

# wwPDB X-ray Structure Validation Summary Report (i)

#### Nov 13, 2023 – 12:21 am GMT

PDB ID : 8OY3

Title : Time-resolved SFX structure of the class II photolyase complexed with a

thymine dimer (3 picosecond pump-probe delay)

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Mashhour, A.R.; Galchenkova, M.; Gunther, S.; Reinke, P.; Kremling, V.; Oberthuer, D.; Henkel, A.; Sprenger, J.; Scheer, T.E.S.; Lange, E.; Yefanov, O.N.; Middendorf, P.; Sellberg, J.A.; Schubert, R.; Fadini, A.; Cirelli, C.; Beale, E.V.; Johnson, P.; Dworkowski, F.; Ozerov, D.; Bertrand, Q.; Wranik,

M.; Zitter, E.D.; Turk, D.; Bajt, S.; Chapman, H.; Bacellar, C.

Deposited on : 2023-05-03

Resolution : 2.16 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 types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

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

MolProbity: 4.02b-467

Mogul: 1.8.4, CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.36

buster-report : 1.1.7 (2018)

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

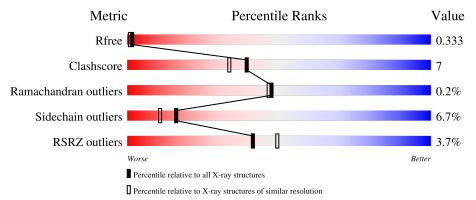
Refmac : 5.8.0158

## 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 2.16 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}({\rm \AA})) \end{array}$
$R_{free}$	130704	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	A	498	76%		13%	• 9%
1	В	498	65%	18%	•	14%

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CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001)

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

Validation Pipeline (wwPDB-VP) : 2.36



Mol	Chain	Length	Quality of chain						
2	С	13	62%		31%	8%			
2	Е	13	23%	62%		15%			
3	D	14	57%		43%				
3	F	14	43%	50%		7%			



## 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 8614 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 Deoxyribodipyrimidine photo-lyase.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	A	453	Total 3701	C 2377	N 623	O 685	S 16	0	3	0
1	В	427	Total 3486	C 2240	N 588	O 643	S 15	0	1	0

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q8PYK9
A	-18	GLY	-	expression tag	UNP Q8PYK9
A	-17	SER	-	expression tag	UNP Q8PYK9
A	-16	SER	-	expression tag	UNP Q8PYK9
A	-15	HIS	-	expression tag	UNP Q8PYK9
A	-14	HIS	-	expression tag	UNP Q8PYK9
A	-13	HIS	-	expression tag	UNP Q8PYK9
A	-12	HIS	-	expression tag	UNP Q8PYK9
A	-11	HIS	-	expression tag	UNP Q8PYK9
A	-10	HIS	-	expression tag	UNP Q8PYK9
A	-9	SER	-	expression tag	UNP Q8PYK9
A	-8	SER	-	expression tag	UNP Q8PYK9
A	-7	GLY	-	expression tag	UNP Q8PYK9
A	-6	LEU	-	expression tag	UNP Q8PYK9
A	-5	VAL	-	expression tag	UNP Q8PYK9
A	-4	PRO	-	expression tag	UNP Q8PYK9
A	-3	ARG	-	expression tag	UNP Q8PYK9
A	-2	GLY	-	expression tag	UNP Q8PYK9
A	-1	SER	-	expression tag	UNP Q8PYK9
A	0	HIS	-	expression tag	UNP Q8PYK9
A	465	ASP	-	expression tag	UNP Q8PYK9
A	466	LYS	-	expression tag	UNP Q8PYK9
A	467	LEU	-	expression tag	UNP Q8PYK9
A	468	ALA	-	expression tag	UNP Q8PYK9
A	469	ALA	-	expression tag	UNP Q8PYK9



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Chain	Residue	Modelled  Modelled	Actual	Comment	Reference
A	470	ALA	-	expression tag	UNP Q8PYK9
A	471	LEU	-	expression tag	UNP Q8PYK9
A	472	GLU	-	expression tag	UNP Q8PYK9
A	473	HIS	-	expression tag	UNP Q8PYK9
A	474	HIS	-	expression tag	UNP Q8PYK9
A	475	HIS	-	expression tag	UNP Q8PYK9
A	476	HIS	-	expression tag	UNP Q8PYK9
A	477	HIS	-	expression tag	UNP Q8PYK9
A	478	HIS	-	expression tag	UNP Q8PYK9
В	-19	MET	-	initiating methionine	UNP Q8PYK9
В	-18	GLY	-	expression tag	UNP Q8PYK9
В	-17	SER	-	expression tag	UNP Q8PYK9
В	-16	SER	-	expression tag	UNP Q8PYK9
В	-15	HIS	-	expression tag	UNP Q8PYK9
В	-14	HIS	-	expression tag	UNP Q8PYK9
В	-13	HIS	-	expression tag	UNP Q8PYK9
В	-12	HIS	-	expression tag	UNP Q8PYK9
В	-11	HIS	-	expression tag	UNP Q8PYK9
В	-10	HIS	-	expression tag	UNP Q8PYK9
В	-9	SER	-	expression tag	UNP Q8PYK9
В	-8	SER	-	expression tag	UNP Q8PYK9
В	-7	GLY	-	expression tag	UNP Q8PYK9
В	-6	LEU	-	expression tag	UNP Q8PYK9
В	-5	VAL	-	expression tag	UNP Q8PYK9
В	-4	PRO	-	expression tag	UNP Q8PYK9
В	-3	ARG	-	expression tag	UNP Q8PYK9
В	-2	GLY	_	expression tag	UNP Q8PYK9
В	-1	SER	-	expression tag	UNP Q8PYK9
В	0	HIS	_	expression tag	UNP Q8PYK9
В	465	ASP	-	expression tag	UNP Q8PYK9
В	466	LYS	-	expression tag	UNP Q8PYK9
В	467	LEU	_	expression tag	UNP Q8PYK9
В	468	ALA	-	expression tag	UNP Q8PYK9
В	469	ALA	-	expression tag	UNP Q8PYK9
В	470	ALA	-	expression tag	UNP Q8PYK9
В	471	LEU	-	expression tag	UNP Q8PYK9
В	472	GLU	-	expression tag	UNP Q8PYK9
В	473	HIS	-	expression tag	UNP Q8PYK9
В	474	HIS	-	expression tag	UNP Q8PYK9
В	475	HIS	-	expression tag	UNP Q8PYK9
В	476	HIS	-	expression tag	UNP Q8PYK9
В	477	HIS	-	expression tag	UNP Q8PYK9



Chain	Residue	Modelled	Actual	Comment	Reference
В	478	HIS	-	expression tag	UNP Q8PYK9

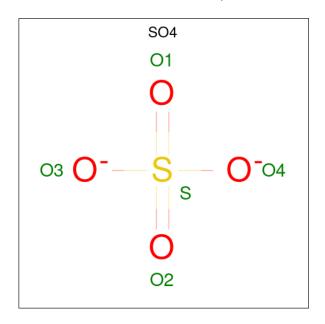
• Molecule 2 is a DNA chain called CPD-COMPRISING OLIGONUCLEOTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	С	13	Total	С	N	О	Р	0	0	0
		10	282	135	51	83	13	U	0	
9	E	11	Total	С	N	О	Р	0	0	0
	E	11	244	115	44	73	12	0	U	U

• Molecule 3 is a DNA chain called COUNTERSTRAND-OLIGONUCLEOTIDE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	D	1.4	Total	С	N	О	Р	0	0	0
3	3 D	14	286	136	56	81	13	U	U	
9	E	19	Total	С	N	О	Р	0	0	0
3	3 F	13	265	126	51	76	12	0	U	

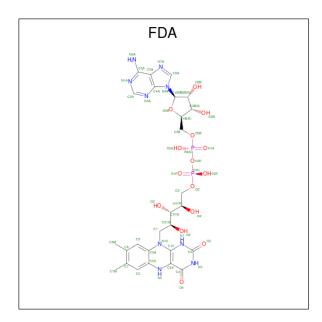
• Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 5	O 4	S 1	0	0

• Molecule 5 is DIHYDROFLAVINE-ADENINE DINUCLEOTIDE (three-letter code: FDA) (formula:  $C_{27}H_{35}N_9O_{15}P_2$ ).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
5	Λ	1	Total	С	N	О	Р	0	0	
9	3 A	1	53	27	9	15	2	0	0	
5	D	1	Total	С	N	О	Р	0	0	
9	5 B	1	53	27	9	15	2	U		

### • Molecule 6 is water.

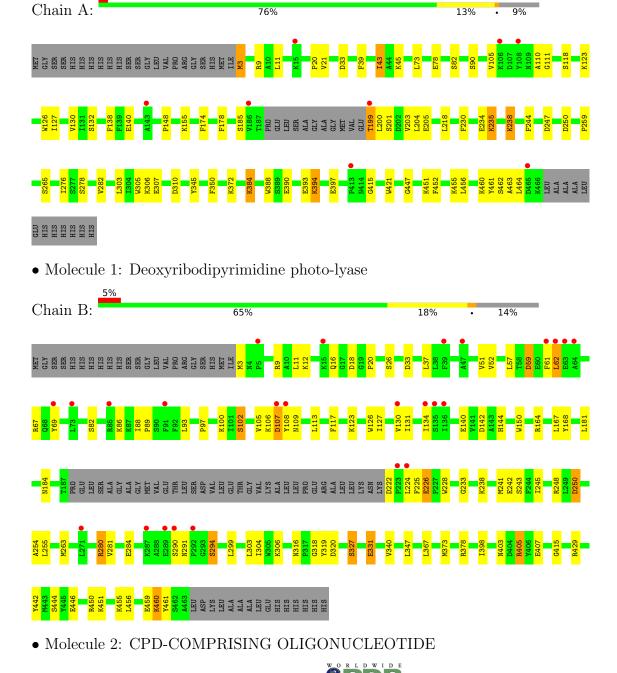
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	140	Total O 140 140	0	0
6	В	65	Total O 65 65	0	0
6	С	14	Total O 14 14	0	0
6	D	12	Total O 12 12	0	0
6	Е	1	Total O 1 1	0	0
6	F	7	Total O 7 7	0	0

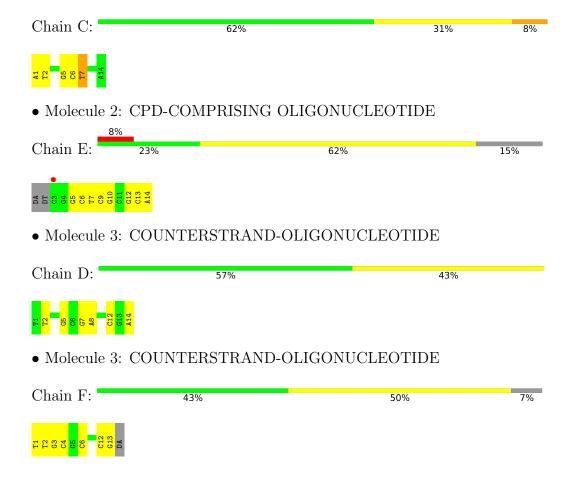


## 3 Residue-property plots (i)

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: Deoxyribodipyrimidine photo-lyase







## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	70.20Å 117.76Å 170.47Å	Donositon
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.29 - 2.16	Depositor
Resolution (A)	31.29 - 2.16	EDS
% Data completeness	79.9 (31.29-2.16)	Depositor
(in resolution range)	79.9 (31.29-2.16)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.57 (at 2.16Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
D D.	0.273 , 0.335	Depositor
$R, R_{free}$	0.275 , $0.333$	DCC
$R_{free}$ test set	3070 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.8	Xtriage
Anisotropy	0.586	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.31 , 61.6	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.52, < L^2>=0.36$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8614	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.19% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical 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.



<sup>&</sup>lt;sup>1</sup>Intensities 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: SO4, TTD, FDA

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	Bo	ond angles
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5
1	A	0.55	0/3802	0.65	0/5139
1	В	0.47	0/3583	0.61	0/4843
2	С	0.97	0/270	0.96	0/412
2	Е	0.87	0/227	0.74	0/345
3	D	0.87	0/321	1.03	1/494~(0.2%)
3	F	0.87	0/297	0.90	0/457
All	All	0.58	0/8500	0.69	1/11690 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
3	D	12	DC	OP2-P-O3'	5.16	116.55	105.20

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3701	0	3625	34	0
1	В	3486	0	3389	63	0
2	С	282	0	159	7	0
2	Е	244	0	135	8	0



n previous	paae
	n previous

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	286	0	158	4	0
3	F	265	0	147	7	0
4	A	5	0	0	0	0
5	A	53	0	33	1	0
5	В	53	0	33	2	0
6	A	140	0	0	3	0
6	В	65	0	0	3	0
6	С	14	0	0	0	0
6	D	12	0	0	2	0
6	Ε	1	0	0	0	0
6	F	7	0	0	0	0
All	All	8614	0	7679	117	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 7.

The worst 5 of 117 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:A:390:GLU:OE2	1:A:394:LYS:HD3	1.88	0.73
1:B:243:SER:OG	1:B:248:ARG:NH2	2.27	0.68
1:A:21:VAL:HG21	1:A:43:ILE:HG22	1.78	0.65
1:B:61:PHE:HB3	1:B:69:TYR:HD1	1.63	0.64
1:B:12:LYS:HB2	1:B:131:ILE:HG23	1.80	0.63

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	452/498 (91%)	429 (95%)	22 (5%)	1 (0%)	47 46



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	В	424/498 (85%)	406 (96%)	17 (4%)	1 (0%)	47 46
All	All	876/996 (88%)	835 (95%)	39 (4%)	2 (0%)	47 46

#### All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	319	TYR
1	A	463	ALA

#### 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	397/429 (92%)	373 (94%)	24 (6%)	19 14
1	В	372/429 (87%)	345 (93%)	27 (7%)	14 9
All	All	769/858 (90%)	718 (93%)	51 (7%)	16 11

5 of 51 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	57	LEU
1	В	144	HIS
1	В	456	LEU
1	В	59	ASP
1	В	93	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	В	272	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)

2 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).

Mol	Trme	Chain	Dec	Link Bond lengths			В	ond ang	les	
IVIOI	Type	Chain	Res	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	TTD	С	7	2	42,45,46	0.53	1 (2%)	62,74,77	0.70	0
2	TTD	Е	7	2	42,45,46	0.37	0	62,74,77	0.66	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
2	TTD	С	7	2	-	11/22/109/110	0/5/6/6
2	TTD	E	7	2	-	11/22/109/110	0/5/6/6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(\text{\AA})$
2	С	7	TTD	PB-O3R	2.33	1.66	1.60

There are no bond angle outliers.

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	С	7	TTD	C2'-C1'-N1-C6
2	Е	7	TTD	C2'-C1'-N1-C6
2	Е	7	TTD	C2'-C1'-N1-C2



Mol	Chain	Res	Type	Atoms
2	Е	7	TTD	O4R-C4'-C5R-O5R
2	Е	7	TTD	O4'-C1'-N1-C6

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	С	7	TTD	5	0
2	Е	7	TTD	2	0

### 5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry (i)

3 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	Type Chain	Chain	Res	es Link Bond lengths				В	ond ang	cles
MIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2	
5	FDA	В	501	-	52,58,58	0.65	0	60,89,89	0.80	3 (5%)	
5	FDA	A	502	-	52,58,58	0.63	0	60,89,89	0.82	3 (5%)	
4	SO4	A	501	-	4,4,4	0.14	0	6,6,6	0.64	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
5	FDA	В	501	-	-	5/30/50/50	0/6/6/6
5	FDA	A	502	-	-	4/30/50/50	0/6/6/6



There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	${f Z}$	$Observed(^{o})$	$Ideal(^{o})$
5	A	502	FDA	C5A-C6A-N6A	2.64	124.36	120.35
5	В	501	FDA	N3-C2-N1	2.42	119.70	115.80
5	A	502	FDA	N3-C2-N1	2.30	119.50	115.80
5	A	502	FDA	O2-C2-N3	-2.19	117.68	121.82
5	В	501	FDA	O2-C2-N3	-2.13	117.80	121.82

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	502	FDA	C5B-O5B-PA-O1A
5	В	501	FDA	C5B-O5B-PA-O1A
5	В	501	FDA	N10-C1'-C2'-C3'
5	В	501	FDA	O4'-C4'-C5'-O5'
5	В	501	FDA	C5B-O5B-PA-O3P

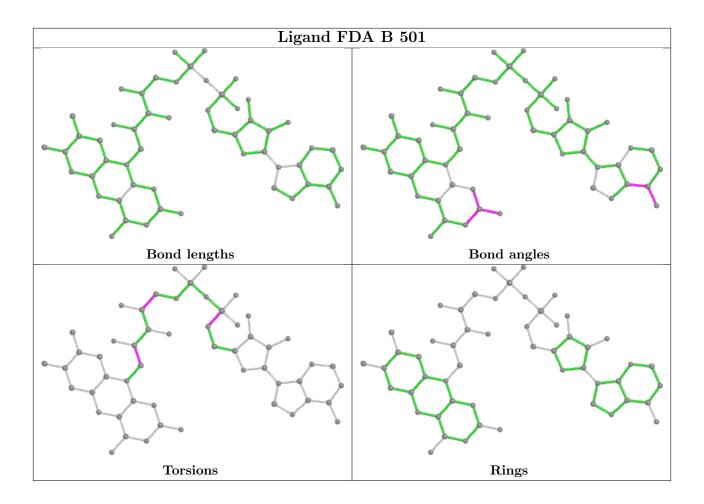
There are no ring outliers.

2 monomers are involved in 3 short contacts:

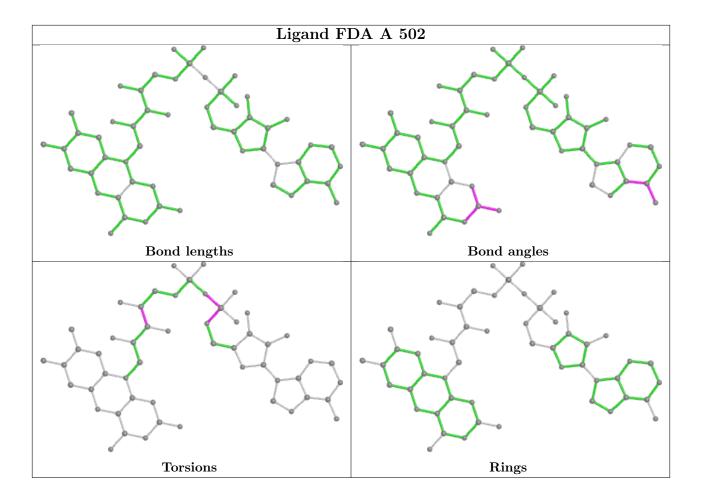
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	В	501	FDA	2	0
5	A	502	FDA	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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></rsrz>	$\# \mathrm{RSRZ} {>} 2$	$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q < 0.9
1	A	453/498 (90%)	0.03	8 (1%) 68 75	25, 39, 70, 109	0
1	В	427/498~(85%)	0.51	25 (5%) 22 30	32, 58, 91, 141	0
2	С	12/13 (92%)	0.39	0 100 100	37, 70, 88, 93	0
2	E	10/13~(76%)	0.46	1 (10%) 7 11	52, 76, 110, 131	0
3	D	14/14 (100%)	0.04	0 100 100	50, 67, 90, 95	0
3	F	13/14 (92%)	0.54	0 100 100	71, 79, 109, 109	0
All	All	929/1050~(88%)	0.27	34 (3%) 41 49	25, 49, 88, 141	0

The worst 5 of 34 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	136	ILE	5.0
1	В	61	PHE	4.6
1	В	64	ALA	4.2
1	В	290	SER	4.2
1	В	135	SER	4.0

### 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	TTD	Ε	7	40/41	0.96	0.15	39,47,67,74	0
2	TTD	С	7	40/41	0.97	0.14	25,34,49,56	0



## 6.3 Carbohydrates (i)

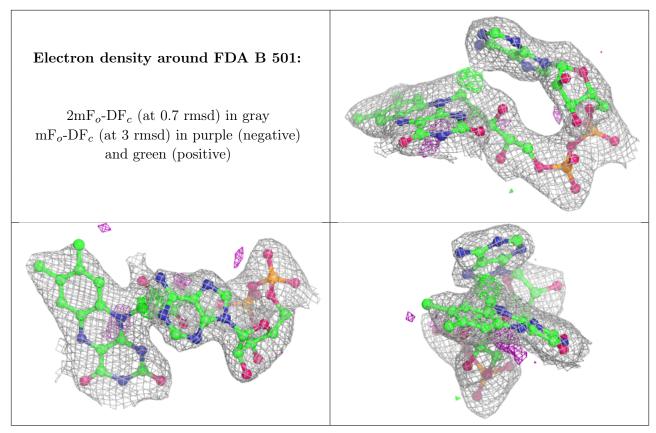
There are no monosaccharides in this entry.

## 6.4 Ligands (i)

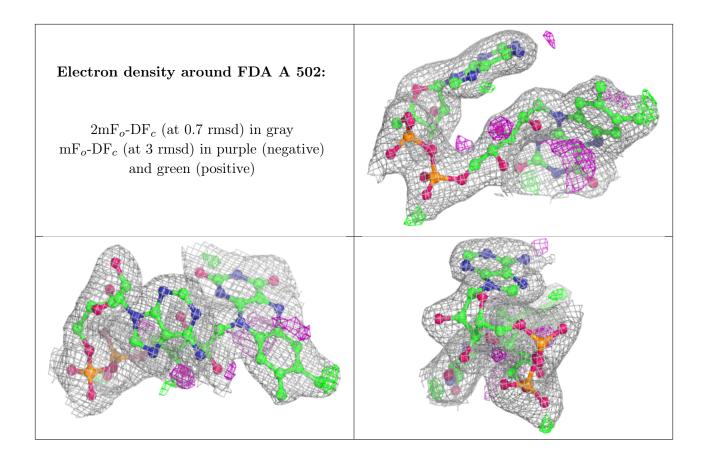
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median,  $95^{th}$  percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
5	FDA	В	501	53/53	0.95	0.17	36,44,53,57	0
5	FDA	A	502	53/53	0.96	0.16	23,31,38,42	0
4	SO4	A	501	5/5	0.99	0.10	36,37,46,46	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.

