



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 08:42 PM UTC

PDB ID : 9FXQ / pdb_00009fxq
Title : Ancestral Prenylcysteine Oxidase 1 (PCYOX1)
Authors : Barone, M.; Mattevi, A.
Deposited on : 2024-07-02
Resolution : 3.10 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

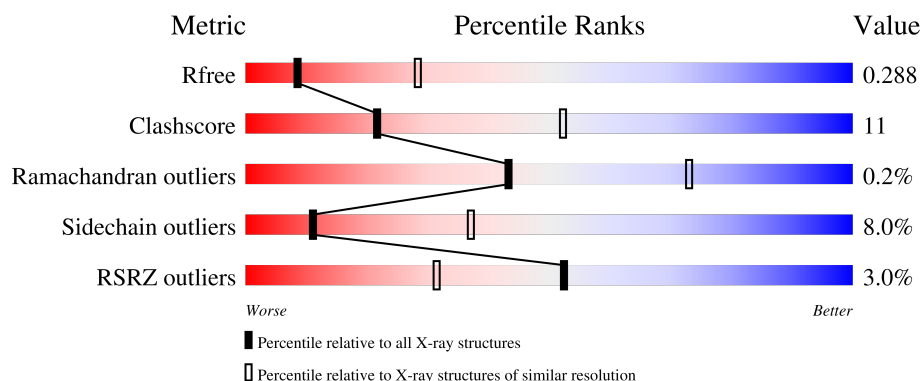
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

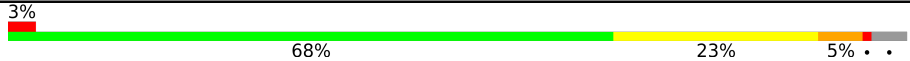
The reported resolution of this entry is 3.10 Å.

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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	482	

2 Entry composition [i](#)

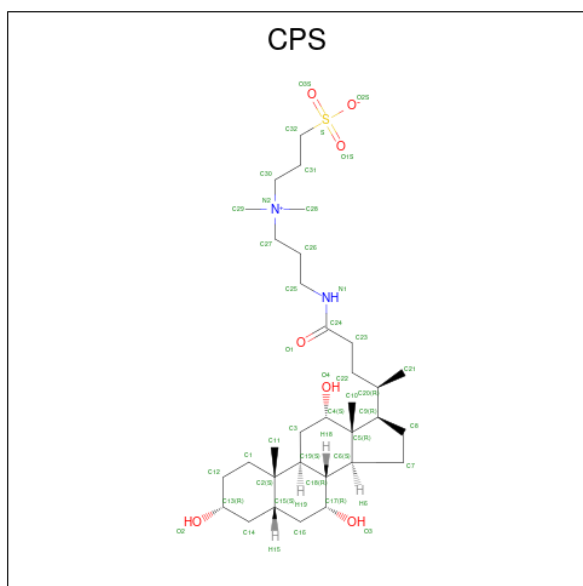
There are 3 unique types of molecules in this entry. The entry contains 3729 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ancestral Prenylcysteine Oxidase 1 (PCYOX1).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3654	2354	613	673	14			

- Molecule 2 is 3-[(3-CHOLAMIDOPROPYL)DIMETHYLAMMONIO]-1-PROPANESULFONATE (CCD ID: CPS) (formula: $C_{32}H_{58}N_2O_7S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			22	19	3		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



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

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	172.66Å 172.66Å 172.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	122.39 – 3.10 122.09 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (122.39-3.10) 99.8 (122.09-3.10)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.07Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, R_{free}	0.225 , 0.275 0.244 , 0.288	Depositor DCC
R_{free} test set	1648 reflections (9.94%)	wwPDB-VP
Wilson B-factor (Å ²)	124.1	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 143.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3729	wwPDB-VP
Average B, all atoms (Å ²)	153.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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: CPS, 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	0.66	0/3744	1.26	27/5069 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	6

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	333	HIS	CB-CA-C	9.68	124.36	109.84
1	A	333	HIS	N-CA-C	-9.00	97.01	110.24
1	A	330	PRO	CB-CA-C	-7.43	100.00	111.22
1	A	431	LEU	N-CA-C	-7.01	95.88	110.80
1	A	240	ASP	CA-CB-CG	6.57	119.17	112.60
1	A	316	LYS	CB-CG-CD	6.38	125.98	111.30
1	A	294	GLN	CB-CA-C	6.26	119.99	109.48
1	A	145	PHE	CA-CB-CG	-6.14	107.66	113.80
1	A	495	TYR	CB-CA-C	-6.03	100.42	110.68
1	A	495	TYR	N-CA-C	5.95	118.53	111.33
1	A	417	LEU	N-CA-C	-5.92	105.32	112.54
1	A	292	THR	CA-CB-OG1	-5.79	100.92	109.60
1	A	287	LYS	N-CA-C	5.77	118.20	107.99
1	A	305	ASP	CA-CB-CG	5.73	118.33	112.60
1	A	287	LYS	N-CA-CB	-5.55	101.26	110.47
1	A	294	GLN	N-CA-CB	-5.54	101.37	110.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	492	GLU	CB-CG-CD	5.53	122.00	112.60
1	A	387	GLU	CB-CA-C	5.49	118.92	110.14
1	A	334	LYS	N-CA-CB	-5.48	103.23	111.56
1	A	295	THR	CA-CB-OG1	-5.46	101.41	109.60
1	A	165	TYR	N-CA-CB	5.31	118.51	110.28
1	A	332	PHE	N-CA-C	-5.31	106.48	113.17
1	A	334	LYS	CB-CA-C	5.23	119.21	110.43
1	A	250	ASN	N-CA-C	-5.14	105.86	111.82
1	A	327	PRO	N-CA-CB	-5.06	97.03	102.60
1	A	418	LEU	CA-C-N	-5.06	114.86	122.81
1	A	418	LEU	C-N-CA	-5.06	114.86	122.81

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	164	ARG	Sidechain
1	A	189	ARG	Sidechain
1	A	27	ARG	Sidechain
1	A	278	ARG	Sidechain
1	A	438	PRO	Peptide
1	A	70	ARG	Sidechain

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	3654	0	3598	80	0
2	A	22	0	30	0	0
3	A	53	0	31	0	0
All	All	3729	0	3659	80	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 11.

All (80) 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:71:LEU:O	1:A:251:LYS:HB3	1.83	0.79
1:A:34:ALA:HB2	1:A:304:TYR:CD2	2.18	0.78
1:A:76:VAL:HG22	1:A:426:LYS:HG3	1.66	0.77
1:A:334:LYS:HB3	1:A:459:TRP:CZ2	2.23	0.74
1:A:108:GLN:HG2	1:A:244:TRP:HB3	1.73	0.70
1:A:70:ARG:O	1:A:84:GLY:HA3	1.95	0.66
1:A:355:TYR:HB3	1:A:358:PRO:HB3	1.77	0.66
1:A:34:ALA:HB2	1:A:304:TYR:CE2	2.32	0.65
1:A:286:VAL:HA	1:A:287:LYS:HZ3	1.63	0.64
1:A:309:VAL:HG11	1:A:313:LEU:HD21	1.79	0.63
1:A:499:LYS:O	1:A:500:THR:C	2.41	0.63
1:A:332:PHE:O	1:A:333:HIS:C	2.41	0.62
1:A:228:ASN:HB2	1:A:231:VAL:HG23	1.84	0.59
1:A:314:ASN:OD1	1:A:316:LYS:HG2	2.04	0.58
1:A:37:GLY:O	1:A:42:GLY:HA3	2.04	0.57
1:A:196:LEU:HD12	1:A:209:LEU:HD13	1.86	0.57
1:A:337:HIS:NE2	1:A:407:GLN:OE1	2.38	0.56
1:A:217:MET:HE1	1:A:235:SER:HB3	1.86	0.56
1:A:81:TYR:OH	1:A:401:VAL:HG21	2.06	0.56
1:A:174:SER:N	1:A:445:ILE:HD12	2.21	0.56
1:A:453:TYR:CE1	1:A:455:ASN:HB2	2.41	0.55
1:A:436:TYR:HE1	1:A:459:TRP:HA	1.72	0.55
1:A:280:LYS:O	1:A:281:GLN:C	2.50	0.54
1:A:130:PHE:HA	1:A:133:ASN:HD22	1.75	0.52
1:A:178:LEU:C	1:A:178:LEU:HD23	2.34	0.52
1:A:333:HIS:O	1:A:334:LYS:CD	2.58	0.52
1:A:314:ASN:HB3	1:A:317:MET:HG3	1.91	0.52
1:A:227:ILE:HG22	1:A:436:TYR:HB2	1.92	0.51
1:A:164:ARG:HD3	1:A:498:LEU:HB3	1.93	0.51
1:A:151:TRP:O	1:A:155:ILE:HG12	2.12	0.50
1:A:447:LEU:HD11	1:A:453:TYR:HB2	1.94	0.50
1:A:319:ASN:OD1	1:A:319:ASN:N	2.43	0.50
1:A:221:TYR:O	1:A:222:GLY:C	2.54	0.49
1:A:173:SER:C	1:A:445:ILE:HD12	2.38	0.49
1:A:79:GLN:NE2	1:A:388:LYS:HD2	2.28	0.49
1:A:82:GLU:HG2	1:A:385:VAL:HG12	1.95	0.48
1:A:461:ALA:HB3	1:A:466:MET:HG3	1.95	0.48
1:A:347:ARG:HG3	1:A:420:ALA:HB3	1.96	0.48
1:A:494:LEU:HA	1:A:497:LYS:HD2	1.95	0.48
1:A:38:ALA:HB2	1:A:63:GLU:HB2	1.96	0.47
1:A:81:TYR:CD1	1:A:384:PRO:HA	2.50	0.47
1:A:414:GLN:O	1:A:418:LEU:HB2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:368:THR:HB	1:A:374:LEU:HD13	1.96	0.47
1:A:317:MET:HG2	1:A:431:LEU:HD12	1.95	0.46
1:A:333:HIS:O	1:A:334:LYS:NZ	2.40	0.46
1:A:377:ASN:HD21	1:A:407:GLN:HE21	1.61	0.46
1:A:234:VAL:HG11	1:A:461:ALA:HB1	1.98	0.46
1:A:53:PHE:HB3	1:A:57:VAL:HG21	1.99	0.45
1:A:371:ASN:OD1	1:A:371:ASN:C	2.59	0.45
1:A:333:HIS:O	1:A:334:LYS:HD2	2.17	0.45
1:A:333:HIS:O	1:A:334:LYS:CG	2.65	0.45
1:A:495:TYR:O	1:A:499:LYS:HB2	2.16	0.45
1:A:334:LYS:HB3	1:A:459:TRP:HZ2	1.76	0.45
1:A:334:LYS:HG2	1:A:459:TRP:CH2	2.53	0.43
1:A:485:ASN:HB3	1:A:487:GLU:OE1	2.17	0.43
1:A:214:THR:HB	1:A:215:PRO:HD3	2.01	0.43
1:A:437:LYS:HD2	1:A:437:LYS:HA	1.56	0.43
1:A:217:MET:SD	1:A:227:ILE:HD11	2.59	0.43
1:A:364:THR:O	1:A:380:GLY:HA2	2.18	0.43
1:A:346:GLY:C	1:A:399:SER:HA	2.43	0.42
1:A:113:LEU:HD22	1:A:113:LEU:HA	1.74	0.42
1:A:146:LEU:HG	1:A:150:MET:HE2	2.02	0.42
1:A:398:GLY:O	1:A:399:SER:C	2.62	0.42
1:A:102:LEU:HD21	1:A:256:GLY:C	2.45	0.41
1:A:165:TYR:CZ	1:A:494:LEU:HB2	2.55	0.41
1:A:386:LYS:HA	1:A:386:LYS:HD2	1.79	0.41
1:A:314:ASN:CG	1:A:316:LYS:HG3	2.46	0.41
1:A:327:PRO:HB2	1:A:328:PRO:HD2	2.02	0.41
1:A:314:ASN:OD1	1:A:316:LYS:CG	2.68	0.41
1:A:404:VAL:HG11	1:A:410:LEU:HD11	2.03	0.41
1:A:53:PHE:CZ	1:A:479:TYR:HA	2.55	0.41
1:A:159:PHE:O	1:A:162:ILE:HG22	2.20	0.41
1:A:172:PHE:O	1:A:473:ASN:ND2	2.52	0.41
1:A:287:LYS:H	1:A:287:LYS:HE2	1.85	0.41
1:A:231:VAL:HG11	1:A:436:TYR:CE2	2.56	0.40
1:A:280:LYS:C	1:A:282:THR:N	2.79	0.40
1:A:358:PRO:O	1:A:359:ASP:C	2.64	0.40
1:A:79:GLN:OE1	1:A:384:PRO:HG3	2.21	0.40
1:A:222:GLY:HA2	1:A:377:ASN:HD22	1.86	0.40
1:A:258:LEU:HD12	1:A:258:LEU:HA	1.93	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	458/482 (95%)	427 (93%)	30 (7%)	1 (0%)	43	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	358	PRO

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	386/403 (96%)	355 (92%)	31 (8%)	11	37

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	LEU
1	A	121	THR
1	A	145	PHE
1	A	155	ILE
1	A	224	SER
1	A	227	ILE
1	A	280	LYS
1	A	285	THR
1	A	287	LYS
1	A	297	SER

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Mol	Chain	Res	Type
1	A	315	ARG
1	A	316	LYS
1	A	319	ASN
1	A	327	PRO
1	A	334	LYS
1	A	347	ARG
1	A	356	LYS
1	A	386	LYS
1	A	387	GLU
1	A	399	SER
1	A	424	SER
1	A	425	VAL
1	A	433	TYR
1	A	437	LYS
1	A	441	LYS
1	A	450	ARG
1	A	462	SER
1	A	486	THR
1	A	497	LYS
1	A	498	LEU
1	A	499	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	180	HIS
1	A	377	ASN
1	A	455	ASN

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](#)

2 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	FAD	A	602	-	58,58,58	0.52	0	85,89,89	0.50	0
2	CPS	A	601	-	25,25,45	0.51	0	39,41,70	0.57	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
3	FAD	A	602	-	-	12/34/50/50	0/6/6/6
2	CPS	A	601	-	-	-	0/4/4/4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	FAD	C5B-O5B-PA-O1A
3	A	602	FAD	C5B-O5B-PA-O2A
3	A	602	FAD	C5B-O5B-PA-O3P
3	A	602	FAD	C1'-C2'-C3'-C4'
3	A	602	FAD	C2'-C3'-C4'-O4'
3	A	602	FAD	O3'-C3'-C4'-O4'
3	A	602	FAD	O3'-C3'-C4'-C5'

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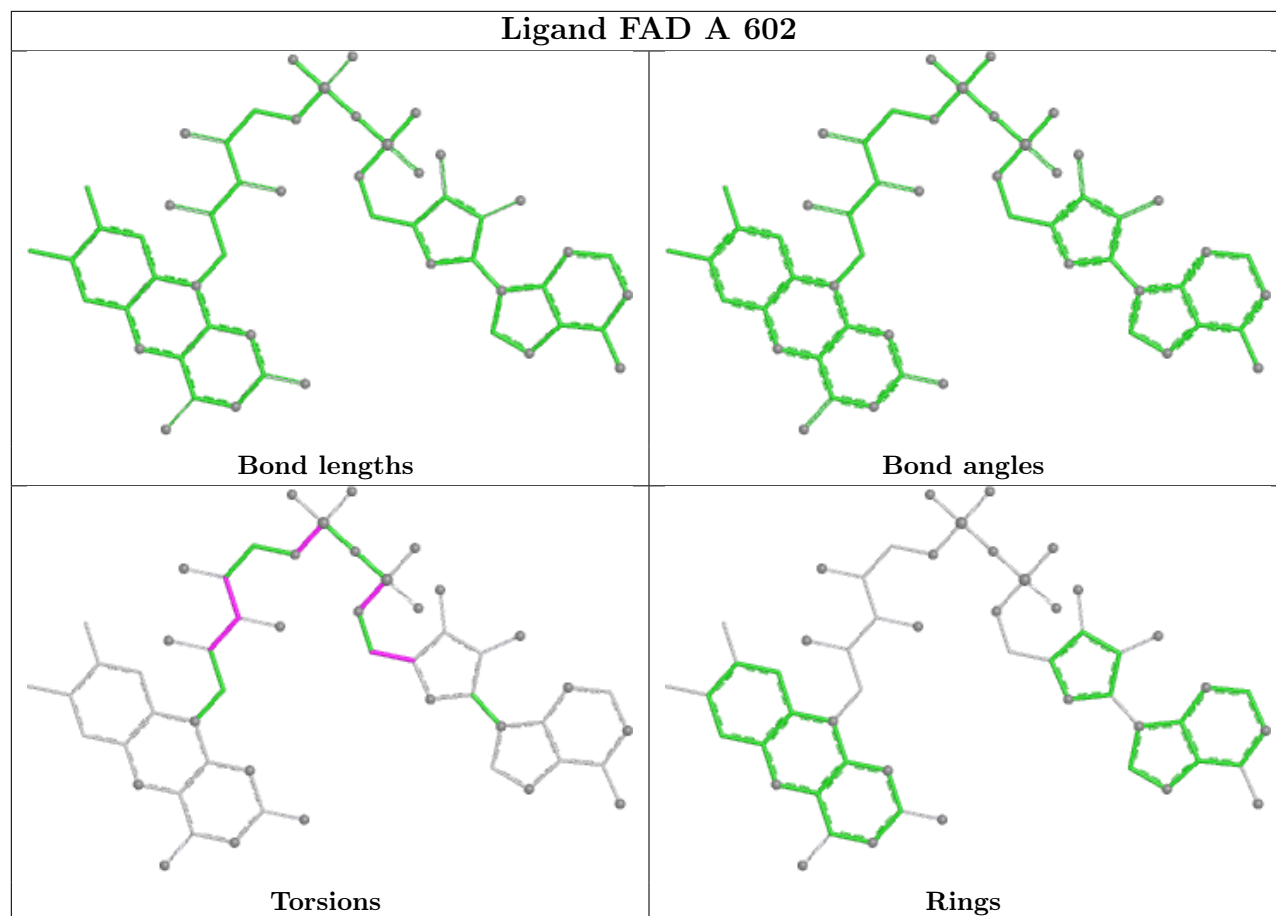
Mol	Chain	Res	Type	Atoms
3	A	602	FAD	C5'-O5'-P-O3P
3	A	602	FAD	C2'-C3'-C4'-C5'
3	A	602	FAD	O4B-C4B-C5B-O5B
3	A	602	FAD	C3B-C4B-C5B-O5B
3	A	602	FAD	C5'-O5'-P-O1P

There are no ring outliers.

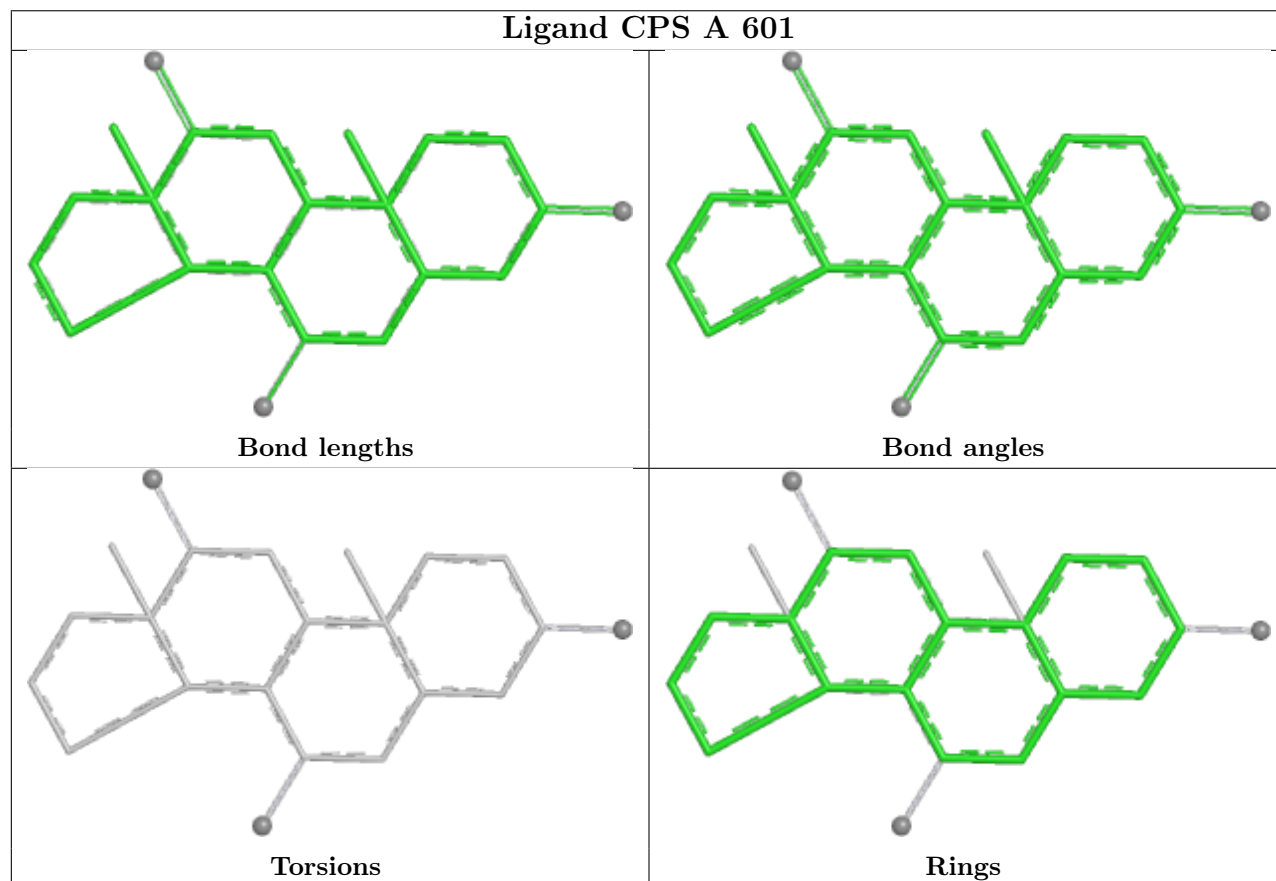
No monomer is involved in short contacts.

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 FAD A 602



Ligand CPS A 601



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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	464/482 (96%)	-0.07	14 (3%)	52 31	100, 147, 214, 258	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	331	GLU	3.9
1	A	281	GLN	3.7
1	A	114	MET	3.5
1	A	280	LYS	3.0
1	A	123	VAL	3.0
1	A	417	LEU	2.8
1	A	343	PHE	2.5
1	A	283	GLY	2.4
1	A	126	GLU	2.4
1	A	124	PHE	2.3
1	A	492	GLU	2.3
1	A	315	ARG	2.3
1	A	357	ALA	2.0
1	A	361	PHE	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](#)

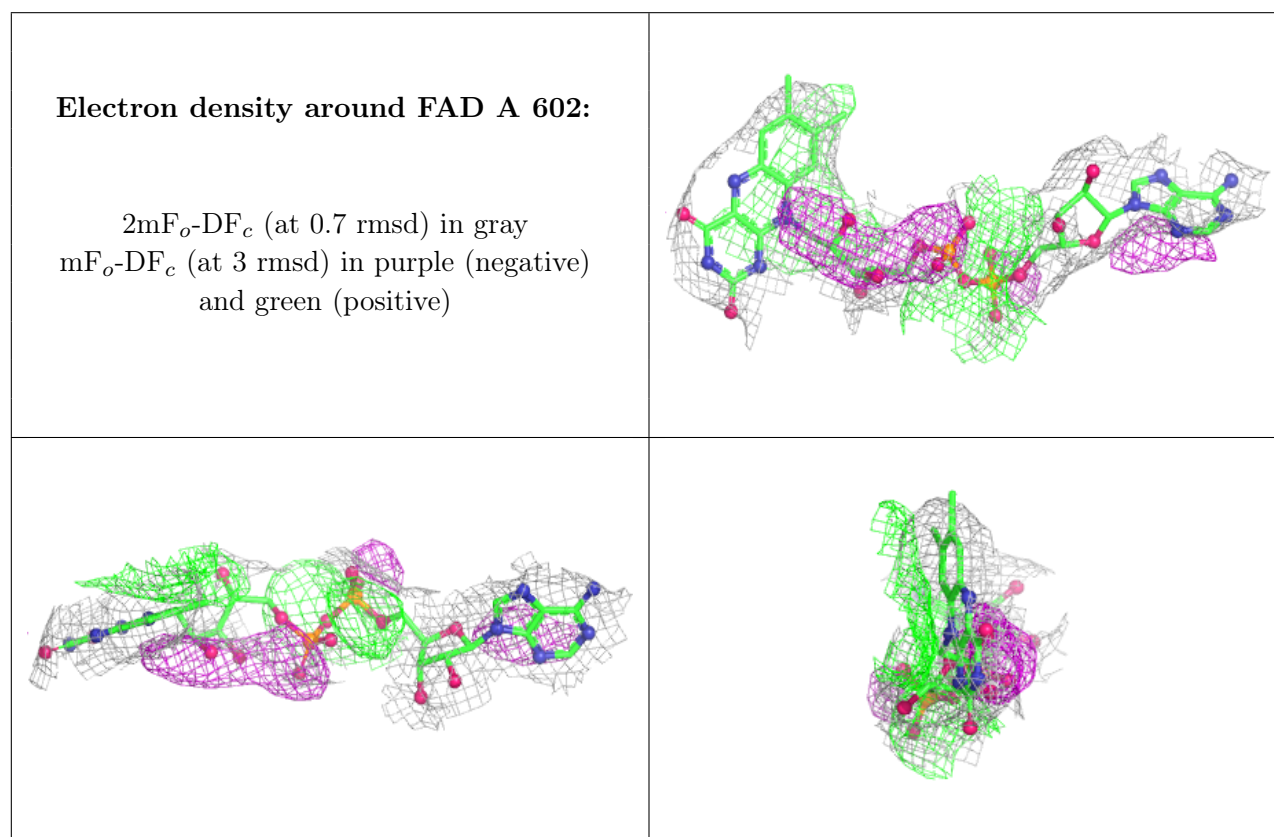
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

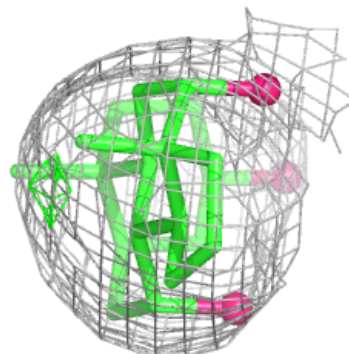
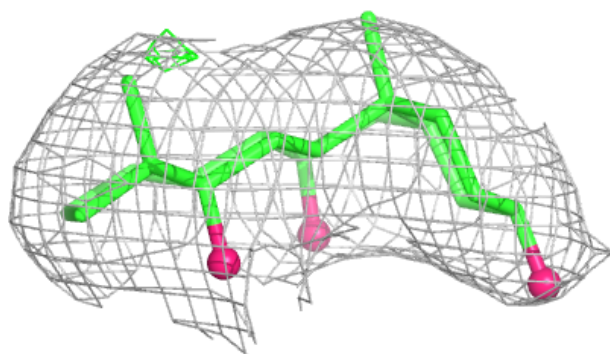
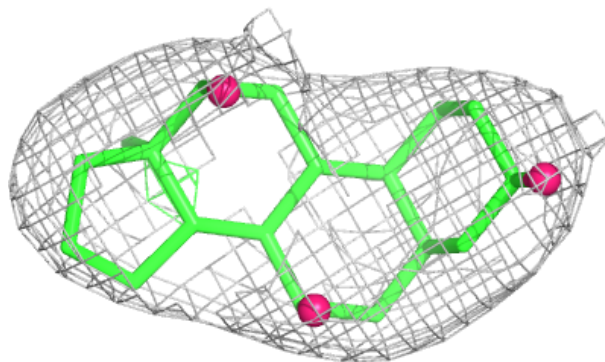
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	FAD	A	602	53/53	0.77	0.16	94,131,181,200	0
2	CPS	A	601	22/42	0.94	0.13	104,133,160,170	0

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 CPS A 601:

$2mF_o-DF_c$ (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.