



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:54 PM GMT

PDB ID : 4QBI
Title : Crystal structure of a stable adenylate kinase variant AKlse6
Authors : Moon, S.; Bae, E.
Deposited on : 2014-05-08
Resolution : 1.67 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
<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

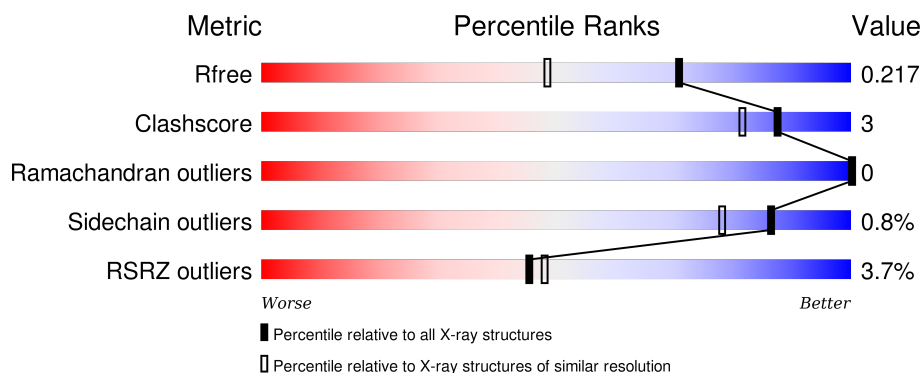
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.67 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	4802 (1.70-1.66)
Clashscore	102246	5317 (1.70-1.66)
Ramachandran outliers	100387	5225 (1.70-1.66)
Sidechain outliers	100360	5224 (1.70-1.66)
RSRZ outliers	91569	4813 (1.70-1.66)

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	217	<div> <div>2%</div> <div>93%</div> <div>6%</div> </div>
1	B	217	<div> <div>5%</div> <div>90%</div> <div>8%</div> </div>

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	ZN	A	301	-	-	-	X
2	ZN	B	301	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4038 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 Adenylate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1706	1067	305	321	13			
1	B	214	Total	C	N	O	S	0	0	0
			1677	1049	298	317	13			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	GLU	ALA	ENGINEERED MUTATION	UNP P27142
A	73	GLY	SER	ENGINEERED MUTATION	UNP P27142
A	78	GLU	GLN	ENGINEERED MUTATION	UNP P27142
A	79	ARG	ASN	ENGINEERED MUTATION	UNP P27142
A	98	GLU	THR	ENGINEERED MUTATION	UNP P27142
A	99	ILE	MET	ENGINEERED MUTATION	UNP P27142
A	101	GLU	ALA	ENGINEERED MUTATION	UNP P27142
A	107	ILE	LEU	ENGINEERED MUTATION	UNP P27142
A	166	GLU	ALA	ENGINEERED MUTATION	UNP P27142
A	170	LYS	ASN	ENGINEERED MUTATION	UNP P27142
A	183	LEU	VAL	ENGINEERED MUTATION	UNP P27142
A	195	VAL	ILE	ENGINEERED MUTATION	UNP P27142
A	201	ILE	MET	ENGINEERED MUTATION	UNP P27142
A	208	VAL	ILE	ENGINEERED MUTATION	UNP P27142
A	216	LYS	ALA	ENGINEERED MUTATION	UNP P27142
B	22	GLU	ALA	ENGINEERED MUTATION	UNP P27142
B	73	GLY	SER	ENGINEERED MUTATION	UNP P27142
B	78	GLU	GLN	ENGINEERED MUTATION	UNP P27142
B	79	ARG	ASN	ENGINEERED MUTATION	UNP P27142
B	98	GLU	THR	ENGINEERED MUTATION	UNP P27142
B	99	ILE	MET	ENGINEERED MUTATION	UNP P27142
B	101	GLU	ALA	ENGINEERED MUTATION	UNP P27142
B	107	ILE	LEU	ENGINEERED MUTATION	UNP P27142
B	166	GLU	ALA	ENGINEERED MUTATION	UNP P27142
B	170	LYS	ASN	ENGINEERED MUTATION	UNP P27142

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Chain	Residue	Modelled	Actual	Comment	Reference
B	183	LEU	VAL	ENGINEERED MUTATION	UNP P27142
B	195	VAL	ILE	ENGINEERED MUTATION	UNP P27142
B	201	ILE	MET	ENGINEERED MUTATION	UNP P27142
B	208	VAL	ILE	ENGINEERED MUTATION	UNP P27142
B	216	LYS	ALA	ENGINEERED MUTATION	UNP P27142

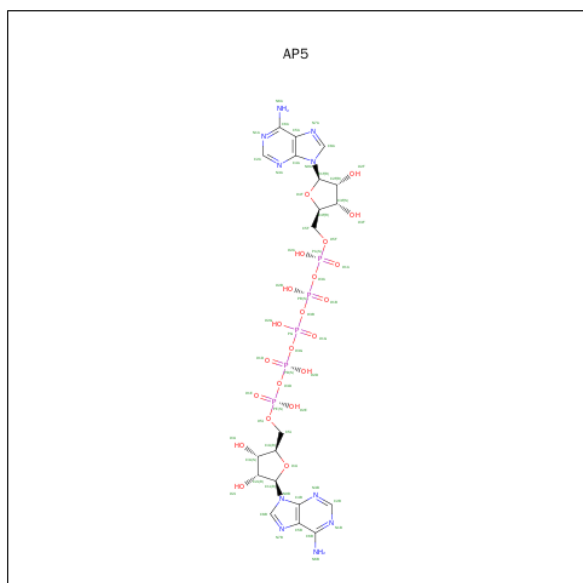
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Zn 1 1	0	0
2	A	1	Total Zn 1 1	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mg 1 1	0	0
3	A	1	Total Mg 1 1	0	0

- Molecule 4 is BIS(ADENOSINE)-5'-PENTAPHOSPHATE (three-letter code: AP5) (formula: C₂₀H₂₉N₁₀O₂₂P₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			57	20	10	22	5		
4	B	1	Total	C	N	O	P	0	0
			57	20	10	22	5		

- Molecule 5 is water.

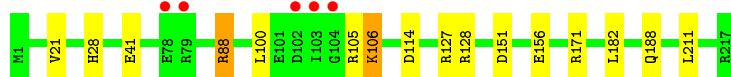
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	280	Total	O	0	0
			280	280		
5	B	257	Total	O	0	0
			257	257		

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: Adenylate kinase

Chain A: 



• Molecule 1: Adenylate kinase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	36.81Å 76.51Å 85.13Å 90.00° 90.20° 90.00°	Depositor
Resolution (Å)	50.00 – 1.67 30.88 – 1.67	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-1.67) 99.9 (30.88-1.67)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 1.67Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.168 , 0.208 0.181 , 0.217	Depositor DCC
R_{free} test set	2779 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 66.5	EDS
Estimated twinning fraction	0.033 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 54799 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4038	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.84% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MG, AP5

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	1.20	0/1731	1.07	8/2324 (0.3%)
1	B	1.22	3/1702 (0.2%)	1.10	11/2288 (0.5%)
All	All	1.21	3/3433 (0.1%)	1.09	19/4612 (0.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	156	GLU	CD-OE2	5.80	1.32	1.25
1	B	158	TYR	CE1-CZ	5.34	1.45	1.38
1	B	137	TYR	CZ-OH	5.12	1.46	1.37

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	209	ARG	NE-CZ-NH1	7.64	124.12	120.30
1	B	118	ASP	CB-CG-OD1	7.48	125.03	118.30
1	A	105	ARG	NE-CZ-NH1	7.34	123.97	120.30
1	A	171	ARG	NE-CZ-NH1	7.28	123.94	120.30
1	B	105	ARG	NE-CZ-NH2	-7.22	116.69	120.30
1	B	151	ASP	CB-CG-OD1	6.83	124.44	118.30
1	B	193	ARG	NE-CZ-NH1	6.38	123.49	120.30
1	B	114	ASP	CB-CG-OD1	6.10	123.79	118.30
1	A	211	LEU	CA-CB-CG	6.02	129.14	115.30
1	A	151	ASP	CB-CG-OD1	5.82	123.53	118.30
1	B	121	MET	CB-CG-SD	5.57	129.11	112.40
1	B	200	ASP	CB-CG-OD1	5.52	123.27	118.30
1	B	128	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	A	88	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	A	114	ASP	CB-CG-OD2	-5.35	113.48	118.30
1	B	209	ARG	NE-CZ-NH2	-5.33	117.63	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	53	MET	CG-SD-CE	-5.33	91.67	100.20
1	A	182	LEU	CB-CG-CD2	-5.14	102.27	111.00
1	A	128	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	1706	0	1716	12	0
1	B	1677	0	1679	6	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	57	0	24	2	0
4	B	57	0	24	1	0
5	A	280	0	0	5	0
5	B	257	0	0	1	0
All	All	4038	0	3443	18	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 (18) 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:104:GLY:O	1:B:105:ARG:HG2	1.85	0.75
1:A:106:LYS:HE2	5:A:503:HOH:O	1.95	0.66
1:B:104:GLY:C	1:B:105:ARG:HG2	2.20	0.61
1:A:41:GLU:HG2	5:A:654:HOH:O	2.01	0.59
1:A:188:GLN:NE2	5:A:664:HOH:O	2.35	0.59
1:A:127:ARG:HH12	4:A:303:AP5:PG	2.28	0.57
1:A:88:ARG:NH2	4:A:303:AP5:O1D	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:LEU:HB2	1:A:106:LYS:HE3	1.89	0.53
1:A:106:LYS:CE	1:A:106:LYS:HA	2.38	0.53
1:A:41:GLU:CG	5:A:654:HOH:O	2.56	0.53
1:A:100:LEU:CB	1:A:106:LYS:HE3	2.41	0.50
1:A:21:VAL:HG21	1:A:28:HIS:HB2	1.94	0.50
1:A:106:LYS:HE2	1:A:106:LYS:HA	1.96	0.46
1:B:104:GLY:HA2	5:B:637:HOH:O	2.15	0.46
1:B:88:ARG:NH2	4:B:303:AP5:O1D	2.49	0.45
1:B:104:GLY:C	1:B:105:ARG:CG	2.85	0.43
1:B:21:VAL:HG11	1:B:28:HIS:HB2	2.02	0.42
1:A:156:GLU:HG3	5:A:479:HOH:O	2.21	0.41

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	215/217 (99%)	210 (98%)	5 (2%)	0	100	100
1	B	212/217 (98%)	210 (99%)	2 (1%)	0	100	100
All	All	427/434 (98%)	420 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	180/180 (100%)	179 (99%)	1 (1%)	90	84
1	B	177/180 (98%)	175 (99%)	2 (1%)	80	67
All	All	357/360 (99%)	354 (99%)	3 (1%)	86	78

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	LYS
1	B	51	GLN
1	B	106	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](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	AP5	A	303	3	44,62,62	1.89	11 (25%)	55,98,98	2.26	14 (25%)
4	AP5	B	303	3	44,62,62	1.04	3 (6%)	55,98,98	2.11	11 (20%)

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	AP5	A	303	3	-	0/36/76/76	0/6/6/6
4	AP5	B	303	3	-	0/36/76/76	0/6/6/6

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	303	AP5	C2J-C3J	2.03	1.58	1.53
4	B	303	AP5	O4F-C1F	2.09	1.43	1.41
4	A	303	AP5	PG-O2G	2.27	1.64	1.54
4	B	303	AP5	C5B-C4B	2.38	1.45	1.40
4	A	303	AP5	O3J-C3J	2.51	1.49	1.43
4	A	303	AP5	C4B-N3B	2.56	1.39	1.35
4	A	303	AP5	C5A-C4A	2.63	1.46	1.40
4	A	303	AP5	C5B-C4B	2.72	1.46	1.40
4	A	303	AP5	C8A-N7A	2.87	1.40	1.34
4	B	303	AP5	O4J-C1J	3.32	1.45	1.41
4	A	303	AP5	C2B-N3B	3.33	1.38	1.32
4	A	303	AP5	C2A-N3A	3.65	1.38	1.32
4	A	303	AP5	O4F-C1F	5.42	1.48	1.41
4	A	303	AP5	O4J-C1J	6.23	1.49	1.41

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	303	AP5	PE-O3D-PD	-10.04	104.53	132.73
4	B	303	AP5	PE-O3D-PD	-5.94	116.05	132.73
4	B	303	AP5	N3A-C2A-N1A	-5.66	124.56	128.89
4	B	303	AP5	PD-O3G-PG	-4.82	119.19	132.73
4	A	303	AP5	N3B-C2B-N1B	-3.82	125.97	128.89
4	B	303	AP5	N3B-C2B-N1B	-3.73	126.04	128.89
4	A	303	AP5	N3A-C2A-N1A	-3.65	126.10	128.89
4	B	303	AP5	O2D-PD-O1D	-3.12	95.63	112.53
4	A	303	AP5	PD-O3G-PG	-3.10	124.03	132.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	303	AP5	C4B-C5B-N7B	-2.96	106.75	109.48
4	B	303	AP5	C2J-C1J-N9B	-2.77	110.06	114.29
4	A	303	AP5	C1F-N9A-C4A	-2.72	122.83	126.94
4	B	303	AP5	C4A-C5A-N7A	-2.49	107.19	109.48
4	B	303	AP5	C1F-N9A-C4A	-2.37	123.36	126.94
4	A	303	AP5	O5J-PE-O1E	-2.12	101.37	109.62
4	A	303	AP5	O2G-PG-O3B	2.03	114.31	105.09
4	A	303	AP5	O3J-C3J-C4J	2.08	117.29	111.05
4	A	303	AP5	O2D-PD-O1D	2.35	125.25	112.53
4	B	303	AP5	O2D-PD-O3D	2.38	115.87	105.09
4	A	303	AP5	C4J-O4J-C1J	2.66	112.64	109.72
4	A	303	AP5	O2D-PD-O3G	2.93	118.39	105.09
4	A	303	AP5	O2G-PG-O1G	3.37	130.78	112.53
4	A	303	AP5	O3D-PE-O5J	4.89	115.92	102.94
4	A	303	AP5	O2E-PE-O3D	5.84	131.59	105.09
4	B	303	AP5	O2D-PD-O3G	6.98	136.75	105.09

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	303	AP5	2	0
4	B	303	AP5	1	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	217/217 (100%)	-0.05	5 (2%) 64 67	16, 23, 40, 63	0
1	B	214/217 (98%)	0.12	11 (5%) 32 32	16, 24, 47, 65	0
All	All	431/434 (99%)	0.04	16 (3%) 45 48	16, 23, 43, 65	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	79	ARG	3.7
1	B	214	GLY	3.6
1	A	79	ARG	3.5
1	B	24	TYR	3.3
1	B	104	GLY	3.2
1	B	213	GLY	3.1
1	B	22	GLU	2.7
1	B	101	GLU	2.7
1	B	103	ILE	2.7
1	A	103	ILE	2.5
1	A	104	GLY	2.5
1	A	78	GLU	2.3
1	B	78	GLU	2.3
1	A	102	ASP	2.3
1	B	102	ASP	2.2
1	B	188	GLN	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. 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(Å ²)	Q<0.9
2	ZN	B	301	1/1	1.00	0.12	3.40	27,27,27,27	0
2	ZN	A	301	1/1	0.99	0.13	3.31	29,29,29,29	0
4	AP5	B	303	57/57	0.96	0.10	0.04	14,17,28,36	8
4	AP5	A	303	57/57	0.97	0.09	-0.17	15,19,32,43	8
3	MG	B	302	1/1	0.99	0.10	-	31,31,31,31	0
3	MG	A	302	1/1	0.97	0.11	-	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.