



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:20 PM UTC

PDB ID : 3BI5 / pdb_00003bi5
Title : Crystal structures of fms1 in complex with its inhibitors
Authors : Huang, Q.; Hao, Q.
Deposited on : 2007-11-29
Resolution : 2.50 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

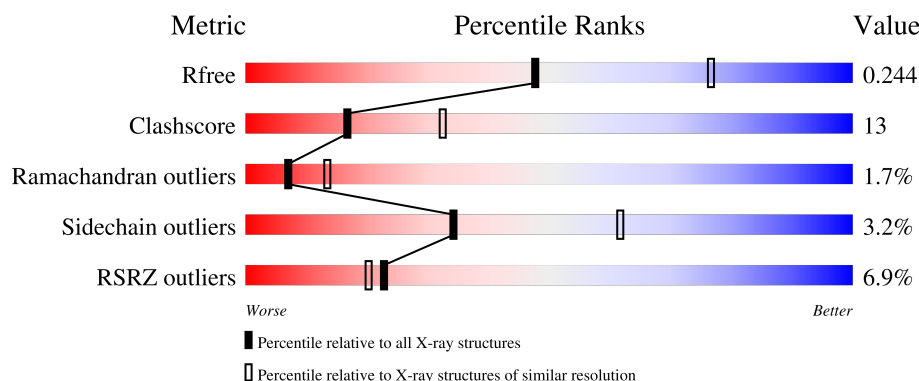
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	4% 73% 21% • •
1	B	516	9% 63% 30% • •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7975 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	497	Total	C	N	O	S	0	0	0
			3890	2462	681	725	22			
1	B	499	Total	C	N	O	S	0	0	0
			3842	2422	668	730	22			

There are 16 discrepancies between the modelled and reference sequences:

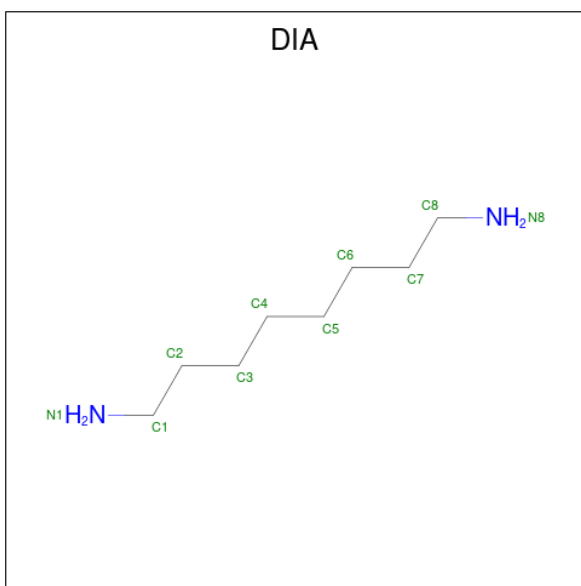
Chain	Residue	Modelled	Actual	Comment	Reference
A	509	LEU	-	expression tag	UNP P50264
A	510	GLU	-	expression tag	UNP P50264
A	511	HIS	-	expression tag	UNP P50264
A	512	HIS	-	expression tag	UNP P50264
A	513	HIS	-	expression tag	UNP P50264
A	514	HIS	-	expression tag	UNP P50264
A	515	HIS	-	expression tag	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 OCTANE 1,8-DIAMINE (CCD ID: DIA) (formula: $\text{C}_8\text{H}_{20}\text{N}_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N 10 8 2	0	0
3	B	1	Total C N 10 8 2	0	0

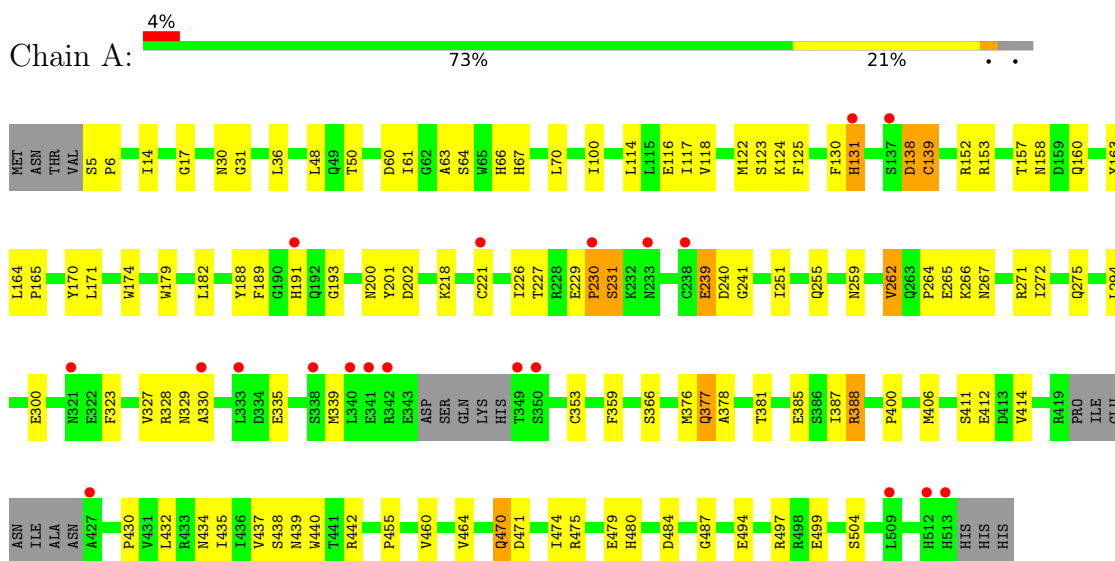
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	69	Total 69	O 69	0	0
4	B	48	Total 48	O 48	0	0

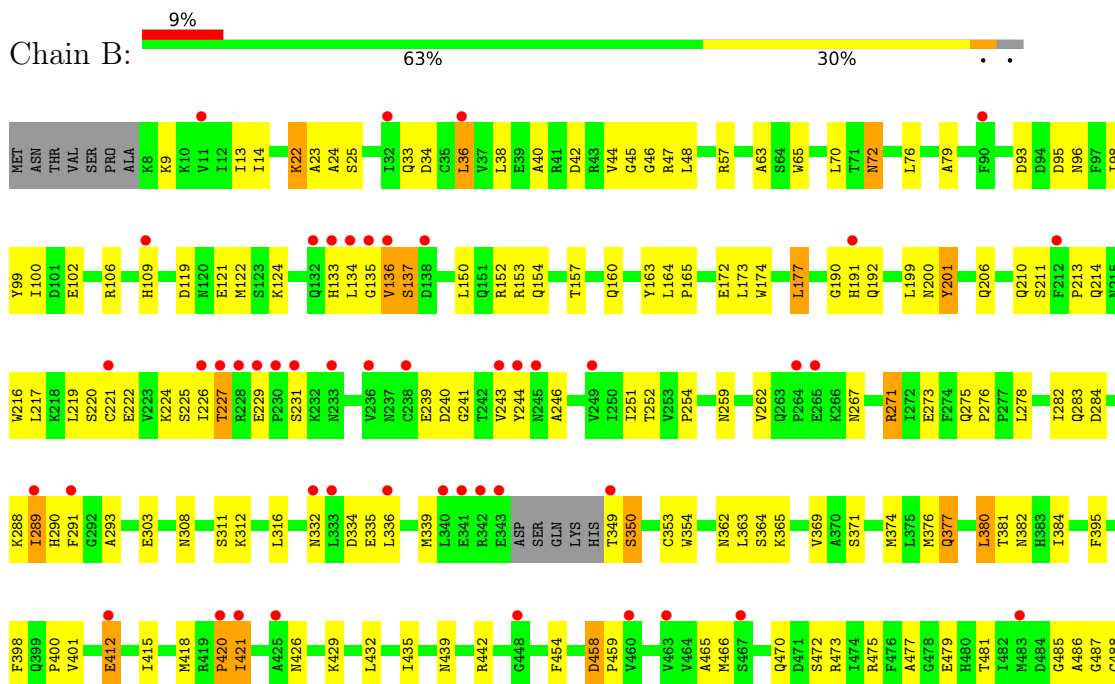
3 Residue-property plots

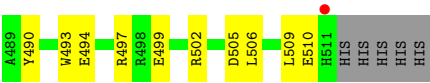
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	164.42Å 103.00Å 77.24Å 90.00° 94.30° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	93.4 (50.00-2.50) 93.4 (50.00-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.06 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.228 , 0.290 0.232 , 0.244	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	27.6	Xtriage
Anisotropy	0.408	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7975	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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, DIA

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	0.48	0/3972	0.92	15/5373 (0.3%)
1	B	0.45	0/3920	0.93	13/5306 (0.2%)
All	All	0.47	0/7892	0.93	28/10679 (0.3%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	454	PHE	N-CA-C	-7.95	99.98	110.40
1	A	138	ASP	N-CA-C	7.08	120.37	111.24
1	A	366	SER	N-CA-C	6.64	118.60	111.36
1	B	267	ASN	N-CA-C	-6.60	106.44	114.75
1	B	472	SER	N-CA-C	6.59	118.55	111.36
1	A	66	HIS	N-CA-C	-6.35	98.39	108.73
1	A	100	ILE	N-CA-C	6.33	117.24	108.12
1	A	163	TYR	N-CA-C	6.24	120.09	112.23
1	A	470	GLN	N-CA-C	-6.12	104.53	111.14
1	B	201	TYR	N-CA-C	-5.93	105.30	112.54
1	A	440	TRP	N-CA-C	5.86	117.67	111.28
1	A	174	TRP	N-CA-C	-5.86	106.27	113.41
1	B	65	TRP	N-CA-C	5.86	119.26	109.95
1	A	294	LEU	N-CA-C	-5.62	102.57	110.50
1	A	182	LEU	N-CA-C	5.53	118.94	110.20
1	B	293	ALA	N-CA-C	5.47	117.88	110.06
1	B	488	CYS	N-CA-C	5.44	118.57	110.14
1	A	179	TRP	N-CA-C	5.30	119.22	112.34
1	A	170	TYR	N-CA-C	-5.28	106.09	112.54
1	A	378	ALA	N-CA-C	-5.28	103.52	110.39
1	B	163	TYR	N-CA-C	5.20	118.78	112.23
1	B	100	ILE	N-CA-C	5.18	115.36	108.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	31	GLY	N-CA-C	5.07	122.59	115.30
1	A	17	GLY	N-CA-C	-5.06	104.86	112.60
1	B	288	LYS	N-CA-C	-5.05	105.86	112.72
1	B	289	ILE	N-CA-C	5.02	119.78	109.34
1	B	458	ASP	CA-C-N	5.01	124.79	119.28
1	B	458	ASP	C-N-CA	5.01	124.79	119.28

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	3890	0	3743	68	0
1	B	3842	0	3633	132	0
2	A	53	0	31	1	0
2	B	53	0	31	2	0
3	A	10	0	20	2	0
3	B	10	0	20	2	0
4	A	69	0	0	2	0
4	B	48	0	0	1	0
All	All	7975	0	7478	196	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 13.

All (196) 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:B:133:HIS:HB3	1:B:136:VAL:HG13	1.36	1.08
1:B:278:LEU:HA	1:B:470:GLN:HE22	1.14	1.06
1:B:227:THR:HG23	1:B:275:GLN:HB3	1.49	0.91
1:B:282:ILE:HD13	1:B:465:ALA:HB1	1.55	0.85
1:B:278:LEU:HA	1:B:470:GLN:NE2	1.94	0.82
1:A:353:CYS:SG	1:A:400:PRO:HG2	2.21	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:PRO:O	1:A:265:GLU:HB3	1.81	0.79
1:B:353:CYS:SG	1:B:400:PRO:HG2	2.24	0.78
1:A:377:GLN:NE2	1:A:377:GLN:H	1.84	0.76
1:B:380:LEU:HD22	1:B:384:ILE:HG12	1.66	0.75
1:B:308:ASN:O	1:B:365:LYS:HD2	1.88	0.74
1:B:506:LEU:O	1:B:510:GLU:HG3	1.87	0.73
1:B:479:GLU:OE1	1:B:487:GLY:HA2	1.88	0.73
1:A:118:VAL:HG23	1:A:164:LEU:HD13	1.70	0.72
1:B:213:PRO:HG2	1:B:216:TRP:CE2	2.24	0.72
1:B:439:ASN:HD21	1:B:442:ARG:HB2	1.55	0.71
1:B:275:GLN:NE2	1:B:276:PRO:HA	2.06	0.71
1:B:213:PRO:HG2	1:B:216:TRP:CD2	2.25	0.70
1:A:406:MET:HE2	1:A:406:MET:HA	1.73	0.69
1:B:72:ASN:C	1:B:72:ASN:HD22	2.01	0.69
1:B:252:THR:HG22	1:B:477:ALA:HB3	1.76	0.67
1:A:330:ALA:CA	1:A:339:MET:HE1	2.24	0.67
1:B:48:LEU:CD2	1:B:63:ALA:HB3	2.26	0.66
1:A:239:GLU:HG3	1:A:240:ASP:N	2.10	0.66
1:B:362:ASN:ND2	1:B:364:SER:H	1.93	0.66
1:A:193:GLY:HA2	1:B:121:GLU:HG3	1.78	0.65
1:B:282:ILE:CD1	1:B:465:ALA:HB1	2.26	0.65
1:A:259:ASN:O	1:A:262:VAL:HG22	1.97	0.64
1:B:172:GLU:HG2	1:B:177:LEU:O	1.97	0.64
1:B:251:ILE:HD12	1:B:251:ILE:N	2.13	0.62
1:A:229:GLU:O	1:A:231:SER:N	2.32	0.62
1:A:414:VAL:HA	1:A:430:PRO:HG2	1.82	0.61
1:B:239:GLU:HA	1:B:239:GLU:OE1	1.99	0.61
1:A:227:THR:HG23	1:A:275:GLN:HB2	1.82	0.61
1:A:48:LEU:CD2	1:A:63:ALA:HB3	2.32	0.60
1:A:479:GLU:OE1	1:A:487:GLY:HA2	2.02	0.60
1:B:473:ARG:HH11	1:B:473:ARG:HG3	1.65	0.60
1:B:439:ASN:ND2	1:B:442:ARG:HB2	2.15	0.60
1:A:5:SER:N	1:A:6:PRO:CD	2.65	0.60
1:B:271:ARG:HH11	1:B:271:ARG:HG3	1.65	0.60
1:B:311:SER:C	1:B:312:LYS:HD3	2.28	0.59
1:B:412:GLU:CD	1:B:429:LYS:HD2	2.28	0.58
1:B:502:ARG:HH11	1:B:502:ARG:HG2	1.68	0.58
1:B:362:ASN:HD21	1:B:364:SER:CB	2.17	0.58
1:B:133:HIS:O	1:B:134:LEU:HB2	2.04	0.57
1:A:158:ASN:HD22	1:A:328:ARG:HH21	1.51	0.57
1:B:376:MET:HG3	1:B:381:THR:OG1	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213:PRO:O	1:B:216:TRP:HB2	2.05	0.57
1:A:385:GLU:OE2	1:A:438:SER:HB2	2.04	0.57
1:B:395:PHE:HB2	1:B:418:MET:HE1	1.87	0.56
1:A:152:ARG:CZ	1:B:70:LEU:HD23	2.35	0.56
1:B:362:ASN:ND2	1:B:362:ASN:C	2.61	0.56
1:A:226:ILE:HD12	1:A:272:ILE:HG21	1.88	0.56
1:B:415:ILE:HD13	1:B:426:ASN:OD1	2.06	0.56
1:B:14:ILE:CD1	1:B:226:ILE:HD11	2.35	0.56
1:A:262:VAL:O	1:A:262:VAL:HG23	2.06	0.55
1:A:376:MET:HG3	1:A:381:THR:OG1	2.06	0.55
1:A:60:ASP:HB3	4:A:2850:HOH:O	2.05	0.55
1:B:63:ALA:HA	2:B:802:FAD:N5	2.21	0.55
1:B:271:ARG:HD3	1:B:271:ARG:C	2.32	0.55
1:B:95:ASP:OD1	1:B:96:ASN:N	2.39	0.54
1:A:475:ARG:HB3	1:A:499:GLU:OE1	2.08	0.54
1:A:359:PHE:CE2	3:A:2841:DIA:HC71	2.43	0.54
1:B:225:SER:HB3	1:B:273:GLU:HB3	1.89	0.54
1:B:45:GLY:O	1:B:48:LEU:HB2	2.08	0.53
1:A:157:THR:HG23	1:A:160:GLN:OE1	2.09	0.53
1:A:164:LEU:N	1:A:165:PRO:HD2	2.24	0.53
1:B:76:LEU:O	1:B:79:ALA:HB3	2.08	0.53
4:A:2862:HOH:O	1:B:109:HIS:HE1	1.92	0.53
1:B:382:ASN:ND2	1:B:382:ASN:H	2.06	0.52
1:B:72:ASN:C	1:B:72:ASN:ND2	2.68	0.52
1:B:332:ASN:ND2	1:B:334:ASP:H	2.06	0.52
1:B:48:LEU:HD23	1:B:63:ALA:HB3	1.91	0.52
1:B:502:ARG:HG2	1:B:502:ARG:NH1	2.25	0.52
1:B:251:ILE:HD12	1:B:251:ILE:H	1.74	0.52
1:B:439:ASN:ND2	1:B:442:ARG:HD3	2.25	0.52
1:A:124:LYS:HE3	1:B:191:HIS:O	2.11	0.51
1:B:174:TRP:CE3	3:B:2842:DIA:HC51	2.44	0.51
1:B:9:LYS:O	1:B:246:ALA:HA	2.11	0.51
1:B:157:THR:OG1	1:B:160:GLN:HG3	2.10	0.51
1:B:225:SER:CB	1:B:273:GLU:HB3	2.40	0.51
1:B:48:LEU:HD21	1:B:63:ALA:HB3	1.92	0.51
1:B:262:VAL:HG12	1:B:283:GLN:HG2	1.91	0.51
1:B:190:GLY:C	1:B:192:GLN:HE21	2.19	0.51
1:A:63:ALA:HA	2:A:801:FAD:N5	2.26	0.50
1:B:362:ASN:C	1:B:362:ASN:HD22	2.17	0.50
1:B:219:LEU:O	1:B:220:SER:HB2	2.11	0.50
1:B:164:LEU:N	1:B:165:PRO:HD2	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:437:VAL:HG12	1:A:438:SER:N	2.27	0.50
1:B:135:GLY:O	1:B:137:SER:N	2.44	0.49
1:A:470:GLN:HB3	1:A:474:ILE:HB	1.95	0.49
1:B:214:GLN:C	1:B:216:TRP:H	2.20	0.49
1:B:432:LEU:HD21	1:B:435:ILE:HD11	1.94	0.49
1:B:475:ARG:HB3	1:B:499:GLU:OE1	2.13	0.49
1:A:48:LEU:HD22	1:A:63:ALA:HB3	1.95	0.49
1:B:36:LEU:HD22	1:B:38:LEU:CD2	2.43	0.49
1:A:5:SER:N	1:A:6:PRO:HD2	2.28	0.48
1:A:30:ASN:HD22	1:A:504:SER:CB	2.26	0.48
1:B:13:ILE:HD13	1:B:23:ALA:HB3	1.96	0.48
1:B:311:SER:O	1:B:312:LYS:HD3	2.13	0.48
1:B:439:ASN:HD21	1:B:442:ARG:HD3	1.79	0.48
1:A:377:GLN:H	1:A:377:GLN:HE21	1.59	0.48
1:B:494:GLU:OE2	1:B:497:ARG:NH2	2.47	0.48
1:B:284:ASP:HA	4:B:2855:HOH:O	2.14	0.48
1:A:122:MET:O	1:A:125:PHE:HB3	2.13	0.48
1:B:353:CYS:SG	1:B:401:VAL:HG13	2.54	0.47
1:A:153:ARG:NH2	1:A:327:VAL:O	2.47	0.47
1:B:57:ARG:HD3	1:B:369:VAL:CG1	2.44	0.47
1:A:189:PHE:CZ	1:A:191:HIS:NE2	2.83	0.47
1:A:239:GLU:C	1:A:241:GLY:H	2.21	0.47
1:B:22:LYS:HG3	1:B:493:TRP:CD1	2.49	0.47
1:B:44:VAL:CG1	1:B:217:LEU:HD21	2.45	0.47
1:A:158:ASN:ND2	1:A:328:ARG:HH21	2.13	0.47
1:B:206:GLN:O	1:B:210:GLN:HG3	2.15	0.46
1:B:473:ARG:HG3	1:B:473:ARG:NH1	2.29	0.46
1:A:385:GLU:O	1:A:388:ARG:HG2	2.15	0.46
1:B:420:PRO:O	1:B:421:ILE:HG12	2.15	0.46
1:B:243:VAL:HG12	1:B:244:TYR:N	2.31	0.46
1:B:335:GLU:HG2	1:B:339:MET:HE3	1.97	0.46
1:B:395:PHE:HB2	1:B:418:MET:CE	2.46	0.46
1:A:255:GLN:HB2	1:A:480:HIS:CD2	2.51	0.46
1:B:36:LEU:HD22	1:B:38:LEU:HD23	1.97	0.46
1:B:259:ASN:O	1:B:262:VAL:HG22	2.16	0.46
1:B:229:GLU:C	1:B:231:SER:N	2.74	0.45
1:B:229:GLU:C	1:B:231:SER:H	2.23	0.45
1:B:199:LEU:HD23	1:B:199:LEU:HA	1.82	0.45
1:B:42:ASP:OD1	1:B:42:ASP:C	2.59	0.45
1:B:133:HIS:C	1:B:135:GLY:H	2.25	0.45
1:B:99:TYR:HB3	1:B:316:LEU:HD11	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:ASP:OD1	1:B:241:GLY:N	2.50	0.45
1:A:114:LEU:HD22	1:A:117:ILE:HD12	1.98	0.44
1:B:102:GLU:HB2	1:B:354:TRP:CZ2	2.52	0.44
1:B:213:PRO:CG	1:B:216:TRP:CE2	3.00	0.44
1:B:485:GLY:O	1:B:486:ALA:C	2.60	0.44
1:B:98:ILE:HG23	1:B:106:ARG:NE	2.32	0.44
1:B:119:ASP:O	1:B:122:MET:HB3	2.17	0.44
1:B:150:LEU:CD1	1:B:153:ARG:HH21	2.30	0.44
1:B:362:ASN:HD21	1:B:364:SER:HB3	1.82	0.44
1:B:46:GLY:C	1:B:48:LEU:H	2.25	0.44
1:B:93:ASP:O	1:B:312:LYS:HE3	2.17	0.44
1:A:50:THR:OG1	1:A:202:ASP:HB3	2.16	0.44
1:A:226:ILE:HD13	1:A:251:ILE:HD11	1.99	0.44
1:B:377:GLN:NE2	1:B:377:GLN:H	2.16	0.44
1:A:61:ILE:HA	1:A:61:ILE:HD12	1.85	0.44
1:A:471:ASP:OD1	1:A:471:ASP:C	2.60	0.44
1:B:275:GLN:HA	1:B:276:PRO:C	2.42	0.44
1:A:494:GLU:OE1	1:A:497:ARG:NH2	2.50	0.43
1:A:70:LEU:HD23	1:B:152:ARG:CZ	2.49	0.43
1:B:133:HIS:HB3	1:B:136:VAL:CG1	2.27	0.43
1:B:374:MET:HE1	1:B:398:PHE:CE2	2.53	0.43
1:B:505:ASP:O	1:B:509:LEU:HG	2.18	0.43
1:B:136:VAL:O	1:B:136:VAL:HG23	2.19	0.43
1:B:362:ASN:HD21	1:B:364:SER:HB2	1.84	0.43
1:A:266:LYS:HA	1:A:271:ARG:HD3	2.00	0.43
1:B:153:ARG:HG3	1:B:154:GLN:N	2.33	0.43
1:A:240:ASP:OD1	1:A:240:ASP:C	2.62	0.42
1:B:222:GLU:O	1:B:239:GLU:N	2.52	0.42
1:A:434:ASN:ND2	1:A:435:ILE:H	2.17	0.42
1:B:33:GLN:O	1:B:34:ASP:HB2	2.19	0.42
1:A:460:VAL:O	1:A:464:VAL:HG23	2.18	0.42
1:A:434:ASN:ND2	1:A:435:ILE:N	2.68	0.42
1:A:439:ASN:ND2	1:A:442:ARG:HB2	2.33	0.42
1:B:40:ALA:O	1:B:220:SER:HA	2.20	0.42
1:B:479:GLU:HG2	2:B:802:FAD:O3'	2.19	0.42
1:A:123:SER:HB3	1:A:188:TYR:CE2	2.54	0.42
1:A:138:ASP:O	1:A:139:CYS:SG	2.77	0.42
1:A:255:GLN:HB2	1:A:480:HIS:CG	2.54	0.42
1:B:135:GLY:O	1:B:136:VAL:C	2.62	0.42
1:B:271:ARG:HG3	1:B:271:ARG:NH1	2.33	0.42
1:A:411:SER:OG	1:A:412:GLU:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:GLU:HA	1:A:230:PRO:HD2	1.92	0.42
1:B:99:TYR:O	1:B:106:ARG:HA	2.20	0.42
1:B:174:TRP:CE2	3:B:2842:DIA:HC71	2.55	0.42
1:A:171:LEU:HD23	1:A:171:LEU:HA	1.86	0.41
1:A:130:PHE:O	1:A:131:HIS:C	2.63	0.41
1:A:14:ILE:HD12	1:A:14:ILE:N	2.35	0.41
1:A:323:PHE:O	1:A:327:VAL:HG23	2.19	0.41
1:A:116:GLU:H	1:A:116:GLU:CD	2.28	0.41
1:B:36:LEU:HD23	1:B:216:TRP:O	2.21	0.41
1:B:466:MET:HG3	1:B:481:THR:HG22	2.03	0.41
1:A:329:ASN:O	1:A:330:ALA:C	2.64	0.41
1:A:300:GLU:O	1:A:432:LEU:HD12	2.21	0.40
1:B:47:ARG:NH2	1:B:254:PRO:HB3	2.36	0.40
1:B:363:LEU:HB2	1:B:371:SER:O	2.21	0.40
1:B:72:ASN:HD21	1:B:490:TYR:HB2	1.87	0.40
1:B:275:GLN:HE21	1:B:276:PRO:HA	1.80	0.40
1:B:376:MET:SD	1:B:381:THR:HA	2.61	0.40
1:B:24:ALA:O	1:B:25:SER:C	2.64	0.40
1:B:458:ASP:HA	1:B:459:PRO:HD2	1.91	0.40
1:B:466:MET:CG	1:B:481:THR:HG22	2.51	0.40
1:A:67:HIS:NE2	3:A:2841:DIA:HC12	2.36	0.40
1:B:109:HIS:O	1:B:109:HIS:CG	2.74	0.40
1:B:224:LYS:HB3	1:B:239:GLU:OE1	2.22	0.40
1:B:251:ILE:H	1:B:251:ILE:CD1	2.35	0.40
1:A:138:ASP:OD2	1:A:455:PRO:HA	2.21	0.40
1:B:349:THR:O	1:B:350:SER:CB	2.70	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	491/516 (95%)	450 (92%)	34 (7%)	7 (1%)	9	17
1	B	495/516 (96%)	430 (87%)	55 (11%)	10 (2%)	6	10
All	All	986/1032 (96%)	880 (89%)	89 (9%)	17 (2%)	7	13

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	200	ASN
1	A	230	PRO
1	A	231	SER
1	B	136	VAL
1	B	200	ASN
1	B	289	ILE
1	B	290	HIS
1	B	137	SER
1	B	421	ILE
1	A	131	HIS
1	B	211	SER
1	B	420	PRO
1	A	262	VAL
1	B	291	PHE
1	A	139	CYS
1	A	387	ILE
1	B	350	SER

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%)	403 (97%)	11 (3%)	39	67
1	B	403/457 (88%)	388 (96%)	15 (4%)	30	57
All	All	817/914 (89%)	791 (97%)	26 (3%)	34	62

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LEU
1	A	64	SER
1	A	201	TYR
1	A	218	LYS
1	A	221	CYS
1	A	239	GLU
1	A	267	ASN
1	A	335	GLU
1	A	377	GLN
1	A	388	ARG
1	A	484	ASP
1	B	22	LYS
1	B	36	LEU
1	B	72	ASN
1	B	124	LYS
1	B	173	LEU
1	B	177	LEU
1	B	201	TYR
1	B	221	CYS
1	B	227	THR
1	B	271	ARG
1	B	303	GLU
1	B	336	LEU
1	B	377	GLN
1	B	380	LEU
1	B	412	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	30	ASN
1	A	158	ASN
1	A	192	GLN
1	A	214	GLN
1	A	267	ASN
1	A	362	ASN
1	A	377	GLN
1	A	434	ASN
1	A	439	ASN
1	B	28	HIS
1	B	29	GLN
1	B	72	ASN

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Mol	Chain	Res	Type
1	B	96	ASN
1	B	109	HIS
1	B	191	HIS
1	B	192	GLN
1	B	215	ASN
1	B	259	ASN
1	B	275	GLN
1	B	329	ASN
1	B	362	ASN
1	B	377	GLN
1	B	382	ASN
1	B	434	ASN
1	B	439	ASN
1	B	470	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	DIA	A	2841	-	9,9,9	0.22	0	8,8,8	0.98	0
2	FAD	B	802	-	58,58,58	1.45	8 (13%)	85,89,89	2.20	9 (10%)
3	DIA	B	2842	-	9,9,9	0.22	0	8,8,8	0.85	0
2	FAD	A	801	-	58,58,58	1.50	10 (17%)	85,89,89	2.21	10 (11%)

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	DIA	A	2841	-	-	0/7/7/7	-
2	FAD	B	802	-	-	6/34/50/50	0/6/6/6
3	DIA	B	2842	-	-	0/7/7/7	-
2	FAD	A	801	-	-	5/34/50/50	0/6/6/6

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	FAD	PA-O3P	4.77	1.64	1.59
2	B	802	FAD	C4X-N5	4.11	1.39	1.30
2	A	801	FAD	C4X-N5	4.00	1.39	1.30
2	B	802	FAD	PA-O3P	3.76	1.63	1.59
2	A	801	FAD	C5B-C4B	2.85	1.60	1.51
2	A	801	FAD	C2A-N1A	2.73	1.38	1.33
2	B	802	FAD	C10-N1	2.72	1.38	1.33
2	B	802	FAD	C2A-N1A	2.65	1.38	1.33
2	A	801	FAD	P-O3P	2.61	1.62	1.59
2	B	802	FAD	C9A-C5X	2.48	1.45	1.41
2	B	802	FAD	P-O3P	2.44	1.62	1.59
2	A	801	FAD	C5'-C4'	2.40	1.55	1.51
2	B	802	FAD	C9-C8	2.31	1.42	1.39
2	A	801	FAD	C9A-C5X	2.20	1.44	1.41
2	A	801	FAD	C10-N1	2.20	1.37	1.33
2	B	802	FAD	C5B-C4B	2.19	1.58	1.51
2	A	801	FAD	O2'-C2'	-2.18	1.38	1.43
2	A	801	FAD	C2A-N3A	2.08	1.37	1.33

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FAD	O3P-PA-O1A	-12.56	72.93	110.70
2	B	802	FAD	O3P-PA-O1A	-12.25	73.85	110.70
2	A	801	FAD	O2A-PA-O5B	-7.62	73.04	107.57
2	A	801	FAD	O5B-PA-O1A	7.12	137.15	108.94
2	B	802	FAD	O2A-PA-O5B	-7.07	75.50	107.57
2	B	802	FAD	O5B-PA-O1A	6.87	136.16	108.94
2	B	802	FAD	O2A-PA-O3P	6.26	124.19	107.27
2	A	801	FAD	O2A-PA-O1A	-5.96	84.73	112.44
2	B	802	FAD	O2A-PA-O1A	-5.57	86.52	112.44
2	A	801	FAD	O2A-PA-O3P	4.72	120.03	107.27
2	B	802	FAD	O3'-C3'-C4'	-2.78	102.62	108.93
2	A	801	FAD	C9A-C5X-N5	-2.73	119.56	122.45
2	A	801	FAD	O4B-C1B-C2B	-2.69	100.86	106.62
2	B	802	FAD	C9A-C5X-N5	-2.60	119.70	122.45
2	B	802	FAD	O4B-C1B-C2B	-2.57	101.11	106.62
2	B	802	FAD	C4'-C3'-C2'	2.50	117.73	113.57
2	A	801	FAD	C4'-C3'-C2'	2.29	117.38	113.57
2	A	801	FAD	O3'-C3'-C4'	-2.26	103.79	108.93
2	A	801	FAD	N3A-C2A-N1A	-2.18	125.29	128.58

There are no chirality outliers.

All (11) 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	B	802	FAD	O4B-C4B-C5B-O5B
2	A	801	FAD	O4B-C4B-C5B-O5B
2	A	801	FAD	C3B-C4B-C5B-O5B
2	B	802	FAD	C3B-C4B-C5B-O5B
2	A	801	FAD	PA-O3P-P-O5'
2	B	802	FAD	PA-O3P-P-O5'
2	A	801	FAD	C4B-C5B-O5B-PA
2	B	802	FAD	C4B-C5B-O5B-PA
2	B	802	FAD	P-O3P-PA-O2A

There are no ring outliers.

4 monomers are involved in 7 short contacts:

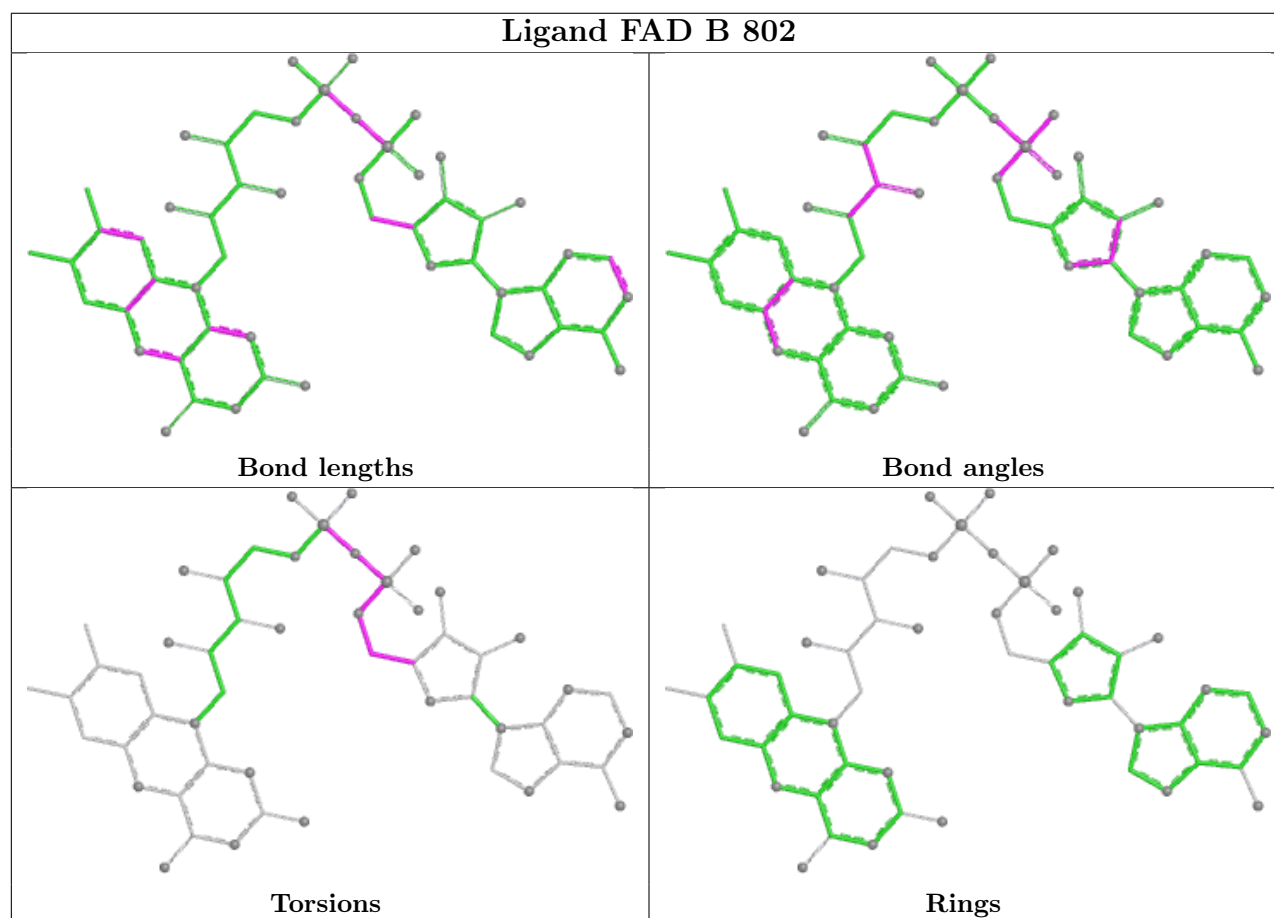
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2841	DIA	2	0
2	B	802	FAD	2	0

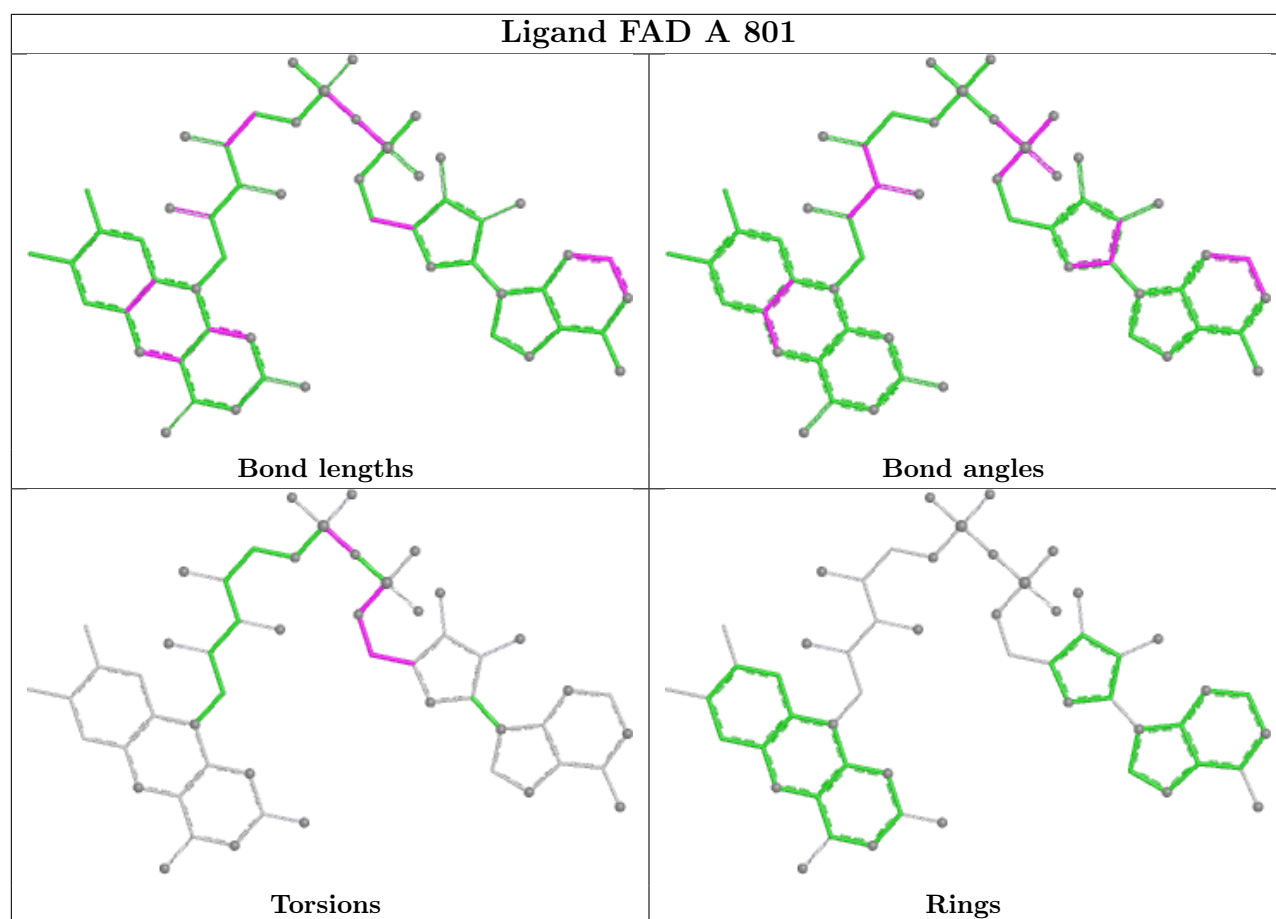
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2842	DIA	2	0
2	A	801	FAD	1	0

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	497/516 (96%)	0.34	20 (4%)	42 38	22, 39, 65, 90	0
1	B	499/516 (96%)	0.72	49 (9%)	13 11	27, 46, 74, 97	0
All	All	996/1032 (96%)	0.53	69 (6%)	23 20	22, 43, 73, 97	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	342	ARG	4.8
1	A	341	GLU	4.7
1	B	289	ILE	4.5
1	B	233	ASN	4.5
1	A	512	HIS	4.4
1	B	236	VAL	3.9
1	B	264	PRO	3.8
1	B	342	ARG	3.7
1	B	134	LEU	3.6
1	B	136	VAL	3.6
1	B	425	ALA	3.6
1	A	221	CYS	3.6
1	A	349	THR	3.6
1	A	509	LEU	3.6
1	B	238	CYS	3.5
1	B	231	SER	3.5
1	B	11	VAL	3.5
1	B	243	VAL	3.4
1	B	483	MET	3.3
1	B	229	GLU	3.3
1	A	238	CYS	3.2
1	B	221	CYS	3.2
1	B	349	THR	3.2
1	B	109	HIS	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	513	HIS	3.1
1	A	131	HIS	3.1
1	B	460	VAL	3.1
1	B	228	ARG	2.9
1	B	467	SER	2.9
1	B	412	GLU	2.9
1	A	333	LEU	2.8
1	A	230	PRO	2.8
1	B	245	ASN	2.8
1	B	249	VAL	2.8
1	B	265	GLU	2.8
1	A	340	LEU	2.7
1	B	340	LEU	2.7
1	B	191	HIS	2.7
1	A	338	SER	2.6
1	B	230	PRO	2.6
1	B	291	PHE	2.6
1	B	133	HIS	2.6
1	B	343	GLU	2.5
1	A	137	SER	2.5
1	A	330	ALA	2.5
1	B	511	HIS	2.5
1	A	427	ALA	2.5
1	B	36	LEU	2.4
1	B	332	ASN	2.4
1	B	226	ILE	2.4
1	B	90	PHE	2.4
1	A	350	SER	2.3
1	B	132	GLN	2.3
1	B	138	ASP	2.3
1	B	448	GLY	2.3
1	A	321	ASN	2.2
1	A	233	ASN	2.2
1	B	333	LEU	2.2
1	B	32	ILE	2.2
1	B	341	GLU	2.2
1	B	227	THR	2.2
1	B	420	PRO	2.1
1	A	191	HIS	2.1
1	B	212	PHE	2.1
1	B	463	VAL	2.1
1	B	244	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	336	LEU	2.0
1	B	135	GLY	2.0
1	B	421	ILE	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](#)

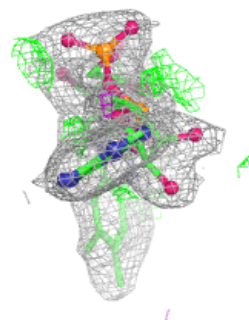
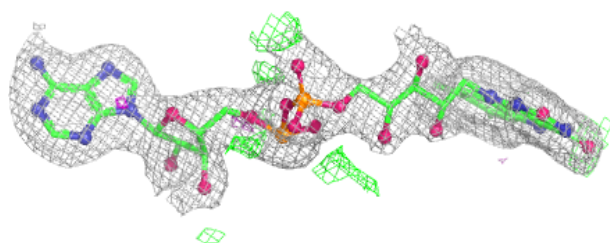
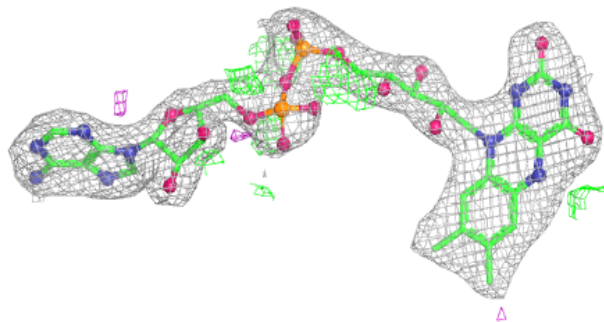
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	DIA	B	2842	10/10	0.79	0.26	72,75,78,78	0
3	DIA	A	2841	10/10	0.85	0.23	49,59,65,66	0
2	FAD	B	802	53/53	0.92	0.11	32,40,45,50	0
2	FAD	A	801	53/53	0.95	0.08	16,25,32,40	0

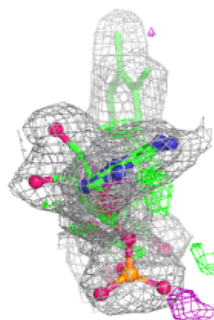
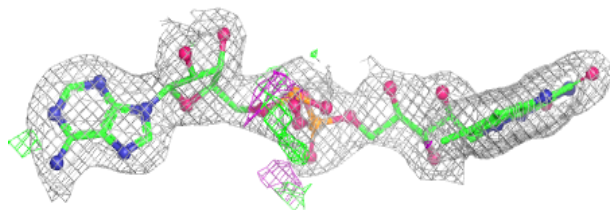
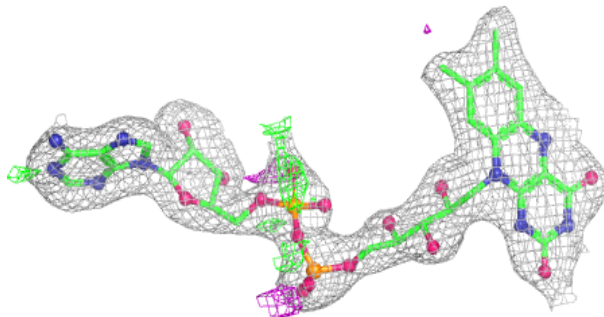
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD B 802:

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

**Electron density around FAD A 801:**

$2mF_o-DF_c$ (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.