



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:06 AM UTC

PDB ID : 3BI2 / pdb_00003bi2
Title : Crystal structures of fms1 in complex with its inhibitors
Authors : Huang, Q.; Hao, Q.
Deposited on : 2007-11-29
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

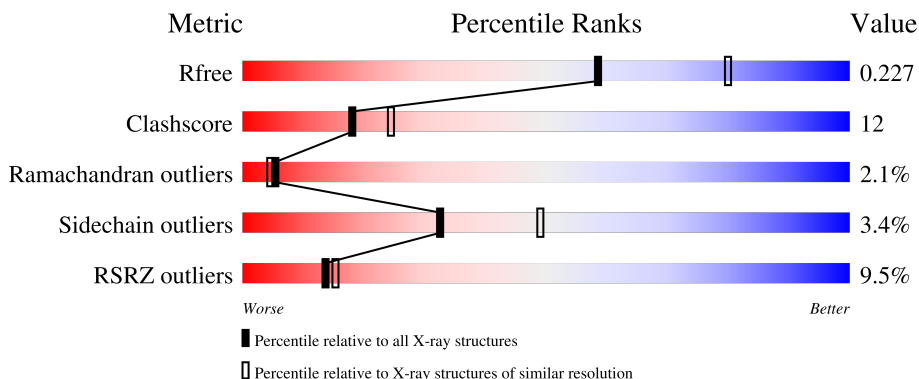
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	516	7% 72% 20% • •
1	B	516	11% 70% 23% • •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7979 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 Polyamine oxidase FMS1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	494	Total	C	N	O	S	0	0	0
			3868	2449	673	724	22			
1	B	499	Total	C	N	O	S	0	0	0
			3853	2429	672	730	22			

There are 16 discrepancies between the modelled and reference sequences:

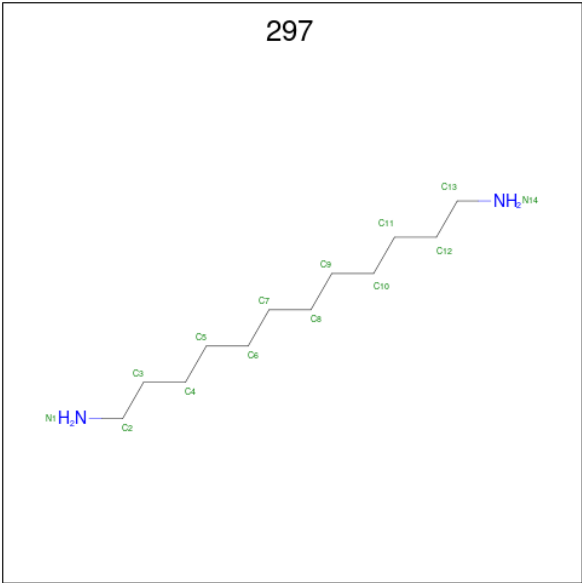
Chain	Residue	Modelled	Actual	Comment	Reference
A	509	LEU	-	expression tag	UNP P50264
A	510	GLU	-	expression tag	UNP P50264
A	511	HIS	-	expression tag	UNP P50264
A	512	HIS	-	expression tag	UNP P50264
A	513	HIS	-	expression tag	UNP P50264
A	514	HIS	-	expression tag	UNP P50264
A	515	HIS	-	expression tag	UNP P50264
A	516	HIS	-	expression tag	UNP P50264
B	509	LEU	-	expression tag	UNP P50264
B	510	GLU	-	expression tag	UNP P50264
B	511	HIS	-	expression tag	UNP P50264
B	512	HIS	-	expression tag	UNP P50264
B	513	HIS	-	expression tag	UNP P50264
B	514	HIS	-	expression tag	UNP P50264
B	515	HIS	-	expression tag	UNP P50264
B	516	HIS	-	expression tag	UNP P50264

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 3 is dodecane-1,12-diamine (CCD ID: 297) (formula: C₁₂H₂₈N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			14	12	2		
3	B	1	Total	C	N	0	0
			14	12	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	61	Total 61	O 61	0	0
4	B	63	Total 63	O 63	0	0

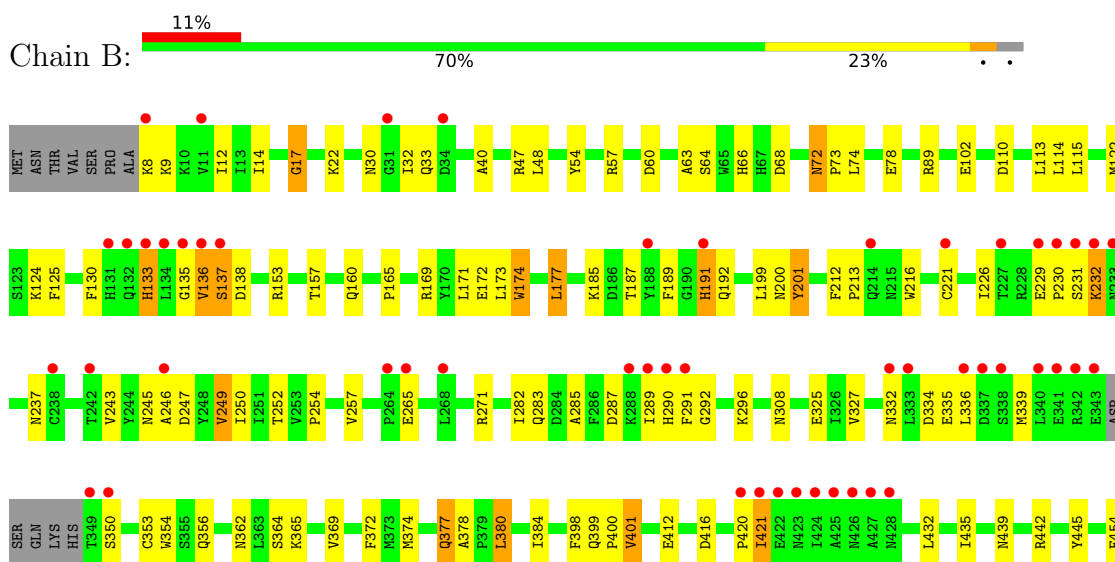
3 Residue-property plots

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: Polyamine oxidase FMS1



• Molecule 1: Polyamine oxidase FMS1





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	102.20Å 215.32Å 116.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 50.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.7 (50.00-2.30) 87.2 (50.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.222 , 0.273 0.226 , 0.227	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.876	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.8	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	0.93	EDS
Total number of atoms	7979	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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, 297

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	0.44	0/3946	0.92	11/5335 (0.2%)
1	B	0.42	0/3931	0.91	11/5319 (0.2%)
All	All	0.43	0/7877	0.91	22/10654 (0.2%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	138	ASP	N-CA-C	8.30	121.90	111.69
1	A	454	PHE	N-CA-C	-7.05	101.17	110.40
1	A	201	TYR	N-CA-C	-6.78	103.69	112.23
1	A	174	TRP	N-CA-C	-6.70	105.23	113.41
1	A	100	ILE	N-CA-C	6.53	117.25	108.11
1	B	470	GLN	N-CA-C	-6.46	104.32	111.36
1	B	201	TYR	N-CA-C	-6.07	105.14	112.54
1	A	170	TYR	N-CA-C	-5.90	105.18	112.38
1	B	454	PHE	N-CA-C	-5.90	102.15	110.29
1	B	458	ASP	CA-C-N	5.64	125.49	119.28
1	B	458	ASP	C-N-CA	5.64	125.49	119.28
1	A	182	LEU	N-CA-C	5.61	118.78	109.46
1	A	440	TRP	N-CA-C	5.60	117.39	111.28
1	A	470	GLN	N-CA-C	-5.57	105.29	111.36
1	B	66	HIS	N-CA-C	-5.56	98.70	108.20
1	B	125	PHE	N-CA-C	-5.53	105.15	111.07
1	B	491	GLY	N-CA-C	-5.52	105.81	112.49
1	A	17	GLY	N-CA-C	-5.48	103.79	112.66
1	B	292	GLY	N-CA-C	-5.42	104.02	112.58
1	A	203	SER	N-CA-C	-5.29	105.57	112.23
1	B	378	ALA	N-CA-C	-5.17	102.95	110.08
1	B	17	GLY	N-CA-C	-5.16	104.10	112.31

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3868	0	3746	89	1
1	B	3853	0	3656	97	0
2	A	53	0	31	4	0
2	B	53	0	31	3	0
3	A	14	0	28	5	0
3	B	14	0	28	0	0
4	A	61	0	0	1	0
4	B	63	0	0	0	0
All	All	7979	0	7520	188	1

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 12.

All (188) 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)
3:A:1001:297:H7	3:A:1001:297:H3A	1.46	0.97
1:B:470:GLN:HB3	1:B:474:ILE:HB	1.52	0.89
1:A:298:ILE:HD13	1:A:298:ILE:H	1.37	0.89
1:B:353:CYS:SG	1:B:400:PRO:HG2	2.14	0.87
1:A:68:ASP:HB3	1:A:192:GLN:HB2	1.58	0.85
1:A:298:ILE:HD11	1:A:436:ILE:HB	1.58	0.84
1:A:377:GLN:NE2	1:A:377:GLN:H	1.77	0.83
1:A:122:MET:HE1	1:A:145:VAL:HG22	1.59	0.82
1:B:282:ILE:HD13	1:B:465:ALA:HB1	1.63	0.81
1:A:298:ILE:CG1	1:A:436:ILE:HB	2.13	0.79
1:A:298:ILE:CD1	1:A:436:ILE:HB	2.14	0.78
1:B:68:ASP:HB3	1:B:192:GLN:HB2	1.71	0.73
1:A:298:ILE:HD13	1:A:298:ILE:N	2.04	0.73
1:A:48:LEU:CD2	1:A:63:ALA:HB3	2.20	0.72
1:B:332:ASN:HD21	1:B:334:ASP:HB2	1.53	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:72:ASN:C	1:B:72:ASN:HD22	1.98	0.70
1:A:48:LEU:HD23	1:A:63:ALA:HB3	1.74	0.70
1:B:377:GLN:NE2	1:B:377:GLN:H	1.91	0.69
1:B:335:GLU:O	1:B:339:MET:HB2	1.92	0.69
1:B:252:THR:HG22	1:B:477:ALA:HB3	1.75	0.69
1:A:157:THR:OG1	1:A:160:GLN:HG3	1.93	0.68
1:B:157:THR:OG1	1:B:160:GLN:HG3	1.95	0.67
1:B:439:ASN:HD21	1:B:442:ARG:HH11	1.41	0.66
1:A:150:LEU:O	1:A:153:ARG:HD2	1.96	0.66
1:A:450:TYR:CE1	3:A:1001:297:H6A	2.31	0.66
1:A:298:ILE:HD11	1:A:436:ILE:CB	2.25	0.66
1:B:138:ASP:HB2	1:B:185:LYS:HE3	1.78	0.65
1:B:130:PHE:CD2	1:B:185:LYS:HG2	2.33	0.64
1:B:48:LEU:CD2	1:B:63:ALA:HB3	2.27	0.64
1:A:5:SER:N	1:A:6:PRO:HD2	2.14	0.63
1:B:136:VAL:O	1:B:137:SER:HB3	2.00	0.61
1:A:103:GLU:HG3	1:A:104:ARG:H	1.65	0.61
1:B:135:GLY:O	1:B:137:SER:N	2.32	0.61
1:B:135:GLY:HA3	1:B:185:LYS:NZ	2.16	0.61
1:B:110:ASP:HB3	1:B:113:LEU:HB2	1.83	0.61
1:A:293:ALA:HB3	1:A:378:ALA:HB2	1.82	0.60
1:B:353:CYS:SG	1:B:401:VAL:HG22	2.41	0.60
1:A:252:THR:HG22	1:A:477:ALA:HB3	1.84	0.59
1:A:298:ILE:HG12	1:A:436:ILE:HB	1.83	0.59
1:A:362:ASN:HD21	1:A:364:SER:HB3	1.67	0.59
1:B:247:ASP:O	1:B:473:ARG:HD2	2.01	0.59
1:B:133:HIS:O	1:B:136:VAL:HG13	2.03	0.59
1:A:278:LEU:HA	1:A:470:GLN:HE22	1.68	0.59
1:A:142:PHE:CE1	1:A:146:MET:HE3	2.38	0.58
1:B:237:ASN:ND2	1:B:243:VAL:HG22	2.17	0.58
1:B:232:LYS:C	1:B:473:ARG:HH12	2.12	0.58
1:B:54:TYR:O	1:B:57:ARG:HG3	2.03	0.58
1:B:30:ASN:HD22	1:B:504:SER:CB	2.17	0.57
1:A:118:VAL:HG23	1:A:164:LEU:HD13	1.87	0.57
1:B:14:ILE:CD1	1:B:226:ILE:HD11	2.34	0.57
1:B:362:ASN:HA	1:B:372:PHE:CD1	2.40	0.57
1:A:353:CYS:SG	1:A:400:PRO:HB2	2.46	0.56
1:B:475:ARG:HB3	1:B:499:GLU:OE1	2.05	0.56
1:B:362:ASN:HA	1:B:372:PHE:HD1	1.71	0.56
3:A:1001:297:H7	3:A:1001:297:C3	2.29	0.56
1:B:271:ARG:HH11	1:B:271:ARG:HG3	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:358:LEU:HD22	1:A:374:MET:HE3	1.89	0.55
1:A:450:TYR:HE1	3:A:1001:297:H6A	1.71	0.55
1:B:72:ASN:C	1:B:72:ASN:ND2	2.61	0.55
1:B:153:ARG:NH1	1:B:327:VAL:O	2.40	0.55
1:B:213:PRO:O	1:B:216:TRP:HB2	2.08	0.54
1:A:504:SER:O	1:A:508:LYS:HB2	2.08	0.54
1:B:398:PHE:O	1:B:401:VAL:HG23	2.07	0.54
1:B:494:GLU:OE1	1:B:497:ARG:NH2	2.41	0.54
1:A:217:LEU:HD11	1:A:219:LEU:CD2	2.38	0.54
1:B:57:ARG:HD3	1:B:369:VAL:HG13	1.90	0.53
1:A:229:GLU:O	1:A:231:SER:N	2.42	0.53
1:A:332:ASN:ND2	1:A:335:GLU:H	2.06	0.53
1:B:12:ILE:HB	1:B:249:VAL:HG23	1.90	0.53
1:B:72:ASN:HD22	1:B:73:PRO:N	2.07	0.52
1:A:115:LEU:HB3	1:A:167:LEU:HD13	1.92	0.52
1:A:167:LEU:C	1:A:167:LEU:HD23	2.35	0.52
1:A:395:PHE:HD1	1:A:432:LEU:HD23	1.75	0.52
1:B:72:ASN:ND2	1:B:74:LEU:H	2.08	0.52
1:A:439:ASN:ND2	1:A:442:ARG:HD3	2.26	0.51
1:A:138:ASP:O	1:A:139:CYS:O	2.28	0.51
1:B:63:ALA:HA	2:B:802:FAD:N5	2.25	0.51
1:A:278:LEU:HA	1:A:470:GLN:NE2	2.25	0.51
1:B:68:ASP:CB	1:B:192:GLN:HB2	2.40	0.51
1:A:224:LYS:HD3	1:A:239:GLU:OE2	2.11	0.50
1:A:41:ARG:HD2	4:A:1034:HOH:O	2.11	0.50
1:A:41:ARG:NH1	1:A:443:ASP:OD2	2.39	0.50
1:A:150:LEU:HD13	1:A:336:LEU:HD22	1.94	0.50
1:B:254:PRO:HG2	1:B:257:VAL:HG23	1.93	0.49
1:B:362:ASN:HD21	1:B:364:SER:HB3	1.76	0.49
1:B:380:LEU:HD22	1:B:384:ILE:HG12	1.92	0.49
1:A:217:LEU:HD11	1:A:219:LEU:HD21	1.95	0.49
1:B:135:GLY:HA3	1:B:185:LYS:HZ3	1.78	0.49
1:A:17:GLY:HA3	2:A:801:FAD:H52A	1.94	0.48
1:A:237:ASN:OD1	1:A:243:VAL:HG22	2.13	0.48
1:B:135:GLY:CA	1:B:185:LYS:NZ	2.76	0.48
1:B:57:ARG:HD3	1:B:369:VAL:CG1	2.44	0.48
1:B:336:LEU:C	1:B:336:LEU:HD23	2.38	0.48
1:A:5:SER:N	1:A:6:PRO:CD	2.77	0.48
1:A:153:ARG:NH2	1:A:327:VAL:O	2.44	0.48
1:A:158:ASN:ND2	1:A:328:ARG:HH21	2.12	0.48
1:B:308:ASN:O	1:B:365:LYS:HD2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:GLY:HA2	1:B:185:LYS:HE2	1.96	0.48
1:A:298:ILE:HD11	1:A:436:ILE:CG2	2.43	0.47
1:A:377:GLN:H	1:A:377:GLN:HE21	1.56	0.47
1:A:150:LEU:HD22	1:A:333:LEU:HD22	1.95	0.47
1:B:199:LEU:HD23	1:B:199:LEU:HA	1.75	0.47
1:A:63:ALA:HA	2:A:801:FAD:N5	2.30	0.47
1:B:102:GLU:HB2	1:B:354:TRP:CE2	2.50	0.47
1:A:297:VAL:HG23	1:A:376:MET:HE2	1.95	0.47
1:A:412:GLU:OE1	1:A:429:LYS:HE3	2.14	0.47
1:B:229:GLU:O	1:B:231:SER:N	2.48	0.46
1:B:172:GLU:HG2	1:B:177:LEU:O	2.15	0.46
1:A:281:VAL:HG12	1:A:465:ALA:HB2	1.98	0.46
1:A:158:ASN:HD22	1:A:328:ARG:HH21	1.63	0.46
1:A:381:THR:O	1:A:385:GLU:HG3	2.16	0.46
1:B:113:LEU:HB3	1:B:115:LEU:HG	1.97	0.46
1:A:330:ALA:CA	1:A:339:MET:HE1	2.45	0.46
1:A:68:ASP:CB	1:A:192:GLN:HB2	2.37	0.46
1:A:11:VAL:HG22	1:A:248:TYR:HB2	1.98	0.46
1:A:260:LEU:HD22	1:A:268:LEU:HD13	1.97	0.46
1:A:356:GLN:CD	1:A:356:GLN:N	2.74	0.46
1:A:9:LYS:HG3	1:A:36:LEU:HD22	1.97	0.45
1:B:32:ILE:HG22	1:B:33:GLN:N	2.31	0.45
1:A:36:LEU:HD21	1:A:244:TYR:CE2	2.51	0.45
1:A:298:ILE:CD1	1:A:298:ILE:N	2.75	0.45
1:A:502:ARG:O	1:A:506:LEU:HG	2.17	0.45
1:B:308:ASN:OD1	1:B:364:SER:HB3	2.16	0.45
1:B:421:ILE:HG22	1:B:421:ILE:O	2.16	0.45
1:B:412:GLU:OE2	1:B:412:GLU:HA	2.16	0.45
1:A:57:ARG:HD3	1:A:369:VAL:CG1	2.47	0.45
1:B:171:LEU:HD13	1:B:187:THR:CG2	2.47	0.45
1:A:47:ARG:NE	2:A:801:FAD:O2A	2.39	0.45
1:A:376:MET:HG3	1:A:381:THR:OG1	2.16	0.45
1:B:252:THR:HA	1:B:477:ALA:O	2.17	0.45
1:B:271:ARG:HG3	1:B:271:ARG:NH1	2.31	0.45
1:A:240:ASP:OD1	1:A:242:THR:HG23	2.17	0.44
1:A:442:ARG:HH11	1:A:442:ARG:HG2	1.82	0.44
1:B:9:LYS:O	1:B:246:ALA:HA	2.17	0.44
1:A:54:TYR:CD2	1:A:55:GLN:HG3	2.53	0.44
1:B:377:GLN:H	1:B:377:GLN:HE21	1.64	0.43
1:A:164:LEU:N	1:A:165:PRO:HD2	2.33	0.43
1:A:218:LYS:HE2	1:A:218:LYS:HB3	1.85	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:VAL:O	1:B:137:SER:CB	2.66	0.43
1:B:48:LEU:HD23	1:B:63:ALA:HB3	2.01	0.43
1:A:63:ALA:HA	2:A:801:FAD:C4X	2.48	0.43
1:B:229:GLU:C	1:B:231:SER:H	2.27	0.43
1:B:285:ALA:C	1:B:287:ASP:H	2.26	0.43
1:A:335:GLU:HG2	1:A:339:MET:HE2	2.01	0.43
1:B:332:ASN:ND2	1:B:334:ASP:H	2.16	0.43
1:B:432:LEU:HD21	1:B:435:ILE:HD11	2.00	0.43
1:B:471:ASP:OD2	1:B:471:ASP:C	2.61	0.43
1:A:451:SER:O	3:A:1001:297:H4A	2.19	0.43
1:B:8:LYS:N	1:B:245:ASN:O	2.52	0.42
1:B:174:TRP:CZ2	1:B:191:HIS:CE1	3.07	0.42
1:B:189:PHE:CD1	1:B:189:PHE:C	2.97	0.42
1:A:395:PHE:CD1	1:A:432:LEU:HD23	2.54	0.42
1:A:399:GLN:HB3	1:A:400:PRO:HD3	2.00	0.42
1:A:442:ARG:HG2	1:A:442:ARG:NH1	2.34	0.42
1:B:63:ALA:HA	2:B:802:FAD:C4X	2.49	0.42
1:B:174:TRP:CD1	1:B:189:PHE:HZ	2.37	0.42
1:A:133:HIS:O	1:A:136:VAL:HG23	2.19	0.42
1:B:296:LYS:HA	1:B:374:MET:O	2.20	0.42
1:B:133:HIS:HB3	1:B:136:VAL:HG22	2.02	0.42
1:B:171:LEU:HD13	1:B:187:THR:HG22	2.02	0.42
1:A:415:ILE:HB	1:A:431:VAL:HG22	2.01	0.42
1:B:283:GLN:HE21	1:B:283:GLN:HB3	1.65	0.42
1:B:14:ILE:HD12	1:B:226:ILE:HD11	2.00	0.42
1:A:264:PRO:O	1:A:265:GLU:HB2	2.20	0.42
1:B:212:PHE:HB2	1:B:213:PRO:HD2	2.02	0.42
1:B:362:ASN:ND2	1:B:364:SER:H	2.18	0.42
1:B:399:GLN:N	1:B:400:PRO:HD2	2.35	0.41
1:A:224:LYS:HB3	1:A:239:GLU:OE1	2.20	0.41
1:B:47:ARG:NH2	1:B:254:PRO:HB3	2.36	0.41
1:B:78:GLU:OE1	1:B:89:ARG:HD3	2.20	0.41
1:B:356:GLN:N	1:B:356:GLN:CD	2.78	0.41
1:B:110:ASP:O	1:B:114:LEU:HD23	2.19	0.41
1:A:234:VAL:HG11	1:A:249:VAL:HG11	2.02	0.41
1:B:165:PRO:O	1:B:169:ARG:HG3	2.20	0.41
1:B:212:PHE:HB2	1:B:213:PRO:CD	2.51	0.41
1:B:250:ILE:HD11	1:B:503:ILE:HD12	2.02	0.41
1:A:122:MET:HG2	1:A:148:TYR:CD2	2.56	0.41
1:B:17:GLY:HA3	2:B:802:FAD:O1A	2.21	0.41
1:B:135:GLY:HA3	1:B:185:LYS:HZ1	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:ALA:HB1	1:B:445:TYR:CZ	2.55	0.41
1:B:60:ASP:HB3	1:B:63:ALA:O	2.20	0.41
1:A:439:ASN:HD21	1:A:442:ARG:HD3	1.86	0.40
1:B:416:ASP:OD1	1:B:416:ASP:C	2.63	0.40
1:B:336:LEU:HD23	1:B:336:LEU:O	2.21	0.40
1:A:9:LYS:HE3	1:A:36:LEU:HD22	2.04	0.40
1:A:104:ARG:HG2	1:A:107:VAL:HG12	2.03	0.40
1:A:122:MET:O	1:A:125:PHE:HB3	2.21	0.40
1:A:240:ASP:OD1	1:A:240:ASP:C	2.64	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:ASN:ND2	1:A:267:ASN:ND2[4_556]	1.83	0.37

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	488/516 (95%)	454 (93%)	27 (6%)	7 (1%)	9	9
1	B	495/516 (96%)	451 (91%)	30 (6%)	14 (3%)	4	2
All	All	983/1032 (95%)	905 (92%)	57 (6%)	21 (2%)	5	4

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	139	CYS
1	A	200	ASN
1	A	232	LYS
1	A	265	GLU

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Mol	Chain	Res	Type
1	B	200	ASN
1	B	289	ILE
1	B	290	HIS
1	B	291	PHE
1	A	230	PRO
1	A	231	SER
1	B	136	VAL
1	B	232	LYS
1	B	265	GLU
1	B	230	PRO
1	B	420	PRO
1	B	510	GLU
1	A	131	HIS
1	B	133	HIS
1	B	137	SER
1	B	350	SER
1	B	421	ILE

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	414/457 (91%)	402 (97%)	12 (3%)	37	55
1	B	405/457 (89%)	389 (96%)	16 (4%)	28	42
All	All	819/914 (90%)	791 (97%)	28 (3%)	32	49

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LEU
1	A	122	MET
1	A	150	LEU
1	A	177	LEU
1	A	201	TYR
1	A	221	CYS

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Mol	Chain	Res	Type
1	A	267	ASN
1	A	298	ILE
1	A	332	ASN
1	A	333	LEU
1	A	377	GLN
1	A	388	ARG
1	B	22	LYS
1	B	64	SER
1	B	72	ASN
1	B	122	MET
1	B	124	LYS
1	B	173	LEU
1	B	174	TRP
1	B	177	LEU
1	B	191	HIS
1	B	201	TYR
1	B	221	CYS
1	B	249	VAL
1	B	325	GLU
1	B	377	GLN
1	B	380	LEU
1	B	401	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	49	GLN
1	A	158	ASN
1	A	259	ASN
1	A	263	GLN
1	A	267	ASN
1	A	332	ASN
1	A	362	ASN
1	A	377	GLN
1	A	434	ASN
1	A	439	ASN
1	B	29	GLN
1	B	30	ASN
1	B	72	ASN
1	B	109	HIS
1	B	237	ASN

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Mol	Chain	Res	Type
1	B	283	GLN
1	B	329	ASN
1	B	332	ASN
1	B	362	ASN
1	B	377	GLN
1	B	434	ASN
1	B	439	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	297	B	1002	-	13,13,13	0.33	0	12,12,12	0.60	0
2	FAD	B	802	-	58,58,58	1.57	7 (12%)	85,89,89	1.86	12 (14%)
2	FAD	A	801	-	58,58,58	1.63	5 (8%)	85,89,89	1.85	11 (12%)
3	297	A	1001	-	13,13,13	0.34	0	12,12,12	0.45	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	297	B	1002	-	-	0/11/11/11	-
2	FAD	B	802	-	-	3/34/50/50	0/6/6/6
2	FAD	A	801	-	-	9/34/50/50	0/6/6/6
3	297	A	1001	-	-	0/11/11/11	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	FAD	PA-O3P	7.24	1.67	1.59
2	B	802	FAD	PA-O3P	6.92	1.67	1.59
2	A	801	FAD	C4X-N5	4.44	1.40	1.30
2	B	802	FAD	C4X-N5	4.20	1.39	1.30
2	A	801	FAD	P-O3P	4.14	1.64	1.59
2	B	802	FAD	P-O3P	3.83	1.63	1.59
2	A	801	FAD	C2A-N1A	2.84	1.39	1.33
2	B	802	FAD	C2A-N1A	2.64	1.38	1.33
2	A	801	FAD	C10-N1	2.60	1.38	1.33
2	B	802	FAD	C9A-C5X	2.28	1.44	1.41
2	B	802	FAD	C9-C8	2.14	1.42	1.39
2	B	802	FAD	C10-N1	2.10	1.37	1.33

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FAD	O3P-PA-O1A	-10.50	79.11	110.70
2	B	802	FAD	O3P-PA-O1A	-9.57	81.91	110.70
2	B	802	FAD	O2A-PA-O3P	-8.75	83.62	107.27
2	A	801	FAD	O2A-PA-O3P	-8.06	85.49	107.27
2	B	802	FAD	O2A-PA-O1A	4.48	133.28	112.44
2	A	801	FAD	O2A-PA-O1A	4.46	133.20	112.44
2	B	802	FAD	C4B-O4B-C1B	-3.11	102.60	109.47
2	B	802	FAD	O4B-C1B-C2B	-2.91	100.39	106.62
2	A	801	FAD	C9A-C5X-N5	-2.81	119.48	122.45
2	A	801	FAD	O3'-C3'-C4'	-2.74	102.70	108.93
2	B	802	FAD	C9A-C5X-N5	-2.63	119.67	122.45
2	B	802	FAD	O3'-C3'-C4'	-2.58	103.07	108.93
2	A	801	FAD	C4B-O4B-C1B	-2.33	104.33	109.47
2	B	802	FAD	C4'-C3'-C2'	2.26	117.34	113.57
2	A	801	FAD	O4B-C1B-C2B	-2.24	101.83	106.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	802	FAD	N3A-C2A-N1A	-2.20	125.25	128.58
2	A	801	FAD	C4-C4X-N5	2.13	121.15	118.21
2	B	802	FAD	C10-C4X-N5	-2.10	120.52	124.81
2	A	801	FAD	N3A-C2A-N1A	-2.09	125.42	128.58
2	A	801	FAD	C5'-C4'-C3'	-2.05	108.35	112.22
2	B	802	FAD	C4-C4X-N5	2.05	121.04	118.21
2	B	802	FAD	C4X-C10-N10	2.03	119.38	116.48
2	A	801	FAD	C10-C4X-N5	-2.00	120.72	124.81

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	FAD	C5B-O5B-PA-O1A
2	B	802	FAD	C5B-O5B-PA-O1A
2	A	801	FAD	C2'-C3'-C4'-O4'
2	A	801	FAD	P-O3P-PA-O1A
2	A	801	FAD	C5B-O5B-PA-O3P
2	B	802	FAD	C5B-O5B-PA-O3P
2	A	801	FAD	O3'-C3'-C4'-O4'
2	A	801	FAD	C4B-C5B-O5B-PA
2	A	801	FAD	O3'-C3'-C4'-C5'
2	B	802	FAD	C4B-C5B-O5B-PA
2	A	801	FAD	C2'-C3'-C4'-C5'
2	A	801	FAD	P-O3P-PA-O2A

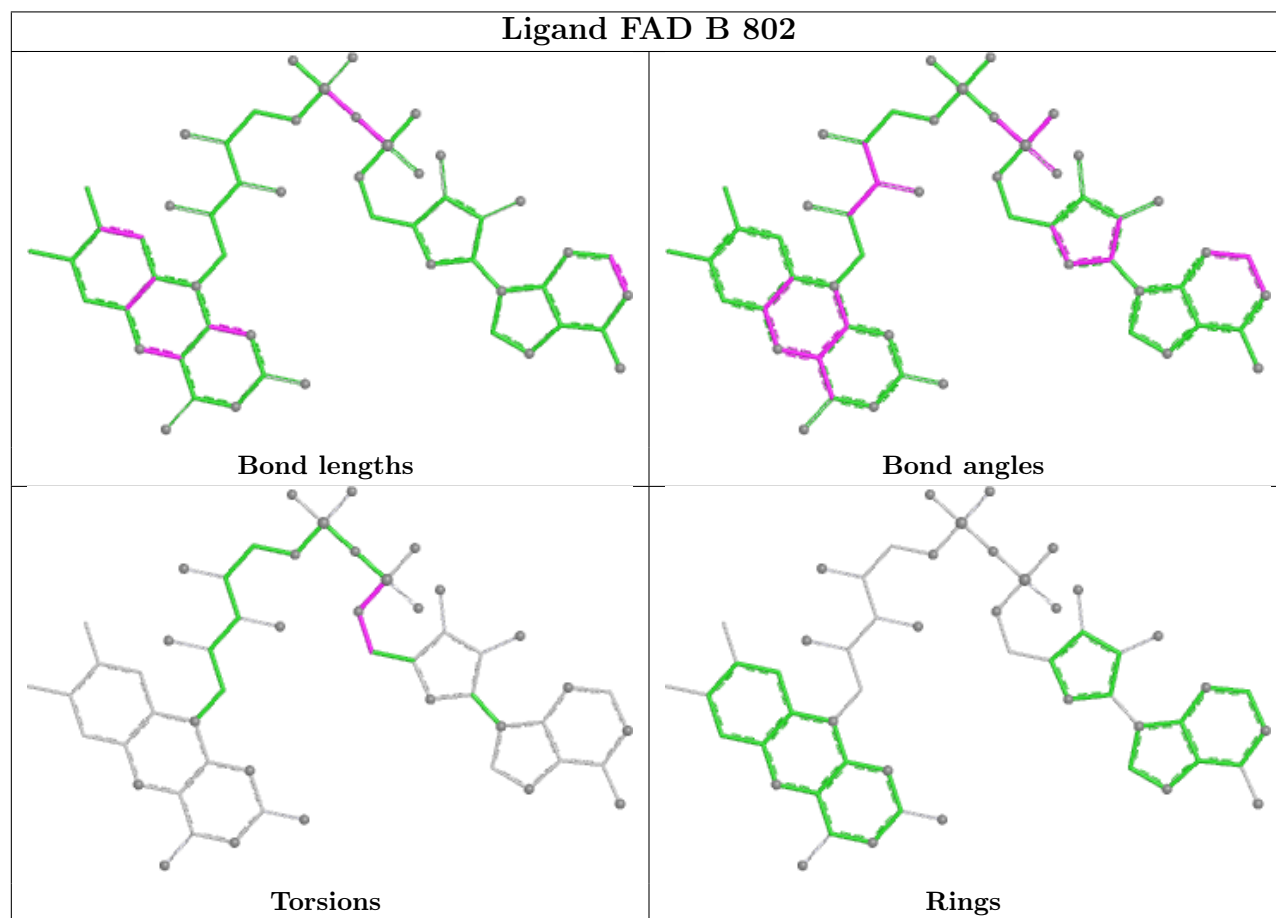
There are no ring outliers.

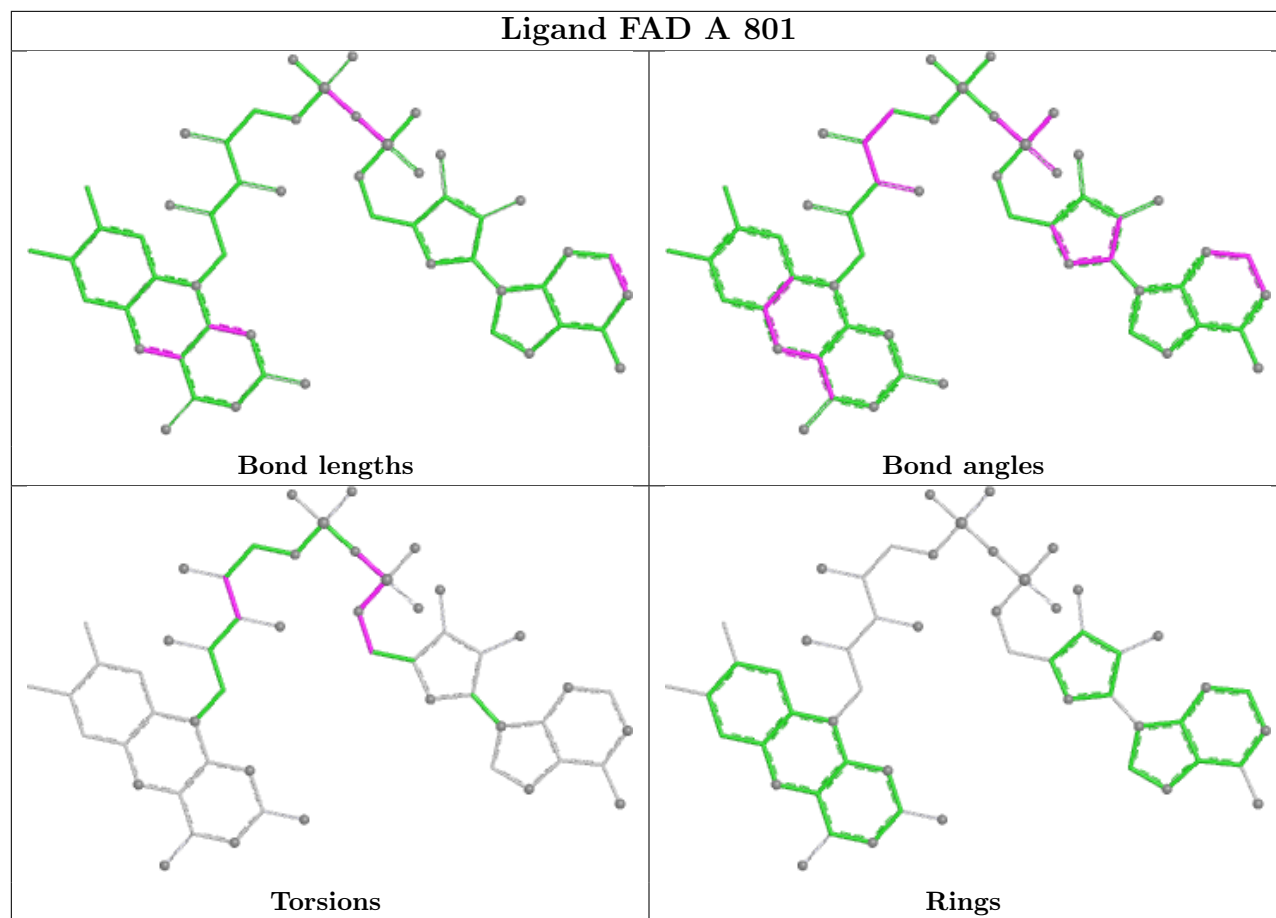
3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	802	FAD	3	0
2	A	801	FAD	4	0
3	A	1001	297	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	494/516 (95%)	0.54	35 (7%)	22 24	25, 44, 66, 86	0
1	B	499/516 (96%)	0.63	59 (11%)	9 10	27, 43, 78, 100	0
All	All	993/1032 (96%)	0.58	94 (9%)	14 15	25, 43, 73, 100	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	289	ILE	7.2
1	B	238	CYS	5.8
1	B	132	GLN	5.5
1	B	134	LEU	5.5
1	A	343	GLU	5.4
1	B	136	VAL	5.3
1	B	264	PRO	5.1
1	B	135	GLY	5.1
1	B	233	ASN	4.9
1	B	509	LEU	4.9
1	A	510	GLU	4.8
1	B	230	PRO	4.7
1	B	349	THR	4.5
1	A	230	PRO	4.2
1	B	288	LYS	4.1
1	A	349	THR	4.0
1	B	350	SER	4.0
1	B	341	GLU	3.9
1	B	427	ALA	3.8
1	B	421	ILE	3.8
1	A	238	CYS	3.7
1	B	229	GLU	3.7
1	B	265	GLU	3.7
1	A	334	ASP	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	509	LEU	3.6
1	B	424	ILE	3.6
1	B	337	ASP	3.6
1	A	221	CYS	3.6
1	A	342	ARG	3.5
1	B	333	LEU	3.4
1	B	8	LYS	3.4
1	B	420	PRO	3.3
1	B	343	GLU	3.3
1	B	231	SER	3.2
1	B	133	HIS	3.1
1	B	290	HIS	3.1
1	B	506	LEU	3.1
1	A	229	GLU	3.1
1	A	431	VAL	2.9
1	B	342	ARG	2.9
1	A	7	ALA	2.9
1	A	264	PRO	2.9
1	B	191	HIS	2.9
1	B	232	LYS	2.8
1	A	428	ASN	2.7
1	A	233	ASN	2.7
1	A	429	LYS	2.7
1	B	291	PHE	2.6
1	A	298	ILE	2.6
1	A	392	GLU	2.6
1	A	5	SER	2.6
1	B	338	SER	2.6
1	B	214	GLN	2.5
1	A	239	GLU	2.5
1	B	340	LEU	2.5
1	A	135	GLY	2.5
1	A	415	ILE	2.5
1	B	428	ASN	2.5
1	B	268	LEU	2.5
1	B	336	LEU	2.5
1	B	507	LEU	2.5
1	B	11	VAL	2.4
1	B	423	ASN	2.4
1	A	350	SER	2.4
1	B	221	CYS	2.4
1	A	263	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	332	ASN	2.3
1	B	469	GLY	2.3
1	B	425	ALA	2.3
1	A	133	HIS	2.3
1	B	426	ASN	2.2
1	A	31	GLY	2.2
1	B	137	SER	2.2
1	A	217	LEU	2.2
1	B	503	ILE	2.2
1	A	339	MET	2.2
1	A	341	GLU	2.2
1	B	242	THR	2.2
1	A	340	LEU	2.1
1	B	34	ASP	2.1
1	B	246	ALA	2.1
1	B	505	ASP	2.1
1	A	372	PHE	2.1
1	A	32	ILE	2.1
1	A	104	ARG	2.1
1	B	227	THR	2.1
1	B	504	SER	2.1
1	B	508	LYS	2.1
1	B	422	GLU	2.1
1	A	206	GLN	2.1
1	A	427	ALA	2.0
1	B	188	TYR	2.0
1	B	31	GLY	2.0
1	B	131	HIS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

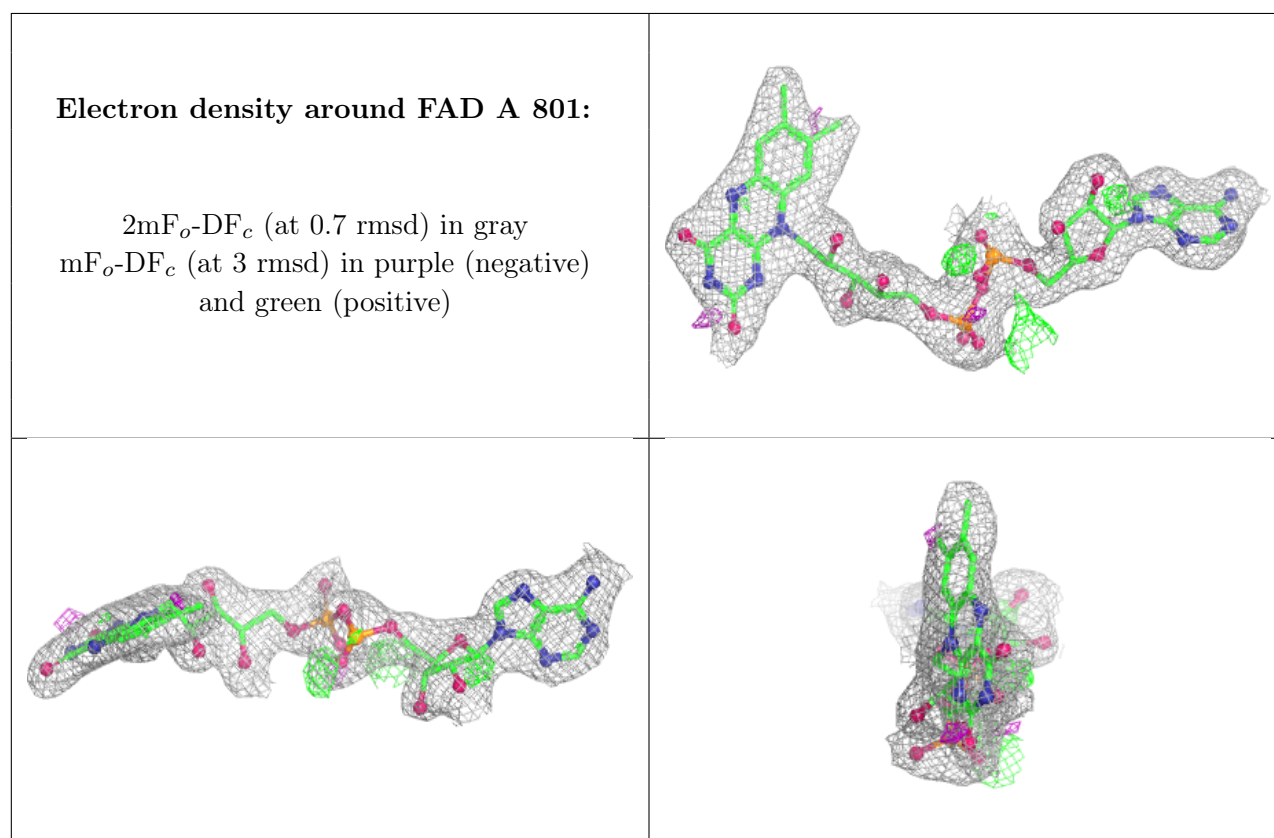
6.4 Ligands [i](#)

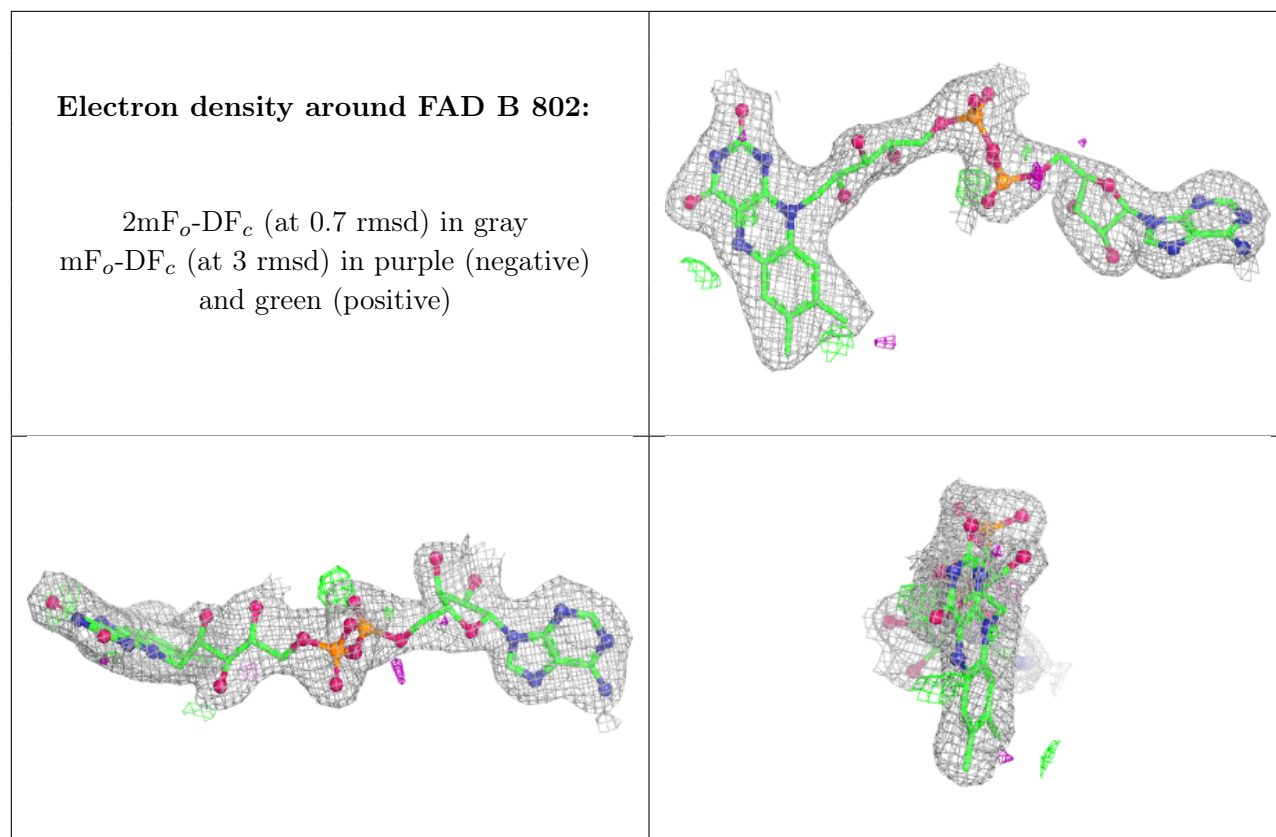
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	297	A	1001	14/14	0.79	0.23	65,65,68,68	0
3	297	B	1002	14/14	0.84	0.20	48,59,72,73	0
2	FAD	A	801	53/53	0.93	0.09	28,33,39,43	0
2	FAD	B	802	53/53	0.95	0.08	25,35,41,42	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.