



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:47 PM UTC

PDB ID : 4G6L / pdb_00004g6l
Title : Crystal structure of human CDK8/CYCC in the DMG-in conformation
Authors : Schneider, E.V.; Blaesse, M.; Huber, R.; Maskos, K.
Deposited on : 2012-07-19
Resolution : 2.70 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)
Validation Pipeline (wwPDB-VP) : 2.49

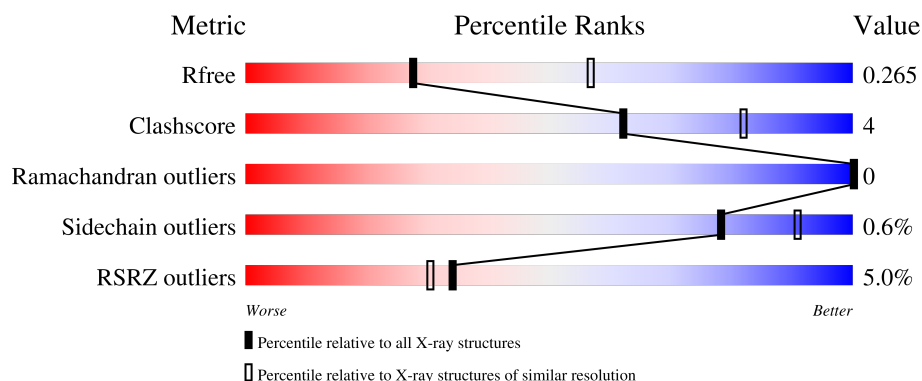
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	406	5% 76% 7% 17%
2	B	286	3% 84% 9% 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5047 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 Cyclin-dependent kinase 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	S	138	0	0
			2779	1791	483	488	17			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ASP	-	expression tag	UNP P49336
A	-1	ASP	-	expression tag	UNP P49336
A	0	LYS	-	expression tag	UNP P49336

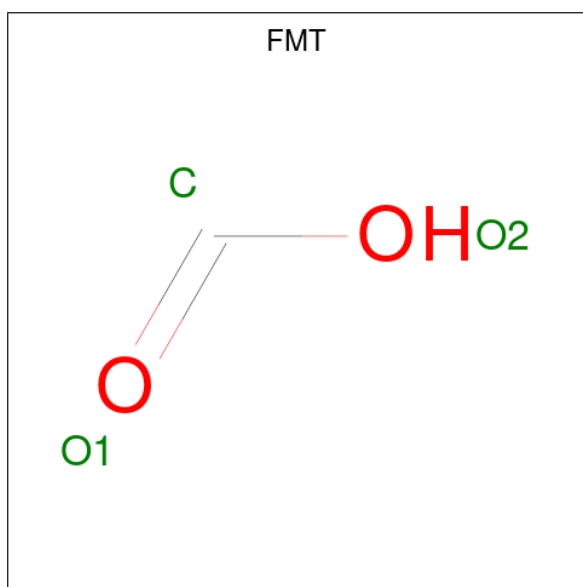
- Molecule 2 is a protein called Cyclin-C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	267	Total	C	N	O	S	72	0	0
			2225	1460	363	385	17			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	ASP	-	expression tag	UNP P24863
B	-1	LYS	-	expression tag	UNP P24863
B	0	ALA	-	expression tag	UNP P24863

- Molecule 3 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 3 1 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	19	Total O 19 19	0	0
4	B	21	Total O 21 21	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.18Å 70.94Å 170.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	85.21 – 2.70 85.21 – 2.70	Depositor EDS
% Data completeness (in resolution range)	(Not available) (85.21-2.70) 99.9 (85.21-2.70)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.212 , 0.266 0.214 , 0.265	Depositor DCC
R_{free} test set	844 reflections (3.50%)	wwPDB-VP
Wilson B-factor (Å ²)	47.4	Xtriage
Anisotropy	0.323	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.027 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5047	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% 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

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

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	0.66	0/2852	0.82	0/3839
2	B	0.65	0/2284	0.85	0/3094
All	All	0.65	0/5136	0.83	0/6933

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2779	0	2768	21	0
2	B	2225	0	2261	20	0
3	A	3	0	1	0	0
4	A	19	0	0	0	0
4	B	21	0	0	0	0
All	All	5047	0	5030	39	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 (39) 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)
2:B:72:PHE:HA	2:B:145:LEU:HD23	1.70	0.74
2:B:193:PRO:HD2	2:B:196:MET:HE2	1.76	0.67
2:B:72:PHE:HA	2:B:145:LEU:CD2	2.29	0.62
1:A:139:ILE:HD11	1:A:152:LEU:HD11	1.84	0.60
1:A:132:LEU:HD22	1:A:224:PHE:CD1	2.43	0.53
2:B:40:LEU:HD23	2:B:250:MET:HE1	1.90	0.53
1:A:149:HIS:CD2	1:A:174:MET:HG3	2.45	0.51
1:A:104:LEU:HD21	1:A:131:LEU:HD22	1.92	0.50
1:A:139:ILE:HG13	1:A:220:ILE:HD11	1.93	0.50
1:A:127:MET:HG3	1:A:345:VAL:HG13	1.95	0.49
1:A:123:LEU:HD23	1:A:127:MET:HE3	1.95	0.48
1:A:139:ILE:CD1	1:A:152:LEU:HD11	2.42	0.48
2:B:166:MET:HE3	2:B:202:LEU:HD21	1.94	0.48
2:B:76:TYR:CD2	2:B:145:LEU:HD22	2.47	0.48
2:B:76:TYR:CE2	2:B:145:LEU:HD13	2.48	0.48
1:A:100:ALA:HB3	1:A:158:LEU:HD13	1.96	0.47
1:A:3:TYR:O	1:A:7:VAL:HG22	2.14	0.47
1:A:280:MET:HA	1:A:280:MET:HE2	1.96	0.47
1:A:279:LEU:HD12	1:A:283:PHE:CD2	2.50	0.47
1:A:61:MET:HE3	1:A:177:ALA:HB1	1.96	0.46
2:B:179:ILE:O	2:B:183:THR:HG23	2.15	0.46
1:A:69:LEU:HD21	2:B:0:ALA:HB1	1.97	0.46
1:A:104:LEU:HD21	1:A:131:LEU:HD13	1.97	0.46
1:A:233:ILE:HG23	1:A:234:PHE:CD2	2.51	0.45
2:B:72:PHE:CZ	2:B:81:ILE:HD12	2.50	0.44
1:A:129:LYS:NZ	1:A:337:GLU:O	2.50	0.44
1:A:91:ARG:NH2	2:B:136:LEU:HD13	2.33	0.44
1:A:127:MET:HG3	1:A:345:VAL:CG1	2.48	0.44
2:B:85:LEU:HD12	2:B:135:ILE:HD13	1.99	0.44
2:B:150:LEU:O	2:B:152:VAL:HG23	2.18	0.43
2:B:51:LEU:HD21	2:B:112:ALA:HA	2.01	0.43
1:A:132:LEU:HD22	1:A:224:PHE:CG	2.55	0.42
2:B:40:LEU:HD23	2:B:250:MET:CE	2.48	0.42
2:B:172:LEU:HD23	2:B:209:GLN:OE1	2.18	0.41
2:B:72:PHE:CG	2:B:86:MET:HE3	2.55	0.41
2:B:166:MET:HE2	2:B:216:TRP:CE3	2.56	0.41
2:B:76:TYR:HE2	2:B:145:LEU:HD13	1.84	0.41
2:B:85:LEU:CD1	2:B:135:ILE:HD13	2.50	0.41
1:A:62:SER:HA	1:A:177:ALA:HB2	2.04	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	328/406 (81%)	317 (97%)	11 (3%)	0	100	100
2	B	265/286 (93%)	263 (99%)	2 (1%)	0	100	100
All	All	593/692 (86%)	580 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	299/361 (83%)	298 (100%)	1 (0%)	86	94
2	B	244/260 (94%)	242 (99%)	2 (1%)	73	88
All	All	543/621 (87%)	540 (99%)	3 (1%)	78	91

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	CYS
2	B	137	GLU
2	B	160	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	HIS
1	A	313	GLN
2	B	4	ASN
2	B	7	GLN
2	B	164	GLN

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](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	FMT	A	501	-	2,2,2	0.59	0	1,1,1	0.29	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	336/406 (82%)	0.32	22 (6%)	25 22	25, 46, 78, 111	34 (10%)
2	B	267/286 (93%)	0.10	8 (2%)	52 49	24, 46, 71, 91	20 (7%)
All	All	603/692 (87%)	0.22	30 (4%)	34 30	24, 46, 76, 111	54 (8%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	246	PRO	4.8
1	A	34	HIS	4.6
1	A	43	GLY	4.3
1	A	114	SER	4.2
1	A	359	LEU	4.1
1	A	350	GLN	3.6
2	B	38	TRP	3.6
2	B	220	LEU	3.3
1	A	32	TYR	3.2
1	A	-2	ASP	3.1
1	A	29	ARG	3.0
1	A	-1	ASP	3.0
1	A	179	LEU	3.0
1	A	180	PHE	2.9
1	A	124	PRO	2.9
1	A	239	GLU	2.8
2	B	215	GLN	2.6
2	B	-2	ASP	2.4
1	A	358	PHE	2.4
1	A	0	LYS	2.4
2	B	102	VAL	2.4
1	A	31	THR	2.3
2	B	221	SER	2.3
2	B	264	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	44	LYS	2.2
1	A	61	MET	2.2
1	A	47	LYS	2.2
1	A	123	LEU	2.1
2	B	99	GLU	2.1
1	A	286	ASN	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	FMT	A	501	3/3	0.77	0.22	51,51,51,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.