

wwPDB X-ray Structure Validation Summary Report (i)

Feb 23, 2022 – 03:24 pm GMT

PDB ID	:	7Q86
Title	:	Crystal structure of human peroxisomal acyl-Co-A oxidase 1a, FAD-bound-
		form
Authors	:	Sonani, R.R.; Blat, A.; Dubin, G.
Deposited on	:	2021-11-10
Resolution	:	2.09 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

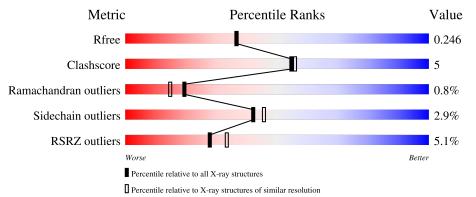
Xtriage (Phenix) EDS buster-report	: : :	1.8.4, CSD as541be (2020) 1.13 2.26 1.1.7 (2018)
		20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.26

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.09 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	5197(2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 <=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	А	667	8%	13%	8%
1	В	667	2% 8 1%	9%	• 9%



2 Entry composition (i)

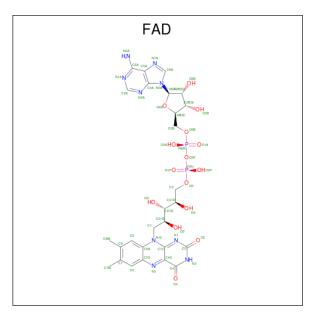
There are 3 unique types of molecules in this entry. The entry contains 10219 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Isoform 2 of Peroxisomal acyl-coenzyme A oxidase 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Δ	612	Total	С	Ν	0	\mathbf{S}	0	0	0
		012	4877	3107	846	895	29			
1	Р	607	Total	С	Ν	0	S	0	0	0
	D	007	4841	3086	839	887	29	0		

• Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
0	۸	1	Total	С	Ν	Ο	Р	0	0
		1	53	27	9	15	2	0	0
0	2 B	1	Total	С	Ν	Ο	Р	0	0
		1	53	27	9	15	2	0	0

• Molecule 3 is water.

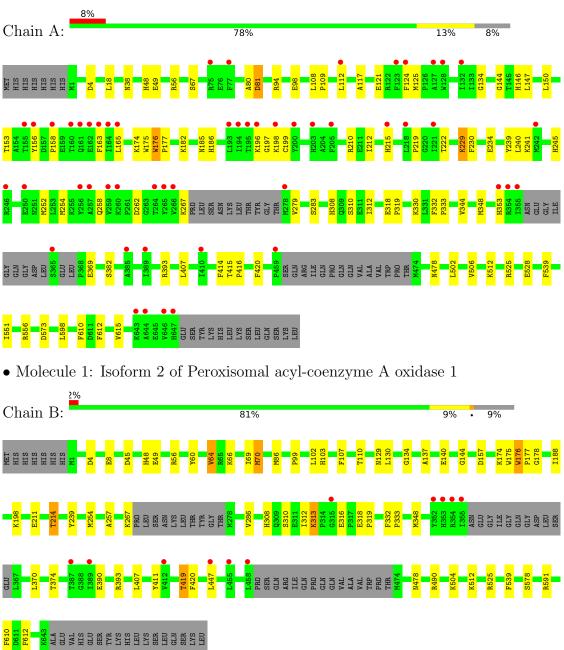


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	188	Total O 188 188	0	0
3	В	207	Total O 207 207	0	0



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: Isoform 2 of Peroxisomal acyl-coenzyme A oxidase 1

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	74.96Å 124.06Å 142.91Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.84 - 2.09	Depositor
Resolution (A)	47.79 - 2.09	EDS
% Data completeness	99.9 (47.84-2.09)	Depositor
(in resolution range)	99.9 (47.79-2.09)	EDS
R _{merge}	0.10	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.68 (at 2.08 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
D D.	0.187 , 0.240	Depositor
R, R_{free}	0.199 , 0.246	DCC
R_{free} test set	4054 reflections $(5.10%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	32.7	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	(Not available), (Not available)	EDS
L-test for twinning ²	$ L > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10219	wwPDB-VP
Average B, all atoms $(Å^2)$	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.22% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: CSO, 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	А	0.80	0/4970	0.91	0/6720	
1	В	0.84	0/4933	0.96	0/6671	
All	All	0.82	0/9903	0.93	0/13391	

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	А	4877	0	4858	62	0
1	В	4841	0	4830	41	0
2	А	53	0	31	0	0
2	В	53	0	31	0	0
3	А	188	0	0	7	0
3	В	207	0	0	2	0
All	All	10219	0	9750	102	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 5.

The worst 5 of 102 close contacts within the same asymmetric unit are listed below, sorted by



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:TYR:CE2	1:A:158:PRO:HA	1.93	1.03
1:A:156:TYR:CE2	1:A:158:PRO:CA	2.58	0.86
1:A:108:LEU:HD22	1:A:124:PHE:CE1	2.16	0.79
1:A:112:LEU:HD21	1:A:124:PHE:CZ	2.21	0.76
1:A:156:TYR:CZ	1:A:158:PRO:HA	2.25	0.71

their clash magnitude.

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	А	601/667~(90%)	576~(96%)	18 (3%)	7 (1%)	13 8
1	В	597/667~(90%)	581 (97%)	14 (2%)	2(0%)	41 41
All	All	1198/1334 (90%)	1157 (97%)	32 (3%)	9(1%)	19 15

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	176	TRP
1	В	176	TRP
1	А	197	GLY
1	А	279	VAL
1	А	81	ASP

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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.



Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
1	А	526/576~(91%)	509~(97%)	17 (3%)	39	41	
1	В	522/576~(91%)	509~(98%)	13 (2%)	47	52	
All	All	1048/1152~(91%)	1018 (97%)	30 (3%)	42	46	

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

 $5~{\rm of}~30$ residues with a non-rotameric side chain are listed below:

Mol	Chain	\mathbf{Res}	Type
1	А	528	GLU
1	В	504	LYS
1	В	56	ARG
1	В	578	SER
1	В	419	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	В	48	HIS
1	В	146	HIS
1	В	627	ASN
1	В	103	HIS
1	В	203	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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).



Mal	Mol Type Chain Res		Res	Res Link	B	Bond lengths			Bond angles		
10101	туре	Unam	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2	
1	CSO	А	199	1	3,6,7	1.32	1 (33%)	$0,\!6,\!8$	-	-	
1	CSO	В	449	1	3,6,7	0.60	0	$0,\!6,\!8$	-	-	
1	CSO	А	449	1	3,6,7	0.79	0	$0,\!6,\!8$	-	-	
1	CSO	В	199	1	3,6,7	1.12	0	$0,\!6,\!8$	-	-	

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
1	CSO	А	199	1	-	0/1/5/7	-
1	CSO	В	449	1	-	0/1/5/7	-
1	CSO	А	449	1	-	1/1/5/7	-
1	CSO	В	199	1	-	0/1/5/7	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
1	А	199	CSO	O-C	2.22	1.28	1.19

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	А	449	CSO	N-CA-CB-SG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

2 ligands are modelled in this entry.



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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 Cha	Chain	Chain Res	Res Link	Bo	Bond lengths			Bond angles		
10101	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	FAD	А	701	-	51,58,58	1.16	3 (5%)	60,89,89	1.79	9 (15%)
2	FAD	В	701	-	51,58,58	1.21	4 (7%)	60,89,89	2.09	8 (13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	А	701	-	-	1/30/50/50	0/6/6/6
2	FAD	В	701	-	-	2/30/50/50	0/6/6/6

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\operatorname{Observed}(\operatorname{\AA})$	$\mathrm{Ideal}(\mathrm{\AA})$
2	А	701	FAD	C4X-C10	5.88	1.44	1.38
2	В	701	FAD	C4X-C10	5.41	1.44	1.38
2	А	701	FAD	C4-N3	2.58	1.37	1.33
2	В	701	FAD	C4X-N5	-2.49	1.29	1.33
2	В	701	FAD	C4-N3	2.44	1.37	1.33

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
2	В	701	FAD	C2-N3-C4	9.66	123.30	115.14
2	А	701	FAD	C2-N3-C4	7.53	121.50	115.14
2	В	701	FAD	C10-C4X-N5	6.20	125.55	121.26
2	А	701	FAD	C10-C4X-C4	-5.20	116.51	119.95
2	В	701	FAD	C4X-C10-N10	-5.09	115.07	120.30

There are no chirality outliers.

All (3) torsion outliers are listed below:

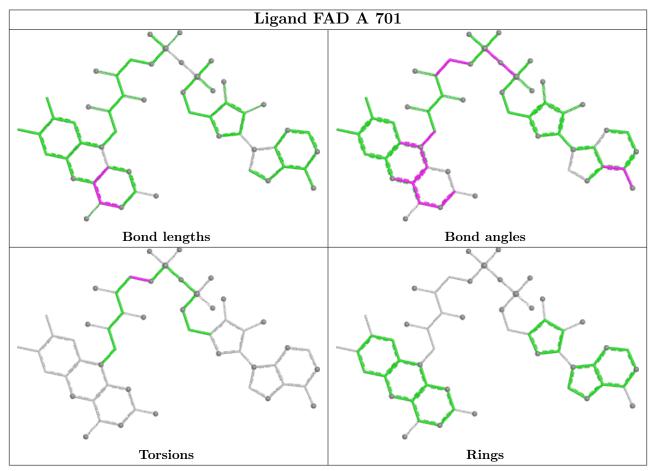


Mol	Chain	Res	Type	Atoms
2	В	701	FAD	O4B-C4B-C5B-O5B
2	А	701	FAD	C4'-C5'-O5'-P
2	В	701	FAD	C3B-C4B-C5B-O5B

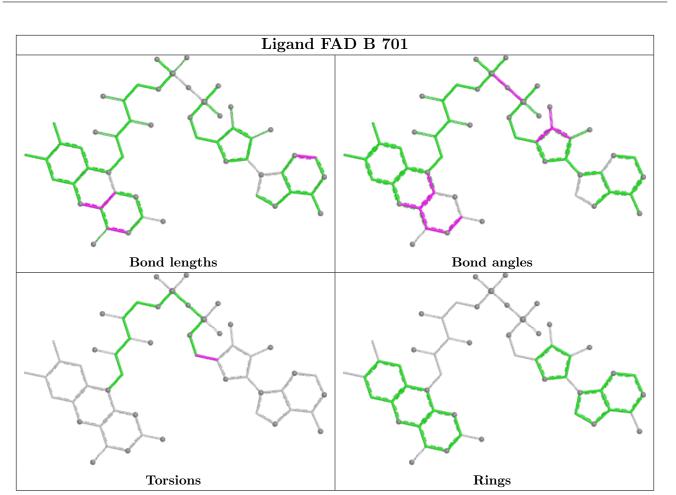
There are no ring outliers.

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 then 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 sufficient 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 (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, 95^{th} 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	$\langle RSRZ \rangle$	# RSRZ > 2		$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	$Q{<}0.9$	
1	А	610/667~(91%)	0.26	51 (8%)	11	14	22, 48, 101, 135	0
1	В	605/667~(90%)	-0.11	11 (1%)	68	72	21, 41, 73, 133	0
All	All	1215/1334 (91%)	0.08	62 (5%)	28	33	21, 43, 93, 135	0

The worst 5 of 62 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	128	TRP	11.3
1	А	256	TYR	7.7
1	А	156	TYR	6.0
1	А	124	PHE	5.8
1	А	200	TYR	5.6

6.2 Non-standard residues in protein, DNA, RNA chains (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, 95^{th} 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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
1	CSO	А	199	7/8	0.84	0.40	20,20,20,20	0
1	CSO	В	449	7/8	0.84	0.16	$63,\!65,\!78,\!82$	0
1	CSO	А	449	7/8	0.90	0.11	$62,\!68,\!77,\!88$	0
1	CSO	В	199	7/8	0.93	0.20	20,20,20,20	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

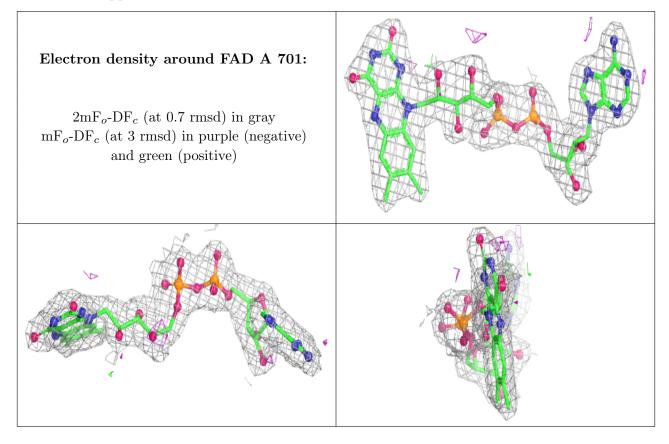


6.4 Ligands (i)

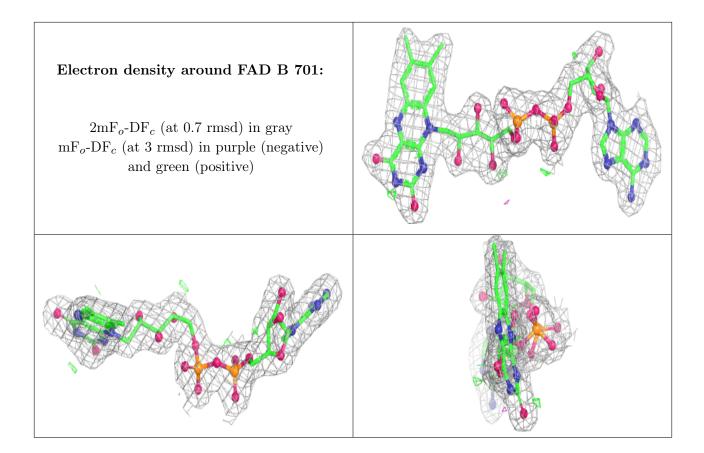
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
2	FAD	А	701	53/53	0.97	0.11	$35,\!42,\!49,\!51$	0
2	FAD	В	701	53/53	0.98	0.09	22,26,30,33	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.

