



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:50 PM UTC

PDB ID : 7CDF / pdb_00007cdf
Title : Crystal structure of LSD1-CoREST in complex with PRSFLVRRK peptide
Authors : Kikuchi, M.; Kitagawa, H.; Sato, S.; Umezawa, N.; Higuchi, T.; Umehara, T.
Deposited on : 2020-06-19
Resolution : 2.68 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

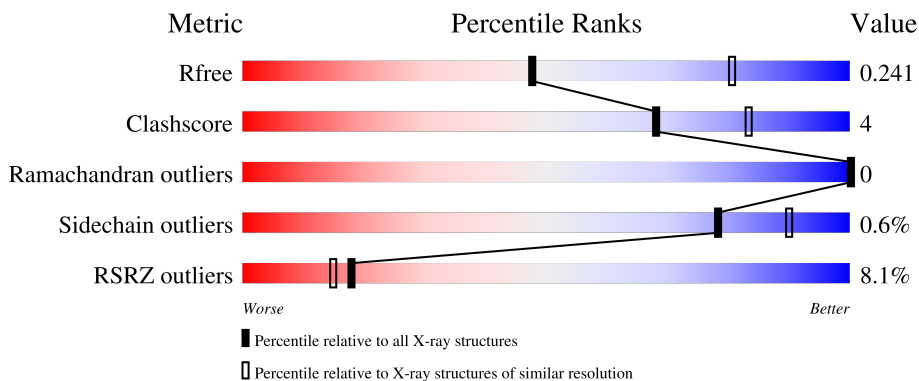
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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	669	 5% 91% 7% ..
2	B	140	 22% 84% 9% 6%
3	C	9	 22% 44% 56%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6459 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
			Total	C	N	O	S			
1	A	663	5185	3300	903	962	20	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	165	GLY	-	expression tag	UNP O60341
A	166	PRO	-	expression tag	UNP O60341
A	167	LEU	-	expression tag	UNP O60341
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
			Total	C	N	O	S			
2	B	131	1019	639	178	199	3	0	0	0

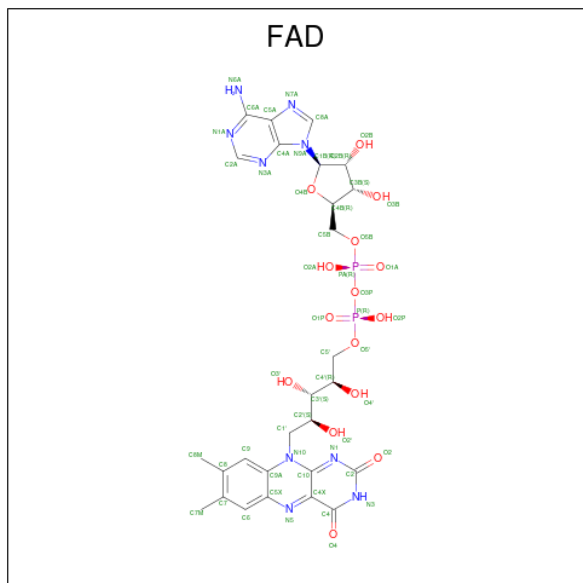
There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	301	GLY	-	expression tag	UNP Q9UKL0
B	302	SER	-	expression tag	UNP Q9UKL0
B	303	SER	-	expression tag	UNP Q9UKL0
B	304	GLY	-	expression tag	UNP Q9UKL0
B	305	SER	-	expression tag	UNP Q9UKL0
B	306	ALA	-	expression tag	UNP Q9UKL0
B	307	SER	-	expression tag	UNP Q9UKL0

- Molecule 3 is a protein called PRO-ARG-SER-PHE-LEU-VAL-ARG-ARG-LYS.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	9	81	52	19	10	0	0	0

- 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
			Total	C	N	O	P		
4	A	1	53	27	9	15	2	0	0

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

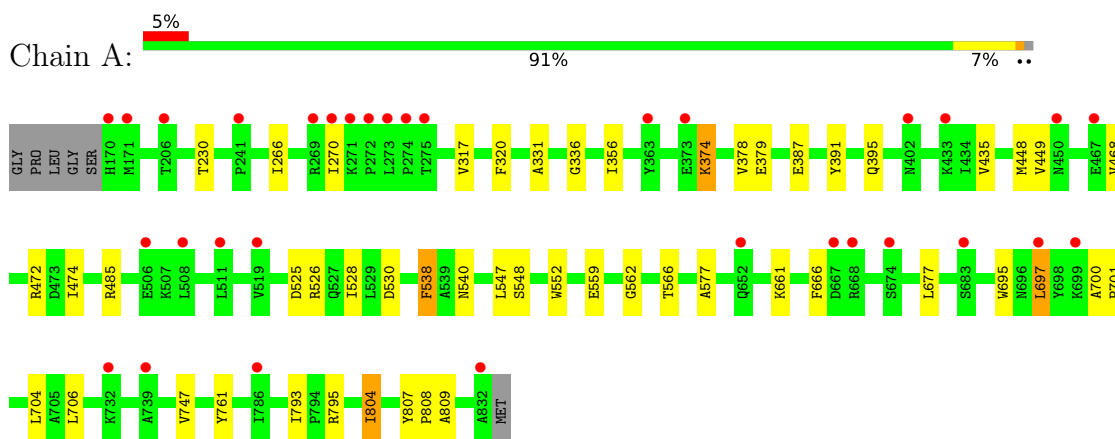
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	94	Total O 94 94	0	0
6	B	5	Total O 5 5	0	0
6	C	4	Total O 4 4	0	0

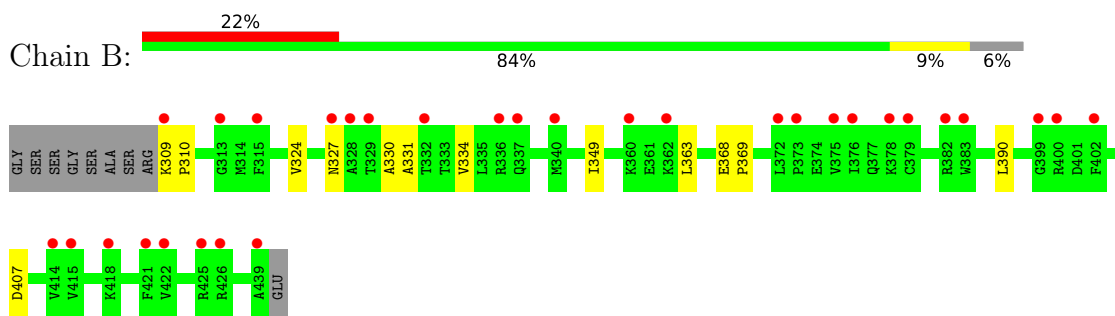
3 Residue-property plots [i](#)

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.

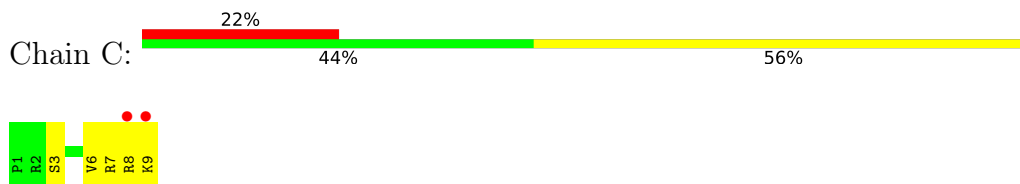
- Molecule 1: Lysine-specific histone demethylase 1A



- Molecule 2: REST corepressor 1



- Molecule 3: PRO-ARG-SER-PHE-LEU-VAL-ARG-ARG-LYS



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.72Å 180.23Å 232.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.00 – 2.68 45.00 – 2.68	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.00-2.68) 100.0 (45.00-2.68)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.69Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.214 , 0.241 0.216 , 0.241	Depositor DCC
R_{free} test set	1505 reflections (2.10%)	wwPDB-VP
Wilson B-factor (Å ²)	70.7	Xtrriage
Anisotropy	0.586	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6459	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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: FAD, GOL

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	0.98	0/5298	1.47	4/7189 (0.1%)
2	B	1.02	0/1034	1.62	0/1404
3	C	0.92	0/82	1.38	0/106
All	All	0.99	0/6414	1.49	4/8699 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	804	ILE	CA-C-N	5.71	128.20	120.38
1	A	804	ILE	C-N-CA	5.71	128.20	120.38
1	A	336	GLY	CA-C-O	-5.28	117.64	122.24
1	A	562	GLY	CA-C-O	-5.10	118.71	122.23

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	5185	0	5202	39	0
2	B	1019	0	979	11	0
3	C	81	0	95	5	0
4	A	53	0	31	2	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	18	0	24	0	0
6	A	94	0	0	0	0
6	B	5	0	0	0	0
6	C	4	0	0	0	0
All	All	6459	0	6331	45	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 4.

All (45) 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:695:TRP:HB3	1:A:697:LEU:HD11	1.74	0.69
1:A:379:GLU:OE1	3:C:7:ARG:NH1	2.27	0.68
1:A:448:MET:HE3	2:B:363:LEU:HD13	1.80	0.64
1:A:468:VAL:O	1:A:472:ARG:NH1	2.35	0.60
2:B:327:ASN:HB3	2:B:330:ALA:HB2	1.86	0.57
1:A:387:GLU:OE2	3:C:9:LYS:NZ	2.39	0.55
1:A:695:TRP:CE3	1:A:697:LEU:HD21	2.42	0.54
1:A:331:ALA:HA	4:A:901:FAD:N5	2.23	0.54
1:A:331:ALA:HA	4:A:901:FAD:C4X	2.38	0.53
2:B:309:LYS:N	2:B:310:PRO:CD	2.72	0.52
2:B:324:VAL:HG13	2:B:331:ALA:HB2	1.91	0.52
2:B:309:LYS:N	2:B:310:PRO:HD2	2.23	0.52
1:A:230:THR:HG23	1:A:270:ILE:HD12	1.93	0.50
1:A:449:VAL:HA	2:B:363:LEU:HD21	1.93	0.50
1:A:526:ARG:NH1	1:A:530:ASP:OD1	2.43	0.49
1:A:356:ILE:HG12	1:A:697:LEU:HD23	1.94	0.48
1:A:761:TYR:CD1	1:A:809:ALA:HB1	2.48	0.48
1:A:695:TRP:HE3	1:A:697:LEU:HD21	1.79	0.48
1:A:793:ILE:HD11	1:A:795:ARG:CZ	2.46	0.46
1:A:317:VAL:HG22	1:A:331:ALA:HB3	1.97	0.46
2:B:368:GLU:N	2:B:369:PRO:CD	2.80	0.45
1:A:485:ARG:HG2	2:B:407:ASP:HB3	1.99	0.45
1:A:695:TRP:HB2	1:A:704:LEU:HB3	1.99	0.45
1:A:540:ASN:HB3	1:A:547:LEU:HD21	1.98	0.44
1:A:320:PHE:CD1	1:A:747:VAL:HG21	2.52	0.44
1:A:548:SER:O	1:A:552:TRP:HB3	2.18	0.43
1:A:677:LEU:HD13	3:C:6:VAL:CG1	2.48	0.43
1:A:435:VAL:HG12	2:B:349:ILE:HG13	2.01	0.42
1:A:804:ILE:O	1:A:804:ILE:HG23	2.19	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:538:PHE:CD1	1:A:706:LEU:HD13	2.54	0.42
1:A:378:VAL:HG11	1:A:528:ILE:HG22	2.00	0.42
1:A:266:ILE:CD1	1:A:577:ALA:HB1	2.49	0.42
1:A:661:LYS:HB3	1:A:704:LEU:HD21	2.01	0.42
2:B:330:ALA:O	2:B:334:VAL:HG23	2.20	0.41
1:A:474:ILE:HD11	2:B:390:LEU:HG	2.02	0.41
1:A:695:TRP:HE1	1:A:706:LEU:HD11	1.85	0.41
1:A:807:TYR:N	1:A:808:PRO:CD	2.83	0.41
1:A:374:LYS:HE3	1:A:525:ASP:OD1	2.21	0.41
1:A:566:THR:HG21	1:A:697:LEU:HB3	2.03	0.41
1:A:391:TYR:CD1	1:A:395:GLN:HG3	2.56	0.41
1:A:559:GLU:OE1	3:C:8:ARG:NH1	2.54	0.41
1:A:677:LEU:HD13	3:C:6:VAL:HG13	2.03	0.41
1:A:666:PHE:O	1:A:701:PRO:HG2	2.21	0.40
1:A:695:TRP:HB3	1:A:697:LEU:CD1	2.48	0.40
1:A:700:ALA:HB1	1:A:701:PRO:HD2	2.04	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	661/669 (99%)	647 (98%)	14 (2%)	0	100	100
2	B	129/140 (92%)	126 (98%)	3 (2%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
All	All	797/818 (97%)	780 (98%)	17 (2%)	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	561/567 (99%)	558 (100%)	3 (0%)	81	91
2	B	106/121 (88%)	106 (100%)	0	100	100
3	C	9/9 (100%)	8 (89%)	1 (11%)	6	13
All	All	676/697 (97%)	672 (99%)	4 (1%)	78	90

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	374	LYS
1	A	538	PHE
1	A	697	LEU
3	C	3	SER

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

Mol	Chain	Res	Type
1	A	324	ASN
1	A	742	GLN
2	B	423	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry

4 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
4	FAD	A	901	-	58,58,58	0.55	0	85,89,89	0.66	0
5	GOL	A	903	-	5,5,5	0.09	0	5,5,5	0.25	0
5	GOL	A	904	-	5,5,5	0.12	0	5,5,5	0.32	0
5	GOL	A	902	-	5,5,5	0.08	0	5,5,5	0.28	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
4	FAD	A	901	-	-	2/34/50/50	0/6/6/6
5	GOL	A	903	-	-	0/4/4/4	-
5	GOL	A	904	-	-	2/4/4/4	-
5	GOL	A	902	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

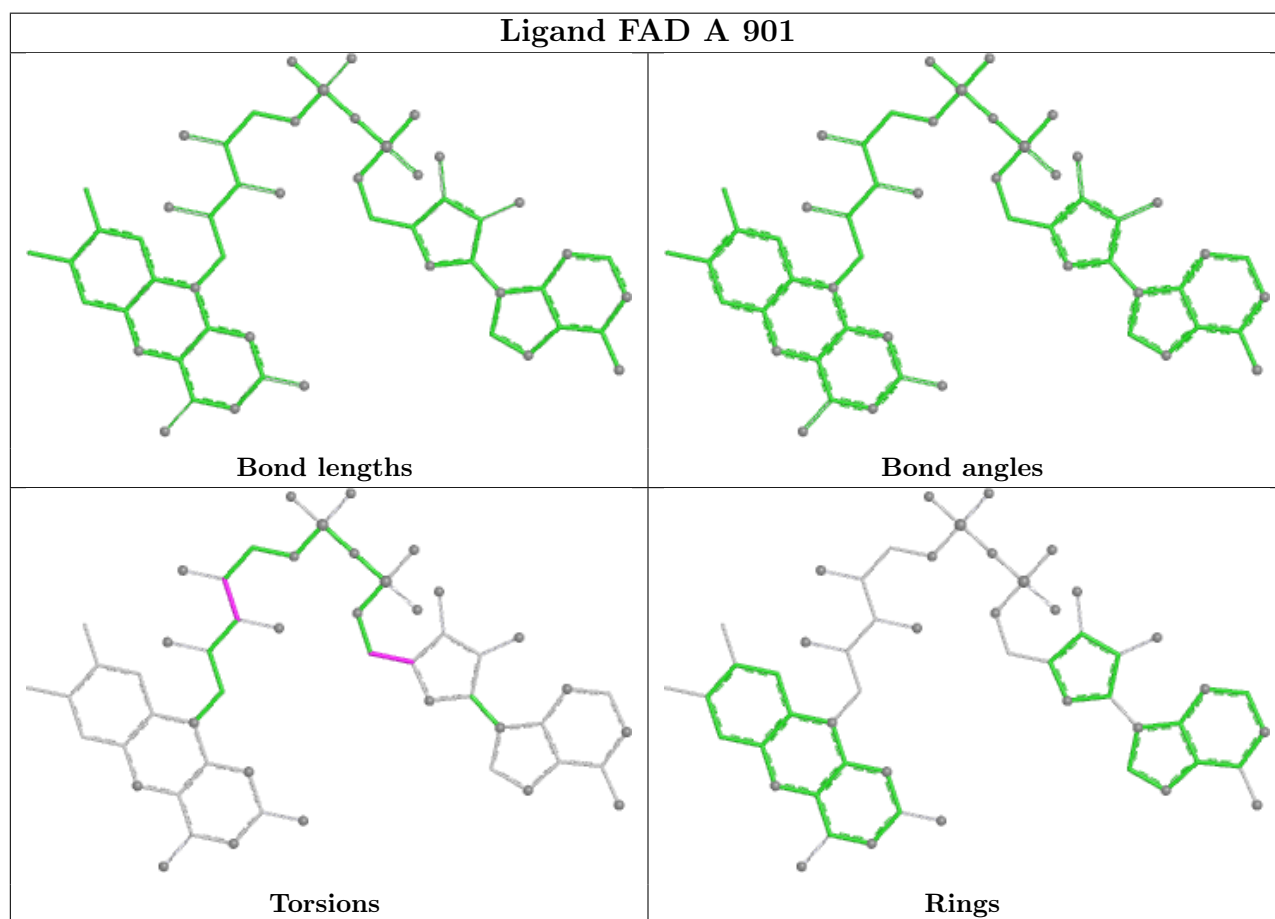
Mol	Chain	Res	Type	Atoms
5	A	904	GOL	C1-C2-C3-O3
5	A	904	GOL	O2-C2-C3-O3
4	A	901	FAD	O4B-C4B-C5B-O5B
4	A	901	FAD	C2'-C3'-C4'-O4'

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

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

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	663/669 (99%)	0.50	32 (4%) 35 31	53, 83, 119, 147	0
2	B	131/140 (93%)	1.38	31 (23%) 2 1	80, 116, 136, 151	0
3	C	9/9 (100%)	1.68	2 (22%) 2 1	70, 74, 120, 132	0
All	All	803/818 (98%)	0.66	65 (8%) 18 15	53, 90, 128, 151	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	170	HIS	5.9
1	A	832	ALA	5.7
1	A	171	MET	5.1
2	B	309	LYS	4.5
2	B	375	VAL	4.5
2	B	426	ARG	4.3
3	C	9	LYS	4.3
2	B	382	ARG	3.8
1	A	269	ARG	3.7
2	B	376	ILE	3.7
1	A	241	PRO	3.6
1	A	271	LYS	3.5
3	C	8	ARG	3.2
1	A	272	PRO	3.2
2	B	332	THR	3.1
1	A	652	GLN	3.1
2	B	383	TRP	3.1
2	B	313	GLY	3.0
2	B	414	VAL	3.0
1	A	275	THR	3.0
2	B	328	ALA	2.9
1	A	273	LEU	2.9
1	A	270	ILE	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	379	CYS	2.8
2	B	315	PHE	2.8
2	B	340	MET	2.8
1	A	467	GLU	2.7
1	A	511	LEU	2.7
2	B	336	ARG	2.7
1	A	274	PRO	2.6
1	A	683	SER	2.6
2	B	327	ASN	2.6
2	B	329	THR	2.6
1	A	363	TYR	2.5
2	B	402	PHE	2.5
1	A	667	ASP	2.5
2	B	422	VAL	2.5
2	B	362	LYS	2.5
2	B	378	LYS	2.5
2	B	425	ARG	2.4
2	B	415	VAL	2.4
2	B	400	ARG	2.4
2	B	337	GLN	2.4
2	B	439	ALA	2.4
1	A	786	ILE	2.4
1	A	508	LEU	2.3
2	B	372	LEU	2.3
1	A	519	VAL	2.3
1	A	697	LEU	2.3
1	A	206	THR	2.2
1	A	506	GLU	2.2
2	B	421	PHE	2.2
1	A	668	ARG	2.2
1	A	373	GLU	2.2
1	A	450	ASN	2.2
2	B	373	PRO	2.2
1	A	699	LYS	2.2
1	A	674	SER	2.1
1	A	739	ALA	2.1
2	B	399	GLY	2.1
1	A	433	LYS	2.1
2	B	360	LYS	2.1
2	B	418	LYS	2.0
1	A	402	ASN	2.0
1	A	732	LYS	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](#)

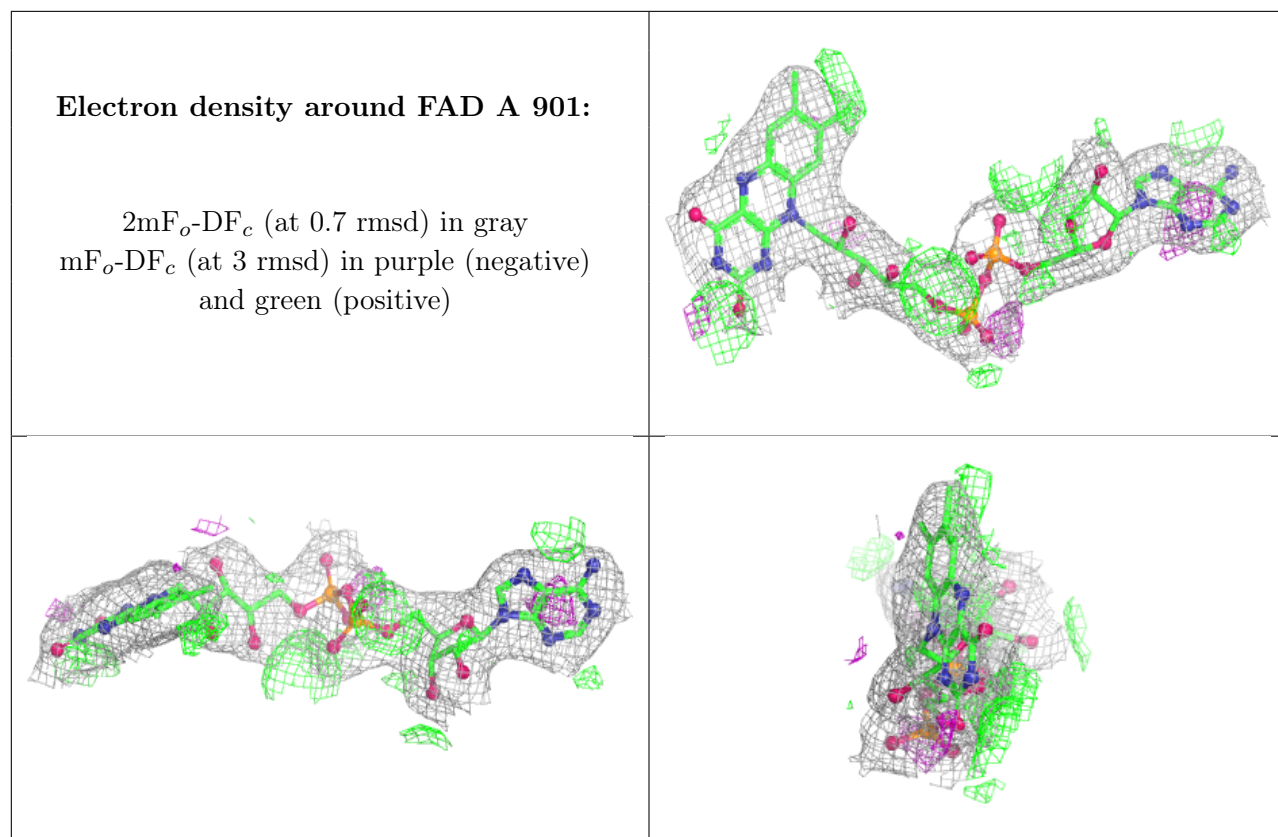
There are no oligosaccharides in this entry.

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	GOL	A	903	6/6	0.77	0.27	133,137,140,140	0
5	GOL	A	904	6/6	0.93	0.18	91,94,99,100	0
5	GOL	A	902	6/6	0.95	0.13	77,82,83,84	0
4	FAD	A	901	53/53	0.97	0.07	51,59,68,69	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 [i](#)

There are no such residues in this entry.