



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:19 PM UTC

PDB ID : 7P98 / pdb_00007p98
Title : Cyclohex-1-ene-1-carboxyl-CoA dehydrogenase in a substrate-free state
Authors : Ermler, U.; Weidenweber, S.; Boll, M.
Deposited on : 2021-07-26
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

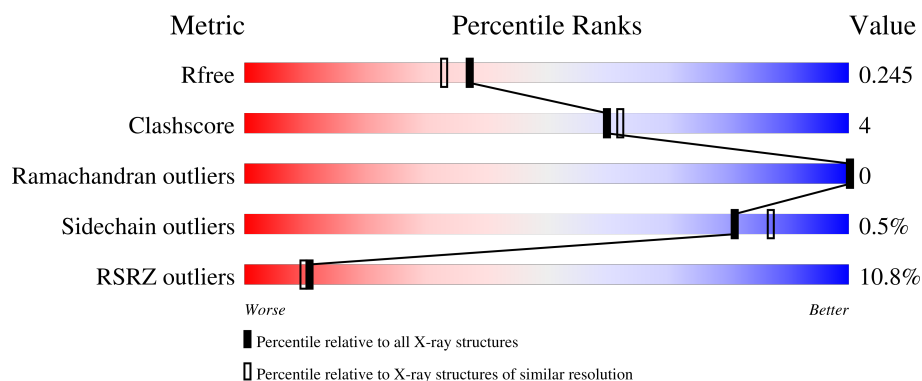
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	412	10% 82% 9% 8%
1	B	412	10% 83% 8% 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6063 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 Short-chain acyl-CoA dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	4	0
			2890	1817	500	551	22			
1	B	378	Total	C	N	O	S	0	2	0
			2868	1803	499	545	21			

There are 64 discrepancies between the modelled and reference sequences:

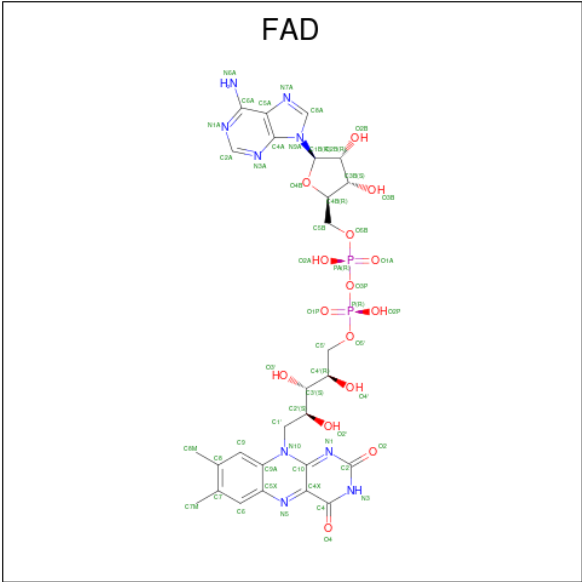
Chain	Residue	Modelled	Actual	Comment	Reference
A	381	LYS	-	expression tag	UNP Q39QF5
A	382	GLY	-	expression tag	UNP Q39QF5
A	383	GLU	-	expression tag	UNP Q39QF5
A	384	LEU	-	expression tag	UNP Q39QF5
A	385	ASN	-	expression tag	UNP Q39QF5
A	386	SER	-	expression tag	UNP Q39QF5
A	387	LYS	-	expression tag	UNP Q39QF5
A	388	LEU	-	expression tag	UNP Q39QF5
A	389	GLU	-	expression tag	UNP Q39QF5
A	390	GLY	-	expression tag	UNP Q39QF5
A	391	LYS	-	expression tag	UNP Q39QF5
A	392	PRO	-	expression tag	UNP Q39QF5
A	393	ILE	-	expression tag	UNP Q39QF5
A	394	PRO	-	expression tag	UNP Q39QF5
A	395	ASN	-	expression tag	UNP Q39QF5
A	396	PRO	-	expression tag	UNP Q39QF5
A	397	LEU	-	expression tag	UNP Q39QF5
A	398	LEU	-	expression tag	UNP Q39QF5
A	399	GLY	-	expression tag	UNP Q39QF5
A	400	LEU	-	expression tag	UNP Q39QF5
A	401	ASP	-	expression tag	UNP Q39QF5
A	402	SER	-	expression tag	UNP Q39QF5
A	403	THR	-	expression tag	UNP Q39QF5
A	404	ARG	-	expression tag	UNP Q39QF5
A	405	THR	-	expression tag	UNP Q39QF5

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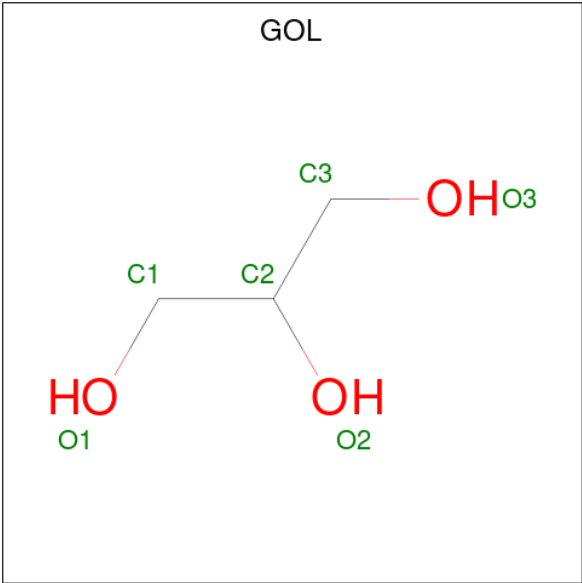
Chain	Residue	Modelled	Actual	Comment	Reference
A	406	GLY	-	expression tag	UNP Q39QF5
A	407	HIS	-	expression tag	UNP Q39QF5
A	408	HIS	-	expression tag	UNP Q39QF5
A	409	HIS	-	expression tag	UNP Q39QF5
A	410	HIS	-	expression tag	UNP Q39QF5
A	411	HIS	-	expression tag	UNP Q39QF5
A	412	HIS	-	expression tag	UNP Q39QF5
B	381	LYS	-	expression tag	UNP Q39QF5
B	382	GLY	-	expression tag	UNP Q39QF5
B	383	GLU	-	expression tag	UNP Q39QF5
B	384	LEU	-	expression tag	UNP Q39QF5
B	385	ASN	-	expression tag	UNP Q39QF5
B	386	SER	-	expression tag	UNP Q39QF5
B	387	LYS	-	expression tag	UNP Q39QF5
B	388	LEU	-	expression tag	UNP Q39QF5
B	389	GLU	-	expression tag	UNP Q39QF5
B	390	GLY	-	expression tag	UNP Q39QF5
B	391	LYS	-	expression tag	UNP Q39QF5
B	392	PRO	-	expression tag	UNP Q39QF5
B	393	ILE	-	expression tag	UNP Q39QF5
B	394	PRO	-	expression tag	UNP Q39QF5
B	395	ASN	-	expression tag	UNP Q39QF5
B	396	PRO	-	expression tag	UNP Q39QF5
B	397	LEU	-	expression tag	UNP Q39QF5
B	398	LEU	-	expression tag	UNP Q39QF5
B	399	GLY	-	expression tag	UNP Q39QF5
B	400	LEU	-	expression tag	UNP Q39QF5
B	401	ASP	-	expression tag	UNP Q39QF5
B	402	SER	-	expression tag	UNP Q39QF5
B	403	THR	-	expression tag	UNP Q39QF5
B	404	ARG	-	expression tag	UNP Q39QF5
B	405	THR	-	expression tag	UNP Q39QF5
B	406	GLY	-	expression tag	UNP Q39QF5
B	407	HIS	-	expression tag	UNP Q39QF5
B	408	HIS	-	expression tag	UNP Q39QF5
B	409	HIS	-	expression tag	UNP Q39QF5
B	410	HIS	-	expression tag	UNP Q39QF5
B	411	HIS	-	expression tag	UNP Q39QF5
B	412	HIS	-	expression tag	UNP Q39QF5

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		

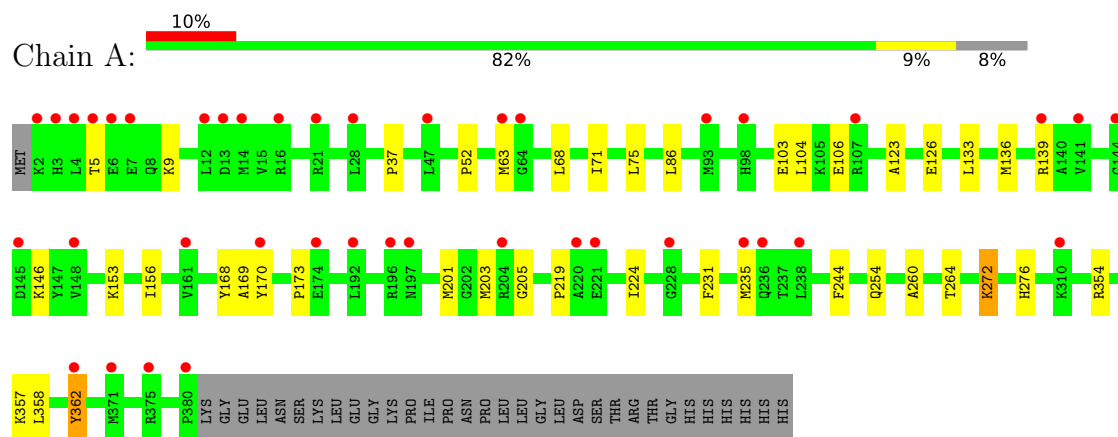
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	106	Total	O	0	0
			106	106		
4	B	74	Total	O	0	1
			75	75		

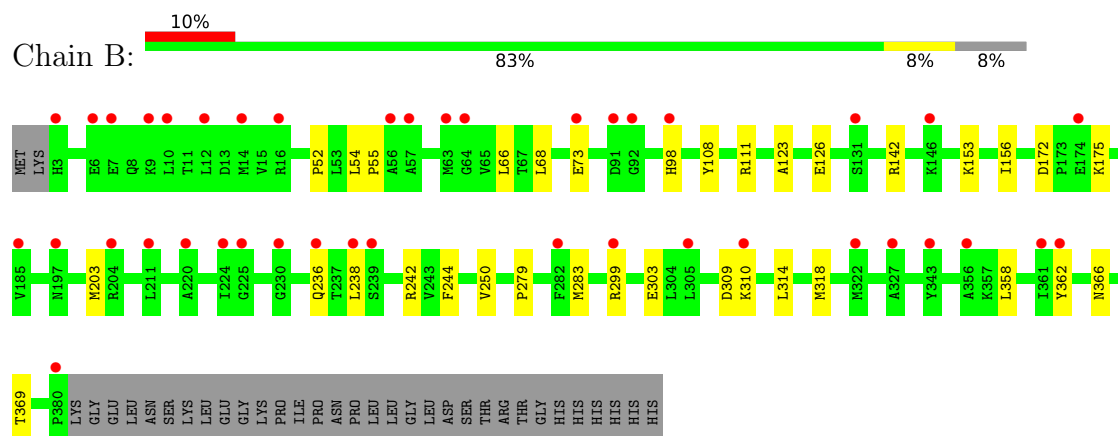
3 Residue-property plots [i](#)

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: Short-chain acyl-CoA dehydrogenase



- Molecule 1: Short-chain acyl-CoA dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	76.32Å 172.20Å 331.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.44 – 2.00 45.44 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.44-2.00) 99.8 (45.44-2.00)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.214 , 0.243 0.218 , 0.245	Depositor DCC
R_{free} test set	3637 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6063	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% 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, GOL

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	0.38	0/2928	0.53	1/3953 (0.0%)
1	B	0.31	0/2906	0.50	0/3924
All	All	0.34	0/5834	0.52	1/7877 (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	A	362	TYR	CB-CA-C	5.60	121.17	109.68

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	2890	0	2954	30	0
1	B	2868	0	2934	21	0
2	A	53	0	29	0	0
2	B	53	0	29	1	0
3	A	12	0	16	3	0
3	B	6	0	8	0	0
4	A	106	0	0	0	0
4	B	75	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6063	0	5970	51	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 (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:A:254:GLN:HE22	3:A:502:GOL:H11	1.49	0.76
1:A:203:MET:HG3	1:A:358:LEU:HD23	1.73	0.70
1:A:201:MET:HE3	1:A:354:ARG:HA	1.72	0.70
1:B:314:LEU:HD11	1:B:318:MET:HE3	1.74	0.69
1:B:203:MET:HG3	1:B:358:LEU:HD23	1.76	0.66
1:A:136:MET:HE1	1:A:168:TYR:O	1.99	0.62
1:A:133:LEU:O	1:A:136:MET:HG2	1.99	0.62
1:A:254:GLN:NE2	3:A:502:GOL:H11	2.18	0.57
1:A:68:LEU:HD23	1:A:244:PHE:CZ	2.42	0.55
1:A:63[B]:MET:HE1	1:A:71:ILE:HD12	1.88	0.54
1:B:126:GLU:HG2	1:B:153:LYS:HD3	1.90	0.52
1:B:238:LEU:O	1:B:242:ARG:HG3	2.11	0.51
1:B:123:ALA:HA	1:B:156:ILE:HD12	1.91	0.51
1:B:172:ASP:HB3	1:B:175:LYS:HD2	1.93	0.50
1:B:68:LEU:HD23	1:B:244:PHE:CZ	2.47	0.50
1:A:139:ARG:NH1	1:A:173:PRO:HG3	2.28	0.48
1:A:52:PRO:HG3	1:A:68:LEU:HD13	1.93	0.48
1:B:98:HIS:HB3	1:B:236:GLN:HE22	1.78	0.48
1:A:123:ALA:HB1	1:A:153:LYS:HG3	1.96	0.48
1:B:66:LEU:HD22	1:B:303:GLU:HG2	1.96	0.48
1:A:123:ALA:HA	1:A:156:ILE:HD12	1.97	0.47
1:A:126:GLU:HG2	1:A:153:LYS:HD3	1.97	0.47
1:A:254:GLN:HE22	3:A:502:GOL:C1	2.22	0.47
1:A:146:LYS:HD3	1:A:219:PRO:HA	1.97	0.46
1:B:299[A]:ARG:O	1:B:303:GLU:HG3	2.15	0.46
1:B:309:ASP:OD1	1:B:310:LYS:N	2.49	0.46
1:B:54:LEU:HD12	1:B:55:PRO:HD2	1.99	0.45
1:B:366:ASN:HA	1:B:369:THR:OG1	2.17	0.45
1:A:103:GLU:H	1:A:103:GLU:CD	2.25	0.44
1:A:75:LEU:CD1	1:A:86:LEU:HD22	2.48	0.44
1:A:203:MET:HG3	1:A:358:LEU:CD2	2.45	0.44
1:B:203:MET:HG3	1:B:358:LEU:CD2	2.45	0.44
1:A:104:LEU:HD21	1:A:224:ILE:HG12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:279:PRO:O	1:B:283:MET:HG3	2.18	0.43
1:A:106:GLU:HG2	1:B:142:ARG:O	2.18	0.43
1:A:260:ALA:O	1:A:264:THR:HG23	2.19	0.43
1:A:231:PHE:CE1	1:A:235:MET:HE3	2.54	0.42
1:A:136:MET:HE1	1:A:169:ALA:HA	2.02	0.42
1:B:52:PRO:HG3	1:B:68:LEU:HD13	2.02	0.42
1:A:136:MET:HE3	1:A:136:MET:HB2	1.77	0.41
1:A:272:LYS:HG3	1:A:276:HIS:HB2	2.01	0.41
1:A:170:TYR:CD1	1:A:173:PRO:HA	2.54	0.41
1:B:299[B]:ARG:O	1:B:303:GLU:HG3	2.20	0.41
1:A:201:MET:HE1	1:A:357:LYS:HB3	2.03	0.41
1:B:73:GLU:HG3	1:B:250:VAL:HG12	2.03	0.41
1:B:108:TYR:O	1:B:111:ARG:HG2	2.21	0.41
1:A:37:PRO:HD3	1:A:205:GLY:O	2.21	0.41
1:A:201:MET:HE1	1:A:357:LYS:CB	2.50	0.41
2:B:501:FAD:H2'	2:B:501:FAD:C9	2.51	0.41
1:B:142:ARG:NH2	4:B:608:HOH:O	2.53	0.40
1:A:5:THR:O	1:A:9:LYS:HG3	2.21	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	381/412 (92%)	379 (100%)	2 (0%)	0	100	100
1	B	378/412 (92%)	377 (100%)	1 (0%)	0	100	100
All	All	759/824 (92%)	756 (100%)	3 (0%)	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	303/328 (92%)	301 (99%)	2 (1%)	76	82
1	B	300/328 (92%)	299 (100%)	1 (0%)	86	91
All	All	603/656 (92%)	600 (100%)	3 (0%)	81	87

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	272	LYS
1	A	362	TYR
1	B	362	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	269	GLN
1	A	276	HIS
1	B	143	GLN
1	B	236	GLN
1	B	367	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

5 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	$\# Z > 2$
3	GOL	A	502	-	5,5,5	0.85	0	5,5,5	1.06	0
2	FAD	A	501	-	58,58,58	1.57	11 (18%)	85,89,89	1.82	22 (25%)
3	GOL	A	503	-	5,5,5	1.42	1 (20%)	5,5,5	0.69	0
3	GOL	B	502	-	5,5,5	1.05	0	5,5,5	0.89	0
2	FAD	B	501	-	58,58,58	1.60	13 (22%)	85,89,89	1.77	22 (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
3	GOL	A	502	-	-	2/4/4/4	-
2	FAD	A	501	-	-	7/34/50/50	0/6/6/6
3	GOL	A	503	-	-	0/4/4/4	-
3	GOL	B	502	-	-	0/4/4/4	-
2	FAD	B	501	-	-	9/34/50/50	0/6/6/6

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	FAD	C5A-C4A	4.87	1.47	1.39
2	A	501	FAD	O2'-C2'	-4.78	1.33	1.43
2	B	501	FAD	O2'-C2'	-4.73	1.33	1.43
2	A	501	FAD	C5A-C4A	4.33	1.46	1.39
2	B	501	FAD	O3'-C3'	-3.96	1.33	1.43
2	A	501	FAD	O3'-C3'	-3.88	1.33	1.43
2	A	501	FAD	C10-N1	3.27	1.39	1.33
2	B	501	FAD	C10-N1	3.24	1.39	1.33
2	B	501	FAD	C4X-N5	2.83	1.36	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	FAD	P-O3P	2.69	1.62	1.59
2	B	501	FAD	C5A-C6A	2.49	1.47	1.41
2	B	501	FAD	C4'-C3'	-2.49	1.49	1.53
2	B	501	FAD	C5A-N7A	-2.47	1.34	1.39
2	A	501	FAD	C5A-C6A	2.44	1.47	1.41
3	A	503	GOL	C1-C2	2.38	1.60	1.51
2	A	501	FAD	C5A-N7A	-2.24	1.35	1.39
2	B	501	FAD	C8A-N7A	2.23	1.36	1.31
2	A	501	FAD	C10-N10	2.13	1.41	1.37
2	B	501	FAD	C4A-N9A	-2.11	1.33	1.37
2	B	501	FAD	PA-O3P	2.08	1.61	1.59
2	B	501	FAD	P-O3P	2.07	1.61	1.59
2	A	501	FAD	C4A-N9A	-2.06	1.33	1.37
2	B	501	FAD	C5'-C4'	2.06	1.54	1.51
2	A	501	FAD	PA-O3P	2.05	1.61	1.59
2	A	501	FAD	C8A-N7A	2.02	1.35	1.31

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	FAD	C5A-C4A-N3A	-5.24	119.50	126.72
2	A	501	FAD	C5A-C4A-N3A	-5.01	119.82	126.72
2	A	501	FAD	N3A-C4A-N9A	4.53	134.86	127.17
2	B	501	FAD	N3A-C4A-N9A	4.41	134.66	127.17
2	B	501	FAD	C10-C4X-N5	-4.40	115.83	124.81
2	A	501	FAD	C9A-N10-C10	-4.10	114.49	120.75
2	A	501	FAD	O3'-C3'-C2'	3.98	117.98	108.93
2	A	501	FAD	C4A-N9A-C8A	3.79	109.72	105.74
2	B	501	FAD	C9A-N10-C10	-3.62	115.23	120.75
2	B	501	FAD	O3'-C3'-C2'	3.58	117.06	108.93
2	B	501	FAD	C2A-N3A-C4A	3.33	119.98	111.83
2	B	501	FAD	C4A-C5A-N7A	-3.30	106.80	110.58
2	A	501	FAD	C2A-N3A-C4A	3.29	119.86	111.83
2	A	501	FAD	N3A-C2A-N1A	-3.23	123.69	128.58
2	B	501	FAD	C4A-N9A-C8A	3.21	109.11	105.74
2	A	501	FAD	C9A-C5X-N5	-3.14	119.12	122.45
2	A	501	FAD	C10-C4X-N5	-3.11	118.47	124.81
2	A	501	FAD	O3'-C3'-C4'	3.10	115.96	108.93
2	B	501	FAD	C5X-N5-C4X	3.01	122.95	118.09
2	A	501	FAD	C5X-N5-C4X	3.00	122.95	118.09
2	A	501	FAD	C4A-C5A-N7A	-2.95	107.21	110.58
2	B	501	FAD	N3A-C2A-N1A	-2.91	124.18	128.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	FAD	O2'-C2'-C3'	2.90	116.04	109.25
2	B	501	FAD	C4-N3-C2	-2.86	120.56	125.64
2	B	501	FAD	C8M-C8-C9	-2.82	114.60	119.57
2	A	501	FAD	O3P-P-O1P	-2.74	102.47	110.70
2	B	501	FAD	C5A-N7A-C8A	2.69	107.67	103.45
2	A	501	FAD	C4-N3-C2	-2.68	120.89	125.64
2	A	501	FAD	N9A-C8A-N7A	-2.65	110.17	113.94
2	A	501	FAD	C5A-N7A-C8A	2.63	107.58	103.45
2	B	501	FAD	C4X-C10-N10	2.57	120.16	116.48
2	B	501	FAD	C9A-C5X-N5	-2.54	119.76	122.45
2	A	501	FAD	C4X-C4-N3	2.51	119.65	113.25
2	B	501	FAD	O4-C4-C4X	-2.47	120.02	126.53
2	B	501	FAD	C4X-C4-N3	2.45	119.49	113.25
2	B	501	FAD	N9A-C8A-N7A	-2.37	110.58	113.94
2	A	501	FAD	O4'-C4'-C5'	2.36	115.20	109.99
2	B	501	FAD	O3P-P-O1P	-2.34	103.67	110.70
2	B	501	FAD	O4'-C4'-C5'	2.23	114.90	109.99
2	A	501	FAD	C8M-C8-C9	-2.23	115.65	119.57
2	B	501	FAD	O2'-C2'-C3'	2.18	114.35	109.25
2	A	501	FAD	O4-C4-C4X	-2.16	120.82	126.53
2	B	501	FAD	C6A-C5A-N7A	2.07	136.08	132.09
2	A	501	FAD	C2A-N1A-C6A	2.05	122.10	118.73

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	FAD	C5B-O5B-PA-O1A
2	A	501	FAD	C2'-C1'-N10-C10
2	A	501	FAD	C1'-C2'-C3'-C4'
2	A	501	FAD	O2'-C2'-C3'-C4'
2	B	501	FAD	C5B-O5B-PA-O1A
2	B	501	FAD	C2'-C1'-N10-C10
2	B	501	FAD	C1'-C2'-C3'-O3'
2	B	501	FAD	C1'-C2'-C3'-C4'
2	B	501	FAD	O2'-C2'-C3'-O3'
2	B	501	FAD	O2'-C2'-C3'-C4'
3	A	502	GOL	O1-C1-C2-C3
2	A	501	FAD	O2'-C2'-C3'-O3'
3	A	502	GOL	O1-C1-C2-O2
2	B	501	FAD	C3'-C4'-C5'-O5'
2	B	501	FAD	C5B-O5B-PA-O3P

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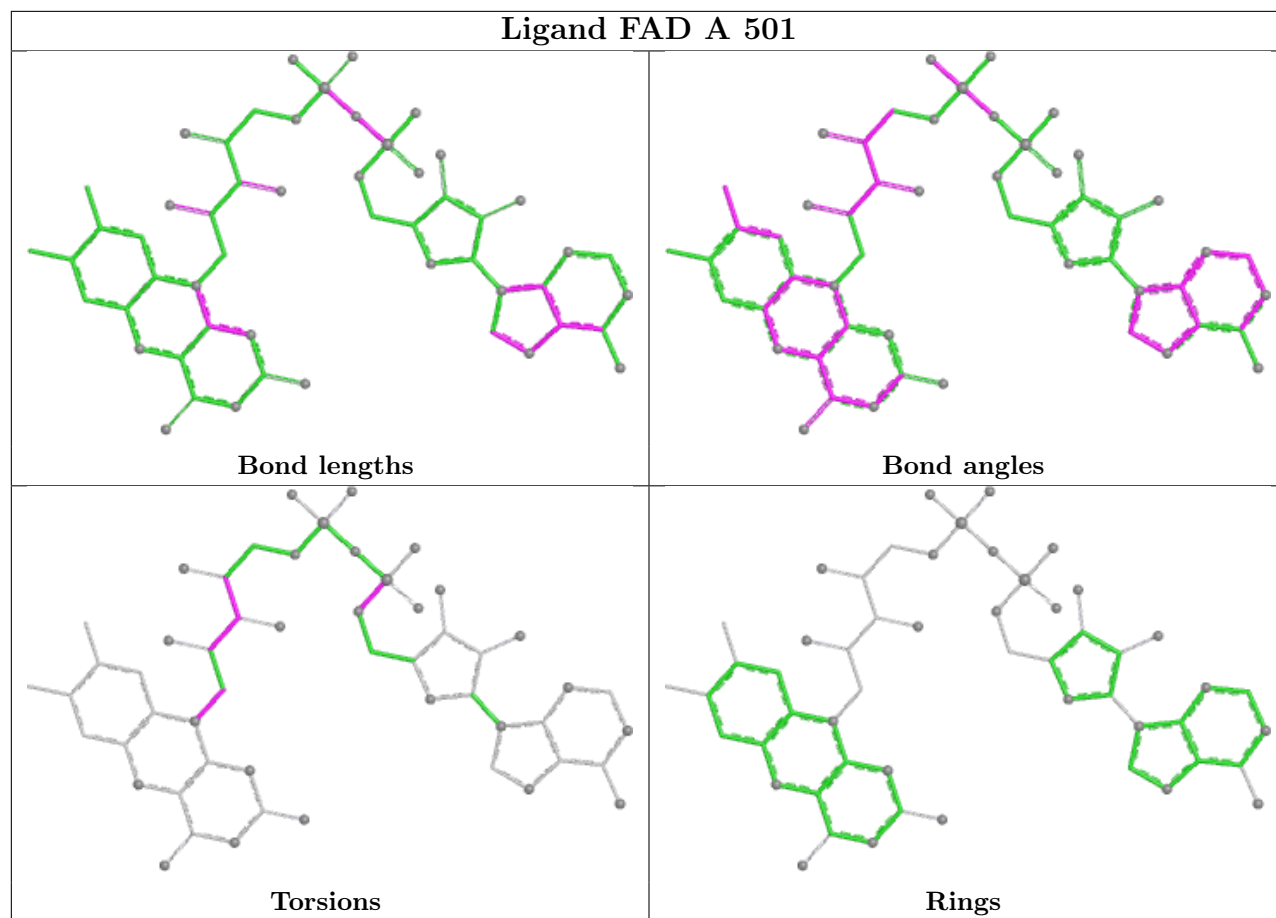
Mol	Chain	Res	Type	Atoms
2	B	501	FAD	C5'-O5'-P-O1P
2	A	501	FAD	O3'-C3'-C4'-C5'
2	A	501	FAD	C1'-C2'-C3'-O3'

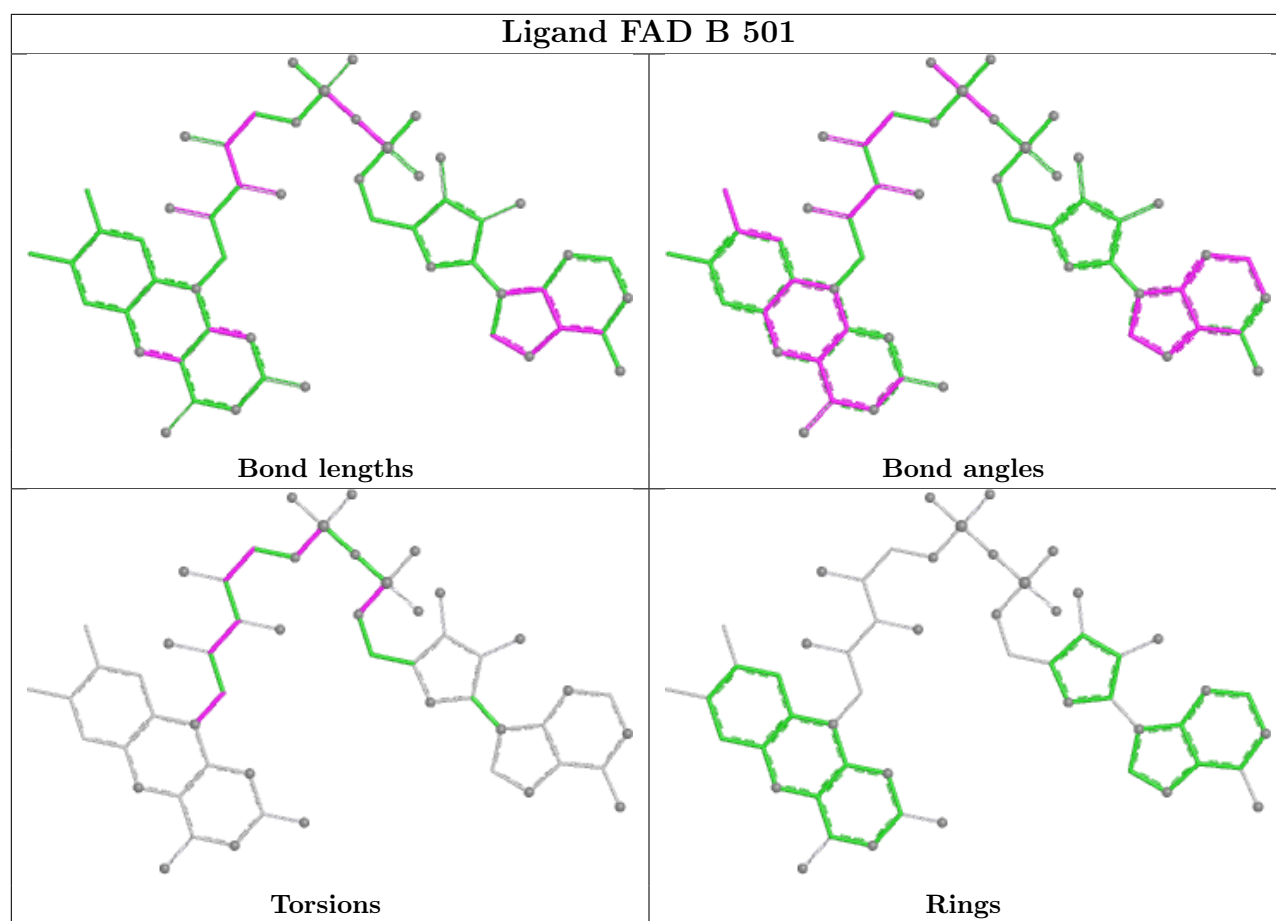
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	GOL	3	0
2	B	501	FAD	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	379/412 (91%)	0.72	41 (10%)	11 10	12, 34, 54, 84	4 (1%)
1	B	378/412 (91%)	0.98	41 (10%)	11 10	16, 39, 58, 87	2 (0%)
All	All	757/824 (91%)	0.85	82 (10%)	11 10	12, 36, 57, 87	6 (0%)

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	63[A]	MET	5.8
1	B	3	HIS	5.3
1	A	2	LYS	4.9
1	A	64	GLY	4.4
1	A	3	HIS	4.3
1	B	10	LEU	4.2
1	B	14	MET	3.9
1	B	239	SER	3.7
1	A	380	PRO	3.6
1	B	356	ALA	3.4
1	B	299[A]	ARG	3.4
1	A	4	LEU	3.3
1	B	185	VAL	3.3
1	A	98	HIS	3.2
1	A	238	LEU	3.2
1	A	16	ARG	3.1
1	B	225	GLY	3.0
1	B	322[A]	MET	3.0
1	B	91	ASP	3.0
1	B	362	TYR	2.9
1	B	73	GLU	2.9
1	A	145	ASP	2.9
1	B	327	ALA	2.9
1	A	174	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	197	ASN	2.7
1	B	9	LYS	2.7
1	A	220	ALA	2.6
1	A	221	GLU	2.6
1	B	98	HIS	2.6
1	B	224	ILE	2.6
1	B	7	GLU	2.6
1	B	56	ALA	2.6
1	A	7	GLU	2.6
1	B	57	ALA	2.6
1	B	238	LEU	2.6
1	B	174	GLU	2.5
1	B	131	SER	2.5
1	B	361	ILE	2.5
1	A	13	ASP	2.5
1	A	12	LEU	2.4
1	A	235	MET	2.4
1	A	21	ARG	2.4
1	A	228	GLY	2.4
1	B	236	GLN	2.4
1	B	220	ALA	2.4
1	B	6	GLU	2.4
1	A	5	THR	2.4
1	B	380	PRO	2.4
1	B	12	LEU	2.3
1	A	371	MET	2.3
1	A	196	ARG	2.3
1	A	6	GLU	2.3
1	A	170	TYR	2.3
1	A	139	ARG	2.3
1	B	305	LEU	2.3
1	A	144	GLY	2.2
1	B	282	PHE	2.2
1	A	141	VAL	2.2
1	B	64	GLY	2.2
1	A	204	ARG	2.2
1	A	310	LYS	2.2
1	A	197	ASN	2.2
1	A	107	ARG	2.1
1	B	16	ARG	2.1
1	B	211	LEU	2.1
1	B	146	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	93	MET	2.1
1	A	148	VAL	2.1
1	A	192	LEU	2.1
1	B	63	MET	2.1
1	A	236	GLN	2.1
1	B	310	LYS	2.1
1	A	47	LEU	2.1
1	A	14	MET	2.1
1	A	362	TYR	2.1
1	A	375	ARG	2.1
1	B	92	GLY	2.1
1	B	230	GLY	2.1
1	A	161	VAL	2.0
1	A	28	LEU	2.0
1	B	204	ARG	2.0
1	B	343	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

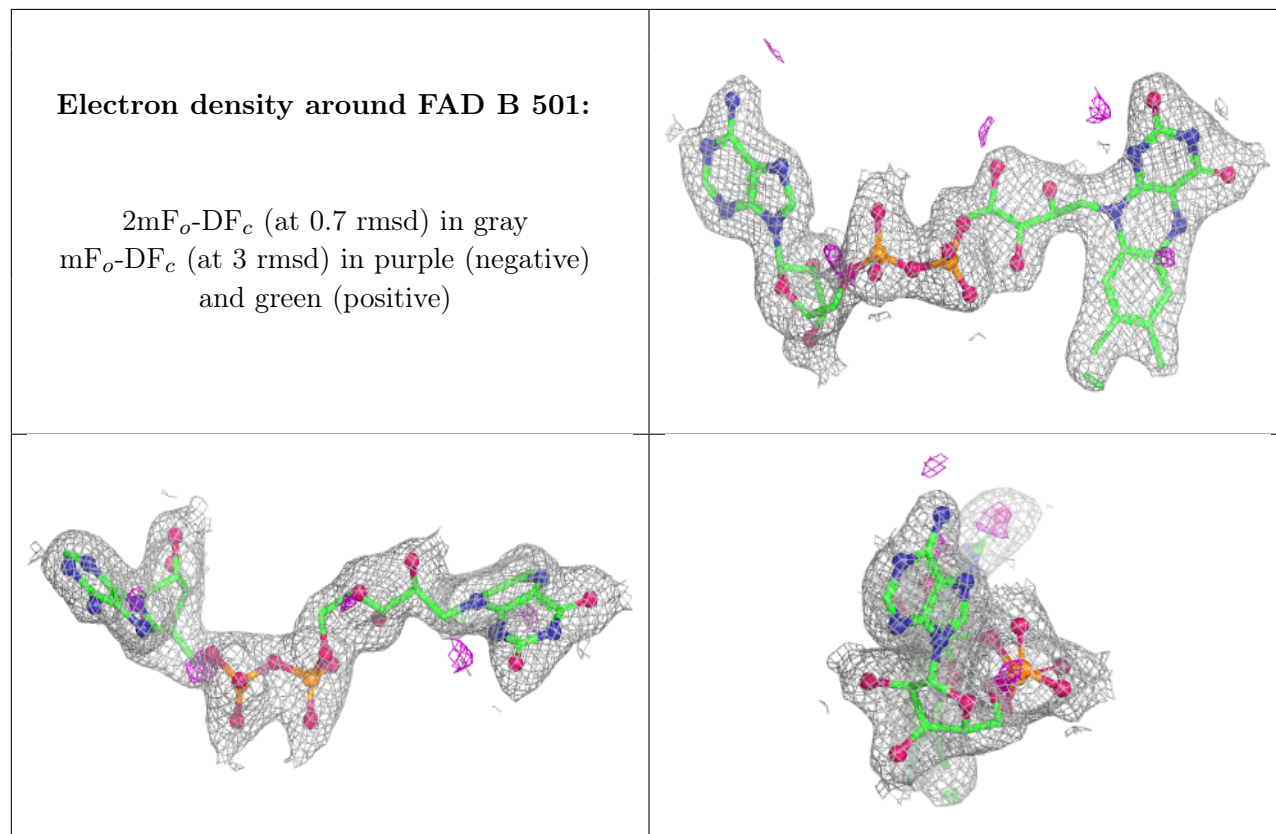
6.4 Ligands [i](#)

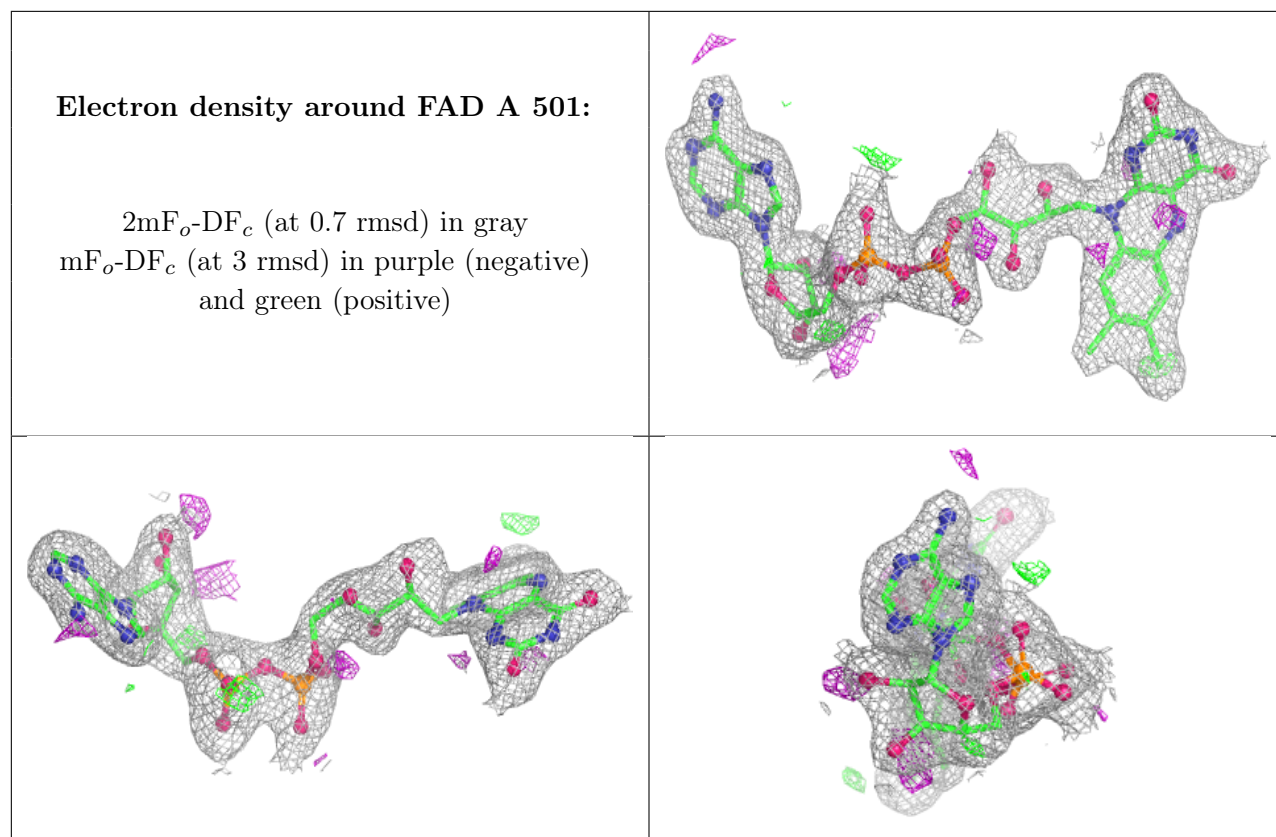
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	B	502	6/6	0.87	0.13	36,39,41,44	0
3	GOL	A	502	6/6	0.88	0.18	49,50,53,55	0
3	GOL	A	503	6/6	0.91	0.11	35,36,37,43	0
2	FAD	B	501	53/53	0.93	0.10	31,37,42,43	0
2	FAD	A	501	53/53	0.94	0.09	29,34,38,41	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.