



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 09:24 PM GMT

PDB ID : 1OSN
Title : Crystal structure of Varicella zoster virus thymidine kinase in complex with BVDU-MP and ADP
Authors : Bird, L.E.; Ren, J.; Wright, A.; Leslie, K.D.; Degreve, B.; Balzarini, J.; Stammers, D.K.
Deposited on : 2003-03-20
Resolution : 3.20 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

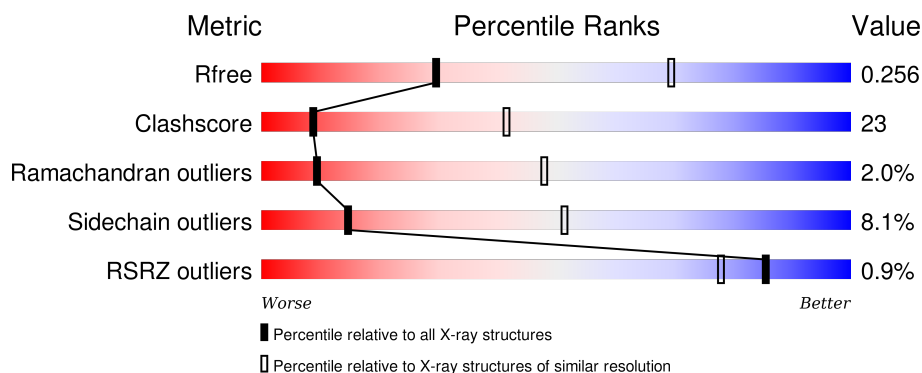
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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))
R_{free}	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	341	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 59% 31% 5% • 5% </div> </div>
1	B	341	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 57% 30% 5% 8% </div> </div>
1	C	341	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 55% 32% 5% 8% </div> </div>
1	D	341	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 57% 29% 6% 8% </div> </div>

2 Entry composition [i](#)

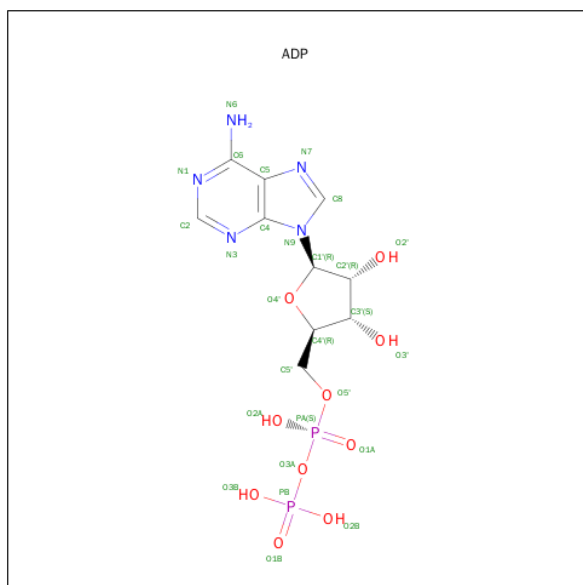
There are 4 unique types of molecules in this entry. The entry contains 10038 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Thymidine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	0	0
			2524	1623	426	456	19			
1	B	315	Total	C	N	O	S	0	0	0
			2446	1574	412	441	19			
1	C	313	Total	C	N	O	S	0	0	0
			2433	1568	411	435	19			
1	D	315	Total	C	N	O	S	0	0	0
			2452	1579	413	441	19			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



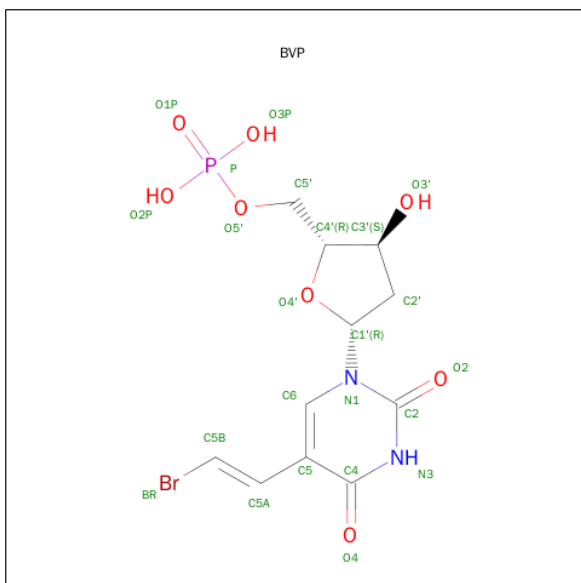
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	O	P		0	0
			8	6	2			
2	B	1	Total	C	N	O	P	0
			27	10	5	10	2	0

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

- Molecule 3 is (E)-5-(2-BROMOVINYL)-2'-DEOXYURIDINE-5'-MONOPHOSPHATE (three-letter code: BVP) (formula: $C_{11}H_{14}BrN_2O_8P$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 23	Br 1	C 11	N 2	O 8	P 1	0	0
3	B	1	Total 23	Br 1	C 11	N 2	O 8	P 1	0	0
3	C	1	Total 23	Br 1	C 11	N 2	O 8	P 1	0	0
3	D	1	Total 23	Br 1	C 11	N 2	O 8	P 1	0	0

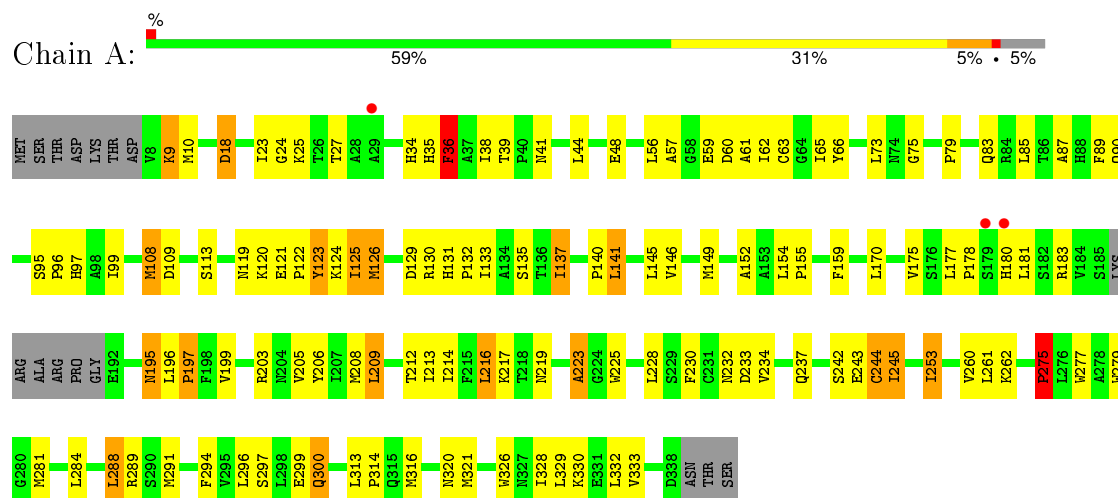
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total	0	0
			2		

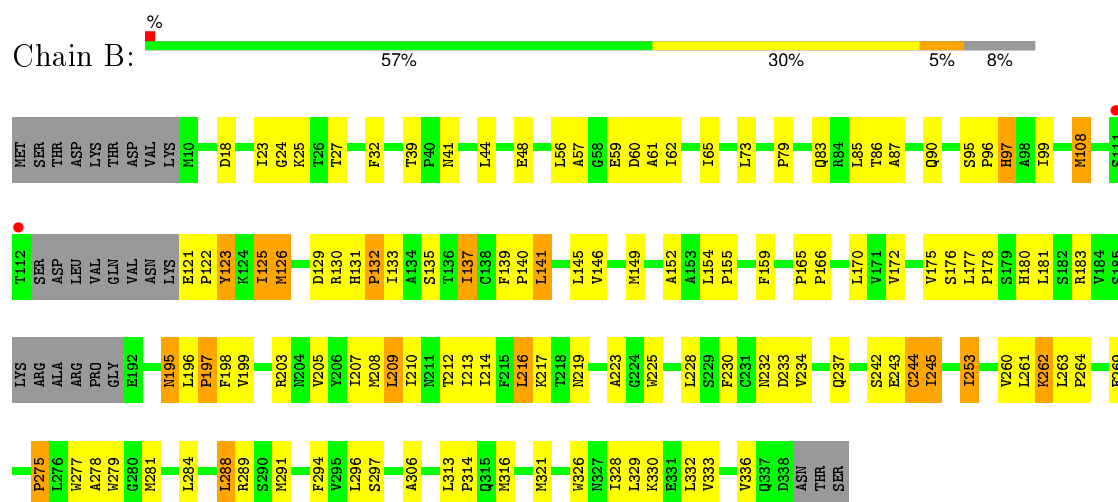
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Thymidine kinase

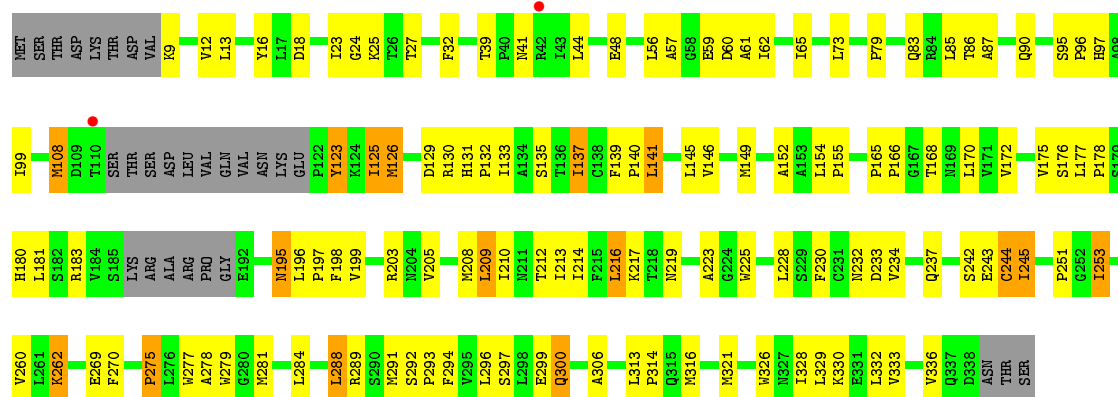


• Molecule 1: Thymidine kinase

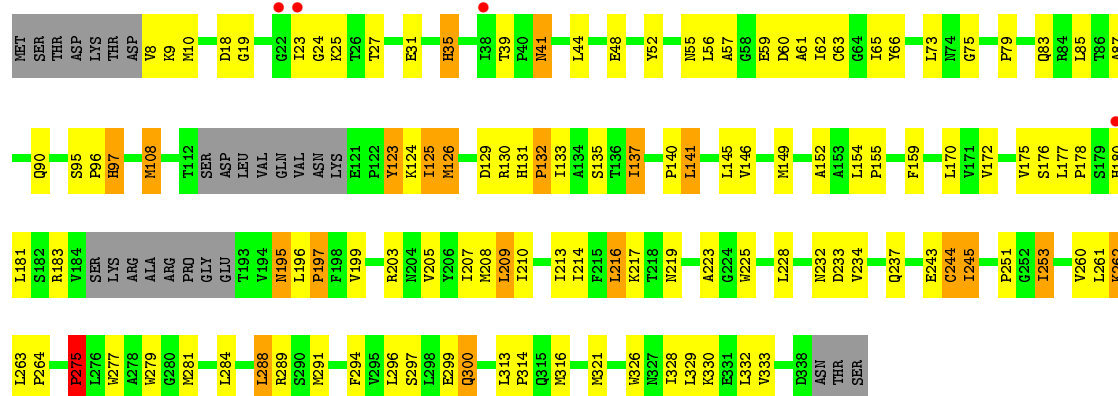


• Molecule 1: Thymidine kinase





● Molecule 1: Thymidine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	99.70 Å 54.20 Å 167.80 Å 90.00° 94.80° 90.00°	Depositor
Resolution (Å)	29.78 – 3.20 29.77 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.78-3.20) 78.2 (29.77-3.20)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 3.18 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.235 , 0.268 0.217 , 0.256	Depositor DCC
R_{free} test set	1490 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.619	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 16.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Outliers	0 of 30033 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10038	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.97% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

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: BVP, 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	0.42	0/2584	0.65	0/3519
1	B	0.41	0/2505	0.63	0/3411
1	C	0.41	0/2492	0.63	0/3391
1	D	0.41	0/2511	0.63	0/3418
All	All	0.41	0/10092	0.64	0/13739

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2524	0	2544	132	0
1	B	2446	0	2456	125	0
1	C	2433	0	2452	127	0
1	D	2452	0	2474	123	0
2	A	8	0	0	2	0
2	B	27	0	12	2	0
2	C	27	0	12	3	0
2	D	27	0	12	2	0
3	A	23	0	10	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	23	0	10	2	0
3	C	23	0	10	3	0
3	D	23	0	10	2	0
4	A	2	0	0	1	0
All	All	10038	0	10002	470	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 23.

The worst 5 of 470 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:63:CYS:HB2	1:D:63:CYS:SG	2.20	0.81
1:A:63:CYS:SG	1:D:63:CYS:HB2	2.21	0.80
1:C:85:LEU:HD13	1:D:85:LEU:HD13	1.63	0.79
1:A:9:LYS:HE3	1:A:109:ASP:OD2	1.83	0.77
1:A:275:PRO:HB3	1:B:336:VAL:HG12	1.66	0.77

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	321/341 (94%)	278 (87%)	35 (11%)	8 (2%)	7	41
1	B	309/341 (91%)	264 (85%)	40 (13%)	5 (2%)	12	54
1	C	307/341 (90%)	267 (87%)	35 (11%)	5 (2%)	12	54
1	D	309/341 (91%)	267 (86%)	35 (11%)	7 (2%)	8	44
All	All	1246/1364 (91%)	1076 (86%)	145 (12%)	25 (2%)	9	48

5 of 25 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	ALA
1	A	245	ILE
1	B	223	ALA
1	B	245	ILE
1	C	223	ALA

5.3.2 Protein sidechains ⓘ

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	280/298 (94%)	257 (92%)	23 (8%)	14	50
1	B	270/298 (91%)	248 (92%)	22 (8%)	15	51
1	C	268/298 (90%)	248 (92%)	20 (8%)	17	55
1	D	272/298 (91%)	249 (92%)	23 (8%)	13	47
All	All	1090/1192 (91%)	1002 (92%)	88 (8%)	15	51

5 of 88 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	219	ASN
1	C	73	LEU
1	D	216	LEU
1	B	253	ILE
1	B	288	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	204	ASN
1	C	74	ASN
1	D	204	ASN
1	B	220	ASN
1	B	300	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 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	ADP	A	400	-	3,7,29	1.63	0	5,10,45	0.75	0
3	BVP	A	500	-	19,24,24	3.68	5 (26%)	23,35,35	2.56	3 (13%)
2	ADP	B	1400	-	22,29,29	1.71	3 (13%)	27,45,45	0.98	1 (3%)
3	BVP	B	1500	-	19,24,24	3.71	5 (26%)	23,35,35	2.54	3 (13%)
2	ADP	C	2400	-	22,29,29	1.75	3 (13%)	27,45,45	1.02	1 (3%)
3	BVP	C	2500	-	19,24,24	3.70	5 (26%)	23,35,35	2.50	3 (13%)
2	ADP	D	3400	-	22,29,29	1.81	4 (18%)	27,45,45	0.98	1 (3%)
3	BVP	D	3500	-	19,24,24	3.65	5 (26%)	23,35,35	2.56	3 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	A	400	-	-	0/2/5/32	0/0/0/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BVP	A	500	-	-	0/9/25/25	0/2/2/2
2	ADP	B	1400	-	-	0/12/32/32	0/3/3/3
3	BVP	B	1500	-	-	0/9/25/25	0/2/2/2
2	ADP	C	2400	-	-	0/12/32/32	0/3/3/3
3	BVP	C	2500	-	-	0/9/25/25	0/2/2/2
2	ADP	D	3400	-	-	0/12/32/32	0/3/3/3
3	BVP	D	3500	-	-	0/9/25/25	0/2/2/2

The worst 5 of 30 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	3500	BVP	C5-C5A	-14.41	1.14	1.46
3	C	2500	BVP	C5-C5A	-14.36	1.14	1.46
3	B	1500	BVP	C5-C5A	-14.18	1.14	1.46
3	A	500	BVP	C5-C5A	-14.18	1.14	1.46
3	C	2500	BVP	P-O3P	-2.30	1.46	1.54

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	3500	BVP	C5-C4-N3	-3.43	117.83	124.04
3	A	500	BVP	C5-C4-N3	-3.37	117.94	124.04
3	C	2500	BVP	C5-C4-N3	-3.32	118.04	124.04
3	B	1500	BVP	C5-C4-N3	-3.29	118.08	124.04
2	B	1400	ADP	PA-O3A-PB	2.19	140.02	132.67

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	400	ADP	2	0
3	A	500	BVP	2	0
2	B	1400	ADP	2	0
3	B	1500	BVP	2	0
2	C	2400	ADP	3	0
3	C	2500	BVP	3	0
2	D	3400	ADP	2	0
3	D	3500	BVP	2	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](#)

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, 95th 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 ²)	Q<0.9
1	A	325/341 (95%)	-0.28	3 (0%) 85 78	5, 26, 88, 105	0
1	B	315/341 (92%)	-0.37	2 (0%) 90 84	5, 30, 87, 103	0
1	C	313/341 (91%)	-0.40	2 (0%) 90 84	5, 30, 87, 102	0
1	D	315/341 (92%)	-0.34	4 (1%) 79 67	7, 29, 89, 107	0
All	All	1268/1364 (92%)	-0.35	11 (0%) 85 78	5, 29, 88, 107	0

The worst 5 of 11 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	179	SER	4.0
1	D	23	ILE	3.1
1	A	29	ALA	2.8
1	D	22	GLY	2.5
1	C	42	ARG	2.4

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	LLDF	B-factors(\AA^2)	Q<0.9
2	ADP	D	3400	27/27	0.92	0.37	0.59	90,113,118,119	0
2	ADP	B	1400	27/27	0.92	0.28	0.30	85,109,116,123	0
2	ADP	C	2400	27/27	0.90	0.25	0.19	84,110,117,122	0
3	BVP	A	500	23/23	0.94	0.22	0.10	13,28,58,77	0
3	BVP	D	3500	23/23	0.95	0.18	-0.35	13,30,58,98	0
3	BVP	C	2500	23/23	0.97	0.17	-0.49	23,31,50,86	0
2	ADP	A	400	8/27	0.90	0.19	-0.49	94,97,105,107	0
3	BVP	B	1500	23/23	0.98	0.17	-0.59	24,35,53,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.