



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:39 AM UTC

PDB ID : 5MK5 / pdb_00005mk5
Title : Structures of DHBN domain of human BLM helicase
Authors : Shi, J.; Chen, W.-F.; Zhang, B.; Fan, S.-H.; Ai, X.; Liu, N.-N.; Rety, S.; Xi, X.-G.
Deposited on : 2016-12-02
Resolution : 2.16 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:

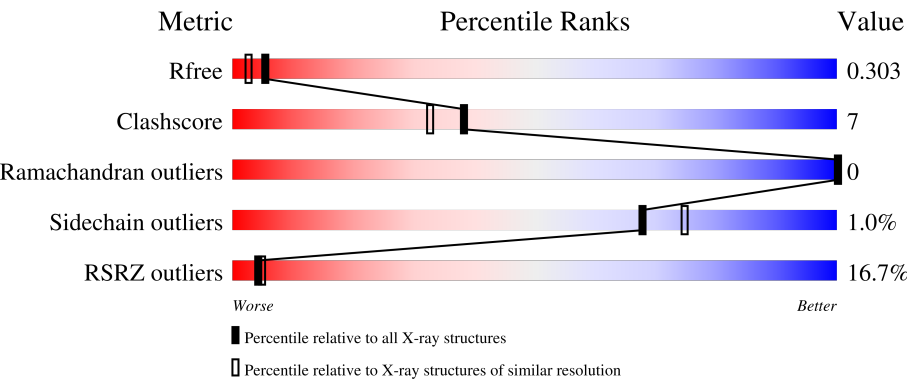
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	54	17% 83%11%6%
1	B	54	11% 89%7%. .
1	C	54	15% 80%15%. . .
1	D	54	20% 81%7%11%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	IOD	D	503	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3615 atoms, of which 1785 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 Bloom syndrome protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	51	Total	C	H	N	O	S	0	0	0
			875	264	453	79	76	3			
1	B	52	Total	C	H	N	O	S	0	0	0
			881	266	453	80	79	3			
1	C	52	Total	C	H	N	O	S	0	0	0
			881	266	453	80	79	3			
1	D	48	Total	C	H	N	O	S	0	0	0
			821	248	426	73	71	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	361	MET	-	initiating methionine	UNP P54132
B	361	MET	-	initiating methionine	UNP P54132
C	361	MET	-	initiating methionine	UNP P54132
D	361	MET	-	initiating methionine	UNP P54132

- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	I	0	0
			5	5		
2	B	4	Total	I	0	0
			4	4		
2	C	7	Total	I	0	0
			7	7		
2	D	4	Total	I	0	0
			4	4		

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

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

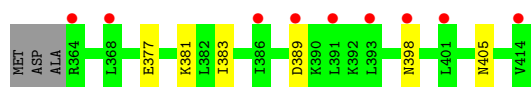
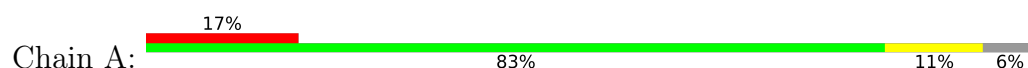
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	39	Total	O	0	0
			39	39		
4	B	35	Total	O	0	0
			35	35		
4	C	31	Total	O	0	0
			31	31		
4	D	29	Total	O	0	0
			29	29		

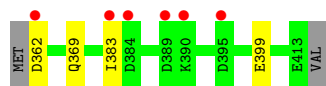
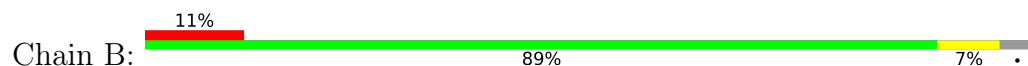
3 Residue-property plots [i](#)

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.

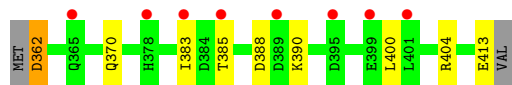
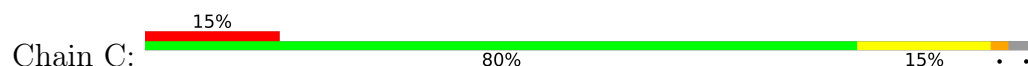
- Molecule 1: Bloom syndrome protein



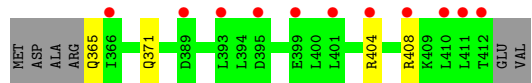
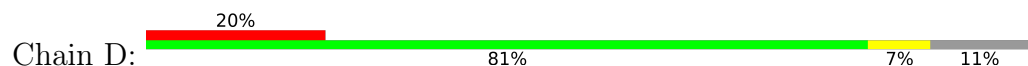
- Molecule 1: Bloom syndrome protein



- Molecule 1: Bloom syndrome protein



- Molecule 1: Bloom syndrome protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	34.08Å 144.73Å 96.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.84 – 2.16 27.84 – 2.16	Depositor EDS
% Data completeness (in resolution range)	91.2 (27.84-2.16) 91.0 (27.84-2.16)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.75 (at 2.16Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.219 , 0.297 0.224 , 0.303	Depositor DCC
R_{free} test set	580 reflections (4.35%)	wwPDB-VP
Wilson B-factor (Å ²)	22.3	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3615	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.29% 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](#)

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

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	0.10	0/424	0.26	0/568
1	B	0.11	0/430	0.22	0/576
1	C	0.11	0/430	0.24	0/576
1	D	0.10	0/397	0.21	0/532
All	All	0.11	0/1681	0.23	0/2252

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 [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	422	453	453	6	0
1	B	428	453	453	5	0
1	C	428	453	453	7	0
1	D	395	426	425	6	0
2	A	5	0	0	0	0
2	B	4	0	0	0	0
2	C	7	0	0	0	0
2	D	4	0	0	1	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	2	0	0	0	0
4	A	39	0	0	5	0
4	B	35	0	0	4	1
4	C	31	0	0	4	0
4	D	29	0	0	7	0
All	All	1830	1785	1784	24	1

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 7.

All (24) 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:C:370:GLN:OE1	4:C:601:HOH:O	1.93	0.86
1:A:398:ASN:ND2	4:A:602:HOH:O	2.09	0.85
1:A:381:LYS:NZ	4:A:603:HOH:O	2.09	0.78
1:B:399:GLU:OE2	4:B:601:HOH:O	2.01	0.78
1:D:408:ARG:NE	4:D:602:HOH:O	2.17	0.78
1:C:413:GLU:OE2	4:C:602:HOH:O	2.05	0.75
1:A:377:GLU:OE2	4:A:601:HOH:O	2.05	0.74
1:B:369:GLN:OE1	4:B:602:HOH:O	2.06	0.73
1:C:413:GLU:OE2	4:C:603:HOH:O	2.07	0.71
2:D:501:IOD:I	4:D:618:HOH:O	2.80	0.69
1:D:371:GLN:OE1	4:D:601:HOH:O	2.12	0.66
1:C:362:ASP:OD1	1:C:362:ASP:N	2.33	0.62
1:B:362:ASP:N	4:B:603:HOH:O	2.36	0.59
1:C:390:LYS:N	4:C:605:HOH:O	2.34	0.59
1:D:404:ARG:NE	4:D:603:HOH:O	2.40	0.54
1:D:365:GLN:N	4:D:604:HOH:O	2.40	0.53
1:B:369:GLN:NE2	4:B:607:HOH:O	2.42	0.51
1:C:383:ILE:HG21	1:C:400:LEU:HD21	1.98	0.46
1:A:383:ILE:HG13	1:B:383:ILE:HD11	1.98	0.45
1:C:388:ASP:OD1	1:C:404:ARG:NH2	2.52	0.43
1:D:408:ARG:CZ	4:D:602:HOH:O	2.64	0.42
1:A:405:ASN:HB2	4:A:610:HOH:O	2.19	0.42
1:A:389:ASP:HB3	4:A:614:HOH:O	2.21	0.40
1:D:404:ARG:CZ	4:D:603:HOH:O	2.68	0.40

All (1) 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	Interatomic distance (Å)	Clash overlap (Å)
4:B:634:HOH:O	4:B:634:HOH:O[3_454]	1.81	0.39

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	49/54 (91%)	47 (96%)	2 (4%)	0	100	100
1	B	50/54 (93%)	50 (100%)	0	0	100	100
1	C	50/54 (93%)	50 (100%)	0	0	100	100
1	D	46/54 (85%)	46 (100%)	0	0	100	100
All	All	195/216 (90%)	193 (99%)	2 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	50/52 (96%)	50 (100%)	0	100	100
1	B	50/52 (96%)	50 (100%)	0	100	100
1	C	50/52 (96%)	48 (96%)	2 (4%)	28	27
1	D	47/52 (90%)	47 (100%)	0	100	100
All	All	197/208 (95%)	195 (99%)	2 (1%)	68	75

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	362	ASP
1	C	385	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	371	GLN
1	A	374	HIS
1	B	365	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](#)

Of 23 ligands modelled in this entry, 23 are monoatomic - leaving 0 for Mogul analysis.

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

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	51/54 (94%)	1.11	9 (17%) 4 4	15, 39, 66, 90	0
1	B	52/54 (96%)	0.95	6 (11%) 9 10	17, 38, 63, 70	0
1	C	52/54 (96%)	0.99	8 (15%) 5 5	21, 39, 61, 74	0
1	D	48/54 (88%)	1.44	11 (22%) 2 2	29, 49, 76, 82	0
All	All	203/216 (93%)	1.12	34 (16%) 4 5	15, 42, 67, 90	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	414	VAL	9.3
1	D	412	THR	5.5
1	C	389	ASP	5.2
1	D	389	ASP	4.5
1	B	384	ASP	3.9
1	D	411	LEU	3.7
1	B	383	ILE	3.3
1	D	408	ARG	3.3
1	C	385	THR	3.1
1	C	401	LEU	3.0
1	A	398	ASN	2.9
1	D	393	LEU	2.9
1	A	368	LEU	2.9
1	B	362	ASP	2.9
1	D	410	LEU	2.8
1	A	389	ASP	2.8
1	B	389	ASP	2.8
1	C	395	ASP	2.8
1	A	364	ARG	2.7
1	D	399	GLU	2.7
1	A	386	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	383	ILE	2.5
1	A	401	LEU	2.5
1	B	390	LYS	2.5
1	D	401	LEU	2.4
1	C	365	GLN	2.3
1	B	395	ASP	2.3
1	D	366	ILE	2.3
1	A	391	LEU	2.3
1	C	378	HIS	2.2
1	C	399	GLU	2.2
1	A	393	LEU	2.2
1	D	404	ARG	2.2
1	D	395	ASP	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 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(Å ²)	Q<0.9
2	IOD	D	503	1/1	0.74	0.41	192,192,192,192	1
3	K	A	506	1/1	0.83	0.14	65,65,65,65	0
2	IOD	B	502	1/1	0.84	0.19	140,140,140,140	0
2	IOD	D	504	1/1	0.86	0.17	137,137,137,137	1
2	IOD	C	508	1/1	0.87	0.18	155,155,155,155	1
2	IOD	A	505	1/1	0.91	0.12	87,87,87,87	1
2	IOD	D	502	1/1	0.92	0.15	143,143,143,143	1
2	IOD	B	503	1/1	0.93	0.15	97,97,97,97	1
3	K	C	502	1/1	0.94	0.14	63,63,63,63	0
3	K	C	501	1/1	0.95	0.06	47,47,47,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	IOD	C	509	1/1	0.95	0.10	149,149,149,149	1
2	IOD	A	504	1/1	0.96	0.07	86,86,86,86	1
2	IOD	B	504	1/1	0.96	0.08	69,69,69,69	1
2	IOD	C	507	1/1	0.97	0.26	74,74,74,74	1
2	IOD	C	506	1/1	0.98	0.07	61,61,61,61	1
2	IOD	D	501	1/1	0.98	0.07	81,81,81,81	0
2	IOD	C	505	1/1	0.99	0.10	51,51,51,51	1
2	IOD	B	501	1/1	0.99	0.02	58,58,58,58	0
2	IOD	A	502	1/1	0.99	0.06	57,57,57,57	1
2	IOD	C	504	1/1	0.99	0.08	48,48,48,48	0
2	IOD	A	503	1/1	1.00	0.03	43,43,43,43	0
2	IOD	C	503	1/1	1.00	0.05	37,37,37,37	0
2	IOD	A	501	1/1	1.00	0.07	39,39,39,39	1

6.5 Other polymers [i](#)

There are no such residues in this entry.