



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2023 – 05:34 AM EDT

PDB ID : 2FMO
Title : Ala177Val mutant of E. coli Methylenetetrahydrofolate Reductase
Authors : Pejchal, R.; Campbell, E.; Guenther, B.D.; Lennon, B.W.; Matthews, R.G.; Ludwig, M.L.
Deposited on : 2006-01-09
Resolution : 2.25 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

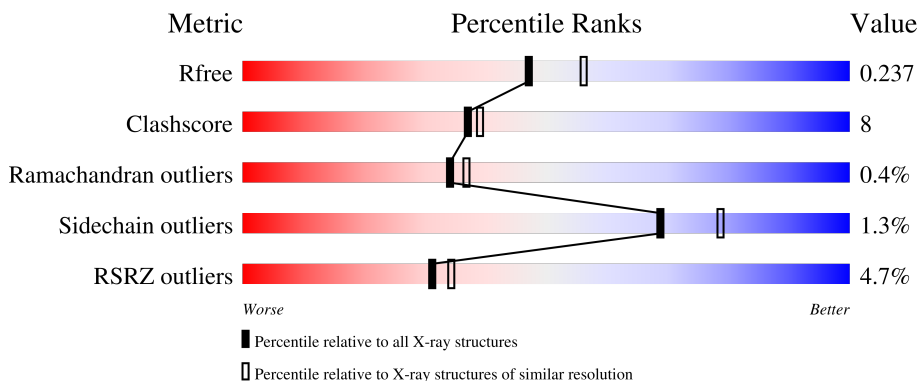
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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}	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	304	 6% 72% 13% • 14%
1	B	304	 5% 74% 17% 9%
1	C	304	 2% 71% 15% • 13%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6563 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 5,10-methylenetetrahydrofolate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	261	1984	1270	334	369	11	0	0	0
1	B	276	2093	1330	358	394	11	0	0	0
1	C	265	2019	1283	345	380	11	0	0	0

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	177	VAL	ALA	engineered mutation	UNP P0AEZ1
A	297	LEU	-	cloning artifact	UNP P0AEZ1
A	298	GLU	-	cloning artifact	UNP P0AEZ1
A	299	HIS	-	expression tag	UNP P0AEZ1
A	300	HIS	-	expression tag	UNP P0AEZ1
A	301	HIS	-	expression tag	UNP P0AEZ1
A	302	HIS	-	expression tag	UNP P0AEZ1
A	303	HIS	-	expression tag	UNP P0AEZ1
A	304	HIS	-	expression tag	UNP P0AEZ1
B	177	VAL	ALA	engineered mutation	UNP P0AEZ1
B	297	LEU	-	cloning artifact	UNP P0AEZ1
B	298	GLU	-	cloning artifact	UNP P0AEZ1
B	299	HIS	-	expression tag	UNP P0AEZ1
B	300	HIS	-	expression tag	UNP P0AEZ1
B	301	HIS	-	expression tag	UNP P0AEZ1
B	302	HIS	-	expression tag	UNP P0AEZ1
B	303	HIS	-	expression tag	UNP P0AEZ1
B	304	HIS	-	expression tag	UNP P0AEZ1
C	177	VAL	ALA	engineered mutation	UNP P0AEZ1
C	297	LEU	-	cloning artifact	UNP P0AEZ1
C	298	GLU	-	cloning artifact	UNP P0AEZ1
C	299	HIS	-	expression tag	UNP P0AEZ1
C	300	HIS	-	expression tag	UNP P0AEZ1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	301	HIS	-	expression tag	UNP P0AEZ1
C	302	HIS	-	expression tag	UNP P0AEZ1
C	303	HIS	-	expression tag	UNP P0AEZ1
C	304	HIS	-	expression tag	UNP P0AEZ1

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 8 6 2	0	0
4	B	1	Total C O 8 6 2	0	0
4	C	1	Total C O 8 6 2	0	0

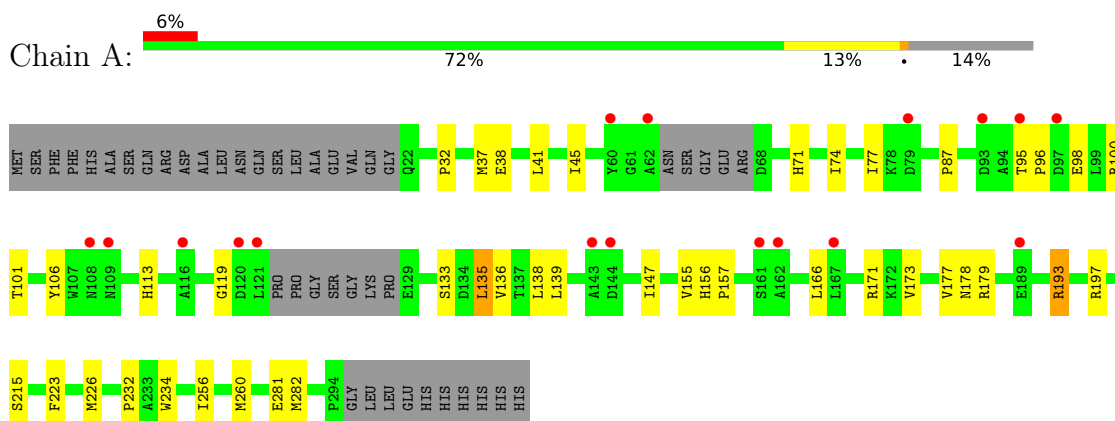
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	50	Total O 50 50	0	0
5	B	105	Total O 105 105	0	0
5	C	104	Total O 104 104	0	0

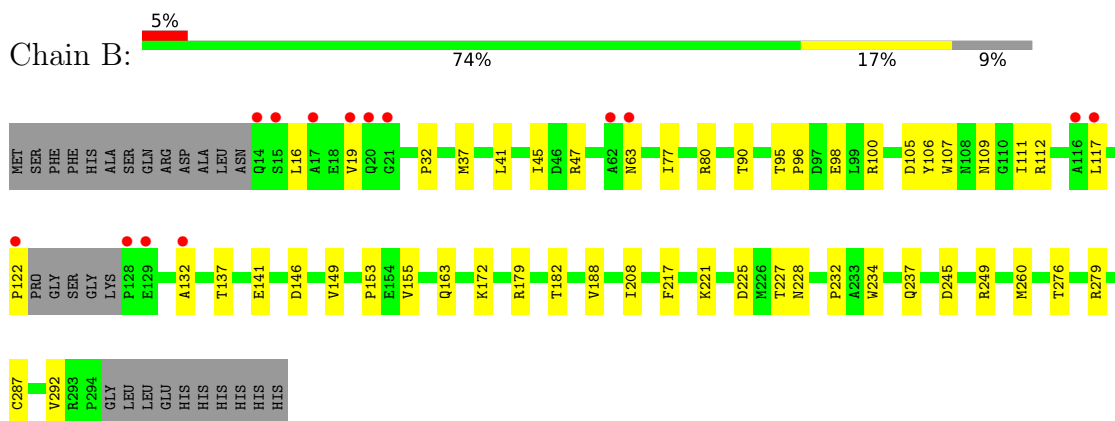
3 Residue-property plots [i](#)

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.

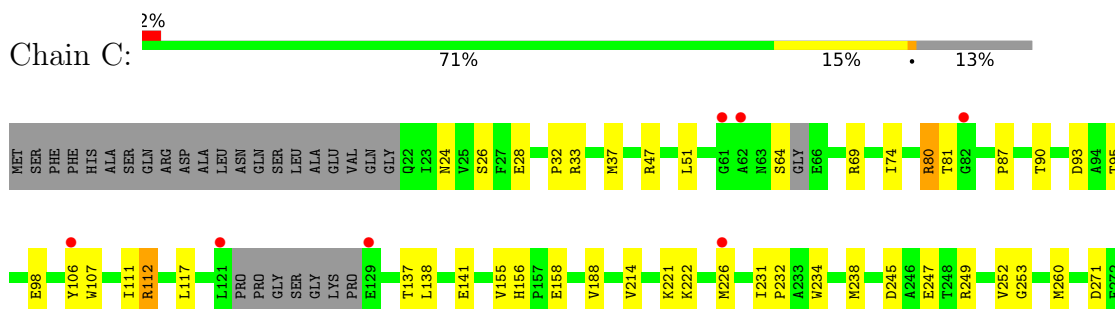
- Molecule 1: 5,10-methylenetetrahydrofolate reductase



- Molecule 1: 5,10-methylenetetrahydrofolate reductase



- Molecule 1: 5,10-methylenetetrahydrofolate reductase



H273	
C287	
V292	
R293	
P294	
GLY	
LEU	
LEU	
GLU	
HIS	
HIS	
HIS	
HIS	
HIS	
HIS	

4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.15Å 128.35Å 97.58Å 90.00° 121.07° 90.00°	Depositor
Resolution (Å)	20.01 – 2.25 20.01 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.01-2.25) 99.7 (20.01-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.09Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.219 , 0.247 0.208 , 0.237	Depositor DCC
R_{free} test set	2583 reflections (4.14%)	wwPDB-VP
Wilson B-factor (Å ²)	30.6	Xtrriage
Anisotropy	0.081	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.014 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6563	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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: FAD, MPD, SO4

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	0.33	0/2025	0.55	0/2755
1	B	0.37	0/2137	0.60	1/2910 (0.0%)
1	C	0.37	0/2060	0.58	0/2802
All	All	0.36	0/6222	0.58	1/8467 (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	146	ASP	N-CA-C	-5.37	96.50	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	1984	0	1901	31	1
1	B	2093	0	1990	32	0
1	C	2019	0	1917	34	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	5	0	0	0	0
3	A	53	0	31	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	53	0	31	0	0
3	C	53	0	31	0	0
4	A	8	0	13	1	0
4	B	8	0	13	0	0
4	C	8	0	13	0	0
5	A	50	0	0	2	0
5	B	105	0	0	1	0
5	C	104	0	0	1	0
All	All	6563	0	5940	96	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 8.

All (96) 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:C:80:ARG:HG3	1:C:81:THR:HG23	1.61	0.82
1:A:223:PHE:HA	1:A:226:MET:HE3	1.61	0.80
1:C:188:VAL:HG22	1:C:260:MET:HG3	1.64	0.77
1:C:32:PRO:HB3	1:C:37:MET:HG3	1.70	0.73
1:C:69:ARG:NH1	5:C:509:HOH:O	2.23	0.72
1:C:188:VAL:CG2	1:C:260:MET:HG3	2.19	0.72
1:C:137:THR:O	1:C:141:GLU:HG3	1.91	0.70
1:A:74:ILE:HG13	1:A:87:PRO:HB3	1.74	0.69
1:B:188:VAL:HG22	1:B:260:MET:HG3	1.75	0.69
1:A:71:HIS:HD1	1:A:106:TYR:HE1	1.39	0.69
1:C:74:ILE:HG13	1:C:87:PRO:HB3	1.76	0.66
1:B:96:PRO:O	1:B:100:ARG:HG3	1.95	0.65
1:A:113:HIS:CE1	1:A:179:ARG:HH21	2.17	0.63
1:C:156:HIS:CD2	1:C:158:GLU:H	2.18	0.62
1:A:96:PRO:O	1:A:100:ARG:HG3	2.00	0.62
1:A:173:VAL:HA	1:A:177:VAL:HG22	1.81	0.61
1:B:16:LEU:O	1:B:19:VAL:HG12	2.00	0.61
1:A:232:PRO:HB2	1:A:234:TRP:CD1	2.36	0.61
1:B:32:PRO:HD3	1:B:41:LEU:HD22	1.83	0.60
1:B:137:THR:O	1:B:141:GLU:HG3	2.01	0.60
1:B:132:ALA:HB3	1:B:172:LYS:HD2	1.83	0.59
1:C:112:ARG:HD2	1:C:112:ARG:N	2.18	0.58
1:C:33:ARG:HD2	1:C:64:SER:HB2	1.84	0.57
1:B:90:THR:HA	1:B:117:LEU:O	2.05	0.56
1:C:106:TYR:HB3	1:C:111:ILE:HB	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:ARG:O	1:B:80:ARG:HG2	2.06	0.55
1:B:188:VAL:CG2	1:B:260:MET:HG3	2.35	0.55
1:B:63:ASN:OD1	1:B:122:PRO:HG3	2.07	0.55
1:C:188:VAL:HG22	1:C:260:MET:CG	2.37	0.54
1:B:132:ALA:CB	1:B:172:LYS:HD2	2.37	0.54
1:C:26:SER:OG	1:C:273:HIS:HD2	1.91	0.53
1:C:238:MET:HG3	1:C:252:VAL:HG11	1.90	0.53
1:A:281:GLU:HG2	1:C:247:GLU:HG2	1.89	0.53
1:B:245:ASP:O	1:B:249:ARG:HG3	2.09	0.53
1:B:287:CYS:HB3	1:B:292:VAL:HB	1.91	0.53
1:B:32:PRO:HB3	1:B:37:MET:HG3	1.92	0.52
1:C:28:GLU:OE1	1:C:273:HIS:HE1	1.93	0.52
1:C:47:ARG:HH21	1:C:47:ARG:HG2	1.75	0.51
1:C:107:TRP:CZ3	1:C:112:ARG:HG3	2.45	0.51
1:A:173:VAL:HG22	1:A:177:VAL:HG21	1.93	0.51
1:A:256:ILE:O	1:A:260:MET:HG3	2.11	0.51
1:A:95:THR:OG1	1:A:98:GLU:HG3	2.11	0.50
1:B:149:VAL:HB	1:B:172:LYS:HE3	1.93	0.49
1:A:173:VAL:HA	1:A:177:VAL:CG2	2.43	0.49
1:C:156:HIS:HD2	1:C:158:GLU:H	1.60	0.49
1:C:238:MET:HG3	1:C:252:VAL:CG1	2.43	0.48
1:B:105:ASP:O	1:B:109:ASN:ND2	2.47	0.48
1:C:95:THR:OG1	1:C:98:GLU:HG3	2.13	0.48
1:C:138:LEU:C	1:C:138:LEU:HD23	2.35	0.47
1:A:45:ILE:HD13	1:A:77:ILE:HG12	1.97	0.47
1:B:112:ARG:HH11	1:B:112:ARG:HG3	1.78	0.47
1:A:119:GLY:H	3:A:395:FAD:H5'1	1.78	0.47
1:C:24:ASN:HB2	1:C:271:ASP:OD1	2.14	0.47
1:B:47:ARG:HH11	1:B:47:ARG:HG2	1.79	0.47
1:A:281:GLU:HG2	1:C:247:GLU:CG	2.46	0.46
1:A:166:LEU:HD22	5:A:602:HOH:O	2.14	0.46
1:C:232:PRO:HB2	1:C:234:TRP:CD1	2.50	0.46
1:A:135:LEU:HD22	1:A:139:LEU:CD1	2.46	0.45
1:B:188:VAL:HG23	5:B:556:HOH:O	2.15	0.45
1:C:28:GLU:HB2	1:C:273:HIS:CE1	2.51	0.45
1:B:107:TRP:CE3	1:B:112:ARG:HD2	2.52	0.45
1:A:223:PHE:CA	1:A:226:MET:HE3	2.39	0.45
1:A:32:PRO:HG2	1:A:38:GLU:OE1	2.17	0.45
1:C:287:CYS:HB3	1:C:292:VAL:HB	1.97	0.45
1:A:173:VAL:HG13	1:A:177:VAL:HG22	1.98	0.45
1:A:232:PRO:HB2	1:A:234:TRP:NE1	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:221:LYS:HA	1:C:231:ILE:HD11	1.99	0.45
1:C:214:VAL:O	1:C:253:GLY:HA3	2.16	0.45
1:C:90:THR:HA	1:C:117:LEU:O	2.17	0.44
1:B:153:PRO:HD3	1:B:182:THR:HB	1.99	0.44
1:B:106:TYR:HB3	1:B:111:ILE:HB	1.99	0.44
1:A:178:ASN:HB2	1:A:179:ARG:NH1	2.33	0.44
1:A:215:SER:HB3	1:A:282:MET:CE	2.48	0.44
1:A:138:LEU:C	1:A:138:LEU:HD23	2.37	0.43
1:B:45:ILE:HD13	1:B:77:ILE:HG12	2.00	0.43
1:B:112:ARG:HG3	1:B:112:ARG:NH1	2.34	0.43
4:A:895:MPD:HM1	5:A:539:HOH:O	2.19	0.43
1:A:32:PRO:HB3	1:A:37:MET:HG3	1.99	0.43
1:A:136:VAL:HG12	1:A:147:ILE:HG21	2.00	0.43
1:A:193:ARG:HB3	1:A:197:ARG:NH1	2.34	0.43
1:B:217:PHE:CE2	1:B:221:LYS:HE2	2.53	0.43
1:B:95:THR:OG1	1:B:98:GLU:HG3	2.19	0.42
1:A:156:HIS:CG	1:A:157:PRO:HD2	2.54	0.42
1:C:245:ASP:O	1:C:249:ARG:HG3	2.20	0.41
1:B:179:ARG:HD3	1:B:208:ILE:HD12	2.02	0.41
1:C:51:LEU:CD2	1:C:294:PRO:HD3	2.50	0.41
1:C:32:PRO:CB	1:C:37:MET:HG3	2.46	0.41
1:B:107:TRP:CZ3	1:B:112:ARG:HD2	2.55	0.41
1:B:96:PRO:CB	1:B:100:ARG:HH12	2.34	0.41
1:B:227:THR:O	1:B:228:ASN:CB	2.69	0.41
1:B:232:PRO:HB2	1:B:234:TRP:CD1	2.56	0.41
1:B:276:THR:HB	1:B:279:ARG:O	2.21	0.41
1:C:222:LYS:O	1:C:226:MET:HG2	2.20	0.41
1:A:133:SER:O	1:A:136:VAL:CG2	2.69	0.40
1:A:32:PRO:HD3	1:A:41:LEU:HD22	2.04	0.40
1:A:98:GLU:O	1:A:101:THR:HB	2.22	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:171:ARG:NH1	1:A:171:ARG:NH1[2_455]	1.98	0.22

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	255/304 (84%)	246 (96%)	8 (3%)	1 (0%)	34	37
1	B	272/304 (90%)	270 (99%)	1 (0%)	1 (0%)	34	37
1	C	259/304 (85%)	253 (98%)	5 (2%)	1 (0%)	34	37
All	All	786/912 (86%)	769 (98%)	14 (2%)	3 (0%)	34	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	155	VAL
1	C	155	VAL
1	A	155	VAL

5.3.2 Protein sidechains [i](#)

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	202/259 (78%)	200 (99%)	2 (1%)	76	84
1	B	213/259 (82%)	210 (99%)	3 (1%)	67	76
1	C	206/259 (80%)	203 (98%)	3 (2%)	65	75
All	All	621/777 (80%)	613 (99%)	8 (1%)	69	79

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	LEU
1	A	193	ARG
1	B	163	GLN
1	B	225	ASP
1	B	237	GLN
1	C	80	ARG
1	C	93	ASP
1	C	112	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	ASN
1	A	163	GLN
1	A	228	ASN
1	B	71	HIS
1	B	109	ASN
1	B	163	GLN
1	C	156	HIS
1	C	168	ASN
1	C	273	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

11 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	SO4	A	901	-	4,4,4	0.26	0	6,6,6	0.09	0
2	SO4	B	903	-	4,4,4	0.31	0	6,6,6	0.10	0
4	MPD	B	896	-	7,7,7	1.25	1 (14%)	9,10,10	0.37	0
4	MPD	C	897	-	7,7,7	1.25	1 (14%)	9,10,10	0.34	0
3	FAD	B	396	-	53,58,58	2.10	16 (30%)	68,89,89	1.68	19 (27%)
2	SO4	C	904	-	4,4,4	0.23	0	6,6,6	0.15	0
2	SO4	B	902	-	4,4,4	0.23	0	6,6,6	0.09	0
3	FAD	A	395	-	53,58,58	1.91	13 (24%)	68,89,89	1.67	18 (26%)
3	FAD	C	397	-	53,58,58	1.75	12 (22%)	68,89,89	1.46	11 (16%)
4	MPD	A	895	-	7,7,7	1.26	1 (14%)	9,10,10	0.35	0
2	SO4	A	900	-	4,4,4	0.28	0	6,6,6	0.09	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
4	MPD	B	896	-	-	0/5/5/5	-
3	FAD	B	396	-	-	7/30/50/50	0/6/6/6
4	MPD	C	897	-	-	0/5/5/5	-
3	FAD	A	395	-	-	9/30/50/50	0/6/6/6
3	FAD	C	397	-	-	1/30/50/50	0/6/6/6
4	MPD	A	895	-	-	1/5/5/5	-

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	395	FAD	C4X-N5	5.79	1.42	1.30
3	B	396	FAD	C10-N10	5.11	1.48	1.37
3	A	395	FAD	O4B-C1B	5.10	1.48	1.41
3	B	396	FAD	C9A-N10	5.09	1.50	1.41
3	B	396	FAD	C4X-N5	4.91	1.40	1.30
3	C	397	FAD	C4X-N5	4.68	1.39	1.30
3	A	395	FAD	C10-N10	4.61	1.47	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	397	FAD	C10-N10	4.59	1.47	1.37
3	C	397	FAD	C9A-N10	4.30	1.48	1.41
3	A	395	FAD	C5'-C4'	4.13	1.57	1.51
3	B	396	FAD	O4B-C1B	3.93	1.46	1.41
3	B	396	FAD	C9-C9A	3.90	1.46	1.39
3	A	395	FAD	C9A-N10	3.85	1.47	1.41
3	C	397	FAD	C9-C9A	3.52	1.45	1.39
3	C	397	FAD	O4B-C1B	3.47	1.45	1.41
3	B	396	FAD	O5B-C5B	-3.46	1.31	1.44
3	B	396	FAD	C6-C5X	3.44	1.45	1.40
3	B	396	FAD	C8A-N7A	-3.38	1.28	1.34
3	A	395	FAD	C9-C9A	3.27	1.44	1.39
4	A	895	MPD	O2-C2	-3.23	1.36	1.44
3	C	397	FAD	C6-C5X	3.23	1.45	1.40
4	B	896	MPD	O2-C2	-3.19	1.36	1.44
3	B	396	FAD	C4'-C3'	3.18	1.59	1.53
4	C	897	MPD	O2-C2	-3.17	1.36	1.44
3	B	396	FAD	C2B-C1B	3.14	1.58	1.53
3	C	397	FAD	C8A-N7A	-3.08	1.29	1.34
3	B	396	FAD	C8-C7	2.86	1.48	1.40
3	A	395	FAD	C6-C5X	2.84	1.44	1.40
3	B	396	FAD	C7M-C7	2.81	1.56	1.51
3	C	397	FAD	C8-C7	2.77	1.47	1.40
3	A	395	FAD	C1'-C2'	2.72	1.56	1.52
3	B	396	FAD	C6-C7	2.61	1.43	1.39
3	C	397	FAD	C1'-C2'	2.59	1.56	1.52
3	A	395	FAD	C8A-N7A	-2.52	1.30	1.34
3	C	397	FAD	C2B-C1B	2.47	1.57	1.53
3	B	396	FAD	C8M-C8	2.42	1.55	1.51
3	A	395	FAD	C8-C7	2.38	1.46	1.40
3	A	395	FAD	C10-N1	2.36	1.38	1.33
3	C	397	FAD	C2B-C3B	-2.34	1.46	1.53
3	C	397	FAD	C8M-C8	2.28	1.55	1.51
3	B	396	FAD	C1'-C2'	2.25	1.55	1.52
3	B	396	FAD	C9A-C5X	2.21	1.44	1.41
3	A	395	FAD	O5B-C5B	-2.08	1.36	1.44
3	A	395	FAD	C9A-C5X	2.01	1.44	1.41

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	396	FAD	C5B-C4B-C3B	-4.94	96.65	115.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	395	FAD	C5'-C4'-C3'	3.79	119.52	112.20
3	B	396	FAD	O2B-C2B-C1B	3.72	124.60	110.85
3	B	396	FAD	C5'-C4'-C3'	3.67	119.29	112.20
3	C	397	FAD	C9-C9A-N10	3.47	126.52	121.84
3	C	397	FAD	O5'-C5'-C4'	3.29	118.14	109.36
3	A	395	FAD	C5B-C4B-C3B	-3.28	102.88	115.18
3	A	395	FAD	O5'-C5'-C4'	3.26	118.06	109.36
3	C	397	FAD	C5X-C9A-N10	-3.22	114.63	117.95
3	B	396	FAD	C9-C9A-N10	3.16	126.11	121.84
3	B	396	FAD	C5X-C9A-N10	-3.08	114.77	117.95
3	A	395	FAD	C9-C9A-N10	3.04	125.94	121.84
3	A	395	FAD	C5X-C9A-N10	-2.99	114.87	117.95
3	A	395	FAD	O2-C2-N3	2.98	124.45	118.65
3	A	395	FAD	O2'-C2'-C3'	2.92	116.19	109.10
3	A	395	FAD	C9A-C5X-N5	2.87	125.55	122.43
3	C	397	FAD	O2B-C2B-C1B	2.86	121.41	110.85
3	C	397	FAD	C8M-C8-C7	2.80	126.48	120.74
3	B	396	FAD	C8M-C8-C7	2.68	126.22	120.74
3	B	396	FAD	O4B-C4B-C5B	2.67	118.15	109.37
3	C	397	FAD	C8M-C8-C9	-2.59	114.70	119.49
3	B	396	FAD	C9A-C5X-N5	2.59	125.24	122.43
3	C	397	FAD	C9A-C5X-N5	2.57	125.22	122.43
3	B	396	FAD	C10-N1-C2	2.50	121.90	116.90
3	B	396	FAD	O2-C2-N3	2.48	123.48	118.65
3	B	396	FAD	C8M-C8-C9	-2.47	114.92	119.49
3	C	397	FAD	C5'-C4'-C3'	2.44	116.92	112.20
3	A	395	FAD	P-O5'-C5'	2.43	135.94	121.68
3	A	395	FAD	N3-C2-N1	-2.40	114.68	119.38
3	A	395	FAD	O2B-C2B-C1B	2.31	119.37	110.85
3	B	396	FAD	O3B-C3B-C2B	2.30	119.26	111.82
3	B	396	FAD	C4A-C5A-N7A	2.27	111.77	109.40
3	B	396	FAD	N3-C2-N1	-2.27	114.92	119.38
3	A	395	FAD	C8M-C8-C7	2.27	125.39	120.74
3	A	395	FAD	C10-N1-C2	2.26	121.43	116.90
3	C	397	FAD	C10-N1-C2	2.25	121.40	116.90
3	C	397	FAD	C3B-C2B-C1B	2.24	104.35	100.98
3	B	396	FAD	O3'-C3'-C4'	2.21	114.15	108.81
3	B	396	FAD	O2A-PA-O5B	2.18	117.85	107.75
3	A	395	FAD	C2B-C3B-C4B	2.14	106.81	102.64
3	B	396	FAD	O2A-PA-O1A	2.13	122.78	112.24
3	B	396	FAD	C9A-C9-C8	2.12	123.57	119.30
3	A	395	FAD	C4'-C3'-C2'	2.08	117.69	113.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	395	FAD	C8M-C8-C9	-2.05	115.69	119.49
3	B	396	FAD	O2P-P-O1P	2.05	122.39	112.24
3	A	395	FAD	O4B-C4B-C5B	2.05	116.11	109.37
3	A	395	FAD	O4B-C1B-C2B	-2.01	103.99	106.93
3	C	397	FAD	N3-C2-N1	-2.00	115.46	119.38

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	396	FAD	C5B-O5B-PA-O1A
3	B	396	FAD	C5B-O5B-PA-O2A
3	A	395	FAD	C3B-C4B-C5B-O5B
3	A	395	FAD	O3'-C3'-C4'-C5'
3	A	395	FAD	O2'-C2'-C3'-C4'
3	B	396	FAD	O2'-C2'-C3'-C4'
3	B	396	FAD	C2'-C3'-C4'-O4'
4	A	895	MPD	O2-C2-C3-C4
3	A	395	FAD	C5B-O5B-PA-O3P
3	A	395	FAD	C2'-C3'-C4'-C5'
3	B	396	FAD	C3B-C4B-C5B-O5B
3	A	395	FAD	O4B-C4B-C5B-O5B
3	A	395	FAD	PA-O3P-P-O1P
3	B	396	FAD	C5B-O5B-PA-O3P
3	A	395	FAD	PA-O3P-P-O2P
3	B	396	FAD	PA-O3P-P-O1P
3	A	395	FAD	C5B-O5B-PA-O1A
3	C	397	FAD	O2'-C2'-C3'-C4'

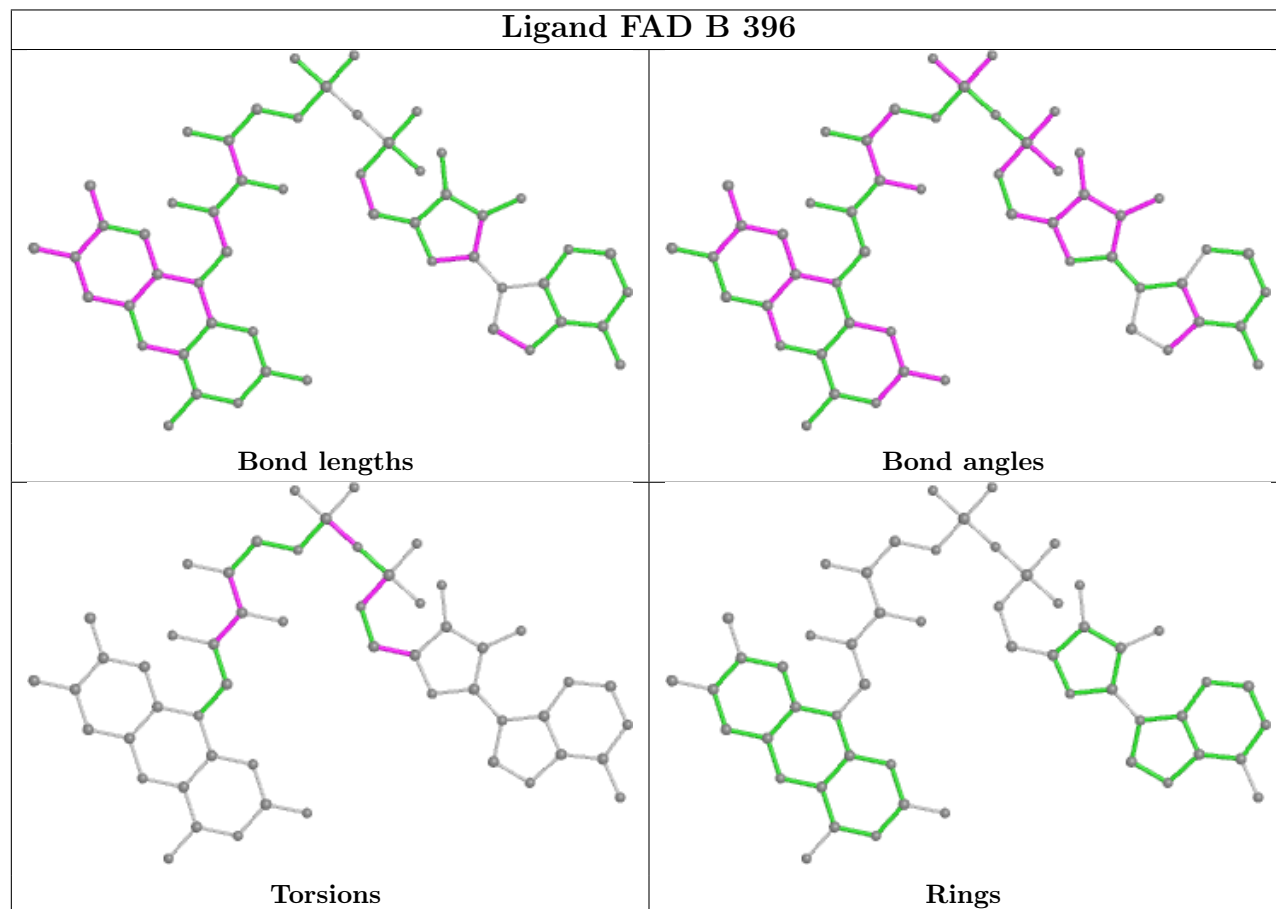
There are no ring outliers.

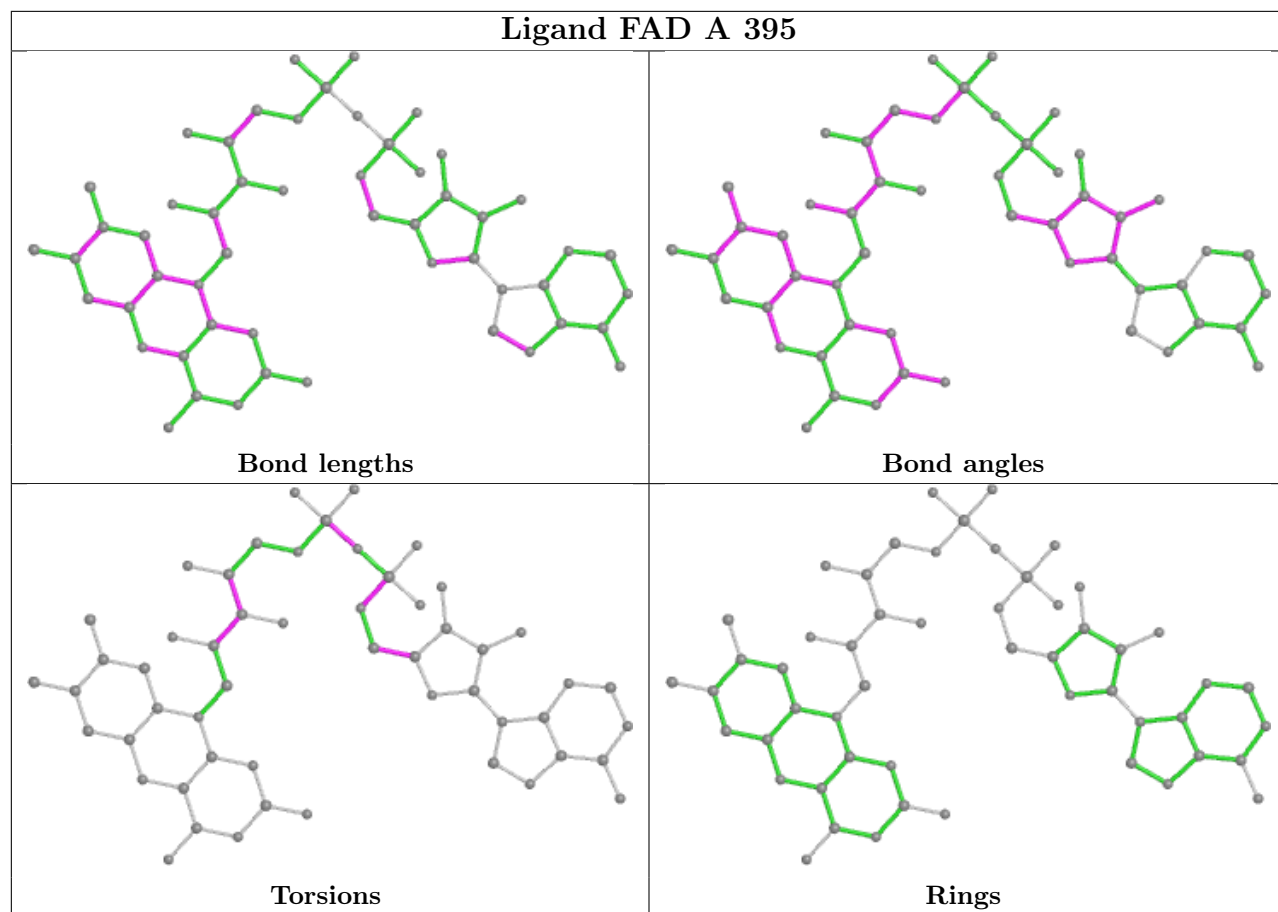
2 monomers are involved in 2 short contacts:

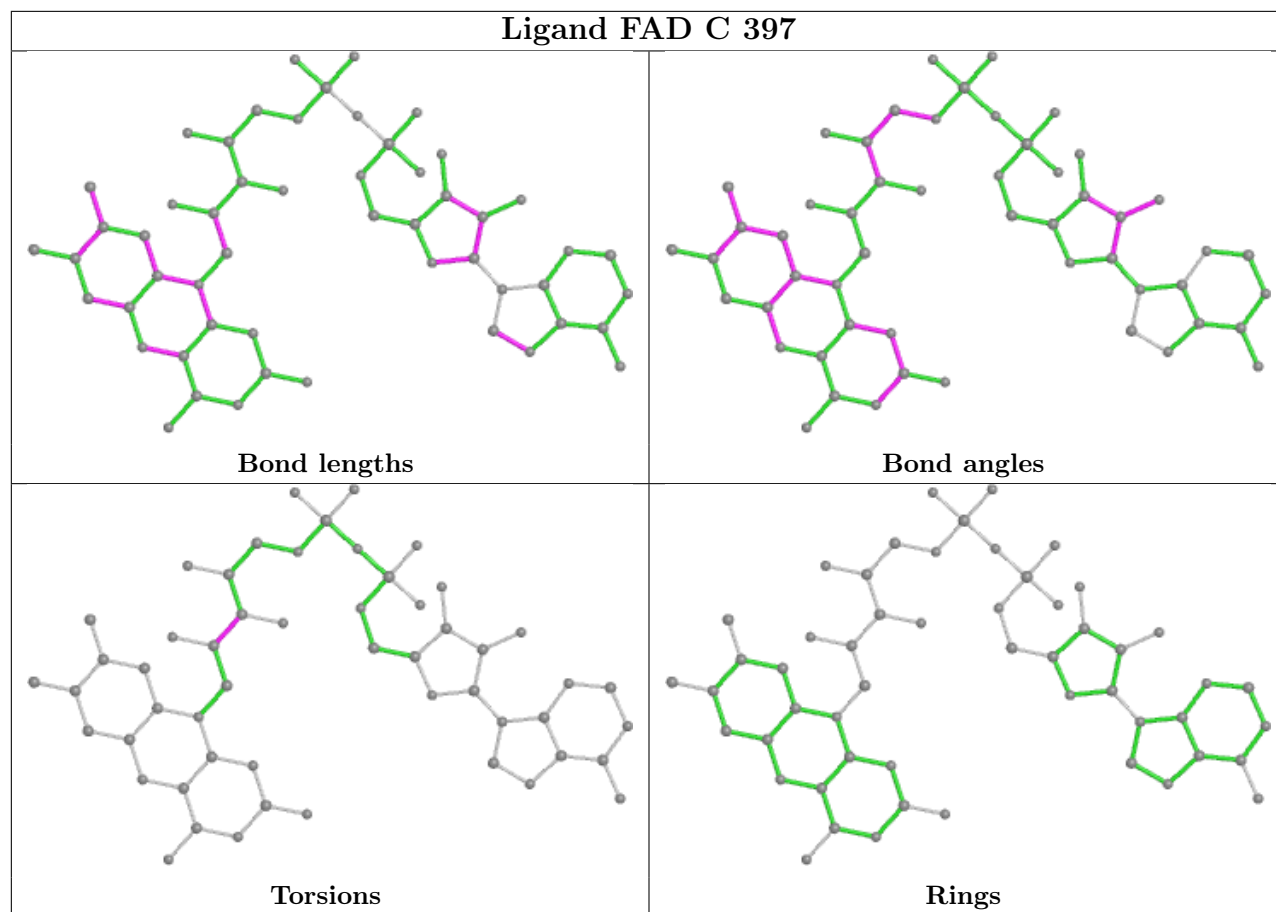
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	395	FAD	1	0
4	A	895	MPD	1	0

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	261/304 (85%)	0.40	17 (6%) 18 20	20, 38, 57, 64	0
1	B	276/304 (90%)	0.08	14 (5%) 28 30	15, 28, 48, 63	0
1	C	265/304 (87%)	-0.13	7 (2%) 56 59	16, 28, 42, 53	0
All	All	802/912 (87%)	0.11	38 (4%) 31 34	15, 31, 53, 64	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	121	LEU	6.2
1	C	121	LEU	4.6
1	A	143	ALA	4.0
1	A	120	ASP	3.9
1	A	93	ASP	3.9
1	B	17	ALA	3.5
1	B	19	VAL	3.5
1	B	14	GLN	3.2
1	A	62	ALA	3.2
1	A	189	GLU	3.0
1	B	122	PRO	2.9
1	B	117	LEU	2.9
1	A	108	ASN	2.9
1	A	144	ASP	2.8
1	B	21	GLY	2.7
1	A	109	ASN	2.6
1	C	106	TYR	2.6
1	C	226	MET	2.6
1	A	95	THR	2.6
1	B	20	GLN	2.5
1	B	116	ALA	2.5
1	B	132	ALA	2.4
1	C	61	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	15	SER	2.4
1	B	63	ASN	2.4
1	A	79	ASP	2.3
1	C	129	GLU	2.3
1	B	62	ALA	2.3
1	A	97	ASP	2.3
1	A	116	ALA	2.3
1	A	162	ALA	2.2
1	B	129	GLU	2.2
1	B	128	PRO	2.2
1	C	82	GLY	2.1
1	A	60	TYR	2.1
1	C	62	ALA	2.1
1	A	161	SER	2.0
1	A	167	LEU	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 monosaccharides 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.

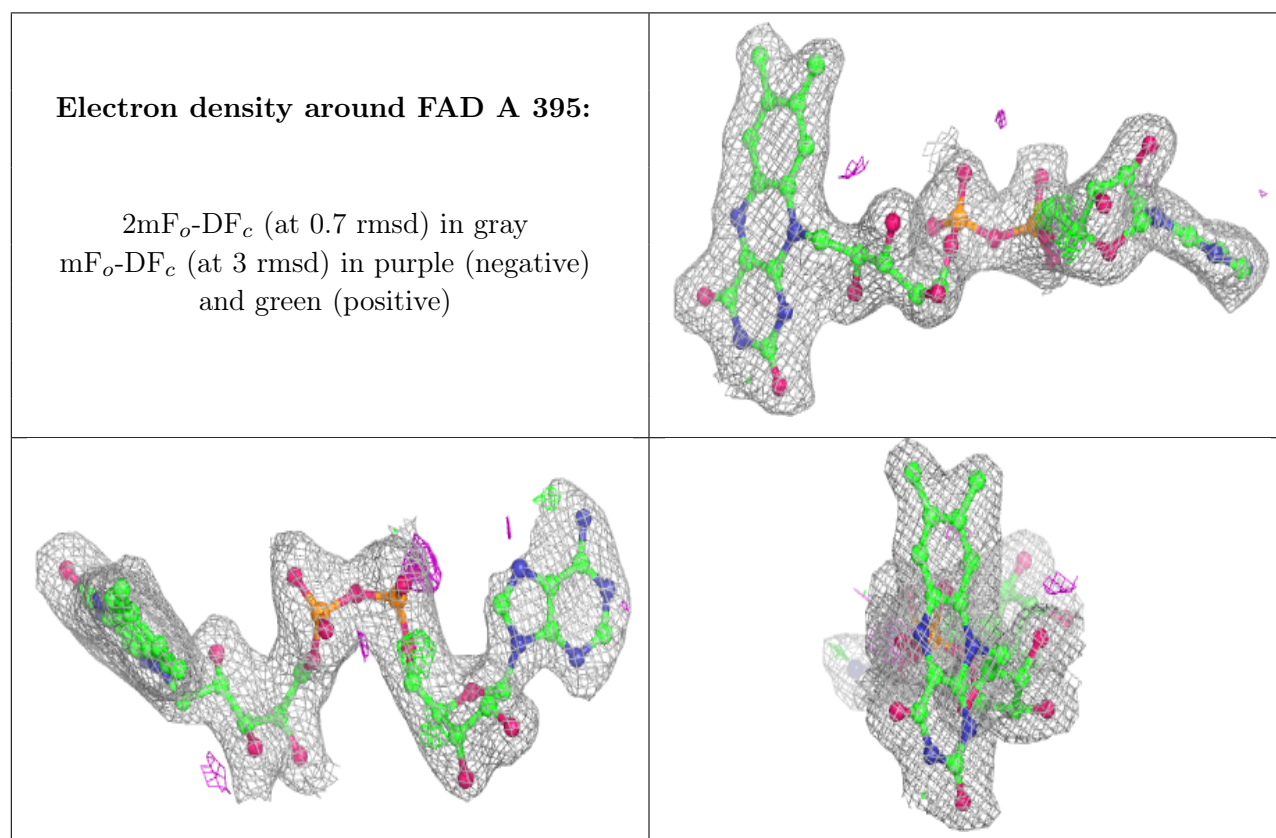
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	A	900	5/5	0.85	0.18	58,59,60,61	5
4	MPD	B	896	8/8	0.86	0.21	39,40,40,41	0
4	MPD	C	897	8/8	0.86	0.18	37,38,40,41	0
3	FAD	A	395	53/53	0.89	0.15	26,35,55,56	0
3	FAD	B	396	53/53	0.93	0.13	19,24,39,41	0
2	SO4	B	903	5/5	0.94	0.22	57,57,58,58	5
4	MPD	A	895	8/8	0.94	0.19	40,42,44,44	0
3	FAD	C	397	53/53	0.95	0.10	19,22,36,38	0

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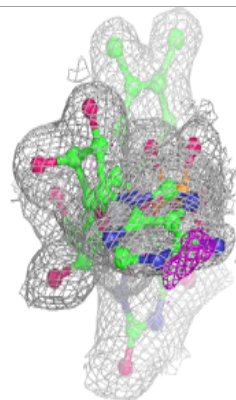
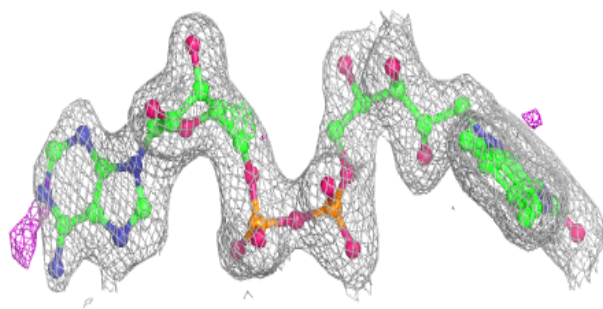
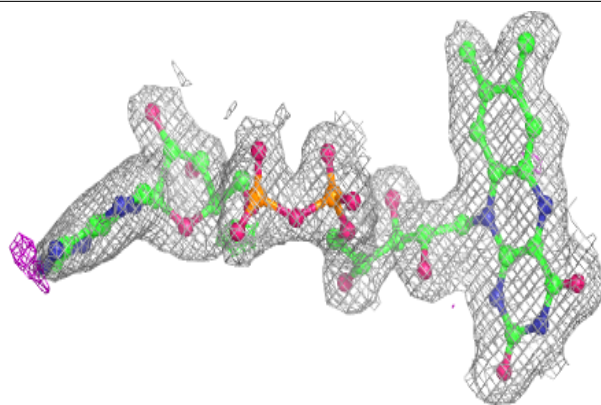
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	A	901	5/5	0.95	0.15	53,54,54,55	5
2	SO4	B	902	5/5	0.97	0.12	39,41,41,43	5
2	SO4	C	904	5/5	0.97	0.13	43,44,45,45	5

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.

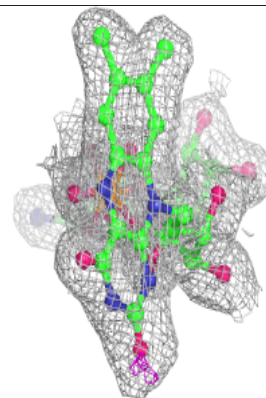
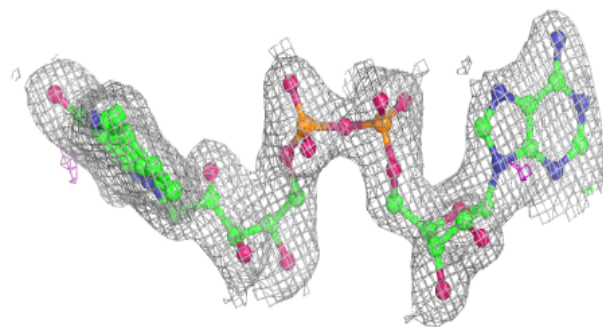
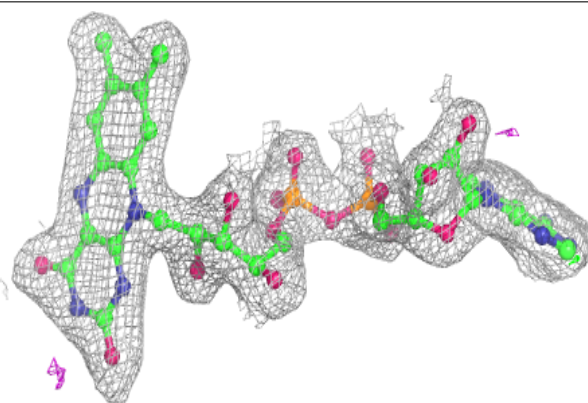


Electron density around FAD B 396:

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

**Electron density around FAD C 397:**

$2mF_o-DF_c$ (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.