



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 06:16 AM BST

PDB ID : 1PFK  
Title : CRYSTAL STRUCTURE OF THE COMPLEX OF PHOSPHOFRUCTOKINASE FROM ESCHERICHIA COLI WITH ITS REACTION PRODUCTS  
Authors : Shirakihara, Y.; Evans, P.R.  
Deposited on : 1988-01-25  
Resolution : 2.40 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](mailto: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.

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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)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
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) : 2.13.1

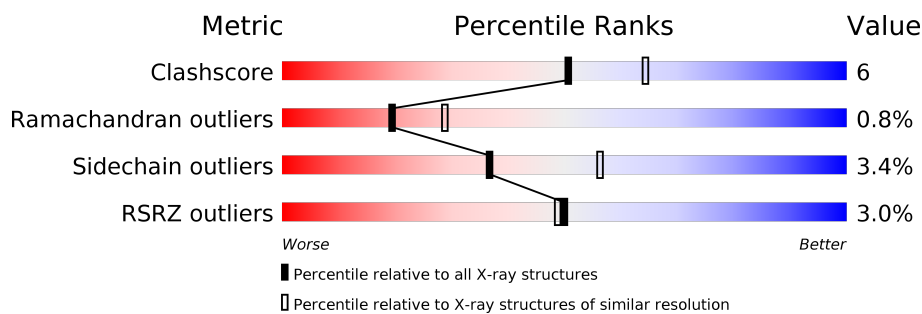
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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  $\geq 3$ , 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	320	<div> <div>4%</div> <div>78%</div> <div>20%</div> <div>..</div> </div>
1	B	320	<div> <div>2%</div> <div>80%</div> <div>18%</div> <div>..</div> </div>

## 2 Entry composition [i](#)

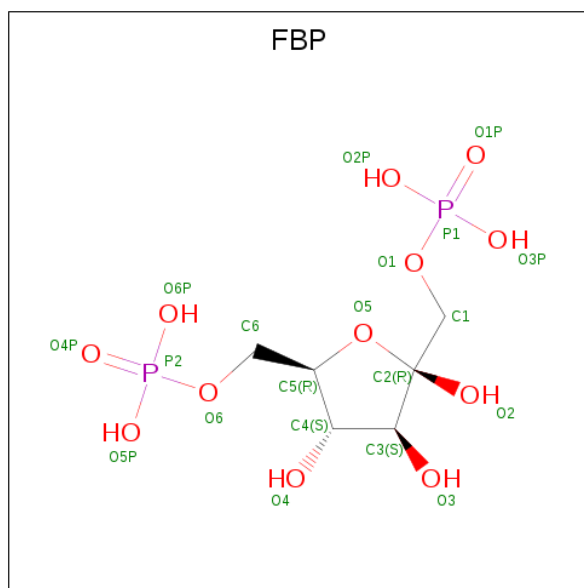
There are 5 unique types of molecules in this entry. The entry contains 5295 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PHOSPHOFRUCTOKINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	0
			2433	1527	426	462	18			
1	B	320	Total	C	N	O	S	0	0	0
			2433	1527	426	462	18			

- Molecule 2 is 1,6-di-O-phosphono-beta-D-fructofuranose (three-letter code: FBP) (formula:  $C_6H_{14}O_{12}P_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			20	6	12	2		
2	B	1	Total	C	O	P	0	0
			20	6	12	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	2	Total Mg 2 2	0	0
3	A	2	Total Mg 2 2	0	0

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O P 27 10 5 10 2	0	0
4	A	1	Total C N O P 27 10 5 10 2	0	0
4	B	1	Total C N O P 27 10 5 10 2	0	0
4	B	1	Total C N O P 27 10 5 10 2	0	0

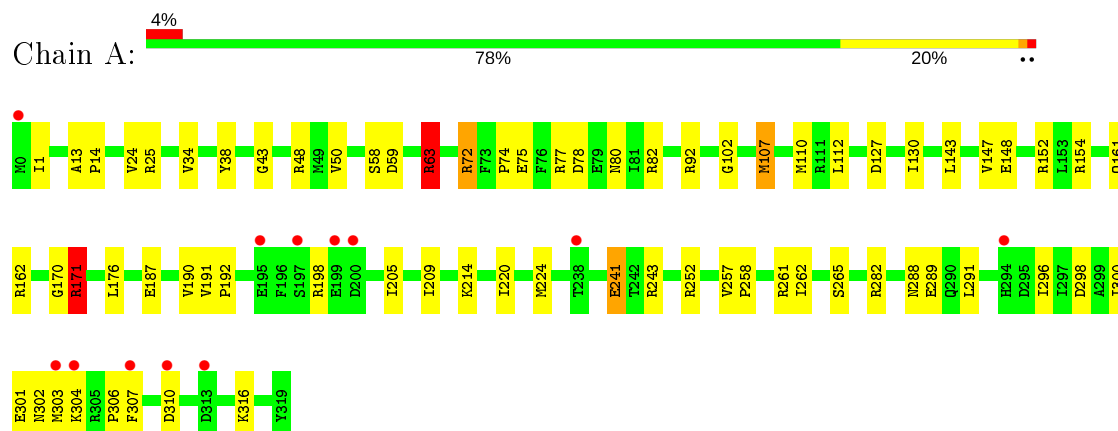
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	141	Total O 141 141	0	0
5	B	136	Total O 136 136	0	0

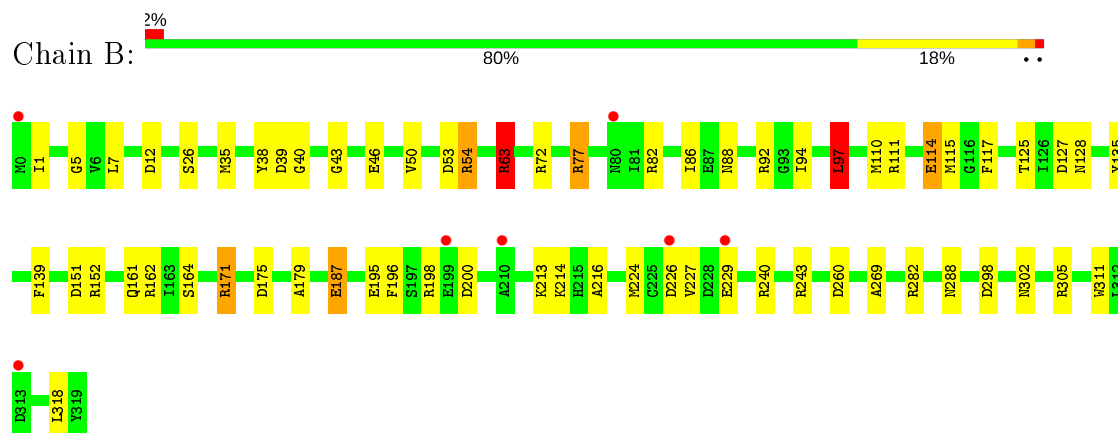
### 3 Residue-property plots [i](#)

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: PHOSPHOFRUCTOKINASE



#### • Molecule 1: PHOSPHOFRUCTOKINASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	112.30 Å 85.40 Å 77.10 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 2.40 22.44 – 2.43	Depositor EDS
% Data completeness (in resolution range)	(Not available) (100.00-2.40) 94.5 (22.44-2.43)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.56 (at 2.44 Å)	Xtriage
Refinement program	PROLSQ	Depositor
R, $R_{free}$	0.165 , (Not available) 0.169 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.9	Xtriage
Anisotropy	0.232	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 56.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5295	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MG, FBP, ADP

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	0.73	0/2470	1.55	32/3329 (1.0%)
1	B	0.75	0/2470	1.49	32/3329 (1.0%)
All	All	0.74	0/4940	1.52	64/6658 (1.0%)

There are no bond length outliers.

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	63	ARG	NE-CZ-NH1	13.04	126.82	120.30
1	A	59	ASP	CB-CG-OD1	11.66	128.79	118.30
1	A	152	ARG	NE-CZ-NH1	11.55	126.07	120.30
1	A	198	ARG	NE-CZ-NH2	-11.50	114.55	120.30
1	A	282	ARG	NE-CZ-NH1	11.16	125.88	120.30
1	A	198	ARG	NE-CZ-NH1	10.05	125.32	120.30
1	B	54	ARG	NE-CZ-NH2	-9.99	115.31	120.30
1	B	54	ARG	NE-CZ-NH1	9.47	125.03	120.30
1	B	243	ARG	NE-CZ-NH1	9.23	124.91	120.30
1	A	77	ARG	NE-CZ-NH1	8.79	124.70	120.30
1	A	63	ARG	CD-NE-CZ	8.78	135.90	123.60
1	A	72	ARG	CD-NE-CZ	8.68	135.75	123.60
1	A	48	ARG	NE-CZ-NH2	-8.59	116.01	120.30
1	A	92	ARG	NE-CZ-NH2	-8.34	116.13	120.30
1	A	59	ASP	CB-CG-OD2	-8.00	111.10	118.30
1	B	162	ARG	NE-CZ-NH1	8.00	124.30	120.30
1	B	77	ARG	NE-CZ-NH2	-7.62	116.49	120.30
1	A	82	ARG	NE-CZ-NH1	7.35	123.97	120.30
1	A	82	ARG	NE-CZ-NH2	-7.32	116.64	120.30
1	A	63	ARG	NE-CZ-NH2	-7.21	116.69	120.30
1	A	77	ARG	NE-CZ-NH2	-7.07	116.77	120.30
1	B	171	ARG	NE-CZ-NH1	6.91	123.76	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	243	ARG	NE-CZ-NH1	6.71	123.66	120.30
1	B	72	ARG	NE-CZ-NH2	-6.69	116.95	120.30
1	A	72	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	A	154	ARG	NE-CZ-NH2	-6.51	117.05	120.30
1	A	252	ARG	NE-CZ-NH1	6.50	123.55	120.30
1	B	162	ARG	NE-CZ-NH2	-6.49	117.06	120.30
1	A	127	ASP	CB-CG-OD1	6.43	124.09	118.30
1	B	12	ASP	CB-CG-OD1	6.42	124.08	118.30
1	B	311	TRP	CA-CB-CG	6.39	125.84	113.70
1	A	78	ASP	CB-CG-OD1	6.32	123.99	118.30
1	A	261	ARG	NE-CZ-NH1	6.31	123.45	120.30
1	B	260	ASP	CB-CG-OD2	-6.21	112.71	118.30
1	A	243	ARG	NE-CZ-NH2	-6.19	117.20	120.30
1	B	53	ASP	CB-CG-OD1	6.16	123.84	118.30
1	A	241	GLU	CA-CB-CG	6.11	126.85	113.40
1	B	63	ARG	NE-CZ-NH2	-6.10	117.25	120.30
1	B	63	ARG	NE-CZ-NH1	5.97	123.28	120.30
1	B	72	ARG	NE-CZ-NH1	5.96	123.28	120.30
1	B	305	ARG	NE-CZ-NH1	5.94	123.27	120.30
1	B	46	GLU	CG-CD-OE1	5.91	130.11	118.30
1	B	282	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	A	152	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	B	12	ASP	CB-CG-OD2	-5.84	113.05	118.30
1	B	97	LEU	CA-CB-CG	5.83	128.70	115.30
1	B	152	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	B	151	ASP	CB-CG-OD1	5.82	123.54	118.30
1	B	171	ARG	CA-CB-CG	5.75	126.05	113.40
1	A	224	MET	CA-CB-CG	-5.67	103.66	113.30
1	A	282	ARG	NE-CZ-NH2	-5.67	117.47	120.30
1	B	39	ASP	CB-CG-OD1	5.57	123.31	118.30
1	B	179	ALA	CB-CA-C	5.53	118.40	110.10
1	A	152	ARG	CD-NE-CZ	5.38	131.13	123.60
1	A	171	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	B	175	ASP	CB-CG-OD1	5.35	123.11	118.30
1	A	298	ASP	CB-CG-OD1	5.33	123.10	118.30
1	A	25	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	B	243	ARG	CD-NE-CZ	5.21	130.90	123.60
1	B	243	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	B	224	MET	CA-CB-CG	-5.20	104.45	113.30
1	B	240	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	B	127	ASP	CB-CG-OD1	5.11	122.90	118.30
1	B	198	ARG	NE-CZ-NH1	5.02	122.81	120.30



There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

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	2433	0	2430	31	0
1	B	2433	0	2430	28	0
2	A	20	0	10	0	0
2	B	20	0	10	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	54	0	24	0	0
4	B	54	0	24	2	0
5	A	141	0	0	3	0
5	B	136	0	0	5	0
All	All	5295	0	4928	57	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 6.

All (57) 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:24:VAL:HG22	1:A:34:VAL:HG11	1.63	0.81
1:A:316:LYS:NZ	5:A:427:HOH:O	2.12	0.81
1:A:161:GLN:HA	1:A:214:LYS:O	1.90	0.71
1:A:162:ARG:HG2	1:A:241:GLU:HG3	1.77	0.66
1:A:102:GLY:HA2	1:A:130:ILE:HD11	1.80	0.63
1:A:107:MET:HE3	1:A:300:ILE:HD13	1.82	0.61
1:B:213:LYS:HE2	4:B:326:ADP:O1A	2.01	0.60
1:B:161:GLN:HA	1:B:214:LYS:O	2.02	0.59
1:B:88:ASN:O	1:B:92:ARG:HG2	2.03	0.59
1:B:187:GLU:HG2	1:B:216:ALA:HA	1.83	0.58
1:A:303:MET:O	1:A:304:LYS:HG3	2.05	0.56
1:B:128:ASN:OD1	1:B:135:TYR:HA	2.05	0.55
1:A:191:VAL:HB	1:A:192:PRO:HD2	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:63:ARG:HH11	1:B:63:ARG:HG2	1.73	0.52
1:A:306:PRO:HG2	5:A:464:HOH:O	2.12	0.50
1:B:38:TYR:O	1:B:43:GLY:HA3	2.11	0.50
1:B:82:ARG:O	1:B:86:ILE:HG13	2.12	0.50
1:A:75:GLU:OE2	1:A:75:GLU:N	2.45	0.50
1:A:205:ILE:O	1:A:209:ILE:HG13	2.12	0.50
1:A:63:ARG:HG2	1:A:63:ARG:HH11	1.76	0.49
1:A:110:MET:HB2	1:A:296:ILE:HG21	1.93	0.49
1:B:54:ARG:HD2	5:B:454:HOH:O	2.12	0.49
1:A:38:TYR:O	1:A:43:GLY:HA3	2.13	0.48
1:B:35:MET:HG3	1:B:94:ILE:HD11	1.95	0.48
1:B:82:ARG:HE	1:B:115:MET:HE1	1.79	0.48
1:B:82:ARG:HE	1:B:115:MET:CE	2.26	0.48
1:B:110:MET:O	1:B:114:GLU:HG2	2.14	0.48
1:B:196:PHE:HA	5:B:408:HOH:O	2.14	0.47
1:A:190:VAL:HB	1:A:220:ILE:HG13	1.97	0.47
1:A:110:MET:HE2	1:A:110:MET:HB3	1.84	0.46
1:A:13:ALA:HA	1:A:14:PRO:HD3	1.74	0.46
1:B:226:ASP:OD1	1:B:229:GLU:HG3	2.16	0.45
1:A:258:PRO:O	1:A:262:ILE:HG13	2.16	0.45
1:A:289:GLU:HB2	1:B:288:ASN:HD22	1.82	0.45
1:B:298:ASP:O	1:B:302:ASN:HB2	2.16	0.45
1:A:107:MET:HA	1:A:107:MET:HE3	1.99	0.44
1:B:82:ARG:NH2	1:B:111:ARG:HD2	2.33	0.44
1:A:265:SER:HB3	1:B:318:LEU:CD1	2.48	0.43
1:A:72:ARG:O	1:A:74:PRO:HD3	2.18	0.43
1:B:86:ILE:HG23	1:B:117:PHE:CE1	2.53	0.43
1:A:58:SER:HB3	4:B:326:ADP:H2'	2.01	0.43
1:B:63:ARG:NH1	1:B:63:ARG:HG2	2.34	0.43
1:A:257:VAL:HB	1:A:258:PRO:HD2	2.00	0.43
1:A:170:GLY:O	1:A:171:ARG:HB3	2.19	0.42
1:A:176:LEU:HA	1:A:176:LEU:HD12	1.82	0.42
1:A:38:TYR:HE2	1:A:50:VAL:HG13	1.84	0.42
1:B:7:LEU:HD13	1:B:40:GLY:HA2	2.02	0.42
1:B:26:SER:OG	1:B:269:ALA:HA	2.19	0.42
1:A:288:ASN:O	1:A:289:GLU:HB2	2.20	0.41
1:B:5:GLY:O	1:B:97:LEU:HA	2.19	0.41
1:B:77:ARG:NH1	5:B:449:HOH:O	2.54	0.41
1:B:128:ASN:HB2	1:B:139:PHE:CD2	2.56	0.41
1:A:143:LEU:O	1:A:147:VAL:HG23	2.19	0.41
1:A:148:GLU:HB2	5:A:393:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:291:LEU:HD23	1:A:291:LEU:HA	1.95	0.41
1:B:114:GLU:HG3	5:B:450:HOH:O	2.21	0.40
1:B:227:VAL:HG22	5:B:379:HOH:O	2.21	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	318/320 (99%)	306 (96%)	10 (3%)	2 (1%)	25	36
1	B	318/320 (99%)	310 (98%)	5 (2%)	3 (1%)	17	25
All	All	636/640 (99%)	616 (97%)	15 (2%)	5 (1%)	19	29

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	171	ARG
1	A	1	ILE
1	B	1	ILE
1	B	171	ARG
1	B	125	THR

### 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	253/255 (99%)	244 (96%)	9 (4%)	35	54
1	B	253/255 (99%)	245 (97%)	8 (3%)	39	59
All	All	506/510 (99%)	489 (97%)	17 (3%)	37	56

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ARG
1	A	80	ASN
1	A	107	MET
1	A	112	LEU
1	A	187	GLU
1	A	301	GLU
1	A	302	ASN
1	A	307	PHE
1	A	310	ASP
1	B	50	VAL
1	B	63	ARG
1	B	97	LEU
1	B	114	GLU
1	B	164	SER
1	B	187	GLU
1	B	195	GLU
1	B	200	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	288	ASN
1	B	160	HIS
1	B	288	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 for Mogul analysis.

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
4	ADP	A	326	3	24,29,29	0.80	0	29,45,45	0.83	1 (3%)
4	ADP	A	324	3	24,29,29	0.88	0	29,45,45	1.06	2 (6%)
2	FBP	B	323	-	18,20,20	1.02	1 (5%)	23,32,32	1.41	4 (17%)
4	ADP	B	326	3	24,29,29	0.79	1 (4%)	29,45,45	1.00	2 (6%)
2	FBP	A	323	3	18,20,20	0.99	0	23,32,32	0.70	0
4	ADP	B	324	3	24,29,29	0.84	0	29,45,45	1.14	2 (6%)

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
4	ADP	A	326	3	-	0/12/32/32	0/3/3/3
4	ADP	A	324	3	-	1/12/32/32	0/3/3/3
2	FBP	B	323	-	-	5/13/32/32	0/1/1/1
4	ADP	B	326	3	-	0/12/32/32	0/3/3/3
2	FBP	A	323	3	-	1/13/32/32	0/1/1/1
4	ADP	B	324	3	-	0/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	323	FBP	P2-O5P	-2.08	1.46	1.54
4	B	326	ADP	C2-N1	2.06	1.37	1.33

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	323	FBP	P1-O1-C1	3.49	127.90	118.30
4	B	324	ADP	C5-C6-N6	3.30	125.37	120.35
4	B	326	ADP	C5-C6-N6	3.25	125.30	120.35
4	A	324	ADP	PA-O3A-PB	2.73	142.21	132.83
4	B	324	ADP	PA-O3A-PB	2.62	141.80	132.83
2	B	323	FBP	O6P-P2-O6	2.61	113.69	106.73
2	B	323	FBP	O1-P1-O1P	2.60	113.77	106.47
4	A	324	ADP	C5-C6-N6	2.40	124.00	120.35
4	A	326	ADP	C4-C5-N7	2.21	111.71	109.40
2	B	323	FBP	O5P-P2-O6	2.17	112.51	106.73
4	B	326	ADP	O2'-C2'-C3'	2.08	118.54	111.82

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	323	FBP	C1-O1-P1-O1P
2	B	323	FBP	C1-O1-P1-O2P
2	B	323	FBP	C1-O1-P1-O3P
2	B	323	FBP	O1-C1-C2-O2
2	B	323	FBP	O1-C1-C2-O5
4	A	324	ADP	PB-O3A-PA-O5'
2	A	323	FBP	C4-C5-C6-O6

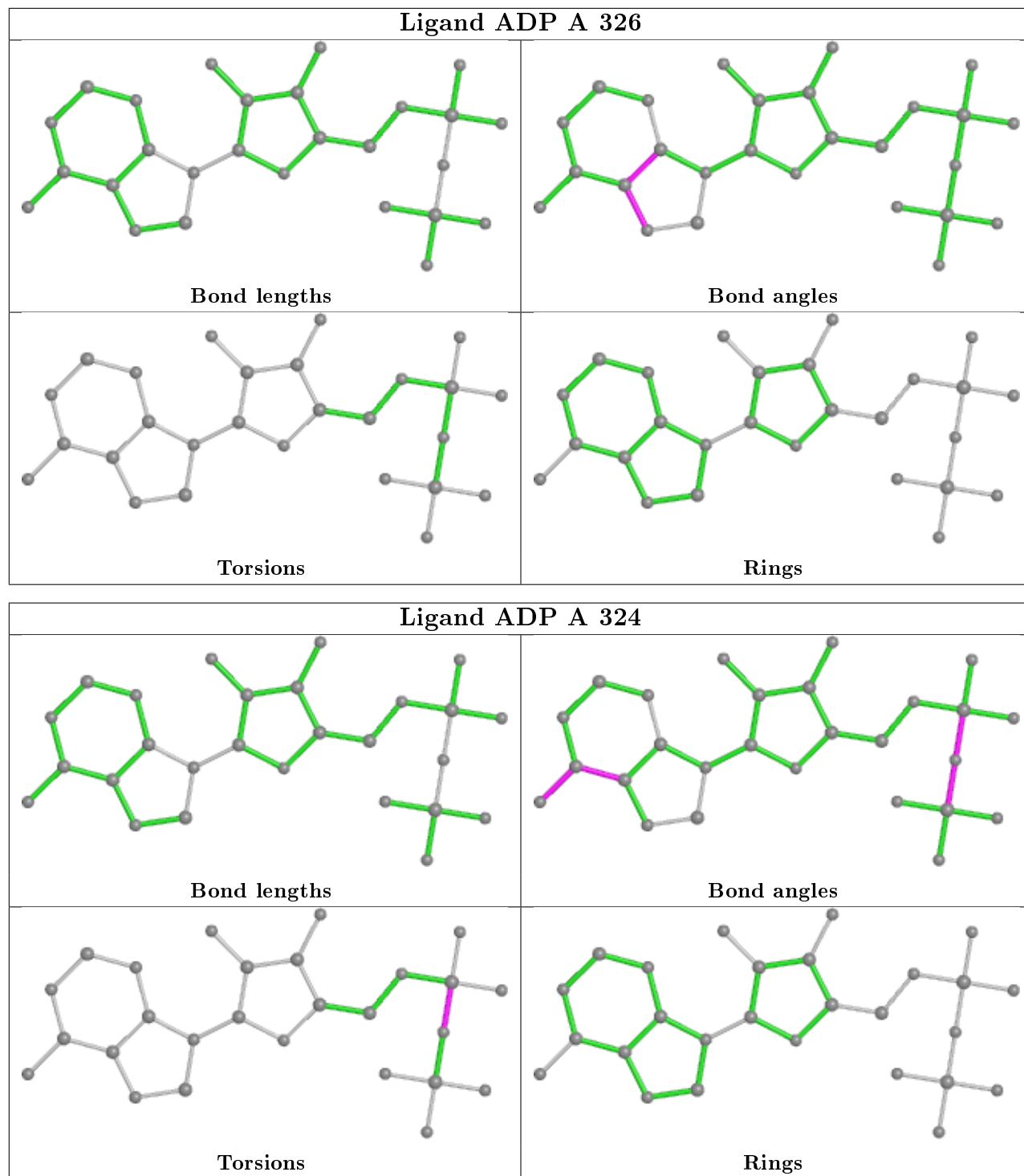
There are no ring outliers.

1 monomer is involved in 2 short contacts:

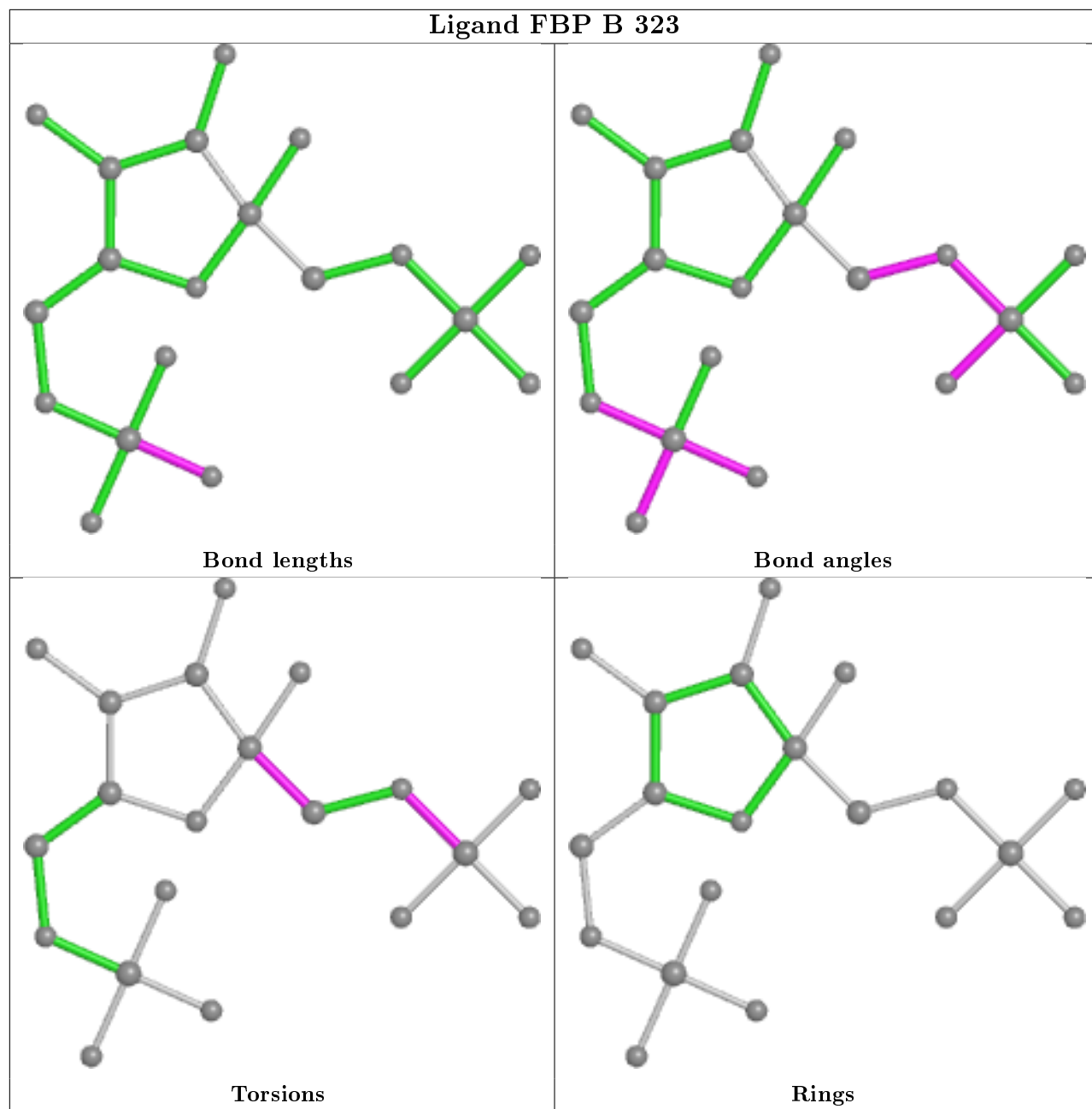
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	326	ADP	2	0

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 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

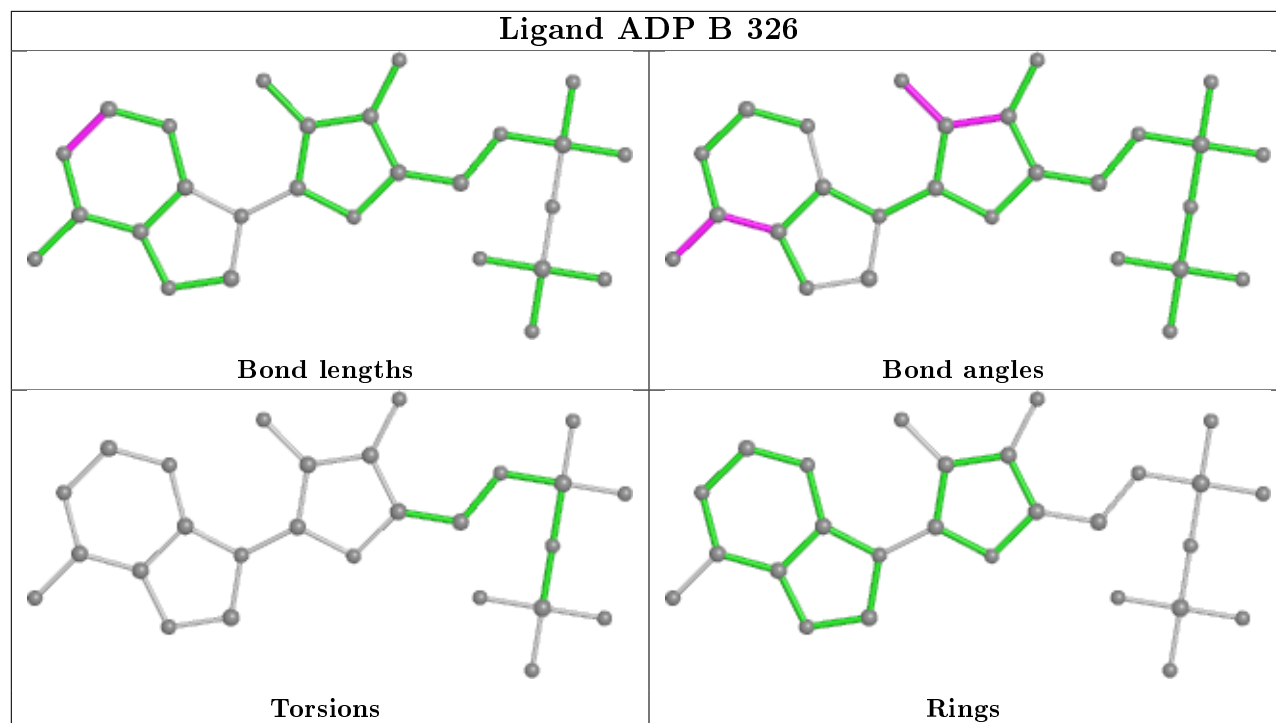
equivalents in the CSD to analyse the geometry.



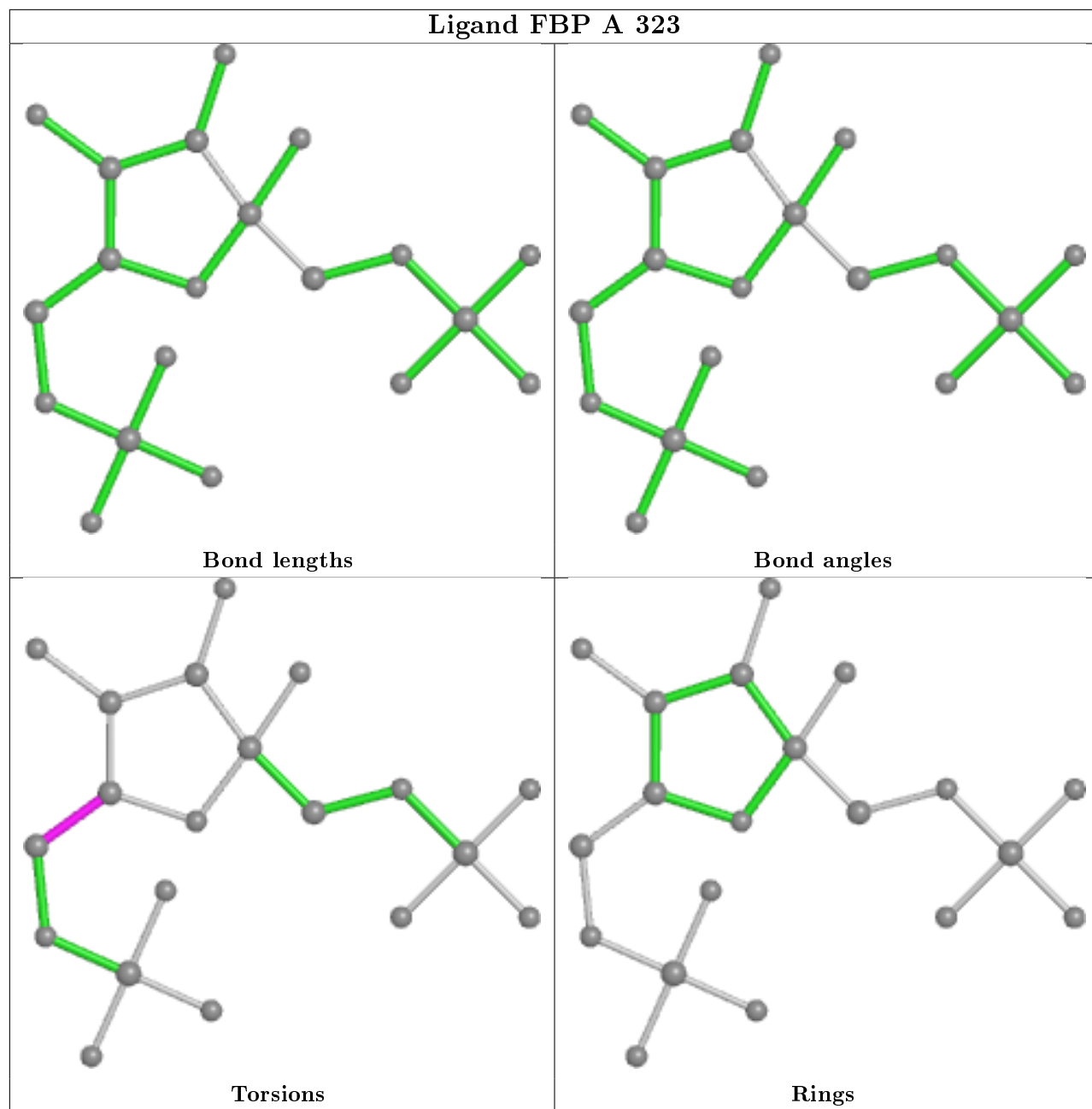
## Ligand FBP B 323

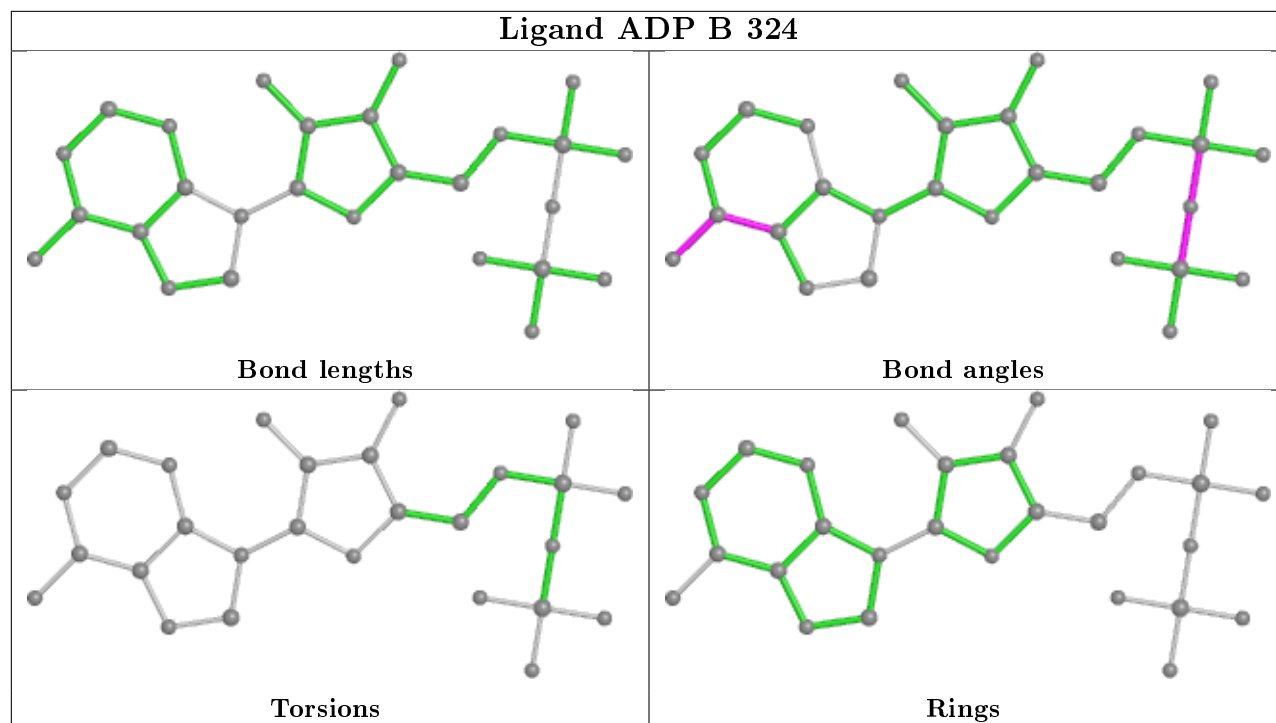






## Ligand FBP A 323





## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	320/320 (100%)	-0.10	12 (3%) 40 39	13, 34, 70, 99	0
1	B	320/320 (100%)	-0.21	7 (2%) 62 60	12, 36, 64, 90	0
All	All	640/640 (100%)	-0.15	19 (2%) 50 49	12, 35, 66, 99	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	0	MET	5.4
1	A	310	ASP	5.0
1	A	200	ASP	3.9
1	A	313	ASP	3.8
1	A	199	GLU	3.4
1	A	303	MET	3.2
1	B	0	MET	2.9
1	B	229	GLU	2.9
1	A	307	PHE	2.7
1	B	199	GLU	2.6
1	B	80	ASN	2.6
1	A	294	HIS	2.6
1	B	313	ASP	2.5
1	A	238	THR	2.4
1	A	195	GLU	2.4
1	A	304	LYS	2.2
1	A	197	SER	2.2
1	B	226	ASP	2.1
1	B	210	ALA	2.1

### 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 ⓘ

There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

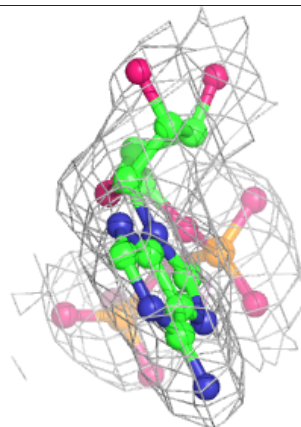
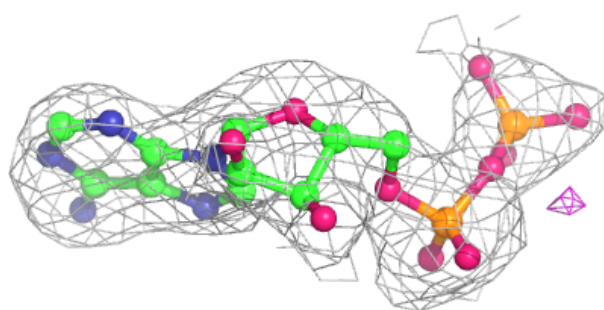
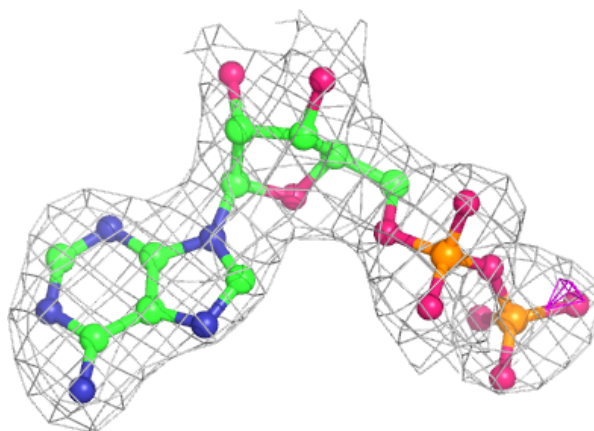
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	MG	B	325	1/1	0.83	0.17	64,64,64,64	0
4	ADP	B	324	27/27	0.92	0.13	49,51,58,58	0
3	MG	A	325	1/1	0.94	0.06	61,61,61,61	0
3	MG	B	327	1/1	0.95	0.05	14,14,14,14	0
4	ADP	A	324	27/27	0.96	0.11	41,46,48,49	0
2	FBP	B	323	20/20	0.96	0.13	28,37,51,51	4
4	ADP	A	326	27/27	0.97	0.11	31,48,51,53	0
4	ADP	B	326	27/27	0.97	0.10	28,35,47,47	0
2	FBP	A	323	20/20	0.98	0.10	27,31,45,45	4
3	MG	A	327	1/1	0.98	0.04	27,27,27,27	0

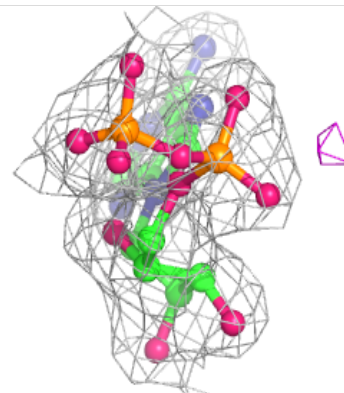
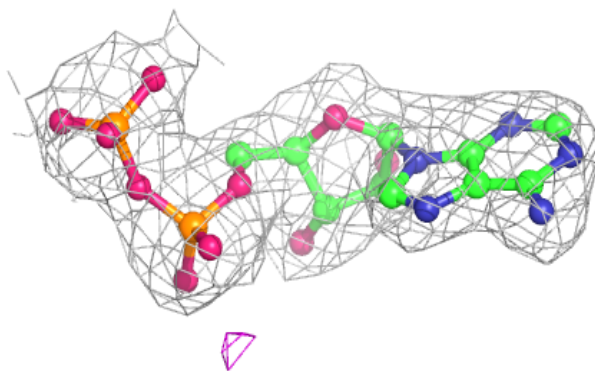
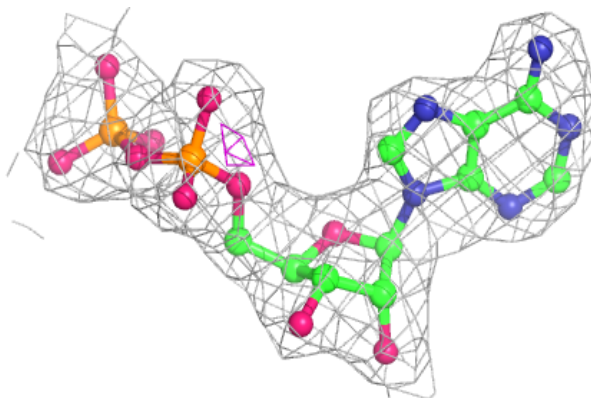
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around ADP B 324:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

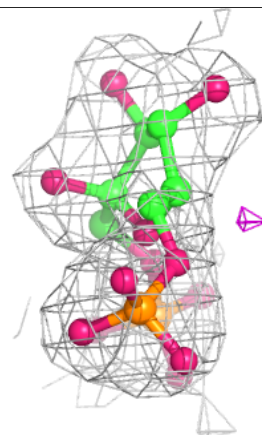
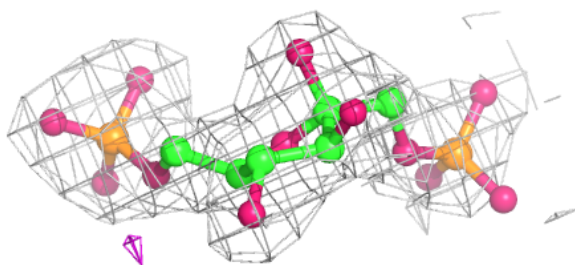
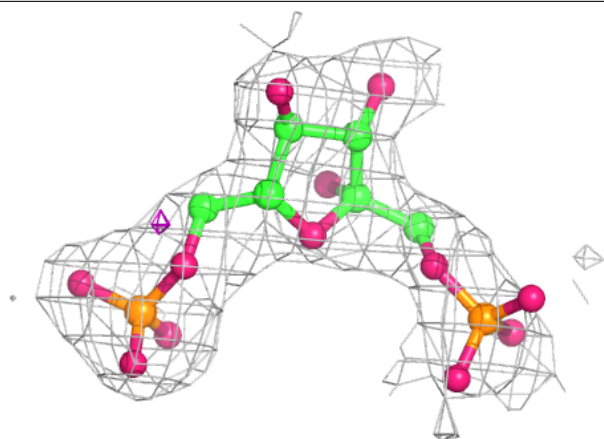
**Electron density around ADP A 324:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

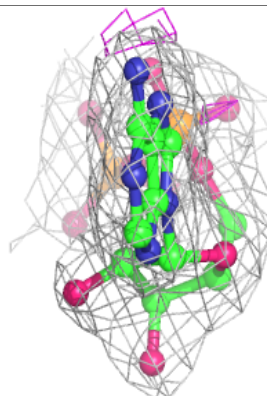
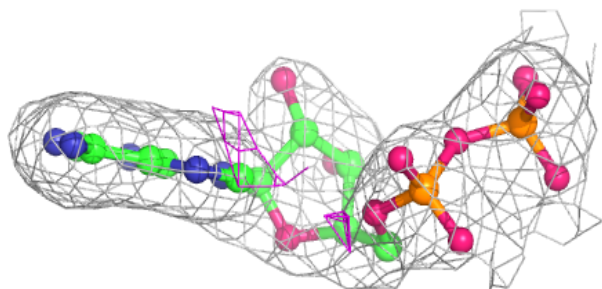
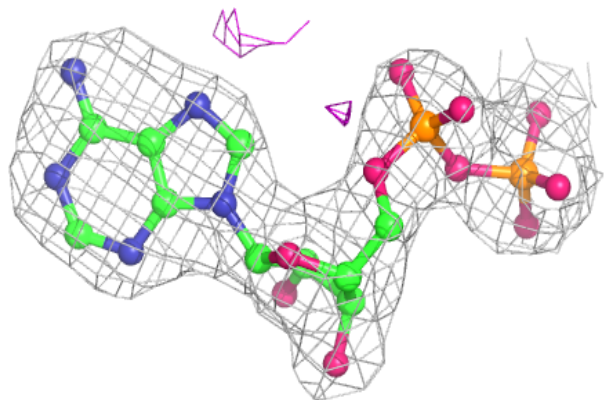


**Electron density around FBP B 323:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

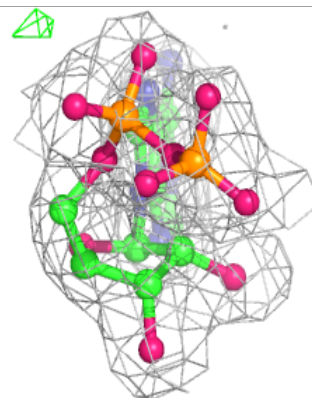
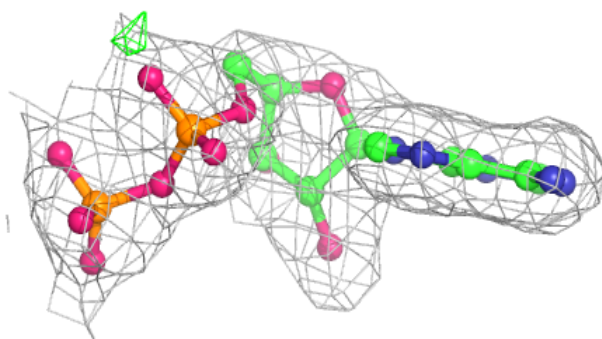
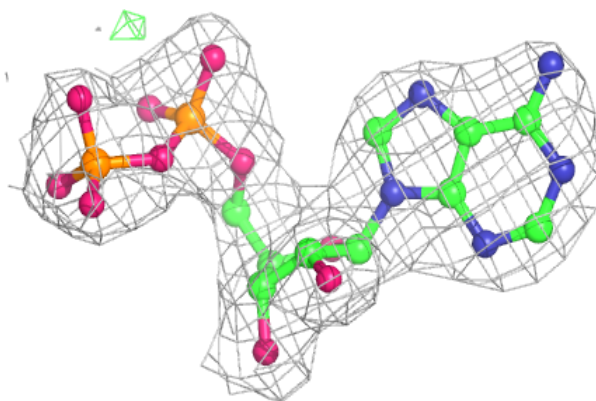
**Electron density around ADP A 326:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

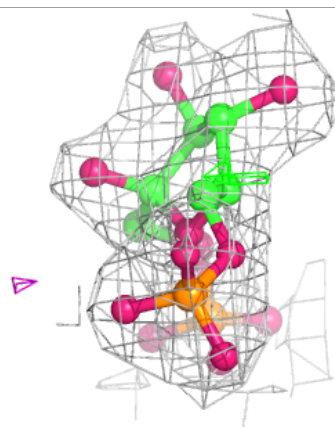
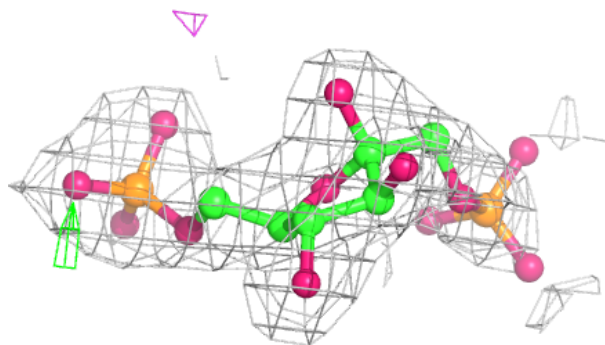
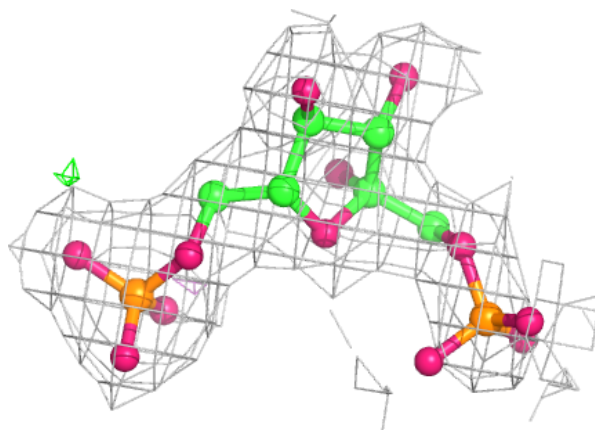


**Electron density around ADP B 326:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around FBP A 323:**

$2mF_o - DF_c$  (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.