



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 25, 2022 – 12:28 PM EDT

PDB ID : 7LY7
Title : Crystal structure of the elongation module of the bacillamide NRPS, BmdB, in complex with the oxidase BmdC
Authors : Fortinez, C.M.; Sharon, I.; Schmeing, T.M.
Deposited on : 2021-03-05
Resolution : 3.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at <http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.29
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) : 2.29

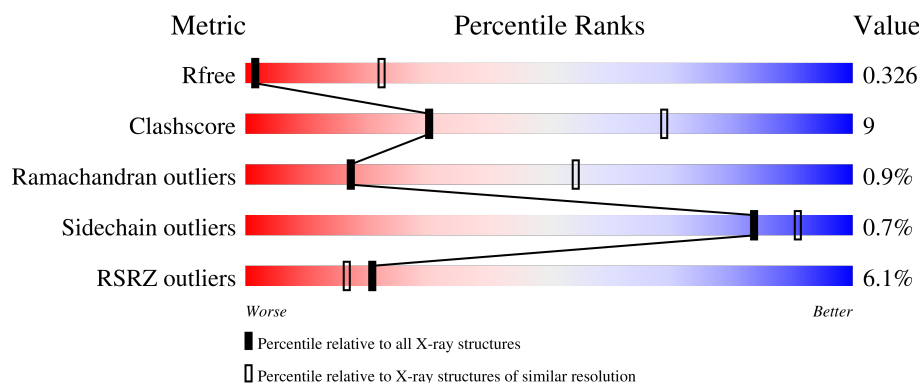
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1212 (4.00-3.60)
Clashscore	141614	1288 (4.00-3.60)
Ramachandran outliers	138981	1243 (4.00-3.60)
Sidechain outliers	138945	1237 (4.00-3.60)
RSRZ outliers	127900	1121 (4.00-3.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	B	335	<div> <div>2%</div> <div>82%</div> <div>15%</div> <div>..</div> </div>
2	A	1027	<div> <div>7%</div> <div>69%</div> <div>18%</div> <div>• 12%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10002 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.

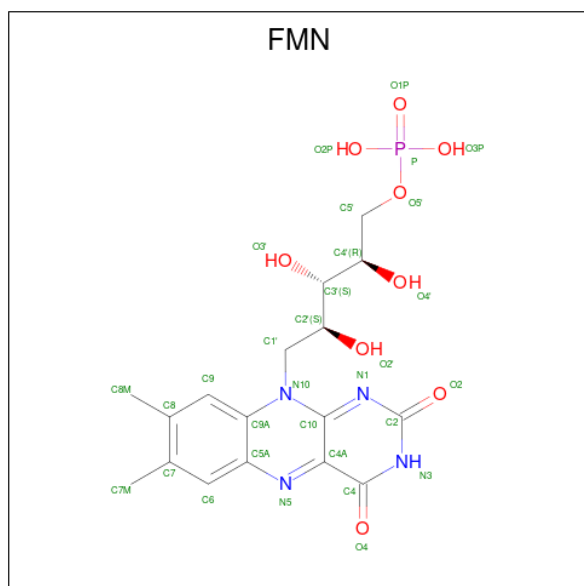
- Molecule 1 is a protein called BmdC, Oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	329	Total	C	N	O	S	0	0	0
			2633	1690	438	495	10			

- Molecule 2 is a protein called BmdB, Bacillamide NRPS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	908	Total	C	N	O	S	0	0	0
			7289	4667	1225	1366	31			

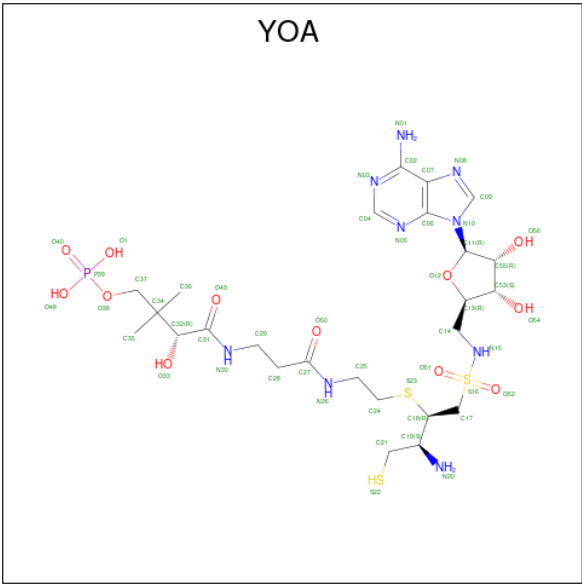
- Molecule 3 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 4 is 5'-{[(2R,3S)-3-amino-2-({2-[(N-{(2R)-4-[(dihydroxyphosphanyl)oxy]-2-hydroxy-3,3-dimethylbutanoyl}-beta-alanyl)amino]ethyl}sulfanyl)-4-sulfanylbutane-1-sulfonyl]a

mino}-5'-deoxyadenosine (three-letter code: YOA) (formula: C₂₅H₄₄N₉O₁₂PS₃) (labeled as "Ligand of Interest" by depositor).

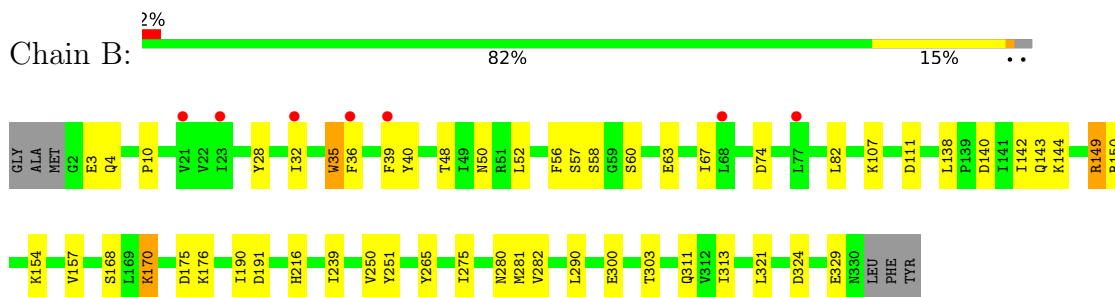


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	0
			49	25	9	11	1	3		

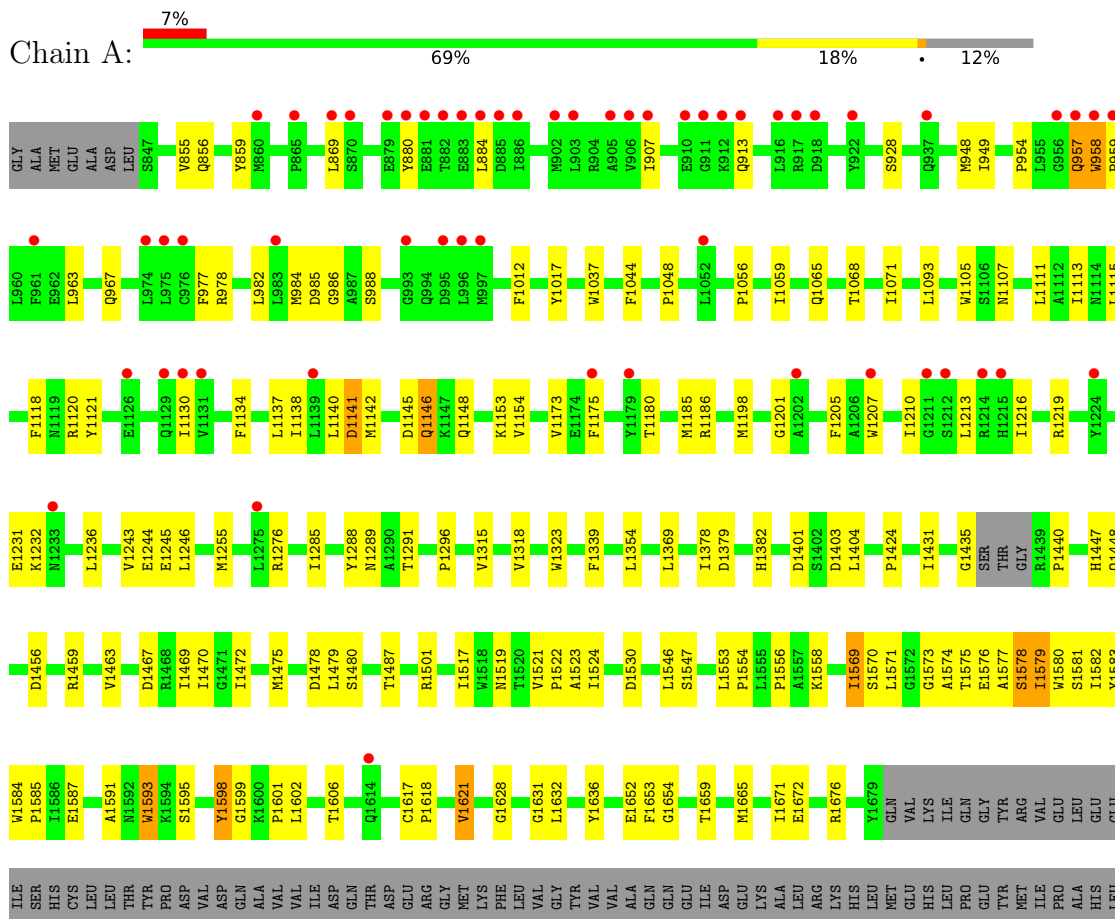
3 Residue-property plots [i](#)

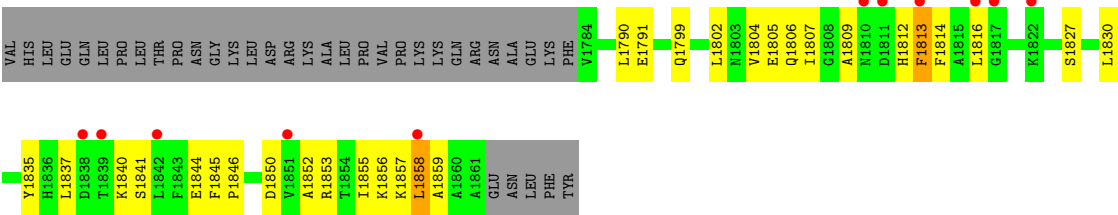
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: BmdC, Oxidase



• Molecule 2: BmdB, Bacillamide NRPS





4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	149.96Å 149.96Å 319.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.28 – 3.80 49.28 – 3.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.28-3.80) 99.9 (49.28-3.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 3.77Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.298 , 0.326 0.298 , 0.326	Depositor DCC
R_{free} test set	2172 reflections (10.01%)	wwPDB-VP
Wilson B-factor (Å ²)	90.3	Xtriage
Anisotropy	0.476	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 87.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.35$, $\langle L^2 \rangle = 0.17$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	10002	wwPDB-VP
Average B, all atoms (Å ²)	182.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% 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: YOA, FMN

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	B	0.30	1/2694 (0.0%)	0.56	1/3648 (0.0%)
2	A	0.26	0/7463	0.46	4/10143 (0.0%)
All	All	0.27	1/10157 (0.0%)	0.49	5/13791 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	35	TRP	C-N	-5.50	1.21	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1141	ASP	CB-CG-OD1	-7.37	111.67	118.30
2	A	1578	SER	N-CA-C	-7.26	91.38	111.00
2	A	1578	SER	CB-CA-C	5.78	121.07	110.10
1	B	149	ARG	NE-CZ-NH2	-5.43	117.59	120.30
2	A	1569	ILE	CG1-CB-CG2	-5.24	99.88	111.40

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	B	2633	0	2572	36	0
2	A	7289	0	7168	144	0
3	B	31	0	19	5	0
4	A	49	0	0	2	0
All	All	10002	0	9759	184	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 9.

All (184) 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)
2:A:958:TRP:HB3	2:A:959:PRO:HD3	1.25	1.15
1:B:300:GLU:O	1:B:303:THR:HG22	1.47	1.14
2:A:958:TRP:HB3	2:A:959:PRO:CD	1.80	1.09
2:A:1617:CYS:HB3	2:A:1621:VAL:HB	1.38	1.01
2:A:1501:ARG:HH12	2:A:1840:LYS:HB3	1.28	0.97
2:A:958:TRP:CB	2:A:959:PRO:CD	2.41	0.95
2:A:1120:ARG:HH22	2:A:1130:ILE:HD11	1.42	0.83
1:B:300:GLU:O	1:B:303:THR:CG2	2.30	0.78
2:A:1501:ARG:NH1	2:A:1840:LYS:HB3	1.97	0.77
2:A:1584:TRP:NE1	2:A:1587:GLU:OE2	2.18	0.75
2:A:958:TRP:CB	2:A:959:PRO:HD2	2.17	0.74
2:A:1059:ILE:HD12	2:A:1245:GLU:HG3	1.70	0.73
2:A:1530:ASP:OD1	2:A:1558:LYS:NZ	2.24	0.70
3:B:401:FMN:H9	3:B:401:FMN:H2'	1.74	0.70
2:A:954:PRO:HB2	2:A:957:GLN:HB2	1.73	0.69
2:A:1115:LEU:HB2	2:A:1140:LEU:HD11	1.74	0.69
2:A:1469:ILE:HG13	2:A:1517:ILE:HB	1.75	0.69
2:A:1571:LEU:HD11	2:A:1584:TRP:H	1.57	0.68
1:B:149:ARG:HD3	3:B:401:FMN:O2	1.94	0.68
2:A:1201:GLY:H	2:A:1231:GLU:HB2	1.59	0.68
1:B:60:SER:O	1:B:63:GLU:HG2	1.94	0.67
2:A:958:TRP:HB2	2:A:959:PRO:HD2	1.75	0.66
1:B:60:SER:HB3	1:B:63:GLU:HB3	1.78	0.66
2:A:1804:VAL:HG11	2:A:1816:LEU:HD22	1.78	0.65
2:A:1570:SER:O	2:A:1571:LEU:HD12	1.98	0.63
2:A:1575:THR:HG23	4:A:1901:YOA:O52	1.98	0.63
2:A:1856:LYS:HG2	2:A:1857:LYS:H	1.64	0.62
2:A:1569:ILE:HG23	2:A:1585:PRO:HA	1.81	0.61
1:B:10:PRO:HD2	1:B:82:LEU:HD11	1.81	0.61
2:A:1569:ILE:HG22	2:A:1571:LEU:CD1	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1618:PRO:O	2:A:1621:VAL:HG23	2.02	0.60
2:A:1478:ASP:HB3	2:A:1579:ILE:HG12	1.83	0.60
2:A:1017:TYR:HB2	2:A:1120:ARG:HH21	1.65	0.59
2:A:1571:LEU:CD1	2:A:1584:TRP:H	2.15	0.59
2:A:928:SER:HA	2:A:967:GLN:HB3	1.85	0.58
1:B:3:GLU:OE2	1:B:50:ASN:ND2	2.27	0.58
2:A:1579:ILE:O	4:A:1901:YOA:S22	2.61	0.58
1:B:149:ARG:HD2	1:B:239:ILE:HD11	1.84	0.58
2:A:1790:LEU:HD13	2:A:1856:LYS:HD2	1.85	0.58
2:A:1037:TRP:HB3	2:A:1175:PHE:HD2	1.69	0.57
2:A:1573:GLY:HA3	2:A:1580:TRP:HA	1.86	0.57
2:A:869:LEU:HD22	2:A:1243:VAL:HG11	1.85	0.57
1:B:32:ILE:HB	1:B:67:ILE:HG21	1.87	0.57
2:A:985:ASP:OD2	2:A:1134:PHE:HB2	2.05	0.56
2:A:1584:TRP:HB2	2:A:1599:GLY:HA3	1.86	0.56
2:A:1288:TYR:CZ	2:A:1601:PRO:HD3	2.40	0.56
1:B:154:LYS:HD3	1:B:329:GLU:HA	1.87	0.56
2:A:1665:MET:HG2	2:A:1671:ILE:HG12	1.85	0.56
2:A:1463:VAL:HA	2:A:1467:ASP:OD2	2.05	0.56
1:B:300:GLU:C	1:B:303:THR:HG22	2.23	0.56
3:B:401:FMN:H9	3:B:401:FMN:C2'	2.34	0.56
2:A:1577:ALA:O	2:A:1579:ILE:N	2.38	0.56
1:B:28:TYR:OH	1:B:74:ASP:OD2	2.23	0.55
2:A:948:MET:HB3	2:A:978:ARG:HD3	1.87	0.55
2:A:1804:VAL:HG21	2:A:1816:LEU:HD13	1.89	0.54
2:A:1146:GLN:CD	2:A:1148:GLN:HB2	2.28	0.54
2:A:1582:ILE:HG13	2:A:1598:TYR:CD2	2.43	0.54
2:A:1068:THR:HG21	2:A:1216:ILE:HB	1.89	0.54
2:A:1093:LEU:HD23	2:A:1154:VAL:HG21	1.90	0.53
2:A:1285:ILE:HA	2:A:1288:TYR:HB3	1.90	0.53
2:A:949:ILE:O	2:A:1219:ARG:NH1	2.37	0.53
1:B:250:VAL:HG22	1:B:313:ILE:HG12	1.90	0.53
2:A:869:LEU:HD11	2:A:1246:LEU:HD11	1.91	0.53
2:A:1469:ILE:HD13	2:A:1487:THR:HG23	1.89	0.53
2:A:1577:ALA:O	2:A:1578:SER:C	2.47	0.53
2:A:1570:SER:C	2:A:1571:LEU:HD12	2.30	0.53
1:B:154:LYS:NZ	1:B:329:GLU:HG2	2.24	0.53
2:A:1805:GLU:N	2:A:1805:GLU:OE2	2.42	0.53
2:A:1807:ILE:HG21	2:A:1816:LEU:HD11	1.90	0.52
2:A:1137:LEU:HD12	2:A:1173:VAL:HB	1.92	0.52
2:A:1276:ARG:NH2	2:A:1652:GLU:OE2	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1289:ASN:HB3	2:A:1606:THR:HG22	1.92	0.52
2:A:1591:ALA:HB1	2:A:1593:TRP:CZ2	2.45	0.52
2:A:907:ILE:HG21	2:A:982:LEU:HD13	1.91	0.51
2:A:1113:ILE:HB	2:A:1140:LEU:HD13	1.92	0.51
2:A:1137:LEU:O	2:A:1138:ILE:HG13	2.11	0.51
1:B:36:PHE:HD2	1:B:40:TYR:CE2	2.28	0.51
2:A:1318:VAL:HG22	2:A:1323:TRP:CG	2.46	0.51
2:A:1628:GLY:HA2	2:A:1632:LEU:HD11	1.93	0.51
2:A:1180:THR:HG23	2:A:1185:MET:HB2	1.93	0.51
1:B:157:VAL:HG22	1:B:321:LEU:HG	1.93	0.51
1:B:280:ASN:HA	1:B:290:LEU:HD12	1.92	0.51
2:A:1571:LEU:HG	2:A:1582:ILE:O	2.12	0.50
3:B:401:FMN:O1P	3:B:401:FMN:H4'	2.12	0.50
2:A:1812:HIS:ND1	2:A:1846:PRO:HB3	2.27	0.50
2:A:1578:SER:O	2:A:1580:TRP:N	2.45	0.50
2:A:1844:GLU:OE1	2:A:1844:GLU:N	2.33	0.50
2:A:855:VAL:O	2:A:859:TYR:N	2.41	0.50
2:A:1044:PHE:CE2	2:A:1141:ASP:OD1	2.65	0.50
1:B:175:ASP:O	1:B:176:LYS:HG3	2.12	0.49
2:A:1205:PHE:CE2	2:A:1207:TRP:HB3	2.47	0.49
2:A:1105:TRP:CD1	2:A:1255:MET:HG2	2.48	0.49
2:A:1205:PHE:CD2	2:A:1232:LYS:HG3	2.47	0.49
2:A:880:TYR:CZ	2:A:1213:LEU:HD13	2.48	0.49
2:A:1118:PHE:CE2	2:A:1198:MET:HG2	2.48	0.49
2:A:1582:ILE:HG13	2:A:1598:TYR:CE2	2.48	0.48
2:A:1571:LEU:HG	2:A:1583:TYR:HA	1.94	0.48
2:A:1659:THR:O	2:A:1676:ARG:NH1	2.34	0.48
2:A:1805:GLU:HG2	2:A:1806:GLN:H	1.78	0.48
2:A:1456:ASP:HA	2:A:1459:ARG:HG2	1.95	0.48
2:A:856:GLN:OE1	2:A:913:GLN:NE2	2.47	0.48
2:A:1837:LEU:HD21	2:A:1858:LEU:HB2	1.95	0.48
2:A:1521:VAL:HG22	2:A:1523:ALA:H	1.79	0.48
1:B:39:PHE:CE2	1:B:52:LEU:HD22	2.49	0.47
3:B:401:FMN:C2'	3:B:401:FMN:C9	2.91	0.47
2:A:1276:ARG:NH1	2:A:1652:GLU:OE2	2.45	0.47
2:A:1522:PRO:HA	2:A:1546:LEU:HD22	1.95	0.47
2:A:1012:PHE:O	2:A:1130:ILE:HG21	2.14	0.47
2:A:1569:ILE:HG22	2:A:1571:LEU:HD13	1.95	0.47
1:B:35:TRP:CE2	1:B:56:PHE:HB3	2.50	0.47
2:A:1424:PRO:HB2	2:A:1448:GLN:HB2	1.96	0.47
1:B:138:LEU:HD22	1:B:142:ILE:HD11	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:ILE:HD12	1:B:275:ILE:HD12	1.97	0.46
2:A:1653:PHE:CG	2:A:1654:GLY:N	2.82	0.46
2:A:963:LEU:HG	2:A:977:PHE:HD1	1.81	0.46
2:A:1017:TYR:HH	2:A:1121:TYR:HD2	1.63	0.46
2:A:1835:TYR:HB3	2:A:1855:ILE:HG13	1.96	0.46
2:A:1841:SER:HB2	2:A:1845:PHE:CE2	2.51	0.46
2:A:1037:TRP:HB3	2:A:1175:PHE:CD2	2.50	0.46
2:A:1120:ARG:HD2	2:A:1134:PHE:CD1	2.51	0.46
1:B:170:LYS:HD2	1:B:216:HIS:HD2	1.81	0.46
2:A:1056:PRO:HA	2:A:1059:ILE:HG12	1.97	0.46
2:A:1339:PHE:CE2	2:A:1404:LEU:HD11	2.50	0.46
2:A:1521:VAL:HA	2:A:1547:SER:O	2.16	0.46
1:B:140:ASP:HB3	1:B:144:LYS:HD2	1.97	0.45
2:A:1120:ARG:HD2	2:A:1134:PHE:CE1	2.50	0.45
2:A:1802:LEU:HD13	2:A:1816:LEU:HB2	1.98	0.45
2:A:1379:ASP:HB3	2:A:1382:HIS:CG	2.51	0.45
2:A:1569:ILE:HG23	2:A:1569:ILE:HD12	1.64	0.45
2:A:1065:GLN:HG3	2:A:1244:GLU:OE1	2.16	0.45
2:A:1583:TYR:HB3	2:A:1602:LEU:HA	1.97	0.45
2:A:1354:LEU:HD23	2:A:1378:ILE:HB	1.98	0.45
2:A:985:ASP:OD1	2:A:986:GLY:N	2.47	0.45
2:A:1553:LEU:O	2:A:1556:PRO:HD2	2.17	0.45
2:A:1369:LEU:HD22	2:A:1447:HIS:CE1	2.52	0.44
2:A:1581:SER:O	2:A:1602:LEU:N	2.35	0.44
2:A:984:MET:HB2	2:A:988:SER:HB2	2.00	0.44
2:A:1435:GLY:HA2	2:A:1440:PRO:HB3	2.00	0.44
2:A:1835:TYR:CE1	2:A:1856:LYS:HG3	2.53	0.43
1:B:36:PHE:CD2	1:B:40:TYR:CE2	3.07	0.43
2:A:1315:VAL:HG11	2:A:1318:VAL:CG2	2.49	0.43
1:B:36:PHE:HD2	1:B:40:TYR:CZ	2.36	0.43
2:A:1791:GLU:OE2	2:A:1809:ALA:HB2	2.19	0.43
1:B:107:LYS:HZ3	1:B:111:ASP:CG	2.22	0.43
2:A:1401:ASP:HB3	2:A:1403:ASP:OD1	2.19	0.43
2:A:1845:PHE:HD1	2:A:1850:ASP:OD2	2.02	0.43
1:B:251:TYR:O	1:B:311:GLN:HA	2.18	0.43
2:A:884:LEU:HD12	2:A:1210:ILE:HG23	2.01	0.43
2:A:1071:ILE:HD12	2:A:1236:LEU:HD23	2.00	0.43
1:B:4:GLN:O	1:B:48:THR:HG22	2.19	0.42
1:B:281:MET:HE2	1:B:281:MET:HB2	1.83	0.42
2:A:1576:GLU:OE2	2:A:1636:TYR:OH	2.23	0.42
2:A:1291:THR:HG21	2:A:1601:PRO:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1501:ARG:NH1	2:A:1840:LYS:HD2	2.34	0.42
2:A:1857:LYS:O	2:A:1859:ALA:N	2.52	0.42
1:B:35:TRP:NE1	1:B:56:PHE:HB3	2.34	0.42
2:A:1813:PHE:CD1	2:A:1813:PHE:O	2.73	0.42
2:A:1431:ILE:HG21	2:A:1478:ASP:OD1	2.20	0.42
1:B:36:PHE:CD2	1:B:40:TYR:CZ	3.07	0.42
2:A:1318:VAL:HG22	2:A:1323:TRP:CB	2.50	0.42
2:A:1107:ASN:HB2	2:A:1621:VAL:CG1	2.50	0.42
2:A:1318:VAL:HG13	2:A:1323:TRP:CE3	2.55	0.42
2:A:1479:LEU:HD11	2:A:1519:ASN:OD1	2.20	0.41
2:A:1802:LEU:HD12	2:A:1816:LEU:HD12	2.02	0.41
1:B:168:SER:HB3	1:B:282:VAL:HG21	2.02	0.41
2:A:1048:PRO:HD3	2:A:1111:LEU:HB2	2.00	0.41
2:A:1553:LEU:HB2	2:A:1554:PRO:HD3	2.02	0.41
2:A:1578:SER:N	2:A:1631:GLY:O	2.52	0.41
2:A:1595:SER:OG	2:A:1672:GLU:OE2	2.27	0.41
1:B:57:SER:OG	1:B:58:SER:N	2.53	0.41
2:A:1017:TYR:CB	2:A:1120:ARG:HH21	2.30	0.41
2:A:1146:GLN:CD	2:A:1153:LYS:HG3	2.41	0.41
2:A:1827:SER:HA	2:A:1830:LEU:HD12	2.02	0.41
2:A:1569:ILE:HG22	2:A:1571:LEU:HD11	2.03	0.41
2:A:1790:LEU:HG	2:A:1852:ALA:HB1	2.02	0.41
2:A:1475:MET:HA	2:A:1480:SER:OG	2.21	0.40
2:A:1850:ASP:HA	2:A:1853:ARG:HD3	2.03	0.40
1:B:142:ILE:HG13	1:B:143:GLN:N	2.36	0.40
2:A:954:PRO:HG2	2:A:957:GLN:HG2	2.04	0.40
2:A:1111:LEU:O	2:A:1142:MET:HB2	2.20	0.40
2:A:1521:VAL:HG13	2:A:1524:ILE:HG12	2.02	0.40
1:B:191:ASP:OD2	1:B:311:GLN:NE2	2.54	0.40
2:A:1470:ILE:HG12	2:A:1472:ILE:HG13	2.04	0.40
2:A:1799:GLN:NE2	2:A:1805:GLU:O	2.54	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	B	327/335 (98%)	309 (94%)	17 (5%)	1 (0%)	41	74
2	A	902/1027 (88%)	834 (92%)	58 (6%)	10 (1%)	14	51
All	All	1229/1362 (90%)	1143 (93%)	75 (6%)	11 (1%)	17	54

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	958	TRP
2	A	1579	ILE
2	A	1186	ARG
2	A	1813	PHE
2	A	1858	LEU
1	B	324	ASP
2	A	1574	ALA
2	A	957	GLN
2	A	1621	VAL
2	A	1145	ASP
2	A	1296	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	B	281/286 (98%)	278 (99%)	3 (1%)	73	85
2	A	788/896 (88%)	784 (100%)	4 (0%)	88	94
All	All	1069/1182 (90%)	1062 (99%)	7 (1%)	84	91

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	150	ARG
1	B	170	LYS

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Mol	Chain	Res	Type
1	B	265	TYR
2	A	1146	GLN
2	A	1593	TRP
2	A	1598	TYR
2	A	1814	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	1156	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	FMN	B	401	-	33,33,33	0.56	0	48,50,50	0.64	1 (2%)
4	YOA	A	1901	2	40,51,52	1.12	4 (10%)	43,73,76	2.17	7 (16%)

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	FMN	B	401	-	-	11/18/18/18	0/3/3/3
4	YOA	A	1901	2	-	22/40/66/67	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1901	YOA	C24-S23	-3.82	1.76	1.81
4	A	1901	YOA	C18-S23	-3.25	1.77	1.82
4	A	1901	YOA	C17-S16	-3.18	1.68	1.77
4	A	1901	YOA	O12-C11	2.03	1.43	1.41

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1901	YOA	O52-S16-O51	-8.57	106.93	119.35
4	A	1901	YOA	C14-N15-S16	-7.96	104.75	120.63
4	A	1901	YOA	N05-C04-N03	-4.44	121.74	128.68
4	A	1901	YOA	O52-S16-C17	3.41	113.26	107.87
4	A	1901	YOA	C29-C28-C27	-2.26	108.59	112.36
4	A	1901	YOA	C35-C34-C32	2.21	112.65	108.82
4	A	1901	YOA	C53-C55-C11	2.06	104.08	100.98
3	B	401	FMN	C4-N3-C2	-2.03	121.89	125.64

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	401	FMN	C2'-C1'-N10-C10
3	B	401	FMN	N10-C1'-C2'-O2'
3	B	401	FMN	N10-C1'-C2'-C3'
3	B	401	FMN	C1'-C2'-C3'-C4'
3	B	401	FMN	C2'-C3'-C4'-O4'
3	B	401	FMN	C2'-C3'-C4'-C5'
3	B	401	FMN	O3'-C3'-C4'-O4'
3	B	401	FMN	O3'-C3'-C4'-C5'
3	B	401	FMN	C4'-C5'-O5'-P
4	A	1901	YOA	C53-C13-C14-N15

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Mol	Chain	Res	Type	Atoms
4	A	1901	YOA	O12-C13-C14-N15
4	A	1901	YOA	S16-C17-C18-C19
4	A	1901	YOA	C17-C18-C19-C21
4	A	1901	YOA	C17-C18-C19-N20
4	A	1901	YOA	S23-C18-C19-C21
4	A	1901	YOA	C17-C18-S23-C24
4	A	1901	YOA	S23-C24-C25-N26
4	A	1901	YOA	C25-C24-S23-C18
4	A	1901	YOA	C27-C28-C29-N30
4	A	1901	YOA	C31-C32-C34-C35
4	A	1901	YOA	C31-C32-C34-C36
4	A	1901	YOA	C31-C32-C34-C37
4	A	1901	YOA	O33-C32-C34-C35
4	A	1901	YOA	O33-C32-C34-C36
4	A	1901	YOA	O33-C32-C34-C37
4	A	1901	YOA	C32-C34-C37-O38
4	A	1901	YOA	C14-N15-S16-C17
4	A	1901	YOA	C14-N15-S16-O51
4	A	1901	YOA	C35-C34-C37-O38
4	A	1901	YOA	C36-C34-C37-O38
3	B	401	FMN	O2'-C2'-C3'-C4'
3	B	401	FMN	O2'-C2'-C3'-O3'
4	A	1901	YOA	O50-C27-C28-C29

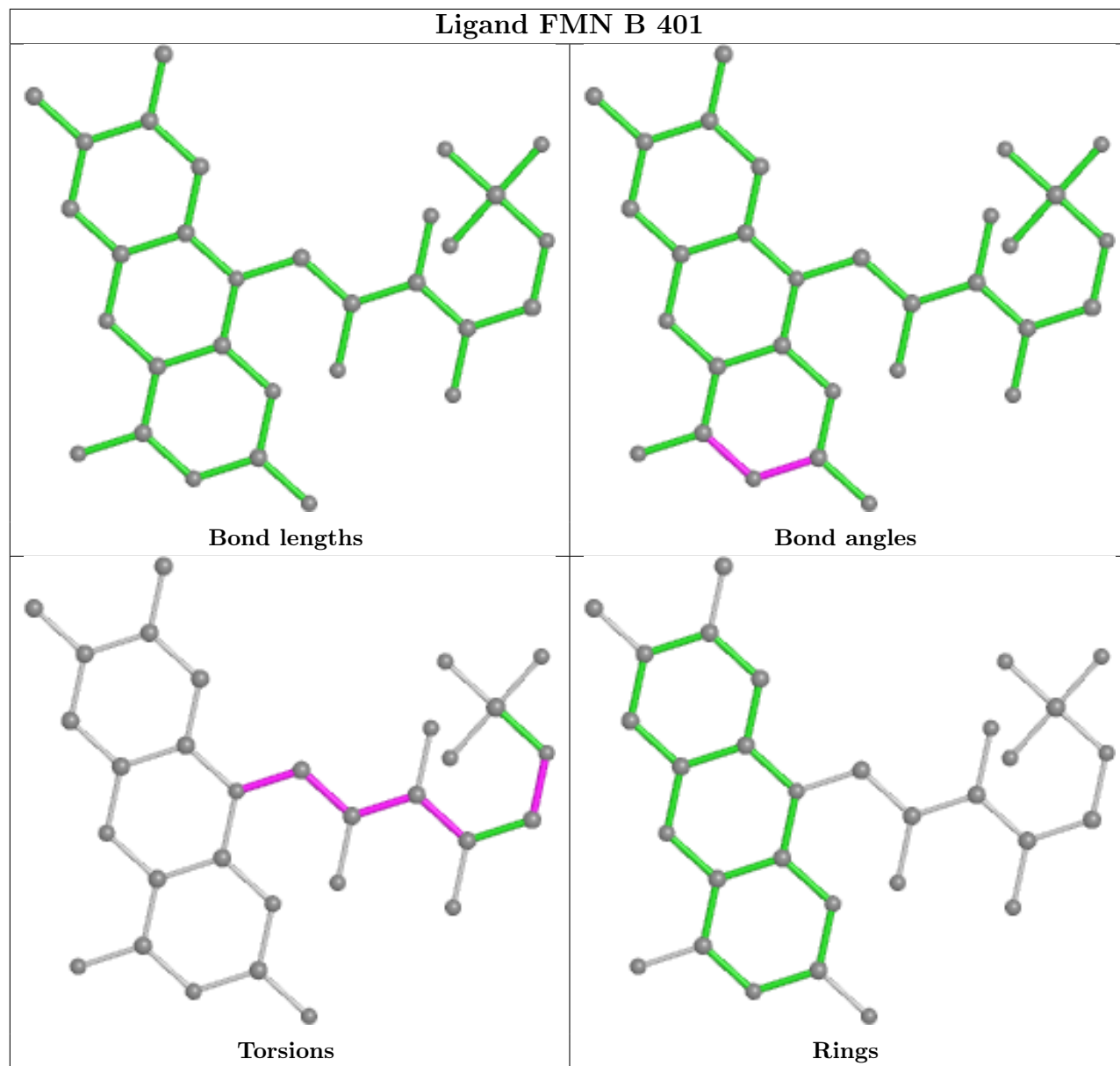
There are no ring outliers.

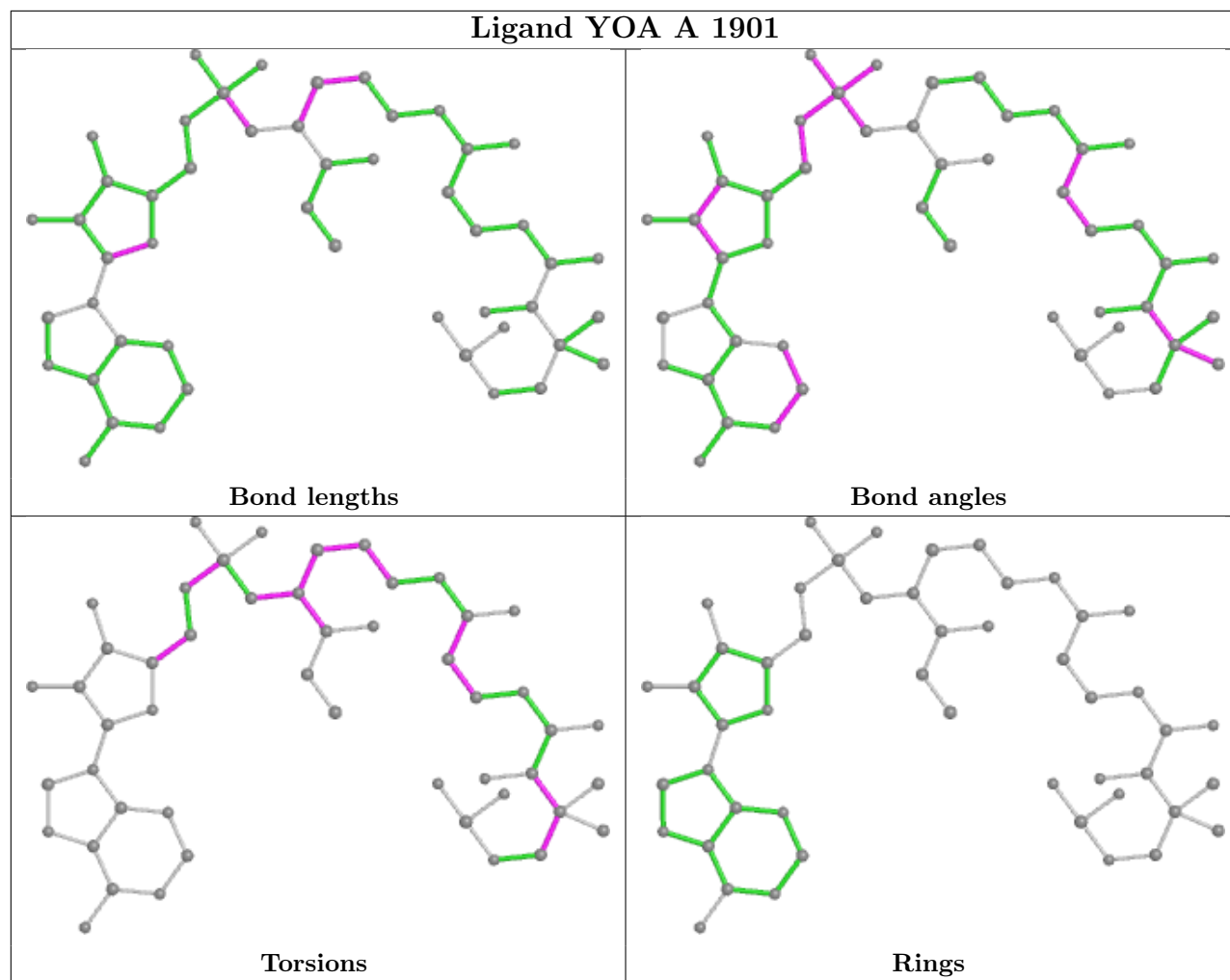
2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	401	FMN	5	0
4	A	1901	YOA	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	B	329/335 (98%)	-0.06	7 (2%) 63 55	50, 105, 257, 297	0
2	A	908/1027 (88%)	0.18	68 (7%) 14 11	71, 187, 365, 460	0
All	All	1237/1362 (90%)	0.12	75 (6%) 21 17	50, 164, 359, 460	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	906	VAL	6.0
2	A	1842	LEU	5.6
1	B	21	VAL	5.4
2	A	880	TYR	5.2
2	A	885	ASP	5.2
2	A	907	ILE	4.8
1	B	39	PHE	4.3
2	A	957	GLN	4.3
2	A	995	ASP	4.2
2	A	911	GLY	4.2
2	A	886	ILE	4.2
2	A	1816	LEU	4.2
2	A	903	LEU	4.0
2	A	1212	SER	4.0
2	A	1207	TRP	4.0
2	A	1839	THR	3.9
2	A	1214	ARG	3.8
2	A	1810	ASN	3.8
2	A	882	THR	3.8
2	A	983	LEU	3.7
2	A	996	LEU	3.7
2	A	997	MET	3.7
2	A	956	GLY	3.6
2	A	881	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
2	A	860	MET	3.4
2	A	1811	ASP	3.3
2	A	902	MET	3.2
2	A	1131	VAL	3.2
2	A	1211	GLY	3.1
2	A	1275	LEU	3.1
1	B	36	PHE	3.1
2	A	884	LEU	3.1
2	A	993	GLY	3.0
2	A	905	ALA	3.0
2	A	922	TYR	3.0
2	A	879	GLU	3.0
2	A	1614	GLN	2.9
2	A	1233	ASN	2.9
2	A	976	CYS	2.9
2	A	917	ARG	2.8
2	A	975	LEU	2.8
2	A	918	ASP	2.7
1	B	23	ILE	2.7
2	A	913	GLN	2.7
2	A	912	LYS	2.7
2	A	910	GLU	2.6
2	A	1202	ALA	2.6
2	A	961	PHE	2.6
2	A	1851	VAL	2.6
2	A	958	TRP	2.6
2	A	1175	PHE	2.5
2	A	869	LEU	2.5
2	A	1858	LEU	2.4
2	A	1215	HIS	2.4
2	A	1179	TYR	2.4
2	A	883	GLU	2.4
1	B	68	LEU	2.3
2	A	1224	TYR	2.3
2	A	1139	LEU	2.3
2	A	1130	ILE	2.3
2	A	1129	GLN	2.3
2	A	1822	LYS	2.3
2	A	1817	GLY	2.3
2	A	959	PRO	2.3
2	A	865	PRO	2.3
2	A	1126	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
2	A	937	GLN	2.2
2	A	1838	ASP	2.2
1	B	77	LEU	2.2
1	B	32	ILE	2.1
2	A	916	LEU	2.1
2	A	974	LEU	2.1
2	A	1813	PHE	2.1
2	A	870	SER	2.1
2	A	1052	LEU	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 monosaccharides in this entry.

6.4 Ligands [i](#)

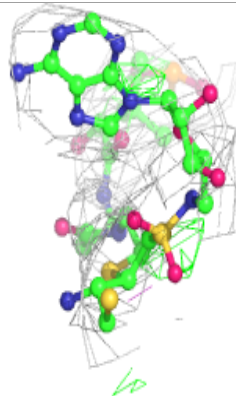
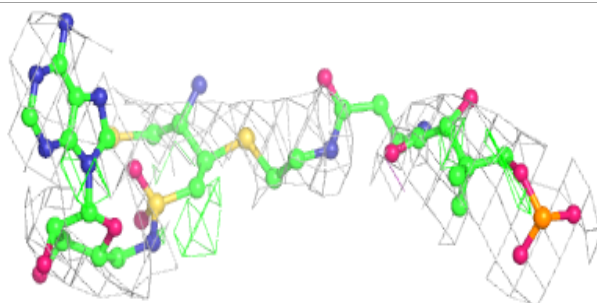
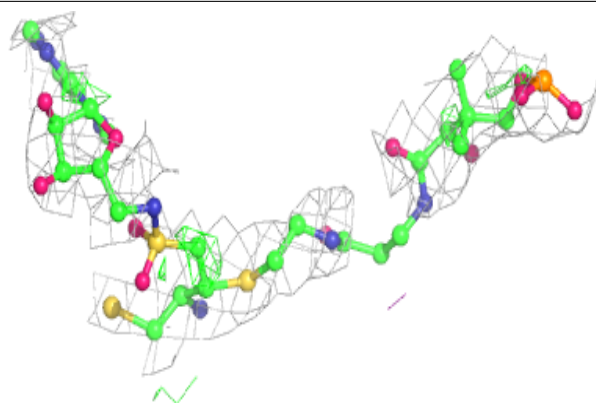
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	YOA	A	1901	49/50	0.84	0.34	203,216,231,234	0
3	FMN	B	401	31/31	0.89	0.31	48,65,93,105	0

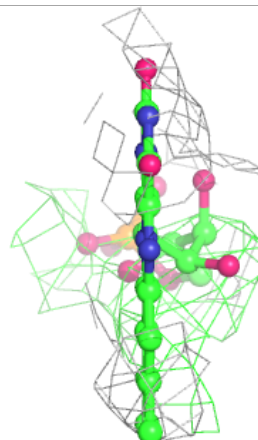
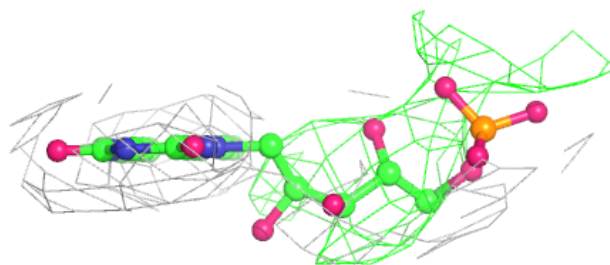
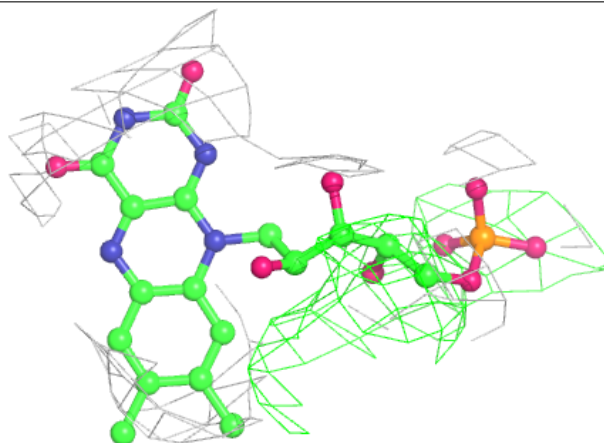
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around YOA A 1901:

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

**Electron density around FMN B 401:**

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



6.5 Other polymers ⓘ

There are no such residues in this entry.