



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 19, 2020 – 07:40 pm BST

PDB ID : 5L3G
Title : LSD1-CoREST1 in complex with polymyxin E (colistin)
Authors : Speranzini, V.; Rotili, D.; Ciossani, G.; Pilotto, S.; Forgione, M.; Lucidi, A.;
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Deposited on : 2016-04-10
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

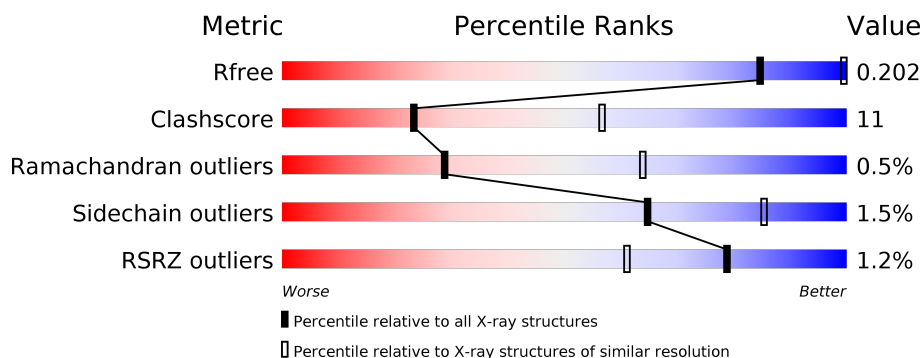
1 Overall quality at a glance

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 $\leq 5\%$. 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	730	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> % <div style="width: 100%; height: 10px; background-color: green;"></div> 9% </div> <div style="display: flex; justify-content: space-between; align-items: center;"> 68% 23% </div> </div>
2	B	178	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> 3% <div style="width: 100%; height: 10px; background-color: green;"></div> 25% </div> <div style="display: flex; justify-content: space-between; align-items: center;"> 56% 19% </div> </div>
3	C	11	<div> <div style="width: 100%; height: 10px; background-color: green;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> 73% 18% 9% </div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6429 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	666	5217	3324	906	967	20	0	0	0

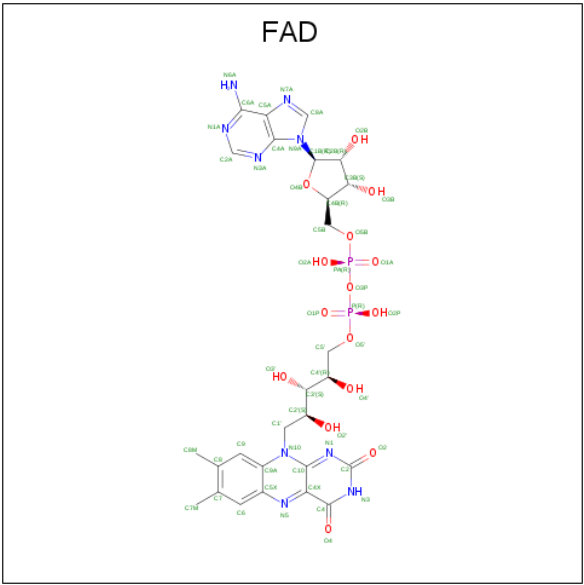
- Molecule 2 is a protein called REST corepressor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	133	1076	676	194	203	3	0	0	0

- Molecule 3 is a protein called polymyxin E.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	11	81	52	16	13	28	0	0

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Na	0	0
			2	2		

Chain C:  73% 18% 9%



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	119.86Å 179.49Å 236.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.78 – 3.10 48.78 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.78-3.10) 99.5 (48.78-3.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.173 , 0.197 0.177 , 0.202	Depositor DCC
R_{free} test set	898 reflections (1.93%)	wwPDB-VP
Wilson B-factor (Å ²)	97.8	Xtriage
Anisotropy	0.617	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 108.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6429	wwPDB-VP
Average B, all atoms (Å ²)	126.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% 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: 4FO, DTH, DAB, NA, 6F5, 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	0.66	1/5331 (0.0%)	0.83	5/7232 (0.1%)
2	B	0.48	0/1091	0.66	0/1471
3	C	2.03	0/15	1.47	0/19
All	All	0.64	1/6437 (0.0%)	0.81	5/8722 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	600	CYS	CB-SG	-5.11	1.73	1.81

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	659	LEU	CA-CB-CG	6.12	129.39	115.30
1	A	697	LEU	CB-CG-CD1	-5.76	101.20	111.00
1	A	693	LEU	CA-CB-CG	5.57	128.10	115.30
1	A	815	LEU	CB-CG-CD1	-5.55	101.57	111.00
1	A	659	LEU	CB-CG-CD1	-5.05	102.42	111.00

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	0	5251	128	0
2	B	1076	0	1091	26	0
3	C	81	0	55	1	0
4	A	53	0	30	6	0
5	A	2	0	0	0	0
All	All	6429	0	6427	145	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 11.

All (145) 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:566:THR:HG21	1:A:697:LEU:HD13	1.49	0.93
1:A:793:ILE:H	1:A:793:ILE:HD12	1.42	0.84
1:A:695:TRP:CE3	1:A:697:LEU:HD11	2.19	0.78
1:A:572:SER:O	1:A:575:PRO:HD2	1.86	0.74
1:A:627:LEU:HD23	1:A:656:PHE:HB2	1.70	0.73
1:A:693:LEU:HD12	1:A:694:PHE:N	2.04	0.72
1:A:319:THR:HB	1:A:572:SER:HB3	1.71	0.71
1:A:625:LEU:HD22	1:A:629:VAL:HG11	1.71	0.71
1:A:321:ARG:HG2	1:A:326:VAL:HG22	1.75	0.69
1:A:356:ILE:HD11	1:A:566:THR:HG23	1.75	0.68
1:A:793:ILE:H	1:A:793:ILE:CD1	2.06	0.67
1:A:801:GLU:HG3	1:A:809:ALA:HA	1.76	0.67
1:A:693:LEU:HD12	1:A:694:PHE:H	1.60	0.67
1:A:468:VAL:O	1:A:472:ARG:NH1	2.31	0.63
1:A:695:TRP:HE3	1:A:697:LEU:HD11	1.60	0.63
1:A:802:HIS:H	1:A:802:HIS:CD2	2.14	0.62
1:A:421:LYS:NZ	2:B:320:ASP:OD2	2.32	0.62
1:A:463:LYS:O	1:A:467:GLU:HG2	2.00	0.61
1:A:534:ALA:HB2	1:A:688:ARG:HG3	1.82	0.61
1:A:374:LYS:HE2	1:A:524:ARG:HH12	1.66	0.60
1:A:313:VAL:HG12	1:A:314:GLY:N	2.17	0.60
1:A:366:ASN:OD1	1:A:367:GLY:N	2.34	0.60
1:A:331:ALA:HA	4:A:901:FAD:N5	2.17	0.60
1:A:189:THR:OG1	1:A:192:GLU:HG3	2.02	0.59
1:A:439:GLU:HG3	2:B:352:ILE:HD13	1.84	0.59
1:A:801:GLU:HG2	1:A:809:ALA:H	1.66	0.59
1:A:238:LEU:HB3	1:A:243:ASN:HB3	1.85	0.58
1:A:230:THR:HG23	1:A:270:ILE:HD12	1.86	0.58
1:A:566:THR:HG21	1:A:697:LEU:CD1	2.30	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:458:LEU:CD2	1:A:487:LEU:HB2	2.35	0.56
1:A:456:LYS:HG2	2:B:370:TYR:CZ	2.41	0.56
1:A:296:GLN:O	1:A:299:SER:HB3	2.06	0.55
1:A:428:ILE:HG12	2:B:342:LEU:HD13	1.88	0.55
2:B:387:GLU:OE1	2:B:412:LYS:NZ	2.39	0.55
1:A:331:ALA:HA	4:A:901:FAD:C4X	2.36	0.55
1:A:374:LYS:HE2	1:A:524:ARG:NH1	2.22	0.55
3:C:6:DAB:O	3:C:7:LEU:HB2	2.07	0.55
1:A:716:GLU:HG2	1:A:750:ARG:HG2	1.90	0.54
2:B:421:PHE:HE2	2:B:434:LEU:HD11	1.72	0.53
2:B:332:THR:O	2:B:336:ARG:HG3	2.08	0.53
1:A:291:LEU:HD11	1:A:313:VAL:CG1	2.40	0.52
2:B:432:GLU:O	2:B:435:GLN:HG2	2.11	0.51
1:A:209:VAL:HG13	1:A:242:TYR:HD1	1.75	0.51
1:A:763:TYR:HE1	1:A:765:ALA:HA	1.75	0.51
2:B:425:ARG:HA	2:B:430:ILE:HD12	1.92	0.51
1:A:262:ILE:HG13	1:A:263:ASN:HD22	1.76	0.51
1:A:537:GLU:HG3	1:A:544:LEU:HG	1.92	0.51
2:B:418:LYS:O	2:B:421:PHE:HB2	2.10	0.50
1:A:341:PRO:HG3	1:A:816:LEU:HD21	1.93	0.50
1:A:646:TRP:CZ3	1:A:647:LYS:HE2	2.46	0.50
1:A:773:TYR:CE2	1:A:808:PRO:HB3	2.46	0.50
1:A:662:VAL:HG13	1:A:748:VAL:HG22	1.94	0.50
1:A:391:TYR:CE2	2:B:310:PRO:HG3	2.46	0.50
1:A:755:PRO:HA	1:A:758:ARG:NH1	2.27	0.50
1:A:647:LYS:O	1:A:651:VAL:HG23	2.12	0.50
1:A:319:THR:CB	1:A:572:SER:HB3	2.40	0.50
1:A:209:VAL:O	1:A:213:ILE:HG13	2.11	0.50
2:B:376:ILE:HG22	2:B:376:ILE:O	2.11	0.50
1:A:346:SER:HB3	1:A:351:MET:CE	2.41	0.49
1:A:346:SER:HB3	1:A:351:MET:HE3	1.93	0.49
1:A:386:LEU:O	1:A:389:THR:OG1	2.24	0.49
1:A:313:VAL:HG12	1:A:314:GLY:H	1.78	0.49
1:A:363:TYR:N	1:A:363:TYR:CD1	2.80	0.49
2:B:383:TRP:CD2	2:B:412:LYS:HE2	2.48	0.49
2:B:327:ASN:HB2	2:B:330:ALA:HB2	1.94	0.49
1:A:801:GLU:CG	1:A:809:ALA:HA	2.43	0.49
1:A:695:TRP:HE1	1:A:706:LEU:HD22	1.77	0.49
1:A:385:LEU:O	1:A:388:ALA:HB3	2.13	0.48
1:A:458:LEU:HD22	1:A:487:LEU:HB2	1.93	0.48
2:B:368:GLU:HB2	2:B:369:PRO:HD3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:GLY:HA3	4:A:901:FAD:O5B	2.14	0.48
1:A:384:ARG:NH2	2:B:312:LYS:O	2.47	0.47
1:A:782:PRO:O	1:A:792:PRO:HG2	2.15	0.47
2:B:340:MET:O	2:B:343:VAL:HG12	2.15	0.47
1:A:724:VAL:HG11	1:A:746:THR:HG21	1.97	0.47
1:A:569:ASN:OD1	1:A:569:ASN:N	2.42	0.47
1:A:437:THR:HG21	1:A:507:LYS:HG2	1.97	0.46
1:A:363:TYR:CE2	1:A:734:ILE:HG23	2.51	0.46
1:A:667:ASP:N	1:A:667:ASP:OD1	2.48	0.46
1:A:538:PHE:CE1	1:A:706:LEU:HD23	2.51	0.46
1:A:475:THR:HA	2:B:393:GLN:OE1	2.16	0.46
1:A:804:ILE:O	1:A:808:PRO:HD3	2.16	0.46
1:A:728:LEU:HD11	1:A:743:PRO:HD3	1.97	0.46
1:A:448:MET:HG3	1:A:497:LEU:HD23	1.97	0.46
1:A:694:PHE:HA	1:A:704:LEU:O	2.15	0.46
1:A:449:VAL:O	1:A:453:GLU:HG2	2.17	0.45
1:A:198:ASP:N	1:A:198:ASP:OD1	2.49	0.45
1:A:297:LEU:HA	1:A:297:LEU:HD23	1.78	0.45
1:A:363:TYR:HD1	1:A:363:TYR:N	2.15	0.45
1:A:627:LEU:HG	1:A:627:LEU:O	2.16	0.45
1:A:258:ARG:NH1	1:A:827:ASP:OD1	2.50	0.45
1:A:485:ARG:HD2	2:B:404:ALA:HA	1.99	0.44
2:B:342:LEU:HD12	2:B:342:LEU:HA	1.79	0.44
1:A:432:LYS:HA	1:A:435:VAL:HG22	1.98	0.44
1:A:329:LEU:HA	1:A:661:LYS:HD2	1.98	0.44
1:A:520:TYR:CD2	1:A:521:LEU:HG	2.53	0.44
1:A:720:ASP:O	1:A:724:VAL:HG23	2.18	0.44
1:A:600:CYS:SG	1:A:795:ARG:HB3	2.58	0.44
1:A:685:THR:O	1:A:688:ARG:HB3	2.18	0.44
1:A:379:GLU:O	1:A:382:PHE:HB3	2.18	0.44
1:A:763:TYR:CE1	1:A:765:ALA:HA	2.52	0.44
1:A:634:PRO:HA	1:A:635:PRO:HD3	1.77	0.44
1:A:807:TYR:N	1:A:808:PRO:CD	2.80	0.44
1:A:760:SER:CB	4:A:901:FAD:HM83	2.47	0.43
1:A:229:LEU:N	1:A:263:ASN:OD1	2.39	0.43
1:A:218:LEU:HD23	1:A:218:LEU:HA	1.80	0.43
1:A:355:LYS:HA	1:A:565:LEU:HD23	2.00	0.43
1:A:269:ARG:HG2	1:A:271:LYS:O	2.19	0.43
1:A:669:VAL:HG13	1:A:671:TRP:CE2	2.54	0.43
1:A:743:PRO:HB2	1:A:746:THR:HG23	2.00	0.43
1:A:463:LYS:HD3	1:A:463:LYS:HA	1.92	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:485:ARG:HG3	2:B:407:ASP:HB2	2.01	0.42
1:A:574:VAL:HB	1:A:575:PRO:HD3	2.00	0.42
2:B:430:ILE:HG22	2:B:434:LEU:HD12	1.99	0.42
1:A:261:LEU:HD23	1:A:261:LEU:HA	1.81	0.42
1:A:716:GLU:CG	1:A:750:ARG:HG2	2.50	0.42
1:A:624:THR:HG22	1:A:799:ALA:HB3	2.01	0.42
1:A:548:SER:O	1:A:552:TRP:HB3	2.19	0.42
1:A:366:ASN:OD1	1:A:368:GLN:N	2.36	0.42
1:A:441:LEU:HD23	2:B:356:ASN:HD22	1.85	0.42
1:A:452:LYS:HE2	1:A:456:LYS:HE3	2.01	0.42
1:A:536:LEU:HA	1:A:536:LEU:HD12	1.82	0.42
1:A:418:LEU:HD11	2:B:321:VAL:HG22	2.00	0.42
1:A:283:ILE:HD12	1:A:294:ALA:HB2	2.01	0.42
1:A:536:LEU:HB3	1:A:544:LEU:HD21	2.02	0.42
1:A:781:THR:HA	1:A:782:PRO:HD2	1.65	0.42
2:B:317:SER:HB3	2:B:320:ASP:OD1	2.19	0.42
1:A:559:GLU:HG2	1:A:560:PHE:N	2.35	0.41
1:A:671:TRP:O	1:A:673:PRO:HD3	2.20	0.41
1:A:188:MET:HG2	1:A:210:PHE:CE2	2.54	0.41
1:A:594:ARG:HA	1:A:640:VAL:O	2.20	0.41
1:A:386:LEU:HA	1:A:386:LEU:HD23	1.84	0.41
1:A:282:ILE:HD13	1:A:282:ILE:HA	1.78	0.41
1:A:784:PRO:HB2	1:A:786:ILE:O	2.20	0.41
1:A:222:LEU:HA	1:A:222:LEU:HD23	1.87	0.41
1:A:760:SER:HB2	4:A:901:FAD:HM83	2.03	0.41
1:A:663:VAL:HG22	1:A:704:LEU:CD2	2.51	0.41
1:A:707:VAL:HG12	1:A:712:ALA:HA	2.03	0.41
4:A:901:FAD:H9	4:A:901:FAD:H1'1	1.58	0.40
1:A:451:LEU:HD23	1:A:494:TYR:HB2	2.03	0.40
1:A:644:PRO:O	1:A:648:THR:HG23	2.22	0.40
1:A:775:LEU:HA	1:A:775:LEU:HD23	1.85	0.40
2:B:401:ASP:O	2:B:405:ILE:HG13	2.21	0.40
1:A:354:ALA:C	1:A:565:LEU:HD23	2.42	0.40
1:A:360:CYS:HB2	1:A:677:LEU:CD1	2.51	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	664/730 (91%)	628 (95%)	33 (5%)	3 (0%)	29	64
2	B	131/178 (74%)	128 (98%)	3 (2%)	0	100	100
3	C	2/11 (18%)	0	1 (50%)	1 (50%)	0	0
All	All	797/919 (87%)	756 (95%)	37 (5%)	4 (0%)	29	64

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	182	ARG
3	C	7	LEU
1	A	365	ALA
1	A	741	PRO

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	566/623 (91%)	559 (99%)	7 (1%)	71	88
2	B	117/156 (75%)	116 (99%)	1 (1%)	78	91
3	C	2/2 (100%)	0	2 (100%)	0	0
All	All	685/781 (88%)	675 (98%)	10 (2%)	65	85

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	191	GLN
1	A	255	TYR
1	A	363	TYR
1	A	571	TYR
1	A	615	ILE
1	A	793	ILE
1	A	801	GLU
2	B	329	THR
3	C	7	LEU
3	C	8	LEU

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

Mol	Chain	Res	Type
1	A	394	HIS
1	A	399	ASN
1	A	802	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

8 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
3	DAB	C	10	3	5,6,7	0.51	0	1,6,8	0.76	0
3	DAB	C	9	3	5,6,7	0.52	0	1,6,8	0.02	0
3	4FO	C	2	3	5,6,7	0.52	0	1,6,8	0.22	0
3	4FO	C	4	3	5,6,7	0.47	0	1,6,8	0.31	0
3	4FO	C	5	3	5,6,7	0.47	0	1,6,8	0.07	0
3	DAB	C	6	3	5,6,7	0.57	0	1,6,8	0.05	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	DAB	C	10	3	-	1/4/5/7	-
3	DAB	C	9	3	-	2/4/5/7	-
3	4FO	C	2	3	-	1/4/5/7	-
3	4FO	C	4	3	-	2/4/5/7	-
3	4FO	C	5	3	-	4/4/5/7	-
3	DAB	C	6	3	-	3/4/5/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	10	DAB	CA-CB-CG-ND
3	C	9	DAB	CA-CB-CG-ND
3	C	4	4FO	O-C-CA-CB
3	C	5	4FO	O-C-CA-CB
3	C	5	4FO	N-CA-CB-CG
3	C	5	4FO	C-CA-CB-CG
3	C	5	4FO	CA-CB-CG-NZ
3	C	2	4FO	CA-CB-CG-NZ
3	C	6	DAB	N-CA-CB-CG
3	C	6	DAB	CA-CB-CG-ND
3	C	9	DAB	N-CA-CB-CG
3	C	4	4FO	CA-CB-CG-NZ
3	C	6	DAB	C-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	6	DAB	1	0

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	FAD	A	901	-	51,58,58	4.27	19 (37%)	60,89,89	2.66	17 (28%)

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	-	-	2/30/50/50	0/6/6/6

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	FAD	C2B-C1B	-17.03	1.27	1.53
4	A	901	FAD	O4B-C1B	13.75	1.60	1.41
4	A	901	FAD	C10-N1	9.67	1.45	1.33
4	A	901	FAD	C5X-N5	7.01	1.46	1.35
4	A	901	FAD	O4B-C4B	-6.42	1.30	1.45
4	A	901	FAD	C4-N3	5.68	1.42	1.33
4	A	901	FAD	C9A-N10	5.25	1.45	1.38
4	A	901	FAD	C2-N1	5.23	1.48	1.38
4	A	901	FAD	C2-N3	4.84	1.47	1.38
4	A	901	FAD	C4X-C10	4.62	1.43	1.38
4	A	901	FAD	C4X-N5	4.52	1.39	1.33
4	A	901	FAD	O2'-C2'	-3.98	1.34	1.43
4	A	901	FAD	C2A-N3A	3.91	1.38	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	FAD	C4-C4X	3.73	1.47	1.41
4	A	901	FAD	C6A-N6A	3.07	1.45	1.34
4	A	901	FAD	C4A-N3A	2.89	1.39	1.35
4	A	901	FAD	C1'-N10	-2.54	1.45	1.48
4	A	901	FAD	O3B-C3B	-2.34	1.37	1.43
4	A	901	FAD	C2A-N1A	2.33	1.38	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	FAD	C5A-C6A-N6A	9.85	135.32	120.35
4	A	901	FAD	C1'-N10-C10	6.74	124.45	118.41
4	A	901	FAD	N6A-C6A-N1A	-6.41	105.28	118.57
4	A	901	FAD	O4B-C1B-C2B	-6.16	97.92	106.93
4	A	901	FAD	C5'-C4'-C3'	-4.75	103.04	112.20
4	A	901	FAD	C4-N3-C2	4.70	119.11	115.14
4	A	901	FAD	C1'-N10-C9A	-3.90	115.23	118.29
4	A	901	FAD	C4X-N5-C5X	3.89	120.66	116.77
4	A	901	FAD	N3A-C2A-N1A	-3.86	122.64	128.68
4	A	901	FAD	C7M-C7-C8	3.77	128.46	120.74
4	A	901	FAD	C7M-C7-C6	-3.58	111.77	120.34
4	A	901	FAD	C5X-C9A-N10	3.43	120.20	117.72
4	A	901	FAD	C4-C4X-C10	3.26	122.11	119.95
4	A	901	FAD	O5'-P-O1P	-3.15	96.74	109.07
4	A	901	FAD	O4'-C4'-C3'	2.69	115.65	109.10
4	A	901	FAD	P-O3P-PA	-2.57	123.99	132.83
4	A	901	FAD	C4X-C4-N3	-2.09	120.57	123.43

There are no chirality outliers.

All (2) torsion outliers are listed below:

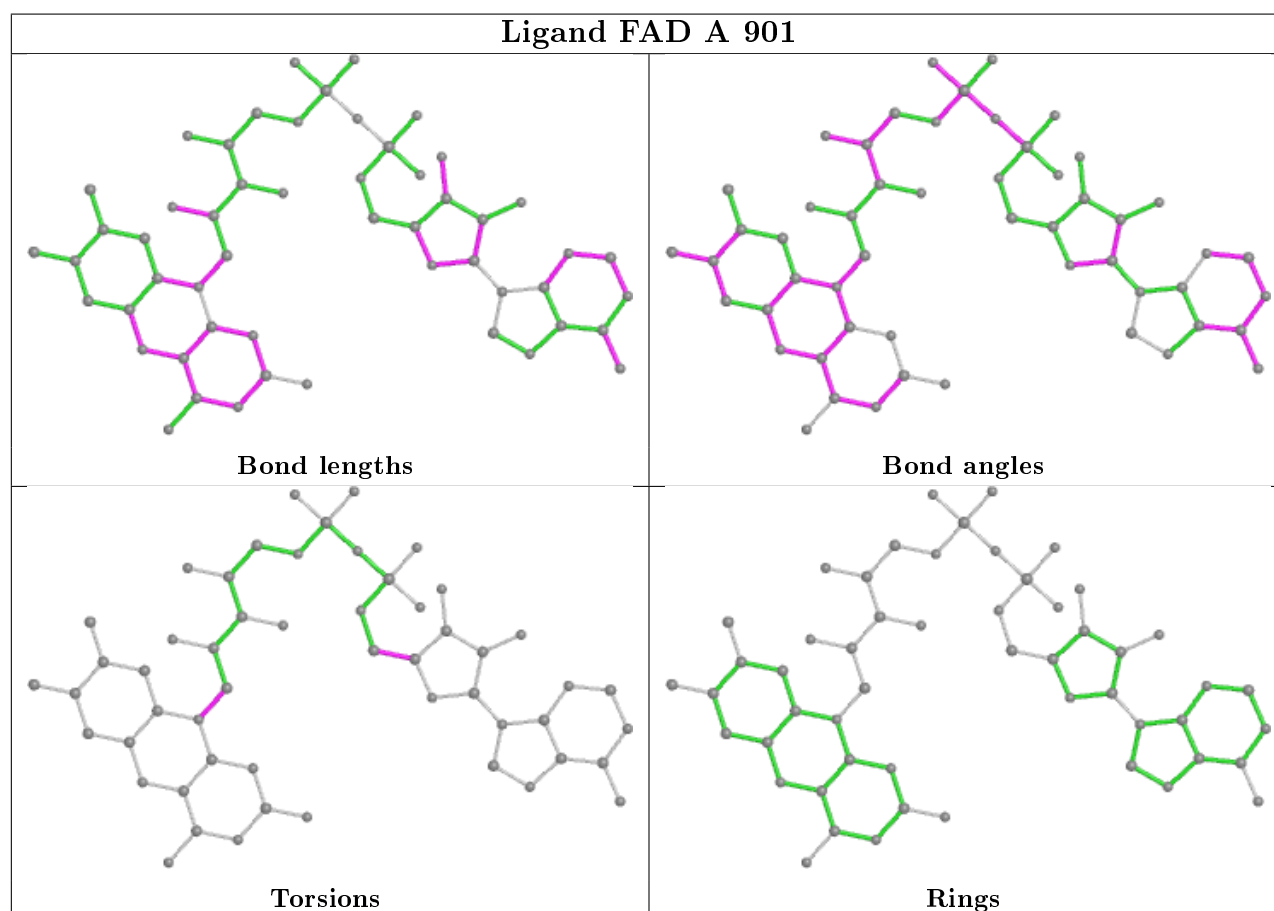
Mol	Chain	Res	Type	Atoms
4	A	901	FAD	C2'-C1'-N10-C9A
4	A	901	FAD	O4B-C4B-C5B-O5B

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	901	FAD	6	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	666/730 (91%)	-0.03	4 (0%) 89 78	71, 115, 171, 213	0
2	B	133/178 (74%)	0.31	6 (4%) 33 16	105, 156, 205, 247	0
3	C	2/11 (18%)	0.32	0 100 100	169, 169, 169, 178	0
All	All	801/919 (87%)	0.03	10 (1%) 79 61	71, 123, 182, 247	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	171	PRO	3.9
2	B	308	ARG	2.8
1	A	398	PHE	2.6
2	B	312	LYS	2.5
1	A	401	LEU	2.4
2	B	309	LYS	2.2
2	B	324	VAL	2.2
1	A	836	LEU	2.2
2	B	316	LEU	2.2
2	B	375	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
3	DAB	C	6	7/8	0.85	0.25	154,184,189,194	0
3	4FO	C	2	7/8	-	-	174,176,177,177	7
3	4FO	C	5	7/8	0.77	0.19	190,197,201,207	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	4FO	C	4	7/8	0.80	0.21	167,171,178,178	5
3	DAB	C	10	7/8	0.83	0.35	146,166,190,191	0
3	DAB	C	9	7/8	0.89	0.22	166,181,196,197	0
3	DTH	C	11	7/8	0.90	0.23	170,173,183,190	0
3	DTH	C	3	7/8	-	-	177,177,179,179	7

6.3 Carbohydrates [i](#)

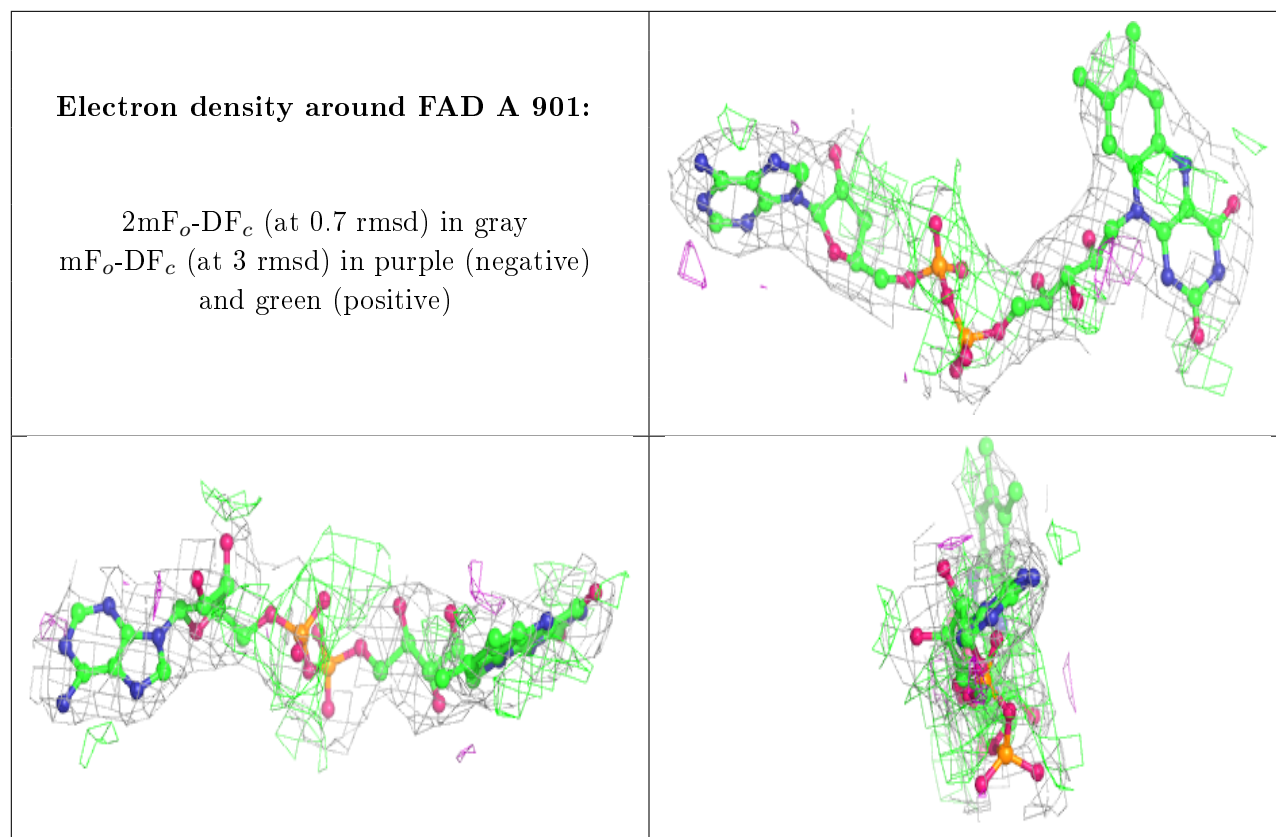
There are no carbohydrates in this entry.

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, 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(\AA^2)	Q<0.9
5	NA	A	903	1/1	0.82	0.79	88,88,88,88	0
5	NA	A	902	1/1	0.92	0.27	73,73,73,73	0
4	FAD	A	901	53/53	0.95	0.27	67,92,113,120	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.