



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 22, 2016 – 01:05 AM EDT

PDB ID : 4Z2M  
Title : Crystal structure of human SPT16 Mid-AID/H3-H4 tetramer FACT Histone complex  
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Deposited on : 2015-03-30  
Resolution : 2.98 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027107  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0122  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027107

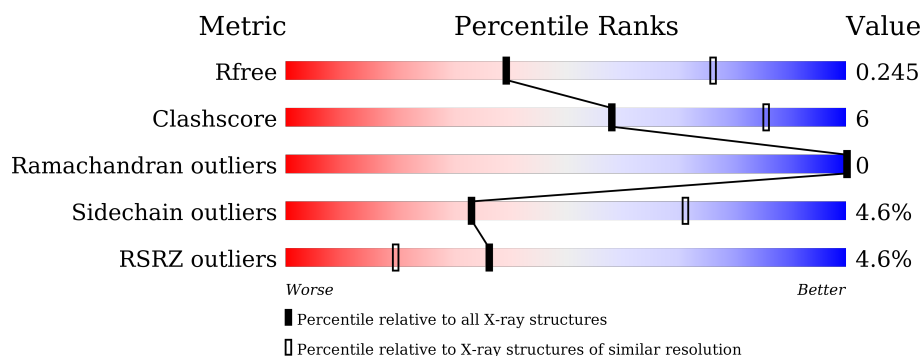
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.98 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))
$R_{free}$	91344	1992 (3.00-2.96)
Clashscore	102246	2349 (3.00-2.96)
Ramachandran outliers	100387	2274 (3.00-2.96)
Sidechain outliers	100360	2277 (3.00-2.96)
RSRZ outliers	91569	2007 (3.00-2.96)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	B	287	<div> <div>5%</div> <div> <div></div> <div>72%</div> <div>21%</div> <div>• 6%</div> </div> </div>
2	G	102	<div> <div>4%</div> <div> <div></div> <div>59%</div> <div>15%</div> <div>• 25%</div> </div> </div>
2	I	102	<div> <div>7%</div> <div> <div></div> <div>63%</div> <div>12%</div> <div>25%</div> </div> </div>
3	H	103	<div> <div>%</div> <div> <div></div> <div>54%</div> <div>14%</div> <div>• 31%</div> </div> </div>
3	J	103	<div> <div>%</div> <div> <div></div> <div>62%</div> <div>• •</div> <div>32%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4594 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 FACT complex subunit SPT16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	271	Total	C	N	O	S	0	0	0
			2220	1429	371	409	11			

- Molecule 2 is a protein called Histone H3.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	77	Total	C	N	O	S	0	0	0
			626	396	116	110	4			
2	I	76	Total	C	N	O	S	0	0	0
			615	390	115	106	4			

- Molecule 3 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	71	Total	C	N	O	S	0	0	0
			572	359	113	99	1			
3	J	70	Total	C	N	O	S	0	0	0
			561	353	109	98	1			





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.44Å 69.78Å 187.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.96 – 2.98 46.96 – 2.98	Depositor EDS
% Data completeness (in resolution range)	97.6 (46.96-2.98) 94.3 (46.96-2.98)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.75 (at 2.96Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.190 , 0.246 0.188 , 0.245	Depositor DCC
$R_{free}$ test set	1723 reflections (10.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	74.0	Xtriage
Anisotropy	0.482	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 42.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 17819 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4594	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	B	0.45	0/2271	0.62	0/3068
2	G	0.45	0/633	0.61	0/849
2	I	0.50	0/622	0.63	0/834
3	H	0.43	0/577	0.57	0/773
3	J	0.43	0/566	0.63	0/759
All	All	0.45	0/4669	0.62	0/6283

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	B	2220	0	2203	33	1
2	G	626	0	656	11	0
2	I	615	0	648	8	1
3	H	572	0	615	9	0
3	J	561	0	602	3	0
All	All	4594	0	4724	55	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 6.

All (55) 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:660:PRO:HG2	1:B:682:ALA:HB3	1.52	0.92
1:B:708:LEU:HB2	1:B:721:HIS:HB3	1.65	0.79
3:H:30:THR:HB	3:H:32:PRO:HD2	1.69	0.73
1:B:659:ASN:ND2	1:B:681:GLU:OE2	2.29	0.65
1:B:688:ARG:HH11	1:B:696:LYS:HD2	1.64	0.63
1:B:915:GLY:HA2	1:B:918:GLU:HG3	1.79	0.63
1:B:664:ASP:O	1:B:742:THR:OG1	2.18	0.62
1:B:691:SER:HB3	1:B:695:ASP:O	1.99	0.61
2:I:79:LYS:N	2:I:80:THR:HA	2.15	0.60
1:B:906:MET:HE1	1:B:909:ILE:HD12	1.84	0.59
1:B:812:ARG:HD2	2:I:116:ARG:O	2.04	0.57
1:B:662:LEU:HB3	1:B:665:LEU:HD11	1.87	0.57
2:G:76:GLN:OE1	3:H:23:ARG:NH1	2.39	0.55
1:B:710:GLN:HB3	1:B:719:VAL:HB	1.91	0.53
1:B:800:PRO:HA	1:B:820:THR:HG22	1.92	0.52
3:J:56:GLY:HA2	3:J:59:LYS:HE2	1.91	0.52
1:B:744:VAL:HG21	1:B:773:ARG:HG2	1.93	0.50
3:J:28:GLY:O	3:J:29:ILE:HD13	2.13	0.48
1:B:668:ARG:HD2	1:B:741:TYR:CE2	2.49	0.48
1:B:883:LYS:NZ	1:B:895:GLU:OE2	2.30	0.48
1:B:892:LYS:HD3	1:B:892:LYS:HA	1.73	0.47
2:G:97:GLU:O	2:G:101:VAL:HG23	2.13	0.47
2:G:134:ARG:NH2	3:H:60:VAL:HG13	2.29	0.47
2:G:128:ARG:NH1	2:G:133:GLU:OE1	2.48	0.47
2:G:70:LEU:O	2:G:74:ILE:HG13	2.15	0.47
1:B:716:MET:SD	1:B:745:GLY:HA3	2.55	0.46
1:B:748:THR:HB	3:H:46:ILE:CG2	2.46	0.46
2:G:73:GLU:HG3	3:H:23:ARG:HB2	1.96	0.45
1:B:664:ASP:HA	1:B:677:GLN:HE21	1.81	0.45
1:B:679:SER:O	1:B:690:THR:N	2.43	0.45
3:H:70:VAL:O	3:H:74:GLU:HG2	2.16	0.45
1:B:777:LYS:HG3	1:B:781:LYS:HE3	1.98	0.45
1:B:700:LEU:O	1:B:703:ASN:N	2.42	0.45
2:I:128:ARG:NH2	2:I:134:ARG:HB2	2.32	0.44
3:H:68:ASP:O	3:H:71:THR:HB	2.18	0.44
2:I:93:GLN:O	2:I:97:GLU:HG3	2.17	0.44
2:G:130:ILE:HG12	2:I:106:ASP:HB3	2.00	0.44
2:I:118:THR:HA	3:J:45:ARG:HB3	1.99	0.43
2:G:100:LEU:HD23	2:G:100:LEU:HA	1.84	0.43
1:B:713:ASP:OD1	1:B:777:LYS:NZ	2.41	0.43
2:I:128:ARG:HG2	2:I:133:GLU:OE1	2.20	0.42
1:B:819:PRO:HB3	1:B:885:TRP:CZ2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:769:GLU:O	1:B:773:ARG:HG3	2.20	0.42
1:B:806:PHE:CG	1:B:882:ILE:HD11	2.54	0.42
1:B:744:VAL:CG2	1:B:773:ARG:HG2	2.50	0.42
1:B:830:TRP:CD2	1:B:832:PRO:HD3	2.55	0.41
1:B:843:ILE:O	1:B:893:TYR:HA	2.20	0.41
1:B:830:TRP:CE3	1:B:832:PRO:HD3	2.55	0.41
3:H:59:LYS:HE2	3:H:63:GLU:OE2	2.19	0.41
1:B:853:LYS:HG3	2:I:105:GLU:OE1	2.21	0.41
2:G:59:GLU:HB2	3:H:40:ARG:HH12	1.87	0.40
1:B:706:HIS:HB2	1:B:723:HIS:HB3	2.03	0.40
1:B:727:ALA:HB2	1:B:736:THR:HG22	2.03	0.40
2:G:125:GLN:O	2:G:129:ARG:HB2	2.22	0.40
2:G:82:LEU:HD23	2:G:82:LEU:HA	1.92	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)
1:B:762:ASP:OD1	2:I:87:SER:OG[1_545]	2.17	0.03

## 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	B	267/287 (93%)	260 (97%)	7 (3%)	0	100	100
2	G	75/102 (74%)	72 (96%)	3 (4%)	0	100	100
2	I	74/102 (72%)	72 (97%)	2 (3%)	0	100	100
3	H	69/103 (67%)	69 (100%)	0	0	100	100
3	J	68/103 (66%)	67 (98%)	1 (2%)	0	100	100
All	All	553/697 (79%)	540 (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	B	247/260 (95%)	235 (95%)	12 (5%)	31	69
2	G	67/88 (76%)	63 (94%)	4 (6%)	24	61
2	I	65/88 (74%)	65 (100%)	0	100	100
3	H	60/79 (76%)	56 (93%)	4 (7%)	20	56
3	J	59/79 (75%)	56 (95%)	3 (5%)	29	67
All	All	498/594 (84%)	475 (95%)	23 (5%)	33	71

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	649	ASP
1	B	659	ASN
1	B	729	MET
1	B	744	VAL
1	B	749	THR
1	B	812	ARG
1	B	813	SER
1	B	825	VAL
1	B	853	LYS
1	B	876	VAL
1	B	878	SER
1	B	894	THR
2	G	76	GLN
2	G	77	ASP
2	G	118	THR
2	G	129	ARG
3	H	29	ILE
3	H	45	ARG
3	H	46	ILE
3	H	47	SER
3	J	45	ARG

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Mol	Chain	Res	Type
3	J	59	LYS
3	J	80	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	735	HIS

### 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 carbohydrates 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

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	B	271/287 (94%)	0.28	13 (4%) 34 19	41, 80, 128, 158	0
2	G	77/102 (75%)	0.27	4 (5%) 31 16	45, 61, 106, 130	0
2	I	76/102 (74%)	0.47	7 (9%) 11 5	45, 71, 126, 144	0
3	H	71/103 (68%)	0.18	1 (1%) 78 57	52, 76, 107, 114	0
3	J	70/103 (67%)	0.36	1 (1%) 78 57	52, 78, 117, 127	0
All	All	565/697 (81%)	0.30	26 (4%) 36 19	41, 76, 124, 158	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	658	SER	6.0
2	I	78	PHE	4.1
3	H	79	LYS	3.8
3	J	77	LYS	3.8
1	B	656	ASN	3.6
1	B	657	ARG	3.5
1	B	732	LYS	3.4
2	I	80	THR	3.3
1	B	729	MET	3.0
2	G	80	THR	3.0
2	G	81	ASP	3.0
2	I	84	PHE	2.8
2	G	78	PHE	2.8
1	B	767	GLN	2.8
1	B	697	VAL	2.4
2	I	81	ASP	2.4
1	B	654	ASN	2.4
1	B	671	ILE	2.4
1	B	655	LEU	2.3
1	B	730	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
2	I	82	LEU	2.2
1	B	647	LYS	2.1
1	B	676	MET	2.1
2	G	58	THR	2.1
2	I	75	ALA	2.0
2	I	72	ARG	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 carbohydrates 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.