



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 26, 2021 – 06:39 PM EDT

PDB ID : 6W4K
Title : Crystal structure of Lysine Specific Demethylase 1 (LSD1) with CC-90011
Authors : Hosfield, D.J.
Deposited on : 2020-03-11
Resolution : 2.93 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 : 2.23.2
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) : 2.23.2

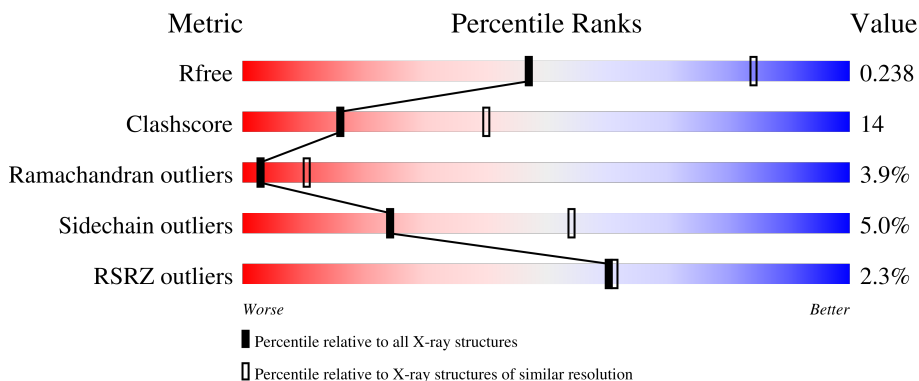
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.93 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	2969 (2.98-2.90)
Clashscore	141614	3218 (2.98-2.90)
Ramachandran outliers	138981	3122 (2.98-2.90)
Sidechain outliers	138945	3124 (2.98-2.90)
RSRZ outliers	127900	2902 (2.98-2.90)

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	659	<div> <div>2%</div> <div>69%</div> <div>26%</div> <div>...</div> </div>
2	B	132	<div> <div>4%</div> <div>61%</div> <div>28%</div> <div>6%</div> <div>5%</div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6223 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

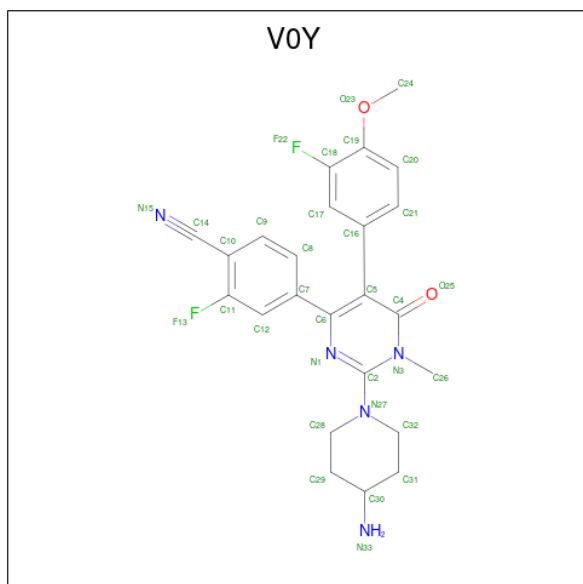
- Molecule 1 is a protein called Lysine-specific histone demethylase 1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	651	Total	C	N	O	S	0	0	0
			5107	3254	888	946	19			

- Molecule 2 is a protein called REST corepressor 1.

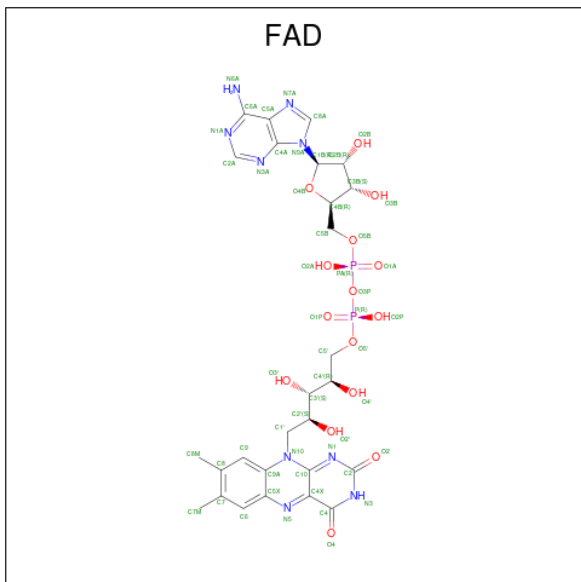
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	125	Total	C	N	O	S	0	0	0
			1008	632	181	192	3			

- Molecule 3 is 4-[2-(4-aminopiperidin-1-yl)-5-(3-fluoro-4-methoxyphenyl)-1-methyl-6-oxo-1,6-dihydropyrimidin-4-yl]-2-fluorobenzonitrile (three-letter code: V0Y) (formula: $C_{24}H_{23}F_2N_5O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			33	24	2	5	2		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $\text{C}_{27}\text{H}_{33}\text{N}_9\text{O}_{15}\text{P}_2$) (labeled as "Ligand of Interest" by depositor).



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

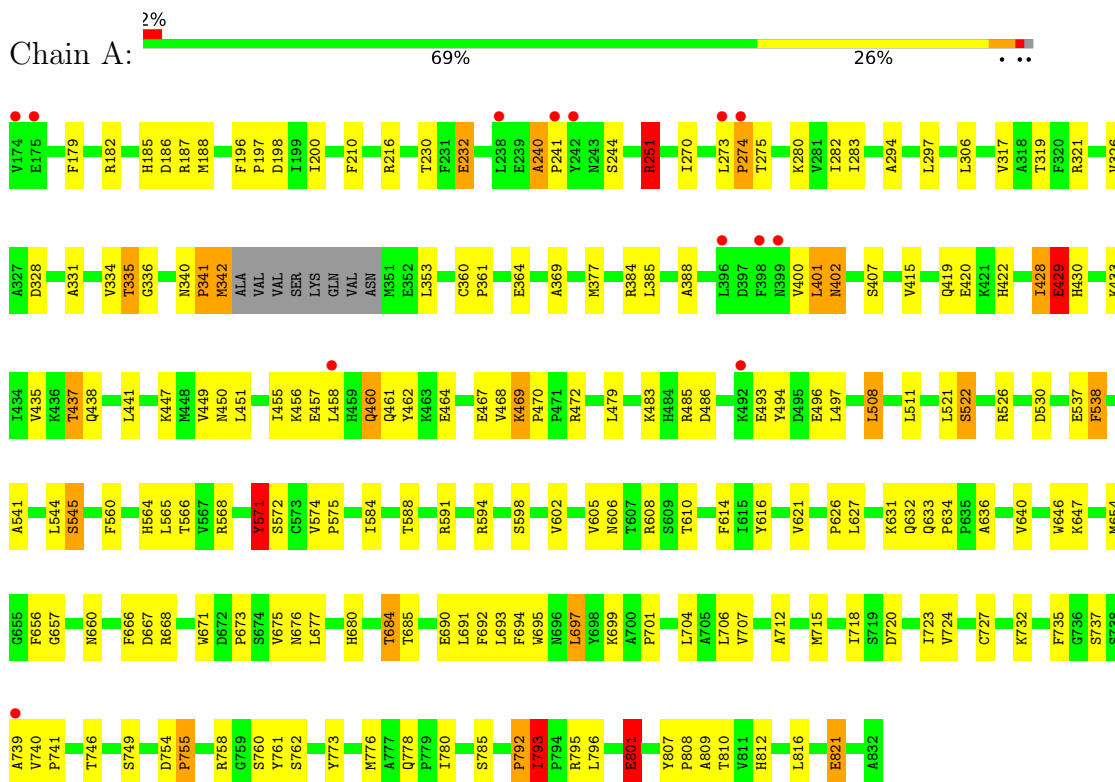
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	22	Total O 22 22	0	0

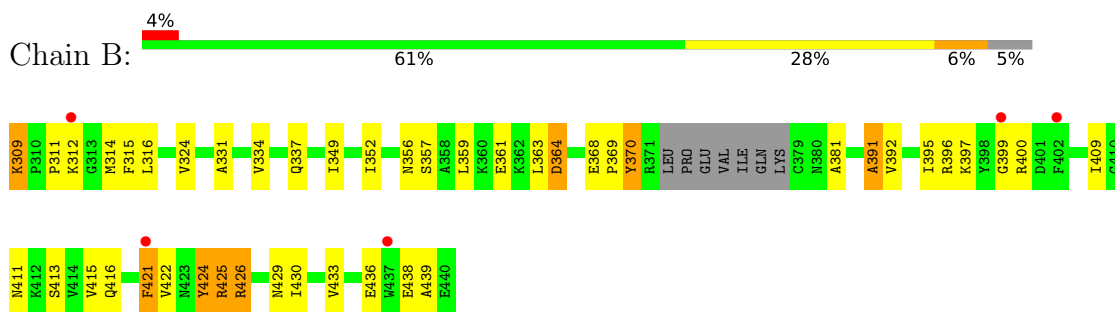
3 Residue-property plots [i](#)

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



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.66Å 178.51Å 236.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.27 – 2.93 45.64 – 2.93	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.27-2.93) 99.4 (45.64-2.93)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.191 , 0.235 0.200 , 0.238	Depositor DCC
R_{free} test set	2723 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	84.2	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 62.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6223	wwPDB-VP
Average B, all atoms (Å ²)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% 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: FAD, V0Y

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	0.78	2/5218 (0.0%)	1.03	8/7077 (0.1%)
2	B	0.79	0/1021	1.05	2/1375 (0.1%)
All	All	0.78	2/6239 (0.0%)	1.03	10/8452 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
2	B	0	1
All	All	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	342	MET	C-O	6.83	1.36	1.23
1	A	821	GLU	CD-OE2	-5.19	1.20	1.25

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	426	ARG	NE-CZ-NH1	9.17	124.88	120.30
1	A	251	ARG	NE-CZ-NH2	7.38	123.99	120.30
1	A	251	ARG	CG-CD-NE	6.13	124.67	111.80
1	A	795	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	A	571	TYR	CB-CG-CD1	5.53	124.32	121.00
1	A	795	ARG	NE-CZ-NH2	-5.38	117.61	120.30
2	B	424	TYR	CB-CA-C	5.31	121.01	110.40

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	496	GLU	CB-CA-C	5.21	120.83	110.40
1	A	216	ARG	CG-CD-NE	-5.09	101.11	111.80
1	A	801	GLU	N-CA-CB	-5.08	101.45	110.60

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	469	LYS	Peptide
1	A	697	LEU	Peptide
1	A	792	PRO	Peptide
2	B	438	GLU	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5107	0	5135	145	0
2	B	1008	0	1012	44	0
3	A	33	0	0	1	0
4	A	53	0	31	4	0
5	A	22	0	0	6	0
All	All	6223	0	6178	179	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 14.

All (179) 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:437:THR:HG22	1:A:508:LEU:HD12	1.62	0.80
1:A:633:GLN:NE2	1:A:634:PRO:CD	2.46	0.78
1:A:633:GLN:HG2	1:A:634:PRO:HD3	1.64	0.78
1:A:633:GLN:NE2	1:A:634:PRO:HD2	2.00	0.76
1:A:606:ASN:HD21	1:A:608:ARG:HH21	1.32	0.75

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:424:TYR:O	2:B:426:ARG:N	2.21	0.74
1:A:240:ALA:HB1	1:A:241:PRO:CD	2.18	0.73
1:A:762:SER:OG	1:A:801:GLU:OE2	2.04	0.72
1:A:437:THR:CG2	1:A:508:LEU:HD12	2.19	0.71
1:A:420:GLU:OE2	1:A:526:ARG:NH1	2.25	0.70
5:A:1012:HOH:O	2:B:352:ILE:HD11	1.94	0.68
2:B:369:PRO:O	2:B:370:TYR:CD1	2.47	0.67
2:B:369:PRO:C	2:B:370:TYR:CD1	2.69	0.67
1:A:807:TYR:N	1:A:808:PRO:HD3	2.10	0.67
2:B:396:ARG:HD3	2:B:436:GLU:OE1	1.95	0.66
2:B:369:PRO:C	2:B:370:TYR:HD1	2.00	0.65
1:A:340:ASN:HB2	1:A:560:PHE:CD2	2.32	0.65
1:A:666:PHE:O	1:A:701:PRO:HG2	1.97	0.65
1:A:633:GLN:CG	1:A:634:PRO:HD3	2.29	0.62
2:B:381:ALA:HA	2:B:416:GLN:HE22	1.63	0.62
1:A:693:LEU:HD12	1:A:694:PHE:N	2.15	0.62
1:A:538:PHE:CE1	1:A:706:LEU:HD23	2.36	0.60
1:A:449:VAL:HG23	2:B:363:LEU:CD2	2.32	0.60
1:A:654:MET:HE1	1:A:773:TYR:CE1	2.36	0.59
2:B:395:ILE:HG22	2:B:433:VAL:CG1	2.32	0.59
1:A:331:ALA:HA	4:A:902:FAD:C4X	2.32	0.59
1:A:530:ASP:OD2	1:A:685:THR:HA	2.02	0.58
1:A:776:MET:HA	1:A:776:MET:HE3	1.84	0.58
1:A:415:VAL:HG22	2:B:316:LEU:HD21	1.84	0.58
2:B:422:VAL:O	2:B:425:ARG:HB2	2.03	0.58
1:A:438:GLN:HB3	1:A:508:LEU:HD11	1.85	0.58
1:A:451:LEU:HD23	1:A:494:TYR:HB2	1.86	0.58
1:A:633:GLN:NE2	1:A:634:PRO:HD3	2.18	0.58
1:A:684:THR:HG22	5:A:1021:HOH:O	2.04	0.58
1:A:654:MET:HE1	1:A:773:TYR:CZ	2.39	0.58
1:A:435:VAL:HG12	2:B:349:ILE:HG13	1.87	0.57
1:A:198:ASP:HB2	1:A:251:ARG:HH22	1.70	0.57
1:A:240:ALA:HB1	1:A:241:PRO:HD2	1.87	0.57
1:A:660:ASN:HA	1:A:749:SER:OG	2.05	0.56
1:A:449:VAL:HG12	1:A:450:ASN:OD1	2.05	0.56
1:A:240:ALA:CB	1:A:241:PRO:CD	2.83	0.56
1:A:185:HIS:CE1	1:A:186:ASP:HB3	2.40	0.56
1:A:739:ALA:O	1:A:741:PRO:HD3	2.05	0.56
3:A:901:V0Y:C28	3:A:901:V0Y:C26	2.84	0.56
1:A:341:PRO:HG3	1:A:816:LEU:HD21	1.88	0.56
2:B:311:PRO:HG2	2:B:314:MET:HG3	1.89	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:ARG:NH2	2:B:312:LYS:O	2.39	0.55
1:A:460:GLN:O	1:A:462:TYR:N	2.40	0.55
1:A:720:ASP:O	1:A:724:VAL:HG23	2.07	0.55
1:A:331:ALA:HA	4:A:902:FAD:N5	2.22	0.55
1:A:537:GLU:OE2	1:A:544:LEU:HG	2.07	0.54
1:A:760:SER:HB2	4:A:902:FAD:HM83	1.89	0.54
1:A:182:ARG:NH1	1:A:341:PRO:HD3	2.23	0.54
1:A:695:TRP:CE3	1:A:697:LEU:HD21	2.42	0.54
1:A:187:ARG:HB2	1:A:187:ARG:NH1	2.21	0.54
1:A:282:ILE:HG21	1:A:602:VAL:HG21	1.90	0.54
1:A:646:TRP:CZ3	1:A:647:LYS:HE2	2.43	0.53
1:A:240:ALA:HB1	1:A:241:PRO:HD3	1.89	0.53
1:A:401:LEU:O	1:A:402:ASN:HB2	2.08	0.53
1:A:319:THR:HB	1:A:572:SER:HB3	1.90	0.53
1:A:521:LEU:O	1:A:522:SER:O	2.26	0.53
1:A:467:GLU:O	1:A:469:LYS:HG3	2.08	0.53
1:A:633:GLN:CD	1:A:634:PRO:HD2	2.28	0.53
2:B:413:SER:OG	2:B:415:VAL:HG12	2.09	0.53
1:A:479:LEU:O	1:A:483:LYS:HG2	2.09	0.52
1:A:447:LYS:HE3	1:A:497:LEU:HD21	1.90	0.52
2:B:395:ILE:HG22	2:B:433:VAL:HG12	1.91	0.52
1:A:485:ARG:HD2	5:A:1014:HOH:O	2.09	0.52
1:A:666:PHE:O	1:A:701:PRO:CG	2.58	0.52
1:A:807:TYR:N	1:A:808:PRO:CD	2.72	0.52
1:A:780:ILE:HB	1:A:796:LEU:HB3	1.92	0.52
1:A:591:ARG:NH1	1:A:605:VAL:HG21	2.25	0.52
1:A:660:ASN:HA	1:A:749:SER:HG	1.75	0.51
1:A:538:PHE:CD1	1:A:706:LEU:HD23	2.46	0.51
2:B:391:ALA:HB2	2:B:409:ILE:HD11	1.93	0.51
2:B:361:GLU:O	2:B:364:ASP:HB2	2.11	0.50
2:B:426:ARG:HH11	2:B:426:ARG:HB2	1.77	0.50
1:A:456:LYS:HA	2:B:370:TYR:HE2	1.77	0.49
1:A:633:GLN:HE21	1:A:634:PRO:HD3	1.76	0.49
1:A:758:ARG:HD2	5:A:1013:HOH:O	2.12	0.49
2:B:426:ARG:HB3	2:B:426:ARG:NH1	2.28	0.49
1:A:633:GLN:CD	1:A:634:PRO:CD	2.81	0.49
1:A:732:LYS:HG2	1:A:740:VAL:HB	1.94	0.49
1:A:566:THR:HG21	1:A:697:LEU:CD2	2.42	0.49
1:A:695:TRP:HE3	1:A:697:LEU:HD21	1.78	0.49
1:A:810:THR:HB	1:A:812:HIS:CE1	2.48	0.49
1:A:428:ILE:O	1:A:430:HIS:N	2.46	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:594:ARG:HA	1:A:640:VAL:O	2.14	0.48
1:A:718:ILE:HG22	1:A:723:ILE:HG13	1.96	0.48
2:B:368:GLU:HB2	2:B:369:PRO:HD3	1.95	0.48
1:A:627:LEU:HD21	1:A:654:MET:HG3	1.95	0.48
1:A:538:PHE:CE1	1:A:706:LEU:CD2	2.97	0.48
1:A:335:THR:HG22	1:A:335:THR:O	2.15	0.47
1:A:632:GLN:OE1	1:A:636:ALA:HB2	2.14	0.47
1:A:633:GLN:HE21	1:A:634:PRO:CD	2.23	0.47
1:A:306:LEU:HD13	1:A:584:ILE:HG12	1.96	0.47
1:A:456:LYS:HA	2:B:370:TYR:CE2	2.49	0.47
1:A:792:PRO:HA	1:A:793:ILE:HD12	1.97	0.47
1:A:441:LEU:HD23	2:B:356:ASN:ND2	2.29	0.47
1:A:707:VAL:HG12	1:A:712:ALA:HA	1.97	0.47
1:A:188:MET:HE2	1:A:210:PHE:CE2	2.49	0.47
1:A:707:VAL:HG11	1:A:715:MET:HG3	1.96	0.47
1:A:821:GLU:OE1	1:A:821:GLU:HA	2.15	0.47
2:B:309:LYS:HE2	2:B:309:LYS:N	2.30	0.46
1:A:470:PRO:O	1:A:472:ARG:N	2.48	0.46
1:A:792:PRO:CA	1:A:793:ILE:HD12	2.46	0.46
1:A:187:ARG:CB	1:A:187:ARG:CZ	2.93	0.46
2:B:426:ARG:HH11	2:B:426:ARG:CB	2.28	0.46
1:A:449:VAL:HG23	2:B:363:LEU:HD23	1.97	0.46
1:A:455:ILE:HG22	2:B:370:TYR:HD2	1.80	0.46
2:B:395:ILE:HG22	2:B:433:VAL:HG11	1.96	0.46
1:A:385:LEU:O	1:A:388:ALA:HB3	2.15	0.46
1:A:724:VAL:HG11	1:A:746:THR:HG21	1.97	0.46
1:A:198:ASP:HB2	1:A:251:ARG:NH2	2.29	0.45
1:A:694:PHE:HA	1:A:704:LEU:O	2.16	0.45
1:A:407:SER:CB	1:A:545:SER:O	2.65	0.45
1:A:568:ARG:HH21	1:A:699:LYS:HD2	1.81	0.45
1:A:428:ILE:HG22	1:A:429:GLU:N	2.32	0.45
1:A:273:LEU:O	1:A:274:PRO:C	2.55	0.45
1:A:564:HIS:N	1:A:564:HIS:CD2	2.84	0.45
1:A:419:GLN:HE22	2:B:315:PHE:H	1.66	0.44
1:A:541:ALA:O	1:A:657:GLY:HA3	2.16	0.44
1:A:283:ILE:HD12	1:A:294:ALA:HB2	1.99	0.44
1:A:451:LEU:HD21	1:A:493:GLU:HG2	2.00	0.44
2:B:324:VAL:O	2:B:324:VAL:HG22	2.18	0.44
2:B:352:ILE:O	2:B:352:ILE:HG22	2.17	0.44
1:A:328:ASP:HB3	5:A:1002:HOH:O	2.17	0.44
1:A:574:VAL:HB	1:A:575:PRO:HD3	2.00	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:359:LEU:HD23	2:B:359:LEU:N	2.32	0.44
1:A:627:LEU:CD2	1:A:654:MET:HG3	2.48	0.43
1:A:317:VAL:HG13	1:A:571:TYR:HB3	1.99	0.43
1:A:511:LEU:HD23	1:A:511:LEU:N	2.33	0.43
2:B:369:PRO:HG2	2:B:370:TYR:HE1	1.83	0.43
1:A:187:ARG:CZ	1:A:187:ARG:HB3	2.49	0.43
1:A:422:HIS:O	1:A:422:HIS:CG	2.72	0.43
1:A:400:VAL:HG12	1:A:401:LEU:N	2.34	0.43
1:A:407:SER:HB3	1:A:545:SER:O	2.18	0.43
2:B:413:SER:OG	2:B:416:GLN:HG3	2.19	0.43
1:A:240:ALA:CB	1:A:241:PRO:HD2	2.47	0.42
1:A:353:LEU:HD13	1:A:565:LEU:HD12	2.01	0.42
1:A:627:LEU:HG	1:A:631:LYS:HD2	2.02	0.42
1:A:633:GLN:CG	1:A:634:PRO:CD	2.95	0.42
1:A:677:LEU:HD12	1:A:693:LEU:HD11	2.01	0.42
2:B:426:ARG:HB3	2:B:426:ARG:CZ	2.49	0.42
1:A:297:LEU:HD23	1:A:297:LEU:HA	1.86	0.42
1:A:334:VAL:O	1:A:336:GLY:N	2.49	0.42
1:A:455:ILE:HG22	2:B:370:TYR:CD2	2.53	0.42
1:A:761:TYR:CD1	1:A:809:ALA:HB1	2.53	0.42
1:A:282:ILE:O	1:A:621:VAL:HA	2.20	0.42
1:A:321:ARG:HG2	1:A:326:VAL:HG22	2.02	0.42
1:A:690:GLU:O	1:A:691:LEU:C	2.58	0.42
1:A:196:PHE:N	1:A:197:PRO:CD	2.82	0.42
1:A:656:PHE:CE1	1:A:761:TYR:HA	2.55	0.42
1:A:778:GLN:HG3	5:A:1019:HOH:O	2.19	0.42
1:A:796:LEU:HD12	1:A:796:LEU:HA	1.87	0.41
2:B:426:ARG:NH1	2:B:426:ARG:CB	2.83	0.41
1:A:754:ASP:OD1	1:A:755:PRO:HD2	2.19	0.41
2:B:334:VAL:HA	2:B:337:GLN:NE2	2.35	0.41
2:B:421:PHE:O	2:B:425:ARG:HG3	2.20	0.41
1:A:360:CYS:HA	1:A:361:PRO:HD2	1.92	0.41
1:A:460:GLN:HE21	1:A:464:GLU:HG3	1.86	0.41
1:A:754:ASP:OD1	1:A:754:ASP:C	2.59	0.41
1:A:626:PRO:HD3	4:A:902:FAD:H51A	2.03	0.41
2:B:411:ASN:CG	2:B:411:ASN:O	2.58	0.41
1:A:660:ASN:CA	1:A:749:SER:OG	2.69	0.41
1:A:671:TRP:HA	1:A:735:PHE:CE1	2.55	0.41
1:A:187:ARG:HB2	1:A:187:ARG:HH11	1.84	0.41
1:A:230:THR:HG23	1:A:270:ILE:HD12	2.03	0.41
1:A:230:THR:OG1	1:A:232:GLU:HB3	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:808:PRO:O	1:A:810:THR:HG23	2.20	0.41
1:A:282:ILE:HD13	1:A:282:ILE:HA	1.93	0.41
1:A:671:TRP:O	1:A:673:PRO:HD3	2.21	0.41
2:B:311:PRO:HG2	2:B:314:MET:CG	2.50	0.41
1:A:614:PHE:HB3	1:A:616:TYR:CE1	2.56	0.40
2:B:391:ALA:CB	2:B:409:ILE:HD11	2.51	0.40
2:B:425:ARG:HA	2:B:430:ILE:HG13	2.02	0.40
1:A:691:LEU:HD22	1:A:727:CYS:SG	2.61	0.40
1:A:198:ASP:CB	1:A:251:ARG:HH22	2.34	0.40
1:A:437:THR:HG22	1:A:508:LEU:CD1	2.43	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	647/659 (98%)	542 (84%)	82 (13%)	23 (4%)	3	12
2	B	121/132 (92%)	96 (79%)	18 (15%)	7 (6%)	1	4
All	All	768/791 (97%)	638 (83%)	100 (13%)	30 (4%)	3	10

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	335	THR
1	A	522	SER
1	A	610	THR
1	A	793	ILE
2	B	331	ALA
2	B	425	ARG
2	B	439	ALA
1	A	232	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	240	ALA
1	A	244	SER
1	A	275	THR
1	A	364	GLU
1	A	402	ASN
1	A	428	ILE
1	A	429	GLU
1	A	460	GLN
1	A	461	GLN
1	A	468	VAL
1	A	801	GLU
2	B	391	ALA
2	B	399	GLY
1	A	274	PRO
1	A	341	PRO
1	A	369	ALA
2	B	429	ASN
1	A	401	LEU
1	A	457	GLU
1	A	486	ASP
1	A	737	SER
2	B	392	VAL

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	553/560 (99%)	527 (95%)	26 (5%)	26	57
2	B	109/116 (94%)	102 (94%)	7 (6%)	17	43
All	All	662/676 (98%)	629 (95%)	33 (5%)	24	54

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	179	PHE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	200	ILE
1	A	251	ARG
1	A	280	LYS
1	A	342	MET
1	A	377	MET
1	A	429	GLU
1	A	433	LYS
1	A	437	THR
1	A	458	LEU
1	A	508	LEU
1	A	538	PHE
1	A	545	SER
1	A	571	TYR
1	A	588	THR
1	A	598	SER
1	A	667	ASP
1	A	668	ARG
1	A	675	VAL
1	A	676	ASN
1	A	680	HIS
1	A	684	THR
1	A	692	PHE
1	A	755	PRO
1	A	785	SER
1	A	793	ILE
2	B	309	LYS
2	B	357	SER
2	B	364	ASP
2	B	370	TYR
2	B	397	LYS
2	B	400	ARG
2	B	421	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	191	GLN
1	A	395	GLN
1	A	422	HIS
1	A	460	GLN
1	A	564	HIS
1	A	680	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	812	HIS
1	A	828	GLN
2	B	416	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
3	V0Y	A	901	-	34,36,36	2.32	10 (29%)	42,52,52	3.35	21 (50%)
4	FAD	A	902	-	51,58,58	1.13	3 (5%)	60,89,89	2.07	9 (15%)

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	V0Y	A	901	-	-	3/16/26/26	0/4/4/4

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FAD	A	902	-	-	13/30/50/50	0/6/6/6

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	901	V0Y	C10-C14	-8.58	1.31	1.44
4	A	902	FAD	C4X-C10	5.20	1.44	1.38
3	A	901	V0Y	C32-N27	-4.71	1.39	1.46
3	A	901	V0Y	C30-N33	-3.67	1.34	1.46
3	A	901	V0Y	C5-C16	3.38	1.54	1.49
3	A	901	V0Y	C28-N27	-3.33	1.41	1.46
3	A	901	V0Y	C17-C18	2.76	1.42	1.37
4	A	902	FAD	C4-N3	2.73	1.37	1.33
3	A	901	V0Y	C20-C19	-2.72	1.33	1.39
3	A	901	V0Y	C5-C6	2.58	1.45	1.41
4	A	902	FAD	C4-C4X	-2.49	1.37	1.41
3	A	901	V0Y	C4-C5	-2.29	1.38	1.43
3	A	901	V0Y	C4-N3	2.19	1.41	1.38

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	V0Y	O23-C19-C18	15.41	126.95	116.26
4	A	902	FAD	C4-N3-C2	9.32	123.01	115.14
4	A	902	FAD	C10-C4X-N5	6.49	125.74	121.26
3	A	901	V0Y	O23-C19-C20	-6.31	113.55	124.37
3	A	901	V0Y	C8-C7-C6	-5.01	112.67	120.61
4	A	902	FAD	C4-C4X-C10	-4.95	116.68	119.95
4	A	902	FAD	C4X-C10-N10	-4.79	115.38	120.30
3	A	901	V0Y	C17-C16-C5	4.30	128.37	120.06
4	A	902	FAD	C4X-C4-N3	-4.22	117.65	123.43
3	A	901	V0Y	C12-C7-C6	3.88	127.62	120.15
3	A	901	V0Y	C16-C17-C18	3.87	122.88	119.59
3	A	901	V0Y	C10-C14-N15	-3.60	171.94	177.88
4	A	902	FAD	C1'-N10-C9A	3.41	120.98	118.29
3	A	901	V0Y	C17-C18-C19	-3.24	118.05	122.81
3	A	901	V0Y	F13-C11-C10	-3.00	113.70	117.67
4	A	902	FAD	O4B-C1B-C2B	-2.72	102.94	106.93
3	A	901	V0Y	C9-C10-C11	2.72	120.11	117.49
3	A	901	V0Y	C21-C16-C17	-2.65	114.42	118.16
4	A	902	FAD	O2P-P-O1P	2.54	124.78	112.24
3	A	901	V0Y	C7-C12-C11	-2.51	117.45	119.59

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	V0Y	F13-C11-C12	2.46	123.51	118.61
3	A	901	V0Y	C32-N27-C2	-2.38	116.94	122.03
3	A	901	V0Y	C21-C16-C5	-2.37	117.18	120.77
3	A	901	V0Y	C4-C5-C16	-2.27	119.53	122.44
3	A	901	V0Y	C7-C6-N1	-2.23	112.13	115.09
3	A	901	V0Y	C24-O23-C19	-2.22	114.18	117.53
3	A	901	V0Y	C9-C10-C14	2.21	123.25	119.40
3	A	901	V0Y	C6-N1-C2	2.15	120.26	117.14
3	A	901	V0Y	C16-C5-C6	2.13	124.58	121.85
4	A	902	FAD	C5A-C6A-N6A	2.05	123.46	120.35

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	901	V0Y	C18-C19-O23-C24
4	A	902	FAD	C5B-O5B-PA-O1A
4	A	902	FAD	O4'-C4'-C5'-O5'
4	A	902	FAD	C5'-O5'-P-O3P
3	A	901	V0Y	N1-C2-N27-C28
3	A	901	V0Y	C20-C19-O23-C24
4	A	902	FAD	C2'-C3'-C4'-O4'
4	A	902	FAD	C3'-C4'-C5'-O5'
4	A	902	FAD	C5B-O5B-PA-O3P
4	A	902	FAD	C5B-O5B-PA-O2A
4	A	902	FAD	C5'-O5'-P-O1P
4	A	902	FAD	C5'-O5'-P-O2P
4	A	902	FAD	C2'-C3'-C4'-C5'
4	A	902	FAD	O3'-C3'-C4'-O4'
4	A	902	FAD	O4B-C4B-C5B-O5B
4	A	902	FAD	O3'-C3'-C4'-C5'

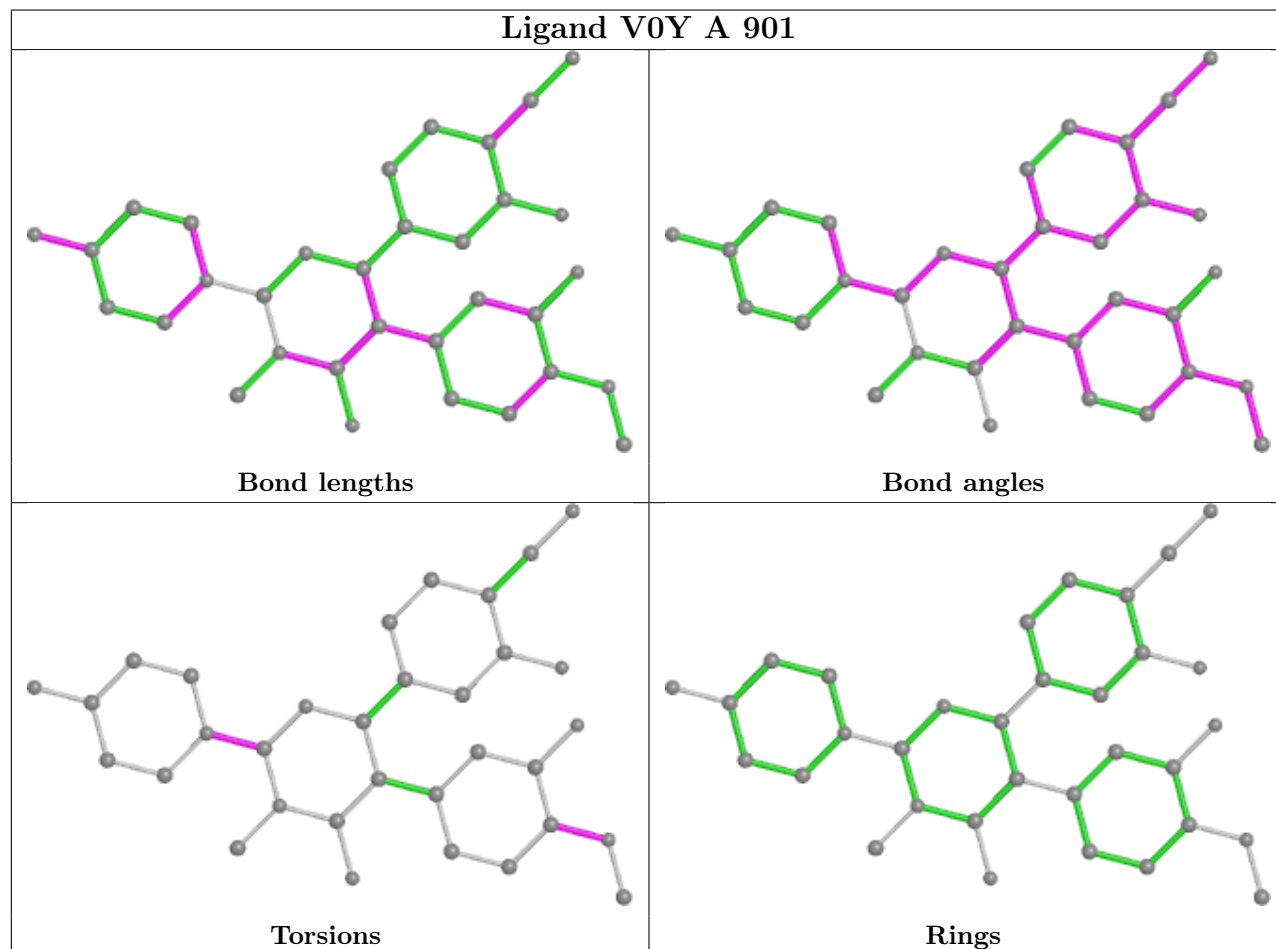
There are no ring outliers.

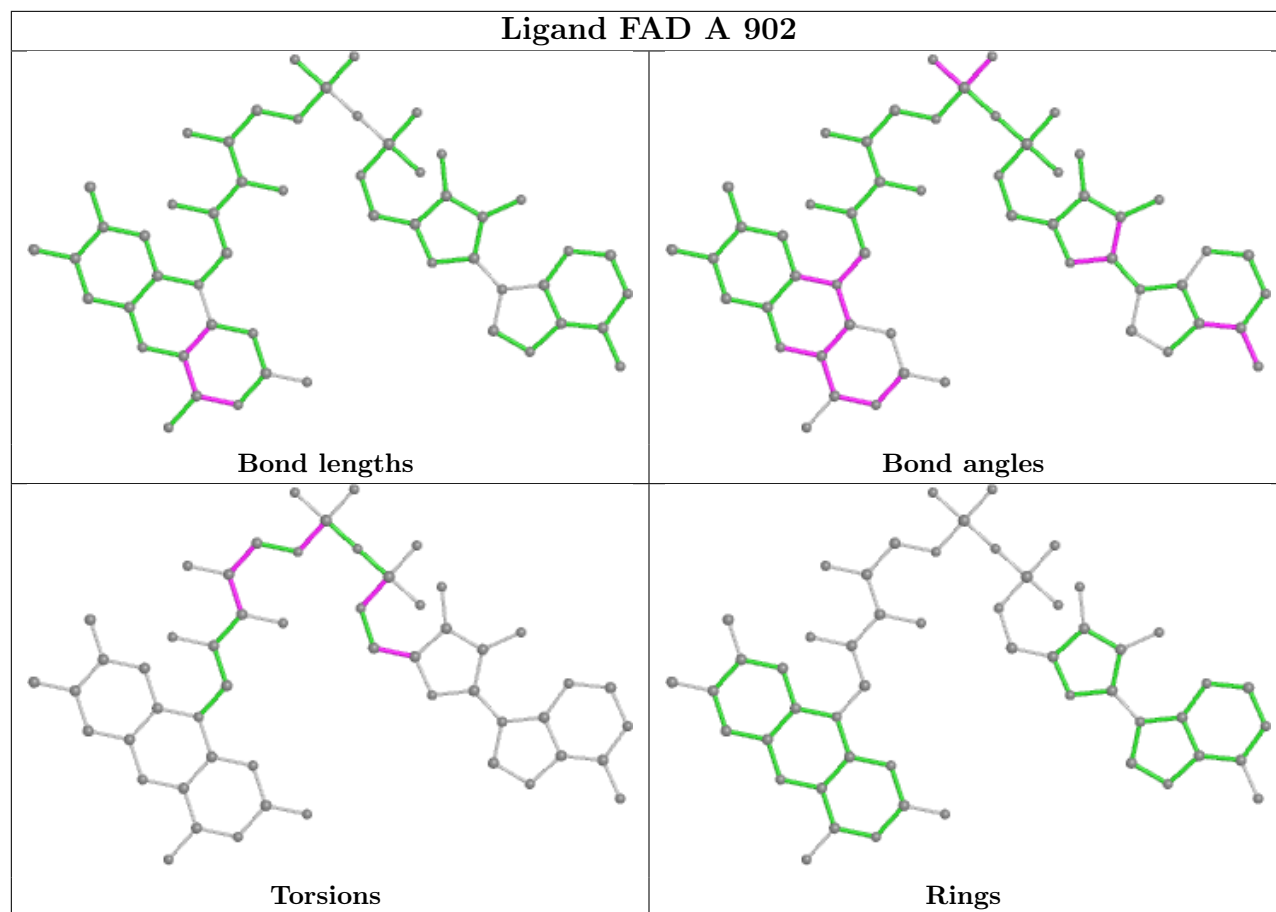
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	V0Y	1	0
4	A	902	FAD	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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	651/659 (98%)	0.15	13 (1%) 65 66	57, 89, 134, 165	0
2	B	125/132 (94%)	0.28	5 (4%) 38 37	88, 120, 142, 179	0
All	All	776/791 (98%)	0.17	18 (2%) 60 61	57, 95, 136, 179	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	174	VAL	3.9
1	A	241	PRO	3.6
1	A	274	PRO	3.3
1	A	175	GLU	3.1
1	A	242	TYR	3.1
2	B	421	PHE	2.7
1	A	399	ASN	2.6
1	A	398	PHE	2.5
2	B	312	LYS	2.5
1	A	238	LEU	2.5
1	A	273	LEU	2.5
2	B	402	PHE	2.4
1	A	492	LYS	2.4
2	B	399	GLY	2.3
1	A	396	LEU	2.2
2	B	437	TRP	2.2
1	A	739	ALA	2.1
1	A	458	LEU	2.1

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 monosaccharides in this entry.

6.4 Ligands [i](#)

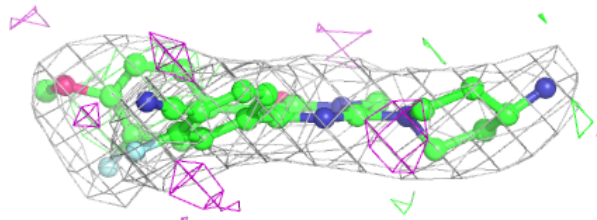
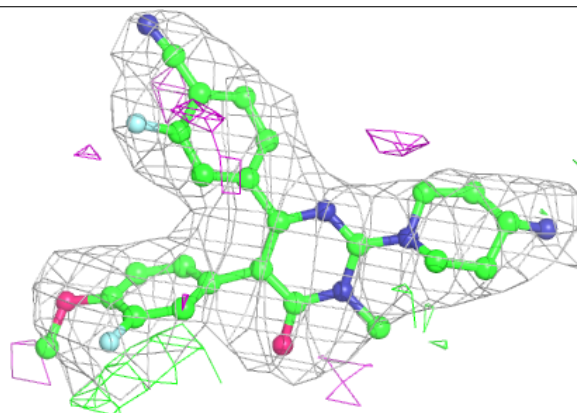
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

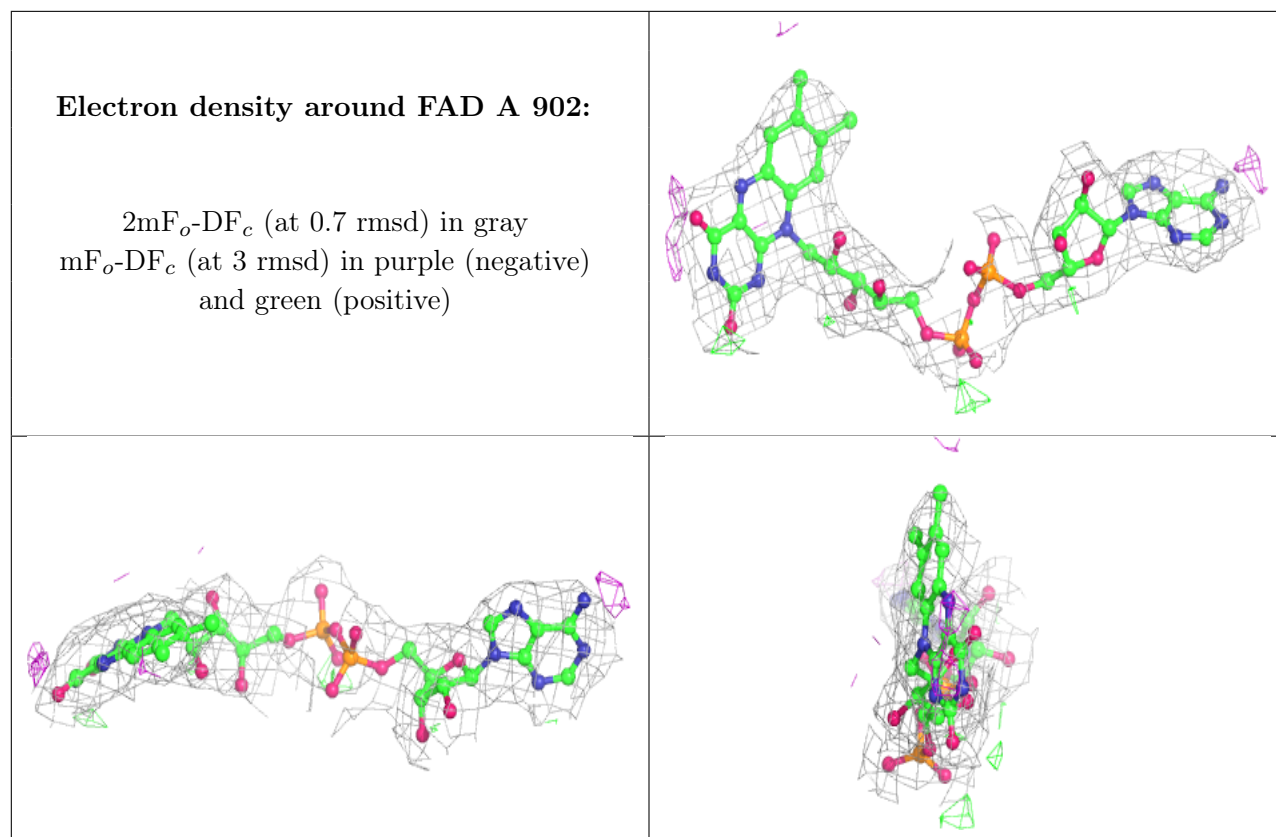
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	V0Y	A	901	33/33	0.96	0.21	51,73,92,106	0
4	FAD	A	902	53/53	0.98	0.22	52,66,84,89	0

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

Electron density around V0Y A 901:

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





6.5 Other polymers [i](#)

There are no such residues in this entry.