



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 25, 2026 – 04:18 PM EDT

PDB ID : 3F8I / pdb_00003f8i
Title : Mouse UHRF1 SRA domain bound with hemi-methylated CpG, crystal structure in space group P21
Authors : Hashimoto, H.; Horton, J.R.; Zhang, X.; Cheng, X.
Deposited on : 2008-11-12
Resolution : 2.29 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

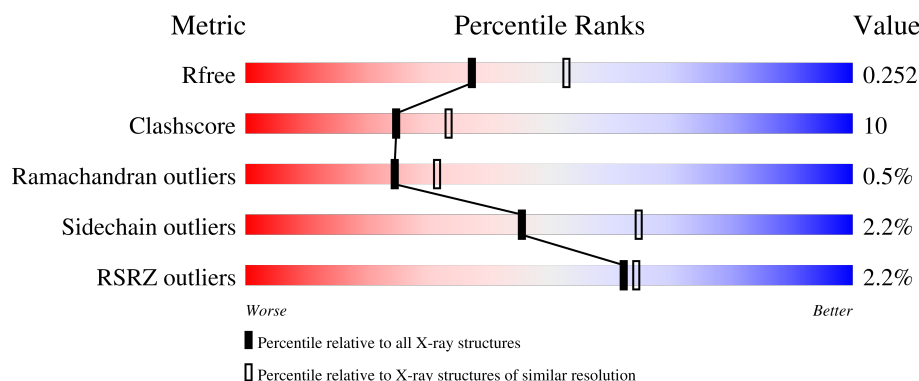
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

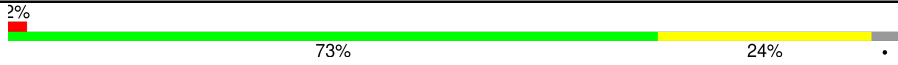
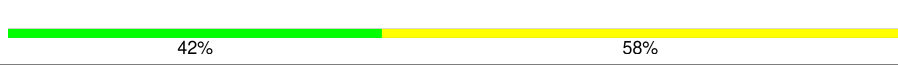
The reported resolution of this entry is 2.29 Å.

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.



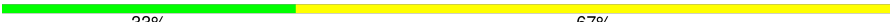
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	212	
1	B	212	
2	D	12	
2	F	12	
3	E	12	

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Mol	Chain	Length	Quality of chain
3	G	12	 A horizontal bar chart showing the quality of chain G. The bar is divided into two segments: a green segment on the left representing 33% and a yellow segment on the right representing 67%. The percentages are labeled below the bar.

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4313 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 E3 ubiquitin-protein ligase UHRF1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	S	0	0	0
			1606	999	308	294	5			
1	B	206	Total	C	N	O	S	0	0	0
			1598	996	304	293	5			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	417	HIS	-	expression tag	UNP Q8VDF2
A	418	MET	VAL	conflict	UNP Q8VDF2
B	417	HIS	-	expression tag	UNP Q8VDF2
B	418	MET	VAL	conflict	UNP Q8VDF2

- Molecule 2 is a DNA chain called 5'-D(*DCP*DCP*DAP*DTP*DGP*(5CM)P*DGP*DCP*DTP*DGP*DAP*DC)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	12	Total	C	N	O	P	0	0	0
			241	116	44	70	11			
2	F	12	Total	C	N	O	P	0	0	0
			241	116	44	70	11			

- Molecule 3 is a DNA chain called 5'-D(*DGP*DTP*DCP*DAP*DGP*DCP*DGP*DCP*DAP*DTP*DGP*DG)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	12	Total	C	N	O	P	0	0	0
			246	117	48	70	11			
3	G	12	Total	C	N	O	P	0	0	0
			246	117	48	70	11			

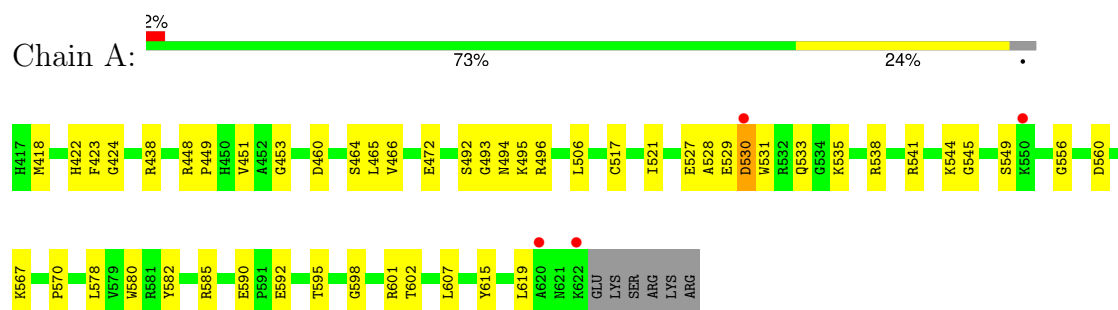
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	44	Total 44	O 44	0	0
4	B	64	Total 64	O 64	0	0
4	D	7	Total 7	O 7	0	0
4	E	4	Total 4	O 4	0	0
4	F	8	Total 8	O 8	0	0
4	G	8	Total 8	O 8	0	0

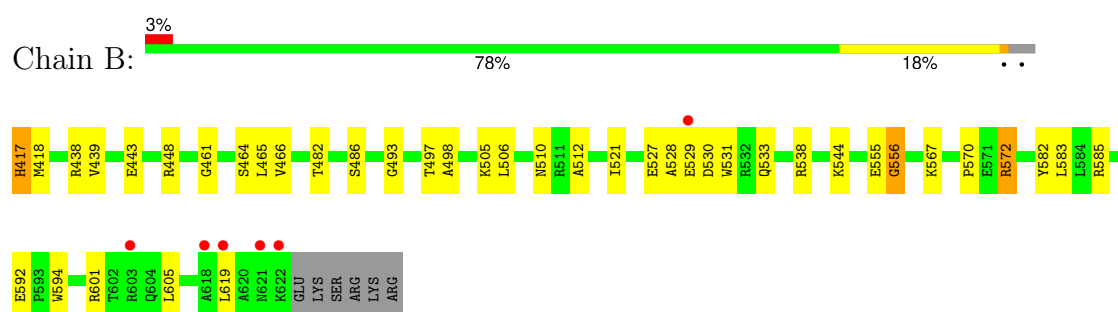
3 Residue-property plots [i](#)

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: E3 ubiquitin-protein ligase UHRF1



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- Molecule 2: 5'-D(*DCP*DCP*DAP*DTP*DGP*(5CM)P*DGP*DCP*DTP*DGP*DAP*DC)-3'

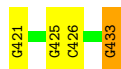


- Molecule 2: 5'-D(*DCP*DCP*DAP*DTP*DGP*(5CM)P*DGP*DCP*DTP*DGP*DAP*DC)-3'

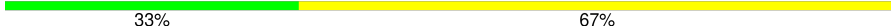


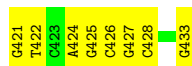
- Molecule 3: 5'-D(*DGP*DTP*DCP*DAP*DGP*DCP*DGP*DCP*DAP*DTP*DGP*DG)-3',

Chain E:  67% 25% 8%



- Molecule 3: 5'-D(*DGP*DTP*DCP*DAP*DGP*DCP*DGP*DCP*DAP*DTP*DGP*DG)-3',

Chain G:  33% 67%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.08Å 62.42Å 113.45Å 90.00° 98.92° 90.00°	Depositor
Resolution (Å)	34.53 – 2.29 34.53 – 2.29	Depositor EDS
% Data completeness (in resolution range)	88.5 (34.53-2.29) 88.5 (34.53-2.29)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.29Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.223 , 0.268 0.205 , 0.252	Depositor DCC
R_{free} test set	1487 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.925	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4313	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.64% 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: 5CM

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	0.38	0/1649	0.88	6/2233 (0.3%)
1	B	0.43	0/1641	0.89	3/2223 (0.1%)
2	D	0.29	0/246	0.68	0/375
2	F	0.28	0/246	0.68	0/375
3	E	0.28	0/276	0.75	1/425 (0.2%)
3	G	0.31	0/276	0.71	0/425
All	All	0.38	0/4334	0.84	10/6056 (0.2%)

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
3	E	0	1
3	G	0	1
All	All	0	2

There are no bond length outliers.

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	433	DG	C2'-C3'-O3'	-6.73	101.41	111.50
1	B	556	GLY	N-CA-C	5.74	121.49	111.50
1	B	493	GLY	N-CA-C	-5.58	106.80	115.67
1	A	496	ARG	N-CA-C	-5.51	107.56	114.56
1	A	556	GLY	N-CA-C	5.47	121.02	111.50

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	E	433	DG	Sidechain
3	G	433	DG	Sidechain

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	1606	0	1526	31	0
1	B	1598	0	1513	29	0
2	D	241	0	138	6	0
2	F	241	0	138	5	0
3	E	246	0	136	4	0
3	G	246	0	136	8	0
4	A	44	0	0	0	0
4	B	64	0	0	2	0
4	D	7	0	0	0	0
4	E	4	0	0	0	0
4	F	8	0	0	0	0
4	G	8	0	0	0	0
All	All	4313	0	3587	76	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 10.

The worst 5 of 76 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:B:486:SER:H	1:B:510:ASN:HD21	1.10	1.00
1:B:572:ARG:HH11	1:B:572:ARG:HB3	1.40	0.86
1:A:590:GLU:CD	1:A:601:ARG:HH12	1.97	0.70
1:A:544:LYS:HE3	2:D:408:DG:OP2	1.92	0.69
1:B:486:SER:N	1:B:510:ASN:HD21	1.89	0.69

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	204/212 (96%)	194 (95%)	9 (4%)	1 (0%)	24	31
1	B	204/212 (96%)	193 (95%)	10 (5%)	1 (0%)	24	31
All	All	408/424 (96%)	387 (95%)	19 (5%)	2 (0%)	24	31

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	530	ASP
1	B	570	PRO

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	161/173 (93%)	158 (98%)	3 (2%)	50	69
1	B	159/173 (92%)	155 (98%)	4 (2%)	42	60
All	All	320/346 (92%)	313 (98%)	7 (2%)	45	65

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	417	HIS
1	B	466	VAL
1	B	572	ARG
1	B	505	LYS

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Mol	Chain	Res	Type
1	A	619	LEU

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

Mol	Chain	Res	Type
1	A	479	ASN
1	A	533	GLN
1	B	479	ASN
1	B	510	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are 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
2	5CM	D	407	2	18,21,22	0.67	1 (5%)	24,30,33	0.86	2 (8%)
2	5CM	F	407	2	18,21,22	0.81	1 (5%)	24,30,33	1.00	1 (4%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	5CM	D	407	2	-	0/7/21/22	0/2/2/2
2	5CM	F	407	2	-	0/7/21/22	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	407	5CM	C5A-C5	-2.49	1.44	1.50
2	D	407	5CM	C5A-C5	-2.13	1.45	1.50

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	407	5CM	C5A-C5-C6	-2.92	118.91	122.85
2	D	407	5CM	C5A-C5-C6	-2.03	120.11	122.85
2	D	407	5CM	C5-C6-N1	-2.02	121.12	123.31

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

There are no chain breaks in this entry.

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	206/212 (97%)	0.29	4 (1%) 66 68	23, 36, 51, 62	0
1	B	206/212 (97%)	0.05	6 (2%) 53 56	14, 27, 44, 73	0
2	D	11/12 (91%)	0.23	0 100 100	31, 43, 54, 58	0
2	F	11/12 (91%)	0.12	0 100 100	26, 39, 52, 56	0
3	E	12/12 (100%)	0.23	0 100 100	33, 43, 52, 52	0
3	G	12/12 (100%)	0.15	0 100 100	21, 40, 50, 51	0
All	All	458/472 (97%)	0.17	10 (2%) 62 64	14, 33, 52, 73	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	622	LYS	3.6
1	B	621	ASN	3.3
1	B	603	ARG	2.6
1	A	622	LYS	2.5
1	B	618	ALA	2.4

6.2 Non-standard residues in protein, DNA, RNA chains [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	5CM	D	407	20/21	0.95	0.08	27,29,33,34	0
2	5CM	F	407	20/21	0.96	0.07	14,21,25,31	0

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.