



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 03:44 pm BST

PDB ID : 3CND
Title : Crystal structure of fms1 in complex with N1-AcSpermine
Authors : Huang, Q.; Hao, Q.
Deposited on : 2008-03-25
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 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 : 2.11
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) : 2.11

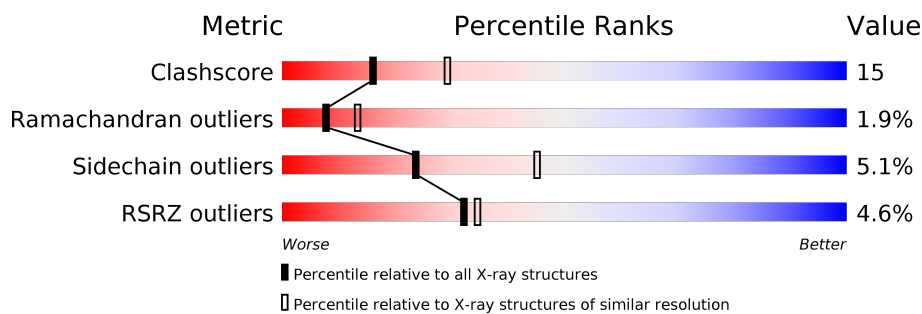
1 Overall quality at a glance

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))
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (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 on 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	<div> <div>2%</div> <div>72%</div> <div>22%</div> <div>• •</div> </div>
1	B	516	<div> <div>7%</div> <div>64%</div> <div>29%</div> <div>• •</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7906 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	B	499	Total	C	N	O	S	0	0	0
			3861	2434	671	734	22			
1	A	494	Total	C	N	O	S	0	0	0
			3863	2445	673	723	22			

There are 16 discrepancies between the modelled and reference sequences:

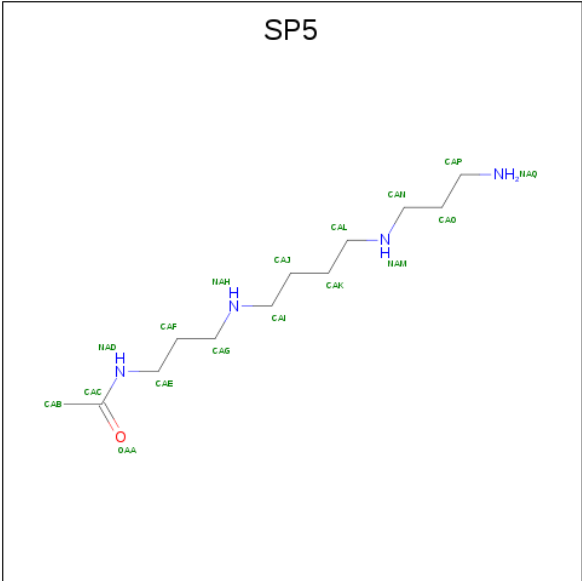
Chain	Residue	Modelled	Actual	Comment	Reference
B	509	LEU	-	EXPRESSION TAG	UNP P50264
B	510	GLU	-	EXPRESSION TAG	UNP P50264
B	511	HIS	-	EXPRESSION TAG	UNP P50264
B	512	HIS	-	EXPRESSION TAG	UNP P50264
B	513	HIS	-	EXPRESSION TAG	UNP P50264
B	514	HIS	-	EXPRESSION TAG	UNP P50264
B	515	HIS	-	EXPRESSION TAG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

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



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

- Molecule 3 is N-[3-({4-[(3-aminopropyl)amino]butyl}amino)propyl]acetamide (three-letter code: SP5) (formula: C₁₂H₂₈N₄O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			17	12	4	1		

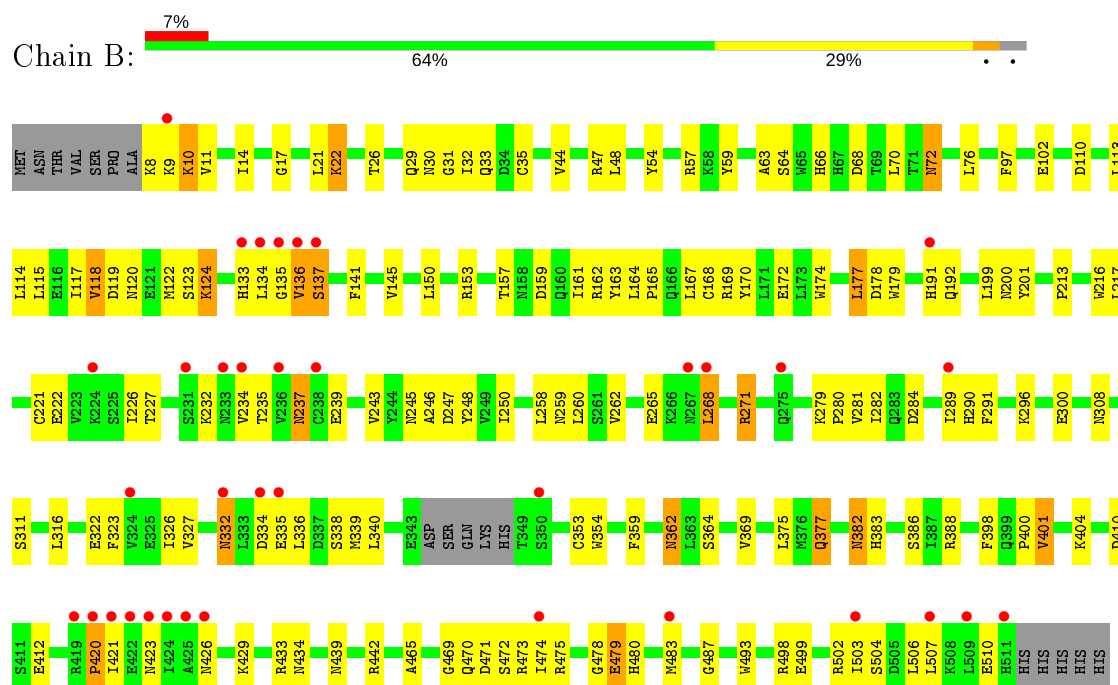
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	25	Total 25	O 25	0	0
4	A	34	Total 34	O 34	0	0

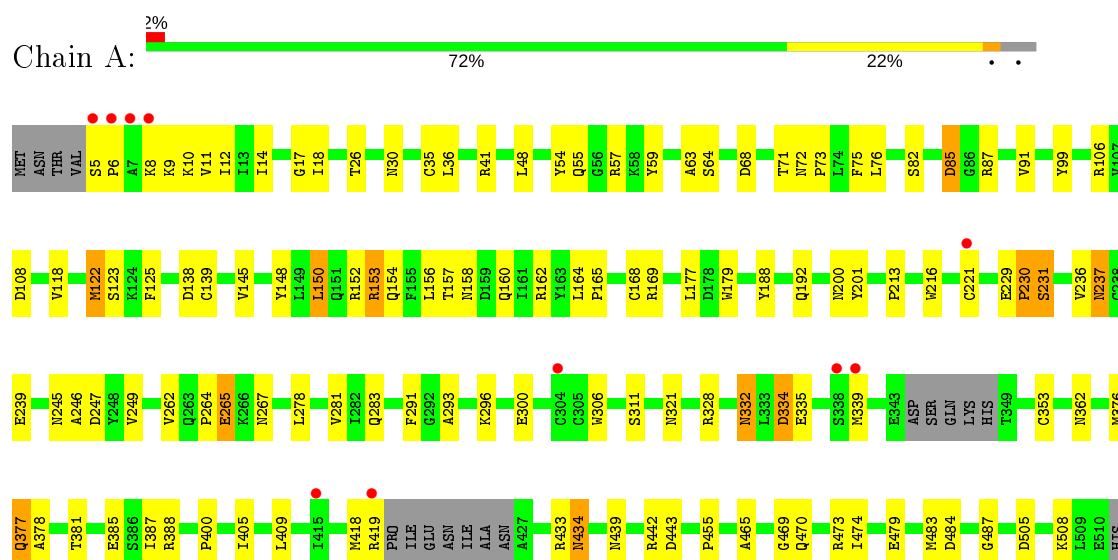
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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



HIS
HIS
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4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	102.58Å 215.85Å 118.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 41.84 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (50.00-2.50) 94.8 (41.84-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.78 (at 2.51Å)	Xtriage
Refinement program	CNS, REFMAC	Depositor
R, R_{free}	0.227 , 0.277 0.232 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	45.9	Xtriage
Anisotropy	0.709	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7906	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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: SP5, FAD

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	0.39	0/3942	0.62	0/5332
1	B	0.38	0/3939	0.61	1/5329 (0.0%)
All	All	0.38	0/7881	0.62	1/10661 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	66	HIS	N-CA-C	-5.17	97.03	111.00

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	A	3863	0	3726	93	0
1	B	3861	0	3669	130	0
2	A	53	0	31	2	0
2	B	53	0	31	3	0
3	B	17	0	28	5	0
4	A	34	0	0	0	0
4	B	25	0	0	1	0
All	All	7906	0	7485	223	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 15.

All (223) 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:353:CYS:SG	1:B:400:PRO:HG2	2.10	0.92
1:A:278:LEU:HA	1:A:470:GLN:HE22	1.36	0.88
1:A:85:ASP:OD2	1:A:87:ARG:HB2	1.74	0.87
1:A:236:VAL:C	1:A:237:ASN:HD22	1.79	0.85
1:B:382:ASN:ND2	1:B:382:ASN:H	1.71	0.85
1:A:377:GLN:NE2	1:A:377:GLN:H	1.79	0.81
1:B:68:ASP:HB3	1:B:192:GLN:HB2	1.64	0.79
1:B:11:VAL:HG22	1:B:248:TYR:HB2	1.65	0.79
1:B:282:ILE:HD13	1:B:465:ALA:HB1	1.68	0.76
1:A:278:LEU:HA	1:A:470:GLN:NE2	2.01	0.76
1:A:353:CYS:SG	1:A:400:PRO:HG2	2.24	0.76
1:A:158:ASN:ND2	1:A:328:ARG:HH21	1.85	0.75
1:B:470:GLN:HB3	1:B:474:ILE:HB	1.69	0.74
1:B:213:PRO:HB2	1:B:216:TRP:CD1	2.22	0.73
1:A:72:ASN:O	1:A:76:LEU:HG	1.91	0.70
1:B:237:ASN:HD22	1:B:243:VAL:HG22	1.56	0.70
1:B:48:LEU:CD2	1:B:63:ALA:HB3	2.23	0.68
1:B:141:PHE:O	1:B:145:VAL:HG23	1.94	0.68
1:B:375:LEU:HD21	3:B:803:SP5:HAL	1.76	0.67
1:A:229:GLU:O	1:A:231:SER:N	2.27	0.66
1:A:237:ASN:HD22	1:A:237:ASN:N	1.92	0.66
1:A:72:ASN:HB3	1:A:75:PHE:HB3	1.78	0.66
1:B:35:CYS:HB2	1:B:216:TRP:CZ3	2.31	0.66
1:B:382:ASN:H	1:B:382:ASN:HD22	1.41	0.66
1:A:54:TYR:CD2	1:A:55:GLN:HG3	2.34	0.62
1:A:48:LEU:CD2	1:A:63:ALA:HB3	2.29	0.62
1:A:157:THR:OG1	1:A:160:GLN:HG3	2.00	0.61
1:A:332:ASN:ND2	1:A:335:GLU:H	1.96	0.61
1:A:158:ASN:HD21	1:A:328:ARG:HE	1.47	0.61
1:A:311:SER:HA	1:A:362:ASN:HB3	1.81	0.61
1:A:162:ARG:NH1	1:A:321:ASN:OD1	2.30	0.60
1:B:135:GLY:O	1:B:137:SER:N	2.34	0.60
1:A:9:LYS:HE3	1:A:36:LEU:HD22	1.83	0.60
1:A:17:GLY:HA3	2:A:801:FAD:O2A	2.02	0.60
1:B:17:GLY:O	1:B:21:LEU:HG	2.01	0.60
1:B:9:LYS:O	1:B:246:ALA:HA	2.02	0.60
1:B:237:ASN:HD22	1:B:243:VAL:CG2	2.14	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:412:GLU:HB2	1:B:429:LYS:HD2	1.83	0.59
1:B:375:LEU:CD2	3:B:803:SP5:HAL	2.31	0.59
1:A:505:ASP:HA	1:A:508:LYS:HE2	1.84	0.59
1:B:32:ILE:HD12	1:B:507:LEU:HD13	1.84	0.59
1:A:306:TRP:HE1	1:A:362:ASN:HD21	1.51	0.59
1:B:322:GLU:O	1:B:326:ILE:HG13	2.02	0.58
1:B:377:GLN:NE2	1:B:377:GLN:H	2.02	0.58
1:B:507:LEU:HD23	1:B:510:GLU:OE2	2.04	0.58
1:A:377:GLN:HE21	1:A:377:GLN:H	1.49	0.58
1:B:8:LYS:N	1:B:245:ASN:HD21	2.02	0.57
1:B:412:GLU:HG3	1:B:429:LYS:CD	2.35	0.57
1:B:507:LEU:HD23	1:B:510:GLU:CD	2.25	0.57
1:B:133:HIS:O	1:B:134:LEU:HB2	2.03	0.57
1:B:281:VAL:HG13	1:B:282:ILE:N	2.20	0.57
1:A:213:PRO:O	1:A:216:TRP:HB2	2.04	0.57
1:A:41:ARG:NH1	1:A:443:ASP:OD2	2.38	0.56
1:A:213:PRO:HB2	1:A:216:TRP:CD1	2.40	0.56
1:B:237:ASN:ND2	1:B:243:VAL:HG22	2.20	0.56
1:A:68:ASP:HB3	1:A:192:GLN:HB2	1.87	0.56
1:A:150:LEU:O	1:A:153:ARG:HD2	2.05	0.56
1:B:118:VAL:HB	1:B:164:LEU:HD22	1.88	0.56
1:B:172:GLU:HG2	1:B:177:LEU:O	2.06	0.56
1:B:68:ASP:CB	1:B:192:GLN:HB2	2.35	0.56
1:B:469:GLY:HA3	1:B:475:ARG:NH1	2.21	0.56
1:B:271:ARG:HH11	1:B:271:ARG:HG3	1.71	0.55
1:B:133:HIS:HB3	1:B:136:VAL:HG13	1.88	0.55
1:B:323:PHE:O	1:B:327:VAL:HG23	2.06	0.55
1:B:63:ALA:HA	2:B:802:FAD:N5	2.21	0.55
1:A:64:SER:OG	1:A:296:LYS:NZ	2.39	0.55
1:B:165:PRO:O	1:B:169:ARG:HG3	2.08	0.54
1:A:278:LEU:CA	1:A:470:GLN:HE22	2.16	0.54
1:B:247:ASP:O	1:B:473:ARG:HD2	2.08	0.54
1:A:332:ASN:HD21	1:A:334:ASP:HB2	1.73	0.54
1:A:470:GLN:HB3	1:A:474:ILE:HB	1.89	0.54
1:A:5:SER:HB2	1:A:6:PRO:CD	2.38	0.54
1:A:5:SER:HB2	1:A:6:PRO:HD3	1.88	0.54
1:A:293:ALA:HB3	1:A:378:ALA:HB2	1.90	0.53
1:B:174:TRP:HZ2	3:B:803:SP5:HNAQ	1.55	0.53
1:B:72:ASN:C	1:B:72:ASN:HD22	2.12	0.53
1:B:30:ASN:ND2	1:B:504:SER:OG	2.40	0.53
1:A:26:THR:HG22	1:A:30:ASN:HD21	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:LYS:HE2	1:A:192:GLN:OE1	2.10	0.52
1:A:418:MET:O	1:A:419:ARG:HG3	2.09	0.52
1:B:26:THR:O	1:B:29:GLN:N	2.42	0.52
1:A:153:ARG:HG2	1:A:154:GLN:N	2.24	0.52
1:B:423:ASN:HB3	1:B:426:ASN:HD22	1.75	0.52
1:B:110:ASP:O	1:B:114:LEU:HD23	2.10	0.51
1:A:388:ARG:NH1	1:A:439:ASN:HB2	2.25	0.51
1:B:281:VAL:HG13	1:B:282:ILE:H	1.75	0.51
1:A:12:ILE:O	1:A:249:VAL:HA	2.10	0.51
1:B:311:SER:HA	1:B:362:ASN:HB3	1.93	0.51
1:A:8:LYS:HD2	1:A:245:ASN:HB3	1.93	0.51
1:A:332:ASN:ND2	1:A:334:ASP:HB2	2.24	0.51
1:B:227:THR:HB	1:B:235:THR:HB	1.92	0.51
1:B:326:ILE:HG23	1:B:339:MET:SD	2.51	0.51
1:B:412:GLU:CB	1:B:429:LYS:HD2	2.40	0.51
1:B:412:GLU:HG3	1:B:429:LYS:HD2	1.91	0.51
1:A:9:LYS:HG3	1:A:12:ILE:HD11	1.93	0.50
1:B:115:LEU:HB3	1:B:167:LEU:HD13	1.94	0.50
1:B:97:PHE:CE2	1:B:359:PHE:HE1	2.29	0.50
1:B:10:LYS:HG3	1:B:33:GLN:O	2.11	0.50
1:A:158:ASN:HD22	1:A:328:ARG:HH21	1.60	0.50
1:B:47:ARG:NE	2:B:802:FAD:O1A	2.42	0.49
1:A:237:ASN:ND2	1:A:237:ASN:N	2.60	0.49
1:B:439:ASN:HD21	1:B:442:ARG:NH1	2.09	0.49
1:B:22:LYS:HG2	1:B:493:TRP:CD1	2.48	0.49
1:B:412:GLU:HG3	1:B:429:LYS:HE3	1.93	0.49
1:A:387:ILE:HG13	1:A:387:ILE:O	2.13	0.49
1:A:164:LEU:N	1:A:165:PRO:HD2	2.27	0.49
1:A:54:TYR:O	1:A:57:ARG:HG3	2.12	0.49
1:B:102:GLU:O	1:B:404:LYS:HE2	2.13	0.49
1:B:332:ASN:OD1	1:B:334:ASP:HB2	2.12	0.49
1:B:338:SER:C	1:B:340:LEU:H	2.15	0.48
1:A:264:PRO:O	1:A:265:GLU:CB	2.60	0.48
1:B:338:SER:C	1:B:340:LEU:N	2.67	0.48
1:B:433:ARG:O	1:B:434:ASN:HB2	2.14	0.48
1:A:381:THR:O	1:A:385:GLU:HG3	2.14	0.48
1:B:157:THR:O	1:B:161:ILE:HG13	2.14	0.48
1:B:271:ARG:HG3	1:B:271:ARG:NH1	2.28	0.48
1:B:57:ARG:HD3	1:B:369:VAL:CG1	2.43	0.48
1:A:122:MET:HG2	1:A:148:TYR:CG	2.49	0.47
1:A:91:VAL:HG13	1:A:91:VAL:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:ASP:O	1:B:122:MET:HB3	2.14	0.47
1:A:63:ALA:HA	2:A:801:FAD:N5	2.29	0.47
1:A:18:ILE:HD12	1:A:48:LEU:HD11	1.97	0.47
1:A:138:ASP:O	1:A:139:CYS:SG	2.73	0.47
1:A:99:TYR:HD1	1:A:108:ASP:HB3	1.79	0.47
1:B:475:ARG:HB3	1:B:499:GLU:OE1	2.14	0.47
1:B:110:ASP:HB3	1:B:113:LEU:HB2	1.95	0.47
1:A:11:VAL:HB	1:A:35:CYS:SG	2.55	0.46
1:B:280:PRO:O	1:B:284:ASP:OD2	2.33	0.46
1:B:507:LEU:HA	1:B:510:GLU:HG3	1.96	0.46
1:A:479:GLU:OE1	1:A:487:GLY:HA2	2.16	0.46
1:B:17:GLY:HA3	2:B:802:FAD:H52A	1.98	0.46
1:B:117:ILE:HD12	1:B:117:ILE:N	2.31	0.46
1:B:507:LEU:HA	1:B:510:GLU:CG	2.45	0.46
1:B:159:ASP:O	1:B:163:TYR:HD1	1.99	0.46
1:A:14:ILE:N	1:A:14:ILE:HD12	2.30	0.46
1:A:418:MET:HG2	1:A:433:ARG:O	2.16	0.46
1:B:222:GLU:HB3	1:B:239:GLU:HB2	1.97	0.46
1:B:507:LEU:HA	1:B:510:GLU:CD	2.37	0.46
1:B:59:TYR:OH	1:B:300:GLU:HG2	2.16	0.46
1:B:439:ASN:HD21	1:B:442:ARG:HH11	1.64	0.45
1:A:376:MET:HG3	1:A:381:THR:OG1	2.17	0.45
1:A:138:ASP:OD2	1:A:455:PRO:HA	2.17	0.45
1:B:11:VAL:HA	1:B:248:TYR:O	2.17	0.45
1:B:308:ASN:OD1	1:B:364:SER:HB3	2.17	0.45
1:A:300:GLU:OE1	1:A:434:ASN:HB3	2.16	0.44
1:A:122:MET:CE	1:A:168:CYS:SG	3.05	0.44
1:B:10:LYS:HA	1:B:247:ASP:OD1	2.16	0.44
1:A:123:SER:HB3	1:A:188:TYR:CE2	2.53	0.44
1:A:474:ILE:N	1:A:474:ILE:HD12	2.32	0.44
1:B:279:LYS:HE2	1:B:470:GLN:HA	1.99	0.44
1:A:12:ILE:HB	1:A:249:VAL:HG12	1.98	0.44
1:A:9:LYS:O	1:A:246:ALA:HA	2.18	0.44
1:B:479:GLU:HB2	1:B:487:GLY:HA2	1.99	0.44
1:A:165:PRO:O	1:A:169:ARG:HG3	2.18	0.44
1:B:502:ARG:O	1:B:506:LEU:HD12	2.18	0.44
1:A:118:VAL:HG23	1:A:164:LEU:HD13	1.99	0.44
1:B:162:ARG:HD3	1:B:163:TYR:CE1	2.52	0.44
1:B:479:GLU:HB2	1:B:487:GLY:CA	2.48	0.44
1:A:334:ASP:N	1:A:334:ASP:OD1	2.51	0.44
1:B:439:ASN:ND2	1:B:442:ARG:HH11	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:338:SER:O	1:B:340:LEU:N	2.51	0.43
1:B:70:LEU:HB2	1:B:192:GLN:O	2.18	0.43
1:A:122:MET:CE	1:A:145:VAL:HG22	2.48	0.43
1:B:234:VAL:N	1:B:246:ALA:O	2.51	0.43
1:B:388:ARG:NE	1:B:439:ASN:HB2	2.33	0.43
1:B:120:ASN:O	1:B:123:SER:HB3	2.18	0.43
1:B:281:VAL:HG13	1:B:282:ILE:HG12	1.99	0.43
1:B:48:LEU:HD23	1:B:63:ALA:HB3	1.99	0.43
1:A:281:VAL:HG12	1:A:465:ALA:HB2	1.99	0.43
1:B:153:ARG:NH1	1:B:327:VAL:O	2.52	0.43
1:B:250:ILE:HD11	1:B:503:ILE:HD12	2.00	0.43
1:B:72:ASN:O	1:B:76:LEU:HG	2.19	0.43
1:B:222:GLU:O	1:B:239:GLU:N	2.52	0.43
1:A:247:ASP:O	1:A:473:ARG:HD2	2.19	0.43
1:A:59:TYR:OH	1:A:300:GLU:HG2	2.19	0.43
1:A:82:SER:HA	1:A:85:ASP:OD1	2.19	0.43
1:B:262:VAL:O	1:B:262:VAL:HG23	2.18	0.43
1:B:420:PRO:O	1:B:421:ILE:HD13	2.18	0.43
1:B:44:VAL:CG1	1:B:217:LEU:HD21	2.49	0.43
1:A:473:ARG:NH1	1:A:473:ARG:HG3	2.34	0.42
1:B:335:GLU:O	1:B:339:MET:HB2	2.19	0.42
1:B:383:HIS:O	1:B:386:SER:HB2	2.19	0.42
1:B:473:ARG:HG3	1:B:473:ARG:HH11	1.84	0.42
1:B:35:CYS:HB2	1:B:216:TRP:CH2	2.53	0.42
1:B:362:ASN:ND2	1:B:364:SER:H	2.17	0.42
1:B:164:LEU:N	1:B:165:PRO:HD2	2.34	0.42
1:B:362:ASN:C	1:B:362:ASN:ND2	2.71	0.42
1:A:230:PRO:O	1:A:231:SER:HB3	2.20	0.42
1:A:26:THR:HG22	1:A:30:ASN:ND2	2.34	0.42
1:A:377:GLN:N	1:A:377:GLN:NE2	2.59	0.42
1:B:150:LEU:HD13	1:B:336:LEU:HD12	2.02	0.42
1:B:199:LEU:HD23	1:B:199:LEU:HA	1.83	0.42
1:B:499:GLU:O	1:B:503:ILE:HG13	2.20	0.42
1:A:262:VAL:CG1	1:A:283:GLN:HG2	2.49	0.42
1:B:8:LYS:N	1:B:245:ASN:ND2	2.68	0.42
1:A:150:LEU:HD12	1:A:150:LEU:HA	1.80	0.42
1:B:478:GLY:O	1:B:480:HIS:N	2.53	0.42
1:A:122:MET:HE3	1:A:168:CYS:SG	2.60	0.42
1:B:248:TYR:CE2	1:B:472:SER:O	2.73	0.42
1:A:335:GLU:HG2	1:A:339:MET:CE	2.50	0.41
1:A:405:ILE:HG22	1:A:409:LEU:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:TYR:C	1:B:170:TYR:CD1	2.94	0.41
1:B:191:HIS:HB2	4:B:804:HOH:O	2.19	0.41
3:B:803:SP5:HALA	3:B:803:SP5:HAO	1.74	0.41
1:A:335:GLU:HG2	1:A:339:MET:HE1	2.02	0.41
1:B:14:ILE:CD1	1:B:226:ILE:HD11	2.50	0.41
1:A:152:ARG:O	1:A:156:LEU:HG	2.21	0.41
1:A:264:PRO:O	1:A:265:GLU:HB2	2.20	0.41
1:A:99:TYR:O	1:A:106:ARG:HA	2.21	0.41
1:B:258:LEU:C	1:B:260:LEU:H	2.24	0.41
1:B:54:TYR:O	1:B:57:ARG:HG3	2.19	0.41
1:A:473:ARG:HH11	1:A:473:ARG:HG3	1.86	0.41
1:B:174:TRP:CZ2	3:B:803:SP5:NAQ	2.88	0.41
1:A:71:THR:O	1:A:73:PRO:HD3	2.21	0.41
1:B:32:ILE:HD12	1:B:507:LEU:CD1	2.50	0.41
1:B:398:PHE:O	1:B:401:VAL:HG23	2.21	0.41
1:B:479:GLU:OE1	1:B:487:GLY:HA2	2.21	0.41
1:A:122:MET:O	1:A:125:PHE:HB3	2.21	0.41
1:B:336:LEU:C	1:B:338:SER:N	2.73	0.41
1:B:22:LYS:HG2	1:B:493:TRP:NE1	2.35	0.41
1:B:64:SER:OG	1:B:296:LYS:NZ	2.54	0.41
1:B:102:GLU:HB2	1:B:354:TRP:CE2	2.56	0.40
1:B:268:LEU:HA	1:B:268:LEU:HD23	1.85	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	488/516 (95%)	444 (91%)	39 (8%)	5 (1%)	15	28
1	B	495/516 (96%)	437 (88%)	44 (9%)	14 (3%)	5	7
All	All	983/1032 (95%)	881 (90%)	83 (8%)	19 (2%)	8	13

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	136	VAL
1	B	200	ASN
1	B	232	LYS
1	B	265	GLU
1	B	289	ILE
1	B	291	PHE
1	A	231	SER
1	A	265	GLU
1	B	498	ARG
1	A	230	PRO
1	B	31	GLY
1	B	479	GLU
1	A	469	GLY
1	B	137	SER
1	B	259	ASN
1	B	290	HIS
1	B	420	PRO
1	B	471	ASP
1	A	200	ASN

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	412/457 (90%)	392 (95%)	20 (5%)	25	47
1	B	408/457 (89%)	386 (95%)	22 (5%)	22	42
All	All	820/914 (90%)	778 (95%)	42 (5%)	24	45

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	10	LYS
1	B	22	LYS
1	B	72	ASN
1	B	118	VAL

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Mol	Chain	Res	Type
1	B	124	LYS
1	B	168	CYS
1	B	177	LEU
1	B	178	ASP
1	B	179	TRP
1	B	201	TYR
1	B	221	CYS
1	B	237	ASN
1	B	268	LEU
1	B	271	ARG
1	B	316	LEU
1	B	332	ASN
1	B	362	ASN
1	B	377	GLN
1	B	382	ASN
1	B	401	VAL
1	B	410	ASP
1	B	483	MET
1	A	10	LYS
1	A	85	ASP
1	A	122	MET
1	A	150	LEU
1	A	153	ARG
1	A	177	LEU
1	A	179	TRP
1	A	201	TYR
1	A	221	CYS
1	A	237	ASN
1	A	239	GLU
1	A	267	ASN
1	A	291	PHE
1	A	332	ASN
1	A	334	ASP
1	A	377	GLN
1	A	434	ASN
1	A	442	ARG
1	A	483	MET
1	A	484	ASP

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

Mol	Chain	Res	Type
1	B	29	GLN
1	B	30	ASN
1	B	72	ASN
1	B	191	HIS
1	B	210	GLN
1	B	237	ASN
1	B	255	GLN
1	B	283	GLN
1	B	329	ASN
1	B	362	ASN
1	B	377	GLN
1	B	382	ASN
1	B	426	ASN
1	B	434	ASN
1	B	439	ASN
1	A	28	HIS
1	A	30	ASN
1	A	33	GLN
1	A	109	HIS
1	A	158	ASN
1	A	237	ASN
1	A	267	ASN
1	A	332	ASN
1	A	362	ASN
1	A	377	GLN
1	A	434	ASN
1	A	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 carbohydrates in this entry.

5.6 Ligand geometry

3 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
2	FAD	B	802	-	51,58,58	1.61	8 (15%)	60,89,89	2.16	11 (18%)
2	FAD	A	801	-	51,58,58	1.62	10 (19%)	60,89,89	2.14	10 (16%)
3	SP5	B	803	-	16,16,16	0.50	0	16,16,16	0.78	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
2	FAD	B	802	-	-	4/30/50/50	0/6/6/6
2	FAD	A	801	-	-	10/30/50/50	0/6/6/6
3	SP5	B	803	-	-	8/14/14/14	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	802	FAD	C4X-N5	5.03	1.40	1.33
2	B	802	FAD	C4X-C10	4.46	1.43	1.38
2	A	801	FAD	C4-N3	4.39	1.40	1.33
2	B	802	FAD	C9A-N10	3.96	1.43	1.38
2	A	801	FAD	C9A-N10	3.96	1.43	1.38
2	B	802	FAD	C10-N1	3.89	1.38	1.33
2	A	801	FAD	C4X-N5	3.86	1.38	1.33
2	A	801	FAD	C4X-C10	3.54	1.42	1.38
2	A	801	FAD	C10-N1	3.21	1.37	1.33
2	B	802	FAD	C4-N3	2.78	1.37	1.33
2	B	802	FAD	C2A-N1A	2.54	1.38	1.33
2	A	801	FAD	C2A-N3A	2.54	1.36	1.32
2	A	801	FAD	C2A-N1A	2.49	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	802	FAD	C2A-N3A	2.42	1.36	1.32
2	A	801	FAD	C5X-N5	2.41	1.39	1.35
2	A	801	FAD	C1'-N10	2.40	1.50	1.48
2	A	801	FAD	PA-O2A	-2.21	1.45	1.55
2	B	802	FAD	C5X-N5	2.12	1.38	1.35

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	802	FAD	C4-N3-C2	11.05	124.47	115.14
2	A	801	FAD	C4-N3-C2	10.25	123.80	115.14
2	A	801	FAD	C1'-N10-C9A	4.72	122.00	118.29
2	A	801	FAD	C4X-C4-N3	-4.70	117.00	123.43
2	B	802	FAD	C4X-N5-C5X	4.64	121.40	116.77
2	B	802	FAD	O2A-PA-O1A	4.56	134.76	112.24
2	B	802	FAD	C4X-C4-N3	-4.48	117.31	123.43
2	A	801	FAD	O4B-C1B-C2B	-4.46	100.40	106.93
2	A	801	FAD	O2A-PA-O1A	4.26	133.30	112.24
2	B	802	FAD	O4B-C1B-C2B	-4.25	100.71	106.93
2	A	801	FAD	C4X-N5-C5X	4.21	120.97	116.77
2	B	802	FAD	C1'-N10-C9A	3.18	120.79	118.29
2	A	801	FAD	O3'-C3'-C4'	-2.55	102.66	108.81
2	A	801	FAD	C4-C4X-C10	-2.47	118.31	119.95
2	A	801	FAD	C5A-C6A-N6A	2.38	123.97	120.35
2	B	802	FAD	C5A-C6A-N6A	2.35	123.92	120.35
2	B	802	FAD	O3'-C3'-C4'	-2.31	103.22	108.81
2	B	802	FAD	P-O3P-PA	2.14	140.18	132.83
2	A	801	FAD	N3A-C2A-N1A	-2.12	125.37	128.68
2	B	802	FAD	N3A-C2A-N1A	-2.07	125.44	128.68
2	B	802	FAD	C4-C4X-C10	-2.01	118.62	119.95

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	802	FAD	C5B-O5B-PA-O2A
2	A	801	FAD	C5B-O5B-PA-O2A
3	B	803	SP5	CAB-CAC-NAD-CAE
3	B	803	SP5	OAA-CAC-NAD-CAE
3	B	803	SP5	CAJ-CAK-CAL-NAM
3	B	803	SP5	CAE-CAF-CAG-NAH
3	B	803	SP5	NAH-CAI-CAJ-CAK

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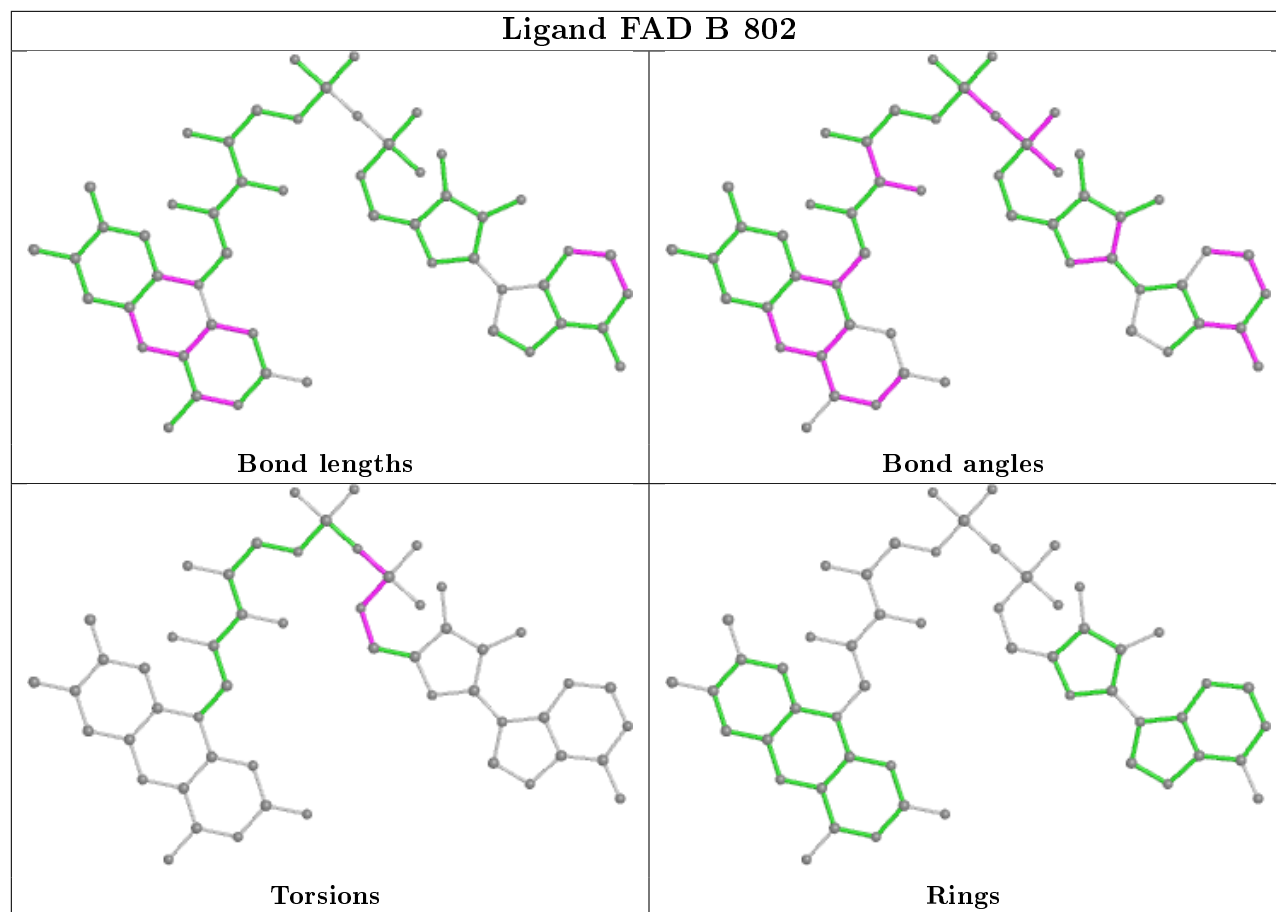
Mol	Chain	Res	Type	Atoms
2	A	801	FAD	C2'-C3'-C4'-O4'
2	A	801	FAD	O4B-C4B-C5B-O5B
3	B	803	SP5	CAJ-CAI-NAH-CAG
2	B	802	FAD	C5B-O5B-PA-O3P
2	A	801	FAD	C5B-O5B-PA-O3P
2	A	801	FAD	C4B-C5B-O5B-PA
3	B	803	SP5	CAF-CAG-NAH-CAI
2	B	802	FAD	P-O3P-PA-O2A
2	B	802	FAD	C4B-C5B-O5B-PA
3	B	803	SP5	CAN-CAO-CAP-NAQ
2	A	801	FAD	C3B-C4B-C5B-O5B
2	A	801	FAD	C2'-C3'-C4'-C5'
2	A	801	FAD	P-O3P-PA-O2A
2	A	801	FAD	O3'-C3'-C4'-C5'
2	A	801	FAD	O3'-C3'-C4'-O4'

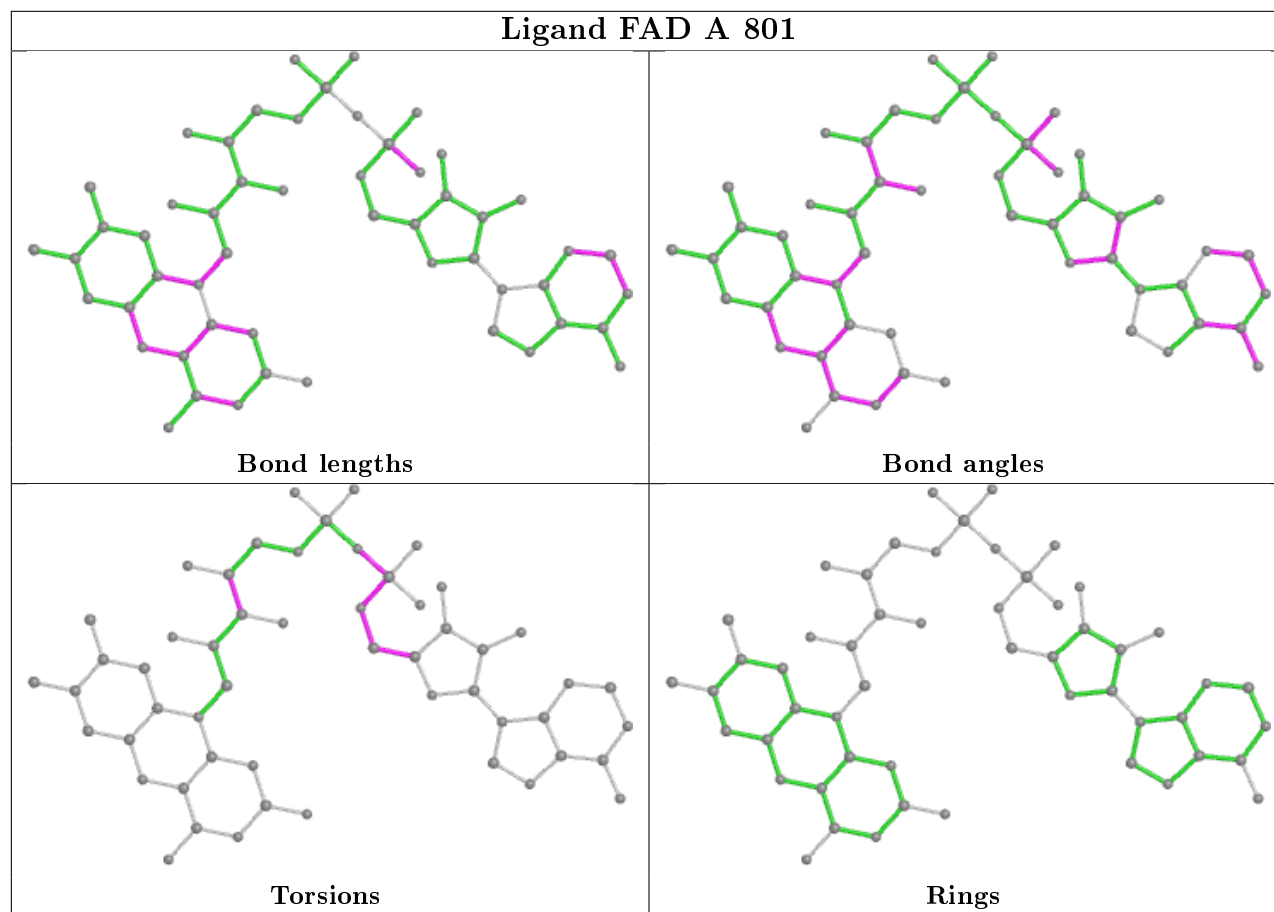
There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	802	FAD	3	0
2	A	801	FAD	2	0
3	B	803	SP5	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%)	0.03	10 (2%) 65 68	27, 46, 69, 88	0
1	B	499/516 (96%)	0.22	36 (7%) 15 16	27, 50, 87, 105	0
All	All	993/1032 (96%)	0.13	46 (4%) 32 34	27, 48, 82, 105	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	135	GLY	9.4
1	B	134	LEU	8.5
1	B	425	ALA	6.5
1	B	136	VAL	6.4
1	B	423	ASN	6.4
1	A	5	SER	5.5
1	B	426	ASN	5.4
1	B	424	ILE	5.2
1	B	511	HIS	5.1
1	B	422	GLU	5.0
1	B	509	LEU	5.0
1	A	6	PRO	4.6
1	B	507	LEU	4.2
1	B	335	GLU	3.7
1	B	133	HIS	3.5
1	A	415	ILE	3.1
1	B	421	ILE	3.0
1	A	419	ARG	3.0
1	A	7	ALA	2.9
1	B	238	CYS	2.7
1	B	275	GLN	2.7
1	B	289	ILE	2.7
1	B	268	LEU	2.6
1	B	324	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	334	ASP	2.5
1	B	191	HIS	2.5
1	B	420	PRO	2.5
1	B	474	ILE	2.4
1	B	350	SER	2.3
1	B	231	SER	2.3
1	B	224	LYS	2.3
1	B	234	VAL	2.3
1	B	267	ASN	2.2
1	A	338	SER	2.2
1	B	137	SER	2.2
1	B	236	VAL	2.2
1	A	8	LYS	2.2
1	A	221	CYS	2.2
1	B	419	ARG	2.1
1	B	332	ASN	2.1
1	B	233	ASN	2.0
1	B	9	LYS	2.0
1	A	304	CYS	2.0
1	A	339	MET	2.0
1	B	503	ILE	2.0
1	B	483	MET	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 carbohydrates in this entry.

6.4 Ligands [i](#)

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

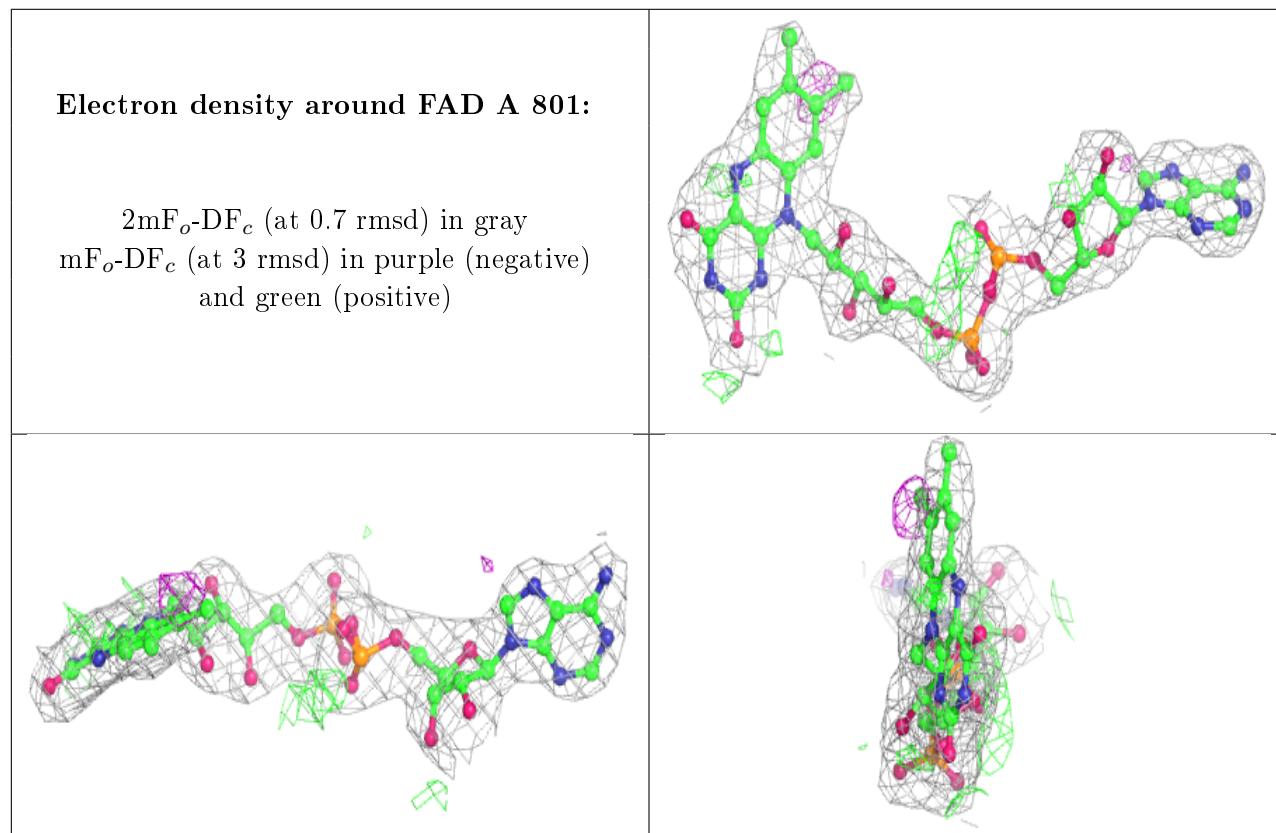
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SP5	B	803	17/17	0.83	0.31	56,61,70,70	0

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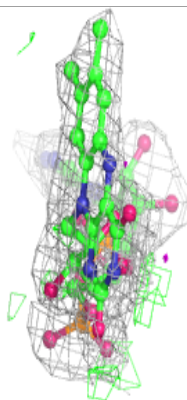
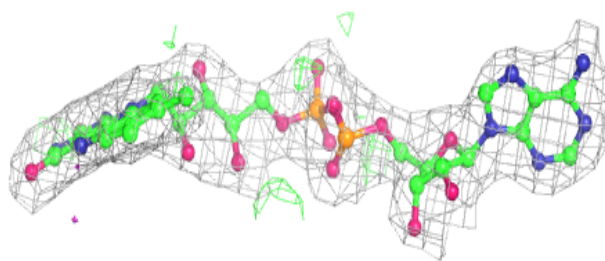
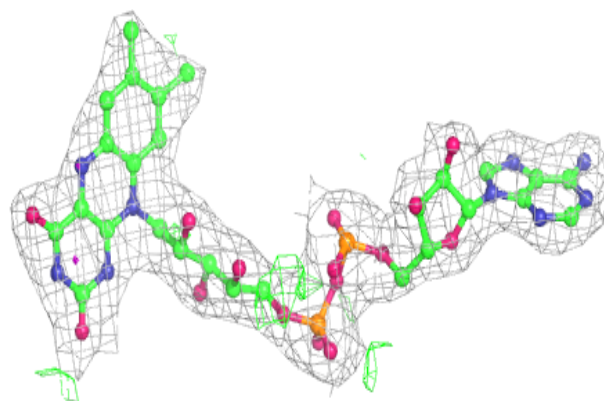
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	FAD	A	801	53/53	0.96	0.19	28,35,42,44	0
2	FAD	B	802	53/53	0.96	0.15	29,39,47,48	0

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)



6.5 Other polymers [i](#)

There are no such residues in this entry.