



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:49 PM UTC

PDB ID : 4XBF / pdb_00004xbf
Title : Structure of LSD1:CoREST in complex with ssRNA
Authors : Luka, Z.; Loukachevitch, L.V.; Martin, W.J.; Wagner, C.; Reiter, N.J.
Deposited on : 2014-12-16
Resolution : 2.80 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 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

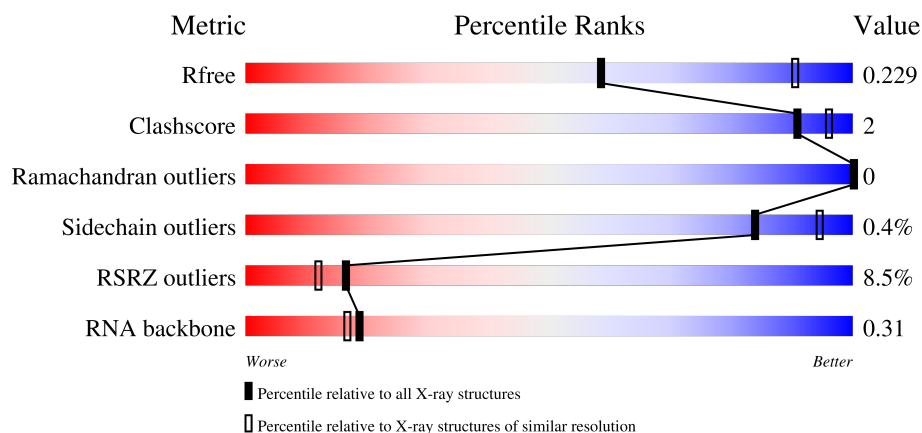
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)
RNA backbone	3983	1114 (3.00-2.60)

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	666	6% 96% .
2	B	133	19% 96% ..
3	D	5	20% 100%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 12911 atoms, of which 6425 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	666	Total	C	H	N	O	S	0	0	0
			10469	3324	5252	906	967	20			

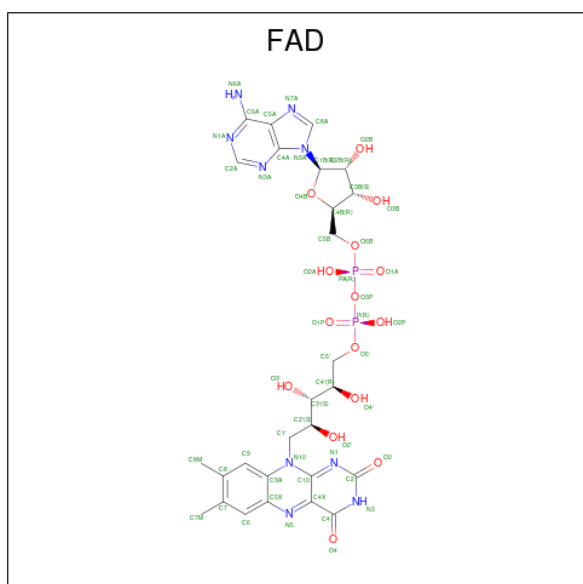
- Molecule 2 is a protein called REST corepressor 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	133	Total	C	H	N	O	S	0	0	0
			2167	676	1091	194	203	3			

- Molecule 3 is a RNA chain called RNA (5'-R(*UP*UP*AP*GP*G)-3').

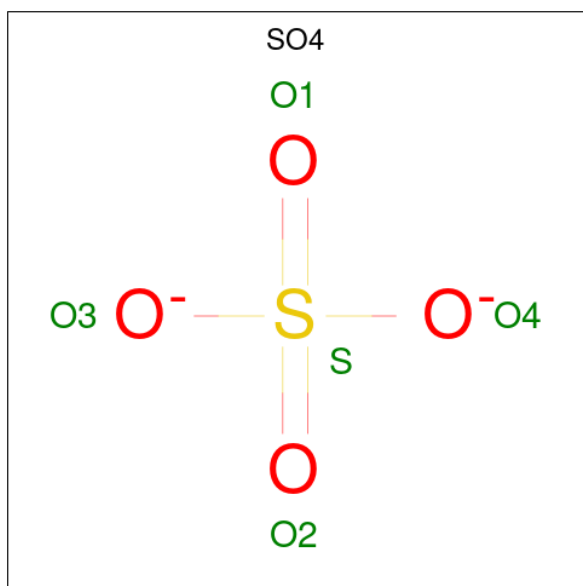
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	D	5	Total	C	H	N	O	P	0	0	0
			129	38	43	14	30	4			

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



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	P	0	0
			84	27	31	9	15	2		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	H	O	0	0
			14	3	8	3		

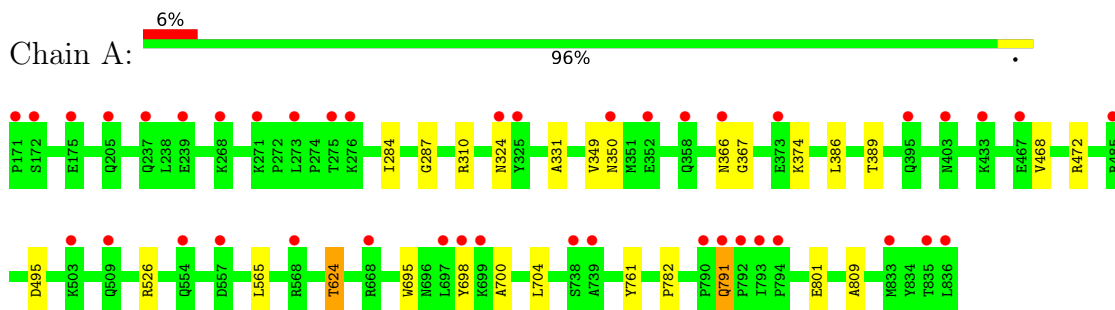
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	20	Total	O	0	0
			20	20		
7	B	2	Total	O	0	0
			2	2		
7	D	1	Total	O	0	0
			1	1		

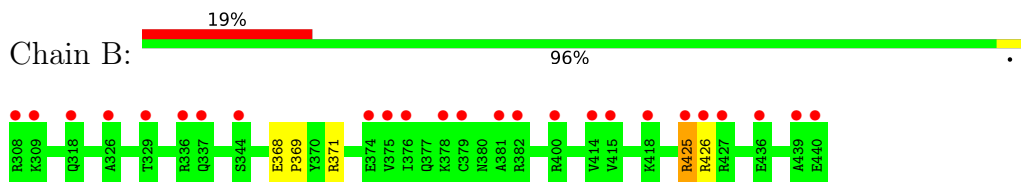
3 Residue-property plots [i](#)

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.

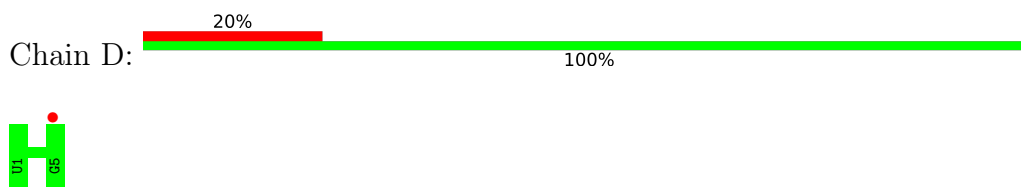
- Molecule 1: Lysine-specific histone demethylase 1A



- Molecule 2: REST corepressor 1



- Molecule 3: RNA (5'-R(*UP*UP*AP*GP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.07Å 178.41Å 234.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.02 – 2.80 49.02 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.02-2.80) 93.9 (49.02-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1839)	Depositor
R, R_{free}	0.201 , 0.227 0.203 , 0.229	Depositor DCC
R_{free} test set	2000 reflections (3.20%)	wwPDB-VP
Wilson B-factor (Å ²)	49.5	Xtriage
Anisotropy	0.692	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 48.3	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	12911	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, GOL, 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	0.33	0/5331	0.65	2/7232 (0.0%)
2	B	0.31	0/1091	0.73	1/1471 (0.1%)
3	D	0.10	0/95	0.33	0/147
All	All	0.32	0/6517	0.66	3/8850 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	425	ARG	N-CA-C	6.13	118.75	111.33
1	A	791	GLN	CA-C-N	5.57	124.92	118.85
1	A	791	GLN	C-N-CA	5.57	124.92	118.85

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	5252	5252	19	0
2	B	1076	1091	1091	4	0
3	D	86	43	43	0	0
4	A	53	31	30	1	0
5	A	25	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	6	8	8	2	0
7	A	20	0	0	2	0
7	B	2	0	0	0	0
7	D	1	0	0	0	0
All	All	6486	6425	6424	22	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 2.

All (22) 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:526:ARG:NH2	6:A:907:GOL:O2	2.22	0.72
1:A:782:PRO:O	1:A:791:GLN:HG2	1.94	0.66
1:A:287:GLY:N	7:A:1016:HOH:O	2.35	0.59
2:B:368:GLU:OE1	2:B:371:ARG:NH1	2.36	0.59
1:A:526:ARG:NH2	6:A:907:GOL:HO2	2.01	0.57
1:A:310:ARG:NH1	5:A:903:SO4:O2	2.42	0.52
2:B:425:ARG:HG3	2:B:426:ARG:N	2.25	0.52
1:A:698:TYR:O	1:A:700:ALA:N	2.42	0.50
1:A:495:ASP:OD1	2:B:371:ARG:NH2	2.46	0.49
1:A:624:THR:HG21	7:A:1016:HOH:O	2.15	0.46
1:A:324:ASN:OD1	1:A:324:ASN:N	2.45	0.46
1:A:801:GLU:HG2	1:A:809:ALA:H	1.80	0.46
1:A:331:ALA:HA	4:A:901:FAD:N5	2.31	0.45
1:A:349:VAL:HG12	1:A:350:ASN:N	2.32	0.44
1:A:791:GLN:O	1:A:791:GLN:HG3	2.18	0.44
1:A:284:ILE:O	1:A:624:THR:HB	2.19	0.43
1:A:468:VAL:O	1:A:472:ARG:NH1	2.49	0.43
1:A:695:TRP:HB2	1:A:704:LEU:HB3	1.99	0.43
1:A:761:TYR:CD1	1:A:809:ALA:HB1	2.53	0.42
1:A:366:ASN:OD1	1:A:367:GLY:N	2.53	0.41
2:B:368:GLU:N	2:B:369:PRO:CD	2.83	0.41
1:A:386:LEU:O	1:A:389:THR:OG1	2.35	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	664/666 (100%)	641 (96%)	23 (4%)	0	100	100
2	B	131/133 (98%)	126 (96%)	5 (4%)	0	100	100
All	All	795/799 (100%)	767 (96%)	28 (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	566/566 (100%)	563 (100%)	3 (0%)	81	93
2	B	117/117 (100%)	117 (100%)	0	100	100
All	All	683/683 (100%)	680 (100%)	3 (0%)	84	94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	374	LYS
1	A	565	LEU
1	A	624	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	185	HIS
1	A	438	GLN
1	A	535	ASN
2	B	318	GLN
2	B	423	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	D	3/5 (60%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

7 ligands 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	$\# Z > 2$
5	SO4	A	906	-	4,4,4	0.24	0	6,6,6	0.09	0
5	SO4	A	903	-	4,4,4	0.22	0	6,6,6	0.09	0
5	SO4	A	904	-	4,4,4	0.24	0	6,6,6	0.07	0
5	SO4	A	905	-	4,4,4	0.23	0	6,6,6	0.11	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GOL	A	907	-	5,5,5	0.38	0	5,5,5	0.47	0
5	SO4	A	902	-	4,4,4	0.25	0	6,6,6	0.07	0
4	FAD	A	901	-	58,58,58	3.31	20 (34%)	85,89,89	1.85	22 (25%)

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
6	GOL	A	907	-	-	2/4/4/4	-
4	FAD	A	901	-	-	1/34/50/50	0/6/6/6

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	FAD	O4B-C1B	8.84	1.62	1.42
4	A	901	FAD	C4X-N5	8.42	1.48	1.30
4	A	901	FAD	C2B-C1B	-7.43	1.30	1.53
4	A	901	FAD	C9A-N10	6.82	1.53	1.41
4	A	901	FAD	PA-O3P	6.69	1.66	1.59
4	A	901	FAD	O4B-C4B	-6.39	1.30	1.45
4	A	901	FAD	C10-N1	6.38	1.46	1.33
4	A	901	FAD	C5X-N5	5.88	1.50	1.39
4	A	901	FAD	C2-N1	5.35	1.48	1.36
4	A	901	FAD	C2-N3	5.17	1.50	1.39
4	A	901	FAD	C6A-N6A	4.61	1.46	1.34
4	A	901	FAD	C10-N10	4.58	1.47	1.37
4	A	901	FAD	C5A-C4A	-3.90	1.32	1.39
4	A	901	FAD	C4-N3	3.63	1.45	1.38
4	A	901	FAD	O2B-C2B	3.60	1.51	1.43
4	A	901	FAD	P-O3P	3.22	1.63	1.59
4	A	901	FAD	O3B-C3B	-2.86	1.35	1.43
4	A	901	FAD	O2'-C2'	-2.71	1.37	1.43
4	A	901	FAD	O2-C2	-2.34	1.19	1.24
4	A	901	FAD	C8A-N9A	-2.27	1.33	1.37

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	FAD	N3A-C2A-N1A	-5.67	120.00	128.58
4	A	901	FAD	N6A-C6A-N1A	-4.94	107.39	118.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	FAD	C5A-C4A-N3A	-4.48	120.55	126.72
4	A	901	FAD	N9A-C8A-N7A	-4.21	107.96	113.94
4	A	901	FAD	C7M-C7-C6	-3.99	112.55	119.57
4	A	901	FAD	C5A-C6A-N6A	3.86	132.85	123.29
4	A	901	FAD	C5'-C4'-C3'	-3.55	105.53	112.22
4	A	901	FAD	C4-N3-C2	-3.43	119.56	125.64
4	A	901	FAD	C7M-C7-C8	3.26	127.42	120.76
4	A	901	FAD	C2A-N3A-C4A	3.21	119.67	111.83
4	A	901	FAD	C8M-C8-C9	-3.13	114.05	119.57
4	A	901	FAD	O4-C4-C4X	-2.79	119.17	126.53
4	A	901	FAD	C5A-N7A-C8A	2.70	107.70	103.45
4	A	901	FAD	N3A-C4A-N9A	2.67	131.70	127.17
4	A	901	FAD	C4X-C4-N3	2.61	119.89	113.25
4	A	901	FAD	C4A-N9A-C8A	2.46	108.32	105.74
4	A	901	FAD	C8M-C8-C7	2.39	125.65	120.76
4	A	901	FAD	C4X-C10-N10	2.32	119.80	116.48
4	A	901	FAD	C9A-C5X-N5	-2.28	120.03	122.45
4	A	901	FAD	C5X-C9A-N10	2.27	120.02	117.97
4	A	901	FAD	C10-C4X-N5	-2.13	120.46	124.81
4	A	901	FAD	C4A-C5A-N7A	-2.02	108.27	110.58

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	907	GOL	O1-C1-C2-C3
6	A	907	GOL	O1-C1-C2-O2
4	A	901	FAD	C2B-C1B-N9A-C8A

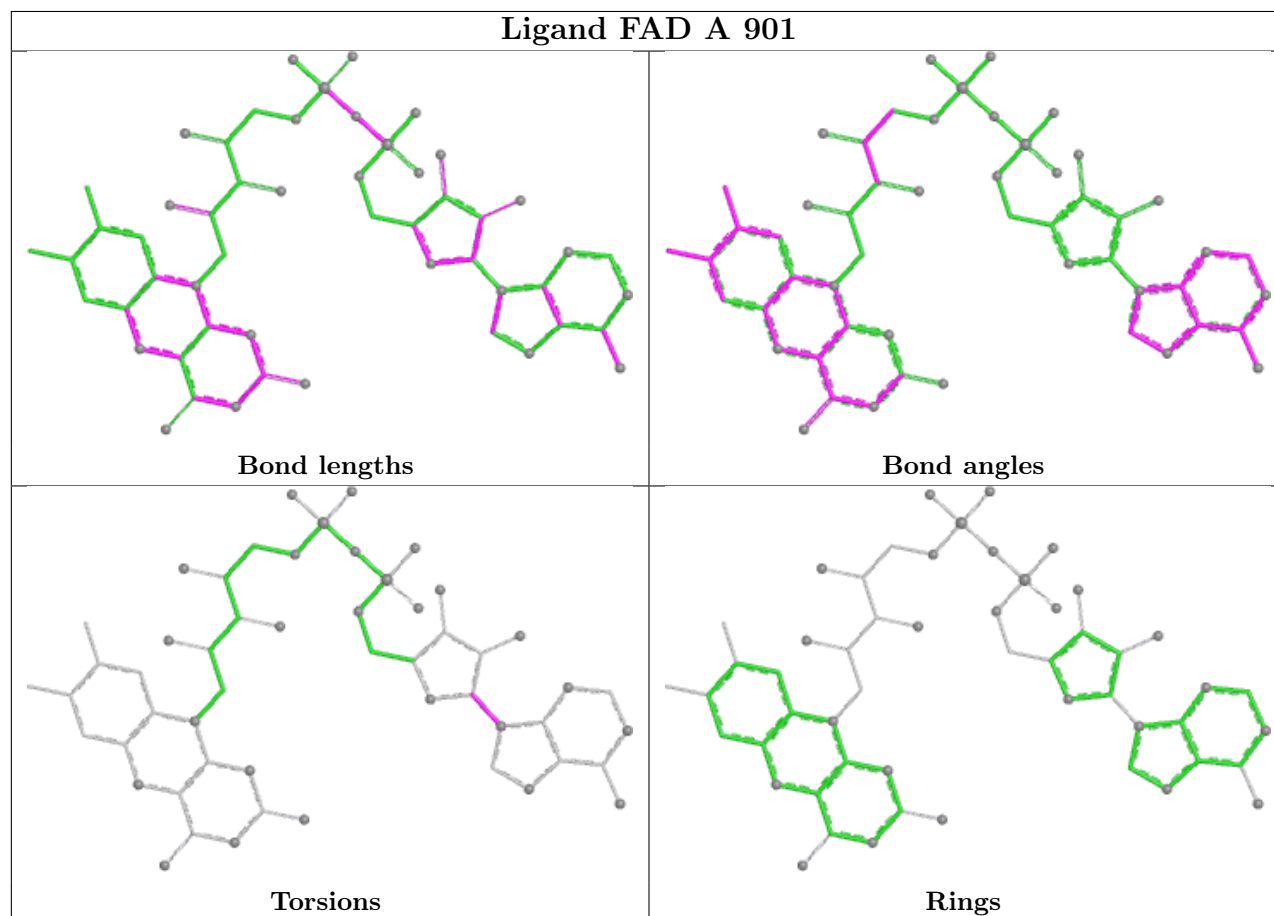
There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	903	SO4	1	0
6	A	907	GOL	2	0
4	A	901	FAD	1	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 than 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.



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	666/666 (100%)	0.17	42 (6%) 26 19	24, 60, 102, 140	0
2	B	133/133 (100%)	1.09	25 (18%) 3 2	52, 94, 124, 135	0
3	D	5/5 (100%)	0.98	1 (20%) 3 2	65, 67, 75, 108	0
All	All	804/804 (100%)	0.33	68 (8%) 16 12	24, 67, 110, 140	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	836	LEU	8.3
2	B	376	ILE	6.5
1	A	699	LYS	6.0
1	A	697	LEU	5.5
1	A	171	PRO	5.5
1	A	791	GLN	5.4
1	A	325	TYR	5.2
2	B	418	LYS	4.5
1	A	698	TYR	4.4
1	A	172	SER	4.4
1	A	793	ILE	4.4
2	B	308	ARG	4.1
2	B	375	VAL	4.1
1	A	237	GLN	4.1
1	A	792	PRO	4.1
3	D	5	G	4.1
1	A	835	THR	4.0
1	A	350	ASN	3.9
1	A	433	LYS	3.5
1	A	568	ARG	3.5
2	B	425	ARG	3.5
1	A	833	MET	3.3
2	B	318	GLN	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	374	GLU	3.2
1	A	395	GLN	3.0
2	B	440	GLU	3.0
1	A	275	THR	2.9
1	A	205	GLN	2.8
1	A	467	GLU	2.8
1	A	268	LYS	2.8
2	B	378	LYS	2.7
2	B	400	ARG	2.7
1	A	554	GLN	2.7
1	A	271	LYS	2.7
1	A	557	ASP	2.6
2	B	439	ALA	2.6
2	B	415	VAL	2.6
2	B	381	ALA	2.6
1	A	324	ASN	2.5
2	B	426	ARG	2.5
1	A	373	GLU	2.4
1	A	273	LEU	2.4
1	A	175	GLU	2.4
1	A	358	GLN	2.3
1	A	790	PRO	2.3
2	B	336	ARG	2.3
1	A	276	LYS	2.3
2	B	344	SER	2.3
1	A	739	ALA	2.2
1	A	794	PRO	2.2
1	A	239	GLU	2.2
1	A	668	ARG	2.2
2	B	382	ARG	2.2
2	B	427	ARG	2.2
1	A	352	GLU	2.2
2	B	309	LYS	2.1
1	A	366	ASN	2.1
1	A	403	ASN	2.1
1	A	485	ARG	2.1
1	A	738	SER	2.1
1	A	509	GLN	2.1
2	B	337	GLN	2.1
2	B	326	ALA	2.1
2	B	379	CYS	2.1
2	B	414	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	503	LYS	2.0
2	B	329	THR	2.0
2	B	436	GLU	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](#)

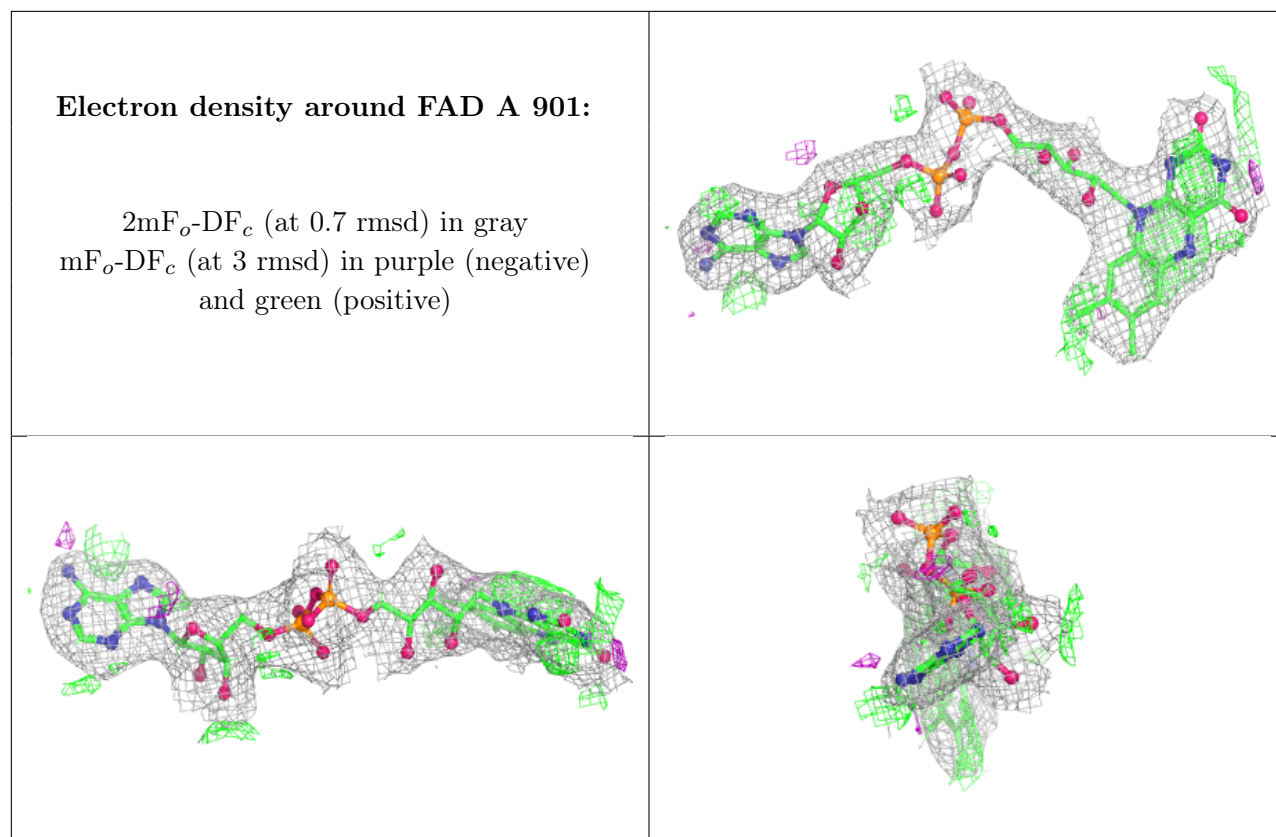
There are no oligosaccharides in this entry.

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, 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
5	SO4	A	906	5/5	0.78	0.35	79,83,88,136	0
5	SO4	A	904	5/5	0.79	0.19	95,100,109,123	0
5	SO4	A	903	5/5	0.82	0.25	73,85,102,106	0
5	SO4	A	905	5/5	0.88	0.21	77,79,92,133	0
5	SO4	A	902	5/5	0.90	0.23	61,74,113,113	0
6	GOL	A	907	6/6	0.91	0.16	59,71,79,85	0
4	FAD	A	901	53/53	0.97	0.08	21,36,55,76	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.