



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 04:42 PM UTC

PDB ID : 4GUS / pdb_00004gus
Title : Crystal structure of LSD2-NPAC with H3 in space group P3221
Authors : Chen, F.; Dong, Z.; Fang, J.; Yang, Y.; Li, Z.; Xu, Y.; Yang, H.; Wang, P.;
Fang, R.; Shi, Y.; Xu, Y.
Deposited on : 2012-08-29
Resolution : 2.23 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

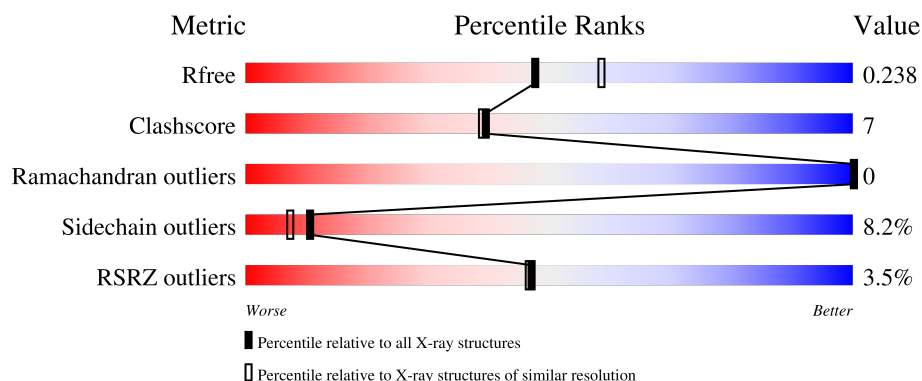
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

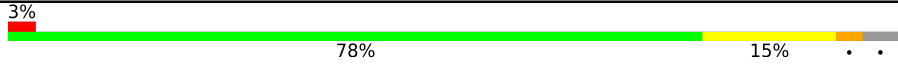
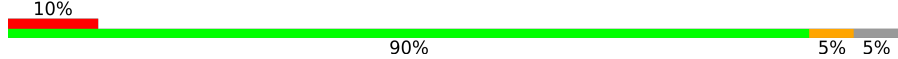
The reported resolution of this entry is 2.23 Å.

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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	776	
2	B	124	
3	C	21	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	IOD	A	904	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6515 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	746	Total	C	N	O	S	0	0	0
			5904	3769	1005	1089	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

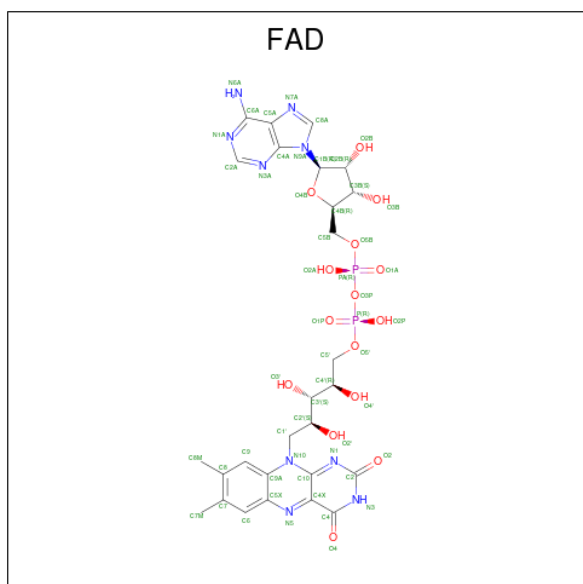
- Molecule 3 is a protein called Histone H3.3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	20	Total	C	N	O	S	0	0	0
			151	90	34	26	1			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	4	MET	LYS	engineered mutation	UNP P84243

- 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 53	C 27	N 9	O 15	P 2	0	0

- Molecule 5 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	6	Total I 6 6	0	0

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



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

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	264	Total	O	0	0
			264	264		
8	B	6	Total	O	0	0
			6	6		
8	C	11	Total	O	0	0
			11	11		

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.08Å 101.08Å 177.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.04 – 2.23 29.04 – 2.23	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.04-2.23) 99.5 (29.04-2.23)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 2.22Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.207 , 0.237 0.209 , 0.238	Depositor DCC
R_{free} test set	2619 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	42.5	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 26.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6515	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: IOD, ZN, 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	0.56	0/6051	0.86	5/8199 (0.1%)
2	B	0.47	0/110	0.77	0/149
3	C	0.64	0/151	0.81	0/198
All	All	0.56	0/6312	0.86	5/8546 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	376	LEU	CA-C-N	5.34	125.33	119.89
1	A	376	LEU	C-N-CA	5.34	125.33	119.89
1	A	760	ASP	CA-C-N	5.29	125.10	119.28
1	A	760	ASP	C-N-CA	5.29	125.10	119.28
1	A	75	LYS	CB-CA-C	-5.08	110.32	117.23

There are no chirality outliers.

There are no planarity outliers.

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	5904	0	5821	80	0
2	B	105	0	91	1	0
3	C	151	0	170	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	53	0	30	3	0
5	A	6	0	0	2	0
6	A	12	0	16	5	0
7	A	3	0	0	0	0
8	A	264	0	0	7	0
8	B	6	0	0	0	0
8	C	11	0	0	0	0
All	All	6515	0	6128	83	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 (83) 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:521:ILE:O	1:A:525:GLY:HA2	1.63	0.98
1:A:521:ILE:O	1:A:525:GLY:CA	2.19	0.91
1:A:521:ILE:O	1:A:525:GLY:N	2.25	0.70
1:A:738:ARG:NH2	6:A:909:GOL:O3	2.26	0.69
1:A:741:PHE:C	1:A:743:GLU:OE2	2.37	0.67
1:A:304:PRO:HB2	1:A:308:LEU:HD22	1.79	0.65
1:A:266:ASN:HD22	1:A:452:GLU:HG2	1.65	0.62
1:A:66:PHE:CE2	1:A:131:MET:HE3	2.34	0.62
1:A:743:GLU:OE1	1:A:744:GLN:NE2	2.33	0.61
1:A:493:VAL:HG11	1:A:516:ILE:HG21	1.84	0.59
1:A:151:ARG:HD2	1:A:169:CYS:SG	2.44	0.58
1:A:114:LYS:O	1:A:118:THR:OG1	2.19	0.57
1:A:498:ARG:HD2	8:A:1170:HOH:O	2.05	0.56
1:A:757:TRP:CD2	6:A:908:GOL:H12	2.41	0.56
1:A:454:LEU:HD21	1:A:585:LYS:HG2	1.87	0.55
1:A:722:ARG:HB3	8:A:1279:HOH:O	2.07	0.55
1:A:194:LEU:HD13	8:A:1183:HOH:O	2.07	0.54
1:A:743:GLU:O	1:A:744:GLN:HB3	2.06	0.54
1:A:127:PRO:O	1:A:131:MET:HG2	2.08	0.53
1:A:51:ARG:NH2	1:A:63:PRO:O	2.42	0.53
1:A:133:ASP:OD1	1:A:212:LYS:NZ	2.41	0.53
1:A:274:GLN:HB2	1:A:277:GLU:HG3	1.90	0.53
1:A:738:ARG:HG2	1:A:746:VAL:HG13	1.91	0.52
1:A:229:GLY:HA3	1:A:309:ALA:HB2	1.90	0.52
1:A:743:GLU:OE2	1:A:743:GLU:N	2.41	0.51
3:C:17:ARG:NE	3:C:17:ARG:H	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:526:ILE:HG12	1:A:526:ILE:O	2.11	0.50
1:A:677:GLY:O	1:A:704:LYS:NZ	2.43	0.50
1:A:422:TRP:CZ3	6:A:908:GOL:H31	2.48	0.49
1:A:741:PHE:HA	1:A:743:GLU:OE2	2.12	0.49
1:A:494:VAL:O	1:A:498:ARG:HG2	2.13	0.49
3:C:17:ARG:H	3:C:17:ARG:HE	1.59	0.49
1:A:618:TYR:OH	8:A:1276:HOH:O	2.19	0.48
1:A:693:LEU:HD22	1:A:694:PHE:CE1	2.47	0.48
1:A:741:PHE:CA	1:A:743:GLU:OE2	2.61	0.48
1:A:766:ALA:HA	4:A:901:FAD:HM82	1.96	0.48
1:A:218:PRO:O	8:A:1213:HOH:O	2.20	0.48
1:A:53:CYS:SG	1:A:84:HIS:HB2	2.54	0.48
1:A:327:THR:OG1	1:A:330:LYS:HG2	2.14	0.47
1:A:127:PRO:HD2	5:A:904:IOD:I	2.84	0.47
1:A:464:ARG:NH2	1:A:676:GLN:HB2	2.29	0.47
1:A:721:VAL:HG22	1:A:729:VAL:HG22	1.96	0.47
1:A:608:VAL:HG21	1:A:623:VAL:HB	1.97	0.47
1:A:300:TYR:OH	1:A:348:GLU:OE2	2.18	0.46
1:A:67:ALA:HB2	1:A:131:MET:HE1	1.96	0.46
1:A:469:GLN:OE1	1:A:473:ARG:HG3	2.14	0.46
1:A:383:LYS:NZ	1:A:822:PHE:O	2.44	0.46
1:A:526:ILE:O	1:A:526:ILE:HG23	2.16	0.46
8:A:1273:HOH:O	2:B:216:HIS:HB2	2.17	0.45
1:A:469:GLN:HB2	1:A:473:ARG:HG2	1.98	0.45
1:A:725:ASP:OD2	1:A:728:GLN:HG3	2.17	0.45
1:A:342:ARG:O	1:A:346:VAL:HG13	2.17	0.45
1:A:67:ALA:HB2	1:A:131:MET:CE	2.47	0.45
1:A:226:ASP:OD1	1:A:226:ASP:N	2.50	0.45
1:A:669:ARG:CG	1:A:707:SER:HB3	2.46	0.45
1:A:298:PRO:O	1:A:301:SER:OG	2.35	0.44
1:A:626:THR:HA	1:A:793:ALA:O	2.17	0.44
1:A:274:GLN:O	1:A:277:GLU:HB2	2.17	0.44
1:A:128:LYS:HG3	5:A:904:IOD:I	2.88	0.44
1:A:136:LEU:HD13	1:A:339:GLY:HA3	2.00	0.44
1:A:266:ASN:ND2	1:A:452:GLU:HG2	2.30	0.43
1:A:552:HIS:O	1:A:552:HIS:CG	2.70	0.43
1:A:734:MET:HG3	6:A:909:GOL:H2	2.01	0.43
1:A:781:ILE:O	1:A:784:GLU:HB2	2.19	0.43
1:A:180:GLU:H	1:A:180:GLU:HG3	1.39	0.43
1:A:493:VAL:HG13	1:A:516:ILE:HD13	2.00	0.43
1:A:396:ALA:HB2	1:A:809:TYR:CD1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:ARG:HB2	8:A:1136:HOH:O	2.19	0.42
1:A:369:VAL:HA	1:A:370:GLY:HA3	1.60	0.42
1:A:757:TRP:CE3	6:A:908:GOL:H12	2.54	0.42
1:A:327:THR:H	1:A:330:LYS:HG3	1.83	0.42
1:A:685:PRO:HB3	1:A:691:ARG:HA	2.02	0.42
1:A:355:PHE:CD2	1:A:356:MET:HE2	2.55	0.42
1:A:685:PRO:HA	1:A:686:PRO:HD3	1.85	0.41
1:A:289:MET:HE3	1:A:307:TYR:CD2	2.55	0.41
1:A:541:SER:HB2	1:A:691:ARG:HG3	2.03	0.41
1:A:692:GLY:HA3	1:A:714:ALA:O	2.21	0.41
1:A:330:LYS:HB3	1:A:330:LYS:HE3	1.88	0.41
1:A:385:VAL:HB	1:A:622:LYS:HB2	2.02	0.41
1:A:436:ALA:HA	4:A:901:FAD:N5	2.36	0.41
1:A:782:ILE:HG22	1:A:797:THR:HG22	2.02	0.41
1:A:803:GLN:O	4:A:901:FAD:O3'	2.39	0.40
1:A:470:GLU:HA	1:A:471:GLY:HA2	1.60	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	742/776 (96%)	713 (96%)	29 (4%)	0	100	100
2	B	10/124 (8%)	10 (100%)	0	0	100	100
3	C	18/21 (86%)	16 (89%)	2 (11%)	0	100	100
All	All	770/921 (84%)	739 (96%)	31 (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	640/662 (97%)	586 (92%)	54 (8%)	10	7
2	B	12/106 (11%)	12 (100%)	0	100	100
3	C	15/15 (100%)	14 (93%)	1 (7%)	15	12
All	All	667/783 (85%)	612 (92%)	55 (8%)	10	7

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	85	LEU
1	A	109	LYS
1	A	124	GLU
1	A	151	ARG
1	A	177	ILE
1	A	180	GLU
1	A	181	THR
1	A	194	LEU
1	A	205	LEU
1	A	285	ARG
1	A	294	LEU
1	A	301	SER
1	A	308	LEU
1	A	330	LYS
1	A	346	VAL
1	A	353	LEU
1	A	361	LEU
1	A	366	VAL
1	A	383	LYS
1	A	385	VAL
1	A	407	LYS
1	A	410	VAL
1	A	414	LYS
1	A	430	VAL
1	A	449	LEU
1	A	457	SER

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Mol	Chain	Res	Type
1	A	464	ARG
1	A	467	LEU
1	A	473	ARG
1	A	493	VAL
1	A	495	SER
1	A	498	ARG
1	A	526	ILE
1	A	527	GLN
1	A	540	LEU
1	A	543	LEU
1	A	553	GLN
1	A	554	VAL
1	A	574	LEU
1	A	585	LYS
1	A	604	SER
1	A	608	VAL
1	A	625	VAL
1	A	680	PHE
1	A	693	LEU
1	A	704	LYS
1	A	718	VAL
1	A	721	VAL
1	A	727	LYS
1	A	730	LEU
1	A	740	LEU
1	A	742	LYS
1	A	743	GLU
1	A	746	VAL
3	C	17	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	141	GLN
1	A	152	GLN
1	A	158	GLN
1	A	184	HIS
1	A	503	GLN
1	A	731	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](#)

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 [i](#)

Of 12 ligands modelled in this entry, 9 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	GOL	A	908	-	5,5,5	0.53	0	5,5,5	0.33	0
4	FAD	A	901	-	58,58,58	2.33	13 (22%)	85,89,89	2.51	23 (27%)
6	GOL	A	909	-	5,5,5	0.41	0	5,5,5	0.87	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
6	GOL	A	908	-	-	3/4/4/4	-
4	FAD	A	901	-	-	2/34/50/50	0/6/6/6
6	GOL	A	909	-	-	2/4/4/4	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	FAD	O2'-C2'	-8.66	1.25	1.43
4	A	901	FAD	O2-C2	8.10	1.40	1.24
4	A	901	FAD	O4-C4	7.35	1.37	1.23
4	A	901	FAD	P-O3P	-3.56	1.55	1.59
4	A	901	FAD	C1'-C2'	-3.37	1.47	1.52
4	A	901	FAD	C9A-N10	-2.64	1.36	1.41
4	A	901	FAD	C5B-C4B	-2.57	1.43	1.51
4	A	901	FAD	PA-O3P	-2.35	1.57	1.59
4	A	901	FAD	C4-N3	-2.30	1.34	1.38
4	A	901	FAD	C3B-C4B	-2.18	1.47	1.53
4	A	901	FAD	C8A-N9A	-2.18	1.33	1.37
4	A	901	FAD	C6A-N6A	2.04	1.39	1.34
4	A	901	FAD	C4X-N5	2.01	1.35	1.30

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	FAD	C4A-N9A-C1B	-9.22	105.08	126.63
4	A	901	FAD	C4A-N9A-C8A	8.32	114.47	105.74
4	A	901	FAD	N9A-C8A-N7A	-6.93	104.10	113.94
4	A	901	FAD	N3A-C4A-N9A	6.37	138.00	127.17
4	A	901	FAD	C1B-N9A-C8A	6.01	140.44	127.09
4	A	901	FAD	C5A-C4A-N3A	-5.58	119.03	126.72
4	A	901	FAD	O2'-C2'-C3'	5.18	121.36	109.25
4	A	901	FAD	C5A-N7A-C8A	4.82	111.02	103.45
4	A	901	FAD	N3A-C2A-N1A	-4.01	122.52	128.58
4	A	901	FAD	C2A-N3A-C4A	3.28	119.83	111.83
4	A	901	FAD	O5'-C5'-C4'	3.02	117.42	109.36
4	A	901	FAD	C4X-C10-N10	2.95	120.70	116.48
4	A	901	FAD	O4B-C1B-C2B	-2.90	100.42	106.62
4	A	901	FAD	C4-N3-C2	-2.89	120.51	125.64
4	A	901	FAD	C1'-C2'-C3'	2.87	117.44	109.66
4	A	901	FAD	O5B-C5B-C4B	2.78	118.45	108.99
4	A	901	FAD	C4A-C5A-N7A	-2.76	107.43	110.58
4	A	901	FAD	C5X-C9A-N10	2.72	120.42	117.97
4	A	901	FAD	C5A-C4A-N9A	-2.60	102.97	105.81
4	A	901	FAD	C3B-C2B-C1B	2.40	106.00	101.46
4	A	901	FAD	O4-C4-C4X	-2.32	120.42	126.53
4	A	901	FAD	C4X-C4-N3	2.24	118.96	113.25
4	A	901	FAD	C10-C4X-N5	-2.14	120.44	124.81

There are no chirality outliers.

All (7) torsion outliers are listed below:

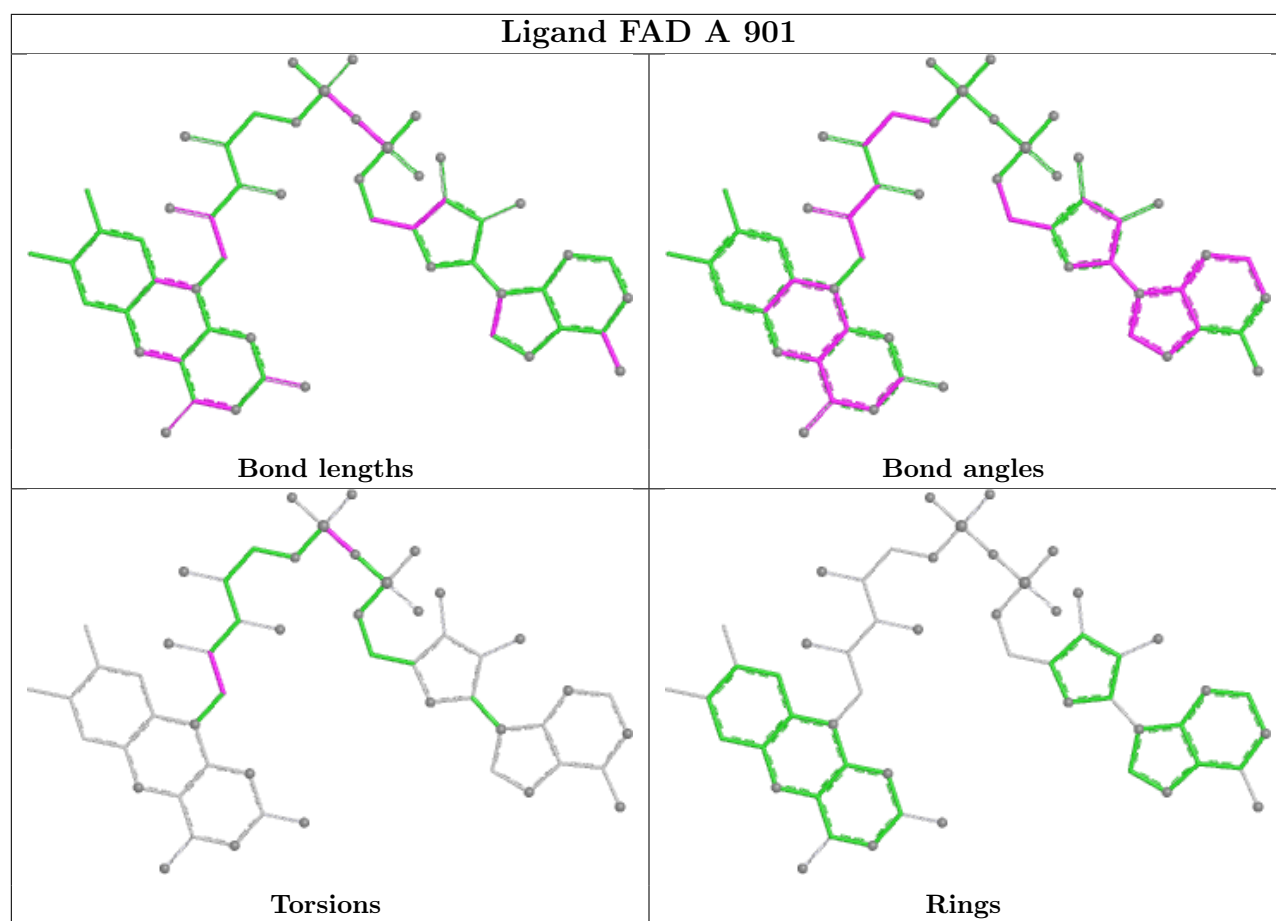
Mol	Chain	Res	Type	Atoms
4	A	901	FAD	N10-C1'-C2'-O2'
6	A	908	GOL	O1-C1-C2-O2
6	A	908	GOL	O1-C1-C2-C3
6	A	909	GOL	O1-C1-C2-C3
6	A	909	GOL	O1-C1-C2-O2
4	A	901	FAD	PA-O3P-P-O5'
6	A	908	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	908	GOL	3	0
4	A	901	FAD	3	0
6	A	909	GOL	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 [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	746/776 (96%)	0.27	24 (3%) 50 50	26, 43, 63, 83	0
2	B	12/124 (9%)	0.47	1 (8%) 17 15	30, 40, 53, 59	0
3	C	20/21 (95%)	0.27	2 (10%) 12 11	30, 37, 66, 67	0
All	All	778/921 (84%)	0.27	27 (3%) 47 46	26, 43, 63, 83	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	526	ILE	4.0
3	C	20	LEU	3.8
1	A	235	THR	3.6
1	A	170	GLY	3.4
1	A	369	VAL	3.4
1	A	371	ALA	3.2
1	A	743	GLU	3.1
1	A	525	GLY	2.8
1	A	177	ILE	2.7
3	C	17	ARG	2.6
1	A	500	ASP	2.6
1	A	174	ASN	2.4
1	A	370	GLY	2.4
1	A	822	PHE	2.4
1	A	78	TYR	2.3
1	A	124	GLU	2.3
2	B	225	THR	2.3
1	A	67	ALA	2.2
1	A	117	TRP	2.2
1	A	49	GLY	2.2
1	A	372	ASP	2.2
1	A	120	ASN	2.1
1	A	473	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	181	THR	2.1
1	A	302	ARG	2.1
1	A	172	LYS	2.0
1	A	524	SER	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](#)

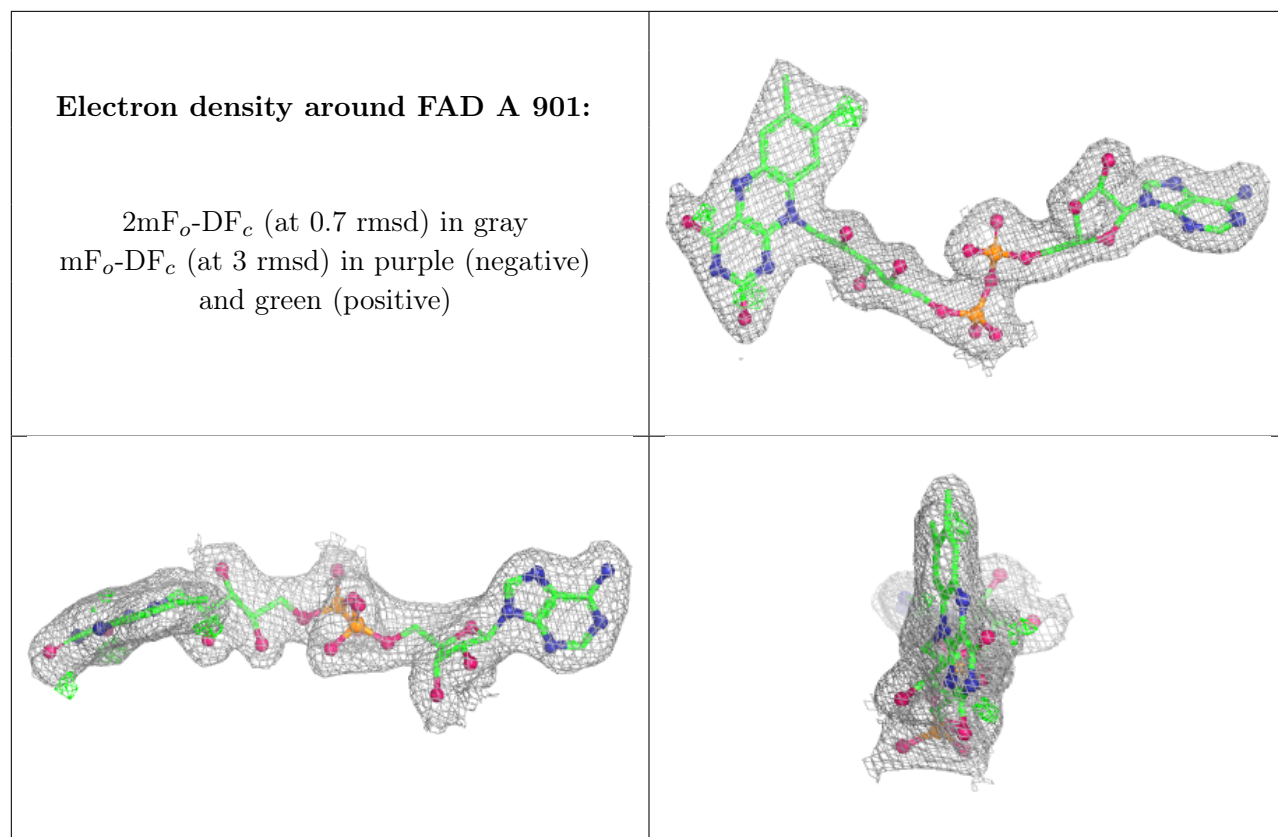
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
6	GOL	A	909	6/6	0.85	0.13	48,49,51,53	0
5	IOD	A	907	1/1	0.86	0.17	131,131,131,131	0
5	IOD	A	905	1/1	0.89	0.19	142,142,142,142	0
6	GOL	A	908	6/6	0.90	0.11	33,38,40,44	0
4	FAD	A	901	53/53	0.97	0.06	26,30,34,38	0
5	IOD	A	902	1/1	0.97	0.08	60,60,60,60	0
5	IOD	A	903	1/1	0.97	0.06	48,48,48,48	0
5	IOD	A	904	1/1	0.97	0.07	72,72,72,72	0
5	IOD	A	906	1/1	0.98	0.13	78,78,78,78	0
7	ZN	A	911	1/1	0.99	0.03	47,47,47,47	0
7	ZN	A	910	1/1	1.00	0.02	42,42,42,42	0
7	ZN	A	912	1/1	1.00	0.02	57,57,57,57	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.