



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:31 PM UTC

PDB ID : 5S9T / pdb_00005s9t
Title : PanDDA analysis group deposition – Crystal Structure of Trypanosoma brucei
Trypanothione reductase in complex with Z1614545742
Authors : Fiorillo, A.; Ilari, A.
Deposited on : 2021-05-18
Resolution : 1.66 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

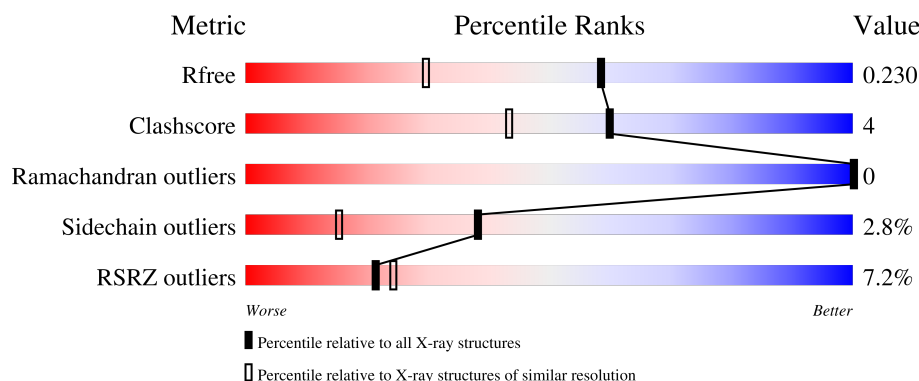
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.66 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	495	10% 86% 12% ..
1	B	495	4% 89% 8% .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	DMS	A	503	-	-	X	-

2 Entry composition ⓘ

There are 7 unique types of molecules in this entry. The entry contains 8140 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.

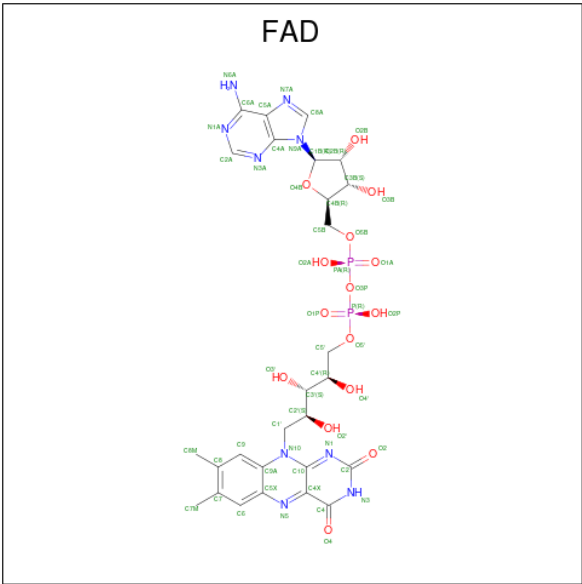
- Molecule 1 is a protein called N(1),N(8)-bis(glutathionyl)spermidine reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	486	Total	C	N	O	S	0	5	0
			3742	2383	637	703	19			
1	B	483	Total	C	N	O	S	0	2	0
			3691	2348	628	696	19			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP A0A3L6KZJ1
A	-1	SER	-	expression tag	UNP A0A3L6KZJ1
A	0	HIS	-	expression tag	UNP A0A3L6KZJ1
B	-2	GLY	-	expression tag	UNP A0A3L6KZJ1
B	-1	SER	-	expression tag	UNP A0A3L6KZJ1
B	0	HIS	-	expression tag	UNP A0A3L6KZJ1

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



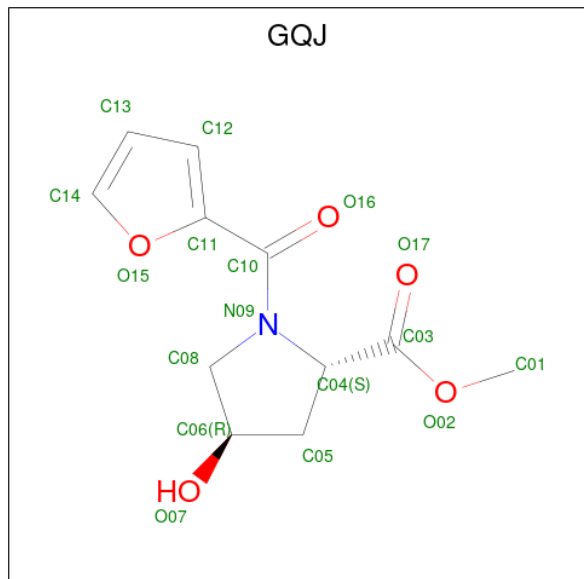
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



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

- Molecule 4 is methyl (2 {S},4 {R})-1-(furan-2-ylcarbonyl)-4-oxidanyl-pyrrolidine-2-carboxylate (CCD ID: GQJ) (formula: C₁₁H₁₃NO₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			17	11	1	5		
4	B	1	Total	C	N	O	0	0
			17	11	1	5		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		
5	B	1	Total	Mg	0	0
			1	1		

- Molecule 6 is BROMIDE ION (CCD ID: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Br	0	0
			1	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	253	Total	O	0	0
			253	253		

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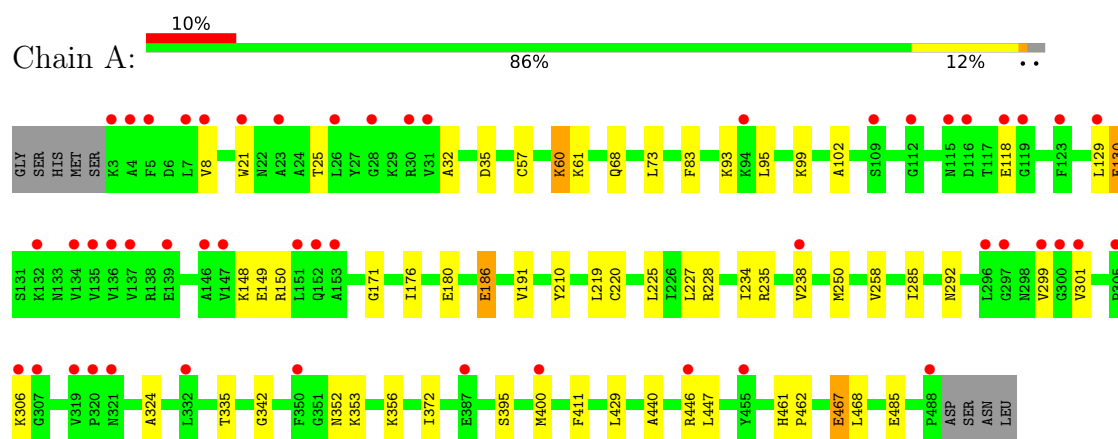
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	271	Total	O	0	0
			271	271		

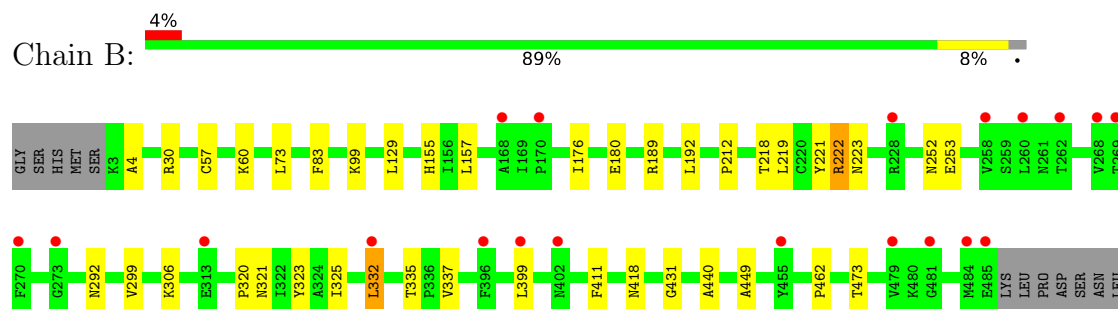
3 Residue-property plots [i](#)

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: N(1),N(8)-bis(glutathionyl)spermidine reductase



- Molecule 1: N(1),N(8)-bis(glutathionyl)spermidine reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.35Å 108.90Å 112.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	78.10 – 1.66 78.10 – 1.66	Depositor EDS
% Data completeness (in resolution range)	99.9 (78.10-1.66) 99.9 (78.10-1.66)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.18 (at 1.66Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.188 , 0.221 0.200 , 0.230	Depositor DCC
R_{free} test set	5707 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 26.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.003 for -h,l,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8140	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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: BR, MG, GQJ, FAD, DMS

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	1.05	2/3822 (0.1%)	1.28	3/5185 (0.1%)
1	B	1.03	0/3769	1.26	2/5113 (0.0%)
All	All	1.04	2/7591 (0.0%)	1.27	5/10298 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	324	ALA	C-O	9.42	1.34	1.23
1	A	342	GLY	C-O	5.17	1.29	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	473	THR	CA-CB-OG1	-6.78	99.42	109.60
1	A	467	GLU	CB-CG-CD	6.15	123.05	112.60
1	B	462	PRO	N-CD-CG	-5.70	96.96	103.80
1	A	342	GLY	CA-C-O	-5.49	115.40	121.05
1	A	342	GLY	O-C-N	5.20	127.28	122.13

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	3742	0	3749	43	0
1	B	3691	0	3692	26	0
2	A	53	0	31	1	0
2	B	53	0	31	0	0
3	A	24	0	36	9	0
3	B	16	0	24	1	0
4	A	17	0	0	1	0
4	B	17	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	B	1	0	0	0	0
7	A	253	0	0	2	0
7	B	271	0	0	1	0
All	All	8140	0	7563	65	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 4.

All (65) 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:320:PRO:O	1:B:321:ASN:OD1	1.70	1.09
1:A:68:GLN:HG2	7:A:840:HOH:O	1.90	0.71
1:A:228:ARG:HE	3:A:503:DMS:H21	1.56	0.70
1:B:253:GLU:OE1	7:B:601:HOH:O	2.09	0.69
1:A:235:ARG:HH22	3:A:503:DMS:H13	1.59	0.66
1:A:238[A]:VAL:HG21	1:A:372:ILE:HD11	1.82	0.61
1:B:129:LEU:CD1	1:B:299:VAL:HG21	2.33	0.59
1:A:21[A]:TRP:CZ2	1:A:25:THR:HG21	2.40	0.57
1:A:225:LEU:H	3:A:503:DMS:H23	1.70	0.56
1:B:320:PRO:O	1:B:321:ASN:CG	2.48	0.55
1:B:176:ILE:HB	1:B:180:GLU:HB2	1.89	0.54
1:A:73:LEU:HD23	1:B:73:LEU:HD22	1.91	0.53
1:B:155:HIS:HD2	1:B:323:TYR:OH	1.91	0.53
1:B:418:ASN:C	1:B:418:ASN:HD22	2.18	0.51
1:B:83:PHE:CE2	3:B:502:DMS:H13	2.47	0.50
1:A:60:LYS:HD2	1:A:60:LYS:C	2.37	0.50
1:A:228:ARG:NE	3:A:503:DMS:H21	2.23	0.50
1:A:234:ILE:O	1:A:238[B]:VAL:HG22	2.12	0.50
1:A:395:SER:HA	1:A:411:PHE:O	2.13	0.49
1:A:129:LEU:HD13	1:A:299:VAL:HG21	1.94	0.49
1:A:176:ILE:HB	1:A:180:GLU:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:LEU:HD13	1:B:299:VAL:HG21	1.95	0.49
1:A:83:PHE:CE2	3:A:502:DMS:H13	2.48	0.48
1:B:129:LEU:HD12	1:B:299:VAL:HG21	1.94	0.48
1:A:130:GLU:OE2	1:A:150:ARG:NE	2.44	0.48
1:A:95:LEU:HD22	1:A:210:TYR:CZ	2.49	0.47
1:A:171:GLY:HA3	1:A:258:VAL:O	2.14	0.47
1:A:225:LEU:O	3:A:503:DMS:H22	2.15	0.47
1:A:99:LYS:HD2	1:A:99:LYS:C	2.39	0.47
1:A:8:VAL:HG22	1:A:32:ALA:HB3	1.95	0.47
1:A:93:LYS:NZ	1:A:186:GLU:OE2	2.32	0.46
1:A:447:LEU:HD11	1:B:449:ALA:HB2	1.97	0.46
1:A:299:VAL:HG23	1:A:301:VAL:HG23	1.97	0.46
1:B:4:ALA:O	1:B:30:ARG:NH1	2.50	0.45
1:A:102:ALA:HB1	1:B:399:LEU:HD11	1.98	0.45
1:A:148:LYS:C	1:A:149:GLU:HG2	2.42	0.45
1:A:440:ALA:HB3	1:B:440:ALA:HB3	1.99	0.45
1:B:411:PHE:CD2	1:B:431:GLY:HA3	2.52	0.45
1:A:228:ARG:HE	3:A:503:DMS:C2	2.24	0.45
1:B:99:LYS:HD2	1:B:99:LYS:C	2.42	0.45
1:B:157:LEU:HD11	1:B:325:ILE:HG12	1.98	0.44
1:B:222:ARG:O	1:B:252:ASN:HA	2.17	0.44
1:A:73:LEU:CD2	1:B:73:LEU:HD22	2.47	0.44
1:A:227:LEU:HD12	1:A:238[B]:VAL:HG21	1.98	0.44
1:A:446[B]:ARG:HH21	1:A:446[B]:ARG:HB2	1.83	0.43
1:B:219:LEU:C	1:B:219:LEU:HD23	2.44	0.43
1:B:418:ASN:C	1:B:418:ASN:ND2	2.77	0.42
1:A:429:LEU:HD21	1:A:468:LEU:HD21	2.01	0.42
1:A:21[A]:TRP:CE2	1:A:25:THR:HG21	2.54	0.42
1:A:219:LEU:C	1:A:219:LEU:HD23	2.44	0.42
4:A:508:GQJ:O15	4:A:508:GQJ:C04	2.67	0.42
1:A:21[B]:TRP:CE2	1:A:25:THR:HG21	2.55	0.42
1:A:461:HIS:HA	1:A:462:PRO:HA	1.87	0.42
1:A:228:ARG:HE	3:A:503:DMS:H11	1.85	0.41
1:B:189:ARG:HA	1:B:212:PRO:HD2	2.02	0.41
1:A:68:GLN:NE2	7:A:618:HOH:O	2.53	0.41
1:A:228:ARG:NE	3:A:503:DMS:H11	2.35	0.41
1:A:352:ASN:HD22	1:A:352:ASN:HA	1.61	0.41
1:A:60:LYS:HD2	1:A:61:LYS:N	2.36	0.40
1:A:148:LYS:O	1:A:149:GLU:HG2	2.22	0.40
1:A:35:ASP:HA	2:A:501:FAD:N3A	2.36	0.40
1:A:220:CYS:HA	1:A:250:MET:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:221:TYR:CE2	1:B:223:ASN:HB2	2.57	0.40
1:B:192:LEU:HA	1:B:218:THR:O	2.20	0.40
1:B:332:LEU:HD13	1:B:337:VAL:HG21	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	489/495 (99%)	477 (98%)	12 (2%)	0	100	100
1	B	483/495 (98%)	474 (98%)	9 (2%)	0	100	100
All	All	972/990 (98%)	951 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	404/407 (99%)	389 (96%)	15 (4%)	30	8
1	B	398/407 (98%)	391 (98%)	7 (2%)	51	30
All	All	802/814 (98%)	780 (97%)	22 (3%)	38	16

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	CYS
1	A	60	LYS
1	A	118	GLU
1	A	130	GLU
1	A	186	GLU
1	A	191	VAL
1	A	285	ILE
1	A	292	ASN
1	A	306	LYS
1	A	335	THR
1	A	353	LYS
1	A	356	LYS
1	A	400	MET
1	A	467	GLU
1	A	485	GLU
1	B	57	CYS
1	B	60	LYS
1	B	222	ARG
1	B	292	ASN
1	B	306	LYS
1	B	332	LEU
1	B	335	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	GLN
1	A	133	ASN
1	A	152	GLN
1	A	174	HIS
1	A	241	GLN
1	A	295	GLN
1	A	352	ASN
1	A	359	HIS
1	B	68	GLN
1	B	155	HIS
1	B	352	ASN
1	B	418	ASN

5.3.3 RNA ⓘ

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 17 ligands modelled in this entry, 3 are monoatomic - leaving 14 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	A	501	-	58,58,58	0.73	0	85,89,89	0.83	2 (2%)
3	DMS	A	503	-	3,3,3	0.36	0	3,3,3	0.41	0
3	DMS	A	504	-	3,3,3	0.28	0	3,3,3	0.16	0
3	DMS	A	502	-	3,3,3	0.28	0	3,3,3	0.29	0
4	GQJ	B	506	-	18,18,18	0.50	0	25,25,25	0.51	0
3	DMS	A	507	-	3,3,3	0.30	0	3,3,3	0.04	0
4	GQJ	A	508	-	18,18,18	0.43	0	25,25,25	0.72	0
2	FAD	B	501	-	58,58,58	1.49	2 (3%)	85,89,89	0.77	1 (1%)
3	DMS	B	505	-	3,3,3	0.20	0	3,3,3	0.15	0
3	DMS	B	504	-	3,3,3	0.33	0	3,3,3	0.13	0
3	DMS	A	506	-	3,3,3	0.31	0	3,3,3	0.08	0
3	DMS	B	502	-	3,3,3	0.37	0	3,3,3	0.12	0
3	DMS	A	505	-	3,3,3	0.31	0	3,3,3	0.16	0
3	DMS	B	503	-	3,3,3	0.42	0	3,3,3	0.26	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	A	501	-	-	5/34/50/50	0/6/6/6
4	GQJ	B	506	-	-	4/14/26/26	0/2/2/2
4	GQJ	A	508	-	-	1/14/26/26	0/2/2/2
2	FAD	B	501	-	-	4/34/50/50	0/6/6/6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	FAD	P-O3P	10.14	1.70	1.59
2	B	501	FAD	C2-N1	-2.08	1.31	1.36

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	FAD	O2P-P-O1P	2.67	124.87	112.44
2	A	501	FAD	C4-N3-C2	-2.60	121.02	125.64
2	B	501	FAD	O2P-P-O1P	2.21	122.73	112.44

There are no chirality outliers.

All (14) torsion outliers are listed below:

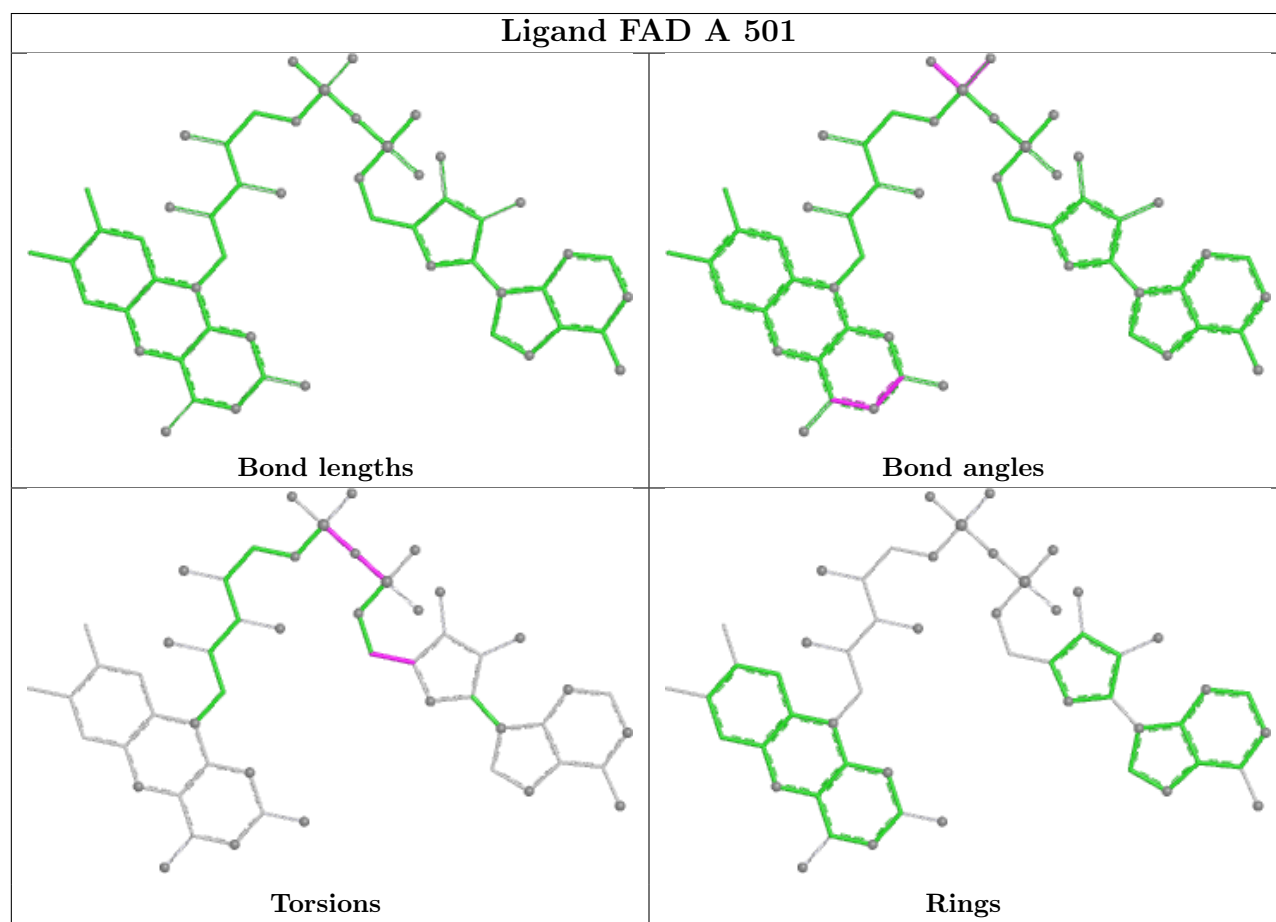
Mol	Chain	Res	Type	Atoms
2	A	501	FAD	PA-O3P-P-O1P
4	B	506	GQJ	O16-C10-C11-C12
2	A	501	FAD	PA-O3P-P-O5'
2	B	501	FAD	PA-O3P-P-O5'
4	B	506	GQJ	N09-C10-C11-O15
2	B	501	FAD	P-O3P-PA-O2A
4	A	508	GQJ	C04-C03-O02-C01
4	B	506	GQJ	N09-C10-C11-C12
2	A	501	FAD	O4B-C4B-C5B-O5B
4	B	506	GQJ	O16-C10-C11-O15
2	B	501	FAD	P-O3P-PA-O1A
2	A	501	FAD	C3B-C4B-C5B-O5B
2	B	501	FAD	O4B-C4B-C5B-O5B
2	A	501	FAD	P-O3P-PA-O2A

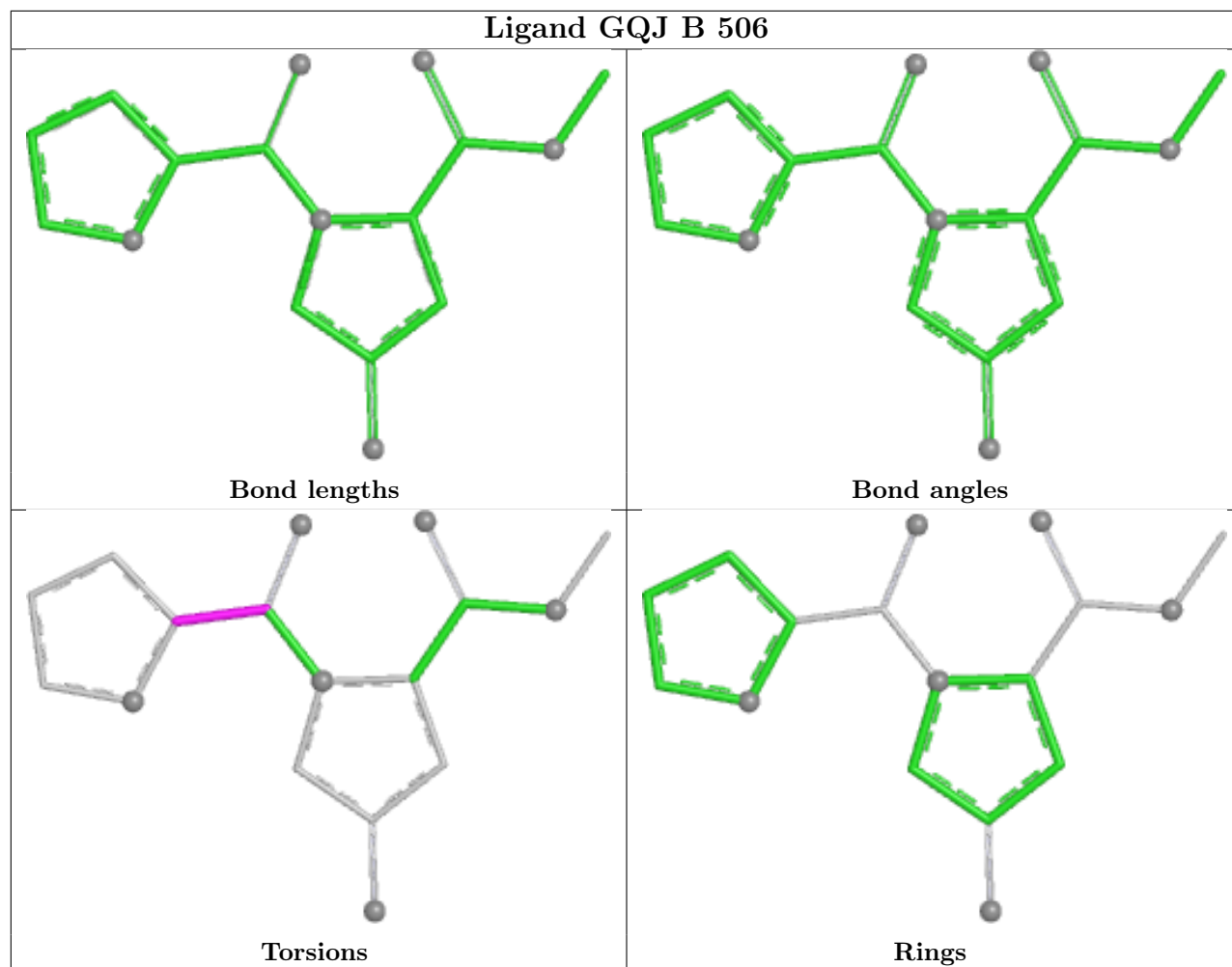
There are no ring outliers.

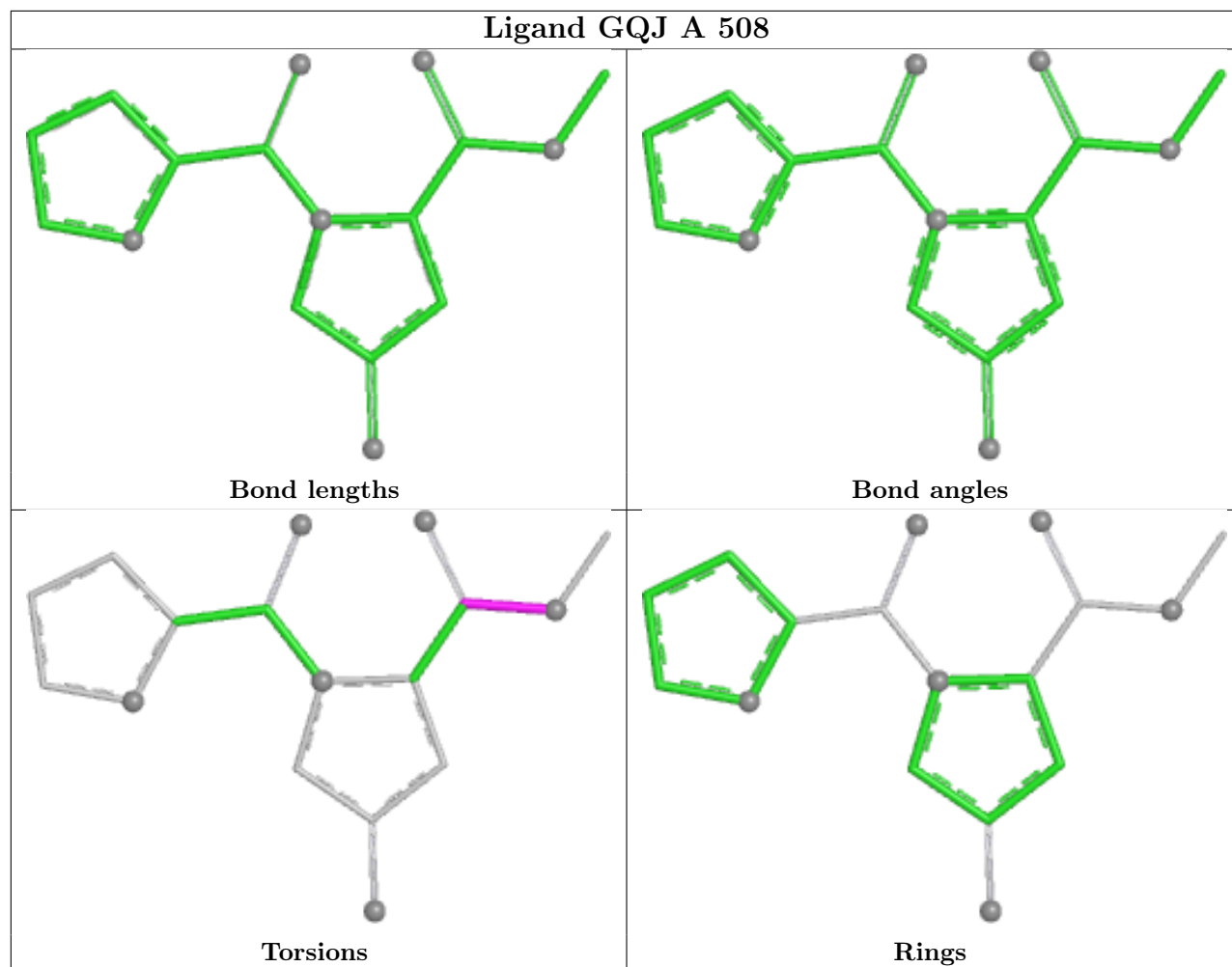
5 monomers are involved in 12 short contacts:

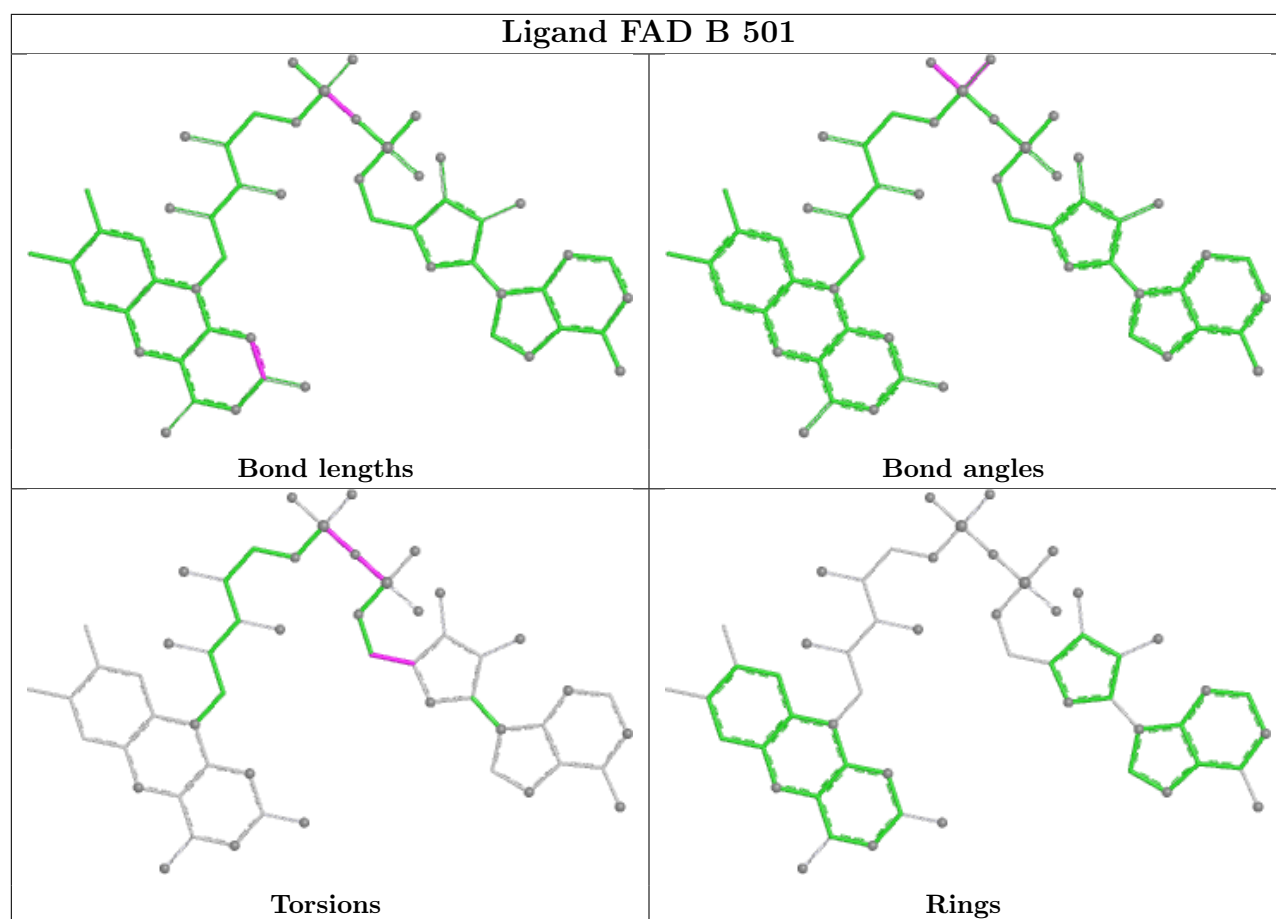
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	FAD	1	0
3	A	503	DMS	8	0
3	A	502	DMS	1	0
4	A	508	GQJ	1	0
3	B	502	DMS	1	0

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	486/495 (98%)	0.53	50 (10%)	12 12	6, 28, 50, 72	7 (1%)
1	B	483/495 (97%)	0.33	20 (4%)	41 45	7, 27, 45, 61	6 (1%)
All	All	969/990 (97%)	0.43	70 (7%)	21 24	6, 27, 49, 72	13 (1%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	455[A]	TYR	16.1
1	A	21[A]	TRP	16.0
1	A	238[A]	VAL	12.3
1	A	94[A]	LYS	10.6
1	A	446[A]	ARG	10.4
1	B	484	MET	10.2
1	B	485	GLU	10.1
1	A	387[A]	GLU	9.9
1	A	400	MET	8.0
1	B	228[A]	ARG	7.1
1	B	332	LEU	6.3
1	B	313	GLU	5.5
1	A	488	PRO	4.7
1	A	332	LEU	4.7
1	A	123	PHE	3.4
1	A	297	GLY	3.4
1	B	399	LEU	3.4
1	B	396	PHE	3.3
1	A	300	GLY	3.3
1	A	31	VAL	3.3
1	A	307	GLY	3.3
1	A	147	VAL	3.3
1	A	305	PRO	3.1
1	B	273	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	151	LEU	3.0
1	A	115	ASN	3.0
1	A	109	SER	3.0
1	B	260	LEU	2.9
1	A	112	GLY	2.9
1	A	153	ALA	2.9
1	A	5	PHE	2.9
1	A	134	VAL	2.9
1	A	3	LYS	2.9
1	A	301	VAL	2.8
1	B	269	THR	2.8
1	A	4	ALA	2.8
1	B	481	GLY	2.7
1	B	168	ALA	2.7
1	A	8	VAL	2.7
1	A	296	LEU	2.5
1	A	321	ASN	2.5
1	A	146	ALA	2.5
1	A	299	VAL	2.5
1	A	26	LEU	2.4
1	A	129	LEU	2.4
1	A	455	TYR	2.4
1	A	152	GLN	2.4
1	A	119	GLY	2.4
1	A	132	LYS	2.4
1	A	118	GLU	2.3
1	A	116	ASP	2.3
1	B	270	PHE	2.3
1	B	262	THR	2.3
1	A	28	GLY	2.3
1	A	135	VAL	2.3
1	B	258	VAL	2.3
1	B	479	VAL	2.2
1	B	402	ASN	2.2
1	A	137	VAL	2.2
1	A	30	ARG	2.2
1	A	23	ALA	2.1
1	A	306	LYS	2.1
1	A	320	PRO	2.1
1	B	170	PRO	2.1
1	A	136	VAL	2.1
1	A	319	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	268	VAL	2.1
1	A	7	LEU	2.1
1	A	350	PHE	2.1
1	A	139	GLU	2.0

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.

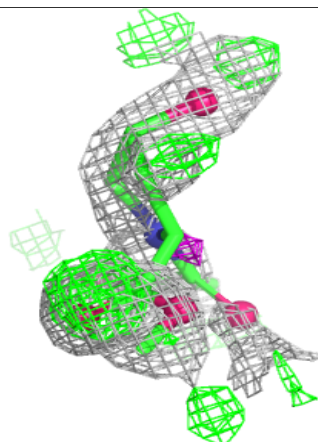
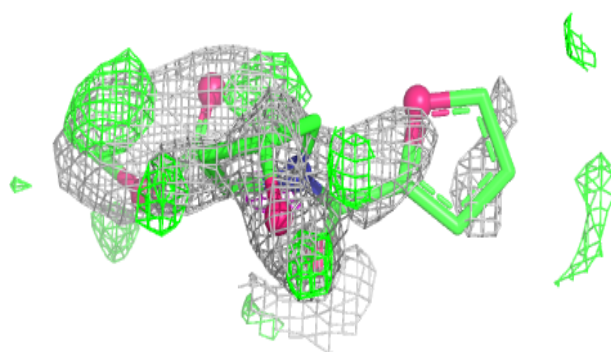
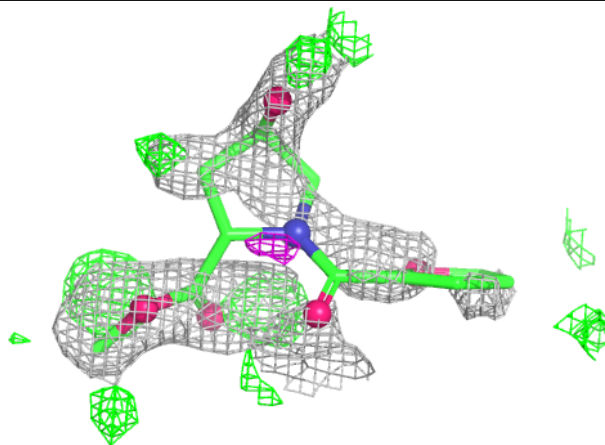
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	DMS	B	504	4/4	0.67	0.25	74,79,79,85	0
3	DMS	A	505	4/4	0.68	0.25	65,67,71,77	0
4	GQJ	B	506	17/17	0.68	0.26	37,44,57,58	17
3	DMS	A	507	4/4	0.70	0.26	80,83,86,91	0
3	DMS	A	506	4/4	0.73	0.28	82,83,83,90	0
3	DMS	A	504	4/4	0.79	0.23	76,86,91,93	0
3	DMS	A	503	4/4	0.81	0.21	55,55,60,64	0
3	DMS	B	503	4/4	0.83	0.20	49,53,61,63	0
5	MG	A	509	1/1	0.84	0.10	48,48,48,48	0
3	DMS	A	502	4/4	0.85	0.17	53,53,55,56	0
4	GQJ	A	508	17/17	0.87	0.12	24,27,29,29	17
3	DMS	B	502	4/4	0.87	0.17	52,58,60,61	0
3	DMS	B	505	4/4	0.87	0.18	70,70,70,77	0
6	BR	B	507	1/1	0.95	0.18	66,66,66,66	0
2	FAD	A	501	53/53	0.98	0.05	17,21,29,31	0
5	MG	B	508	1/1	0.98	0.10	30,30,30,30	0
2	FAD	B	501	53/53	0.98	0.04	17,20,24,29	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.

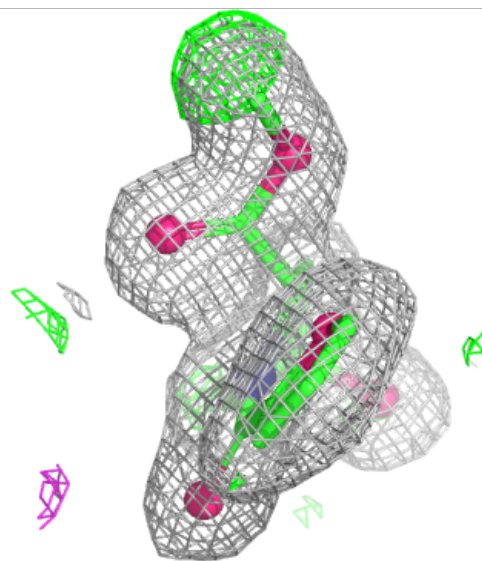
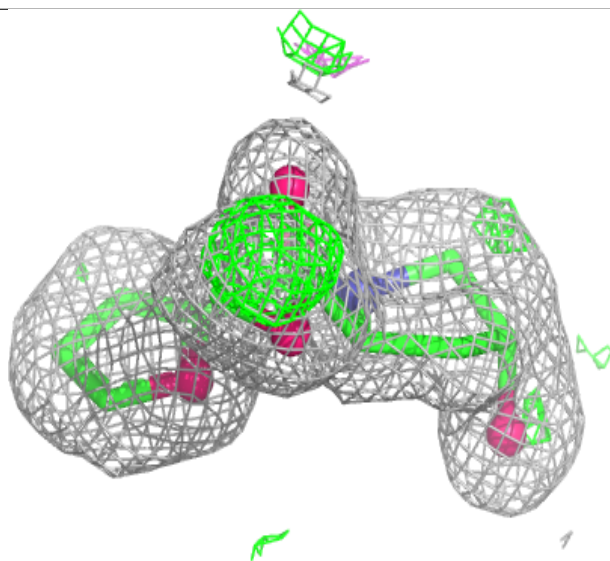
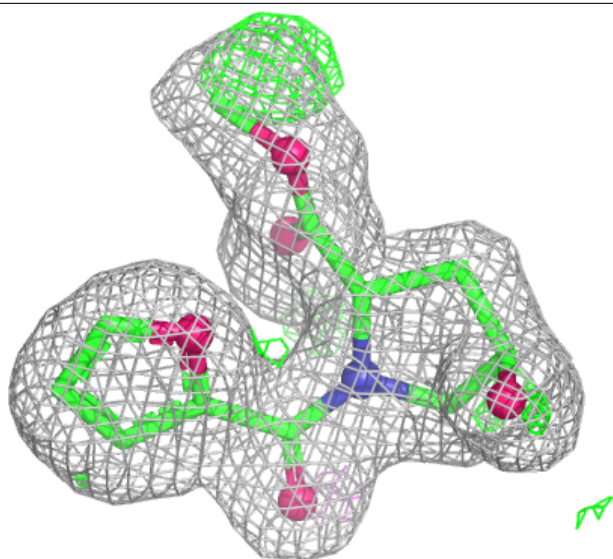
Electron density around GQJ B 506:

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



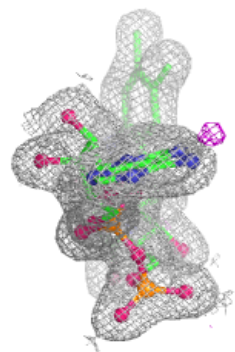
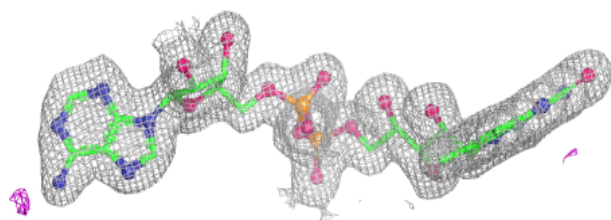
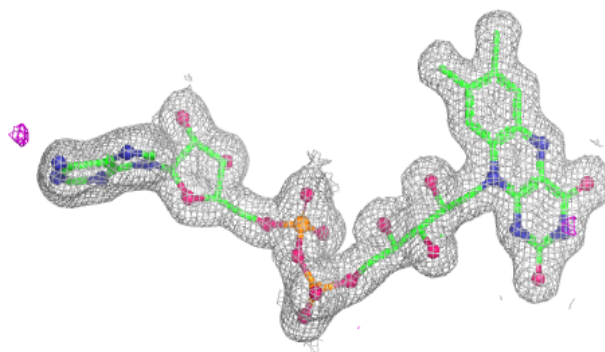
Electron density around GQJ A 508:

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

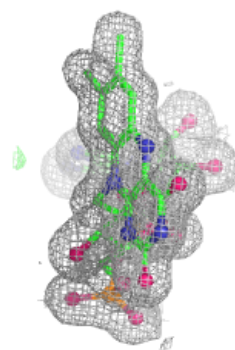
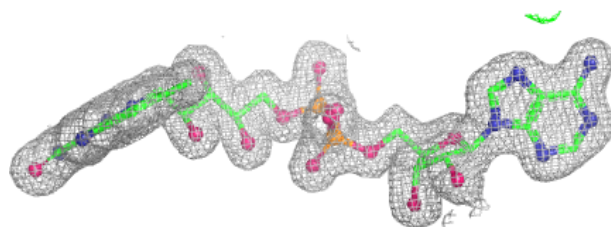
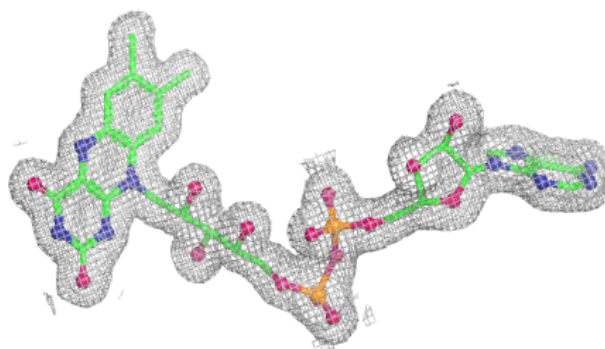


Electron density around FAD A 501:

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

**Electron density around FAD B 501:**

$2mF_o-DF_c$ (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.