



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:49 AM GMT

PDB ID : 2IW5  
Title : STRUCTURAL BASIS FOR COREST-DEPENDENT DEMETHYLATION  
OF NUCLEOSOMES BY THE HUMAN LSD1 HISTONE DEMETHYLASE  
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Otwinowski, Z.; Yu, H.  
Deposited on : 2006-06-26  
Resolution : 2.57 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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

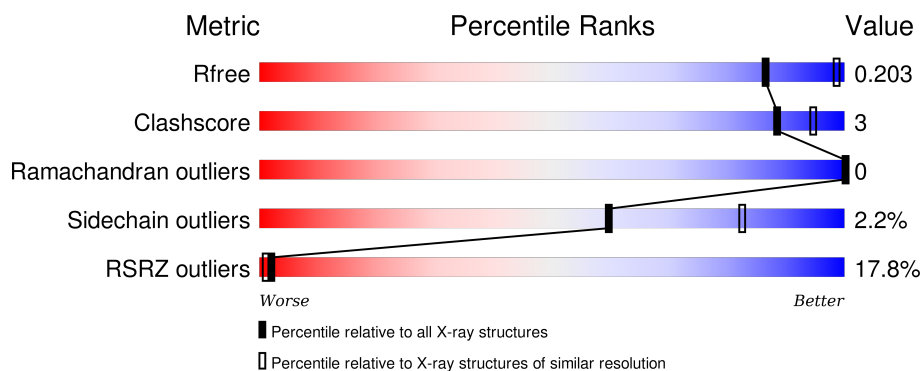
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.57 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	2636 (2.60-2.56)
Clashscore	102246	3003 (2.60-2.56)
Ramachandran outliers	100387	2956 (2.60-2.56)
Sidechain outliers	100360	2956 (2.60-2.56)
RSRZ outliers	91569	2642 (2.60-2.56)

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	A	666	<div> <div>16%</div> <div>93%</div> <div>7%</div> </div>
2	B	235	<div> <div>15%</div> <div>54%</div> <div>43%</div> </div>

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
6	NH4	A	1842	-	-	-	X

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6440 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.

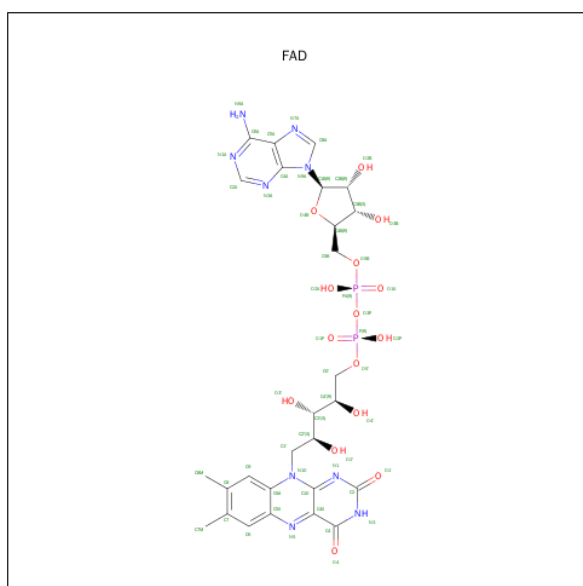
- Molecule 1 is a protein called LYSINE-SPECIFIC HISTONE DEMETHYLASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	666	Total	C	N	O	S	0	0	0
			5217	3324	906	967	20			

- Molecule 2 is a protein called REST COREPRESSOR 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	133	Total	C	N	O	S	0	0	0
			1076	676	194	203	3			

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

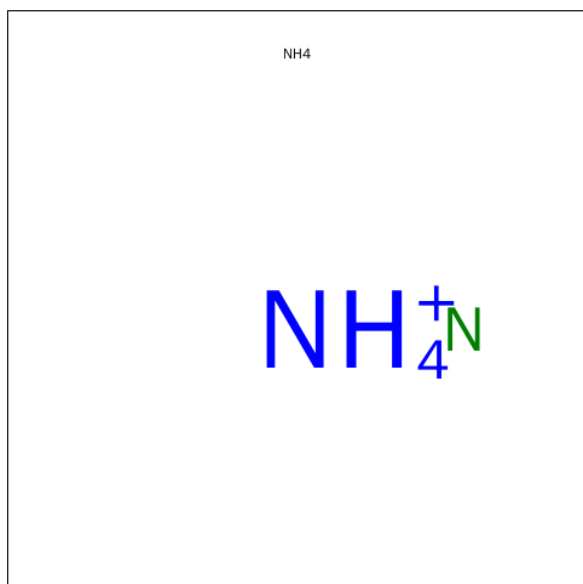
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Cl	0	0
			3	3		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is AMMONIUM ION (three-letter code: NH4) (formula:  $H_4N$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total N 1 1	0	0

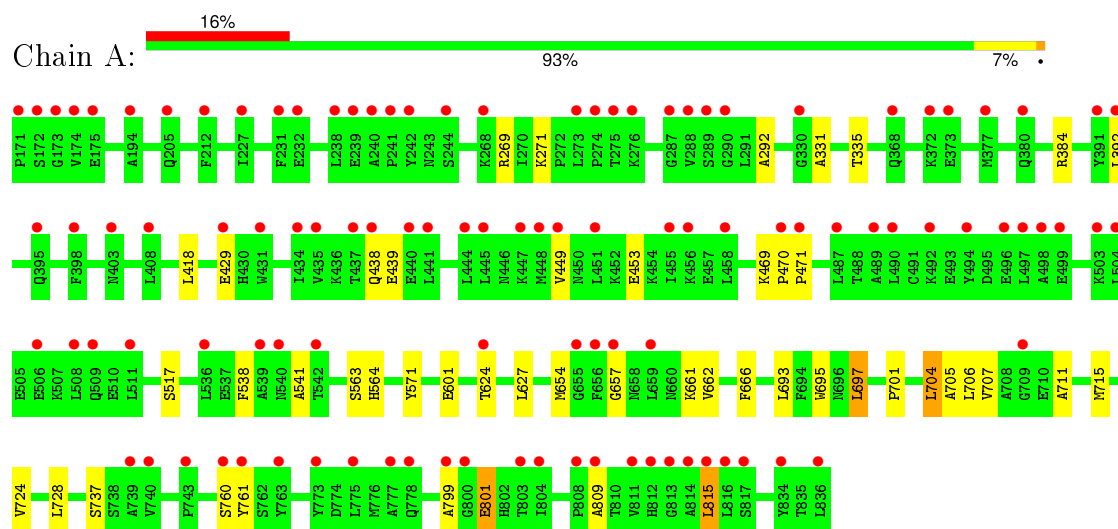
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	77	Total O 77 77	0	0
7	B	1	Total O 1 1	0	0

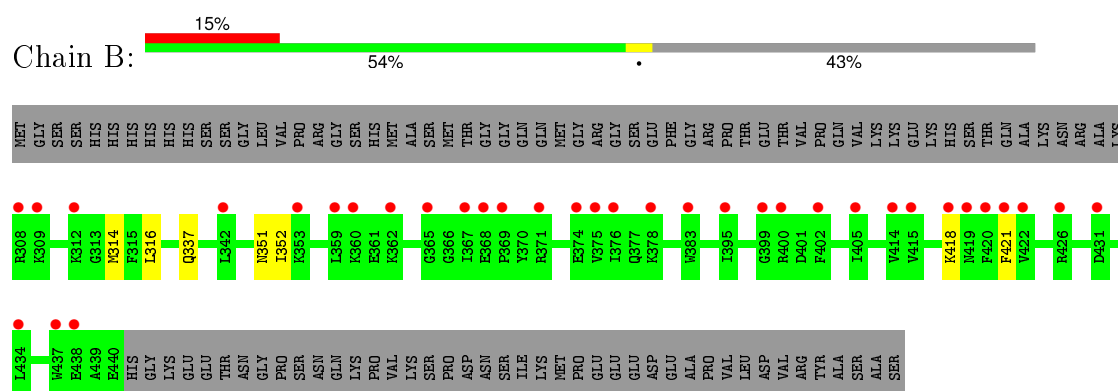
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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: LYSINE-SPECIFIC HISTONE DEMETHYLASE 1



#### • Molecule 2: REST COREPRESSOR 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	120.35Å 178.22Å 234.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.57 49.87 – 2.57	Depositor EDS
% Data completeness (in resolution range)	93.2 (20.00-2.57) 93.2 (49.87-2.57)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.73 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.203 , 0.212 0.193 , 0.203	Depositor DCC
$R_{free}$ test set	1498 reflections (2.04%)	DCC
Wilson B-factor (Å <sup>2</sup> )	62.7	Xtriage
Anisotropy	0.475	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 68.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 75084 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6440	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CL, FAD, NH4

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	0.50	0/5331	0.66	2/7232 (0.0%)
2	B	0.42	0/1091	0.55	0/1471
All	All	0.48	0/6422	0.65	2/8703 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	815	LEU	CA-CB-CG	6.60	130.49	115.30
1	A	697	LEU	N-CA-C	5.30	125.32	111.00

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	5217	0	5252	30	0
2	B	1076	0	1091	5	0
3	A	53	0	31	2	0
4	A	3	0	0	1	0
5	A	12	0	16	0	0
6	A	1	0	0	0	0
7	A	77	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	1	0	0	0	0
All	All	6440	0	6390	32	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 3.

All (32) 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)
4:A:1838:CL:CL	4:A:1843:CL:CL	2.19	1.34
1:A:384:ARG:HB3	2:B:314:MET:HE1	1.59	0.85
1:A:801:GLU:HG2	1:A:809:ALA:H	1.48	0.76
1:A:707:VAL:HG11	1:A:715:MET:HG3	1.75	0.67
1:A:801:GLU:HG3	1:A:809:ALA:HA	1.76	0.66
1:A:384:ARG:HB3	2:B:314:MET:CE	2.27	0.65
1:A:335:THR:HG22	1:A:564:HIS:ND1	2.20	0.57
1:A:761:TYR:CD1	1:A:809:ALA:HB1	2.41	0.55
1:A:695:TRP:CE3	1:A:697:LEU:HD11	2.43	0.54
1:A:470:PRO:HA	1:A:471:PRO:C	2.28	0.53
1:A:711:ALA:O	1:A:715:MET:HG2	2.10	0.52
1:A:624:THR:HG22	1:A:799:ALA:HB3	1.92	0.51
1:A:335:THR:HG22	1:A:564:HIS:CE1	2.47	0.50
1:A:627:LEU:HD22	1:A:654:MET:HE2	1.93	0.50
1:A:661:LYS:HB3	1:A:704:LEU:HD21	1.96	0.47
1:A:661:LYS:HD3	1:A:704:LEU:HD21	1.96	0.47
1:A:707:VAL:CG1	1:A:715:MET:HG3	2.45	0.47
1:A:449:VAL:O	1:A:453:GLU:HG2	2.16	0.46
1:A:801:GLU:HG2	1:A:809:ALA:N	2.26	0.45
1:A:269:ARG:HG2	1:A:271:LYS:O	2.17	0.45
1:A:724:VAL:O	1:A:728:LEU:HG	2.17	0.45
1:A:760:SER:HB2	3:A:900:FAD:HM83	2.00	0.44
1:A:292:ALA:HB2	1:A:815:LEU:HD22	2.01	0.43
1:A:693:LEU:HD23	1:A:706:LEU:HD22	2.00	0.43
1:A:662:VAL:HB	1:A:705:ALA:HB3	2.00	0.43
1:A:438:GLN:HG3	2:B:352:ILE:HG22	2.01	0.43
2:B:418:LYS:O	2:B:421:PHE:HB2	2.18	0.43
1:A:801:GLU:CG	1:A:809:ALA:HA	2.45	0.42
1:A:541:ALA:O	1:A:657:GLY:HA3	2.19	0.42
1:A:666:PHE:O	1:A:701:PRO:HG2	2.20	0.42
1:A:439:GLU:HG3	2:B:352:ILE:HD13	2.02	0.41
1:A:331:ALA:HA	3:A:900:FAD:C4X	2.51	0.40

There are no symmetry-related clashes.

## 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	664/666 (100%)	647 (97%)	17 (3%)	0	100	100
2	B	131/235 (56%)	127 (97%)	4 (3%)	0	100	100
All	All	795/901 (88%)	774 (97%)	21 (3%)	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	566/566 (100%)	554 (98%)	12 (2%)	61	84
2	B	117/203 (58%)	114 (97%)	3 (3%)	54	78
All	All	683/769 (89%)	668 (98%)	15 (2%)	60	82

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	392	LEU
1	A	418	LEU
1	A	429	GLU
1	A	469	LYS
1	A	517	SER

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Mol	Chain	Res	Type
1	A	538	PHE
1	A	563	SER
1	A	571	TYR
1	A	601	GLU
1	A	704	LEU
1	A	737	SER
1	A	801	GLU
2	B	316	LEU
2	B	337	GLN
2	B	351	ASN

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

Mol	Chain	Res	Type
1	A	532	HIS
1	A	633	GLN
2	B	348	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 1 is modelled with single atom and 3 are monoatomic - leaving 3 for Mogul analysis.

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	$\# Z  > 2$
5	GOL	A	1839	-	5,5,5	0.38	0	5,5,5	0.32	0
5	GOL	A	1840	-	5,5,5	0.33	0	5,5,5	0.27	0
3	FAD	A	900	-	48,58,58	1.23	5 (10%)	54,89,89	2.43	8 (14%)

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
5	GOL	A	1839	-	-	0/4/4/4	0/0/0/0
5	GOL	A	1840	-	-	0/4/4/4	0/0/0/0
3	FAD	A	900	-	-	0/30/50/50	0/6/6/6

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	900	FAD	C10-N1	2.06	1.39	1.35
3	A	900	FAD	C2A-N1A	2.67	1.39	1.33
3	A	900	FAD	C4-N3	2.71	1.38	1.33
3	A	900	FAD	C2A-N3A	3.35	1.38	1.32
3	A	900	FAD	C4X-N5	3.58	1.38	1.33

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	900	FAD	N3A-C2A-N1A	-13.88	118.27	128.89
3	A	900	FAD	C1B-N9A-C4A	-3.28	121.99	126.94
3	A	900	FAD	P-O3P-PA	-2.94	124.49	132.73
3	A	900	FAD	C4X-C10-N10	-2.62	118.97	120.52
3	A	900	FAD	C4A-C5A-N7A	-2.18	107.48	109.48
3	A	900	FAD	C6-C5X-C9A	2.11	121.76	118.98
3	A	900	FAD	O3P-P-O5'	2.21	108.79	102.94
3	A	900	FAD	C4-N3-C2	5.88	120.33	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	900	FAD	2	0

## 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	A	666/666 (100%)	1.16	107 (16%) <b>3</b> <b>2</b>	18, 70, 84, 105	0
2	B	133/235 (56%)	1.32	35 (26%) <b>1</b> <b>0</b>	59, 73, 96, 101	0
All	All	799/901 (88%)	1.19	142 (17%) <b>2</b> <b>1</b>	18, 70, 86, 105	0

All (142) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	172	SER	7.1
2	B	422	VAL	5.8
1	A	836	LEU	4.9
1	A	171	PRO	4.9
2	B	375	VAL	4.4
2	B	376	ILE	4.4
1	A	506	GLU	4.3
1	A	174	VAL	4.3
1	A	508	LEU	4.2
1	A	275	THR	4.2
1	A	238	LEU	3.9
1	A	504	LEU	3.8
2	B	402	PHE	3.6
2	B	309	LYS	3.5
1	A	490	LEU	3.5
1	A	242	TYR	3.5
2	B	308	ARG	3.4
1	A	273	LEU	3.4
1	A	431	TRP	3.4
2	B	437	TRP	3.4
1	A	492	LYS	3.3
1	A	241	PRO	3.3
1	A	497	LEU	3.3
1	A	503	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	451	LEU	3.3
2	B	378	LYS	3.3
1	A	173	GLY	3.2
2	B	374	GLU	3.2
2	B	399	GLY	3.2
1	A	403	ASN	3.2
1	A	373	GLU	3.2
1	A	372	LYS	3.1
1	A	231	PHE	3.0
1	A	814	ALA	3.0
1	A	739	ALA	3.0
1	A	441	LEU	2.9
2	B	420	PHE	2.9
1	A	809	ALA	2.9
2	B	365	GLY	2.9
1	A	268	LYS	2.9
2	B	371	ARG	2.9
1	A	496	GLU	2.8
1	A	239	GLU	2.8
2	B	353	LYS	2.8
2	B	360	LYS	2.8
2	B	426	ARG	2.8
2	B	312	LYS	2.8
1	A	494	TYR	2.8
1	A	429	GLU	2.8
2	B	414	VAL	2.8
1	A	232	GLU	2.8
1	A	445	LEU	2.8
1	A	499	GLU	2.8
2	B	415	VAL	2.8
1	A	212	PHE	2.7
1	A	455	ILE	2.7
2	B	359	LEU	2.7
1	A	509	GLN	2.7
1	A	740	VAL	2.7
1	A	444	LEU	2.7
1	A	377	MET	2.7
1	A	777	ALA	2.7
1	A	656	PHE	2.7
2	B	383	TRP	2.7
1	A	487	LEU	2.6
1	A	274	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	438	GLU	2.6
2	B	421	PHE	2.6
1	A	811	VAL	2.6
1	A	813	GLY	2.6
1	A	175	GLU	2.5
2	B	369	PRO	2.5
2	B	342	LEU	2.5
1	A	761	TYR	2.5
1	A	804	ILE	2.5
1	A	539	ALA	2.5
1	A	438	GLN	2.5
1	A	288	VAL	2.5
1	A	511	LEU	2.5
1	A	763	TYR	2.5
1	A	773	TYR	2.5
1	A	834	TYR	2.5
1	A	816	LEU	2.5
1	A	817	SER	2.5
1	A	380	GLN	2.4
1	A	498	ALA	2.4
2	B	418	LYS	2.4
2	B	368	GLU	2.4
1	A	449	VAL	2.4
1	A	803	THR	2.4
1	A	395	GLN	2.4
1	A	448	MET	2.4
1	A	808	PRO	2.3
1	A	471	PRO	2.3
1	A	799	ALA	2.3
1	A	655	GLY	2.2
1	A	775	LEU	2.2
1	A	289	SER	2.2
1	A	489	ALA	2.2
1	A	227	ILE	2.2
1	A	408	LEU	2.2
2	B	395	ILE	2.2
1	A	398	PHE	2.2
1	A	458	LEU	2.2
1	A	815	LEU	2.2
2	B	434	LEU	2.2
1	A	435	VAL	2.2
1	A	440	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	405	ILE	2.2
1	A	659	LEU	2.2
1	A	470	PRO	2.2
1	A	540	ASN	2.2
2	B	431	ASP	2.2
1	A	205	GLN	2.2
1	A	391	TYR	2.1
1	A	536	LEU	2.1
1	A	709	GLY	2.1
1	A	240	ALA	2.1
1	A	542	THR	2.1
1	A	287	GLY	2.1
1	A	194	ALA	2.1
2	B	367	ILE	2.1
2	B	419	ASN	2.1
1	A	368	GLN	2.1
1	A	330	GLY	2.1
1	A	437	THR	2.1
1	A	812	HIS	2.1
1	A	244	SER	2.1
1	A	447	LYS	2.1
2	B	400	ARG	2.1
1	A	290	GLY	2.1
1	A	657	GLY	2.1
1	A	743	PRO	2.1
1	A	778	GLN	2.0
1	A	760	SER	2.0
1	A	276	LYS	2.0
1	A	456	LYS	2.0
2	B	362	LYS	2.0
1	A	800	GLY	2.0
1	A	434	ILE	2.0
1	A	392	LEU	2.0
1	A	624	THR	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
6	NH4	A	1842	1/1	0.72	0.45	9.98	70,70,70,70	0
5	GOL	A	1840	6/6	0.77	0.23	0.25	113,114,115,115	0
5	GOL	A	1839	6/6	0.96	0.14	-2.40	61,62,64,65	0
3	FAD	A	900	53/53	0.98	0.10	-4.28	38,43,46,47	0
4	CL	A	1838	1/1	0.90	0.29	-	53,53,53,53	1
4	CL	A	1843	1/1	0.96	0.32	-	58,58,58,58	1
4	CL	A	1841	1/1	0.83	0.15	-	92,92,92,92	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.