



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:16 PM GMT

PDB ID : 4HSU
Title : Crystal structure of LSD2-NPAC with H3(1-26)in space group P21
Authors : Chen, F.; Dong, Z.; Fang, J.; Xu, Y.
Deposited on : 2012-10-30
Resolution : 1.99 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

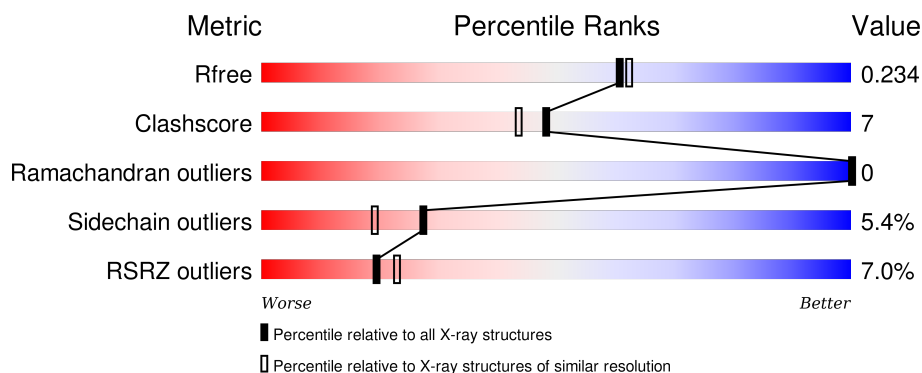
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	8664 (2.00-1.96)
Clashscore	102246	9905 (2.00-1.96)
Ramachandran outliers	100387	9792 (2.00-1.96)
Sidechain outliers	100360	9791 (2.00-1.96)
RSRZ outliers	91569	8679 (2.00-1.96)

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 ≥ 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	776	<div> <div>5%</div> <div>78%</div> <div>15%</div> <div>5%</div> </div>
2	B	124	<div> <div>4%</div> <div>7%</div> <div>90%</div> </div>
3	C	30	<div> <div>27%</div> <div>67%</div> <div>13%</div> <div>7%</div> <div>13%</div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6505 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 1B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	737	Total	C	N	O	S	0	0	0
			5837	3727	994	1075	41			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	PRO	-	EXPRESSION TAG	UNP Q8NB78
A	48	LEU	-	EXPRESSION TAG	UNP Q8NB78
A	49	GLY	-	EXPRESSION TAG	UNP Q8NB78
A	50	SER	-	EXPRESSION TAG	UNP Q8NB78

- Molecule 2 is a protein called Putative oxidoreductase GLYR1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	12	Total	C	N	O	0	0	0
			105	69	19	17			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	145	PRO	-	EXPRESSION TAG	UNP Q49A26
B	146	LEU	-	EXPRESSION TAG	UNP Q49A26
B	147	GLY	-	EXPRESSION TAG	UNP Q49A26
B	148	SER	-	EXPRESSION TAG	UNP Q49A26
B	149	PRO	-	EXPRESSION TAG	UNP Q49A26
B	150	GLU	-	EXPRESSION TAG	UNP Q49A26
B	151	PHE	-	EXPRESSION TAG	UNP Q49A26

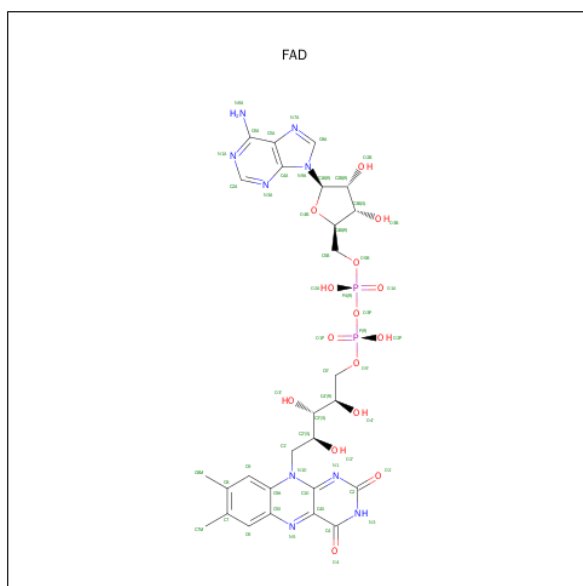
- Molecule 3 is a protein called Histone H3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	26	Total	C	N	O	S	0	0	0
			193	115	44	33	1			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	4	MET	LYS	ENGINEERED MUTATION	UNP Q92133

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	Zn	0	0
			3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	302	Total	O	0	0
			302	302		

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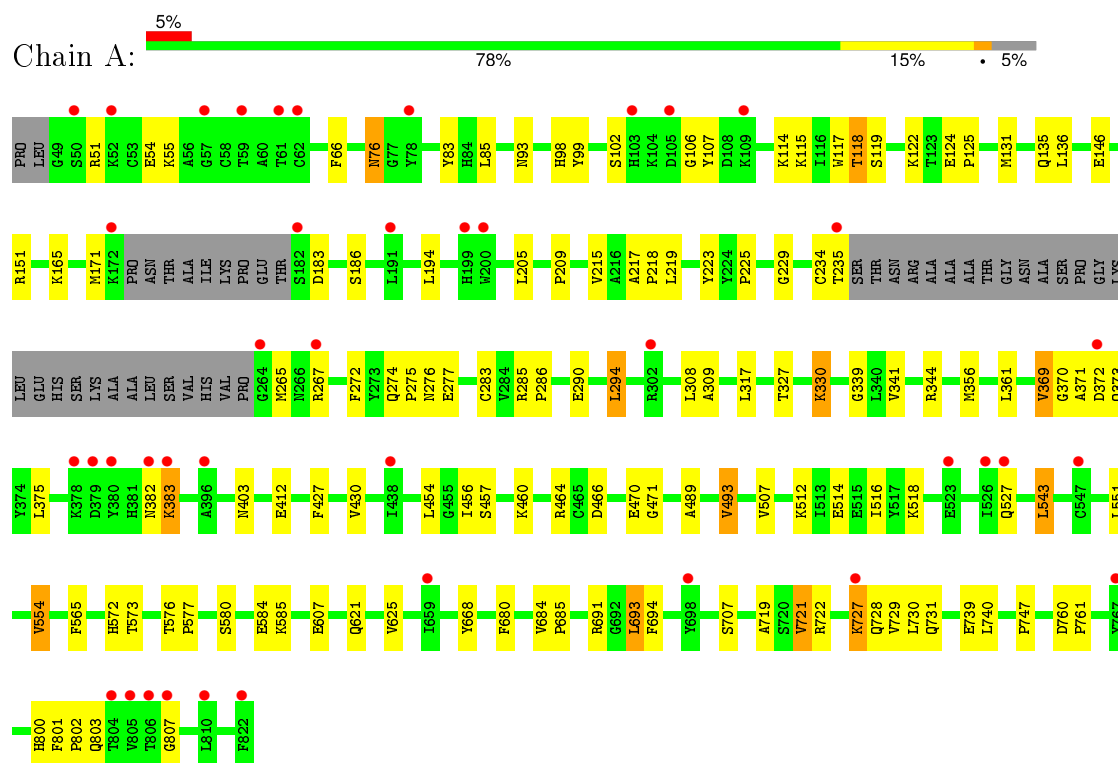
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total	O	0	0
			2	2		
6	C	10	Total	O	0	0
			10	10		

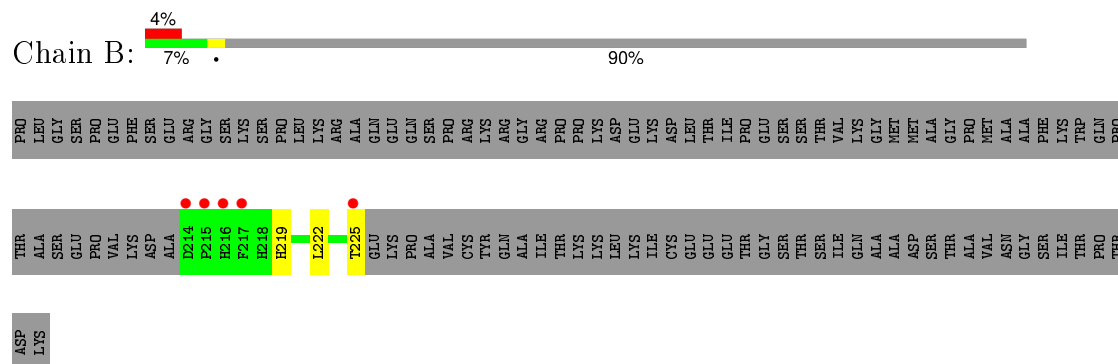
3 Residue-property plots [i](#)

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 ($RSRZ > 2$). 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.

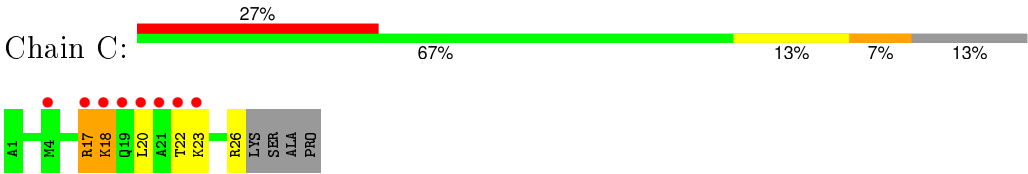
- Molecule 1: Lysine-specific histone demethylase 1B



- Molecule 2: Putative oxidoreductase GLYR1



- Molecule 3: Histone H3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.75Å 89.48Å 88.35Å 90.00° 102.89° 90.00°	Depositor
Resolution (Å)	40.44 – 1.99 40.43 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.0 (40.44-1.99) 99.1 (40.43-1.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.17 (at 1.98Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.199 , 0.232 0.204 , 0.234	Depositor DCC
R_{free} test set	3291 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	29.9	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 64866 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6505	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.87% 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, 0.375 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: ZN, FAD

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	0.31	0/5981	0.46	1/8100 (0.0%)
2	B	0.27	0/110	0.42	0/149
3	C	0.34	0/193	0.45	0/254
All	All	0.31	0/6284	0.46	1/8503 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	151	ARG	NE-CZ-NH1	-5.17	117.71	120.30

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	5837	0	5751	84	0
2	B	105	0	91	2	0
3	C	193	0	218	10	0
4	A	53	0	30	2	0
5	A	3	0	0	0	0
6	A	302	0	0	3	0
6	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	C	10	0	0	1	0
All	All	6505	0	6090	90	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 7.

All (90) 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:A:369:VAL:HG22	1:A:373:GLN:HB3	1.48	0.95
3:C:17:ARG:HD3	3:C:17:ARG:H	1.31	0.95
1:A:580:SER:O	1:A:584:GLU:HG2	1.69	0.92
1:A:727:LYS:HE3	1:A:731:GLN:HE22	1.37	0.90
1:A:727:LYS:HE3	1:A:731:GLN:NE2	1.89	0.87
3:C:17:ARG:HD3	3:C:17:ARG:N	1.91	0.85
1:A:209:PRO:HG3	1:A:341:VAL:HG21	1.60	0.84
1:A:493:VAL:HG22	1:A:516:ILE:HD13	1.63	0.80
1:A:327:THR:OG1	1:A:330:LYS:HG2	1.89	0.72
1:A:209:PRO:HG3	1:A:341:VAL:CG2	2.21	0.69
1:A:276:ASN:HA	3:C:20:LEU:O	1.94	0.67
1:A:76:ASN:HD22	1:A:76:ASN:H	1.41	0.67
1:A:668:TYR:CZ	1:A:747:PRO:HG3	2.32	0.65
1:A:382:ASN:O	1:A:383:LYS:HB3	1.95	0.65
1:A:727:LYS:HD3	1:A:728:GLN:N	2.13	0.64
1:A:719:ALA:O	1:A:722:ARG:HG2	1.98	0.63
1:A:285:ARG:HD3	3:C:20:LEU:HD13	1.84	0.59
1:A:99:TYR:CE2	1:A:131:MET:HE1	2.37	0.59
1:A:693:LEU:HD22	1:A:694:PHE:CE1	2.39	0.58
1:A:76:ASN:HD22	1:A:76:ASN:N	2.02	0.57
1:A:361:LEU:HD11	2:B:219:HIS:CE1	2.39	0.57
1:A:721:VAL:HG22	1:A:729:VAL:HG22	1.87	0.56
1:A:454:LEU:HD21	1:A:585:LYS:HG2	1.88	0.56
3:C:17:ARG:N	3:C:17:ARG:CD	2.61	0.54
1:A:514:GLU:OE1	1:A:518:LYS:HE3	2.07	0.54
1:A:223:TYR:CE2	1:A:275:PRO:HD3	2.44	0.53
1:A:122:LYS:HG2	1:A:122:LYS:O	2.09	0.53
1:A:727:LYS:HD3	1:A:728:GLN:HG3	1.90	0.53
1:A:217:ALA:HB3	1:A:218:PRO:HD3	1.91	0.53
1:A:456:ILE:HG23	1:A:577:PRO:HG2	1.92	0.51
1:A:551:LEU:HA	1:A:554:VAL:HG13	1.92	0.51
1:A:114:LYS:HG3	1:A:125:PRO:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:668:TYR:CE2	1:A:747:PRO:HG3	2.47	0.50
1:A:489:ALA:O	1:A:493:VAL:HG12	2.11	0.50
1:A:131:MET:HE1	1:A:135:GLN:HG3	1.94	0.50
1:A:572:HIS:HB2	6:C:108:HOH:O	2.12	0.49
3:C:18:LYS:NZ	3:C:18:LYS:HB3	2.27	0.49
1:A:576:THR:OG1	1:A:577:PRO:HD3	2.13	0.48
1:A:83:TYR:HE2	1:A:93:ASN:HD22	1.61	0.48
1:A:412:GLU:OE1	4:A:901:FAD:O3B	2.27	0.48
3:C:17:ARG:CD	3:C:17:ARG:H	2.00	0.47
1:A:146:GLU:H	1:A:146:GLU:CD	2.17	0.47
1:A:99:TYR:HD1	1:A:106:GLY:O	1.96	0.47
1:A:512:LYS:HD2	1:A:512:LYS:HA	1.69	0.47
1:A:685:PRO:HB3	1:A:691:ARG:HA	1.96	0.47
1:A:215:VAL:HG21	1:A:317:LEU:HD21	1.97	0.47
1:A:802:PRO:O	1:A:803:GLN:HB2	2.14	0.47
1:A:383:LYS:HA	1:A:621:GLN:OE1	2.15	0.46
1:A:136:LEU:HD13	1:A:339:GLY:HA3	1.98	0.45
1:A:183:ASP:OD2	1:A:186:SER:HB3	2.16	0.45
1:A:131:MET:CE	1:A:135:GLN:HG3	2.47	0.45
1:A:285:ARG:HD3	3:C:20:LEU:CD1	2.45	0.45
1:A:507:VAL:HG21	1:A:512:LYS:HD3	1.98	0.45
1:A:493:VAL:CG2	1:A:516:ILE:HD13	2.43	0.44
1:A:225:PRO:HD3	3:C:26:ARG:NH2	2.32	0.44
1:A:117:TRP:CD1	1:A:125:PRO:HG3	2.53	0.44
1:A:76:ASN:ND2	1:A:76:ASN:H	2.11	0.44
1:A:285:ARG:HA	1:A:286:PRO:HD3	1.86	0.43
3:C:20:LEU:HD23	3:C:20:LEU:HA	1.73	0.43
1:A:470:GLU:HA	1:A:471:GLY:HA2	1.62	0.43
1:A:102:SER:HA	1:A:107:TYR:CD1	2.53	0.43
1:A:370:GLY:O	1:A:373:GLN:HB2	2.19	0.43
1:A:565:PHE:CD1	1:A:800:HIS:CE1	3.06	0.43
1:A:801:PHE:O	1:A:807:GLY:HA3	2.18	0.43
1:A:219:LEU:HD11	1:A:265:MET:HE1	2.01	0.43
1:A:371:ALA:O	1:A:372:ASP:HB2	2.18	0.43
1:A:464:ARG:NH1	1:A:466:ASP:OD2	2.53	0.42
1:A:234:CYS:O	1:A:235:THR:HG22	2.19	0.42
1:A:171:MET:HE1	6:A:1396:HOH:O	2.20	0.42
1:A:290:GLU:O	1:A:294:LEU:HD22	2.19	0.42
1:A:607:GLU:HA	6:A:1175:HOH:O	2.18	0.42
1:A:803:GLN:O	4:A:901:FAD:O3'	2.37	0.41
1:A:471:GLY:HA2	1:A:739:GLU:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:684:VAL:HA	1:A:685:PRO:HD3	1.85	0.41
1:A:272:PHE:HB3	1:A:286:PRO:HG3	2.03	0.41
1:A:456:ILE:CG2	1:A:577:PRO:HG2	2.50	0.41
1:A:290:GLU:HG3	2:B:222:LEU:HD13	2.02	0.41
1:A:427:PHE:HB2	6:A:1155:HOH:O	2.21	0.41
1:A:115:LYS:HA	1:A:118:THR:OG1	2.20	0.41
1:A:375:LEU:HD12	1:A:403:ASN:HB3	2.02	0.41
1:A:98:HIS:CD2	1:A:106:GLY:HA3	2.57	0.40
1:A:54:GLU:HG3	1:A:55:LYS:N	2.36	0.40
1:A:760:ASP:HA	1:A:761:PRO:HD3	1.88	0.40
1:A:693:LEU:HA	1:A:693:LEU:HD23	1.91	0.40
1:A:124:GLU:HA	1:A:125:PRO:HD3	1.86	0.40
1:A:229:GLY:HA3	1:A:309:ALA:HB2	2.02	0.40
1:A:274:GLN:O	1:A:277:GLU:HB2	2.21	0.40
1:A:543:LEU:HB3	1:A:551:LEU:HD21	2.04	0.40
1:A:460:LYS:HA	1:A:573:THR:HG22	2.04	0.40
1:A:66:PHE:CD2	1:A:66:PHE:C	2.94	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	731/776 (94%)	704 (96%)	27 (4%)	0	100	100
2	B	10/124 (8%)	10 (100%)	0	0	100	100
3	C	24/30 (80%)	22 (92%)	2 (8%)	0	100	100
All	All	765/930 (82%)	736 (96%)	29 (4%)	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	A	632/662 (96%)	601 (95%)	31 (5%)	31	22
2	B	12/106 (11%)	11 (92%)	1 (8%)	14	7
3	C	18/21 (86%)	14 (78%)	4 (22%)	1	0
All	All	662/789 (84%)	626 (95%)	36 (5%)	27	19

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	ARG
1	A	76	ASN
1	A	85	LEU
1	A	118	THR
1	A	119	SER
1	A	165	LYS
1	A	194	LEU
1	A	205	LEU
1	A	267	ARG
1	A	283	CYS
1	A	294	LEU
1	A	308	LEU
1	A	330	LYS
1	A	344	ARG
1	A	356	MET
1	A	369	VAL
1	A	383	LYS
1	A	430	VAL
1	A	457	SER
1	A	493	VAL
1	A	527	GLN
1	A	543	LEU
1	A	554	VAL
1	A	625	VAL
1	A	680	PHE
1	A	693	LEU

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Mol	Chain	Res	Type
1	A	707	SER
1	A	721	VAL
1	A	727	LYS
1	A	730	LEU
1	A	740	LEU
2	B	225	THR
3	C	17	ARG
3	C	18	LYS
3	C	22	THR
3	C	23	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	ASN
1	A	93	ASN
1	A	98	HIS
1	A	134	GLN
1	A	276	ASN
1	A	400	GLN
1	A	440	ASN
1	A	527	GLN
1	A	731	GLN
2	B	218	HIS

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 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 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
4	FAD	A	901	-	48,58,58	2.34	23 (47%)	54,89,89	2.15	13 (24%)

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
4	FAD	A	901	-	-	0/30/50/50	0/6/6/6

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	FAD	C10-N10	-5.09	1.33	1.39
4	A	901	FAD	PA-O5B	-4.36	1.39	1.59
4	A	901	FAD	O4B-C1B	-4.19	1.35	1.41
4	A	901	FAD	P-O2P	-4.09	1.37	1.54
4	A	901	FAD	PA-O1A	-3.89	1.37	1.51
4	A	901	FAD	PA-O2A	-3.85	1.38	1.54
4	A	901	FAD	C6-C5X	-3.59	1.36	1.41
4	A	901	FAD	P-O1P	-3.48	1.38	1.51
4	A	901	FAD	C5A-N7A	-2.96	1.29	1.39
4	A	901	FAD	C5A-C4A	-2.90	1.33	1.40
4	A	901	FAD	O3'-C3'	-2.79	1.36	1.43
4	A	901	FAD	C4A-N3A	-2.76	1.31	1.35
4	A	901	FAD	C9A-C5X	-2.73	1.37	1.42
4	A	901	FAD	P-O5'	-2.65	1.46	1.59
4	A	901	FAD	C4-C4X	-2.51	1.36	1.41
4	A	901	FAD	C2A-N1A	-2.44	1.29	1.33
4	A	901	FAD	C4X-C10	-2.42	1.36	1.41
4	A	901	FAD	O5B-C5B	-2.41	1.34	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	FAD	C1'-N10	-2.38	1.45	1.48
4	A	901	FAD	O4'-C4'	-2.32	1.38	1.43
4	A	901	FAD	C8A-N7A	-2.30	1.30	1.34
4	A	901	FAD	C9-C9A	-2.24	1.35	1.40
4	A	901	FAD	C6A-N6A	-2.14	1.28	1.34

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	FAD	N3A-C2A-N1A	-8.06	122.73	128.89
4	A	901	FAD	C4A-C5A-N7A	-3.00	106.72	109.48
4	A	901	FAD	C4X-C4-N3	-2.58	120.06	123.59
4	A	901	FAD	O4'-C4'-C5'	-2.46	104.83	110.19
4	A	901	FAD	C4B-O4B-C1B	-2.30	107.19	109.72
4	A	901	FAD	C4-C4X-C10	-2.25	118.50	119.94
4	A	901	FAD	O2A-PA-O3P	-2.14	95.39	105.09
4	A	901	FAD	C1B-N9A-C4A	-2.07	123.82	126.94
4	A	901	FAD	C1'-N10-C9A	3.72	123.04	118.86
4	A	901	FAD	C5X-C9A-N10	3.90	120.58	117.62
4	A	901	FAD	O3P-PA-O5B	4.33	114.42	102.94
4	A	901	FAD	O3P-P-O5'	5.51	117.56	102.94
4	A	901	FAD	C4-N3-C2	5.77	120.24	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
4	A	901	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, 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	737/776 (94%)	0.29	41 (5%) 28 32	23, 35, 60, 77	0
2	B	12/124 (9%)	1.51	5 (41%) 0 1	34, 45, 62, 67	0
3	C	26/30 (86%)	1.37	8 (30%) 1 1	31, 38, 67, 72	0
All	All	775/930 (83%)	0.35	54 (6%) 19 23	23, 35, 61, 77	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	21	ALA	6.6
1	A	822	PHE	6.0
3	C	20	LEU	5.2
1	A	50	SER	4.9
2	B	225	THR	4.9
1	A	527	GLN	4.7
1	A	382	ASN	4.5
1	A	200	TRP	4.0
1	A	182	SER	3.9
1	A	526	ILE	3.8
3	C	18	LYS	3.6
1	A	199	HIS	3.5
1	A	302	ARG	3.5
3	C	19	GLN	3.4
1	A	804	THR	3.4
3	C	17	ARG	3.3
1	A	103	HIS	3.3
2	B	214	ASP	3.2
1	A	379	ASP	3.2
3	C	23	LYS	3.1
2	B	215	PRO	3.1
1	A	62	CYS	3.1
1	A	264	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	438	ILE	3.1
1	A	383	LYS	3.0
1	A	805	VAL	3.0
1	A	172	LYS	3.0
1	A	523	GLU	3.0
1	A	378	LYS	2.9
1	A	235	THR	2.9
2	B	217	PHE	2.8
1	A	372	ASP	2.8
1	A	59	THR	2.8
3	C	22	THR	2.7
1	A	105	ASP	2.7
1	A	396	ALA	2.6
1	A	767	TYR	2.6
1	A	61	THR	2.6
1	A	806	THR	2.6
1	A	727	LYS	2.5
1	A	109	LYS	2.5
1	A	698	TYR	2.5
1	A	78	TYR	2.5
1	A	57	GLY	2.4
1	A	52	LYS	2.4
2	B	216	HIS	2.4
1	A	810	LEU	2.3
1	A	380	TYR	2.3
1	A	807	GLY	2.3
1	A	191	LEU	2.3
1	A	547	CYS	2.2
3	C	4	MET	2.1
1	A	267	ARG	2.0
1	A	659	ILE	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](#)

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, 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	LLDF	B-factors(Å ²)	Q<0.9
5	ZN	A	904	1/1	0.99	0.10	0.28	39,39,39,39	0
4	FAD	A	901	53/53	0.95	0.14	-0.32	20,25,30,31	0
5	ZN	A	902	1/1	0.99	0.08	-1.08	42,42,42,42	0
5	ZN	A	903	1/1	0.95	0.08	-1.23	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.