



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:41 AM UTC

PDB ID : 6GAS / pdb_00006gas
Title : Crystal structure of oxidised ferredoxin/flavodoxin NADP+ oxidoreductase 2 (FNR2) from Bacillus cereus
Authors : Gudim, I.; Hersleth, H.-P.
Deposited on : 2018-04-12
Resolution : 2.40 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

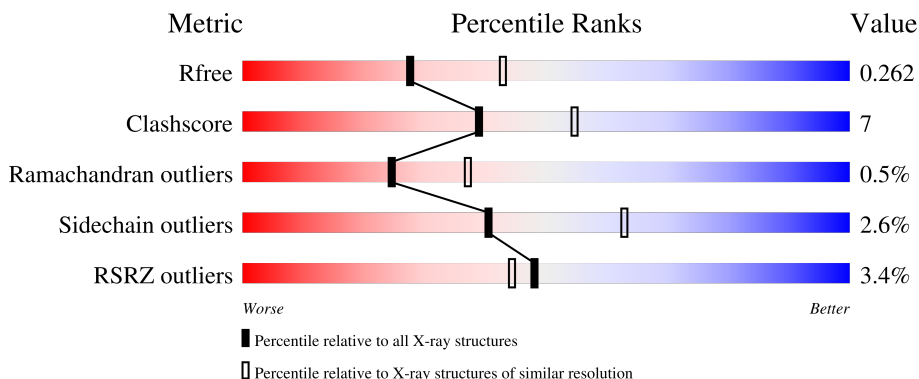
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	331	4% 82% 14% ..
1	B	331	2% 81% 16% ..
1	C	331	5% 83% 15% ..
1	D	331	3% 83% 15% ..

2 Entry composition [i](#)

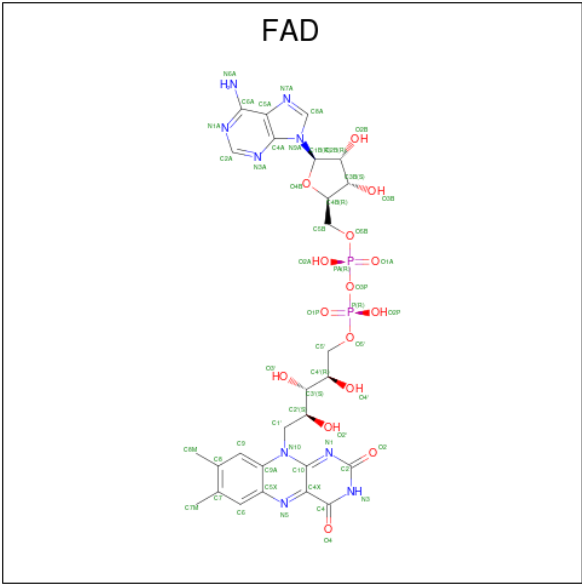
There are 4 unique types of molecules in this entry. The entry contains 10632 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ferredoxin–NADP reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total	C	N	O	S	0	0	0
			2547	1626	426	483	12			
1	B	327	Total	C	N	O	S	0	0	0
			2556	1631	427	486	12			
1	C	327	Total	C	N	O	S	0	0	0
			2556	1631	427	486	12			
1	D	326	Total	C	N	O	S	0	0	0
			2547	1626	426	483	12			

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

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

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		
3	B	1	Total	Na	0	0
			1	1		
3	C	1	Total	Na	0	0
			1	1		
3	D	1	Total	Na	0	0
			1	1		

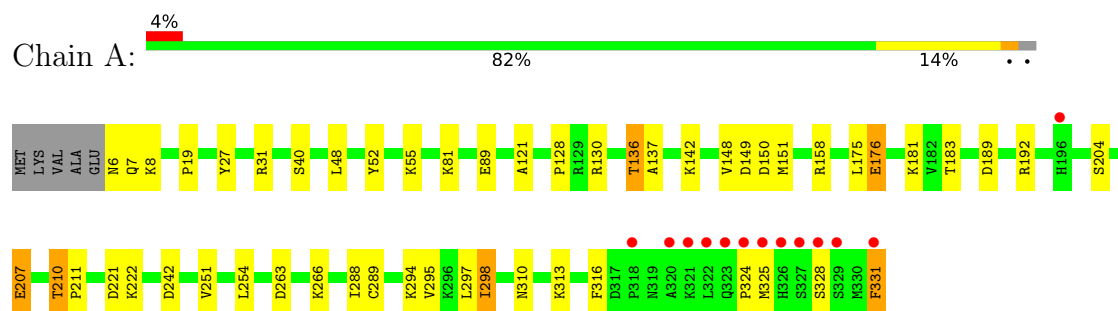
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	59	Total	O	0	0
			59	59		
4	B	62	Total	O	0	0
			62	62		
4	C	45	Total	O	0	0
			45	45		
4	D	44	Total	O	0	0
			44	44		

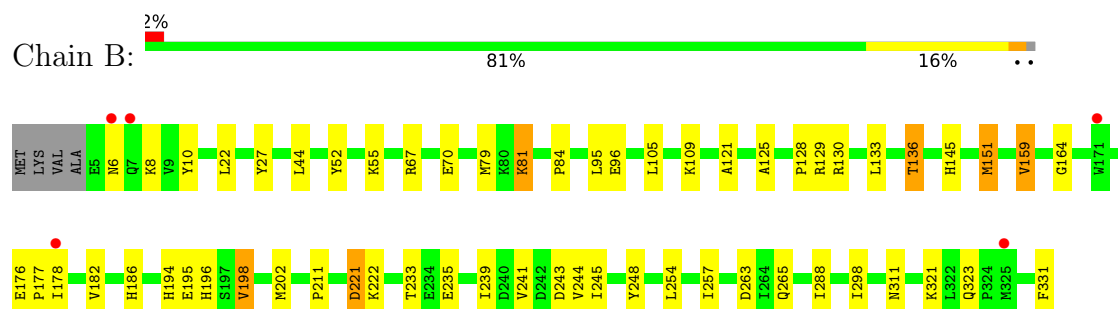
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

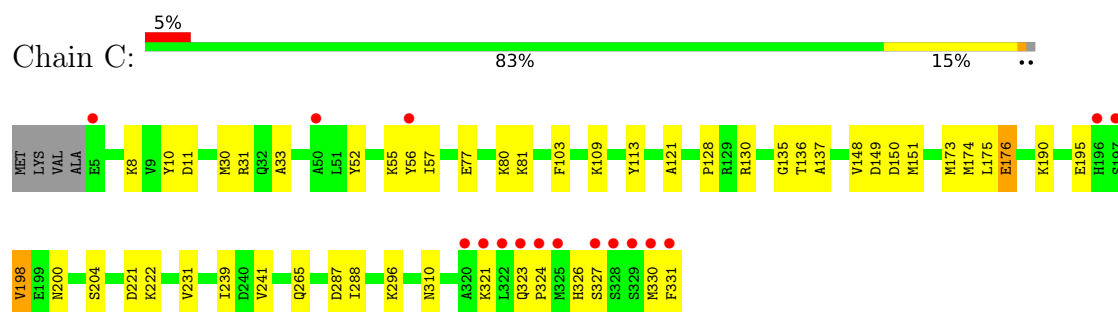
- Molecule 1: Ferredoxin–NADP reductase



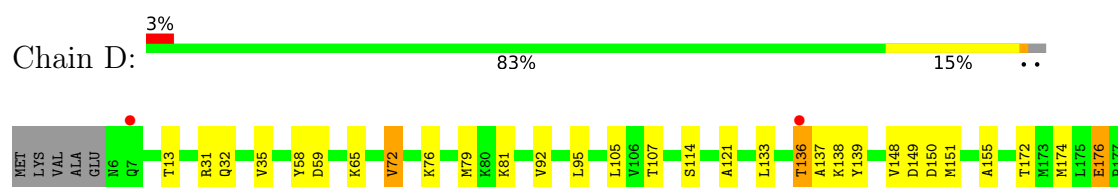
- Molecule 1: Ferredoxin–NADP reductase

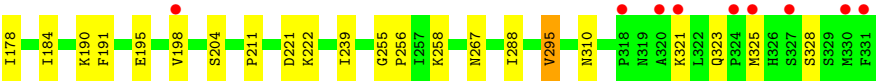


- Molecule 1: Ferredoxin–NADP reductase



- Molecule 1: Ferredoxin–NADP reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	109.66Å 109.66Å 242.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.60 – 2.40 46.60 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.60-2.40) 99.9 (46.60-2.40)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.207 , 0.262 0.210 , 0.262	Depositor DCC
R_{free} test set	3385 reflections (4.18%)	wwPDB-VP
Wilson B-factor (Å ²)	50.1	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 32.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10632	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.25% 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: NA, 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	0.50	0/2598	0.71	0/3511
1	B	0.57	0/2607	0.78	0/3523
1	C	0.48	0/2607	0.68	0/3523
1	D	0.48	0/2598	0.67	0/3511
All	All	0.51	0/10410	0.71	0/14068

There are no bond length outliers.

There are no bond angle outliers.

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	2547	0	2553	33	0
1	B	2556	0	2559	40	0
1	C	2556	0	2559	41	0
1	D	2547	0	2553	32	0
2	A	53	0	31	2	0
2	B	53	0	31	4	0
2	C	53	0	31	1	0
2	D	53	0	31	1	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	59	0	0	4	0
4	B	62	0	0	3	0
4	C	45	0	0	0	0
4	D	44	0	0	1	0
All	All	10632	0	10348	142	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 7.

All (142) 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:159:VAL:HG12	1:B:182:VAL:HG23	1.64	0.79
1:B:151:MET:HG2	1:B:178:ILE:HD12	1.64	0.78
1:B:133:LEU:O	1:B:136:THR:HG23	1.85	0.75
1:C:52:TYR:HE2	1:C:57:ILE:HD11	1.53	0.73
2:B:1001:FAD:O2'	2:B:1001:FAD:O4'	2.04	0.73
1:C:81:LYS:HZ3	1:C:331:PHE:HZ	1.37	0.71
1:B:67:ARG:HB3	1:B:70:GLU:HG2	1.72	0.71
1:A:7:GLN:HE22	1:C:231:VAL:HG12	1.57	0.69
1:B:221:ASP:OD1	1:B:222:LYS:N	2.26	0.69
1:C:151:MET:HE2	1:C:175:LEU:HD21	1.74	0.69
1:A:298:ILE:HG22	2:A:1001:FAD:O2	1.93	0.69
1:C:148:VAL:HG21	1:C:151:MET:HE3	1.75	0.69
1:A:31:ARG:NH1	1:A:310:ASN:OD1	2.26	0.69
1:D:121:ALA:HB3	1:D:288:ILE:HD12	1.73	0.69
1:C:151:MET:HG2	1:C:174:MET:HG2	1.76	0.68
1:B:194:HIS:HD2	1:B:196:HIS:H	1.41	0.67
1:D:76:LYS:HA	1:D:79:MET:HE2	1.77	0.66
1:D:149:ASP:OD1	1:D:150:ASP:N	2.25	0.65
1:A:183:THR:HG23	1:A:207:GLU:HG2	1.80	0.64
1:B:176:GLU:HG3	1:B:177:PRO:HD3	1.79	0.64
1:D:221:ASP:OD1	1:D:222:LYS:N	2.31	0.63
1:C:77:GLU:HA	1:C:80:LYS:HE3	1.78	0.63
1:D:148:VAL:HG11	1:D:151:MET:SD	2.39	0.63
1:A:148:VAL:HG11	1:A:151:MET:SD	2.39	0.61
1:D:136:THR:HG23	1:D:137:ALA:H	1.66	0.61
1:D:321:LYS:HE2	1:D:323:GLN:O	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:326:HIS:O	1:C:330:MET:HG3	2.02	0.60
1:C:30:MET:SD	1:C:330:MET:HE1	2.43	0.59
1:B:194:HIS:CD2	1:B:196:HIS:H	2.19	0.58
1:C:121:ALA:HB3	1:C:288:ILE:HD12	1.85	0.58
1:C:195:GLU:O	1:C:198:VAL:HG22	2.04	0.58
1:C:148:VAL:HG21	1:C:151:MET:CE	2.34	0.58
1:A:189:ASP:OD1	1:A:210:THR:HG23	2.04	0.57
1:D:136:THR:HG23	1:D:137:ALA:N	2.19	0.57
1:A:176:GLU:OE1	1:A:204:SER:HB2	2.05	0.57
1:D:58:TYR:CE1	1:D:65:LYS:HG3	2.40	0.56
1:B:128:PRO:HG2	1:B:130:ARG:HH22	1.70	0.56
1:D:190:LYS:HE2	1:D:191:PHE:O	2.07	0.55
1:D:13:THR:HG23	1:D:114:SER:OG	2.06	0.54
1:C:176:GLU:OE2	1:C:204:SER:HB2	2.06	0.54
1:B:52:TYR:HB2	1:B:55:LYS:HG2	1.89	0.54
1:A:121:ALA:HB3	1:A:288:ILE:HD12	1.89	0.53
1:B:159:VAL:HG22	1:B:243:ASP:HB2	1.90	0.53
1:D:13:THR:HG21	4:D:1104:HOH:O	2.08	0.53
1:B:254:LEU:HD11	1:B:288:ILE:HD11	1.90	0.53
1:D:138:LYS:HE3	1:D:139:TYR:CZ	2.44	0.53
1:C:128:PRO:HG2	1:C:130:ARG:NH2	2.23	0.53
1:B:321:LYS:HE2	1:B:323:GLN:O	2.09	0.53
1:B:233:THR:OG1	1:B:235:GLU:HG2	2.09	0.52
1:A:136:THR:OG1	1:A:137:ALA:N	2.41	0.52
1:A:221:ASP:OD1	1:A:222:LYS:N	2.42	0.52
1:C:221:ASP:OD1	1:C:222:LYS:N	2.41	0.52
1:B:109:LYS:HE2	4:B:1158:HOH:O	2.09	0.52
1:C:149:ASP:OD1	1:C:150:ASP:N	2.36	0.51
1:A:149:ASP:OD1	1:A:150:ASP:N	2.38	0.51
1:A:40:SER:O	1:A:89:GLU:HA	2.10	0.51
1:A:6:ASN:HB3	1:A:8:LYS:H	1.77	0.50
1:B:195:GLU:HA	1:B:198:VAL:HG13	1.93	0.50
1:C:331:PHE:HE2	1:D:58:TYR:CD1	2.29	0.49
1:D:95:LEU:HD13	1:D:105:LEU:HG	1.93	0.49
1:B:244:VAL:C	1:B:245:ILE:HD12	2.37	0.49
1:B:95:LEU:HD13	1:B:105:LEU:HG	1.94	0.49
1:B:79:MET:HE3	1:B:84:PRO:HG3	1.95	0.49
1:B:22:LEU:HD21	1:B:44:LEU:HD13	1.95	0.49
1:C:176:GLU:HG3	1:C:204:SER:HB2	1.95	0.49
1:A:27:TYR:OH	1:A:324:PRO:HG3	2.13	0.48
1:C:31:ARG:CD	1:C:324:PRO:HB3	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:92:VAL:HA	1:D:107:THR:HG22	1.95	0.48
1:A:325:MET:HE3	1:D:321:LYS:HD2	1.94	0.48
1:B:145:HIS:HB2	1:B:245:ILE:HG13	1.95	0.48
1:A:210:THR:HG21	4:A:1109:HOH:O	2.13	0.48
1:B:265:GLN:HG3	1:C:265:GLN:HG3	1.96	0.48
1:A:151:MET:HE3	1:A:175:LEU:HG	1.95	0.47
1:C:321:LYS:HE2	1:C:323:GLN:O	2.13	0.47
1:B:121:ALA:HB3	1:B:288:ILE:HD12	1.95	0.47
1:A:158:ARG:NH1	4:A:1104:HOH:O	2.42	0.47
1:C:128:PRO:HG2	1:C:130:ARG:HH22	1.79	0.47
1:A:19:PRO:HG3	1:A:48:LEU:HD13	1.97	0.47
1:C:135:GLY:O	1:C:137:ALA:N	2.48	0.46
1:A:128:PRO:HG2	1:A:130:ARG:NH2	2.30	0.46
1:A:81:LYS:NZ	4:B:1107:HOH:O	2.49	0.46
1:C:173:MET:HE2	1:C:200:ASN:HB3	1.98	0.46
1:C:190:LYS:HD2	1:C:190:LYS:HA	1.68	0.46
1:B:81:LYS:HD3	1:B:331:PHE:CE1	2.51	0.46
1:B:239:ILE:HG22	1:B:241:VAL:HG13	1.97	0.46
1:B:198:VAL:O	1:B:202:MET:HG2	2.15	0.46
1:A:298:ILE:HG22	2:A:1001:FAD:C2	2.45	0.45
1:B:254:LEU:HD12	1:B:257:ILE:HG21	1.98	0.45
1:C:239:ILE:HG22	1:C:241:VAL:HG13	1.97	0.45
1:D:81:LYS:HE3	1:D:81:LYS:HB2	1.41	0.45
1:B:164:GLY:HA3	1:B:186:HIS:ND1	2.32	0.45
1:B:81:LYS:HB2	1:B:81:LYS:HE3	1.74	0.45
1:C:56:TYR:O	1:D:328:SER:HB3	2.17	0.45
1:C:30:MET:CE	1:D:59:ASP:HB3	2.47	0.45
1:D:176:GLU:OE2	1:D:204:SER:HB2	2.17	0.45
1:C:31:ARG:HD3	1:C:324:PRO:HB3	1.97	0.45
1:A:181:LYS:NZ	1:A:207:GLU:OE1	2.48	0.44
1:A:295:VAL:HG12	1:A:297:LEU:HG	1.98	0.44
1:B:298:ILE:HG22	2:B:1001:FAD:O2	2.17	0.44
1:B:159:VAL:O	1:B:182:VAL:HA	2.18	0.44
1:B:254:LEU:HD12	1:B:257:ILE:CG2	2.47	0.44
1:C:174:MET:HE3	1:C:174:MET:HB2	1.88	0.44
2:C:1001:FAD:H9	2:C:1001:FAD:H1'1	1.86	0.44
1:B:311:ASN:HB3	4:B:1146:HOH:O	2.17	0.44
2:D:1001:FAD:O2'	2:D:1001:FAD:O4'	2.31	0.44
1:B:159:VAL:CG2	1:B:243:ASP:HB2	2.47	0.43
1:D:133:LEU:O	1:D:136:THR:HG22	2.18	0.43
1:C:52:TYR:HD2	1:C:55:LYS:HB2	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:103:PHE:O	1:C:113:TYR:HA	2.18	0.43
1:D:258:LYS:NZ	1:D:267:ASN:OD1	2.48	0.43
1:B:130:ARG:HD3	1:B:130:ARG:HA	1.83	0.43
1:B:321:LYS:HG2	1:B:323:GLN:O	2.18	0.43
1:D:172:THR:OG1	1:D:184:ILE:HD12	2.18	0.43
1:A:81:LYS:HE3	1:A:81:LYS:HB2	1.12	0.42
1:D:31:ARG:NH1	1:D:310:ASN:OD1	2.45	0.42
1:C:310:ASN:ND2	1:D:295:VAL:HG21	2.35	0.42
1:C:81:LYS:NZ	1:C:331:PHE:HZ	2.11	0.42
1:C:81:LYS:HB2	1:C:81:LYS:HE3	1.76	0.42
1:A:242:ASP:HA	4:A:1152:HOH:O	2.20	0.42
1:A:266:LYS:NZ	4:A:1111:HOH:O	2.53	0.42
1:A:313:LYS:HA	1:A:316:PHE:CZ	2.55	0.42
1:C:287:ASP:O	1:C:296:LYS:HE3	2.20	0.42
1:D:155:ALA:HA	1:D:178:ILE:O	2.19	0.42
1:C:331:PHE:CE2	1:D:58:TYR:CG	3.08	0.41
1:A:289:CYS:O	1:A:294:LYS:HE2	2.20	0.41
1:A:81:LYS:HD3	1:A:331:PHE:CZ	2.56	0.41
1:A:142:LYS:HE3	1:A:221:ASP:HB2	2.02	0.41
1:A:331:PHE:N	1:A:331:PHE:CD1	2.89	0.41
1:B:8:LYS:HD3	1:B:10:TYR:CE1	2.56	0.41
2:B:1001:FAD:HI'1	2:B:1001:FAD:H9	1.87	0.41
1:C:327:SER:HA	1:C:330:MET:SD	2.61	0.41
1:D:174:MET:HE2	1:D:174:MET:HB3	1.94	0.40
1:A:52:TYR:HB2	1:A:55:LYS:HG3	2.03	0.40
1:B:52:TYR:HD2	1:B:55:LYS:HB2	1.85	0.40
1:B:125:ALA:HA	2:B:1001:FAD:C8A	2.51	0.40
1:D:195:GLU:HA	1:D:198:VAL:HG22	2.03	0.40
1:C:151:MET:HE3	1:C:151:MET:HA	2.03	0.40
1:D:72:VAL:O	1:D:76:LYS:HG3	2.22	0.40
1:C:8:LYS:HD3	1:C:10:TYR:CE1	2.57	0.40
1:B:129:ARG:HD2	1:B:248:TYR:HB3	2.02	0.40
1:C:11:ASP:HB2	1:C:33:ALA:HB1	2.04	0.40
1:D:255:GLY:HA3	1:D:256:PRO:HD3	1.91	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	324/331 (98%)	308 (95%)	14 (4%)	2 (1%)	21	32
1	B	325/331 (98%)	313 (96%)	10 (3%)	2 (1%)	21	32
1	C	325/331 (98%)	312 (96%)	12 (4%)	1 (0%)	36	50
1	D	324/331 (98%)	309 (95%)	13 (4%)	2 (1%)	21	32
All	All	1298/1324 (98%)	1242 (96%)	49 (4%)	7 (0%)	24	37

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	136	THR
1	C	136	THR
1	D	136	THR
1	B	6	ASN
1	D	211	PRO
1	B	211	PRO
1	A	211	PRO

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	277/281 (99%)	267 (96%)	10 (4%)	31	52
1	B	278/281 (99%)	269 (97%)	9 (3%)	34	56
1	C	278/281 (99%)	275 (99%)	3 (1%)	65	82

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	D	277/281 (99%)	270 (98%)	7 (2%)	42 64
All	All	1110/1124 (99%)	1081 (97%)	29 (3%)	40 63

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	176	GLU
1	A	192	ARG
1	A	207	GLU
1	A	210	THR
1	A	251	VAL
1	A	254	LEU
1	A	263	ASP
1	A	298	ILE
1	A	328	SER
1	A	331	PHE
1	B	27	TYR
1	B	81	LYS
1	B	96	GLU
1	B	136	THR
1	B	151	MET
1	B	159	VAL
1	B	198	VAL
1	B	221	ASP
1	B	263	ASP
1	C	109	LYS
1	C	176	GLU
1	C	198	VAL
1	D	32	GLN
1	D	35	VAL
1	D	72	VAL
1	D	176	GLU
1	D	239	ILE
1	D	295	VAL
1	D	325	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	47	GLN

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Mol	Chain	Res	Type
1	A	259	ASN
1	A	267	ASN
1	B	32	GLN
1	B	259	ASN
1	C	7	GLN
1	C	123	ASN
1	D	32	GLN
1	D	229	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](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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	1001	-	58,58,58	0.50	1 (1%)	85,89,89	0.53	1 (1%)
2	FAD	D	1001	3	58,58,58	0.40	0	85,89,89	0.65	2 (2%)
2	FAD	A	1001	3	58,58,58	0.63	1 (1%)	85,89,89	0.60	0
2	FAD	C	1001	3	58,58,58	0.50	0	85,89,89	0.56	1 (1%)

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	1001	-	-	2/34/50/50	0/6/6/6
2	FAD	D	1001	3	-	1/34/50/50	0/6/6/6
2	FAD	A	1001	3	-	0/34/50/50	0/6/6/6
2	FAD	C	1001	3	-	1/34/50/50	0/6/6/6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	FAD	P-O3P	3.09	1.62	1.59
2	B	1001	FAD	P-O2P	-2.23	1.45	1.55

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1001	FAD	O2A-PA-O1A	2.82	125.58	112.44
2	D	1001	FAD	O2P-P-O1P	2.72	125.09	112.44
2	C	1001	FAD	O2P-P-O1P	2.66	124.82	112.44
2	B	1001	FAD	C4'-C3'-C2'	-2.31	109.72	113.57

There are no chirality outliers.

All (4) torsion outliers are listed below:

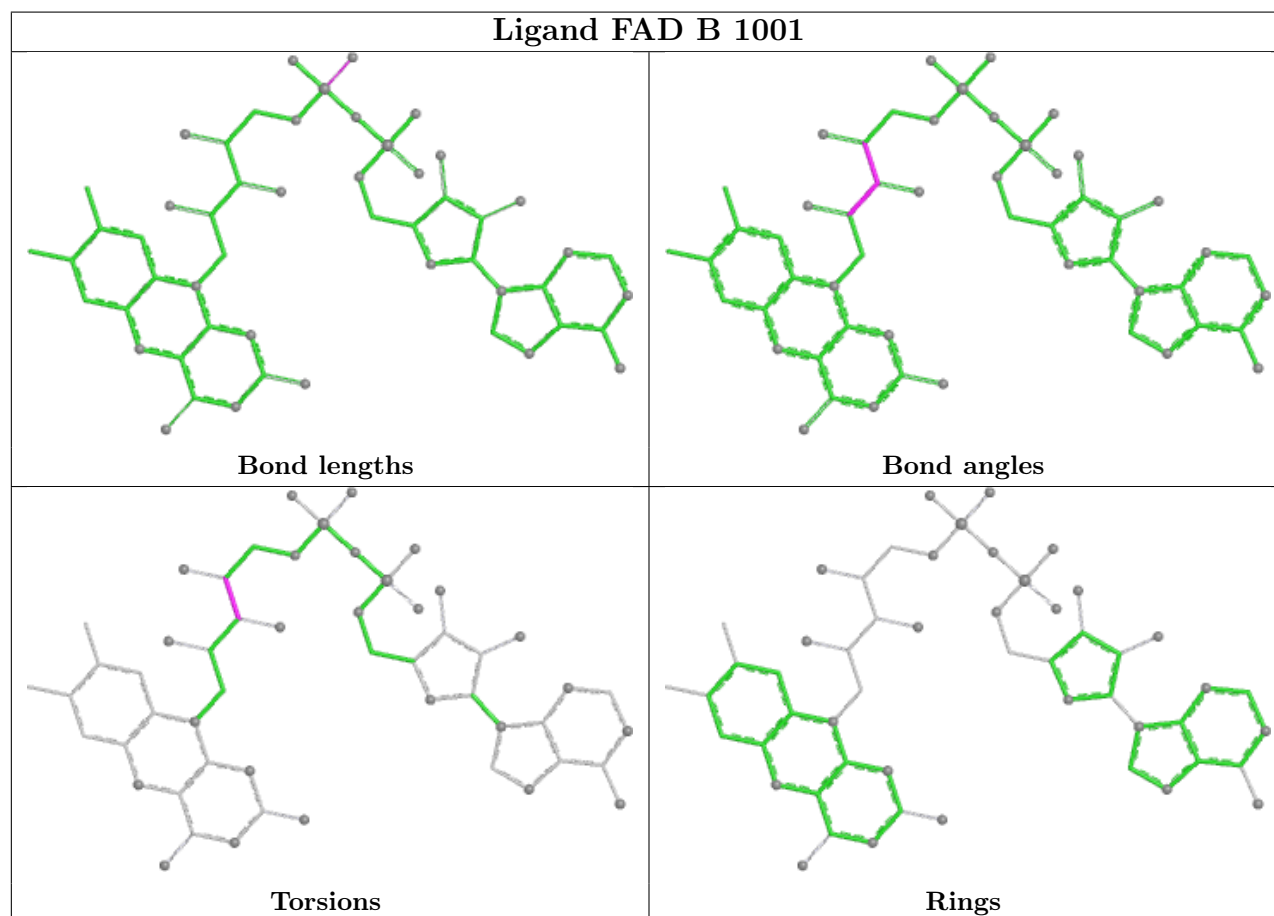
Mol	Chain	Res	Type	Atoms
2	D	1001	FAD	O4B-C4B-C5B-O5B
2	B	1001	FAD	O3'-C3'-C4'-C5'
2	B	1001	FAD	O3'-C3'-C4'-O4'
2	C	1001	FAD	O4B-C4B-C5B-O5B

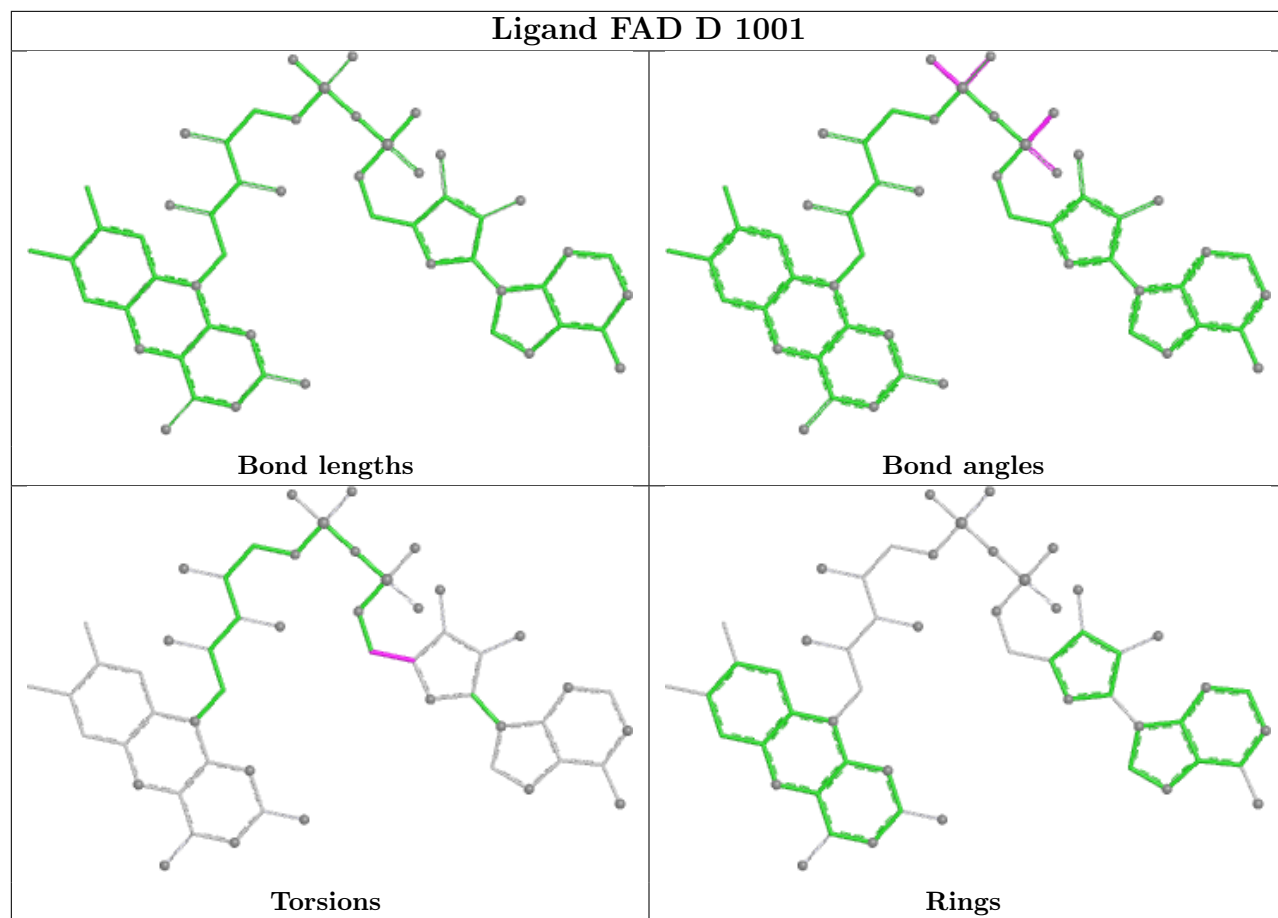
There are no ring outliers.

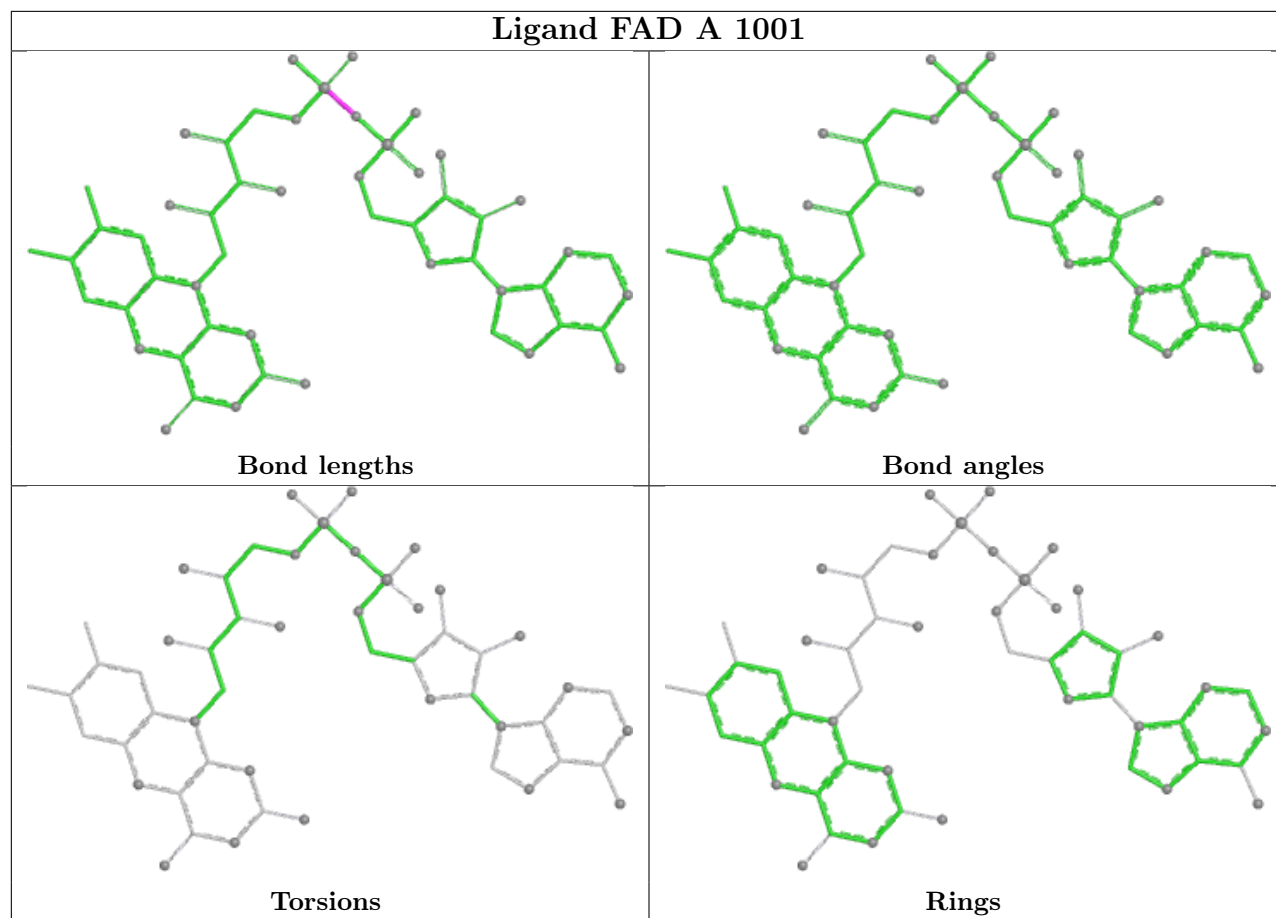
4 monomers are involved in 8 short contacts:

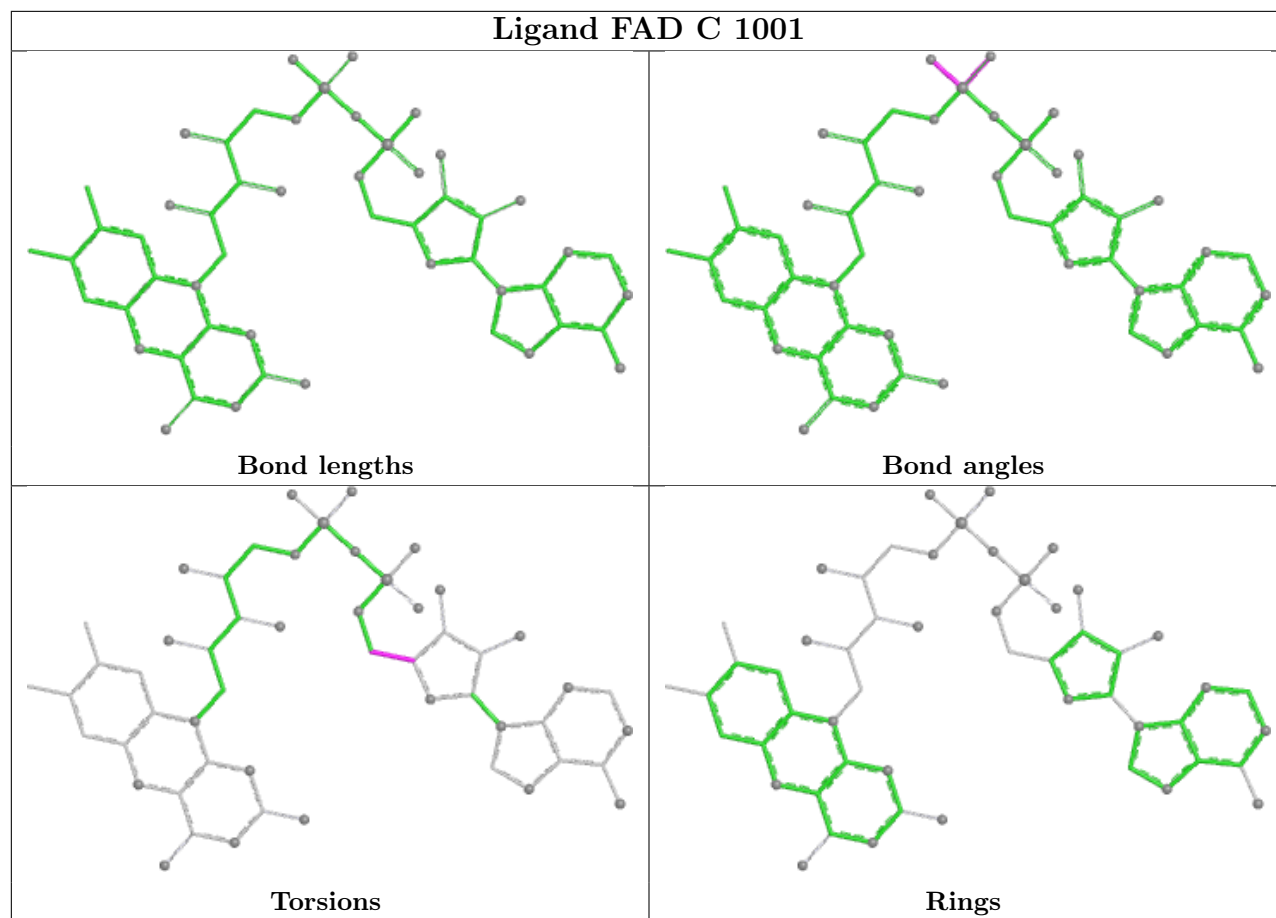
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1001	FAD	4	0
2	D	1001	FAD	1	0
2	A	1001	FAD	2	0
2	C	1001	FAD	1	0

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	326/331 (98%)	0.15	13 (3%) 42 38	33, 50, 85, 101	0
1	B	327/331 (98%)	0.30	5 (1%) 72 68	33, 49, 70, 102	0
1	C	327/331 (98%)	0.35	16 (4%) 35 31	35, 57, 91, 115	0
1	D	326/331 (98%)	0.36	11 (3%) 48 44	35, 57, 91, 114	0
All	All	1306/1324 (98%)	0.29	45 (3%) 48 44	33, 53, 85, 115	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	7	GLN	4.6
1	C	328	SER	4.3
1	A	321	LYS	4.1
1	C	327	SER	4.0
1	D	321	LYS	3.9
1	A	320	ALA	3.7
1	A	324	PRO	3.6
1	D	325	MET	3.6
1	C	330	MET	3.4
1	A	325	MET	3.4
1	C	323	GLN	3.3
1	A	323	GLN	3.1
1	D	330	MET	3.0
1	B	6	ASN	3.0
1	C	321	LYS	2.9
1	C	329	SER	2.9
1	C	320	ALA	2.9
1	B	325	MET	2.9
1	D	331	PHE	2.8
1	A	322	LEU	2.7
1	D	324	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	324	PRO	2.5
1	C	325	MET	2.5
1	C	196	HIS	2.5
1	C	197	SER	2.5
1	A	318	PRO	2.5
1	A	328	SER	2.5
1	D	327	SER	2.5
1	A	196	HIS	2.3
1	B	171	TRP	2.3
1	C	322	LEU	2.3
1	A	327	SER	2.2
1	B	178	ILE	2.2
1	D	318	PRO	2.2
1	D	7	GLN	2.2
1	C	5	GLU	2.2
1	D	136	THR	2.1
1	C	56	TYR	2.1
1	D	198	VAL	2.1
1	C	50	ALA	2.1
1	A	326	HIS	2.1
1	A	331	PHE	2.0
1	C	331	PHE	2.0
1	D	320	ALA	2.0
1	A	329	SER	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](#)

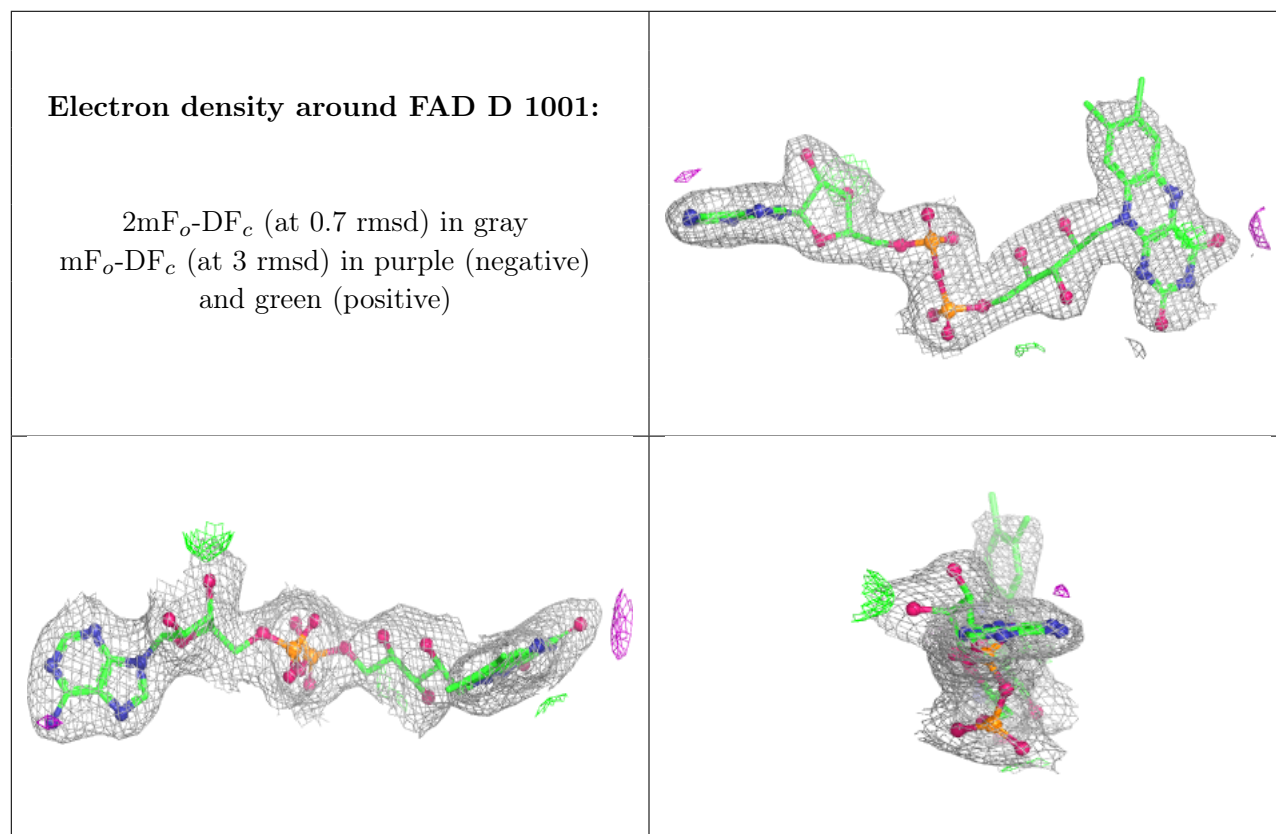
There are no oligosaccharides in this entry.

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.

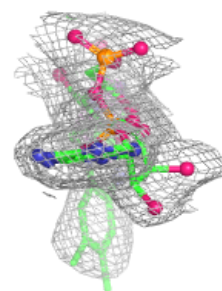
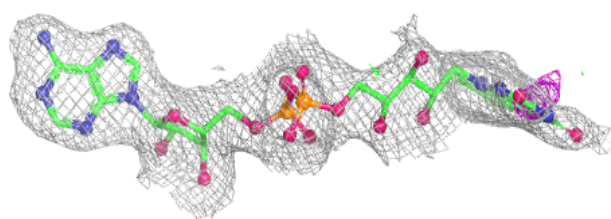
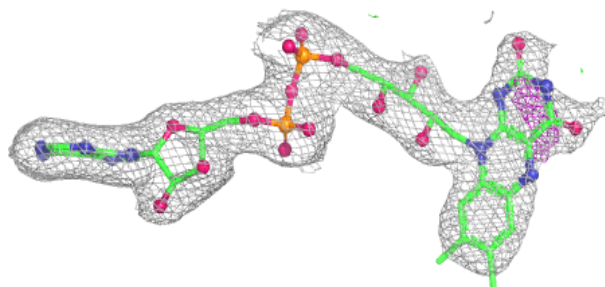
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NA	B	1002	1/1	0.90	0.13	56,56,56,56	0
3	NA	C	1002	1/1	0.92	0.12	54,54,54,54	0
3	NA	D	1002	1/1	0.94	0.17	55,55,55,55	0
2	FAD	D	1001	53/53	0.95	0.09	37,55,82,87	0
2	FAD	C	1001	53/53	0.95	0.10	36,55,83,87	0
2	FAD	B	1001	53/53	0.96	0.10	30,45,79,82	0
3	NA	A	1002	1/1	0.97	0.12	45,45,45,45	0
2	FAD	A	1001	53/53	0.97	0.07	30,42,55,61	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.

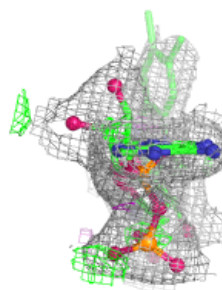
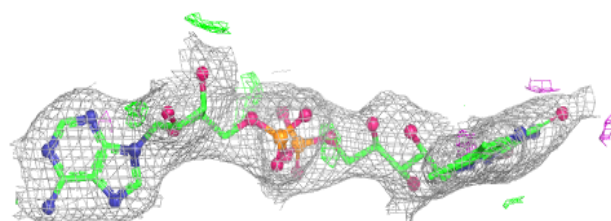
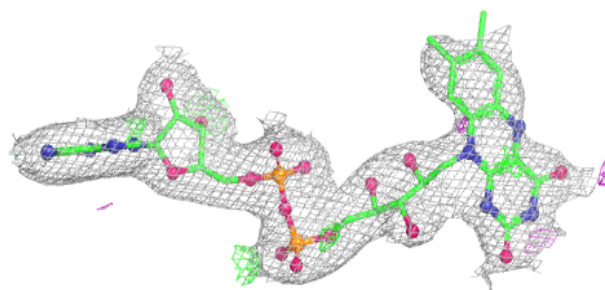


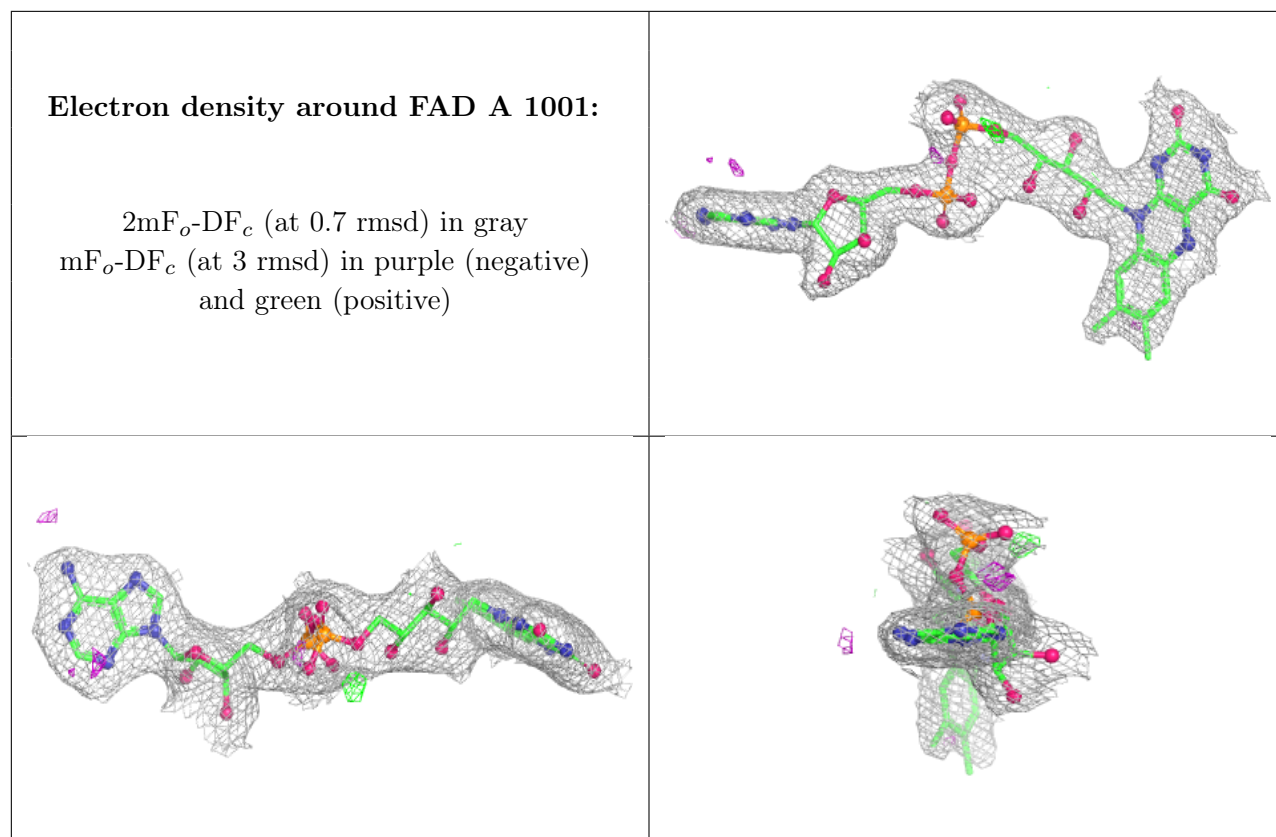
Electron density around FAD C 1001:

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

**Electron density around FAD B 1001:**

$2mF_o-DF_c$ (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.