



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 08:24 PM UTC

PDB ID : 3QSS / pdb_00003qss
Title : Crystal structure for the MSOX.chloride.MTA ternary complex
Authors : Kommoju, P.; Chen, Z.; Bruckner, R.C.; Mathews, F.S.; Jorns, M.S.
Deposited on : 2011-02-21
Resolution : 1.85 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

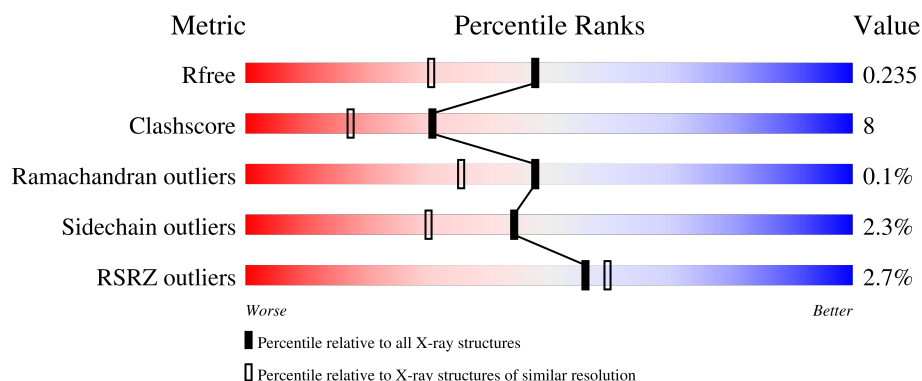
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	389	2% 77% 20% ...
1	B	389	4% 76% 20% ..

2 Entry composition [i](#)

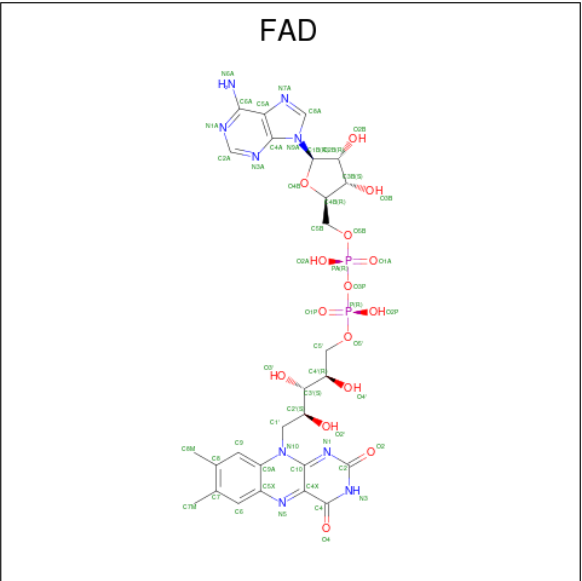
There are 5 unique types of molecules in this entry. The entry contains 6632 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Monomeric sarcosine oxidase.

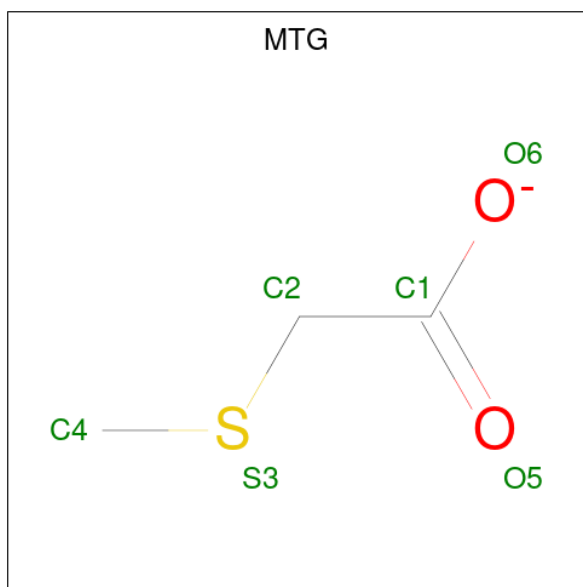
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	383	Total	C	N	O	S	0	0	0
			2991	1905	502	574	10			
1	B	383	Total	C	N	O	S	0	0	0
			2991	1905	502	574	10			

- 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
3	A	2	Total Cl 2 2	0	0
3	B	2	Total Cl 2 2	0	0

- Molecule 4 is [METHYLTHIO]ACETATE (CCD ID: MTG) (formula: $C_3H_5O_2S$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O S 6 3 2 1	0	0
4	B	1	Total C O S 6 3 2 1	0	0

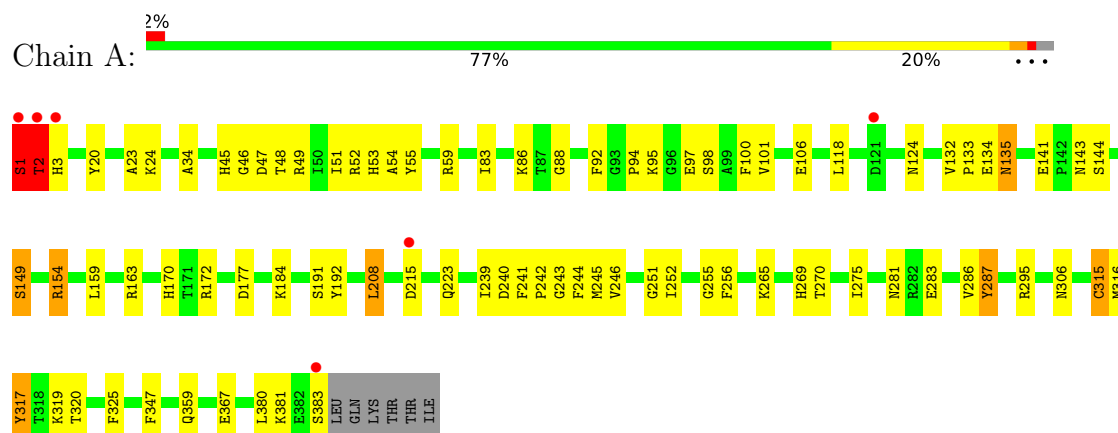
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	275	Total O 275 275	0	0
5	B	253	Total O 253 253	0	0

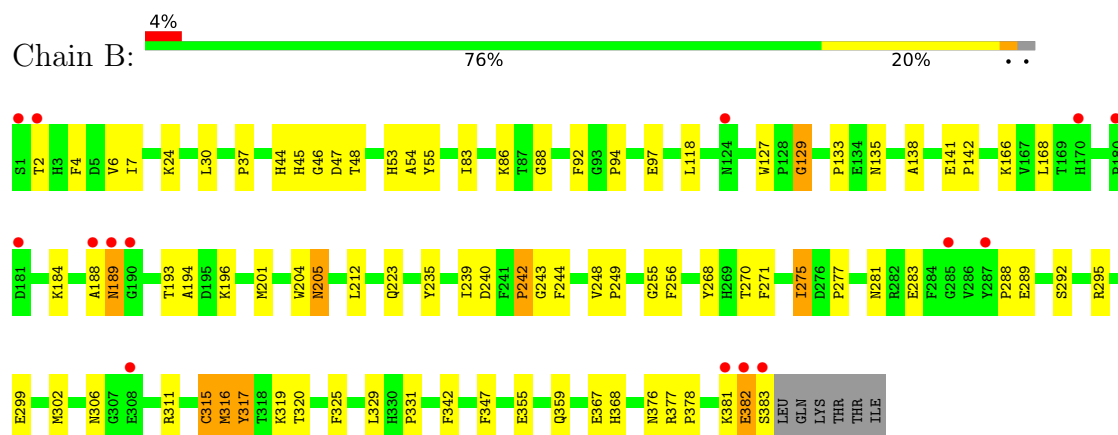
3 Residue-property plots

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: Monomeric sarcosine oxidase



- Molecule 1: Monomeric sarcosine oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.98Å 69.48Å 72.75Å 90.00° 92.33° 90.00°	Depositor
Resolution (Å)	29.73 – 1.85 29.73 – 1.85	Depositor EDS
% Data completeness (in resolution range)	88.4 (29.73-1.85) 88.6 (29.73-1.85)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.02 (at 1.85Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.208 , 0.236 0.208 , 0.235	Depositor DCC
R_{free} test set	2877 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.565	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 40.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.002 for l,k,-h 0.036 for h,-k,-l 0.025 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6632	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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, CL, MTG

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	0.53	5/3066 (0.2%)	1.06	21/4152 (0.5%)
1	B	0.37	0/3066	1.01	24/4152 (0.6%)
All	All	0.46	5/6132 (0.1%)	1.04	45/8304 (0.5%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	134	GLU	C-O	-9.08	1.12	1.24
1	A	1	SER	CA-CB	-8.50	1.36	1.53
1	A	2	THR	C-N	6.70	1.42	1.33
1	A	2	THR	CA-CB	6.26	1.64	1.53
1	A	1	SER	N-CA	-5.89	1.35	1.46

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	243	GLY	N-CA-C	-10.80	97.91	112.82
1	A	243	GLY	N-CA-C	-9.44	99.79	112.82
1	A	2	THR	CA-CB-OG1	9.17	123.35	109.60
1	B	54	ALA	N-CA-C	-7.97	98.09	109.96
1	A	241	PHE	N-CA-C	-7.59	100.19	109.83
1	B	88	GLY	N-CA-C	-7.40	103.36	112.68
1	A	55	TYR	N-CA-C	7.27	120.63	108.20
1	A	149	SER	N-CA-C	7.08	119.60	111.11
1	A	2	THR	CA-C-O	-6.72	110.90	120.51
1	A	269	HIS	N-CA-C	6.63	119.85	111.69
1	B	188	ALA	N-CA-C	-6.48	104.31	111.82
1	B	316	MET	N-CA-C	6.42	119.86	109.40
1	B	240	ASP	N-CA-C	6.25	120.25	112.24
1	A	240	ASP	N-CA-C	6.22	120.20	112.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	GLY	N-CA-C	-6.19	104.88	112.68
1	B	235	TYR	N-CA-C	6.15	120.92	113.17
1	A	54	ALA	N-CA-C	-6.03	100.47	109.15
1	B	48	THR	N-CA-C	6.02	119.35	109.72
1	B	55	TYR	N-CA-C	6.00	118.52	108.02
1	A	208	LEU	N-CA-C	5.97	120.56	113.28
1	A	320	THR	N-CA-C	-5.94	101.66	110.46
1	A	48	THR	N-CA-C	5.76	118.94	109.85
1	B	242	PRO	N-CA-C	5.75	119.69	111.13
1	A	316	MET	N-CA-C	5.74	118.83	109.24
1	B	320	THR	N-CA-C	-5.72	102.00	110.46
1	A	287	TYR	N-CA-C	-5.63	101.64	110.14
1	A	2	THR	CB-CA-C	-5.62	99.24	110.42
1	B	83	ILE	N-CA-C	5.58	116.36	111.56
1	A	317	TYR	N-CA-C	-5.49	101.64	110.14
1	B	292	SER	N-CA-C	5.44	117.29	111.36
1	B	37	PRO	N-CA-C	5.43	117.32	110.70
1	A	83	ILE	N-CA-C	5.41	116.21	111.56
1	B	319	LYS	N-CA-C	5.35	118.82	110.32
1	B	317	TYR	N-CA-C	-5.31	101.91	110.14
1	A	319	LYS	N-CA-C	5.29	118.74	110.32
1	B	275	ILE	N-CA-C	5.29	114.98	108.53
1	B	129	GLY	N-CA-C	-5.27	108.21	114.48
1	A	134	GLU	CG-CD-OE1	5.25	130.48	118.40
1	B	212	LEU	N-CA-C	-5.23	106.21	112.59
1	B	329	LEU	N-CA-C	-5.17	103.57	110.55
1	A	251	GLY	N-CA-C	5.17	118.31	110.97
1	B	83	ILE	CB-CA-C	-5.16	106.54	111.44
1	B	205	ASN	N-CA-C	-5.10	105.84	111.71
1	B	46	GLY	N-CA-C	-5.09	108.04	115.63
1	B	44	HIS	N-CA-C	5.04	120.05	113.30

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	2991	0	2891	50	0
1	B	2991	0	2891	49	0
2	A	53	0	30	1	0
2	B	53	0	29	3	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	6	0	5	1	0
4	B	6	0	5	0	0
5	A	275	0	0	2	0
5	B	253	0	0	3	0
All	All	6632	0	5851	99	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 8.

All (99) 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:A:1:SER:HB3	1:A:191:SER:O	1.44	1.14
1:B:295:ARG:O	1:B:299:GLU:HG2	1.69	0.92
1:A:124:ASN:HD21	1:A:132:VAL:H	1.30	0.78
1:B:30:LEU:HD11	1:B:168:LEU:HD13	1.73	0.70
1:A:244:PHE:CZ	1:A:255:GLY:HA3	2.27	0.68
1:A:53:HIS:HB2	1:A:141:GLU:OE1	1.97	0.65
1:B:189:ASN:HD22	1:B:189:ASN:N	1.97	0.63
1:B:244:PHE:CZ	1:B:255:GLY:HA3	2.33	0.63
1:A:20:TYR:CE1	1:A:24:LYS:HD2	2.33	0.63
1:B:295:ARG:HH11	1:B:306:ASN:ND2	1.98	0.61
1:B:94:PRO:HB2	1:B:97:GLU:HB2	1.83	0.60
1:B:189:ASN:HD22	1:B:189:ASN:H	1.47	0.60
1:A:172:ARG:NH1	5:A:1305:HOH:O	2.35	0.58
1:A:2:THR:HG22	1:A:3:HIS:N	2.18	0.58
1:A:135:ASN:H	1:A:135:ASN:HD22	1.52	0.57
1:A:281:ASN:C	1:A:281:ASN:HD22	2.11	0.57
1:B:47:ASP:OD2	1:B:311:ARG:NH1	2.36	0.57
1:B:4:PHE:CE2	1:B:30:LEU:HD22	2.40	0.56
1:B:382:GLU:HG2	1:B:383:SER:N	2.20	0.55
1:B:204:TRP:CE3	2:B:400:FAD:H8A	2.43	0.54
1:A:380:LEU:O	1:A:381:LYS:HB2	2.09	0.53
1:B:288:PRO:HG2	1:B:289:GLU:OE2	2.09	0.53
1:B:7:ILE:HG13	1:B:194:ALA:HB2	1.92	0.52
1:A:177:ASP:HB3	1:A:184:LYS:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:ARG:HH11	1:A:306:ASN:ND2	2.07	0.51
1:A:20:TYR:CZ	1:A:24:LYS:HD2	2.45	0.51
1:B:281:ASN:HD21	1:B:283:GLU:HB3	1.75	0.51
1:A:100:PHE:CE1	1:A:245:MET:HE2	2.46	0.50
1:B:289:GLU:H	1:B:289:GLU:CD	2.19	0.50
1:B:248:VAL:HB	1:B:249:PRO:HD2	1.93	0.50
1:A:256:PHE:CD2	1:A:265:LYS:HE2	2.47	0.50
1:A:2:THR:HG22	1:A:3:HIS:O	2.12	0.49
1:A:86:LYS:HD2	1:A:143:ASN:HA	1.93	0.49
1:A:215:ASP:HB3	1:A:383:SER:O	2.12	0.49
1:A:51:ILE:O	1:A:144:SER:HB2	2.13	0.48
1:B:239:ILE:O	1:B:239:ILE:HG13	2.14	0.48
1:B:355:GLU:OE2	1:B:368:HIS:HE1	1.96	0.48
1:B:53:HIS:HB2	1:B:141:GLU:OE2	2.13	0.48
1:B:359:GLN:HE22	1:B:367:GLU:HG2	1.78	0.48
1:B:184:LYS:HG3	1:B:193:THR:HG22	1.94	0.48
1:B:381:LYS:HA	1:B:381:LYS:HD3	1.64	0.48
1:B:347:PHE:HB2	2:B:400:FAD:H2'	1.96	0.47
1:B:92:PHE:HB3	1:B:138:ALA:HB2	1.97	0.47
1:B:295:ARG:HH11	1:B:306:ASN:HD21	1.63	0.47
1:A:275:ILE:HD12	1:A:275:ILE:C	2.40	0.47
1:B:315:CYS:SG	1:B:316:MET:N	2.88	0.47
1:A:59:ARG:NH2	1:A:106:GLU:HG2	2.30	0.47
1:B:347:PHE:H	2:B:400:FAD:H2'	1.80	0.46
1:B:325:PHE:N	1:B:325:PHE:CD1	2.82	0.46
1:B:223:GLN:HG3	1:B:317:TYR:CG	2.50	0.46
1:A:49:ARG:HG3	1:A:149:SER:HB3	1.98	0.46
1:A:223:GLN:HG3	1:A:317:TYR:CG	2.51	0.46
1:B:201:MET:H	1:B:205:ASN:HD21	1.64	0.46
1:B:6:VAL:HG22	1:B:196:LYS:HB2	1.98	0.46
1:A:100:PHE:CZ	1:A:245:MET:HE2	2.51	0.45
1:A:1:SER:HA	1:A:192:TYR:CE1	2.51	0.45
1:A:133:PRO:HB2	1:A:135:ASN:ND2	2.32	0.45
1:A:154:ARG:HB2	1:A:154:ARG:HH11	1.80	0.45
1:B:127:TRP:CD1	1:B:242:PRO:HD3	2.51	0.45
1:B:381:LYS:NZ	5:B:1264:HOH:O	2.47	0.45
1:B:189:ASN:N	1:B:189:ASN:ND2	2.64	0.45
1:B:268:TYR:CE2	1:B:271:PHE:HB3	2.52	0.44
1:A:98:SER:OG	1:A:101:VAL:HG23	2.17	0.44
1:B:378:PRO:HD2	5:B:1476:HOH:O	2.17	0.44
1:A:281:ASN:ND2	1:A:283:GLU:H	2.15	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:LYS:HG2	5:B:1368:HOH:O	2.18	0.44
1:A:95:LYS:HB2	1:A:135:ASN:O	2.17	0.44
1:A:159:LEU:O	1:A:163:ARG:HG2	2.18	0.43
1:A:46:GLY:O	1:A:47:ASP:HB2	2.18	0.43
1:A:34:ALA:O	1:A:170:HIS:HA	2.19	0.43
1:B:248:VAL:HB	1:B:249:PRO:CD	2.48	0.43
1:A:239:ILE:O	1:A:239:ILE:HG13	2.19	0.42
1:A:45:HIS:CE1	1:A:315:CYS:HB2	2.54	0.42
1:A:92:PHE:H	1:A:246:VAL:HG12	1.82	0.42
1:B:377:ARG:HA	1:B:378:PRO:HD3	1.85	0.42
1:A:347:PHE:HB2	2:A:400:FAD:H2'	2.00	0.42
1:B:325:PHE:CD1	1:B:342:PHE:HA	2.55	0.42
1:B:256:PHE:CD2	1:B:256:PHE:N	2.87	0.42
1:A:1:SER:O	1:A:2:THR:OG1	2.28	0.42
1:A:281:ASN:HD22	1:A:283:GLU:H	1.68	0.42
1:A:94:PRO:HB2	1:A:97:GLU:HB2	2.01	0.41
1:A:256:PHE:CE2	1:A:265:LYS:HE2	2.55	0.41
1:A:23:ALA:HB1	1:A:163:ARG:HB2	2.02	0.41
1:A:325:PHE:CD1	1:A:325:PHE:N	2.85	0.41
1:A:295:ARG:HG2	1:A:306:ASN:HD21	1.86	0.41
1:A:52:ARG:HH21	4:A:405:MTG:C1	2.34	0.41
1:B:92:PHE:CD1	1:B:92:PHE:C	2.99	0.41
1:A:245:MET:HE3	1:A:252:ILE:HG21	2.03	0.41
1:A:286:VAL:HG23	1:A:287:TYR:CD2	2.56	0.41
1:B:45:HIS:CE1	1:B:315:CYS:HB2	2.55	0.41
1:B:133:PRO:HB2	1:B:135:ASN:ND2	2.36	0.41
1:B:204:TRP:CE2	1:B:277:PRO:HB2	2.56	0.41
1:B:275:ILE:C	1:B:275:ILE:HD12	2.45	0.41
1:B:86:LYS:HE3	1:B:142:PRO:O	2.20	0.41
1:A:154:ARG:HH11	1:A:154:ARG:CB	2.34	0.40
1:A:2:THR:HG23	5:A:1254:HOH:O	2.21	0.40
1:B:129:GLY:O	1:B:302:MET:HA	2.21	0.40
1:B:331:PRO:HB3	1:B:376:ASN:HD21	1.86	0.40
1:A:359:GLN:HE22	1:A:367:GLU:HG2	1.86	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	381/389 (98%)	370 (97%)	10 (3%)	1 (0%)	36	25
1	B	381/389 (98%)	368 (97%)	13 (3%)	0	100	100
All	All	762/778 (98%)	738 (97%)	23 (3%)	1 (0%)	48	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	THR

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	320/326 (98%)	312 (98%)	8 (2%)	42	27
1	B	320/326 (98%)	313 (98%)	7 (2%)	45	32
All	All	640/652 (98%)	625 (98%)	15 (2%)	44	30

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	118	LEU
1	A	135	ASN
1	A	154	ARG
1	A	208	LEU

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Mol	Chain	Res	Type
1	A	242	PRO
1	A	270	THR
1	A	315	CYS
1	B	2	THR
1	B	24	LYS
1	B	118	LEU
1	B	189	ASN
1	B	270	THR
1	B	315	CYS
1	B	382	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	21	GLN
1	A	111	HIS
1	A	124	ASN
1	A	135	ASN
1	A	143	ASN
1	A	189	ASN
1	A	205	ASN
1	A	213	ASN
1	A	219	GLN
1	A	281	ASN
1	A	293	ASN
1	A	306	ASN
1	A	359	GLN
1	A	368	HIS
1	B	21	GLN
1	B	135	ASN
1	B	143	ASN
1	B	189	ASN
1	B	205	ASN
1	B	219	GLN
1	B	281	ASN
1	B	293	ASN
1	B	306	ASN
1	B	333	HIS
1	B	359	GLN
1	B	368	HIS
1	B	376	ASN

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	400	1	58,58,58	2.35	18 (31%)	85,89,89	1.59	13 (15%)
2	FAD	A	400	1	58,58,58	2.40	17 (29%)	85,89,89	1.58	14 (16%)
4	MTG	B	405	-	5,5,5	0.97	0	4,5,5	1.50	0
4	MTG	A	405	-	5,5,5	0.93	0	4,5,5	1.63	1 (25%)

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	400	1	-	1/34/50/50	0/6/6/6
2	FAD	A	400	1	-	1/34/50/50	0/6/6/6
4	MTG	B	405	-	-	1/3/3/3	-
4	MTG	A	405	-	-	0/3/3/3	-

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	400	FAD	P-O3P	9.76	1.70	1.59
2	B	400	FAD	P-O3P	8.18	1.68	1.59
2	A	400	FAD	PA-O3P	7.75	1.67	1.59
2	B	400	FAD	PA-O3P	7.43	1.67	1.59
2	A	400	FAD	C8M-C8	-5.38	1.40	1.51
2	B	400	FAD	C8M-C8	-5.18	1.41	1.51
2	B	400	FAD	C4X-N5	4.44	1.40	1.30
2	B	400	FAD	C6-C5X	4.29	1.46	1.40
2	A	400	FAD	C4X-N5	4.08	1.39	1.30
2	A	400	FAD	C6-C5X	3.89	1.45	1.40
2	B	400	FAD	C4-N3	3.24	1.44	1.38
2	B	400	FAD	O4'-C4'	-3.11	1.36	1.43
2	A	400	FAD	O4'-C4'	-2.98	1.37	1.43
2	B	400	FAD	C1'-N10	2.84	1.54	1.47
2	A	400	FAD	C4-N3	2.83	1.44	1.38
2	B	400	FAD	C2'-C3'	2.79	1.58	1.53
2	B	400	FAD	P-O2P	-2.70	1.42	1.55
2	A	400	FAD	C2A-N3A	2.68	1.38	1.33
2	B	400	FAD	C10-N10	2.56	1.42	1.37
2	B	400	FAD	C2A-N3A	2.55	1.38	1.33
2	B	400	FAD	C9A-N10	2.49	1.45	1.41
2	A	400	FAD	P-O2P	-2.48	1.43	1.55
2	B	400	FAD	C9-C8	-2.45	1.36	1.39
2	A	400	FAD	C9A-N10	2.44	1.45	1.41
2	B	400	FAD	C1'-C2'	2.43	1.56	1.52
2	B	400	FAD	O2'-C2'	2.35	1.48	1.43
2	A	400	FAD	O2'-C2'	2.28	1.48	1.43
2	A	400	FAD	C4'-C3'	2.26	1.57	1.53
2	A	400	FAD	C8-C7	2.22	1.46	1.40
2	A	400	FAD	C9A-C5X	2.22	1.44	1.41
2	A	400	FAD	C1'-N10	2.21	1.53	1.47
2	B	400	FAD	C8-C7	2.20	1.46	1.40
2	A	400	FAD	C9-C8	-2.18	1.36	1.39
2	A	400	FAD	C2'-C3'	2.18	1.57	1.53
2	B	400	FAD	C9A-C5X	2.06	1.44	1.41

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	400	FAD	O4B-C1B-N9A	6.07	119.74	108.09
2	A	400	FAD	O4B-C1B-N9A	5.70	119.04	108.09
2	B	400	FAD	C5'-C4'-C3'	-4.57	103.60	112.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	400	FAD	C5'-C4'-C3'	-4.25	104.19	112.22
2	A	400	FAD	O4'-C4'-C5'	4.20	119.26	109.99
2	B	400	FAD	O4'-C4'-C5'	3.99	118.78	109.99
2	B	400	FAD	O3P-P-O1P	-3.92	98.90	110.70
2	A	400	FAD	C4'-C3'-C2'	-3.88	107.11	113.57
2	A	400	FAD	O3P-P-O1P	-3.70	99.58	110.70
2	B	400	FAD	C4'-C3'-C2'	-3.32	108.05	113.57
2	B	400	FAD	O2P-P-O1P	3.31	127.84	112.44
2	A	400	FAD	O2P-P-O1P	3.30	127.80	112.44
2	B	400	FAD	C8M-C8-C9	-3.24	113.87	119.57
2	B	400	FAD	O2A-PA-O3P	-3.16	98.73	107.27
2	A	400	FAD	O2A-PA-O3P	-3.15	98.75	107.27
2	A	400	FAD	C8M-C8-C9	-3.10	114.11	119.57
2	B	400	FAD	O4'-C4'-C3'	2.42	114.90	109.25
2	B	400	FAD	O2A-PA-O1A	2.35	123.39	112.44
2	A	400	FAD	C9-C8-C7	2.35	123.13	119.69
2	A	400	FAD	O2A-PA-O1A	2.33	123.27	112.44
2	A	400	FAD	C4-C4X-N5	-2.23	115.12	118.21
4	A	405	MTG	C4-S3-C2	-2.19	92.88	100.00
2	A	400	FAD	O4'-C4'-C3'	2.18	114.36	109.25
2	B	400	FAD	C9-C8-C7	2.14	122.83	119.69
2	B	400	FAD	C5X-N5-C4X	-2.13	114.64	118.09
2	A	400	FAD	C5X-N5-C4X	-2.11	114.67	118.09
2	A	400	FAD	C1'-C2'-C3'	-2.03	104.16	109.66
2	B	400	FAD	C1'-C2'-C3'	-2.02	104.17	109.66

There are no chirality outliers.

All (3) torsion outliers are listed below:

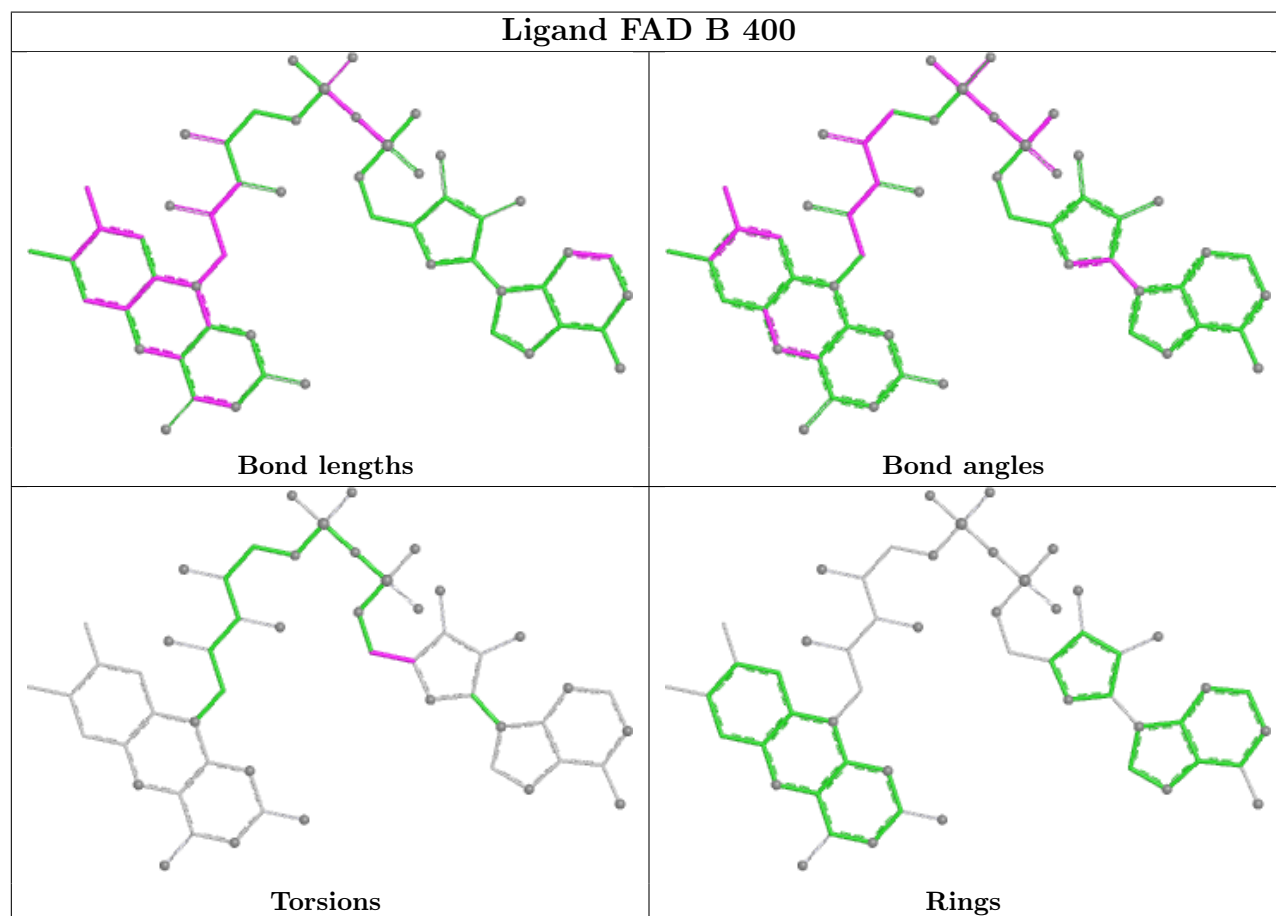
Mol	Chain	Res	Type	Atoms
2	A	400	FAD	O4B-C4B-C5B-O5B
2	B	400	FAD	O4B-C4B-C5B-O5B
4	B	405	MTG	O6-C1-C2-S3

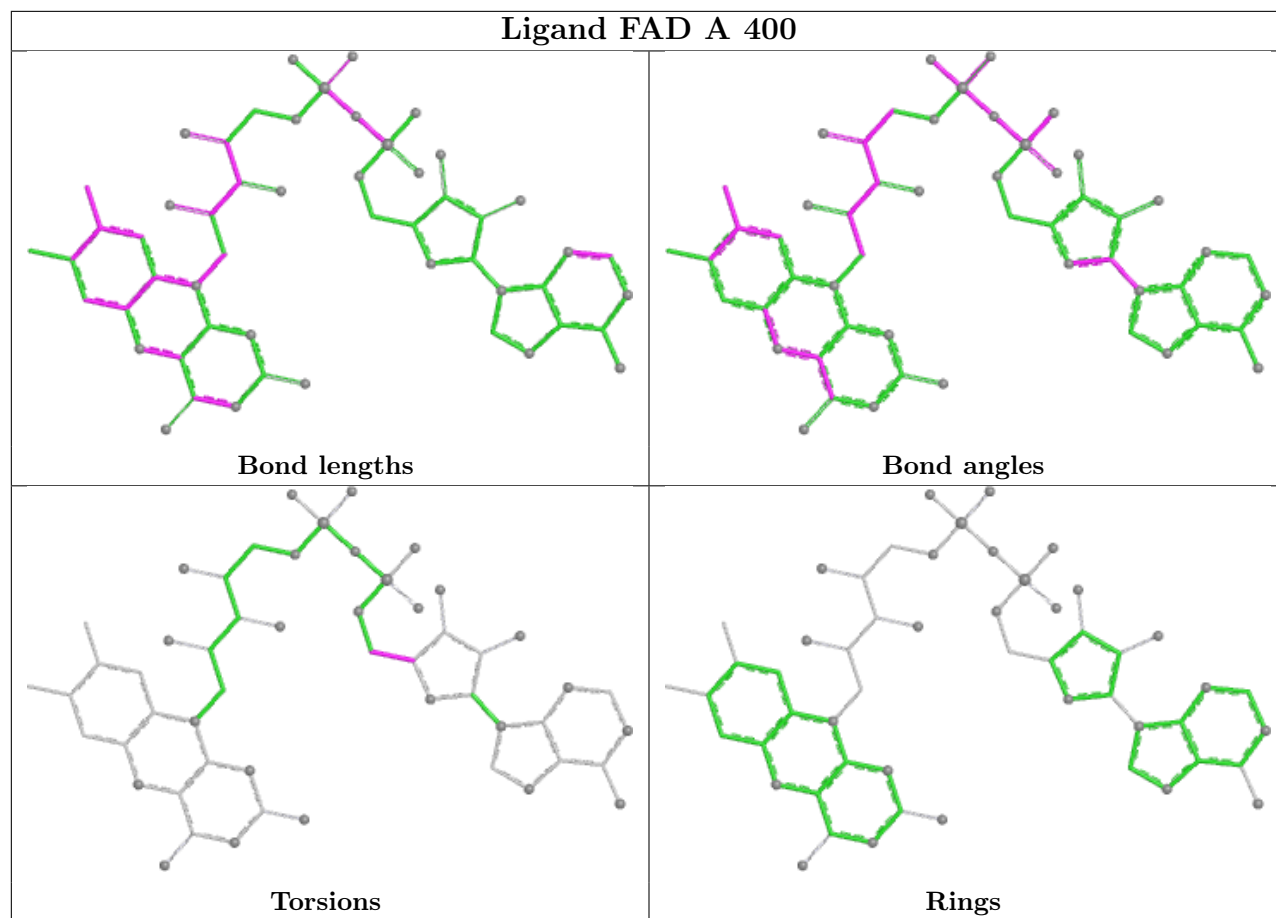
There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	400	FAD	3	0
2	A	400	FAD	1	0
4	A	405	MTG	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	383/389 (98%)	0.24	6 (1%) 70 74	17, 29, 40, 57	0
1	B	383/389 (98%)	0.52	15 (3%) 43 47	19, 31, 45, 59	0
All	All	766/778 (98%)	0.38	21 (2%) 56 59	17, 30, 44, 59	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	383	SER	4.3
1	B	2	THR	4.1
1	B	188	ALA	3.8
1	A	1	SER	3.6
1	B	189	ASN	3.0
1	A	2	THR	3.0
1	B	190	GLY	2.9
1	B	287	TYR	2.8
1	B	381	LYS	2.6
1	B	383	SER	2.5
1	B	382	GLU	2.4
1	B	285	GLY	2.4
1	A	215	ASP	2.3
1	A	3	HIS	2.3
1	B	124	ASN	2.2
1	B	170	HIS	2.2
1	B	1	SER	2.2
1	B	308	GLU	2.1
1	B	180	PRO	2.1
1	B	181	ASP	2.1
1	A	121	ASP	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

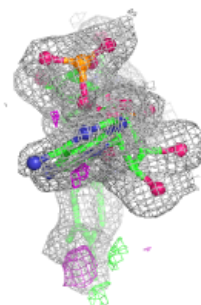
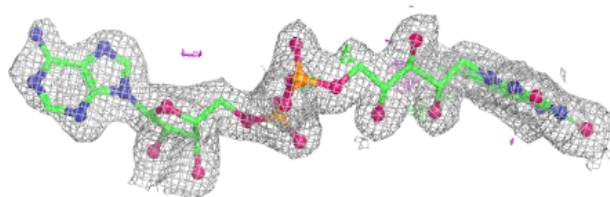
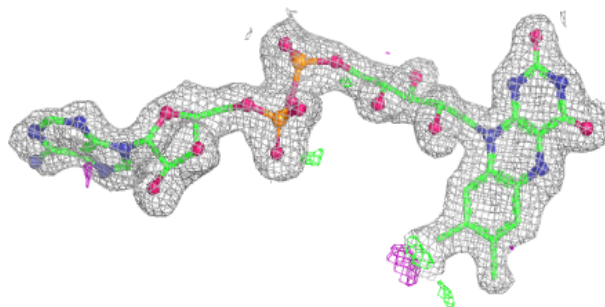
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	MTG	A	405	6/6	0.90	0.15	34,37,38,38	0
4	MTG	B	405	6/6	0.92	0.09	30,33,34,37	0
2	FAD	B	400	53/53	0.95	0.08	20,25,36,38	0
2	FAD	A	400	53/53	0.96	0.07	18,21,24,27	0
3	CL	B	404	1/1	0.96	0.14	45,45,45,45	0
3	CL	A	404	1/1	0.97	0.06	31,31,31,31	0
3	CL	B	403	1/1	0.98	0.05	27,27,27,27	0
3	CL	A	403	1/1	1.00	0.02	23,23,23,23	0

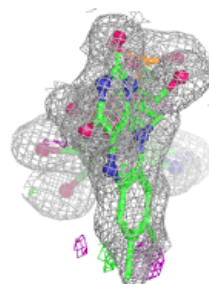
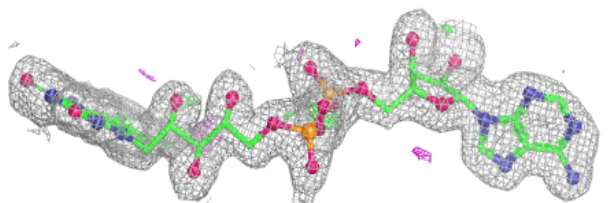
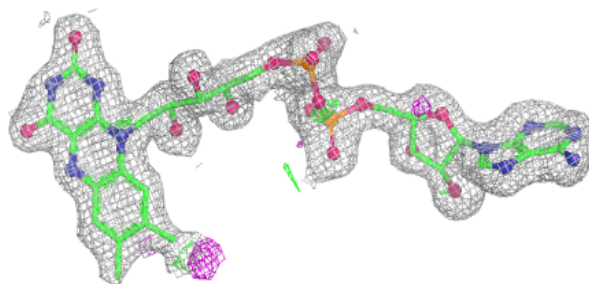
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD B 400:

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

**Electron density around FAD A 400:**

$2mF_o-DF_c$ (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.