



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:25 AM UTC

PDB ID : 8PNS / pdb_00008pns
Title : Crystal structure of the acyl-CoA dehydrogenase PA0506 (FadE1) from *Pseudomonas aeruginosa*
Authors : Wang, M.; Brear, P.; Welch, M.
Deposited on : 2023-06-30
Resolution : 2.08 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

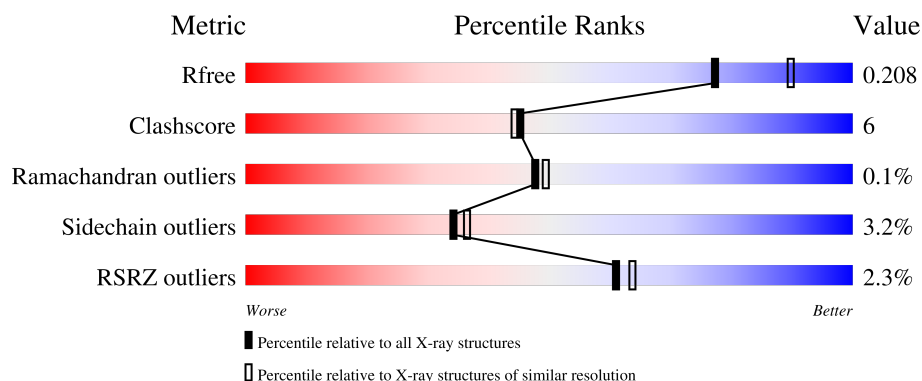
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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	603	2% 86% 12% .
1	B	603	3% 82% 16% ..

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 9545 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.

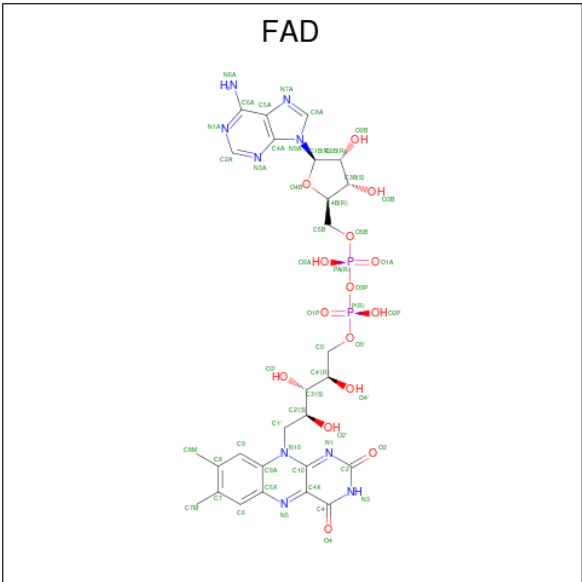
- Molecule 1 is a protein called Acyl-CoA dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	600	Total	C	N	O	S	0	2	0
			4587	2897	794	858	38			
1	B	599	Total	C	N	O	S	0	1	0
			4578	2889	794	858	37			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP A0A072ZR57
A	0	HIS	-	expression tag	UNP A0A072ZR57
B	-1	GLY	-	expression tag	UNP A0A072ZR57
B	0	HIS	-	expression tag	UNP A0A072ZR57

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



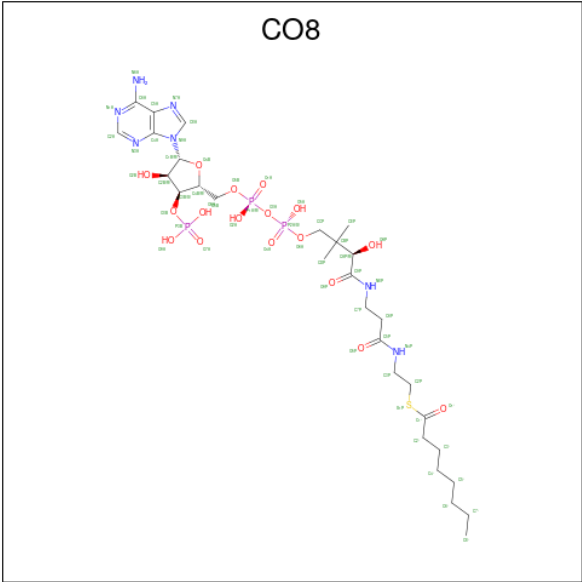
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is OCTANOYL-COENZYME A (CCD ID: CO8) (formula: $C_{29}H_{50}N_7O_{17}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	0
			57	29	7	17	3	1		
4	B	1	Total	C	N	O	P	S	0	0
			57	29	7	17	3	1		

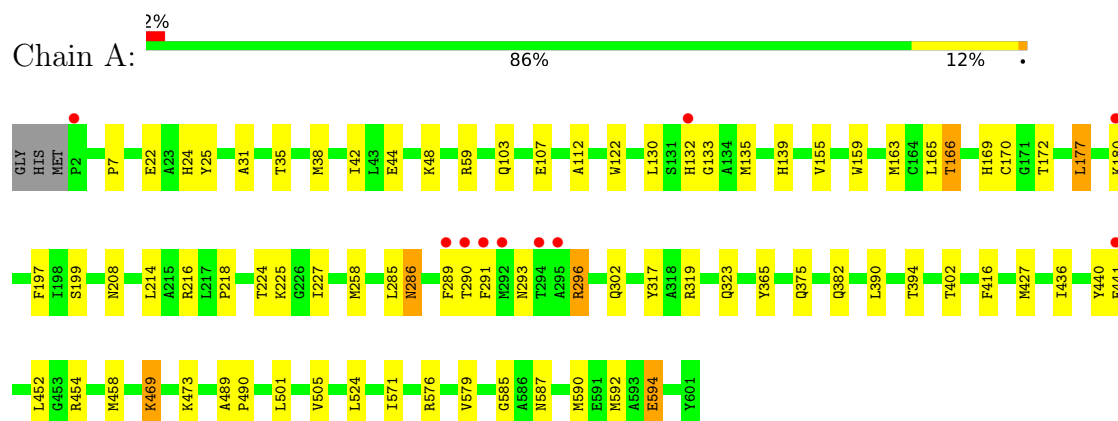
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	78	Total	O	0	0
			78	78		
5	B	68	Total	O	0	0
			68	68		

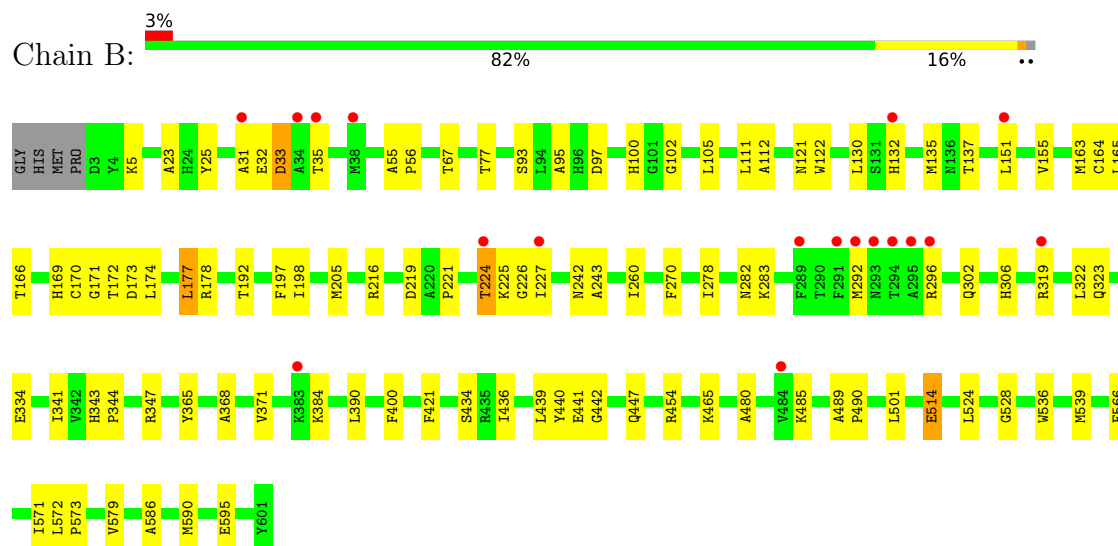
3 Residue-property plots [i](#)

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: Acyl-CoA dehydrogenase



- Molecule 1: Acyl-CoA dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	155.92Å 155.92Å 99.51Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	135.03 – 2.08 135.03 – 2.08	Depositor EDS
% Data completeness (in resolution range)	99.9 (135.03-2.08) 99.9 (135.03-2.08)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.204 , 0.249 (Not available) , 0.208	Depositor DCC
R_{free} test set	4062 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	36.3	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 29.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.026 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9545	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% 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: PEG, FAD, CO8

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	1.10	6/4693 (0.1%)	1.47	19/6340 (0.3%)
1	B	1.09	0/4677	1.51	10/6319 (0.2%)
All	All	1.10	6/9370 (0.1%)	1.49	29/12659 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	427[A]	MET	C-O	6.41	1.32	1.24
1	A	427[B]	MET	C-O	6.41	1.32	1.24
1	A	59	ARG	C-O	5.96	1.31	1.24
1	A	319[A]	ARG	C-O	5.72	1.31	1.24
1	A	319[B]	ARG	C-O	5.72	1.31	1.24
1	A	139	HIS	CE1-NE2	5.40	1.38	1.32

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	319[A]	ARG	CA-C-O	7.51	128.62	119.31
1	A	319[B]	ARG	CA-C-O	7.51	128.62	119.31
1	A	427[A]	MET	CA-C-O	7.17	128.21	119.97
1	A	427[B]	MET	CA-C-O	7.17	128.21	119.97
1	A	166	THR	N-CA-C	5.91	118.91	110.10
1	B	334	GLU	CB-CG-CD	5.84	122.53	112.60
1	A	218	PRO	N-CA-C	5.74	124.30	112.47
1	B	166	THR	N-CA-C	5.72	118.63	110.10
1	A	585	GLY	CA-C-O	-5.70	118.30	122.23
1	B	283	LYS	CA-C-N	5.54	126.09	119.94
1	B	283	LYS	C-N-CA	5.54	126.09	119.94
1	A	7	PRO	CA-C-O	-5.54	117.21	121.31
1	A	290	THR	CA-C-N	5.47	127.55	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	290	THR	C-N-CA	5.47	127.55	120.44
1	B	67	THR	CA-CB-OG1	-5.45	101.42	109.60
1	A	402	THR	CA-CB-OG1	5.43	117.75	109.60
1	A	22	GLU	CA-C-N	5.43	127.55	120.28
1	A	22	GLU	C-N-CA	5.43	127.55	120.28
1	A	382	GLN	CA-C-N	5.30	127.39	120.28
1	A	382	GLN	C-N-CA	5.30	127.39	120.28
1	B	23	ALA	CA-C-N	5.25	127.62	120.54
1	B	23	ALA	C-N-CA	5.25	127.62	120.54
1	A	103	GLN	N-CA-C	-5.24	106.93	113.38
1	B	77	THR	CA-CB-OG1	-5.23	101.76	109.60
1	B	32	GLU	N-CA-C	-5.13	107.67	114.04
1	A	394	THR	CA-CB-OG1	-5.11	101.93	109.60
1	B	292	MET	CA-C-O	-5.10	115.10	120.55
1	A	576	ARG	CA-C-N	5.00	126.94	120.44
1	A	576	ARG	C-N-CA	5.00	126.94	120.44

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	4587	0	4499	52	0
1	B	4578	0	4481	63	0
2	A	53	0	31	8	0
2	B	53	0	31	6	0
3	A	14	0	20	3	0
4	A	57	0	46	5	0
4	B	57	0	46	6	0
5	A	78	0	0	5	0
5	B	68	0	0	2	0
All	All	9545	0	9154	110	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 6.

All (110) 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:436:ILE:HD11	2:A:701:FAD:HM82	1.52	0.92
2:B:701:FAD:C4X	4:B:702:CO8:H3'2	2.24	0.68
1:A:224:THR:HA	1:A:227:ILE:HD12	1.77	0.66
1:A:286:ASN:HD22	1:A:286:ASN:H	1.45	0.65
1:A:285:LEU:HD11	4:A:703:CO8:H142	1.77	0.65
1:A:436:ILE:HD11	2:A:701:FAD:C8M	2.27	0.63
1:B:100:HIS:HE1	5:B:802:HOH:O	1.81	0.62
1:A:323:GLN:HG3	2:B:701:FAD:O2A	2.00	0.62
1:A:285:LEU:HD21	4:A:703:CO8:H121	1.81	0.61
1:B:442:GLY:HA2	4:B:702:CO8:H62	1.82	0.61
1:A:501:LEU:HD11	1:A:579:VAL:HG22	1.81	0.61
1:A:286:ASN:H	1:A:286:ASN:ND2	1.97	0.61
1:B:172:THR:HA	4:B:702:CO8:H71	1.83	0.60
1:B:501:LEU:HD11	1:B:579:VAL:HG22	1.82	0.59
1:B:566:PHE:HE2	1:B:571:ILE:HD11	1.66	0.59
1:A:436:ILE:CD1	2:A:701:FAD:HM82	2.30	0.59
1:A:587:ASN:HB3	5:A:851:HOH:O	2.02	0.59
1:A:38:MET:HE3	1:A:42:ILE:HD11	1.85	0.59
1:B:489:ALA:HB3	1:B:490:PRO:CD	2.33	0.58
1:B:165:LEU:HD12	1:B:165:LEU:N	2.19	0.58
1:A:323:GLN:HE22	1:B:170:CYS:HA	1.70	0.57
1:B:296:ARG:HD3	1:B:441:GLU:HG3	1.84	0.57
1:B:341:ILE:O	1:B:347:ARG:HD2	2.04	0.56
1:B:112:ALA:HA	1:B:365:TYR:OH	2.06	0.56
1:B:524:LEU:C	1:B:524:LEU:HD23	2.31	0.56
1:A:107:GLU:HG3	1:A:375:GLN:OE1	2.06	0.56
1:B:260:ILE:HD11	2:B:701:FAD:HM73	1.87	0.56
2:A:701:FAD:O1A	1:B:323:GLN:HG3	2.06	0.56
1:B:178:ARG:HD3	5:B:866:HOH:O	2.06	0.55
1:B:151:LEU:HD23	1:B:155:VAL:HG23	1.88	0.55
1:B:172:THR:HA	4:B:702:CO8:C7P	2.38	0.54
1:A:132:HIS:HA	1:A:135:MET:HE3	1.90	0.53
1:B:489:ALA:HB3	1:B:490:PRO:HD3	1.90	0.53
1:A:163:MET:HB3	1:A:291:PHE:CE2	2.44	0.53
1:B:566:PHE:CE2	1:B:571:ILE:HD11	2.44	0.53
1:A:122:TRP:CH2	1:A:302:GLN:HB2	2.44	0.52
1:B:197:PHE:HB3	2:B:701:FAD:C9A	2.39	0.52
1:B:33:ASP:N	1:B:33:ASP:OD1	2.43	0.51
1:B:25:TYR:O	1:B:31:ALA:HB3	2.10	0.51
1:B:130:LEU:HD23	1:B:163:MET:HE3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:PHE:O	2:B:701:FAD:C4X	2.58	0.51
1:A:165:LEU:HD12	1:A:165:LEU:N	2.26	0.50
1:A:285:LEU:HD11	4:A:703:CO8:CEP	2.41	0.49
1:B:242:ASN:O	1:B:243:ALA:C	2.55	0.49
1:B:132:HIS:HA	1:B:135:MET:HE3	1.93	0.49
1:A:293:ASN:HB3	3:A:702:PEG:H12	1.95	0.49
1:A:317:TYR:CE2	1:A:416:PHE:HA	2.48	0.48
1:A:590:MET:HA	1:B:347:ARG:HE	1.77	0.48
1:A:169:HIS:O	1:B:323:GLN:NE2	2.47	0.48
1:B:536:TRP:CE3	1:B:539:MET:HE3	2.48	0.48
1:A:130:LEU:HD11	1:A:199:SER:HB2	1.96	0.48
1:A:390:LEU:HD13	3:A:702:PEG:H22	1.96	0.48
1:A:489:ALA:HB3	1:A:490:PRO:CD	2.44	0.48
1:A:172:THR:HA	4:A:703:CO8:H72	1.95	0.47
1:B:151:LEU:HD23	1:B:155:VAL:CG2	2.43	0.47
1:A:197:PHE:O	2:A:701:FAD:C4X	2.62	0.47
1:B:121:ASN:C	1:B:121:ASN:HD22	2.21	0.47
1:A:44:GLU:HG2	5:A:833:HOH:O	2.14	0.46
1:B:164:CYS:HA	1:B:198:ILE:HD12	1.96	0.46
1:B:586:ALA:O	1:B:590:MET:HG2	2.16	0.46
1:B:400:PHE:CD1	1:B:528:GLY:HA3	2.51	0.46
1:B:436:ILE:HD12	1:B:439:LEU:HB2	1.98	0.46
1:A:170:CYS:HA	1:B:323:GLN:HE22	1.81	0.46
1:B:221:PRO:O	1:B:226:GLY:HA3	2.16	0.46
1:A:197:PHE:HB3	2:A:701:FAD:C9A	2.46	0.46
1:A:289:PHE:O	1:A:293:ASN:ND2	2.49	0.45
1:A:133:GLY:HA3	1:A:291:PHE:CE1	2.52	0.45
1:B:480:ALA:HA	1:B:485:LYS:HE3	1.97	0.45
1:B:322:LEU:O	1:B:323:GLN:HG2	2.17	0.45
1:B:55:ALA:HB3	1:B:56:PRO:HD3	1.99	0.45
1:B:572:LEU:N	1:B:573:PRO:CD	2.80	0.45
1:A:177:LEU:O	1:A:216:ARG:HD2	2.17	0.44
1:A:592:MET:O	1:B:347:ARG:NH2	2.49	0.44
1:B:122:TRP:CH2	1:B:302:GLN:HB2	2.54	0.43
1:A:25:TYR:O	1:A:31:ALA:HB3	2.19	0.43
1:B:97:ASP:OD2	1:B:100:HIS:HD2	2.00	0.43
1:B:95:ALA:HA	1:B:105:LEU:O	2.19	0.43
1:A:258:MET:HB3	1:B:421:PHE:CE1	2.54	0.43
1:A:524:LEU:C	1:A:524:LEU:HD23	2.44	0.43
1:A:452:LEU:HD21	1:A:505:VAL:HG11	2.01	0.43
1:B:170:CYS:HB2	1:B:173:ASP:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:296:ARG:HD2	1:B:447:GLN:HE22	1.84	0.43
1:B:371:VAL:HG23	1:B:390:LEU:HD21	2.01	0.43
1:A:159:TRP:CG	1:A:208:ASN:HB3	2.54	0.42
1:A:296:ARG:HD3	5:A:875:HOH:O	2.18	0.42
1:A:469:LYS:HE3	1:A:469:LYS:HA	2.00	0.42
1:A:112:ALA:HA	1:A:365:TYR:OH	2.19	0.42
1:A:489:ALA:HB3	1:A:490:PRO:HD3	2.01	0.42
1:A:293:ASN:HB3	3:A:702:PEG:C1	2.50	0.42
1:A:214:LEU:HD22	1:A:227:ILE:CG2	2.49	0.42
1:B:177:LEU:O	1:B:216:ARG:HD2	2.20	0.42
1:B:224:THR:HA	1:B:227:ILE:HD12	2.00	0.42
1:A:594:GLU:HG3	1:B:319[B]:ARG:CZ	2.50	0.42
1:B:111:LEU:HD22	1:B:368:ALA:HB2	2.02	0.42
1:B:165:LEU:HD22	4:B:702:CO8:H31	2.02	0.42
1:B:343:HIS:HA	1:B:344:PRO:HD3	1.95	0.42
1:B:225:LYS:O	1:B:282:ASN:HB3	2.20	0.42
1:A:35:THR:HG22	5:A:872:HOH:O	2.20	0.41
1:B:93:SER:HB2	1:B:102:GLY:HA2	2.02	0.41
1:B:306:HIS:CD2	1:B:434:SER:HA	2.55	0.41
1:A:24:HIS:HD2	5:A:878:HOH:O	2.03	0.41
1:B:390:LEU:HD23	1:B:390:LEU:HA	1.92	0.41
1:B:514:GLU:HG2	1:B:590:MET:HG3	2.01	0.41
1:A:289:PHE:CE2	4:A:703:CO8:H143	2.56	0.41
2:A:701:FAD:C9	2:A:701:FAD:H2'	2.51	0.41
1:A:323:GLN:NE2	1:B:169:HIS:O	2.54	0.41
1:B:174:LEU:HD12	4:B:702:CO8:H131	2.02	0.41
1:A:166:THR:OG1	2:A:701:FAD:H1'1	2.21	0.40
1:B:192:THR:HA	1:B:270:PHE:O	2.20	0.40
2:B:701:FAD:HM83	2:B:701:FAD:HM71	1.91	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	600/603 (100%)	582 (97%)	18 (3%)	0	100	100
1	B	598/603 (99%)	577 (96%)	20 (3%)	1 (0%)	43	44
All	All	1198/1206 (99%)	1159 (97%)	38 (3%)	1 (0%)	48	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	171	GLY

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	465/465 (100%)	450 (97%)	15 (3%)	34	36
1	B	463/465 (100%)	448 (97%)	15 (3%)	34	36
All	All	928/930 (100%)	898 (97%)	30 (3%)	34	36

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	LYS
1	A	155	VAL
1	A	177	LEU
1	A	180	LYS
1	A	225	LYS
1	A	286	ASN
1	A	296	ARG
1	A	440	TYR
1	A	441	GLU
1	A	454	ARG
1	A	458	MET
1	A	469	LYS
1	A	473	LYS
1	A	571	ILE
1	A	594	GLU

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Mol	Chain	Res	Type
1	B	5	LYS
1	B	33	ASP
1	B	35	THR
1	B	137	THR
1	B	177	LEU
1	B	205	MET
1	B	219	ASP
1	B	224	THR
1	B	278	ILE
1	B	384	LYS
1	B	440	TYR
1	B	454	ARG
1	B	465	LYS
1	B	514	GLU
1	B	595	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	HIS
1	A	286	ASN
1	A	302	GLN
1	A	476	GLN
1	B	100	HIS
1	B	103	GLN
1	B	121	ASN
1	B	136	ASN
1	B	184	GLN
1	B	282	ASN
1	B	323	GLN
1	B	419	HIS
1	B	447	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

6 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	$\# Z > 2$
4	CO8	B	702	-	57,59,59	0.44	0	79,85,85	0.72	1 (1%)
3	PEG	A	704	-	6,6,6	0.52	0	5,5,5	0.31	0
4	CO8	A	703	-	57,59,59	0.44	0	79,85,85	0.74	1 (1%)
2	FAD	B	701	-	58,58,58	0.61	0	85,89,89	0.59	1 (1%)
3	PEG	A	702	-	6,6,6	0.47	0	5,5,5	0.28	0
2	FAD	A	701	-	58,58,58	0.66	1 (1%)	85,89,89	0.60	1 (1%)

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	CO8	B	702	-	-	13/58/74/74	0/3/3/3
3	PEG	A	704	-	-	2/4/4/4	-
4	CO8	A	703	-	-	23/58/74/74	0/3/3/3
2	FAD	B	701	-	-	2/34/50/50	0/6/6/6
3	PEG	A	702	-	-	2/4/4/4	-
2	FAD	A	701	-	-	4/34/50/50	0/6/6/6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	FAD	C9-C8	-2.03	1.36	1.39

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	702	CO8	P3B-O3B-C3B	-4.39	111.70	123.43
4	A	703	CO8	P3B-O3B-C3B	-4.22	112.16	123.43
2	A	701	FAD	C4-N3-C2	-2.07	121.96	125.64
2	B	701	FAD	C4-N3-C2	-2.03	122.03	125.64

There are no chirality outliers.

All (46) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	FAD	C1'-C2'-C3'-O3'
2	A	701	FAD	C1'-C2'-C3'-C4'
4	A	703	CO8	CCP-O6A-P2A-O4A
4	A	703	CO8	CDP-CBP-CCP-O6A
4	A	703	CO8	CEP-CBP-CCP-O6A
4	A	703	CO8	CAP-CBP-CCP-O6A
4	A	703	CO8	C9P-CAP-CBP-CCP
4	A	703	CO8	C9P-CAP-CBP-CDP
4	A	703	CO8	S1P-C2P-C3P-N4P
4	A	703	CO8	O1'-C1'-S1P-C2P
4	A	703	CO8	C2'-C1'-S1P-C2P
4	A	703	CO8	C1'-C2'-C3'-C4'
4	B	702	CO8	CAP-C9P-N8P-C7P
4	B	702	CO8	S1P-C2P-C3P-N4P
4	B	702	CO8	C3P-C2P-S1P-C1'
4	B	702	CO8	O1'-C1'-S1P-C2P
4	B	702	CO8	C2'-C1'-S1P-C2P
4	B	702	CO8	O9P-C9P-N8P-C7P
4	A	703	CO8	O5P-C5P-N4P-C3P
4	A	703	CO8	C6P-C5P-N4P-C3P
2	A	701	FAD	O2'-C2'-C3'-O3'
2	A	701	FAD	O2'-C2'-C3'-C4'
3	A	702	PEG	O1-C1-C2-O2
3	A	702	PEG	O2-C3-C4-O4
3	A	704	PEG	O1-C1-C2-O2
4	A	703	CO8	C2'-C3'-C4'-C5'
4	B	702	CO8	C3'-C4'-C5'-C6'
4	A	703	CO8	C5P-C6P-C7P-N8P
4	A	703	CO8	OAP-CAP-CBP-CDP
4	B	702	CO8	C5'-C6'-C7'-C8'
4	A	703	CO8	C4'-C5'-C6'-C7'
4	B	702	CO8	C6P-C7P-N8P-C9P

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Mol	Chain	Res	Type	Atoms
4	B	702	CO8	C4'-C5'-C6'-C7'
3	A	704	PEG	C1-C2-O2-C3
4	A	703	CO8	OAP-CAP-CBP-CCP
4	A	703	CO8	OAP-CAP-CBP-CEP
4	B	702	CO8	C2'-C3'-C4'-C5'
2	B	701	FAD	O2'-C2'-C3'-C4'
4	A	703	CO8	P2A-O3A-P1A-O1A
4	B	702	CO8	P2A-O3A-P1A-O1A
4	A	703	CO8	C9P-CAP-CBP-CEP
4	A	703	CO8	C3'-C4'-C5'-C6'
4	A	703	CO8	P2A-O3A-P1A-O2A
2	B	701	FAD	O2'-C2'-C3'-O3'
4	A	703	CO8	N8P-C9P-CAP-OAP
4	B	702	CO8	P2A-O3A-P1A-O2A

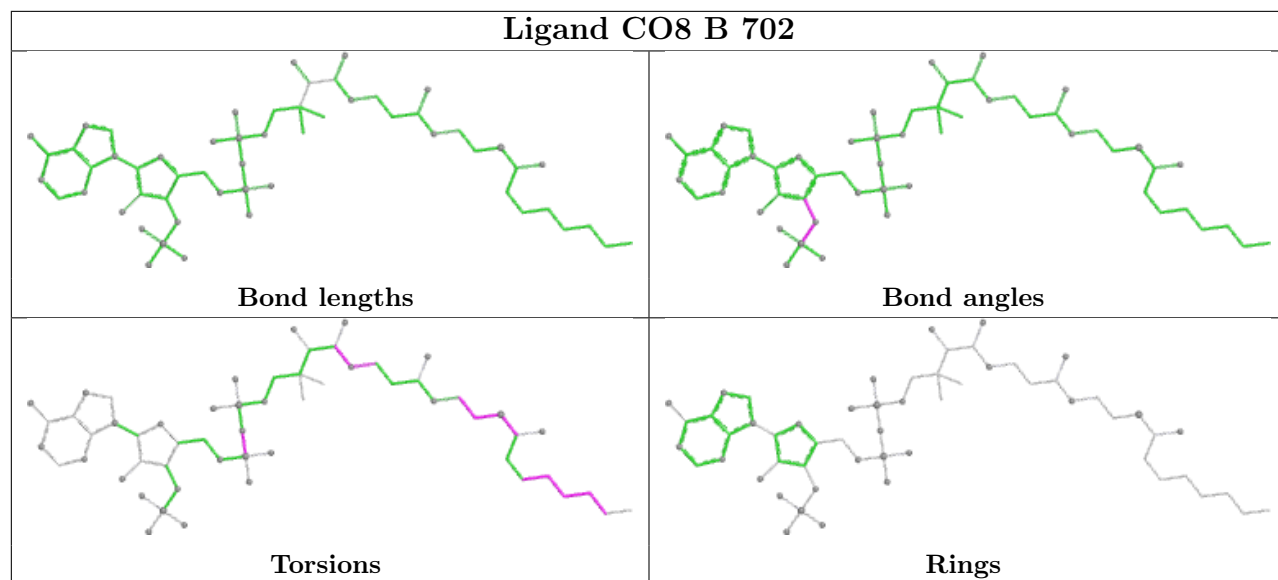
There are no ring outliers.

5 monomers are involved in 27 short contacts:

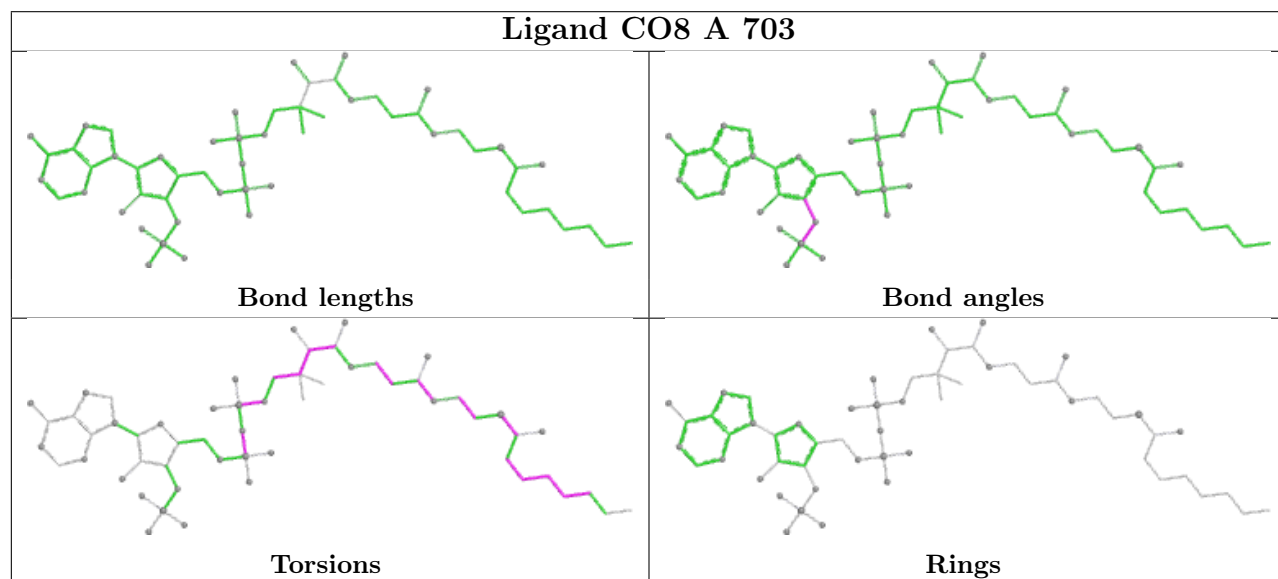
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	702	CO8	6	0
4	A	703	CO8	5	0
2	B	701	FAD	6	0
3	A	702	PEG	3	0
2	A	701	FAD	8	0

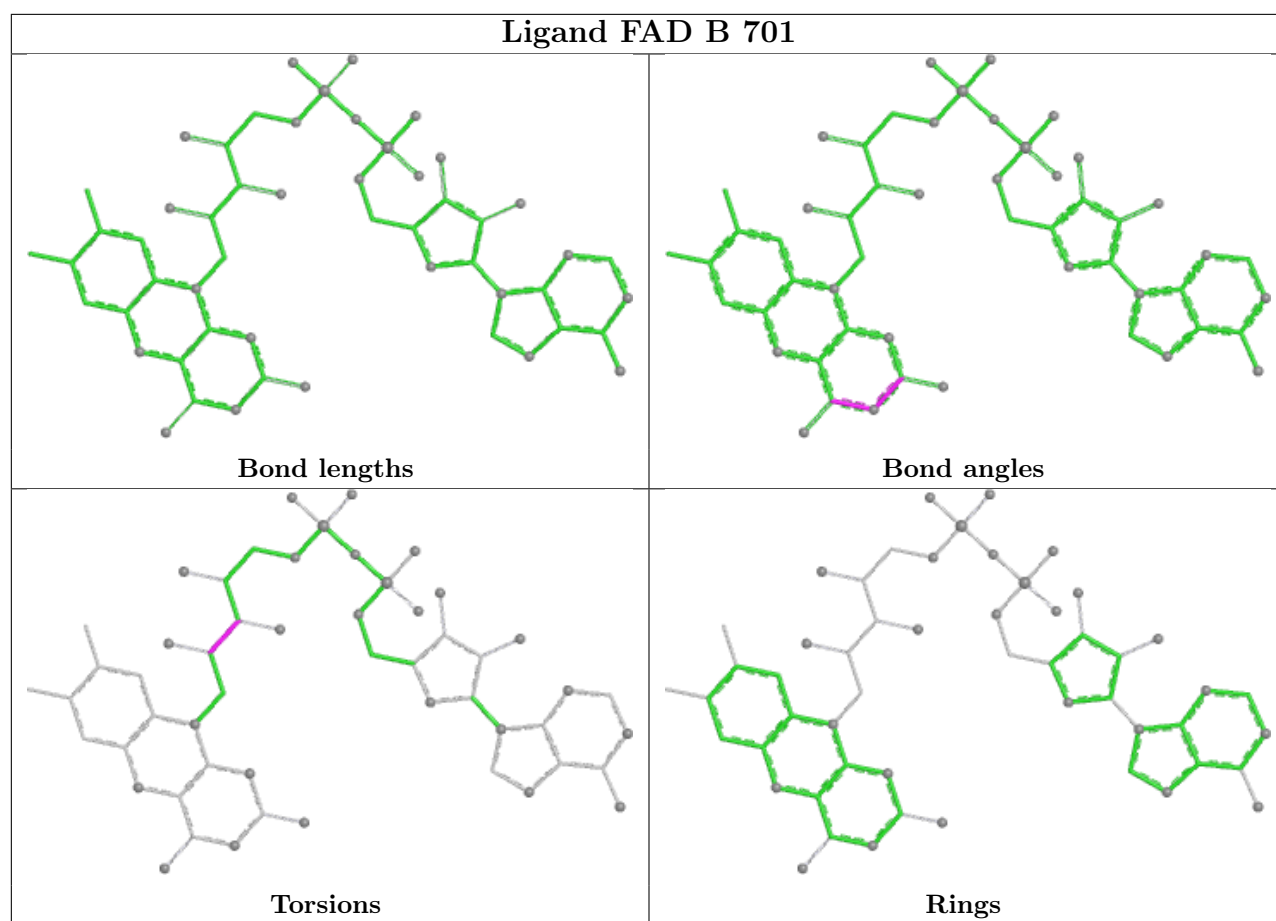
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

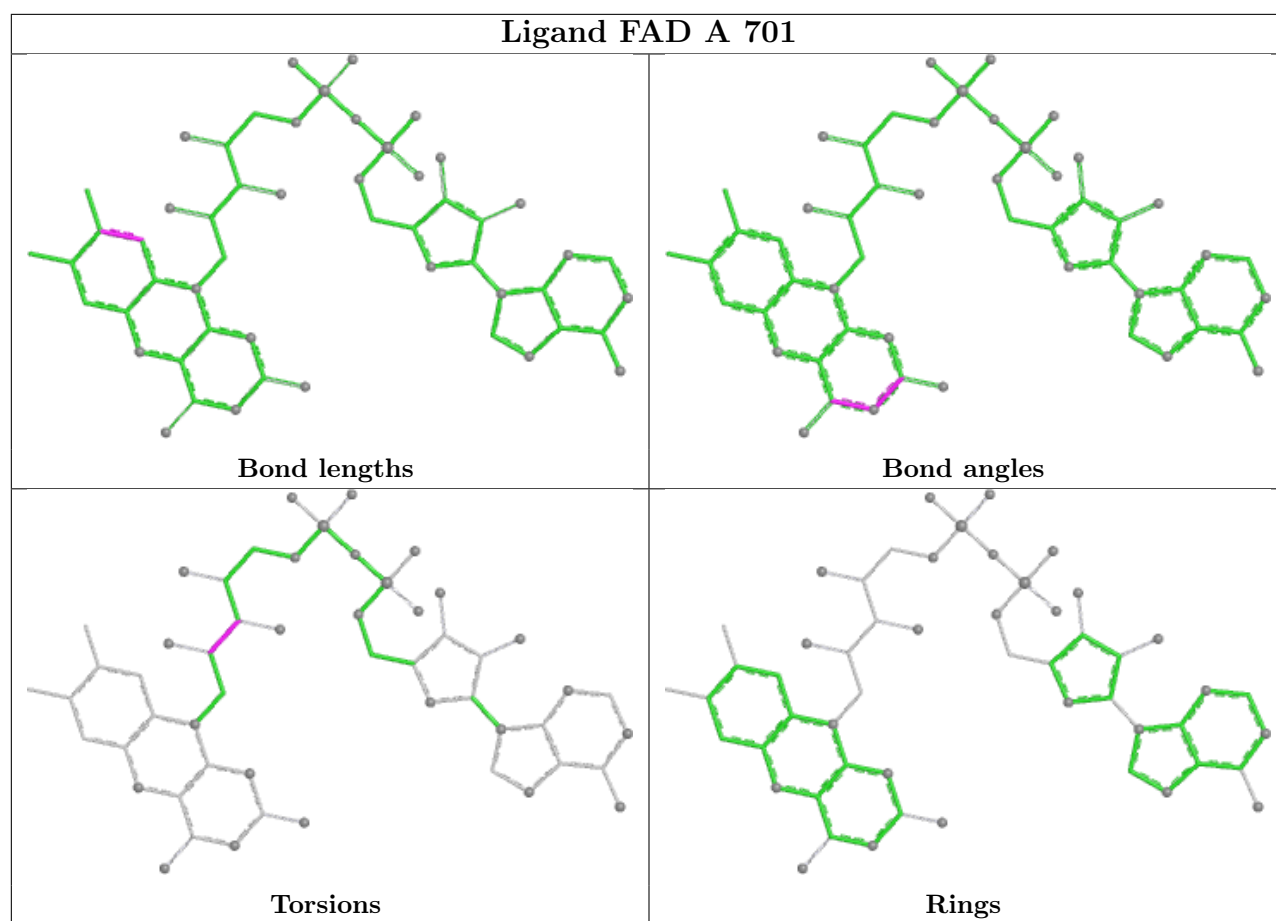
Ligand CO8 B 702



Ligand CO8 A 703







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	600/603 (99%)	0.16	10 (1%) 69 71	18, 39, 61, 78	2 (0%)
1	B	599/603 (99%)	0.30	18 (3%) 52 54	16, 42, 66, 88	1 (0%)
All	All	1199/1206 (99%)	0.23	28 (2%) 61 64	16, 41, 64, 88	3 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	PRO	5.1
1	A	291	PHE	4.6
1	B	291	PHE	4.4
1	A	295	ALA	3.8
1	B	289	PHE	3.7
1	B	132	HIS	3.3
1	A	294	THR	3.2
1	A	180	LYS	3.1
1	B	294	THR	3.1
1	B	295	ALA	3.0
1	B	292	MET	3.0
1	B	31	ALA	2.9
1	B	34	ALA	2.9
1	B	293	ASN	2.9
1	B	296	ARG	2.7
1	A	289	PHE	2.7
1	A	292	MET	2.6
1	B	224	THR	2.5
1	A	441	GLU	2.4
1	B	484	VAL	2.3
1	B	38	MET	2.2
1	B	319[A]	ARG	2.2
1	B	35	THR	2.1
1	B	227	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	151	LEU	2.1
1	A	132	HIS	2.0
1	B	383	LYS	2.0
1	A	290	THR	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](#)

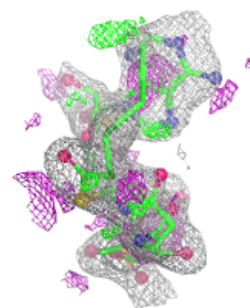
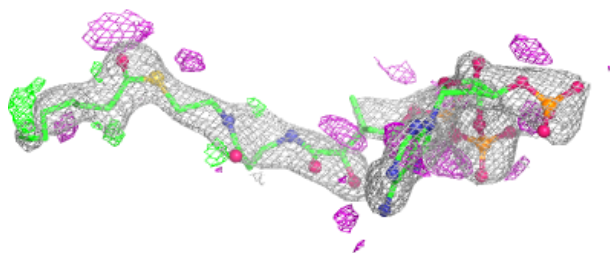
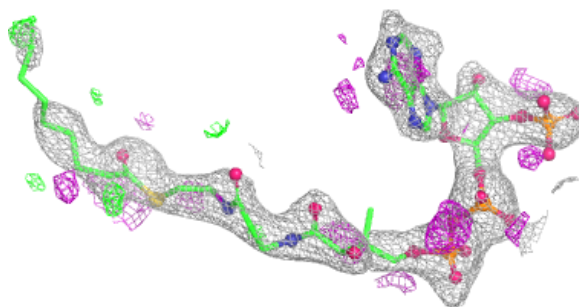
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	PEG	A	702	7/7	0.79	0.20	61,61,69,69	0
4	CO8	B	702	57/57	0.79	0.17	45,87,108,114	0
4	CO8	A	703	57/57	0.82	0.16	48,77,104,126	0
3	PEG	A	704	7/7	0.89	0.15	39,50,55,56	0
2	FAD	A	701	53/53	0.93	0.09	25,31,37,40	0
2	FAD	B	701	53/53	0.94	0.08	27,34,40,41	0

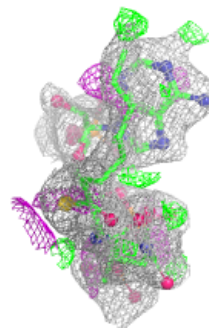
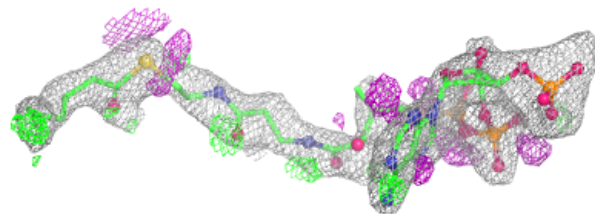
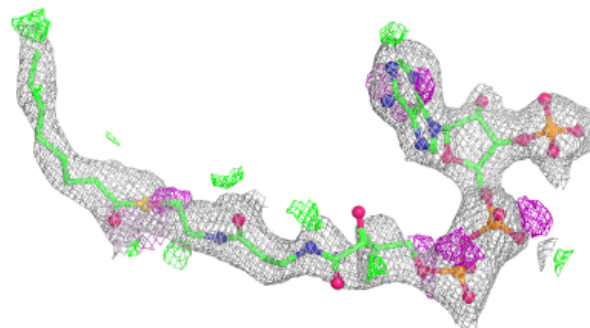
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CO8 B 702:

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

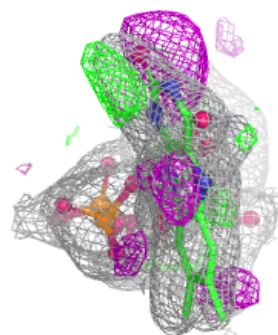
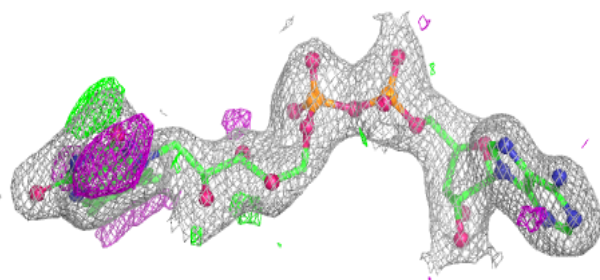
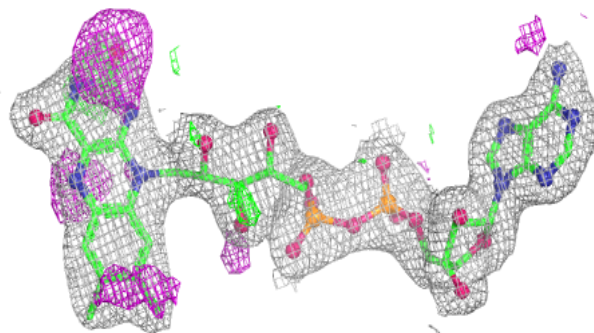
**Electron density around CO8 A 703:**

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

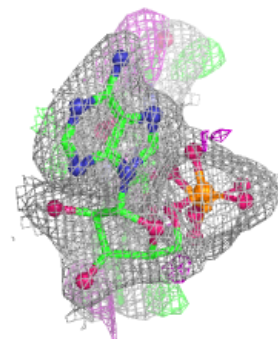
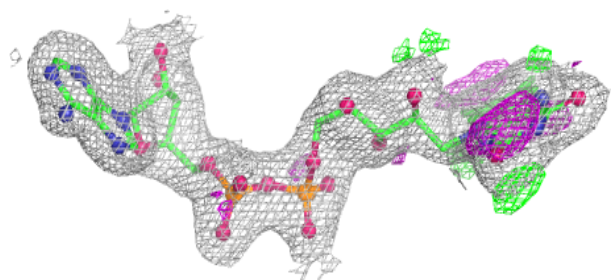
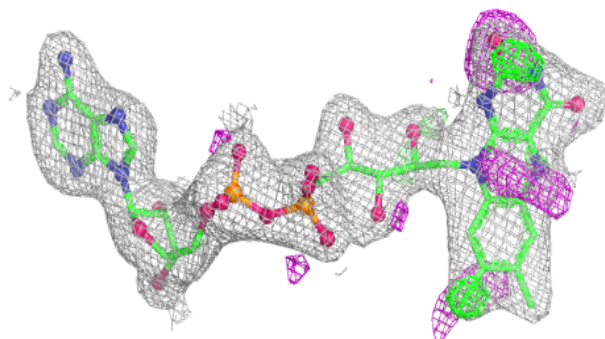


Electron density around FAD A 701:

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

**Electron density around FAD B 701:**

$2mF_o-DF_c$ (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.