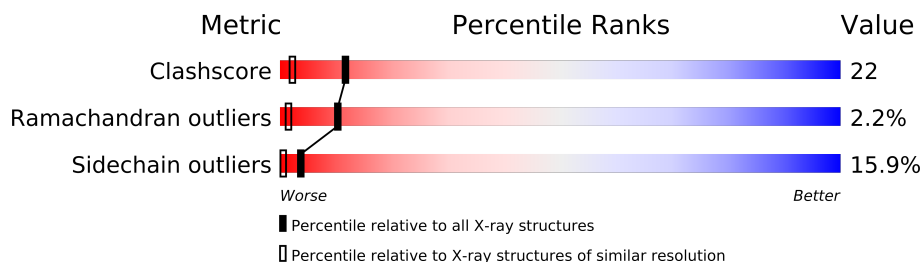
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.80 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))
Clashscore	79885	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (1.80-1.80)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	34	
1	B	34	
1	C	34	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 862 atoms, of which 0 are hydrogen and 0 are deuterium.

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 GCN4P-II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	32	Total	C	N	O	S	0	0	0
			265	170	46	48	1			
1	B	33	Total	C	N	O	S	0	0	0
			270	173	47	49	1			
1	C	34	Total	C	N	O	S	0	0	0
			279	178	48	52	1			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	ILE	LEU	ENGINEERED	UNP P03069
A	9	ILE	VAL	ENGINEERED	UNP P03069
A	12	ILE	LEU	ENGINEERED	UNP P03069
A	16	ILE	ASN	ENGINEERED	UNP P03069
A	19	ILE	LEU	ENGINEERED	UNP P03069
A	23	ILE	VAL	ENGINEERED	UNP P03069
A	26	ILE	LEU	ENGINEERED	UNP P03069
A	30	ILE	VAL	ENGINEERED	UNP P03069
B	5	ILE	LEU	ENGINEERED	UNP P03069
B	9	ILE	VAL	ENGINEERED	UNP P03069
B	12	ILE	LEU	ENGINEERED	UNP P03069
B	16	ILE	ASN	ENGINEERED	UNP P03069
B	19	ILE	LEU	ENGINEERED	UNP P03069
B	23	ILE	VAL	ENGINEERED	UNP P03069
B	26	ILE	LEU	ENGINEERED	UNP P03069
B	30	ILE	VAL	ENGINEERED	UNP P03069
C	5	ILE	LEU	ENGINEERED	UNP P03069
C	9	ILE	VAL	ENGINEERED	UNP P03069
C	12	ILE	LEU	ENGINEERED	UNP P03069
C	16	ILE	ASN	ENGINEERED	UNP P03069
C	19	ILE	LEU	ENGINEERED	UNP P03069
C	23	ILE	VAL	ENGINEERED	UNP P03069
C	26	ILE	LEU	ENGINEERED	UNP P03069

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Chain	Residue	Modelled	Actual	Comment	Reference
C	30	ILE	VAL	ENGINEERED	UNP P03069

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	15	Total 15	O 15	0	0
2	B	11	Total 11	O 11	0	0
2	C	22	Total 22	O 22	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

• Molecule 1: GCN4P-II

Chain A: 



• Molecule 1: GCN4P-II

Chain B: 



• Molecule 1: GCN4P-II

Chain C: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	35.32Å 39.67Å 36.97Å 90.00° 98.38° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.80)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.175 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	862	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	1.11	1/264 (0.4%)	1.41	1/349 (0.3%)
1	B	1.22	6/269 (2.2%)	1.44	3/356 (0.8%)
1	C	1.22	4/278 (1.4%)	1.46	4/368 (1.1%)
All	All	1.19	11/811 (1.4%)	1.43	8/1073 (0.7%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	20	GLU	CD-OE1	8.32	1.34	1.25
1	A	20	GLU	CD-OE1	6.12	1.32	1.25
1	B	6	GLU	CD-OE1	6.12	1.32	1.25
1	C	22	GLU	CD-OE2	5.98	1.32	1.25
1	B	10	GLU	CD-OE2	5.93	1.32	1.25
1	C	10	GLU	CD-OE2	5.60	1.31	1.25
1	C	11	GLU	CD-OE1	5.55	1.31	1.25
1	B	22	GLU	CD-OE2	-5.43	1.19	1.25
1	B	22	GLU	CD-OE1	5.28	1.31	1.25
1	C	20	GLU	CD-OE1	5.27	1.31	1.25
1	B	11	GLU	CD-OE1	5.14	1.31	1.25

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	0	ACE	C-N-CA	10.17	147.12	121.70
1	C	1	ARG	NE-CZ-NH2	-7.68	116.46	120.30
1	C	7	ASP	CB-CG-OD1	-6.69	112.28	118.30
1	C	1	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	B	25	ARG	NE-CZ-NH2	-5.78	117.41	120.30
1	B	25	ARG	NE-CZ-NH1	5.77	123.19	120.30
1	B	0	ACE	C-N-CA	5.67	135.88	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	7	ASP	CB-CG-OD2	5.61	123.35	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	265	0	289	15	2
1	B	270	0	292	22	0
1	C	279	0	298	12	0
2	A	15	0	0	0	0
2	B	11	0	0	3	0
2	C	22	0	0	1	2
All	All	862	0	879	37	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 22.

All (37) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:30:ILE:HG13	1:A:31:GLY:H	1.24	0.98
1:A:1:ARG:HA	1:A:1:ARG:HH11	1.36	0.88
1:A:1:ARG:HA	1:A:1:ARG:NH1	1.97	0.79
1:B:4:GLN:HG2	1:B:8:LYS:NZ	1.98	0.78
1:B:15:LYS:HD2	2:B:43:HOH:O	1.83	0.78
1:B:11:GLU:CD	1:B:15:LYS:HE3	2.05	0.76
1:B:18:HIS:HA	2:B:35:HOH:O	1.85	0.76
1:A:2:MET:O	1:A:6:GLU:HG3	1.86	0.74
1:B:4:GLN:HG2	1:B:8:LYS:HZ3	1.56	0.69
1:C:4:GLN:HE21	1:C:4:GLN:HA	1.58	0.68
2:B:34:HOH:O	1:C:20:GLU:HG2	1.93	0.68
1:B:11:GLU:OE1	1:C:13:LEU:HD21	1.95	0.67
1:B:25:ARG:NE	1:C:32:GLU:OE2	2.28	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:2:MET:HE1	1:B:2:MET:HE3	1.77	0.65
1:B:2:MET:HE2	1:B:2:MET:HA	1.78	0.65
1:A:30:ILE:HG13	1:A:31:GLY:N	2.04	0.62
1:B:2:MET:HE3	1:C:2:MET:CE	2.33	0.59
1:A:30:ILE:CG1	1:A:31:GLY:H	2.08	0.55
1:A:13:LEU:HD11	1:C:8:LYS:HE2	1.90	0.53
1:A:2:MET:HE1	1:B:2:MET:CE	2.38	0.53
1:B:28:LYS:O	1:B:31:GLY:O	2.27	0.53
1:B:25:ARG:O	1:B:29:LEU:HG	2.09	0.53
1:B:5:ILE:HG22	1:B:6:GLU:N	2.25	0.52
1:B:2:MET:HE3	1:C:2:MET:HE3	1.93	0.51
1:A:3:LYS:NZ	1:A:7:ASP:OD2	2.29	0.50
1:B:4:GLN:HG2	1:B:8:LYS:HZ2	1.74	0.50
1:C:4:GLN:NE2	1:C:4:GLN:HA	2.26	0.48
1:A:5:ILE:HG23	1:B:9:ILE:HD11	1.93	0.48
1:A:25:ARG:HH12	1:B:27:LYS:NZ	2.12	0.48
1:B:2:MET:HE3	1:C:2:MET:HE2	1.95	0.47
1:C:20:GLU:OE1	2:C:37:HOH:O	2.20	0.47
1:A:5:ILE:HG23	1:B:9:ILE:CD1	2.46	0.46
1:C:9:ILE:HG21	1:C:9:ILE:HD13	1.70	0.46
1:B:3:LYS:O	1:B:3:LYS:HG3	2.16	0.44
1:A:2:MET:CE	1:B:2:MET:CE	2.98	0.41
1:C:29:LEU:HA	1:C:29:LEU:HD23	1.78	0.41
1:A:1:ARG:NH1	1:A:4:GLN:NE2	2.69	0.41

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:30:ILE:C	2:C:55:HOH:O[1_656]	2.09	0.11
1:A:30:ILE:O	2:C:55:HOH:O[1_656]	2.18	0.02

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	30/34 (88%)	28 (93%)	1 (3%)	1 (3%)	6	0
1	B	31/34 (91%)	29 (94%)	2 (6%)	0	100	100
1	C	32/34 (94%)	30 (94%)	1 (3%)	1 (3%)	7	0
All	All	93/102 (91%)	87 (94%)	4 (4%)	2 (2%)	10	1

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1	ARG
1	A	30	ILE

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	29/31 (94%)	27 (93%)	2 (7%)	22	7
1	B	29/31 (94%)	23 (79%)	6 (21%)	2	0
1	C	30/31 (97%)	24 (80%)	6 (20%)	2	0
All	All	88/93 (95%)	74 (84%)	14 (16%)	4	0

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	15	LYS
1	B	2	MET
1	B	5	ILE
1	B	8	LYS
1	B	10	GLU
1	B	25	ARG
1	B	28	LYS
1	C	1	ARG
1	C	2	MET
1	C	3	LYS
1	C	4	GLN
1	C	8	LYS

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Mol	Chain	Res	Type
1	C	28	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	B	21	ASN
1	C	4	GLN
1	C	21	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.