



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:06 PM UTC

PDB ID : 6QB6 / pdb_00006qb6
Title : Mcl1 in complex with a Fab
Authors : Hargreaves, D.
Deposited on : 2018-12-20
Resolution : 2.24 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

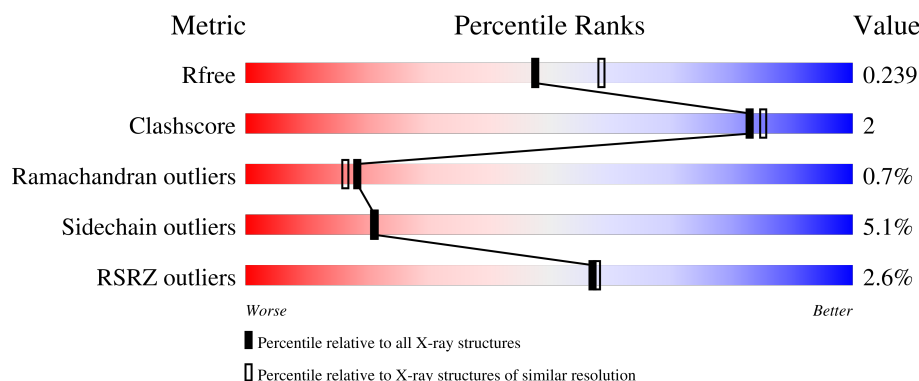
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.24 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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	162	6% 85% 12% ...
2	H	230	2% 80% 11% 9%
3	L	220	89% 7% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4583 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 Induced myeloid leukemia cell differentiation protein Mcl-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	159	Total	C	N	O	S	0	0	0
			1265	793	231	238	3			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	166	GLY	-	expression tag	UNP Q07820
A	167	PRO	-	expression tag	UNP Q07820
A	168	LEU	-	expression tag	UNP Q07820
A	169	GLY	-	expression tag	UNP Q07820
A	170	SER	-	expression tag	UNP Q07820
A	171	GLU	-	expression tag	UNP Q07820
A	172	ASP	-	expression tag	UNP Q07820
A	173	ASP	-	expression tag	UNP Q07820
A	193	SER	ALA	conflict	UNP Q07820
A	196	SER	THR	conflict	UNP Q07820
A	199	LEU	MET	conflict	UNP Q07820
A	201	GLU	ARG	conflict	UNP Q07820
A	202	ALA	SER	conflict	UNP Q07820
A	205	ALA	THR	conflict	UNP Q07820
A	206	GLY	SER	conflict	UNP Q07820
A	208	ARG	LYS	conflict	UNP Q07820

- Molecule 2 is a protein called Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	210	Total	C	N	O	S	0	0	0
			1565	992	260	306	7			

- Molecule 3 is a protein called Fab Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	213	Total 1599	C 990	N 271	O 332	S 6	0	0	0

- Molecule 4 is water.

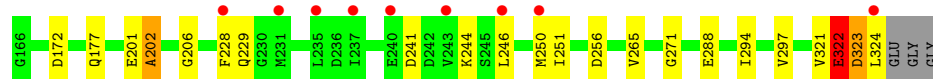
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	29	Total 29	O 29	0	0
4	H	52	Total 52	O 52	0	0
4	L	73	Total 73	O 73	0	0

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.

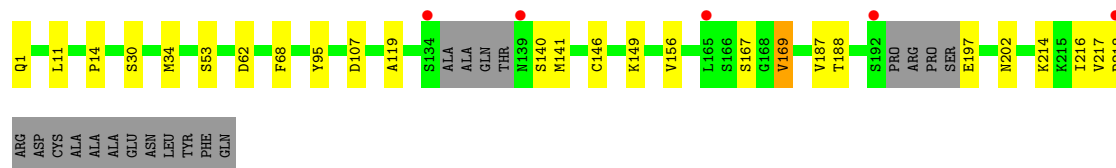
- Molecule 1: Induced myeloid leukemia cell differentiation protein Mcl-1

Chain A: 



- Molecule 2: Fab Heavy Chain

Chain H: 



- Molecule 3: Fab Light Chain

Chain L: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	148.05Å 42.46Å 106.23Å 90.00° 113.19° 90.00°	Depositor
Resolution (Å)	47.71 – 2.24 47.71 – 2.24	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.71-2.24) 99.1 (47.71-2.24)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.24Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.187 , 0.232 0.191 , 0.239	Depositor DCC
R_{free} test set	1497 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.232	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 34.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4583	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.21% 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	0.89	0/1286	1.48	2/1731 (0.1%)
2	H	0.87	1/1602 (0.1%)	1.14	5/2184 (0.2%)
3	L	0.86	0/1636	1.15	1/2233 (0.0%)
All	All	0.87	1/4524 (0.0%)	1.25	8/6148 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	34	MET	SD-CE	-5.10	1.66	1.79

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	172	ASP	CA-CB-CG	8.93	121.53	112.60
3	L	216	ASP	CA-CB-CG	6.13	118.73	112.60
2	H	68	PHE	CA-CB-CG	5.70	119.50	113.80
2	H	107	ASP	CA-CB-CG	5.65	118.25	112.60
2	H	95	TYR	CA-C-N	5.61	130.09	122.19
2	H	95	TYR	C-N-CA	5.61	130.09	122.19
1	A	228	PHE	CA-CB-CG	5.16	118.96	113.80
2	H	149	LYS	N-CA-C	5.00	117.64	109.59

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	1265	0	1263	10	0
2	H	1565	0	1539	6	0
3	L	1599	0	1526	5	0
4	A	29	0	0	0	0
4	H	52	0	0	0	0
4	L	73	0	0	1	0
All	All	4583	0	4328	21	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 2.

All (21) 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)
3:L:35:ASN:HD22	3:L:50:TYR:HA	1.59	0.68
1:A:177:GLN:HG3	1:A:202:ALA:HB2	1.78	0.65
1:A:241:ASP:HA	1:A:244:LYS:HD3	1.80	0.61
1:A:251:ILE:HG13	1:A:297:VAL:HG22	1.85	0.57
3:L:19:THR:HG22	4:L:371:HOH:O	2.06	0.54
2:H:169:VAL:HG12	2:H:187:VAL:HG23	1.89	0.54
2:H:141:MET:HE3	2:H:188:THR:HG22	1.89	0.53
1:A:323:ASP:HB3	1:A:324:LEU:HG	1.94	0.50
2:H:30:SER:O	2:H:53:SER:HB2	2.12	0.49
1:A:271:GLY:HA3	1:A:294:ILE:CD1	2.43	0.49
2:H:14:PRO:HD2	2:H:119:ALA:HB2	1.94	0.48
1:A:202:ALA:HB1	1:A:206:GLY:N	2.29	0.47
1:A:321:VAL:O	1:A:322:GLU:HB2	2.15	0.47
3:L:56:PRO:HD2	3:L:59:VAL:HG21	1.96	0.47
3:L:166:THR:HG23	3:L:167:GLN:O	2.16	0.45
1:A:271:GLY:HA3	1:A:294:ILE:HD11	1.98	0.45
3:L:155:ASP:HA	3:L:194:SER:HB3	2.00	0.44
1:A:202:ALA:HB1	1:A:206:GLY:H	1.83	0.44
1:A:250:MET:HE2	1:A:297:VAL:HG21	2.01	0.42
2:H:146:CYS:SG	2:H:216:ILE:HD11	2.60	0.42
2:H:217:VAL:HA	2:H:218:PRO:HD3	2.00	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	157/162 (97%)	149 (95%)	5 (3%)	3 (2%)	6	2
2	H	204/230 (89%)	196 (96%)	7 (3%)	1 (0%)	24	23
3	L	209/220 (95%)	205 (98%)	4 (2%)	0	100	100
All	All	570/612 (93%)	550 (96%)	16 (3%)	4 (1%)	18	16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	202	ALA
1	A	201	GLU
1	A	322	GLU
2	H	140	SER

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	136/137 (99%)	129 (95%)	7 (5%)	21	21
2	H	176/192 (92%)	167 (95%)	9 (5%)	21	21
3	L	183/189 (97%)	174 (95%)	9 (5%)	22	23
All	All	495/518 (96%)	470 (95%)	25 (5%)	21	21

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	229	GLN
1	A	246	LEU
1	A	256	ASP
1	A	265	VAL
1	A	288	GLU
1	A	322	GLU
1	A	323	ASP
2	H	1	GLN
2	H	11	LEU
2	H	62	ASP
2	H	156	VAL
2	H	167	SER
2	H	169	VAL
2	H	197	GLU
2	H	202	ASN
2	H	214	LYS
3	L	19	THR
3	L	130	GLU
3	L	150	VAL
3	L	163	MET
3	L	166	THR
3	L	169	SER
3	L	184	LEU
3	L	193	SER
3	L	196	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	189	GLN
1	A	221	GLN
2	H	82	GLN
2	H	177	GLN
3	L	35	ASN
3	L	53	ASN
3	L	54	GLN
3	L	80	GLN
3	L	201	HIS

5.3.3 RNA ⓘ

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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	L	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	L	202:GLU	C	205:GLY	N	2.93

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	159/162 (98%)	0.31	9 (5%)	29 27	28, 54, 133, 140	0
2	H	210/230 (91%)	-0.08	5 (2%)	59 61	24, 39, 73, 100	0
3	L	213/220 (96%)	-0.37	1 (0%)	87 88	24, 36, 54, 98	0
All	All	582/612 (95%)	-0.08	15 (2%)	57 58	24, 41, 94, 140	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	324	LEU	3.9
1	A	243	VAL	3.8
1	A	246	LEU	3.6
2	H	192	SER	3.2
3	L	215	ALA	2.9
2	H	165	LEU	2.8
1	A	237	ILE	2.8
2	H	218	PRO	2.7
2	H	139	ASN	2.4
1	A	235	LEU	2.3
1	A	240	GLU	2.3
1	A	228	PHE	2.2
1	A	250	MET	2.2
1	A	231	MET	2.1
2	H	134	SER	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.