



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:07 PM UTC

PDB ID : 7LY7 / pdb_00007ly7
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	:	4-5-2 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

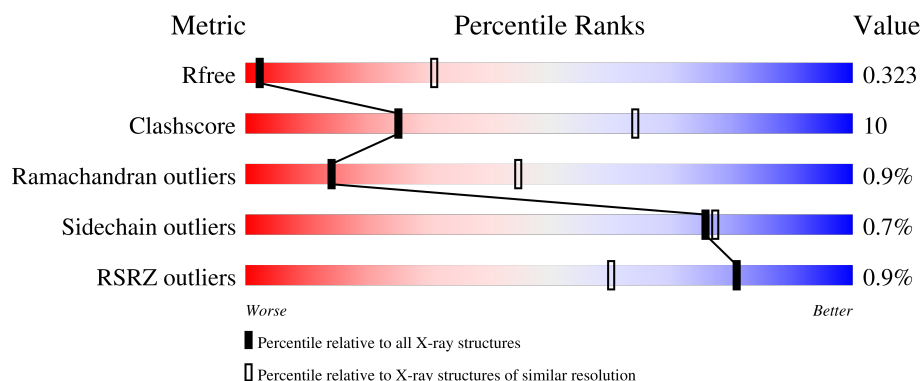
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

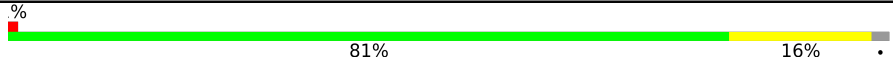
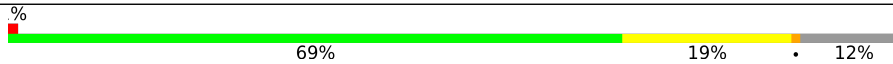
The reported resolution of this entry is 3.80 Å.

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}	180053	1065 (3.96-3.64)
Clashscore	190562	1012 (3.94-3.66)
Ramachandran outliers	187476	1048 (3.96-3.64)
Sidechain outliers	187428	1043 (3.96-3.64)
RSRZ outliers	180081	1064 (3.96-3.64)

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	
2	A	1027	

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.

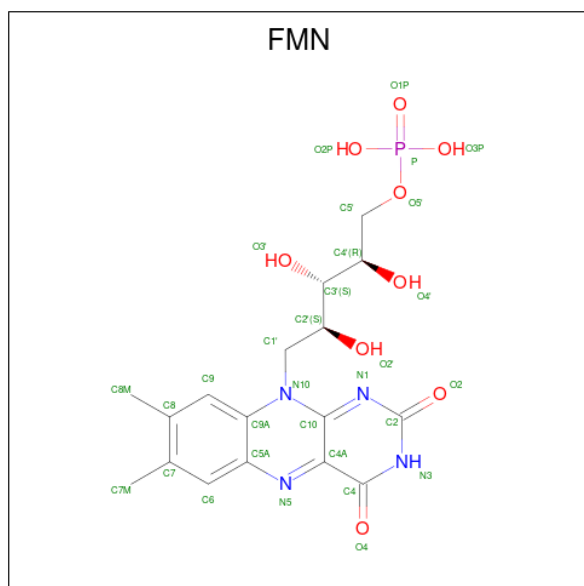
- Molecule 1 is a protein called BmdC, Oxidase.

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

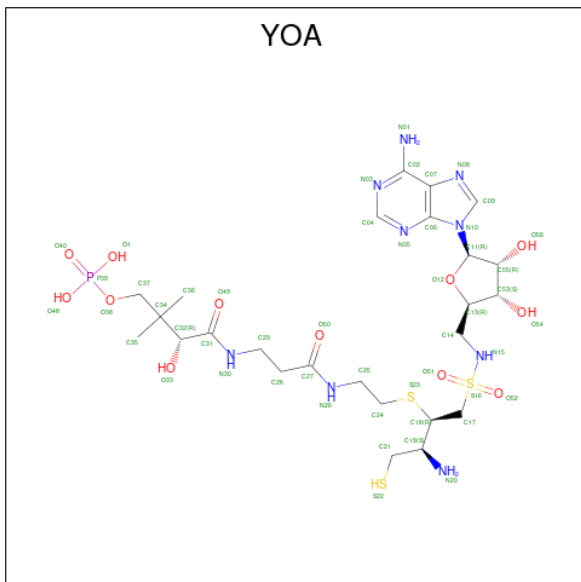
- Molecule 2 is a protein called BmdB, Bacillamide NRPS.

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 (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



no}-5'-deoxyadenosine (CCD ID: YOA) (formula: $C_{25}H_{44}N_9O_{12}PS_3$) (labeled as "Ligand of Interest" by depositor).

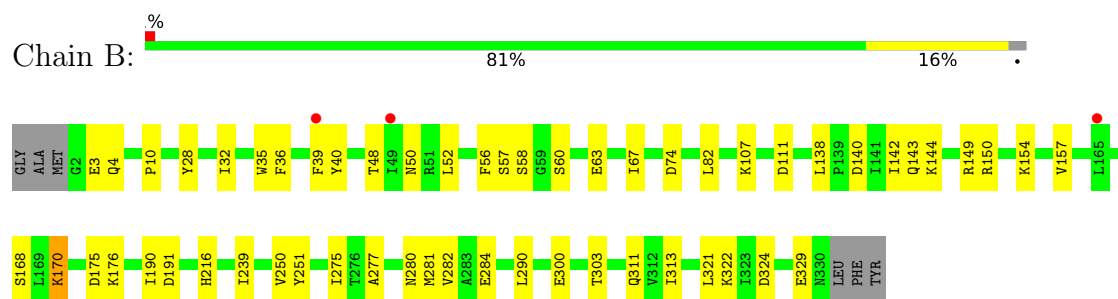


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	0
			49	25	9	11	1	3		

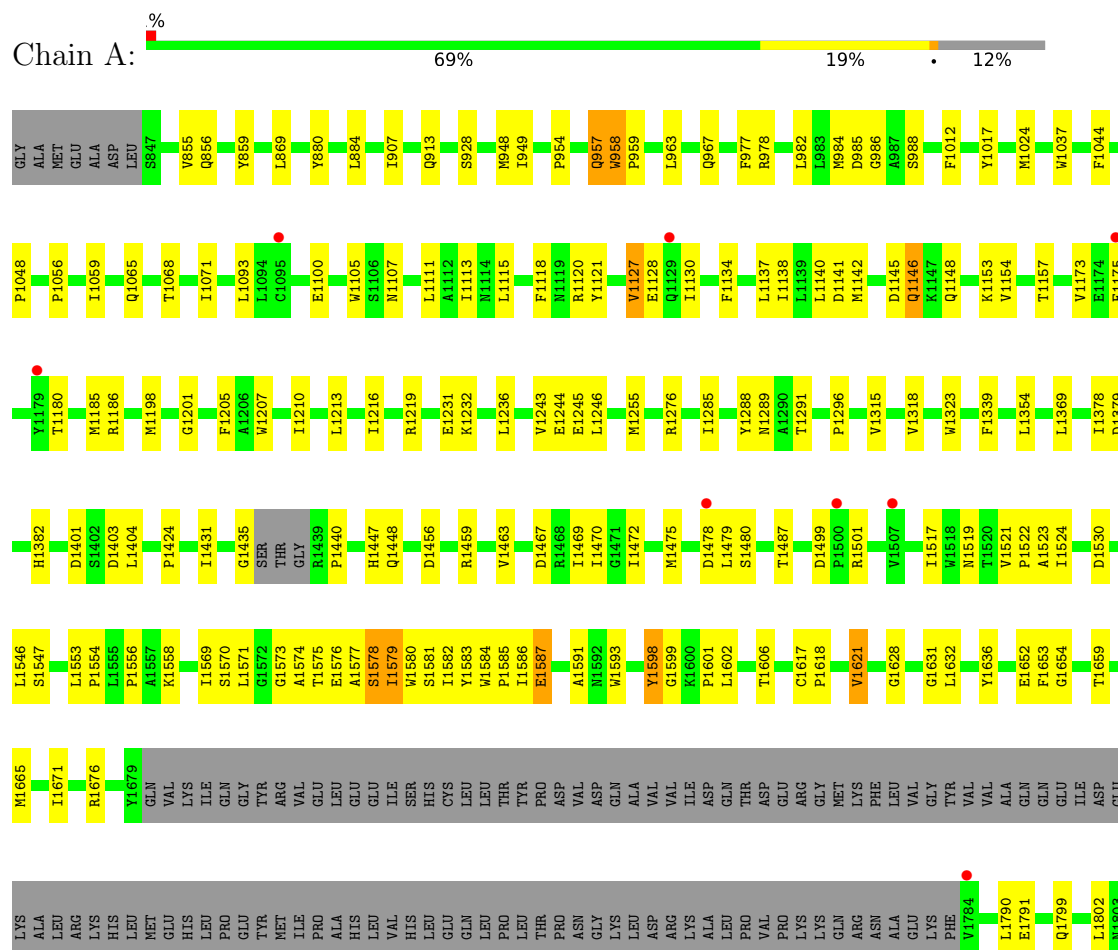
3 Residue-property plots [i](#)

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



Y1804	E1805	Q1806	I1807	G1808	A1809		H1812	F1813	F1814	A1815	L1816		S1827		L1830		Y1835	H1836	L1837		K1840	S1841		E1844	F1845	P1846		D1850	V1851	A1852	R1853	T1854	I1855	K1856	K1857	L1858	A1859	A1860	A1861	GLU	ASN	LEU	PHE	TYR
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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) 99.8 (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 , 0.323	Depositor DCC
R_{free} test set	2174 reflections (10.00%)	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)	0.28 , 111.2	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	0.86	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: FMN, YOA

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	0.18	0/2694	0.54	0/3648
2	A	0.16	0/7463	0.38	3/10143 (0.0%)
All	All	0.17	0/10157	0.43	3/13791 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1578	SER	N-CA-C	-9.12	91.38	110.80
2	A	1578	SER	CB-CA-C	5.35	121.07	110.42
2	A	1586	ILE	N-CA-C	-5.13	98.67	109.34

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	2573	39	0
2	A	7289	0	7168	149	0
3	B	31	0	19	5	0
4	A	49	0	0	2	0
All	All	10002	0	9760	192	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 10.

All (192) 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
1:B:300:GLU:C	1:B:303:THR:HG22	2.23	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
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:1146:GLN:CD	2:A:1148:GLN:HB2	2.28	0.59
2:A:928:SER:HA	2:A:967:GLN:HB3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:1570:SER:C	2:A:1571:LEU:HD12	2.30	0.57
2:A:1577:ALA:O	2:A:1578:SER:C	2.47	0.57
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
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: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
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:154:LYS:HZ3	1:B:329:GLU:HG2	1.77	0.49
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
1:B:277:ALA:O	1:B:281:MET:HE2	2.13	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
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
2:A:1024:MET:HE2	2:A:1024:MET:HB3	1.87	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:1146:GLN:CD	2:A:1153:LYS:HG3	2.41	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
1:B:107:LYS:HZ3	1:B:111:ASP:CG	2.23	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:1024:MET:HE1	2:A:1121:TYR:CE2	2.52	0.45
1:B:281:MET:HE3	1:B:281:MET:HB2	1.70	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:1576:GLU:OE2	2:A:1636:TYR:OH	2.23	0.44
2:A:1581:SER:O	2:A:1602:LEU:N	2.35	0.44
2:A:1499:ASP:OD2	2:A:1840:LYS:NZ	2.45	0.44
2:A:1142:MET:HE1	2:A:1157:THR:OG1	2.18	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
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
2:A:1291:THR:HG21	2:A:1601:PRO:HB2	2.01	0.42
2:A:1501:ARG:NH1	2:A:1840:LYS:HD2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:B:284:GLU:OE1	1:B:322:LYS:NZ	2.47	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:1244:GLU:C	2:A:1246:LEU:H	2.27	0.42
2:A:1107:ASN:HB2	2:A:1621:VAL:CG1	2.50	0.42
2:A:1127:VAL:HG13	2:A:1128:GLU:CD	2.45	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
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: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:1100:GLU:OE2	2:A:1153:LYS:NZ	2.39	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%)	36	67
2	A	902/1027 (88%)	834 (92%)	58 (6%)	10 (1%)	11	41
All	All	1229/1362 (90%)	1143 (93%)	75 (6%)	11 (1%)	14	45

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%)	279 (99%)	2 (1%)	76	77
2	A	788/896 (88%)	783 (99%)	5 (1%)	78	79
All	All	1069/1182 (90%)	1062 (99%)	7 (1%)	76	77

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

Mol	Chain	Res	Type
1	B	150	ARG

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Mol	Chain	Res	Type
1	B	170	LYS
2	A	1127	VAL
2	A	1146	GLN
2	A	1587	GLU
2	A	1598	TYR
2	A	1814	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	1002	GLN
2	A	1156	GLN
2	A	1321	GLN
2	A	1393	ASN
2	A	1448	GLN
2	A	1452	ASN
2	A	1455	GLN
2	A	1810	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [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
4	YOA	A	1901	2	45,51,52	1.12	6 (13%)	57,73,76	2.29	14 (24%)
3	FMN	B	401	-	33,33,33	0.58	0	48,50,50	0.64	1 (2%)

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	YOA	A	1901	2	-	22/44/66/67	0/3/3/3
3	FMN	B	401	-	-	11/18/18/18	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1901	YOA	C24-S23	-3.47	1.76	1.81
4	A	1901	YOA	C18-S23	-3.25	1.77	1.82
4	A	1901	YOA	C17-S16	-3.21	1.68	1.77
4	A	1901	YOA	C07-N08	-2.20	1.35	1.39
4	A	1901	YOA	C09-N08	2.08	1.35	1.31
4	A	1901	YOA	C09-N10	-2.05	1.34	1.37

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1901	YOA	O52-S16-O51	-8.87	106.93	119.34
4	A	1901	YOA	C14-N15-S16	-7.84	104.75	120.79
4	A	1901	YOA	C07-C06-N05	-5.39	119.29	126.72
4	A	1901	YOA	N05-C04-N03	-4.52	121.74	128.58
4	A	1901	YOA	N05-C06-N10	3.44	133.02	127.17
4	A	1901	YOA	O52-S16-C17	3.42	113.26	107.86
4	A	1901	YOA	C04-N05-C06	3.38	120.08	111.83
4	A	1901	YOA	N10-C09-N08	-3.33	109.22	113.94
4	A	1901	YOA	C06-C07-N08	-2.82	107.36	110.58
4	A	1901	YOA	C07-N08-C09	2.58	107.50	103.45
4	A	1901	YOA	C06-N10-C09	2.37	108.23	105.74
4	A	1901	YOA	C29-C28-C27	-2.29	108.59	112.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1901	YOA	C35-C34-C32	2.27	112.65	108.77
4	A	1901	YOA	C55-C11-N10	-2.23	107.78	113.30
3	B	401	FMN	C4-N3-C2	-2.11	121.89	125.64

There are no chirality outliers.

All (33) torsion outliers are listed below:

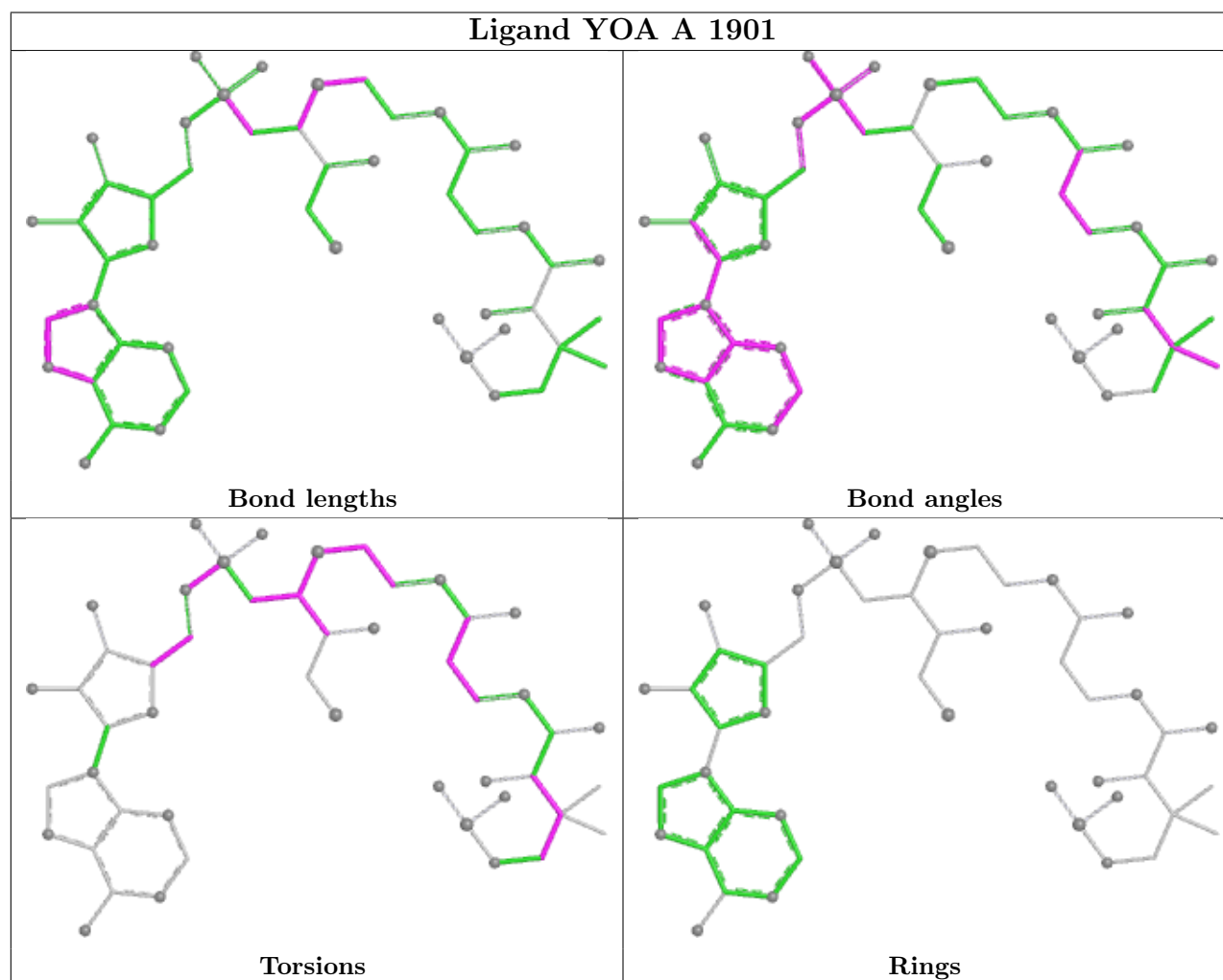
Mol	Chain	Res	Type	Atoms
3	B	401	FMN	C2'-C1'-N10-C10
3	B	401	FMN	N10-C1'-C2'-O2'
3	B	401	FMN	N10-C1'-C2'-C3'
3	B	401	FMN	C1'-C2'-C3'-C4'
3	B	401	FMN	C2'-C3'-C4'-O4'
3	B	401	FMN	C2'-C3'-C4'-C5'
3	B	401	FMN	O3'-C3'-C4'-O4'
3	B	401	FMN	O3'-C3'-C4'-C5'
3	B	401	FMN	C4'-C5'-O5'-P
4	A	1901	YOA	C53-C13-C14-N15
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	S23-C24-C25-N26
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	C35-C34-C37-O38
4	A	1901	YOA	C36-C34-C37-O38
4	A	1901	YOA	C14-N15-S16-C17
4	A	1901	YOA	C14-N15-S16-O51
3	B	401	FMN	O2'-C2'-C3'-C4'
4	A	1901	YOA	C25-C24-S23-C18
3	B	401	FMN	O2'-C2'-C3'-O3'
4	A	1901	YOA	C17-C18-S23-C24
4	A	1901	YOA	O50-C27-C28-C29

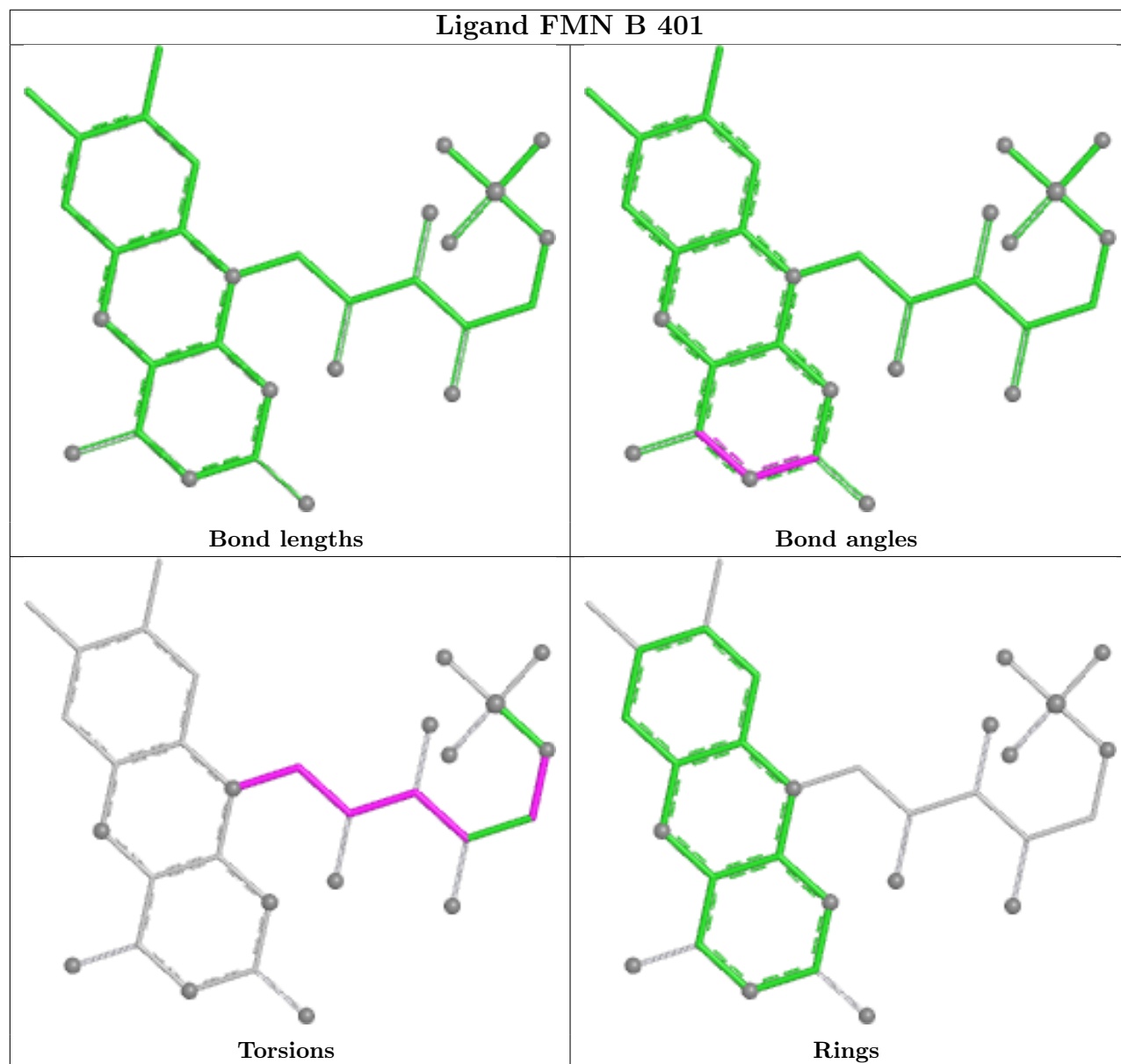
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1901	YOA	2	0
3	B	401	FMN	5	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	B	329/335 (98%)	-0.16	3 (0%) 81 60	50, 105, 257, 297	0
2	A	908/1027 (88%)	-0.14	8 (0%) 81 60	71, 187, 365, 460	0
All	All	1237/1362 (90%)	-0.14	11 (0%) 81 60	50, 164, 359, 460	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	1129	GLN	3.2
2	A	1175	PHE	2.8
2	A	1179	TYR	2.8
2	A	1507	VAL	2.5
2	A	1478	ASP	2.3
1	B	165	LEU	2.3
2	A	1500	PRO	2.2
1	B	49	ILE	2.1
2	A	1095	CYS	2.1
2	A	1784	VAL	2.0
1	B	39	PHE	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 ⓘ

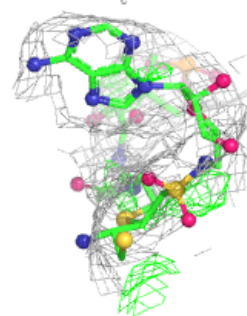
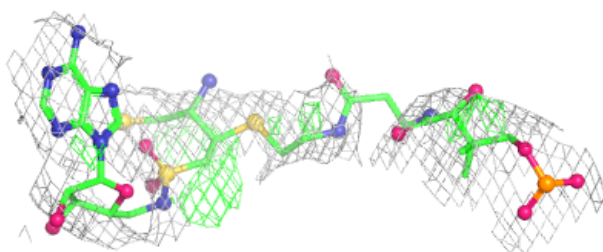
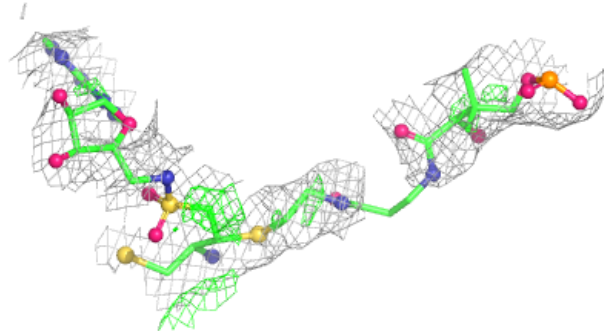
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

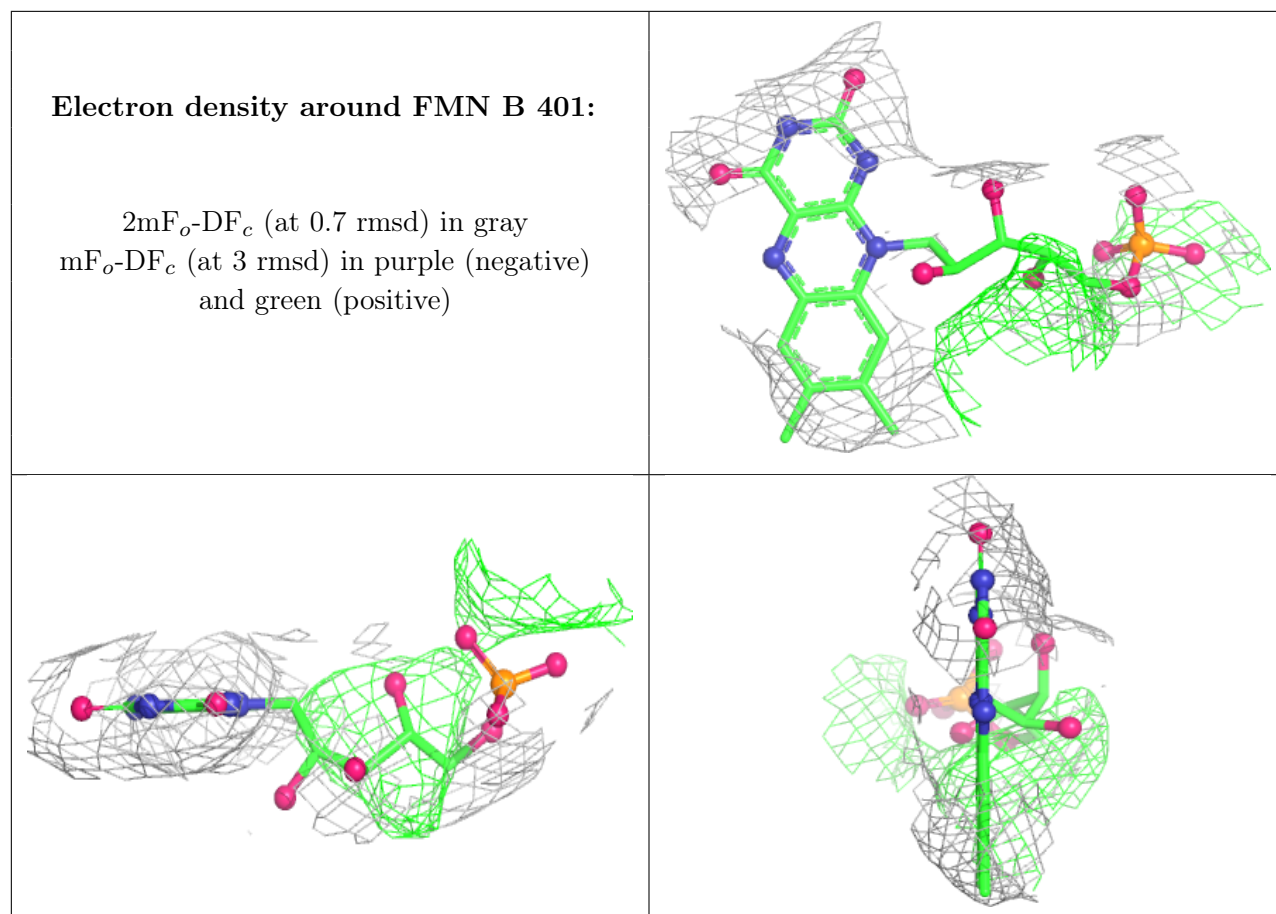
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	YOA	A	1901	49/50	0.87	0.17	203,216,231,234	0
3	FMN	B	401	31/31	0.92	0.09	48,65,93,105	0

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)





6.5 Other polymers [i](#)

There are no such residues in this entry.