



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 23, 2024 – 10:19 AM EDT

PDB ID : 1CI9
Title : DFP-INHIBITED ESTERASE ESTB FROM BURKHOLDERIA GLADIOLI
Authors : Wagner, U.G.; Petersen, E.I.; Schwab, H.; Kratky, C.
Deposited on : 1999-04-08
Resolution : 1.80 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.1

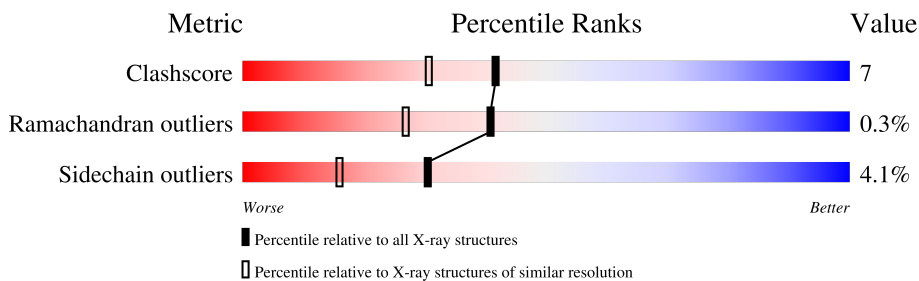
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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))
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	392	
1	B	392	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	DFP	B	401	-	-	X	-

2 Entry composition [i](#)

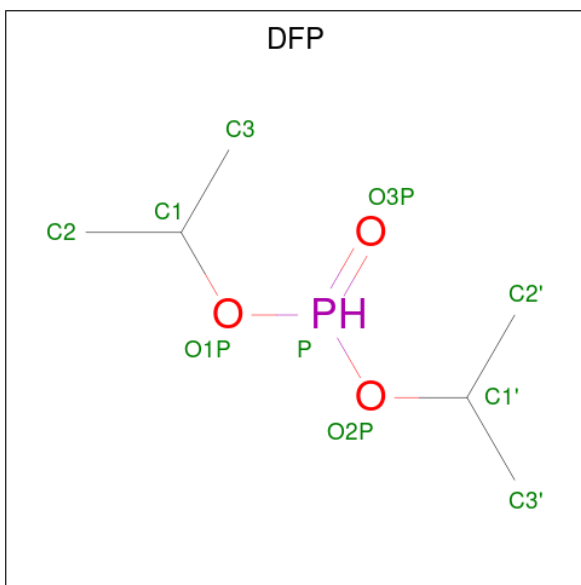
There are 3 unique types of molecules in this entry. The entry contains 6261 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (CARBOXYLESTERASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	377	Total 2835	C 1793	N 514	O 522	S 6	0	0	0
1	B	377	Total 2835	C 1793	N 514	O 522	S 6	0	0	0

- Molecule 2 is DIISOPROPYL PHOSPHONATE (three-letter code: DFP) (formula: C₆H₁₅O₃P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
2	A	1	Total 10	C 6	O 3	P 1	0	0
2	B	1	Total 10	C 6	O 3	P 1	0	0

- Molecule 3 is water.

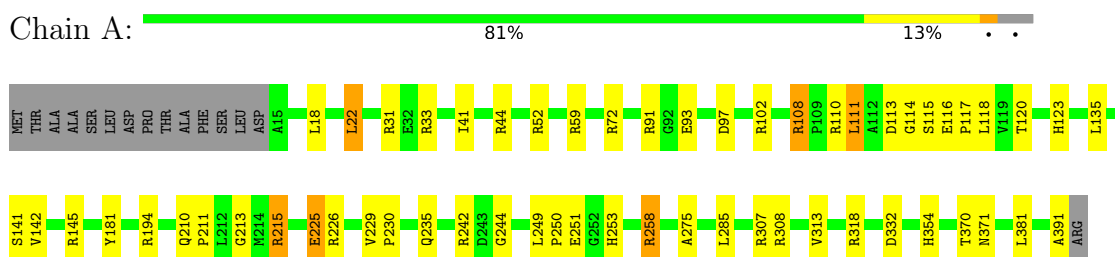
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	306	Total 306	O 306	0	0
3	B	265	Total 265	O 265	0	0

3 Residue-property plots [i](#)

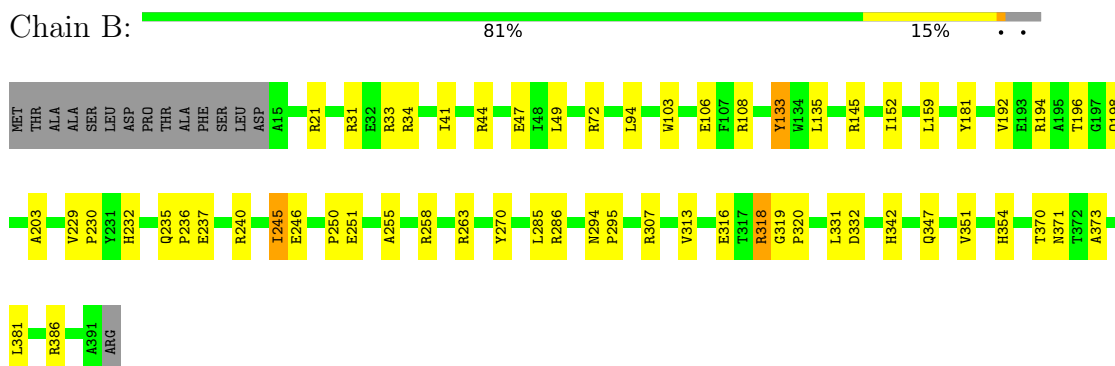
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

- Molecule 1: PROTEIN (CARBOXYLESTERASE)



- Molecule 1: PROTEIN (CARBOXYLESTERASE)



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	83.44Å 83.44Å 194.58Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 1.80	Depositor
% Data completeness (in resolution range)	95.0 (15.00-1.80)	Depositor
R_{merge}	0.02	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.176 , 0.235	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6261	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

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: DFP

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	0.50	1/2902 (0.0%)	1.08	11/3956 (0.3%)
1	B	0.37	0/2902	1.01	9/3956 (0.2%)
All	All	0.44	1/5804 (0.0%)	1.05	20/7912 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	391	ALA	C-O	17.46	1.56	1.23

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	391	ALA	CA-C-O	-14.35	89.96	120.10
1	A	258	ARG	NE-CZ-NH2	-9.61	115.50	120.30
1	B	386	ARG	NE-CZ-NH1	9.02	124.81	120.30
1	A	258	ARG	NE-CZ-NH1	8.56	124.58	120.30
1	A	59	ARG	NE-CZ-NH1	7.42	124.01	120.30
1	A	307	ARG	NE-CZ-NH1	7.19	123.89	120.30
1	B	270	TYR	CB-CG-CD1	6.94	125.17	121.00
1	A	308	ARG	NE-CZ-NH2	6.48	123.54	120.30
1	A	242	ARG	NE-CZ-NH1	-6.01	117.30	120.30
1	B	286	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	B	34	ARG	CD-NE-CZ	5.86	131.81	123.60
1	A	108	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	B	263	ARG	NE-CZ-NH2	5.82	123.21	120.30
1	B	286	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	A	226	ARG	NE-CZ-NH2	-5.65	117.48	120.30
1	A	31	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	B	108	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	A	226	ARG	NE-CZ-NH1	5.37	122.99	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	307	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	B	307	ARG	NE-CZ-NH1	5.18	122.89	120.30

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	2835	0	2798	37	0
1	B	2835	0	2798	42	0
2	A	10	0	14	4	0
2	B	10	0	14	6	0
3	A	306	0	0	4	0
3	B	265	0	0	5	0
All	All	6261	0	5624	79	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.

All (79) 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:B:196:THR:HG23	1:B:198:GLN:H	1.36	0.88
1:A:244:GLY:O	1:A:258:ARG:HD2	1.83	0.79
1:A:141:SER:O	1:A:145:ARG:HG3	1.85	0.76
1:B:313:VAL:HG12	3:B:1394:HOH:O	1.87	0.73
1:B:235:GLN:NE2	1:B:236:PRO:HA	2.03	0.73
1:A:142:VAL:HG22	1:A:145:ARG:NH2	2.06	0.71
1:B:196:THR:HG23	1:B:198:GLN:N	2.06	0.70
1:B:354:HIS:HD2	1:B:370:THR:H	1.40	0.68
1:A:108:ARG:HD3	1:A:116:GLU:OE1	1.94	0.68
1:A:110:ARG:HD3	1:A:114:GLY:O	1.93	0.67
1:B:313:VAL:HG22	1:B:332:ASP:OD2	1.96	0.66
1:A:91:ARG:HB3	1:A:93:GLU:OE1	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:LEU:HD22	2:A:401:DFP:H21	1.79	0.62
1:B:47:GLU:OE2	1:B:49:LEU:HD21	2.00	0.62
1:B:237:GLU:HA	3:B:1423:HOH:O	2.00	0.61
1:A:250:PRO:HG2	1:A:253:HIS:CG	2.37	0.59
1:B:33:ARG:NH2	1:B:236:PRO:HB2	2.18	0.58
1:A:72:ARG:O	1:A:354:HIS:HE1	1.87	0.58
1:B:354:HIS:CD2	1:B:370:THR:H	2.22	0.57
1:A:225:GLU:HB2	3:A:1152:HOH:O	2.04	0.57
1:A:313:VAL:HG12	1:A:332:ASP:OD2	2.05	0.56
1:B:245:ILE:HD13	1:B:246:GLU:N	2.21	0.56
1:B:351:VAL:HG13	2:B:401:DFP:H1'	1.88	0.56
1:A:354:HIS:HD2	1:A:370:THR:H	1.52	0.56
1:B:192:VAL:O	1:B:196:THR:HG22	2.05	0.55
1:A:318:ARG:HB3	1:A:318:ARG:NH1	2.21	0.55
1:B:181:TYR:HE2	2:B:401:DFP:H3'2	1.72	0.55
1:A:142:VAL:HG22	1:A:145:ARG:HH22	1.71	0.55
1:B:196:THR:HG21	1:B:203:ALA:HB2	1.90	0.54
1:B:318:ARG:NH1	1:B:347:GLN:HG2	2.22	0.53
1:B:235:GLN:HE21	1:B:236:PRO:HA	1.71	0.53
1:B:232:HIS:HA	1:B:373:ALA:O	2.08	0.53
1:B:181:TYR:CE2	2:B:401:DFP:H3'2	2.44	0.53
1:A:41:ILE:HG21	1:A:285:LEU:HD22	1.93	0.51
1:B:72:ARG:O	1:B:354:HIS:HE1	1.94	0.51
1:A:250:PRO:O	1:A:253:HIS:HB2	2.12	0.50
1:B:135:LEU:CD2	2:B:401:DFP:H33	2.42	0.50
1:B:235:GLN:CG	1:B:236:PRO:HA	2.41	0.50
1:A:313:VAL:HG12	1:A:332:ASP:CB	2.41	0.50
1:A:44:ARG:HD2	3:A:1382:HOH:O	2.12	0.50
1:A:213:GLY:HA2	3:A:1342:HOH:O	2.13	0.49
1:A:215:ARG:HH11	1:A:215:ARG:HB3	1.77	0.49
1:A:194:ARG:HG2	1:A:194:ARG:HH11	1.77	0.49
1:B:152:ILE:HG13	2:B:401:DFP:H32	1.94	0.49
1:B:316:GLU:HG2	1:B:320:PRO:HA	1.94	0.49
1:A:275:ALA:HB2	2:A:401:DFP:C3	2.43	0.48
1:B:133:TYR:HB3	1:B:135:LEU:HG	1.95	0.48
1:A:120:THR:H	1:A:123:HIS:HD2	1.62	0.48
1:B:196:THR:OG1	1:B:198:GLN:OE1	2.30	0.47
1:B:94:LEU:HD23	1:B:103:TRP:CZ2	2.51	0.46
1:B:235:GLN:CD	1:B:236:PRO:HA	2.36	0.46
1:A:111:LEU:HD11	1:A:117:PRO:HB3	1.97	0.46
1:A:230:PRO:HD2	3:A:1103:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:230:PRO:HD2	3:B:1053:HOH:O	2.15	0.45
1:B:44:ARG:NH1	3:B:1460:HOH:O	2.49	0.45
1:B:294:ASN:N	1:B:295:PRO:HD3	2.31	0.45
1:A:120:THR:H	1:A:123:HIS:CD2	2.34	0.45
1:B:21:ARG:NH1	3:B:1395:HOH:O	2.49	0.45
1:B:229:VAL:O	1:B:371:ASN:HB2	2.17	0.44
1:B:316:GLU:HA	1:B:319:GLY:O	2.17	0.44
1:A:250:PRO:HG2	1:A:253:HIS:HB2	2.00	0.44
1:B:246:GLU:OE1	1:B:258:ARG:NH2	2.51	0.44
1:A:251:GLU:OE1	1:A:251:GLU:N	2.49	0.43
1:B:240:ARG:HH11	1:B:240:ARG:HD2	1.64	0.43
1:A:181:TYR:CE2	2:A:401:DFP:H1'	2.54	0.43
1:A:181:TYR:OH	2:A:401:DFP:H1'	2.19	0.42
1:B:246:GLU:OE1	1:B:258:ARG:NE	2.50	0.42
1:B:235:GLN:HG3	1:B:236:PRO:HA	2.01	0.42
1:A:250:PRO:HG2	1:A:253:HIS:CB	2.49	0.42
1:A:210:GLN:HB2	1:A:211:PRO:HD3	2.02	0.42
1:B:44:ARG:O	1:B:47:GLU:HG2	2.20	0.42
1:A:229:VAL:O	1:A:371:ASN:HB2	2.21	0.41
1:A:111:LEU:HD12	1:A:115:SER:OG	2.21	0.41
1:A:194:ARG:HG2	1:A:194:ARG:NH1	2.36	0.41
1:A:249:LEU:CD2	1:A:250:PRO:HD2	2.51	0.41
1:B:351:VAL:HG21	2:B:401:DFP:H31	2.03	0.41
1:B:145:ARG:HG3	1:B:145:ARG:HH11	1.86	0.40
1:A:18:LEU:HG	1:A:22:LEU:HD22	2.03	0.40
1:B:41:ILE:HG21	1:B:285:LEU:HD22	2.03	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	375/392 (96%)	364 (97%)	11 (3%)	0	100	100
1	B	375/392 (96%)	361 (96%)	12 (3%)	2 (0%)	29	15
All	All	750/784 (96%)	725 (97%)	23 (3%)	2 (0%)	41	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	255	ALA
1	B	250	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	278/290 (96%)	266 (96%)	12 (4%)	29	14
1	B	278/290 (96%)	267 (96%)	11 (4%)	31	16
All	All	556/580 (96%)	533 (96%)	23 (4%)	30	16

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	33	ARG
1	A	52	ARG
1	A	97	ASP
1	A	102	ARG
1	A	111	LEU
1	A	113	ASP
1	A	118	LEU
1	A	215	ARG
1	A	225	GLU
1	A	235	GLN
1	A	381	LEU
1	B	31	ARG
1	B	106	GLU

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Mol	Chain	Res	Type
1	B	133	TYR
1	B	159	LEU
1	B	194	ARG
1	B	245	ILE
1	B	251	GLU
1	B	318	ARG
1	B	331	LEU
1	B	342	HIS
1	B	381	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	123	HIS
1	A	354	HIS
1	B	28	GLN
1	B	235	GLN
1	B	341	GLN
1	B	354	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](#)

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

5.5 Carbohydrates [i](#)

There are no monosaccharides 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
2	DFP	B	401	1	6,9,9	1.27	1 (16%)	6,11,11	0.66	0
2	DFP	A	401	1	6,9,9	1.44	2 (33%)	6,11,11	0.69	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DFP	B	401	1	-	1/4/8/8	-
2	DFP	A	401	1	-	2/4/8/8	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	DFP	O1P-C1	-2.12	1.43	1.45
2	B	401	DFP	O2P-C1'	-2.06	1.43	1.45
2	A	401	DFP	O2P-C1'	-2.05	1.43	1.45

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	DFP	C3-C1-O1P-P
2	B	401	DFP	C2'-C1'-O2P-P
2	A	401	DFP	C2-C1-O1P-P

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	DFP	6	0
2	A	401	DFP	4	0

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](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.