



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 08:55 am BST

PDB ID : 5J60
Title : Structure of a thioredoxin reductase from *Gloeobacter violaceus*
Authors : Buey, R.M.; de Pereda, J.M.; Balsera, M.
Deposited on : 2016-04-04
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

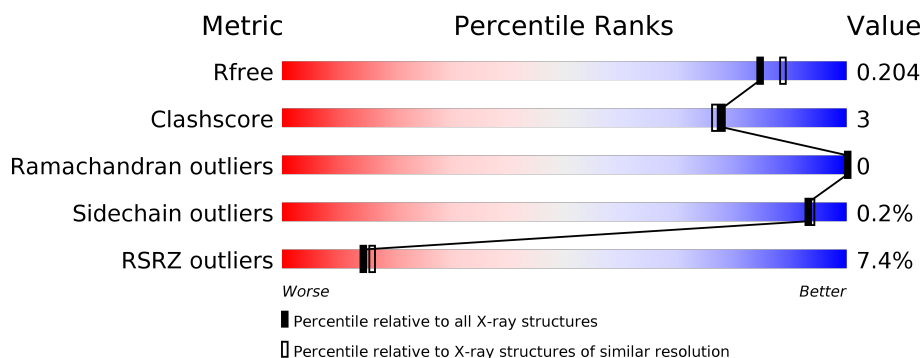
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	320	<div> <div>3%</div> <div> <div></div> <div>95%</div> <div></div> </div> <div></div> </div>
1	B	320	<div> <div>2%</div> <div> <div></div> <div>95%</div> <div></div> </div> <div></div> </div>
1	C	320	<div> <div>15%</div> <div> <div></div> <div>85%</div> <div>8%</div> </div> <div>8%</div> </div>
1	D	320	<div> <div>8%</div> <div> <div></div> <div>88%</div> <div>5%</div> </div> <div>7%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 18562 atoms, of which 8740 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 Thioredoxin reductase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	316	Total	C	H	N	O	S	0	0	0
			4577	1484	2247	398	436	12			
1	B	316	Total	C	H	N	O	S	0	0	0
			4587	1489	2252	397	437	12			
1	C	296	Total	C	H	N	O	S	0	0	0
			4208	1370	2062	371	393	12			
1	D	297	Total	C	H	N	O	S	0	0	0
			4159	1368	2021	365	393	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q7NMP6
A	-1	SER	-	expression tag	UNP Q7NMP6
A	0	HIS	-	expression tag	UNP Q7NMP6
B	-2	GLY	-	expression tag	UNP Q7NMP6
B	-1	SER	-	expression tag	UNP Q7NMP6
B	0	HIS	-	expression tag	UNP Q7NMP6
C	-2	GLY	-	expression tag	UNP Q7NMP6
C	-1	SER	-	expression tag	UNP Q7NMP6
C	0	HIS	-	expression tag	UNP Q7NMP6
D	-2	GLY	-	expression tag	UNP Q7NMP6
D	-1	SER	-	expression tag	UNP Q7NMP6
D	0	HIS	-	expression tag	UNP Q7NMP6

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



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	0	0
			83	27	30	9	15	2		
2	B	1	Total	C	H	N	O	P	0	0
			83	27	30	9	15	2		
2	C	1	Total	C	H	N	O	P	0	0
			84	27	31	9	15	2		
2	D	1	Total	C	H	N	O	P	0	0
			84	27	31	9	15	2		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		
3	A	1	Total	Ca	0	0
			1	1		

- Molecule 4 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			31	8	18	5		
4	B	1	Total	C	H	O	0	0
			31	8	18	5		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	198	Total	O	0	0
			198	198		
5	B	206	Total	O	0	0
			206	206		
5	C	107	Total	O	0	0
			107	107		
5	D	122	Total	O	0	0
			122	122		

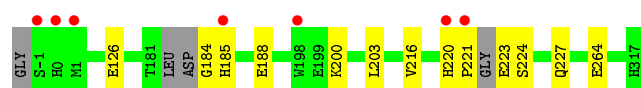
3 Residue-property plots [i](#)

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

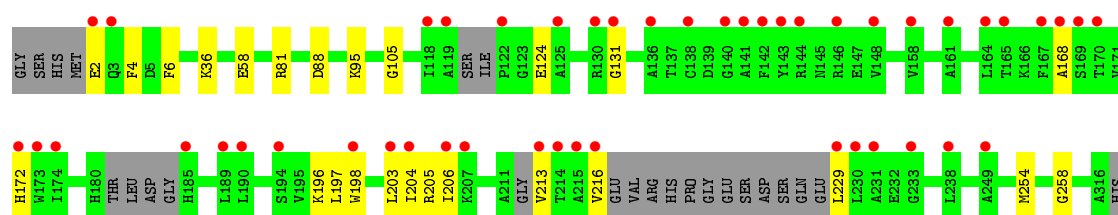
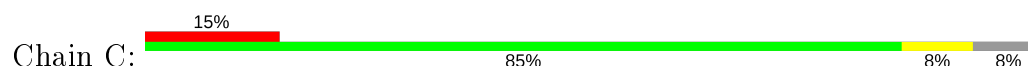
- Molecule 1: Thioredoxin reductase



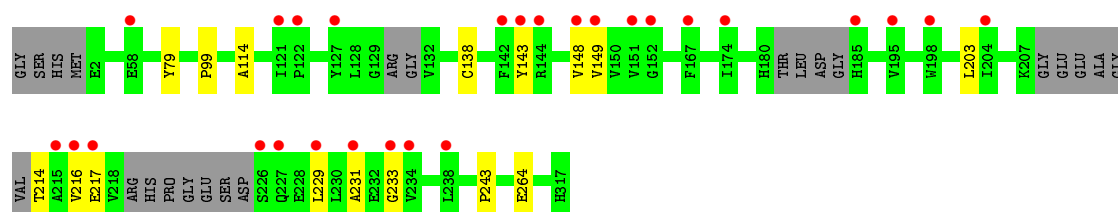
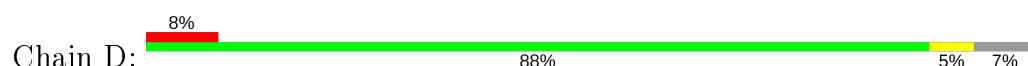
- Molecule 1: Thioredoxin reductase



- Molecule 1: Thioredoxin reductase



- Molecule 1: Thioredoxin reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.28Å 122.04Å 139.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.22 – 1.90 49.22 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.22-1.90) 100.0 (49.22-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 1.90Å)	Xtriage
Refinement program	PHENIX (dev_2341: ???)	Depositor
R, R_{free}	0.177 , 0.202 0.180 , 0.204	Depositor DCC
R_{free} test set	5599 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	33.3	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 52.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	18562	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.87 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.4061e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: CA, PG4, 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	0.42	0/2380	0.56	0/3243
1	B	0.43	0/2386	0.60	0/3251
1	C	0.37	0/2190	0.55	0/2983
1	D	0.38	0/2183	0.56	0/2979
All	All	0.40	0/9139	0.57	0/12456

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	2330	2247	2244	9	0
1	B	2335	2252	2251	9	0
1	C	2146	2062	2062	18	0
1	D	2138	2021	2016	16	0
2	A	53	30	31	0	0
2	B	53	30	31	0	0
2	C	53	31	31	0	0
2	D	53	31	31	0	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
4	A	13	18	17	1	0
4	B	13	18	17	1	0
5	A	198	0	0	5	0
5	B	206	0	0	1	1
5	C	107	0	0	2	0
5	D	122	0	0	0	0
All	All	9822	8740	8731	51	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 3.

All (51) 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:168:ALA:O	5:C:601:HOH:O	1.76	1.03
4:A:403:PG4:O5	1:D:264:GLU:OE2	1.84	0.96
1:A:251:GLN:NE2	5:A:501:HOH:O	2.08	0.84
1:D:99:PRO:HG3	1:D:138:CYS:SG	2.21	0.81
1:D:203:LEU:HD13	1:D:216:VAL:HG13	1.70	0.73
1:A:205:ARG:HD2	5:A:650:HOH:O	1.92	0.68
1:B:220:HIS:NE2	1:B:227:GLN:OE1	2.27	0.67
1:D:216:VAL:N	1:D:229:LEU:O	2.25	0.67
1:C:6:PHE:O	1:C:105:GLY:HA2	1.99	0.62
1:D:214:THR:HA	1:D:231:ALA:O	2.02	0.59
1:D:216:VAL:O	1:D:229:LEU:N	2.29	0.58
1:C:124:GLU:HA	1:C:206:ILE:HD12	1.86	0.58
1:A:254:MET:CE	5:A:506:HOH:O	2.52	0.58
1:B:184:GLY:HA2	1:B:185:HIS:C	2.25	0.57
1:D:149:VAL:HG23	1:D:231:ALA:CB	2.37	0.54
1:D:143:TYR:CD2	1:D:148:VAL:HG12	2.44	0.53
1:C:58:GLU:HA	5:C:638:HOH:O	2.08	0.52
1:A:245:THR:HG21	1:A:254:MET:HE1	1.93	0.51
1:C:196:LYS:NZ	1:C:198:TRP:CB	2.74	0.51
1:C:254:MET:CE	1:C:258:GLY:O	2.59	0.49
1:D:143:TYR:CE2	1:D:233:GLY:HA3	2.47	0.49
1:A:302:ASP:OD2	1:A:308:ARG:NH2	2.27	0.49
1:C:131:GLY:HA3	1:C:213:VAL:HG23	1.94	0.48
1:B:220:HIS:HB2	1:B:221:PRO:HD2	1.94	0.48
1:C:172:HIS:CE1	1:C:229:LEU:HD21	2.49	0.46
1:D:203:LEU:HD13	1:D:216:VAL:CG1	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2:GLU:O	1:C:4:PHE:CE1	2.70	0.45
1:D:143:TYR:HB3	1:D:148:VAL:CG1	2.48	0.44
1:D:99:PRO:HG3	1:D:138:CYS:HG	1.83	0.43
1:B:223:GLU:O	1:B:224:SER:CB	2.66	0.43
1:C:203:LEU:HD12	1:C:204:ILE:N	2.33	0.43
1:D:114:ALA:HA	1:D:243:PRO:HA	2.00	0.43
1:C:203:LEU:HD11	1:C:205:ARG:O	2.19	0.43
1:A:203:LEU:HD11	1:A:216:VAL:HB	2.00	0.43
1:C:36:LYS:O	1:C:81:ARG:HD2	2.18	0.42
1:A:302:ASP:CG	1:A:308:ARG:HH22	2.15	0.42
1:C:131:GLY:HA3	1:C:213:VAL:CG2	2.49	0.42
1:A:309:LYS:HB2	5:A:505:HOH:O	2.20	0.42
1:A:50:ALA:HA	5:A:503:HOH:O	2.19	0.42
1:C:88:ASP:HB3	1:C:95:LYS:HB2	2.02	0.42
1:B:126:GLU:HB2	5:B:612:HOH:O	2.19	0.42
1:B:203:LEU:CD1	1:B:216:VAL:HB	2.50	0.41
1:B:188:GLU:HG2	1:D:79:TYR:HB3	2.02	0.41
1:D:203:LEU:CD1	1:D:216:VAL:HG13	2.46	0.41
1:B:264:GLU:HA	4:B:501:PG4:H31	2.02	0.41
1:C:197:LEU:HD12	1:C:198:TRP:N	2.36	0.41
1:C:172:HIS:NE2	1:C:229:LEU:HD21	2.36	0.41
1:C:203:LEU:HD11	1:C:216:VAL:HB	2.02	0.40
1:B:200:LYS:O	1:B:221:PRO:HD3	2.21	0.40
1:D:216:VAL:HG12	1:D:217:GLU:N	2.36	0.40
1:C:206:ILE:HA	1:C:216:VAL:HG12	2.04	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)
5:B:683:HOH:O	5:B:702:HOH:O[3_555]	2.10	0.10

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	310/320 (97%)	301 (97%)	9 (3%)	0	100	100
1	B	310/320 (97%)	300 (97%)	10 (3%)	0	100	100
1	C	286/320 (89%)	276 (96%)	10 (4%)	0	100	100
1	D	287/320 (90%)	279 (97%)	8 (3%)	0	100	100
All	All	1193/1280 (93%)	1156 (97%)	37 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	228/252 (90%)	226 (99%)	2 (1%)	78	79
1	B	229/252 (91%)	229 (100%)	0	100	100
1	C	205/252 (81%)	205 (100%)	0	100	100
1	D	199/252 (79%)	199 (100%)	0	100	100
All	All	861/1008 (85%)	859 (100%)	2 (0%)	93	94

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	117	ARG
1	A	145	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	$\# Z > 2$
4	PG4	A	403	3	12,12,12	0.50	0	11,11,11	0.31	0
2	FAD	A	401	-	51,58,58	2.05	7 (13%)	60,89,89	2.78	12 (20%)
2	FAD	D	401	-	51,58,58	2.16	10 (19%)	60,89,89	2.82	11 (18%)
2	FAD	C	500	-	51,58,58	2.07	11 (21%)	60,89,89	2.70	12 (20%)
4	PG4	B	501	3	12,12,12	0.48	0	11,11,11	0.32	0
2	FAD	B	500	-	51,58,58	2.03	10 (19%)	60,89,89	2.86	12 (20%)

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	PG4	A	403	3	-	5/10/10/10	-
2	FAD	A	401	-	-	2/30/50/50	0/6/6/6
2	FAD	D	401	-	-	2/30/50/50	0/6/6/6
2	FAD	C	500	-	-	3/30/50/50	0/6/6/6
4	PG4	B	501	3	-	3/10/10/10	-
2	FAD	B	500	-	-	2/30/50/50	0/6/6/6

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	FAD	C4X-C10	8.72	1.47	1.38
2	D	401	FAD	C4X-C10	8.37	1.47	1.38
2	B	500	FAD	C4X-C10	7.94	1.46	1.38
2	C	500	FAD	C4X-C10	7.45	1.46	1.38
2	C	500	FAD	C4-N3	6.88	1.45	1.33
2	D	401	FAD	C4-C4X	5.69	1.51	1.41
2	D	401	FAD	C4-N3	5.66	1.42	1.33
2	A	401	FAD	C4-C4X	5.27	1.50	1.41
2	A	401	FAD	C4-N3	5.04	1.41	1.33
2	B	500	FAD	C4-N3	4.88	1.41	1.33
2	A	401	FAD	C5X-N5	4.30	1.42	1.35
2	B	500	FAD	C4-C4X	4.27	1.48	1.41
2	C	500	FAD	C4-C4X	4.26	1.48	1.41
2	B	500	FAD	C9A-N10	4.00	1.43	1.38
2	C	500	FAD	C5X-N5	3.91	1.41	1.35
2	B	500	FAD	C5X-N5	3.91	1.41	1.35
2	D	401	FAD	C5X-N5	3.19	1.40	1.35
2	C	500	FAD	PA-O1A	3.11	1.61	1.50
2	B	500	FAD	C2-N3	2.94	1.44	1.38
2	D	401	FAD	P-O1P	2.81	1.60	1.50
2	A	401	FAD	C9A-N10	2.76	1.42	1.38
2	C	500	FAD	C9A-N10	2.76	1.42	1.38
2	A	401	FAD	C3B-C4B	2.48	1.59	1.53
2	D	401	FAD	PA-O1A	2.48	1.59	1.50
2	B	500	FAD	O4'-C4'	-2.44	1.38	1.43
2	A	401	FAD	C2-N3	2.43	1.43	1.38
2	C	500	FAD	P-O1P	2.39	1.59	1.50
2	D	401	FAD	C3B-C4B	2.37	1.59	1.53
2	D	401	FAD	C5'-C4'	-2.28	1.48	1.51
2	D	401	FAD	O5'-C5'	-2.24	1.36	1.44
2	B	500	FAD	C2B-C3B	2.20	1.59	1.53
2	C	500	FAD	C3B-C4B	2.15	1.58	1.53
2	C	500	FAD	PA-O2A	-2.13	1.45	1.55
2	B	500	FAD	O4-C4	-2.07	1.19	1.24
2	C	500	FAD	O4-C4	-2.07	1.19	1.24
2	C	500	FAD	C2A-N3A	2.06	1.35	1.32
2	D	401	FAD	O4-C4	-2.05	1.19	1.24
2	B	500	FAD	P-O1P	2.02	1.58	1.50

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	FAD	C4-N3-C2	15.52	128.24	115.14
2	B	500	FAD	C4-N3-C2	15.40	128.15	115.14
2	A	401	FAD	C4-N3-C2	15.10	127.89	115.14
2	C	500	FAD	C4-N3-C2	13.77	126.77	115.14
2	D	401	FAD	C4-C4X-C10	-7.54	114.96	119.95
2	B	500	FAD	C4-C4X-C10	-7.33	115.10	119.95
2	A	401	FAD	C4X-C4-N3	-7.25	113.52	123.43
2	C	500	FAD	C4-C4X-C10	-7.13	115.23	119.95
2	D	401	FAD	C4X-C4-N3	-6.45	114.61	123.43
2	B	500	FAD	C4X-C4-N3	-6.36	114.74	123.43
2	A	401	FAD	C4-C4X-C10	-6.23	115.83	119.95
2	C	500	FAD	C10-C4X-N5	6.08	125.46	121.26
2	C	500	FAD	C4X-C4-N3	-5.95	115.30	123.43
2	B	500	FAD	C10-C4X-N5	5.78	125.25	121.26
2	B	500	FAD	C4X-C10-N10	-5.09	115.07	120.30
2	D	401	FAD	C10-C4X-N5	4.92	124.66	121.26
2	C	500	FAD	C4X-C10-N10	-4.78	115.39	120.30
2	A	401	FAD	C10-C4X-N5	4.56	124.41	121.26
2	A	401	FAD	C4X-C10-N10	-4.49	115.69	120.30
2	D	401	FAD	C4X-C10-N10	-4.24	115.94	120.30
2	C	500	FAD	P-O3P-PA	4.23	147.35	132.83
2	C	500	FAD	C1'-N10-C9A	4.07	121.50	118.29
2	D	401	FAD	C1'-N10-C9A	4.05	121.48	118.29
2	B	500	FAD	P-O3P-PA	3.92	146.29	132.83
2	A	401	FAD	P-O3P-PA	3.62	145.26	132.83
2	D	401	FAD	P-O3P-PA	3.61	145.23	132.83
2	A	401	FAD	C1'-N10-C9A	3.44	121.00	118.29
2	B	500	FAD	C7-C6-C5X	2.86	125.26	121.22
2	B	500	FAD	C1'-N10-C9A	2.63	120.36	118.29
2	B	500	FAD	C6-C5X-N5	2.57	121.88	119.05
2	A	401	FAD	C7-C6-C5X	2.57	124.84	121.22
2	A	401	FAD	C1'-N10-C10	-2.54	116.14	118.41
2	C	500	FAD	C7-C6-C5X	2.52	124.78	121.22
2	C	500	FAD	C8-C9-C9A	2.43	126.72	119.19
2	B	500	FAD	C1'-N10-C10	-2.38	116.27	118.41
2	D	401	FAD	C1'-N10-C10	-2.35	116.30	118.41
2	D	401	FAD	C7-C6-C5X	2.34	124.52	121.22
2	C	500	FAD	C1'-N10-C10	-2.30	116.35	118.41
2	A	401	FAD	C8-C9-C9A	2.29	126.30	119.19
2	C	500	FAD	O2'-C2'-C1'	-2.29	104.08	109.59
2	D	401	FAD	O2'-C2'-C1'	-2.27	104.13	109.59
2	B	500	FAD	C8-C9-C9A	2.25	126.18	119.19
2	D	401	FAD	C8-C9-C9A	2.18	125.95	119.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	FAD	O2A-PA-O1A	2.12	122.74	112.24
2	A	401	FAD	C6-C5X-N5	2.10	121.36	119.05
2	C	500	FAD	C5A-C6A-N6A	2.04	123.46	120.35
2	B	500	FAD	C9-C9A-C5X	-2.01	116.45	119.88

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	FAD	PA-O3P-P-O5'
2	B	500	FAD	PA-O3P-P-O5'
4	A	403	PG4	O1-C1-C2-O2
4	A	403	PG4	O3-C5-C6-O4
4	B	501	PG4	O1-C1-C2-O2
2	D	401	FAD	PA-O3P-P-O5'
2	C	500	FAD	PA-O3P-P-O5'
4	A	403	PG4	C6-C5-O3-C4
2	C	500	FAD	P-O3P-PA-O1A
4	B	501	PG4	C1-C2-O2-C3
2	D	401	FAD	O4B-C4B-C5B-O5B
4	B	501	PG4	C3-C4-O3-C5
2	C	500	FAD	O4B-C4B-C5B-O5B
2	B	500	FAD	O4B-C4B-C5B-O5B
4	A	403	PG4	C5-C6-O4-C7
2	A	401	FAD	O4B-C4B-C5B-O5B
4	A	403	PG4	O4-C7-C8-O5

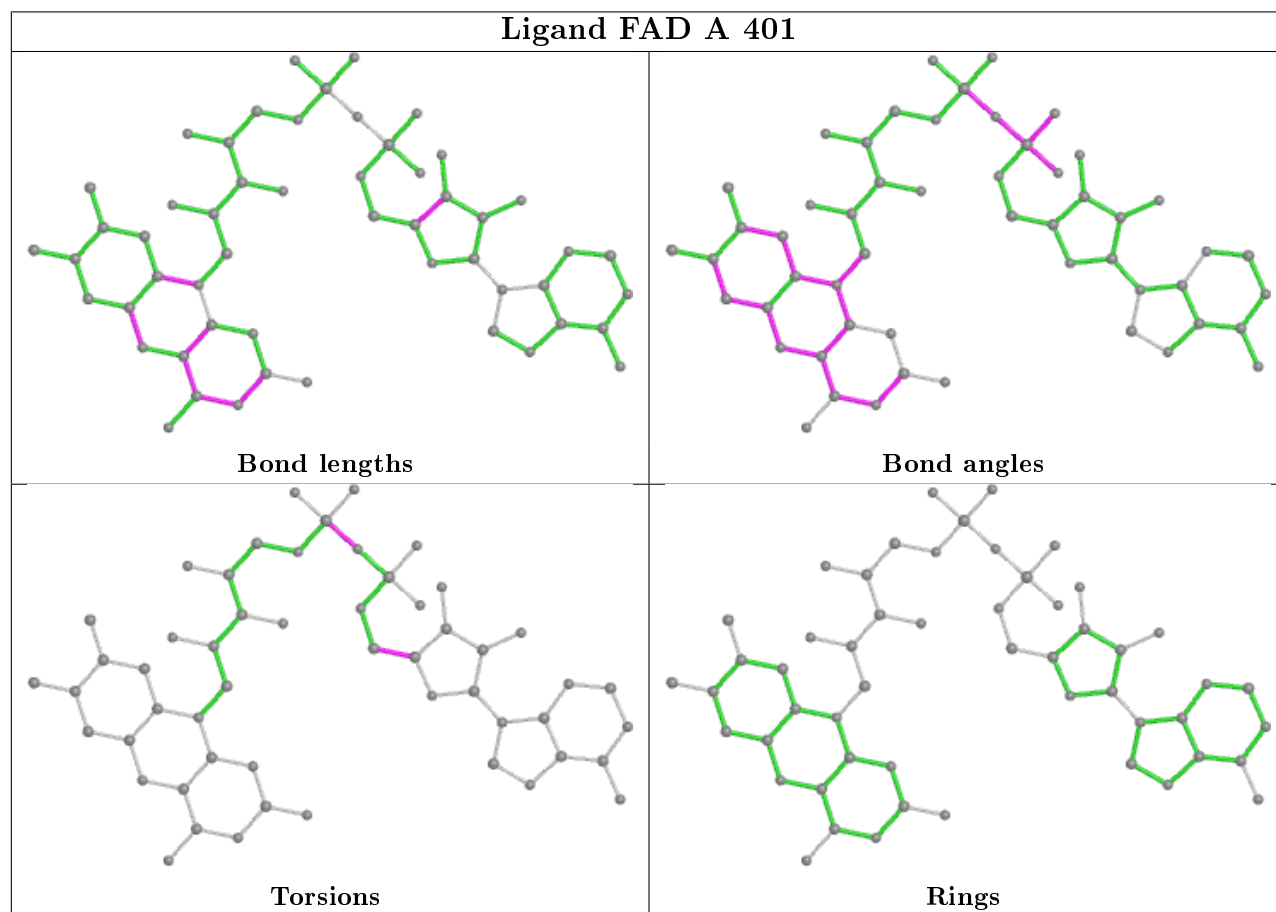
There are no ring outliers.

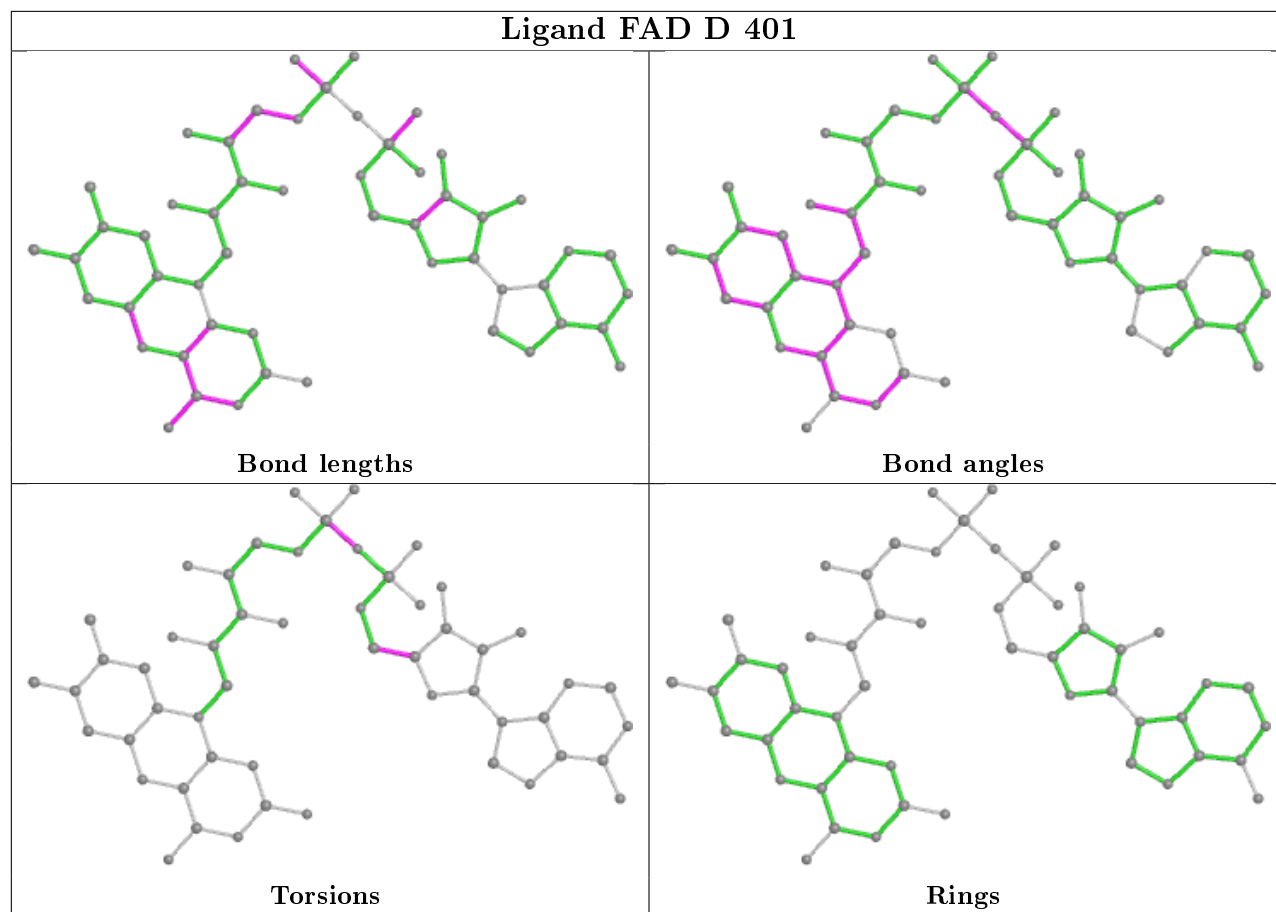
2 monomers are involved in 2 short contacts:

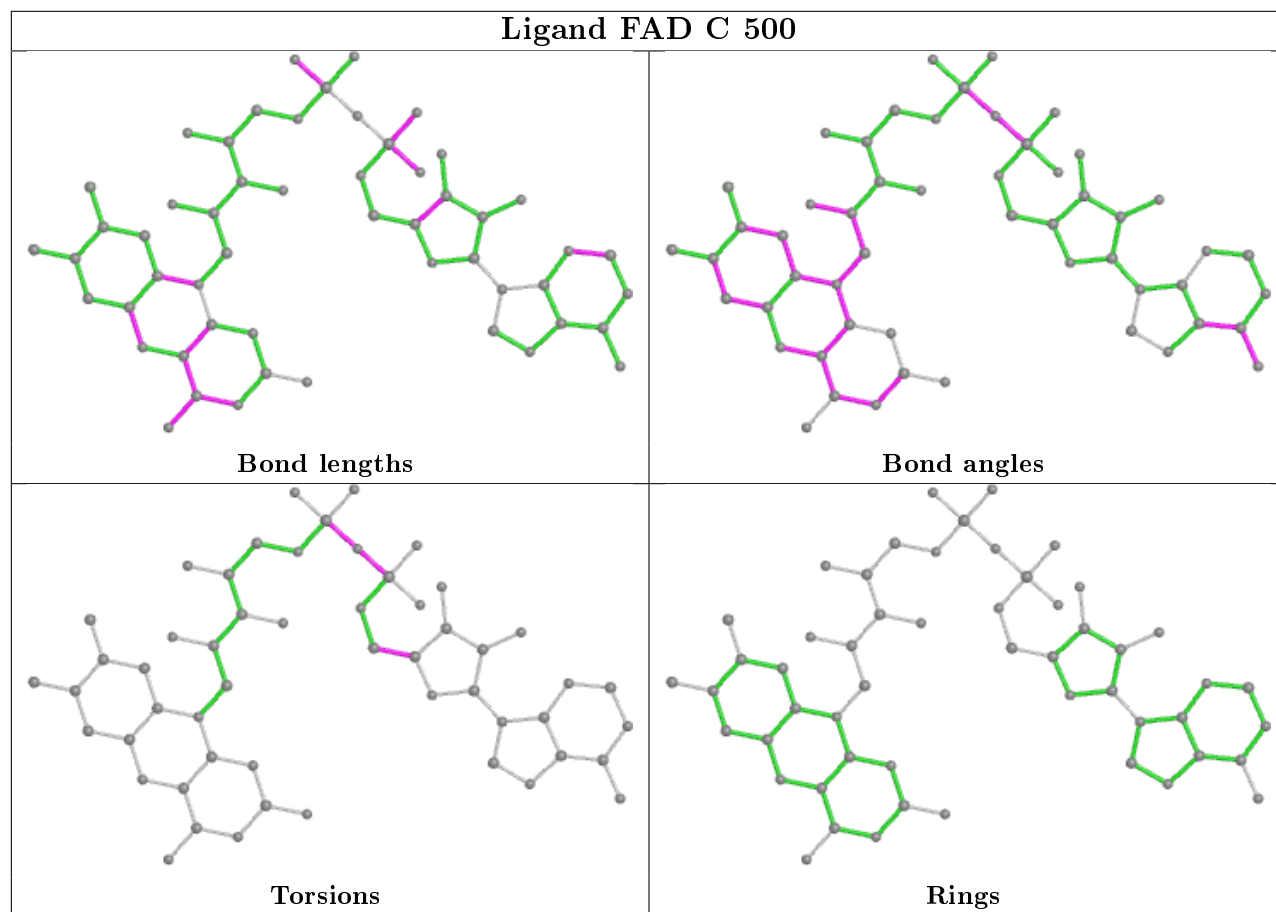
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	403	PG4	1	0
4	B	501	PG4	1	0

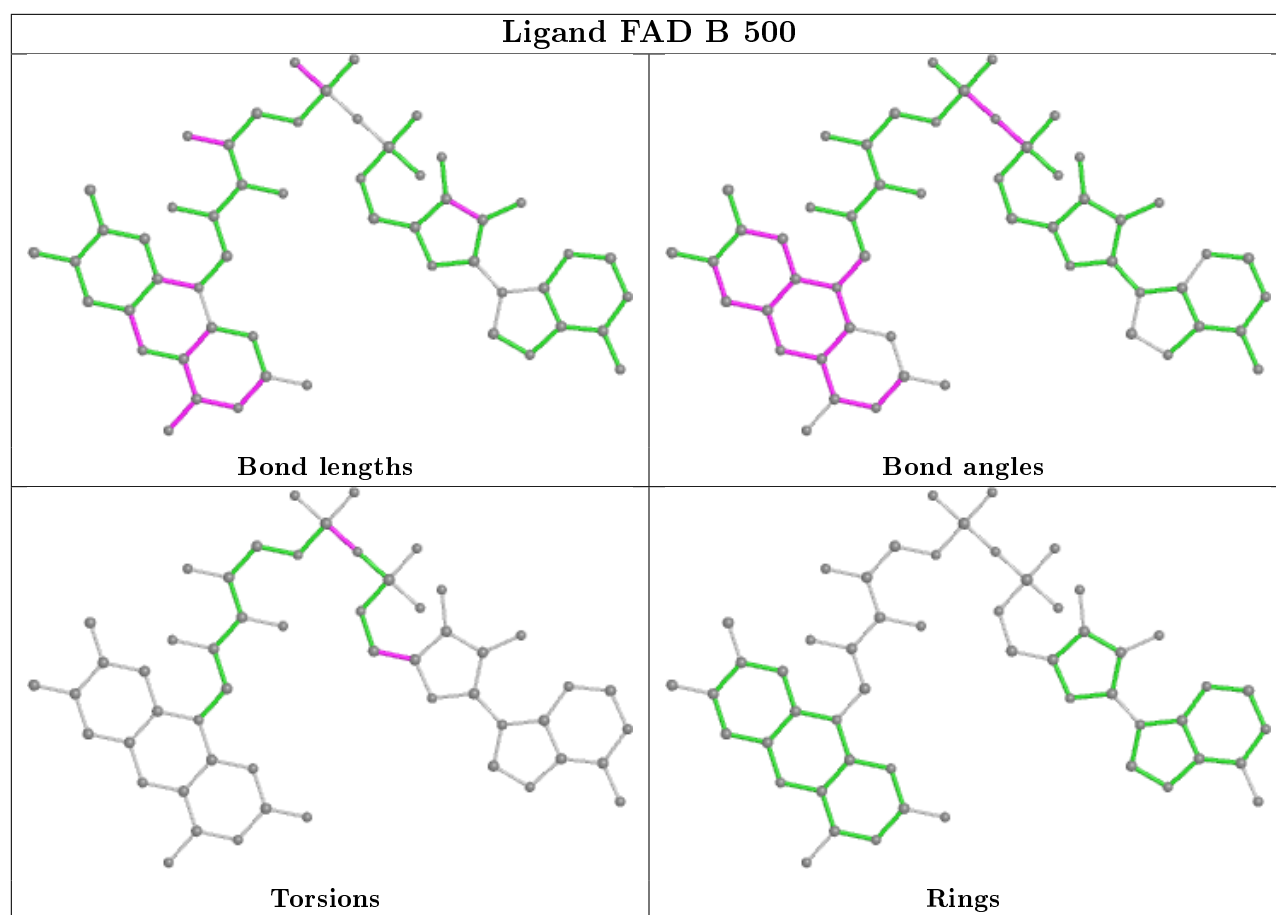
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	316/320 (98%)	0.23	10 (3%) 47 50	21, 36, 70, 117	0
1	B	316/320 (98%)	0.32	7 (2%) 62 64	22, 37, 68, 115	0
1	C	296/320 (92%)	0.95	47 (15%) 1 2	24, 51, 104, 123	0
1	D	297/320 (92%)	0.58	27 (9%) 9 10	24, 47, 99, 145	0
All	All	1225/1280 (95%)	0.51	91 (7%) 14 16	21, 40, 97, 145	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-2	GLY	7.6
1	C	167	PHE	7.3
1	C	204	ILE	6.9
1	B	-1	SER	6.4
1	D	233	GLY	6.1
1	C	174	ILE	5.9
1	D	174	ILE	5.9
1	C	206	ILE	5.8
1	C	142	PHE	5.6
1	C	161	ALA	5.5
1	D	231	ALA	5.4
1	A	-1	SER	5.1
1	B	221	PRO	5.0
1	D	148	VAL	4.7
1	C	203	LEU	4.7
1	D	234	VAL	4.7
1	C	140	GLY	4.6
1	D	215	ALA	4.5
1	B	220	HIS	4.5
1	D	216	VAL	4.5
1	C	190	LEU	4.4

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Mol	Chain	Res	Type	RSRZ
1	C	148	VAL	4.4
1	C	172	HIS	4.3
1	C	231	ALA	4.3
1	C	230	LEU	4.2
1	C	213	VAL	4.2
1	C	229	LEU	4.1
1	C	189	LEU	4.0
1	C	170	THR	4.0
1	C	216	VAL	4.0
1	B	0	HIS	3.8
1	B	1	MET	3.8
1	D	142	PHE	3.7
1	D	151	VAL	3.6
1	A	181	THR	3.6
1	C	164	LEU	3.6
1	C	198	TRP	3.6
1	C	143	TYR	3.6
1	C	215	ALA	3.5
1	D	127	TYR	3.5
1	D	198	TRP	3.5
1	C	3	GLN	3.4
1	D	149	VAL	3.4
1	C	144	ARG	3.3
1	D	143	TYR	3.3
1	C	119	ALA	3.3
1	A	0	HIS	3.3
1	C	169	SER	3.3
1	C	194	SER	3.2
1	C	165	THR	3.2
1	A	1	MET	3.1
1	C	138	CYS	3.1
1	C	185	HIS	3.1
1	D	204	ILE	3.0
1	D	58	GLU	3.0
1	D	167	PHE	2.9
1	C	122	PRO	2.9
1	C	214	THR	2.9
1	D	238	LEU	2.9
1	C	168	ALA	2.8
1	D	121	ILE	2.8
1	D	144	ARG	2.8
1	C	207	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	185	HIS	2.7
1	D	217	GLU	2.7
1	A	317	HIS	2.7
1	A	185	HIS	2.7
1	C	233	GLY	2.6
1	C	249	ALA	2.6
1	C	146	ARG	2.6
1	C	173	TRP	2.5
1	C	131	GLY	2.5
1	D	229	LEU	2.5
1	C	141	ALA	2.4
1	C	158	VAL	2.4
1	C	2	GLU	2.3
1	D	226	SER	2.3
1	C	136	ALA	2.3
1	D	122	PRO	2.3
1	B	185	HIS	2.3
1	A	47	HIS	2.3
1	C	125	ALA	2.2
1	A	57	GLY	2.2
1	D	227	GLN	2.2
1	C	118	ILE	2.2
1	D	195	VAL	2.2
1	B	198	TRP	2.1
1	A	224	SER	2.1
1	C	238	LEU	2.1
1	D	152	GLY	2.0
1	C	130	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

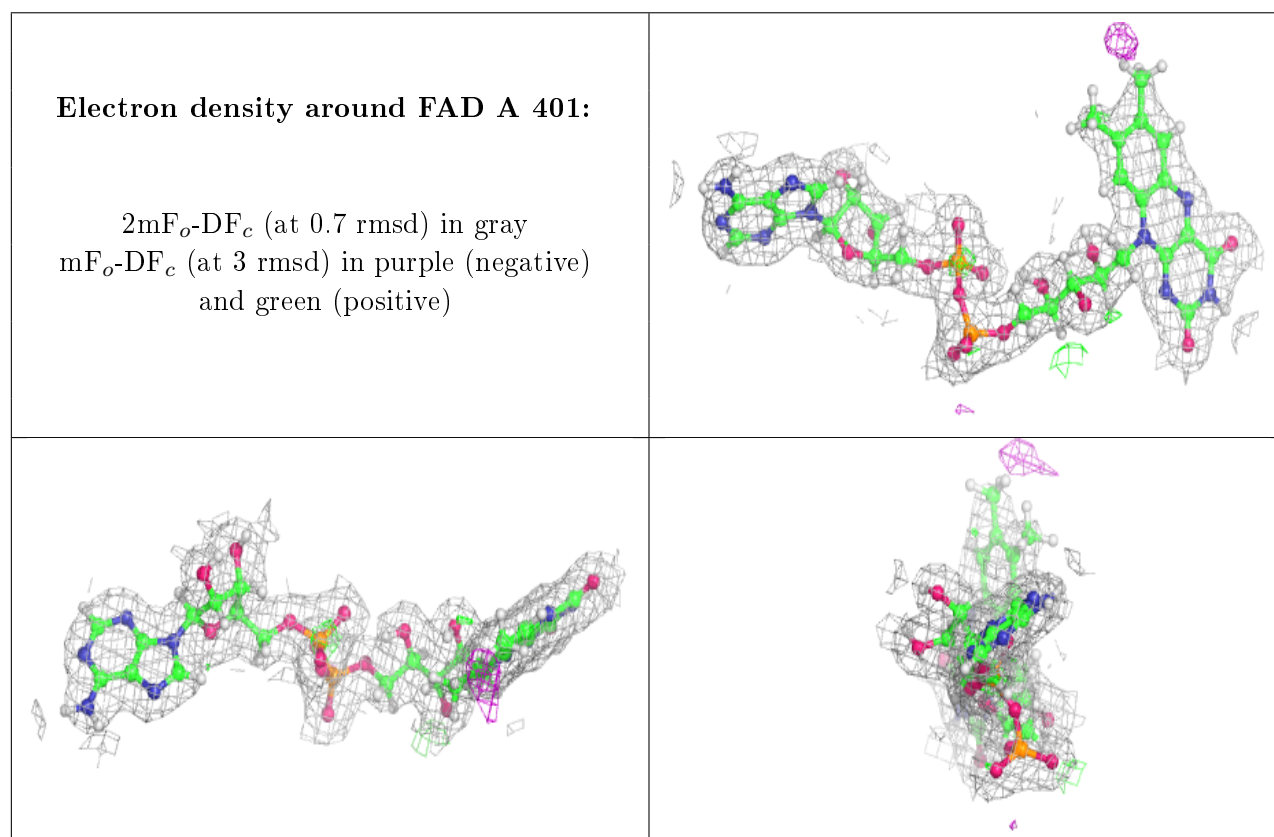
6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

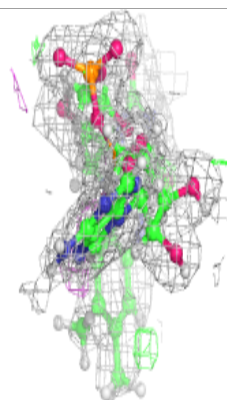
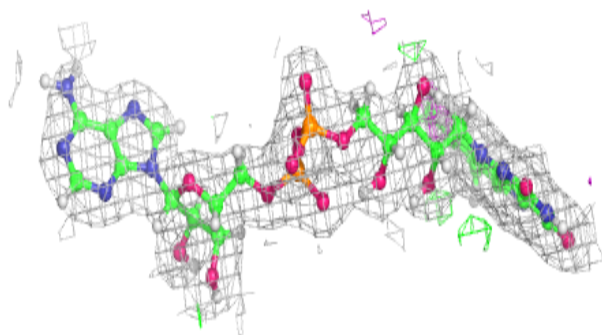
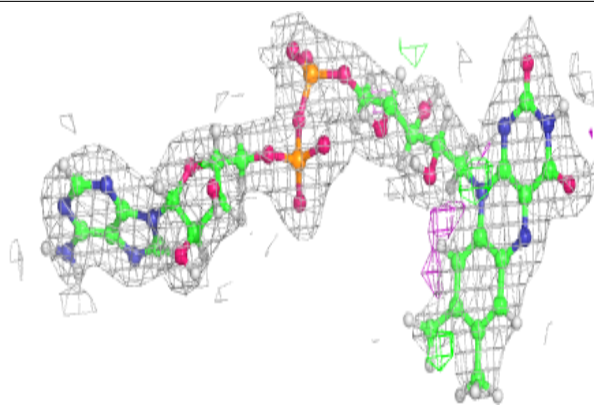
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	PG4	B	501	13/13	0.81	0.16	41,52,62,71	0
4	PG4	A	403	13/13	0.84	0.13	43,54,62,66	0
2	FAD	A	401	53/53	0.98	0.12	18,27,44,50	0
2	FAD	D	401	53/53	0.98	0.12	21,33,47,50	0
2	FAD	C	500	53/53	0.98	0.13	22,34,47,50	0
3	CA	B	502	1/1	0.98	0.06	40,40,40,40	0
2	FAD	B	500	53/53	0.98	0.12	20,30,46,48	0
3	CA	A	402	1/1	0.99	0.06	40,40,40,40	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.

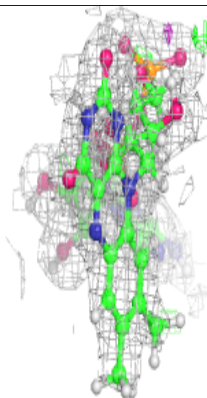
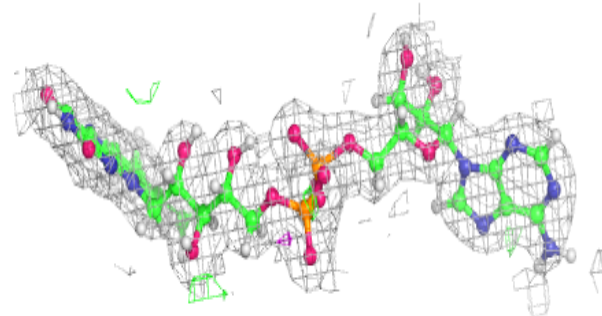
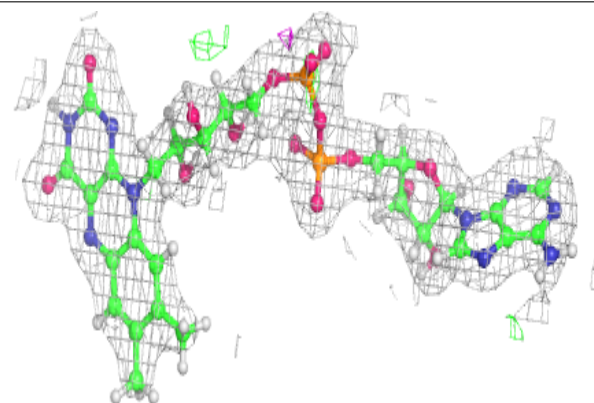


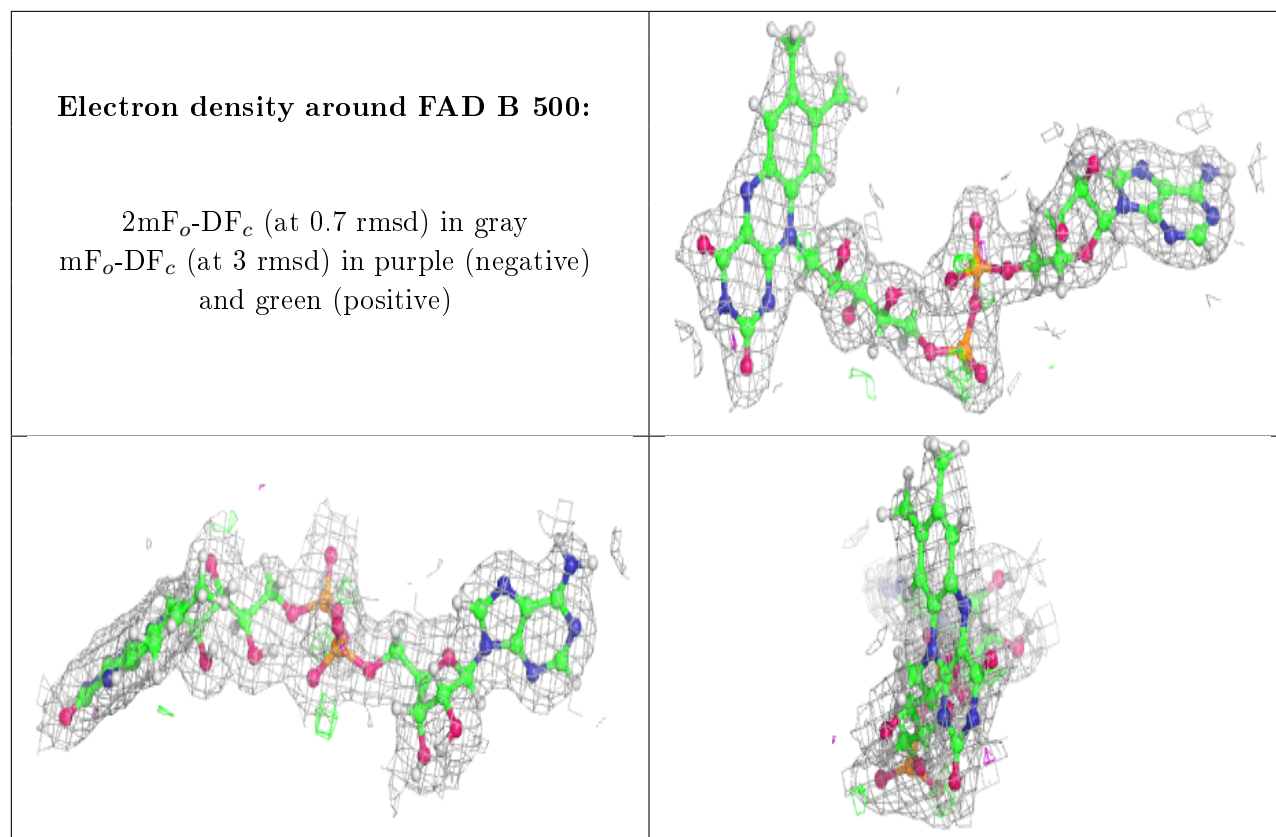
Electron density around FAD D 401:

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

**Electron density around FAD C 500:**

$2mF_o-DF_c$ (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.