Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 03:37 am BST

PDB ID	:	6S35
Title	:	LSD1/CoREST1 complex with macrocyclic peptide inhibitor
Authors	:	Talibov, V.O.; Dobritzsch, D.
Deposited on	:	2019-06-24
Resolution	:	3.10 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 following versions of software and data (see references ⓘ) were used in the production of this report:

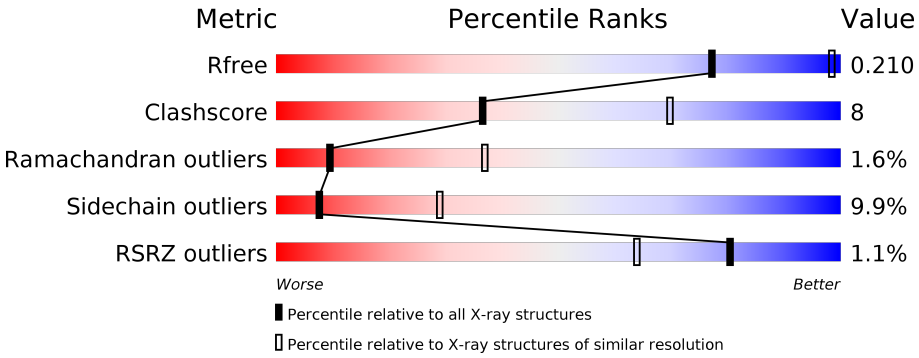
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance

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

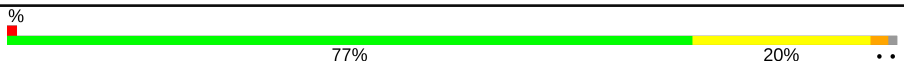
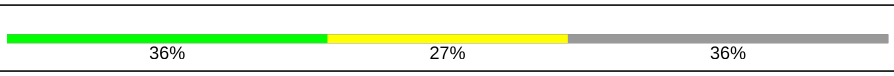
The reported resolution of this entry is 3.10 Å.

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}	130704	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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 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 <=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	
2	B	182	
3	C	11	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6376 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 1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	662	Total	C	N	O	S	0	0	0
			5183	3300	902	961	20			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	168	GLY	-	expression tag	UNP O60341
A	169	SER	-	expression tag	UNP O60341
A	170	HIS	-	expression tag	UNP O60341
A	171	MET	-	expression tag	UNP O60341

- Molecule 2 is a protein called REST corepressor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	134	Total	C	N	O	S	0	0	0
			1085	682	196	204	3			

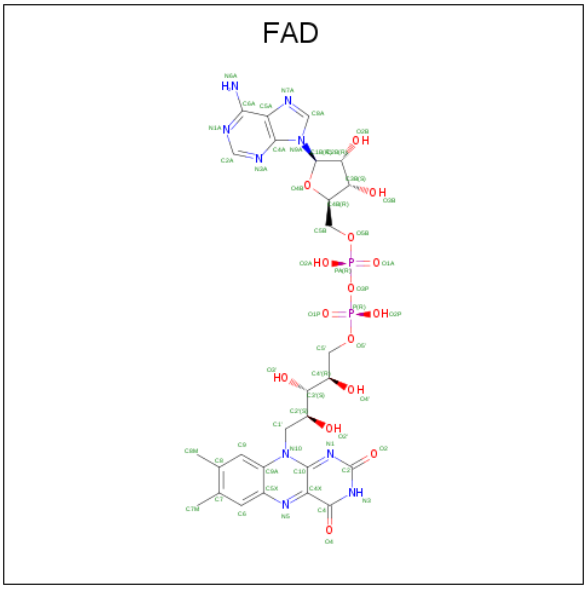
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	304	GLY	-	expression tag	UNP Q9UKL0
B	305	SER	-	expression tag	UNP Q9UKL0
B	306	HIS	-	expression tag	UNP Q9UKL0
B	307	MET	-	expression tag	UNP Q9UKL0

- Molecule 3 is a protein called ALA-ARG-(D)LYS-MET-GLN-GLU-ALA-ARG-LYS-SER-THR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	7	Total	C	N	O	S	0	0	0
			55	33	12	9	1			

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

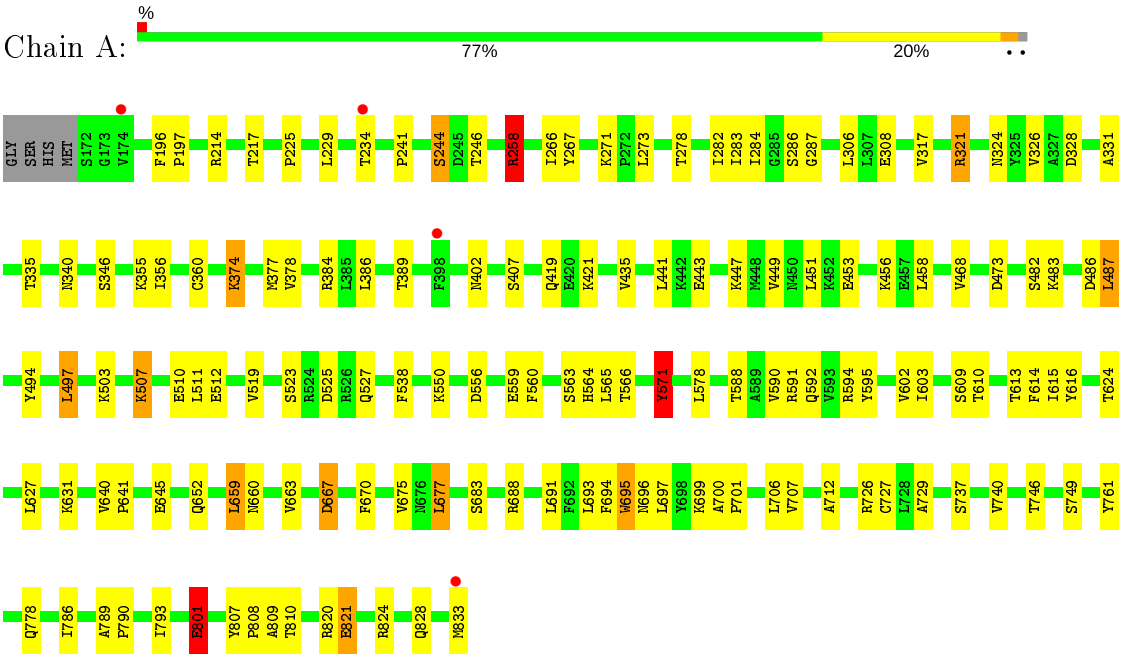


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	A	1	53	27	9	15	2	0	0

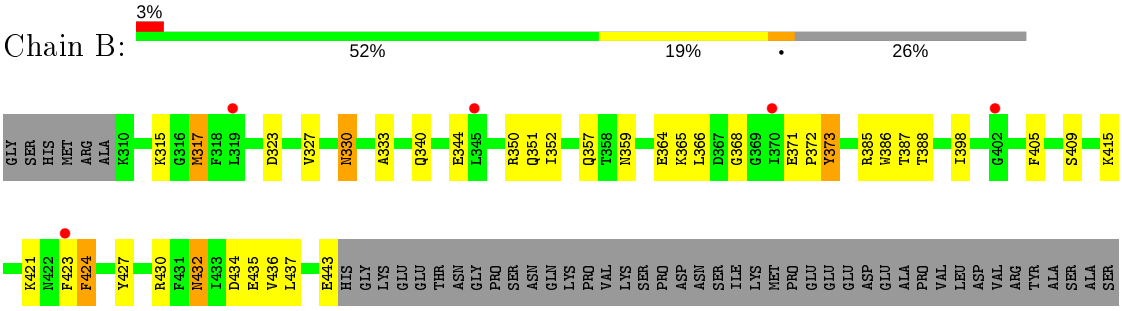
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

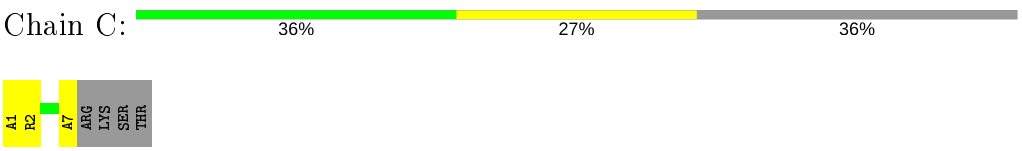
- Molecule 1: Lysine-specific histone demethylase 1A



- Molecule 2: REST corepressor 1



- Molecule 3: ALA-ARG-(D)LYS-MET-GLN-GLU-ALA-ARG-LYS-SER-THR



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	120.55Å 179.50Å 234.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	117.21 – 3.10 117.21 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.5 (117.21-3.10) 99.5 (117.21-3.10)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.71 (at 3.13Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.183 , 0.213 0.185 , 0.210	Depositor DCC
R_{free} test set	2300 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å^2)	91.1	Xtriage
Anisotropy	0.615	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 75.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6376	wwPDB-VP
Average B, all atoms (Å^2)	112.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: DLY, 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	0.72	1/5295 (0.0%)	0.91	4/7182 (0.1%)
2	B	0.67	0/1100	0.87	0/1482
3	C	0.66	0/44	1.00	0/54
All	All	0.71	1/6439 (0.0%)	0.90	4/8718 (0.0%)

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
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	833	MET	C-O	17.10	1.55	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	571	TYR	CB-CG-CD1	6.22	124.73	121.00
1	A	258	ARG	CB-CA-C	-5.67	99.06	110.40
1	A	820	ARG	NE-CZ-NH2	5.60	123.10	120.30
1	A	820	ARG	NE-CZ-NH1	-5.51	117.55	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	695	TRP	Peptide

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	5183	0	5217	82	0
2	B	1085	0	1104	26	0
3	C	55	0	57	2	0
4	A	53	0	31	4	0
All	All	6376	0	6409	101	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 8.

All (101) 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:308:GLU:OE2	4:A:901:FAD:O3B	1.97	0.81
1:A:360:CYS:O	3:C:1:ALA:N	2.19	0.76
1:A:451:LEU:HD23	1:A:494:TYR:HB2	1.68	0.75
1:A:456:LYS:HA	2:B:373:TYR:CE1	2.23	0.73
1:A:456:LYS:HA	2:B:373:TYR:HE1	1.55	0.70
1:A:538:PHE:CE1	1:A:706:LEU:HD13	2.30	0.66
1:A:700:ALA:HB1	1:A:701:PRO:HD2	1.78	0.66
1:A:435:VAL:HG23	2:B:352:ILE:HG13	1.77	0.65
2:B:398:ILE:HG22	2:B:436:VAL:CG1	2.29	0.63
1:A:266:ILE:HD11	1:A:578:LEU:HD23	1.81	0.62
1:A:801:GLU:HG2	1:A:809:ALA:HA	1.83	0.60
2:B:330:ASN:ND2	2:B:333:ALA:HB2	2.15	0.60
1:A:494:TYR:CD2	1:A:494:TYR:O	2.55	0.59
1:A:693:LEU:HD12	1:A:694:PHE:H	1.67	0.59
2:B:327:VAL:O	2:B:327:VAL:HG12	2.02	0.59
1:A:331:ALA:HA	4:A:901:FAD:N5	2.16	0.58
1:A:355:LYS:HA	1:A:565:LEU:HD23	1.85	0.58
2:B:405:PHE:CD2	2:B:421:LYS:HE2	2.38	0.58
1:A:273:LEU:O	1:A:273:LEU:HD12	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:801:GLU:CG	1:A:809:ALA:HA	2.33	0.58
1:A:821:GLU:OE2	1:A:824:ARG:NH1	2.37	0.57
2:B:386:TRP:CZ2	2:B:423:PHE:HB2	2.40	0.57
1:A:331:ALA:HA	4:A:901:FAD:C4X	2.36	0.56
1:A:808:PRO:O	1:A:810:THR:HG23	2.06	0.56
1:A:659:LEU:HD23	1:A:660:ASN:N	2.21	0.55
2:B:398:ILE:HG22	2:B:436:VAL:HG11	1.88	0.55
1:A:435:VAL:HG23	2:B:352:ILE:CG1	2.37	0.55
1:A:419:GLN:NE2	2:B:317:MET:HA	2.23	0.54
1:A:447:LYS:HE3	1:A:497:LEU:HD11	1.90	0.54
1:A:229:LEU:HD21	1:A:234:THR:CG2	2.37	0.53
1:A:595:TYR:CZ	1:A:641:PRO:HD2	2.43	0.53
1:A:360:CYS:HB2	1:A:677:LEU:HD13	1.90	0.53
1:A:229:LEU:HD21	1:A:234:THR:HG22	1.91	0.52
2:B:340:GLN:O	2:B:340:GLN:HG3	2.10	0.52
1:A:707:VAL:HG12	1:A:712:ALA:HA	1.92	0.52
1:A:789:ALA:O	1:A:790:PRO:C	2.48	0.52
1:A:374:LYS:HA	1:A:374:LYS:HE3	1.91	0.52
1:A:441:LEU:HD23	2:B:359:ASN:HD22	1.75	0.51
1:A:695:TRP:HE1	1:A:706:LEU:HD11	1.76	0.51
1:A:473:ASP:OD1	1:A:473:ASP:C	2.50	0.50
1:A:807:TYR:N	1:A:808:PRO:CD	2.75	0.49
2:B:409:SER:OG	2:B:415:LYS:O	2.20	0.49
2:B:424:PHE:HE2	2:B:437:LEU:HD11	1.76	0.49
1:A:603:ILE:HG12	1:A:615:ILE:HD13	1.95	0.49
1:A:507:LYS:O	1:A:511:LEU:HG	2.13	0.49
1:A:258:ARG:HG2	1:A:258:ARG:HH11	1.78	0.48
1:A:793:ILE:HB	1:A:828:GLN:NE2	2.28	0.48
1:A:284:ILE:HG12	1:A:590:VAL:HG21	1.94	0.48
1:A:761:TYR:CD1	1:A:809:ALA:HB1	2.48	0.48
1:A:321:ARG:HG3	1:A:326:VAL:HG22	1.94	0.48
2:B:340:GLN:O	2:B:340:GLN:CG	2.61	0.48
2:B:330:ASN:HD22	2:B:333:ALA:HB2	1.78	0.47
1:A:663:VAL:O	1:A:746:THR:HA	2.14	0.47
1:A:660:ASN:HA	1:A:749:SER:OG	2.13	0.47
1:A:559:GLU:OE1	3:C:7:ALA:HB1	2.15	0.47
1:A:696:ASN:N	1:A:696:ASN:OD1	2.46	0.47
1:A:258:ARG:NH1	1:A:258:ARG:HG2	2.30	0.47
1:A:419:GLN:HE22	2:B:317:MET:HA	1.78	0.46
1:A:670:PHE:CE1	1:A:740:VAL:HG22	2.50	0.46
1:A:196:PHE:N	1:A:197:PRO:CD	2.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:LYS:HE2	1:A:525:ASP:OD1	2.16	0.46
1:A:801:GLU:HG2	1:A:809:ALA:CA	2.44	0.46
1:A:523:SER:O	1:A:527:GLN:NE2	2.49	0.45
1:A:659:LEU:C	1:A:659:LEU:CD2	2.85	0.45
1:A:726:ARG:O	1:A:729:ALA:HB3	2.16	0.45
1:A:627:LEU:O	1:A:631:LYS:HG3	2.16	0.45
1:A:374:LYS:O	1:A:378:VAL:HG23	2.16	0.45
1:A:494:TYR:C	1:A:494:TYR:CD2	2.90	0.45
1:A:592:GLN:HB3	1:A:603:ILE:HD12	1.98	0.45
1:A:282:ILE:HG21	1:A:602:VAL:HG21	1.98	0.45
1:A:667:ASP:N	1:A:667:ASP:OD1	2.48	0.45
1:A:384:ARG:NH2	2:B:315:LYS:O	2.50	0.45
1:A:241:PRO:O	1:A:244:SER:HB3	2.18	0.44
1:A:266:ILE:CD1	1:A:578:LEU:HD23	2.46	0.44
1:A:286:SER:HB3	1:A:308:GLU:OE1	2.18	0.43
1:A:283:ILE:HB	1:A:306:LEU:HD23	1.99	0.43
1:A:458:LEU:HD23	1:A:487:LEU:HD12	2.00	0.43
2:B:386:TRP:O	2:B:387:THR:HG23	2.18	0.43
1:A:786:ILE:HD12	1:A:786:ILE:H	1.83	0.43
1:A:614:PHE:HD1	1:A:616:TYR:HH	1.67	0.43
1:A:821:GLU:OE2	1:A:821:GLU:HA	2.18	0.43
2:B:405:PHE:CE2	2:B:421:LYS:HE2	2.54	0.43
1:A:356:ILE:HD11	1:A:566:THR:HG23	2.00	0.43
1:A:659:LEU:C	1:A:659:LEU:HD23	2.39	0.42
2:B:427:TYR:CD1	2:B:430:ARG:NH2	2.87	0.42
1:A:340:ASN:HB2	1:A:560:PHE:CD2	2.54	0.42
1:A:317:VAL:HG13	1:A:571:TYR:HB3	2.00	0.42
1:A:283:ILE:O	1:A:306:LEU:HA	2.20	0.42
1:A:321:ARG:HH11	1:A:321:ARG:HG2	1.85	0.41
1:A:386:LEU:O	1:A:389:THR:N	2.49	0.41
2:B:398:ILE:HG22	2:B:436:VAL:HG12	2.00	0.41
2:B:371:GLU:HB3	2:B:372:PRO:HD3	2.01	0.41
1:A:217:THR:HG23	1:A:234:THR:HG21	2.02	0.41
1:A:691:LEU:HD22	1:A:727:CYS:SG	2.60	0.41
1:A:693:LEU:HD12	1:A:694:PHE:N	2.32	0.41
2:B:405:PHE:CD2	2:B:421:LYS:CE	3.04	0.41
1:A:321:ARG:CG	1:A:326:VAL:HG22	2.51	0.40
1:A:640:VAL:HA	1:A:641:PRO:HA	1.82	0.40
1:A:386:LEU:HA	1:A:386:LEU:HD23	1.95	0.40
4:A:901:FAD:HO3A	4:A:901:FAD:HO2A	1.68	0.40
2:B:327:VAL:O	2:B:327:VAL:CG1	2.67	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	660/666 (99%)	597 (90%)	54 (8%)	9 (1%)	11 40
2	B	132/182 (72%)	110 (83%)	18 (14%)	4 (3%)	4 23
3	C	4/11 (36%)	4 (100%)	0	0	100 100
All	All	796/859 (93%)	711 (89%)	72 (9%)	13 (2%)	9 37

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	801	GLU
1	A	287	GLY
1	A	468	VAL
1	A	699	LYS
2	B	432	ASN
1	A	483	LYS
2	B	344	GLU
1	A	244	SER
1	A	486	ASP
2	B	434	ASP
1	A	225	PRO
1	A	482	SER
2	B	368	GLY

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	562/565 (100%)	511 (91%)	51 (9%)	9 33
2	B	118/159 (74%)	102 (86%)	16 (14%)	3 16
3	C	4/8 (50%)	3 (75%)	1 (25%)	0 2
All	All	684/732 (93%)	616 (90%)	68 (10%)	8 29

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	214	ARG
1	A	246	THR
1	A	258	ARG
1	A	267	TYR
1	A	271	LYS
1	A	278	THR
1	A	321	ARG
1	A	324	ASN
1	A	328	ASP
1	A	335	THR
1	A	346	SER
1	A	374	LYS
1	A	377	MET
1	A	402	ASN
1	A	407	SER
1	A	421	LYS
1	A	443	GLU
1	A	449	VAL
1	A	453	GLU
1	A	487	LEU
1	A	497	LEU
1	A	503	LYS
1	A	507	LYS
1	A	510	GLU
1	A	512	GLU
1	A	519	VAL
1	A	550	LYS
1	A	556	ASP
1	A	563	SER
1	A	564	HIS
1	A	571	TYR
1	A	588	THR

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Mol	Chain	Res	Type
1	A	591	ARG
1	A	594	ARG
1	A	609	SER
1	A	610	THR
1	A	613	THR
1	A	624	THR
1	A	645	GLU
1	A	652	GLN
1	A	659	LEU
1	A	667	ASP
1	A	675	VAL
1	A	677	LEU
1	A	683	SER
1	A	688	ARG
1	A	697	LEU
1	A	737	SER
1	A	778	GLN
1	A	801	GLU
1	A	821	GLU
2	B	317	MET
2	B	323	ASP
2	B	330	ASN
2	B	350	ARG
2	B	351	GLN
2	B	357	GLN
2	B	364	GLU
2	B	365	LYS
2	B	366	LEU
2	B	373	TYR
2	B	385	ARG
2	B	388	THR
2	B	424	PHE
2	B	432	ASN
2	B	435	GLU
2	B	443	GLU
3	C	2	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	243	ASN
1	A	430	HIS

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Mol	Chain	Res	Type
1	A	527	GLN
1	A	587	ASN
1	A	742	GLN
1	A	771	ASN
1	A	828	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)

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
3	DLY	C	3	3	7,8,9	0.51	0	3,8,10	0.08	0

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
3	DLY	C	3	3	-	1/6/7/9	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	3	DLY	CE-CD-CG-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	$\#\|Z\| > 2$
4	FAD	A	901	-	51,58,58	1.08	2 (3%)	60,89,89	1.89	10 (16%)

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	-	-	6/30/50/50	0/6/6/6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	FAD	C4X-C10	5.53	1.44	1.38
4	A	901	FAD	C4-N3	2.61	1.37	1.33

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	901	FAD	C4-N3-C2	8.16	122.03	115.14
4	A	901	FAD	C10-C4X-N5	5.64	125.16	121.26
4	A	901	FAD	C4X-C10-N10	-4.18	116.01	120.30
4	A	901	FAD	C4-C4X-C10	-4.02	117.29	119.95
4	A	901	FAD	C4X-C4-N3	-3.71	118.36	123.43
4	A	901	FAD	O4B-C1B-C2B	-2.92	102.65	106.93
4	A	901	FAD	O5'-C5'-C4'	2.83	116.91	109.36
4	A	901	FAD	C1'-N10-C9A	2.37	120.16	118.29
4	A	901	FAD	C5'-C4'-C3'	-2.17	108.01	112.20
4	A	901	FAD	O2A-PA-O1A	2.09	122.58	112.24

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	901	FAD	O4'-C4'-C5'-O5'
4	A	901	FAD	PA-O3P-P-O5'
4	A	901	FAD	C3'-C4'-C5'-O5'
4	A	901	FAD	C5'-O5'-P-O1P
4	A	901	FAD	O4B-C4B-C5B-O5B
4	A	901	FAD	C1'-C2'-C3'-O3'

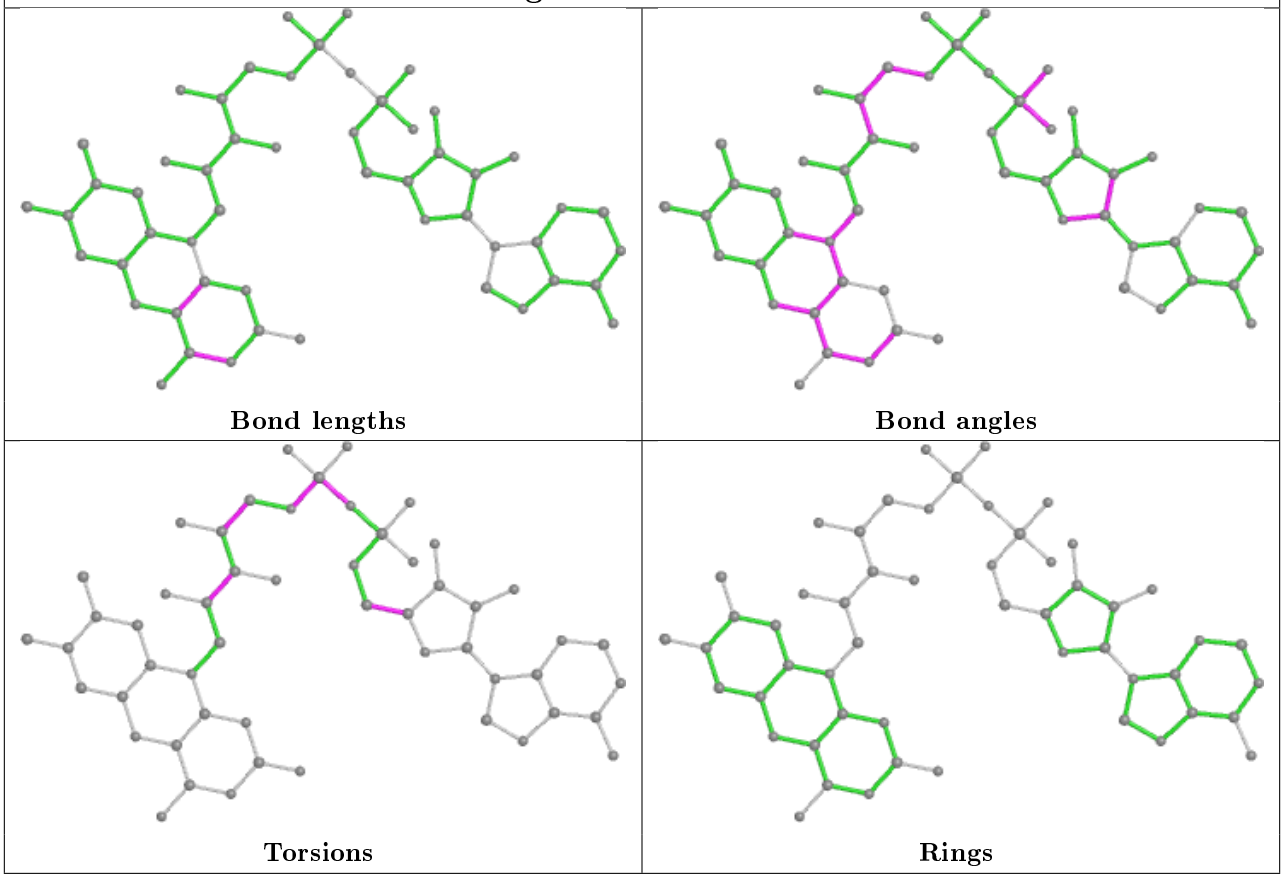
There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	901	FAD	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand FAD A 901



5.7 Other polymers

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, 95^{th} 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	662/666 (99%)	0.57	4 (0%) 89 78	63, 104, 147, 190	0
2	B	134/182 (73%)	0.46	5 (3%) 41 21	99, 136, 166, 187	0
3	C	6/11 (54%)	0.87	0 100 100	121, 135, 143, 146	0
All	All	802/859 (93%)	0.56	9 (1%) 80 64	63, 112, 153, 190	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	174	VAL	3.2
1	A	398	PHE	3.2
2	B	402	GLY	2.3
2	B	319	LEU	2.3
2	B	370	ILE	2.2
2	B	423	PHE	2.1
1	A	833	MET	2.0
1	A	234	THR	2.0
2	B	345	LEU	2.0

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

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, 95^{th} 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
3	DLY	C	3	9/10	0.96	0.40	129,130,134,137	0

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

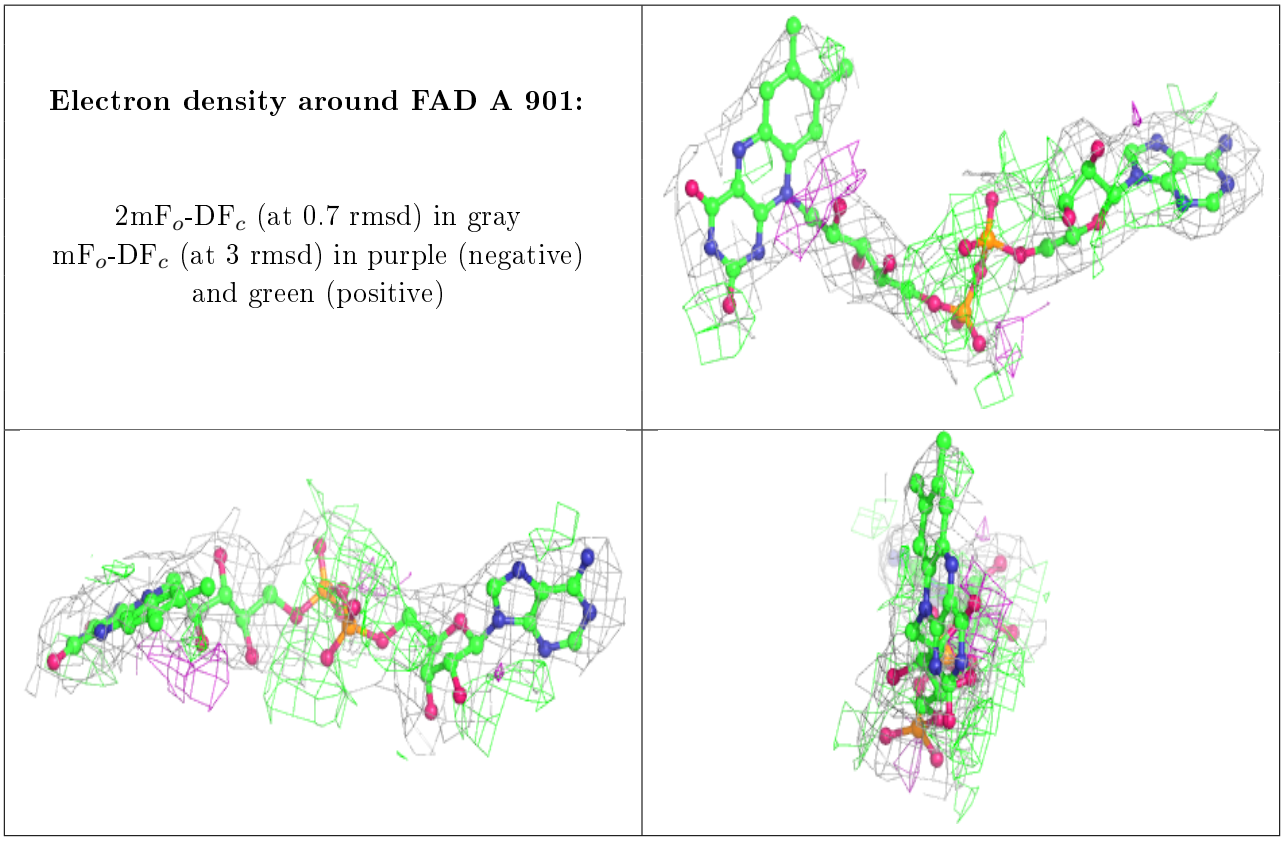
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
4	FAD	A	901	53/53	0.98	0.28	61,78,98,100	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around FAD A 901:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.