

Full wwPDB X-ray Structure Validation Report (i)

May 22, 2020 – 08:25 pm BST

PDB ID : 6C6P

Title: Human squalene epoxidase (SQLE, squalene monooxygenase) structure with

FAD and NB-598

Authors : Padyana, A.K.; Jin, L.

Deposited on : 2018-01-19

Resolution : 2.50 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 (i) symbol.

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

MolProbity : 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.11

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac: 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

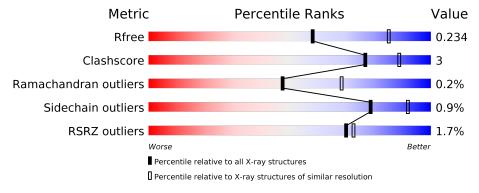
Validation Pipeline (wwPDB-VP) : 2.11

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.50 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	458	91%	7%	-			
1	В	458	88%	10%	-			



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 7583 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.

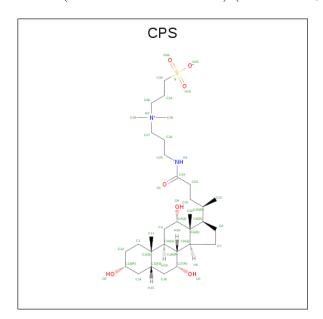
• Molecule 1 is a protein called Squalene monooxygenase.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	449	Total	С	N	О	S	0	9	0
1	A	449	3532	2285	594	638	15	U		0
1	B	449	Total	С	N	О	S	0	0	0
1	Б	449	3518	2276	592	635	15	U	U	

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	117	GLY	_	expression tag	UNP Q14534
В	117	GLY	-	expression tag	UNP Q14534

• Molecule 2 is 3-[(3-CHOLAMIDOPROPYL)DIMETHYLAMMONIO]-1-PROPANESULFO NATE (three-letter code: CPS) (formula: C₃₂H₅₈N₂O₇S).



I	Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
	2	A	1	Total C O 25 22 3	0	0

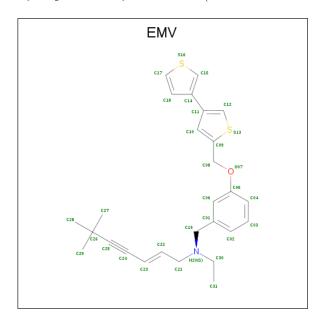
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N O 29 24 1 4	0	0
2	A	1	Total C O 25 22 3	0	0
2	В	1	Total C N O 30 25 1 4	0	0
2	В	1	Total C N O 30 25 1 4	0	0
2	В	1	Total C N O 30 25 1 4	0	0
2	В	1	Total C O 26 23 3	0	0

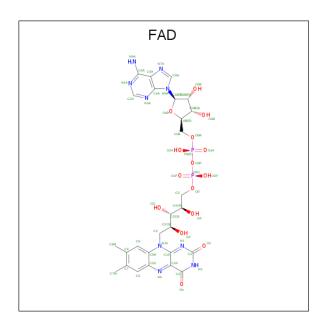
• Molecule 3 is (2E)-N-($\{3-[([3,3'-bithiophen]-5-yl)methoxy|phenyl\}methyl)$ -N-ethyl-6,6-dime thylhept-2-en-4-yn-1-amine (three-letter code: EMV) (formula: $C_{27}H_{31}NOS_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
9	Α	1	Total	С	N	О	S	0	0	
3	A	1	31	27	1	1	2	U	0	
9	D	1	Total	С	N	О	S	0	0	
)	D	1	31	27	1	1	2	U	0	

• Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).





Mol	Chain	Residues	${f Atoms}$					ZeroOcc	AltConf		
1	Α	Λ	Α	1	Total	С	N	О	Р	0	0
4	A	1	53	27	9	15	2	U			
4	D	1	Total	С	N	О	Р	0	0		
4	Б	1	53	27	9	15	2	U			

• Molecule 5 is water.

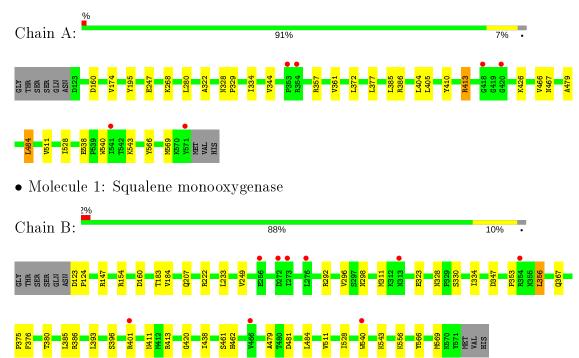
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	87	Total O 87 87	0	0
5	В	83	Total O 83 83	0	0



3 Residue-property plots (i)

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

• Molecule 1: Squalene monooxygenase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants	127.86Å 127.86Å 165.09Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.57 - 2.50	Depositor
Resolution (A)	40.57 - 2.50	EDS
% Data completeness	88.5 (40.57-2.50)	Depositor
(in resolution range)	88.5 (40.57-2.50)	EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	4.86 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575)	Depositor
D.D.	0.193 , 0.234	Depositor
R, R_{free}	0.193 , 0.234	DCC
R_{free} test set	2315 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	56.2	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32 , 42.8	EDS
L-test for twinning ²	$< L >=0.49, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.020 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7583	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.23% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $< L^2>$ 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: CPS, EMV, 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		
WIOI	Chain	RMSZ	# Z >5	RMSZ	# Z > 5	
1	A	0.25	0/3615	0.42	0/4895	
1	В	0.25	0/3601	0.42	0/4876	
All	All	0.25	0/7216	0.42	0/9771	

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)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
1	A	3532	0	3588	19	0
1	В	3518	0	3574	24	0
2	A	79	0	109	1	0
2	В	116	0	157	2	0
3	A	31	0	0	0	0
3	В	31	0	0	0	0
4	A	53	0	31	0	0
4	В	53	0	31	2	0
5	A	87	0	0	0	0
5	В	83	0	0	0	0
All	All	7583	0	7490	45	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 3.

All (45) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
		distance (Å)	overlap (Å)
1:B:393:LEU:O	1:B:411:ASN:ND2	2.27	0.65
1:A:247:GLU:HG2	1:A:268:LYS:HD2	1.79	0.63
1:A:174:VAL:HG21	1:A:426:LYS:HG3	1.80	0.62
1:B:540:TRP:HA	1:B:543:LYS:HB2	1.81	0.62
1:B:147:ARG:NH2	1:B:438:ILE:O	2.34	0.60
1:A:334:ILE:HD13	1:A:372:LEU:HD13	1.83	0.60
1:A:479:ALA:HB1	1:A:484:LEU:HB3	1.85	0.58
1:A:566:TYR:HA	1:A:569:MET:HE2	1.87	0.57
1:A:538:GLU:O	1:A:543:LYS:NZ	2.39	0.55
1:B:311:MET:HE3	1:B:375:PRO:HB2	1.88	0.55
1:A:357:ARG:HG3	1:A:385:LEU:HD22	1.92	0.52
1:B:479:ALA:HB1	1:B:484:LEU:HB3	1.93	0.51
1:B:356:LEU:HB3	1:B:385:LEU:HD12	1.94	0.49
2:A:601:CPS:H15	2:A:602:CPS:H8	1.94	0.49
1:B:481:ASP:HB3	1:B:484:LEU:HB2	1.95	0.48
1:B:184:VAL:HG11	1:B:222:ARG:HG3	1.96	0.48
1:B:330:SER:HB3	1:B:347:ASP:HB3	1.95	0.48
1:A:280:LEU:HD11	1:A:404:LEU:HG	1.96	0.48
1:A:361:VAL:HG13	1:A:377:LEU:HD11	1.96	0.47
1:A:540:TRP:CE3	1:A:543:LYS:HG3	2.50	0.46
1:B:328:ASN:O	1:B:367:GLN:NE2	2.48	0.45
1:B:160:ASP:OD1	1:B:386:ARG:NH2	2.48	0.45
1:A:413:ARG:HD3	1:A:467:ASN:OD1	2.17	0.45
1:B:123:ASP:HB2	1:B:124:PRO:HD2	1.99	0.44
1:A:528:ILE:HG13	2:B:602:CPS:H11A	2.01	0.43
1:A:405:LEU:HG	1:A:410:TYR:HB2	1.99	0.43
1:A:328:ASN:HA	1:A:329:PRO:HA	1.83	0.43
1:B:461:SER:OG	1:B:462:HIS:N	2.49	0.43
1:B:323:GLU:HB2	1:B:334:ILE:HB	2.01	0.43
1:B:183:THR:HG21	1:B:233:LEU:HD23	2.00	0.42
1:B:154:ARG:HG3	1:B:249:VAL:HG22	2.00	0.42
1:B:292:ARG:O	1:B:296:VAL:HG22	2.20	0.42
1:B:376:PHE:O	1:B:380:THR:HG23	2.19	0.42
1:A:195:TYR:HB3	1:A:322:ALA:HB3	2.01	0.42
1:B:528:ILE:HG13	2:B:601:CPS:H11A	2.02	0.42
1:B:420:GLY:HA3	4:B:606:FAD:H1'2	2.01	0.42
1:A:160:ASP:OD1	1:A:386:ARG:NH2	2.48	0.42

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Atom-1	Atom-2	$egin{aligned} & ext{Interatomic} \ & ext{distance} \ & ext{($\mathring{\mathbf{A}}$)} \end{aligned}$	Clash overlap (Å)
1:A:377:LEU:HD12	1:A:377:LEU:HA	1.85	0.42
1:A:413:ARG:HD2	1:A:466:VAL:HG12	2.02	0.42
1:B:566:TYR:HA	1:B:569:MET:HE2	2.01	0.41
4:B:606:FAD:H1'1	4:B:606:FAD:H9	1.87	0.41
1:B:353:PRO:HG2	1:B:356:LEU:HD22	2.03	0.41
1:B:298:ASN:OD1	1:B:396:SER:HB3	2.20	0.41
1:A:334:ILE:HG12	1:A:344:VAL:HG22	2.03	0.41
1:B:569:MET:HE2	1:B:569:MET:HB2	1.94	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	449/458 (98%)	437 (97%)	11 (2%)	1 (0%)	47	68
1	В	447/458 (98%)	437 (98%)	9 (2%)	1 (0%)	47	68
All	All	896/916 (98%)	874 (98%)	20 (2%)	2 (0%)	47	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	Α	511	VAL
1	В	511	VAL

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	Rotameric Outliers	
1	A	$385/391 \ (98\%)$	383 (100%)	2 (0%)	88 96
1	В	383/391 (98%)	378 (99%)	5 (1%)	69 87
All	All	768/782 (98%)	761 (99%)	7 (1%)	78 92

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	413	ARG
1	A	484	LEU
1	В	207	GLN
1	В	356	LEU
1	В	401	ARG
1	В	413	ARG
1	В	556	LYS

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

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

11 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	Т	Chain	n Res Link		Bo	nd leng	ths	Bond angles		
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EMV	A	604	-	32,33,33	0.47	0	35,44,44	1.08	3 (8%)
2	CPS	В	604	-	29,29,45	0.19	0	47,47,70	0.42	0
2	CPS	В	601	_	33,33,45	0.17	0	52,52,70	0.35	0
2	CPS	A	603	-	28,28,45	0.20	0	46,46,70	0.38	0
2	CPS	В	602	-	33,33,45	0.18	0	52,52,70	0.34	0
2	CPS	A	602	_	32,32,45	0.18	0	51,51,70	0.37	0
3	EMV	В	605	-	32,33,33	0.51	0	35,44,44	1.11	3 (8%)
2	CPS	A	601	-	28,28,45	0.18	0	46,46,70	0.36	0
4	FAD	В	606	-	51,58,58	1.21	5 (9%)	60,89,89	2.24	8 (13%)
2	CPS	В	603	-	33,33,45	0.18	0	52,52,70	0.37	0
4	FAD	A	605	_	51,58,58	1.19	5 (9%)	60,89,89	2.23	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
3	EMV	A	604	_	-	3/22/24/24	0/3/3/3
2	CPS	В	604	-	-	5/6/71/90	0/4/4/4
2	CPS	В	601	-	-	0/11/76/90	0/4/4/4
2	CPS	A	603	-	-	0/4/69/90	0/4/4/4
2	CPS	В	602	_	-	2/11/76/90	0/4/4/4
2	CPS	A	602	-	-	0/9/74/90	0/4/4/4
3	EMV	В	605	_	-	4/22/24/24	0/3/3/3
2	CPS	A	601	_	-	0/4/69/90	0/4/4/4
4	FAD	В	606	_	_	9/30/50/50	0/6/6/6
2	CPS	В	603	-	-	1/11/76/90	0/4/4/4
4	FAD	A	605	-	-	2/30/50/50	0/6/6/6

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$Ideal(\AA)$
4	A	605	FAD	C4X-C10	5.48	1.44	1.38
4	В	606	FAD	C4X-C10	5.48	1.44	1.38
4	В	606	FAD	C4-N3	3.12	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	${ m Observed}({ m \AA})$	$\operatorname{Ideal}(ext{\AA})$
4	A	605	FAD	C4-N3	2.95	1.38	1.33
4	A	605	FAD	C4-C4X	2.35	1.45	1.41
4	В	606	FAD	C4-C4X	2.33	1.45	1.41
4	A	605	FAD	C5X-N5	2.26	1.39	1.35
4	В	606	FAD	C5X-N5	2.25	1.39	1.35
4	В	606	FAD	C9A-N10	2.16	1.41	1.38
4	A	605	FAD	C9A-N10	2.14	1.41	1.38

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
4	В	606	FAD	C4-N3-C2	13.04	126.15	115.14
4	A	605	FAD	C4-N3-C2	13.00	126.12	115.14
4	В	606	FAD	C4X-C4-N3	-6.99	113.87	123.43
4	A	605	FAD	C4X-C4-N3	-6.95	113.93	123.43
4	В	606	FAD	C10-C4X-N5	4.73	124.53	121.26
4	A	605	FAD	C10-C4X-N5	4.70	124.50	121.26
4	A	605	FAD	C4-C4X-C10	-3.85	117.40	119.95
4	В	606	FAD	C4-C4X-C10	-3.84	117.41	119.95
3	В	605	EMV	C21-C22-C23	-3.66	118.64	124.68
3	A	604	EMV	C21-C22-C23	-3.62	118.70	124.68
4	В	606	FAD	C4X-C10-N10	-3.50	116.71	120.30
4	A	605	FAD	C4X-C10-N10	-3.46	116.75	120.30
4	A	605	FAD	C1'-N10-C9A	3.15	120.77	118.29
4	В	606	FAD	C1'-N10-C9A	3.04	120.68	118.29
3	В	605	EMV	C15-C14-C11	-2.88	123.66	127.58
4	В	606	FAD	P-O3P-PA	-2.80	123.22	132.83
3	A	604	EMV	C15-C14-C11	-2.78	123.79	127.58
3	A	604	EMV	C12-C11-C14	-2.66	123.95	127.58
3	В	605	EMV	C12-C11-C14	-2.65	123.97	127.58
4	A	605	FAD	P-O3P-PA	-2.43	124.49	132.83
4	A	605	FAD	C5A-C6A-N6A	2.29	123.83	120.35
4	В	606	FAD	C5A-C6A-N6A	2.28	123.82	120.35

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	604	EMV	C22-C23-C24-C25
3	В	605	EMV	C22-C23-C24-C25
4	В	606	FAD	C5B-O5B-PA-O2A
4	A	605	FAD	O4B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
4	В	606	FAD	O4B-C4B-C5B-O5B
4	В	606	FAD	C3B-C4B-C5B-O5B
4	A	605	FAD	C3B-C4B-C5B-O5B
3	A	604	EMV	C23-C24-C25-C26
2	В	602	CPS	C22-C23-C24-N1
2	В	602	CPS	C22-C23-C24-O1
2	В	603	CPS	C20-C22-C23-C24
2	В	604	CPS	C9-C20-C22-C23
4	В	606	FAD	C5B-O5B-PA-O3P
4	В	606	FAD	C5'-O5'-P-O3P
4	В	606	FAD	C5B-O5B-PA-O1A
3	В	605	EMV	C04-C05-O07-C08
3	В	605	EMV	C06-C05-O07-C08
3	В	605	EMV	C23-C24-C25-C26
4	В	606	FAD	PA-O3P-P-O1P
2	В	604	CPS	C22-C20-C9-C8
2	В	604	CPS	C22-C20-C9-C5
2	В	604	CPS	C21-C20-C22-C23
3	A	604	EMV	C12-C11-C14-C15
4	В	606	FAD	PA-O3P-P-O2P
2	В	604	CPS	C21-C20-C9-C5
4	В	606	FAD	C5'-O5'-P-O1P

There are no ring outliers.

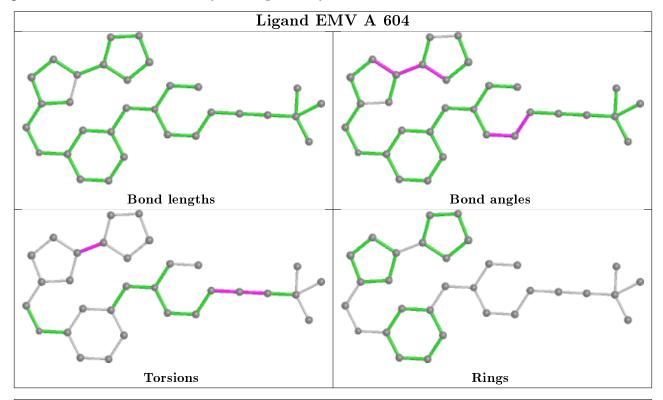
5 monomers are involved in 5 short contacts:

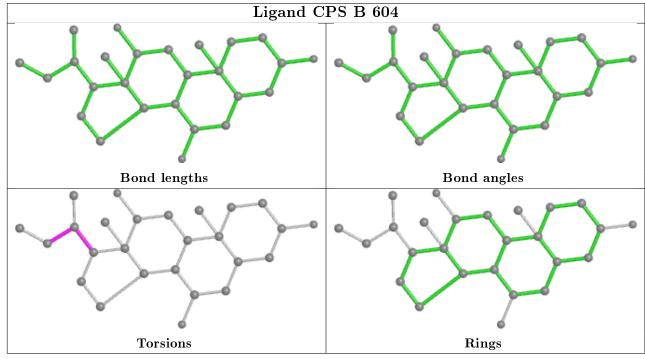
Mol	Chain	${f Res}$	Type	Clashes	Symm-Clashes
2	В	601	CPS	1	0
2	В	602	CPS	1	0
2	A	602	CPS	1	0
2	A	601	CPS	1	0
4	В	606	FAD	2	0

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 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

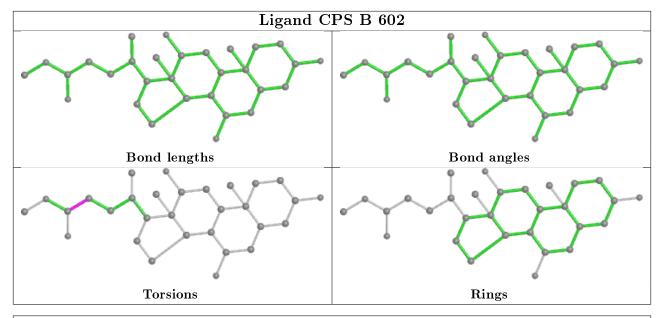


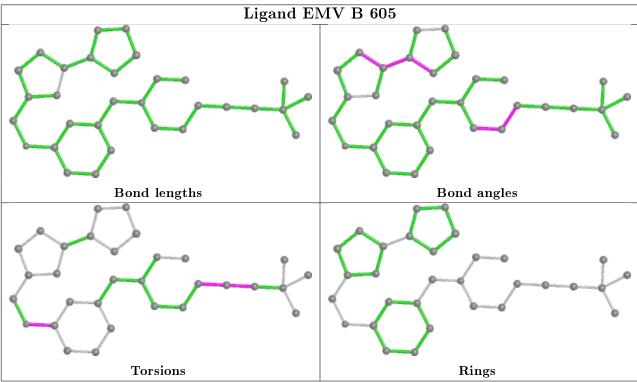
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



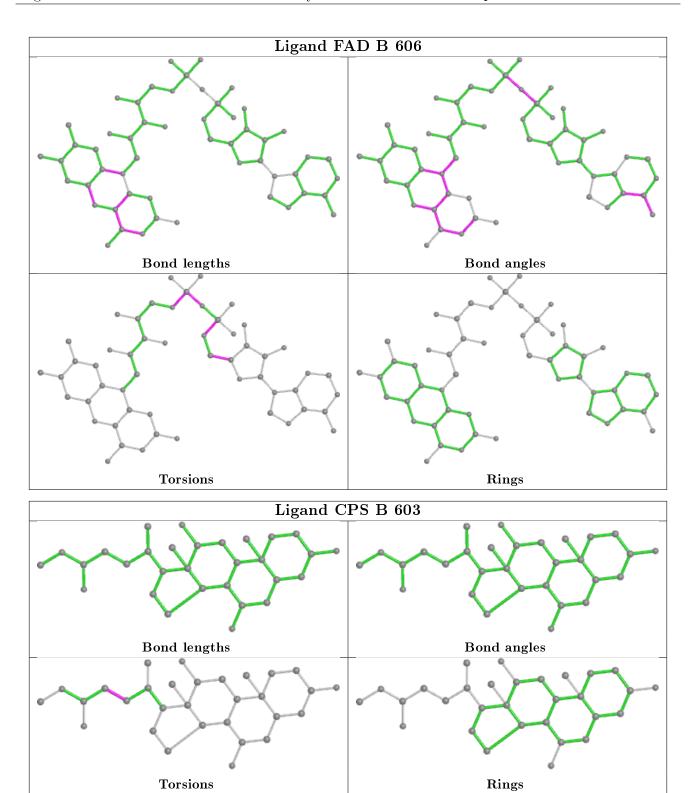




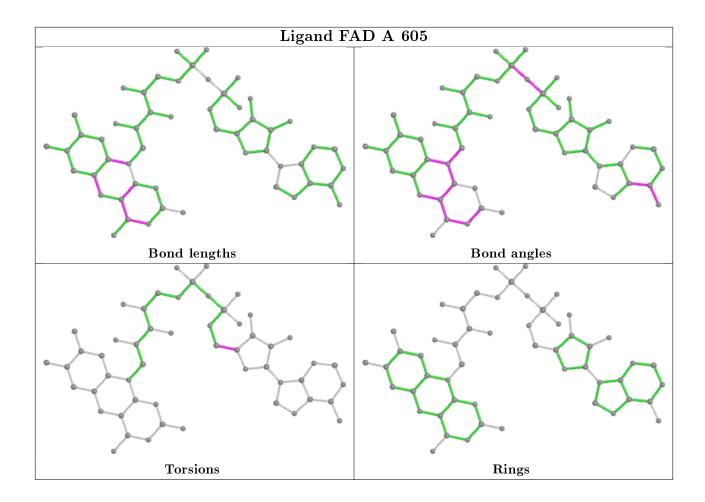












5.7 Other polymers (i)

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 (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	<RSRZ $>$	$\#\mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q < 0.9
1	A	449/458 (98%)	-0.24	6 (1%) 77 79	37, 56, 85, 108	0
1	В	449/458 (98%)	-0.02	9 (2%) 65 68	38, 58, 88, 106	0
All	All	898/916 (98%)	-0.13	15 (1%) 70 72	37, 57, 87, 108	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	541	ILE	4.5
1	A	354	ARG	3.7
1	В	540	TRP	3.5
1	В	272	ASP	3.0
1	В	273	ILE	2.7
1	A	353	PRO	2.5
1	A	571	TYR	2.5
1	В	256	GLU	2.5
1	В	276	LEU	2.4
1	В	354	ARG	2.3
1	В	313	ASN	2.2
1	В	401	ARG	2.2
1	A	420	GLY	2.1
1	В	466	VAL	2.1
1	A	418	GLY	2.0

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

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

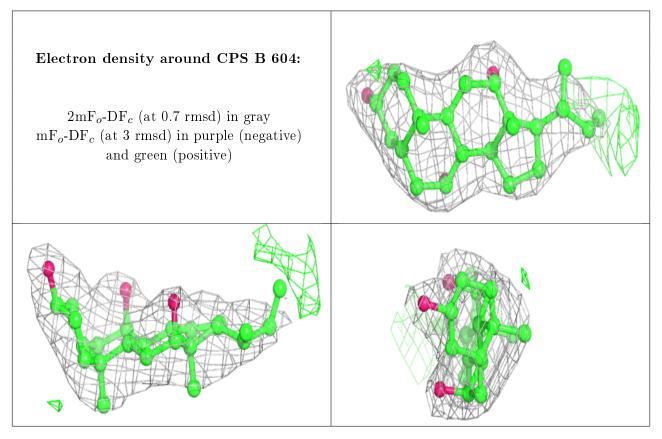


6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
2	CPS	В	604	26/42	0.92	0.23	66,73,81,83	0
2	CPS	A	602	29/42	0.93	0.16	60,70,83,87	0
2	CPS	В	602	30/42	0.94	0.13	40,49,87,93	0
2	CPS	A	603	25/42	0.94	0.16	47,57,73,78	0
3	EMV	В	605	31/31	0.94	0.19	38,54,60,62	0
2	CPS	В	603	30/42	0.94	0.14	60,69,86,92	0
2	CPS	В	601	30/42	0.95	0.13	41,48,83,90	0
3	EMV	A	604	31/31	0.95	0.20	42,53,61,65	0
4	FAD	A	605	53/53	0.95	0.17	46,52,65,67	0
2	CPS	A	601	25/42	0.96	0.21	67,74,80,81	0
4	FAD	В	606	53/53	0.96	0.17	44,55,74,75	0

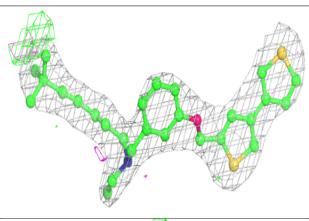
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

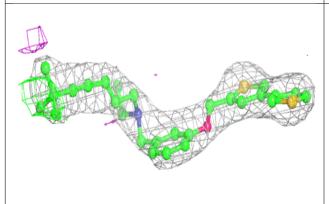


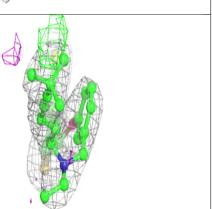


Electron density around CPS B 602: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray mF_o -DF_c (at 3 rmsd) in purple (negative) and green (positive) Electron density around EMV B 605:

 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)









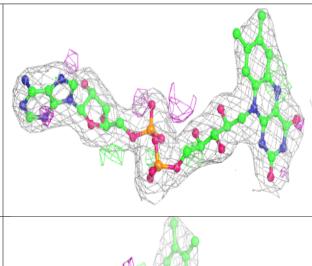
Electron density around CPS B 603: 2mF_o-DF_c (at 0.7 rmsd) in gray mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive) Electron density around EMV A 604:

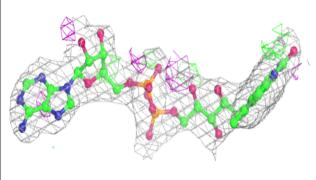
Electron density around EMV A 604: 2mF_o-DF_c (at 0.7 rmsd) in gray mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

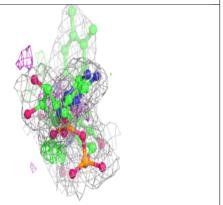


Electron density around FAD B 606:

 $2 {\rm mF}_o\text{-}{\rm DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)









6.5 Other polymers (i)

There are no such residues in this entry.

