



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:23 PM GMT

PDB ID : 4I5V  
Title : Crystal structure of yeast Ap4A phosphorylase Apa2 in complex with Ap4A  
Authors : Jiang, Y.L.; Hou, W.T.; Chen, Y.; Zhou, C.Z.  
Deposited on : 2012-11-29  
Resolution : 2.70 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (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

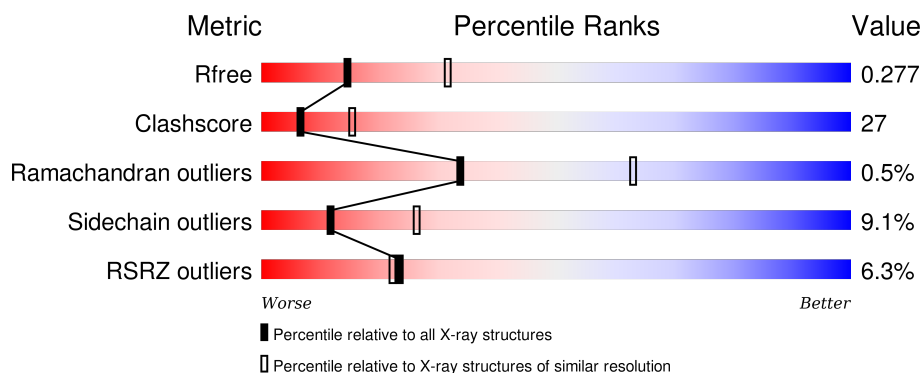
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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. 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	333	
1	B	333	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4843 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.

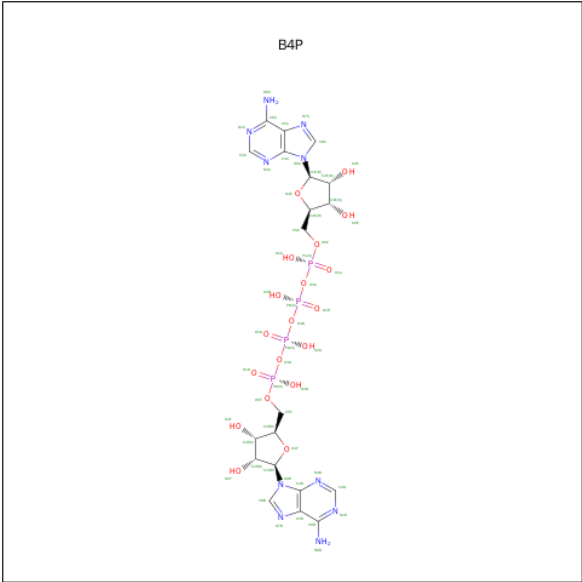
- Molecule 1 is a protein called 5',5'''-P-1,P-4-tetraphosphate phosphorylase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	0	0
			2347	1497	386	448	16			
1	B	295	Total	C	N	O	S	0	0	0
			2338	1492	384	446	16			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	EXPRESSION TAG	UNP P22108
A	-6	GLY	-	EXPRESSION TAG	UNP P22108
A	-5	HIS	-	EXPRESSION TAG	UNP P22108
A	-4	HIS	-	EXPRESSION TAG	UNP P22108
A	-3	HIS	-	EXPRESSION TAG	UNP P22108
A	-2	HIS	-	EXPRESSION TAG	UNP P22108
A	-1	HIS	-	EXPRESSION TAG	UNP P22108
A	0	HIS	-	EXPRESSION TAG	UNP P22108
A	161	ALA	HIS	ENGINEERED MUTATION	UNP P22108
B	-7	MET	-	EXPRESSION TAG	UNP P22108
B	-6	GLY	-	EXPRESSION TAG	UNP P22108
B	-5	HIS	-	EXPRESSION TAG	UNP P22108
B	-4	HIS	-	EXPRESSION TAG	UNP P22108
B	-3	HIS	-	EXPRESSION TAG	UNP P22108
B	-2	HIS	-	EXPRESSION TAG	UNP P22108
B	-1	HIS	-	EXPRESSION TAG	UNP P22108
B	0	HIS	-	EXPRESSION TAG	UNP P22108
B	161	ALA	HIS	ENGINEERED MUTATION	UNP P22108

- Molecule 2 is BIS(ADENOSINE)-5'-TETRAPHOSPHATE (three-letter code: B4P) (formula: C<sub>20</sub>H<sub>28</sub>N<sub>10</sub>O<sub>19</sub>P<sub>4</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	20	10	19	4		
2	B	1	Total	C	N	O	P	0	0
			53	20	10	19	4		

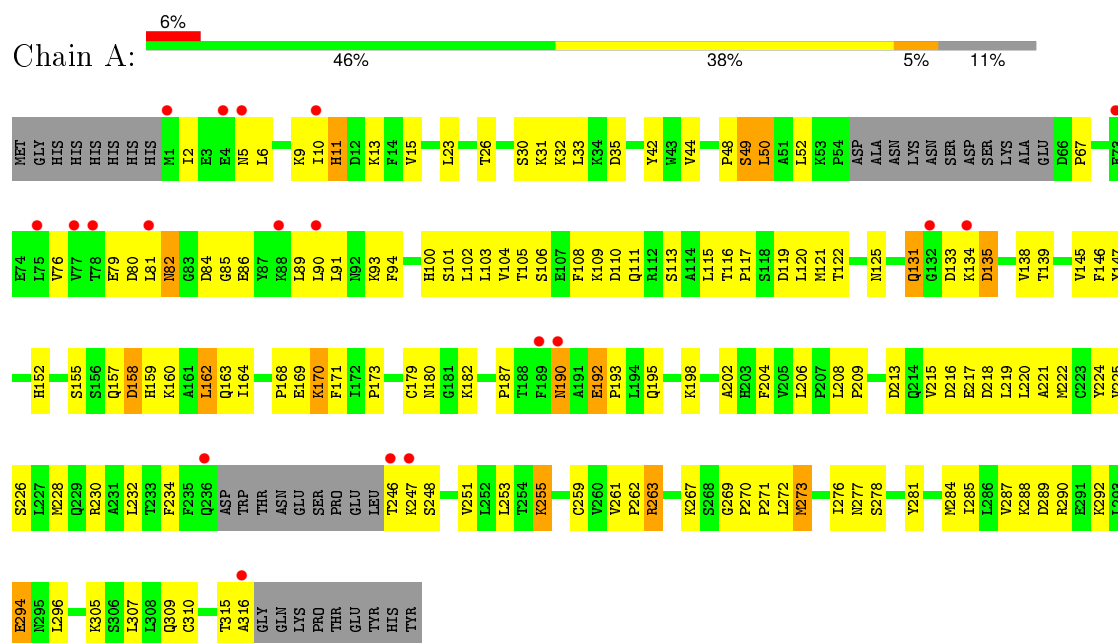
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	37	Total	O	0	0
			37	37		
3	B	15	Total	O	0	0
			15	15		

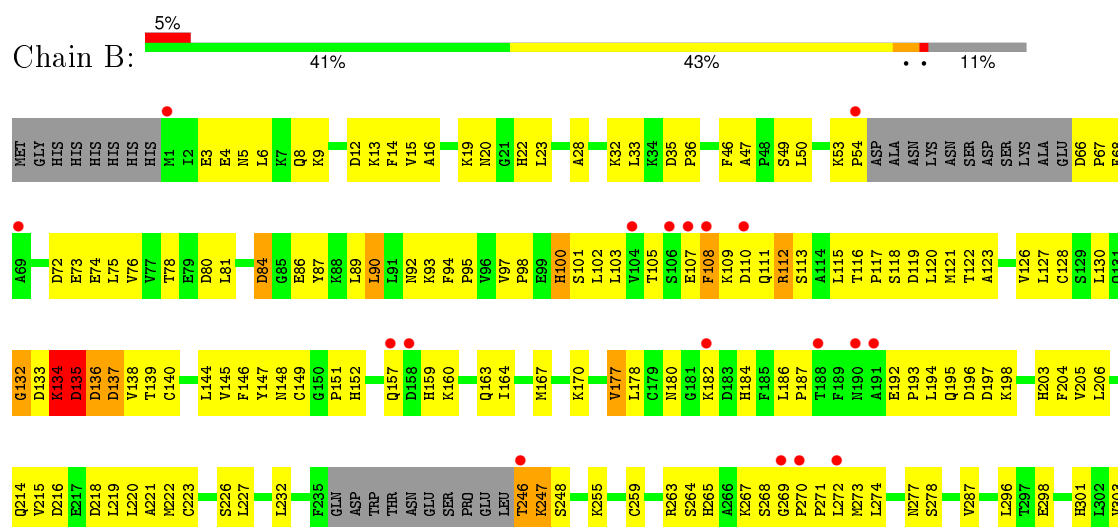
### 3 Residue-property plots

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: 5',5'''-P-1,P-4-tetraphosphate phosphorylase 2



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D304	R305	S306	L307	L308	D309	C310	P313	N314	T315	A316	GLY	GLN	LYS	PRO	THR	GLU	TYR	HIS	TYR
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## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	203.35Å 52.78Å 70.76Å 90.00° 99.92° 90.00°	Depositor
Resolution (Å)	42.07 – 2.70 42.07 – 2.70	Depositor EDS
% Data completeness (in resolution range)	93.7 (42.07-2.70) 93.7 (42.07-2.70)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.91 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, $R_{free}$	0.181 , 0.261 0.210 , 0.277	Depositor DCC
$R_{free}$ test set	988 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	69.5	Xtriage
Anisotropy	0.446	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 61.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 19446 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4843	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% 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, 0.375 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: B4P

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	0.45	0/2401	0.64	0/3255
1	B	0.40	0/2392	0.59	1/3243 (0.0%)
All	All	0.43	0/4793	0.61	1/6498 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	132	GLY	N-CA-C	-5.53	99.27	113.10

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	2347	0	2323	118	1
1	B	2338	0	2315	142	1
2	A	53	0	22	1	0
2	B	53	0	22	0	0
3	A	37	0	0	2	0
3	B	15	0	0	0	0
All	All	4843	0	4682	258	1

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 27.

All (258) 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:67:PRO:HD2	1:A:157:GLN:HE22	0.99	1.08
1:A:192:GLU:HB2	1:A:193:PRO:HD2	1.38	1.02
1:B:84:ASP:HB2	1:B:86:GLU:HG2	1.52	0.92
1:A:67:PRO:HD2	1:A:157:GLN:NE2	1.84	0.92
1:A:263:ARG:NH2	1:A:276:ILE:O	2.02	0.91
1:B:53:LYS:HG3	1:B:54:PRO:HA	1.53	0.90
1:B:116:THR:HG1	1:B:118:SER:HG	1.19	0.89
1:B:187:PRO:HG2	1:B:226:SER:OG	1.73	0.88
1:A:2:ILE:HG12	1:A:82:ASN:HD22	1.40	0.87
1:A:218:ASP:O	1:A:222:MET:HG3	1.76	0.85
1:B:4:GLU:HG2	1:B:5:ASN:H	1.42	0.82
1:B:134:LYS:O	1:B:135:ASP:HB2	1.78	0.82
1:A:2:ILE:HG12	1:A:82:ASN:ND2	1.95	0.82
1:A:232:LEU:HD11	1:A:247:LYS:HE3	1.60	0.81
1:B:136:ASP:O	1:B:138:VAL:N	2.13	0.81
1:B:218:ASP:O	1:B:222:MET:HG2	1.83	0.79
1:B:268:SER:HB3	1:B:274:LEU:H	1.50	0.77
1:B:78:THR:HG21	1:B:81:LEU:HD13	1.67	0.77
1:B:135:ASP:O	1:B:137:ASP:N	2.18	0.76
1:B:136:ASP:O	1:B:139:THR:N	2.18	0.76
1:B:272:LEU:HD12	1:B:306:SER:OG	1.86	0.74
1:A:170:LYS:HE3	1:B:139:THR:O	1.88	0.74
1:B:145:VAL:HG22	1:B:164:ILE:HG12	1.70	0.74
1:A:121:MET:HA	1:A:220:LEU:HD12	1.70	0.73
1:A:192:GLU:HB2	1:A:193:PRO:CD	2.16	0.73
1:B:128:CYS:SG	1:B:255:LYS:NZ	2.61	0.72
1:B:5:ASN:O	1:B:9:LYS:HG3	1.90	0.71
1:A:67:PRO:CD	1:A:157:GLN:HE22	1.92	0.71
1:A:209:PRO:HG2	1:A:215:VAL:HG22	1.71	0.71
1:B:149:CYS:O	1:B:248:SER:HA	1.90	0.70
1:B:89:LEU:HD12	1:B:102:LEU:O	1.92	0.70
1:A:2:ILE:HD12	1:A:125:ASN:HB2	1.74	0.70
1:B:263:ARG:NH1	1:B:265:HIS:O	2.24	0.69
1:A:272:LEU:HD13	1:A:296:LEU:HD21	1.74	0.69
1:B:206:LEU:HD12	1:B:223:CYS:HB3	1.74	0.69
1:B:80:ASP:HB2	1:B:84:ASP:O	1.92	0.68
1:B:263:ARG:NH1	1:B:310:CYS:HB3	2.09	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:ASP:OD2	1:A:86:GLU:HB2	1.95	0.67
1:B:28:ALA:HB2	1:B:47:ALA:HB2	1.75	0.67
1:B:122:THR:O	1:B:126:VAL:HG23	1.93	0.67
1:B:196:ASP:OD1	1:B:198:LYS:HD3	1.95	0.66
1:A:173:PRO:HA	3:A:508:HOH:O	1.95	0.66
1:A:187:PRO:HD3	1:A:206:LEU:HD11	1.77	0.65
1:B:78:THR:HG22	1:B:89:LEU:HB3	1.78	0.65
1:B:90:LEU:N	1:B:90:LEU:HD12	2.11	0.65
1:B:206:LEU:HG	1:B:227:LEU:HD21	1.78	0.65
1:B:90:LEU:O	1:B:101:SER:HB2	1.97	0.65
1:B:216:ASP:OD1	1:B:219:LEU:HB2	1.97	0.65
1:B:195:GLN:HG2	1:B:203:HIS:O	1.97	0.64
1:A:13:LYS:HG3	1:A:91:LEU:HD11	1.79	0.64
1:B:20:ASN:HB2	1:B:22:HIS:ND1	2.12	0.64
1:A:193:PRO:HA	1:A:206:LEU:HD21	1.80	0.64
1:A:116:THR:N	1:A:119:ASP:OD2	2.27	0.63
1:A:120:LEU:HD11	1:A:224:TYR:CD1	2.34	0.62
1:A:30:SER:HA	1:A:44:VAL:O	1.99	0.62
1:B:187:PRO:HG2	1:B:226:SER:HG	1.64	0.62
1:A:26:THR:HG21	1:A:50:LEU:HD22	1.80	0.62
1:B:67:PRO:HG2	1:B:68:PHE:CE1	2.34	0.62
1:A:289:ASP:HB3	1:A:292:LYS:HD2	1.80	0.62
1:B:133:ASP:O	1:B:135:ASP:N	2.33	0.61
1:B:13:LYS:HE3	1:B:74:GLU:O	2.00	0.61
1:A:190:ASN:OD1	1:A:190:ASN:N	2.32	0.61
1:B:67:PRO:HD3	1:B:157:GLN:HE22	1.64	0.61
1:B:268:SER:HB3	1:B:274:LEU:N	2.16	0.60
1:B:264:SER:HB2	1:B:265:HIS:CE1	2.37	0.60
1:A:221:ALA:O	1:A:225:VAL:HG23	2.03	0.59
1:A:290:ARG:NE	1:A:294:GLU:OE1	2.32	0.59
1:B:116:THR:O	1:B:120:LEU:HD13	2.02	0.59
1:B:270:PRO:N	1:B:271:PRO:HD2	2.18	0.59
1:A:2:ILE:CG1	1:A:82:ASN:HD22	2.13	0.59
1:B:126:VAL:O	1:B:130:LEU:HG	2.03	0.58
1:A:115:LEU:CD2	1:A:162:LEU:HD21	2.32	0.58
1:B:136:ASP:C	1:B:138:VAL:N	2.57	0.58
1:B:127:LEU:HD21	1:B:145:VAL:HG23	1.86	0.58
1:A:35:ASP:HB2	1:A:42:TYR:CE2	2.38	0.58
1:B:92:ASN:O	1:B:95:PRO:HD3	2.04	0.57
1:A:2:ILE:HD12	1:A:125:ASN:CB	2.35	0.57
1:B:136:ASP:CG	1:B:137:ASP:N	2.58	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:272:LEU:O	1:A:292:LYS:HE2	2.05	0.56
1:A:267:LYS:HE3	1:A:273:MET:HB3	1.86	0.56
1:B:134:LYS:HG3	1:B:139:THR:HG21	1.87	0.56
1:B:134:LYS:O	1:B:135:ASP:CB	2.49	0.56
1:A:284:MET:C	1:A:285:ILE:HD12	2.26	0.55
1:B:115:LEU:HD21	1:B:147:TYR:CD2	2.41	0.55
1:A:116:THR:OG1	1:A:119:ASP:N	2.38	0.55
1:A:159:HIS:HD2	1:A:160:LYS:H	1.53	0.55
1:B:32:LYS:O	1:B:33:LEU:HD23	2.07	0.55
1:A:159:HIS:HD2	1:A:160:LYS:N	2.04	0.55
1:B:35:ASP:OD2	1:B:304:ASP:HB2	2.07	0.55
1:A:113:SER:O	1:A:147:TYR:OH	2.19	0.55
1:B:269:GLY:C	1:B:271:PRO:HD2	2.27	0.54
1:A:133:ASP:O	1:A:134:LYS:HB3	2.08	0.54
1:A:305:LYS:O	1:A:309:GLN:HG3	2.08	0.54
1:A:116:THR:O	1:A:119:ASP:HB2	2.09	0.54
1:A:285:ILE:N	1:A:285:ILE:CD1	2.71	0.54
1:A:134:LYS:O	1:A:135:ASP:HB2	2.08	0.53
1:A:93:LYS:HD3	1:A:94:PHE:CZ	2.44	0.53
1:B:81:LEU:HD22	1:B:89:LEU:HB2	1.90	0.53
1:B:152:HIS:HE1	1:B:246:THR:HG23	1.73	0.53
1:B:6:LEU:HD11	1:B:89:LEU:HD23	1.91	0.53
1:B:5:ASN:HB3	1:B:9:LYS:HE3	1.91	0.53
1:B:115:LEU:HB3	1:B:120:LEU:CD1	2.39	0.52
1:B:28:ALA:HB1	1:B:46:PHE:O	2.09	0.52
1:B:110:ASP:OD1	1:B:112:ARG:NH1	2.42	0.52
1:B:151:PRO:HG2	1:B:152:HIS:CE1	2.44	0.52
1:B:145:VAL:HA	1:B:163:GLN:O	2.10	0.52
1:A:121:MET:HA	1:A:220:LEU:CD1	2.40	0.52
1:A:285:ILE:N	1:A:285:ILE:HD12	2.25	0.52
1:B:132:GLY:O	1:B:134:LYS:N	2.43	0.51
1:A:105:THR:HB	1:A:109:LYS:HD2	1.92	0.51
1:A:157:GLN:HB3	2:A:401:B4P:O2D	2.09	0.51
1:B:206:LEU:HD12	1:B:223:CYS:CB	2.39	0.51
1:A:115:LEU:HD21	1:A:162:LEU:HD21	1.92	0.51
1:A:103:LEU:HB3	1:A:162:LEU:HB2	1.93	0.51
1:A:228:MET:O	1:A:232:LEU:HB2	2.11	0.51
1:B:272:LEU:HD13	1:B:296:LEU:HD21	1.93	0.51
1:A:49:SER:O	1:A:52:LEU:HD12	2.10	0.51
1:B:147:TYR:CE2	1:B:149:CYS:HB2	2.45	0.51
1:A:168:PRO:HG2	1:A:171:PHE:CB	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:THR:CB	1:A:109:LYS:HD2	2.40	0.50
1:B:264:SER:N	1:B:310:CYS:O	2.44	0.50
1:A:272:LEU:HD13	1:A:296:LEU:CD2	2.42	0.50
1:A:138:VAL:HG13	1:A:139:THR:N	2.26	0.50
1:A:32:LYS:O	1:A:33:LEU:HD23	2.11	0.50
1:A:159:HIS:CD2	1:A:160:LYS:N	2.80	0.50
1:A:195:GLN:NE2	1:A:202:ALA:HA	2.27	0.50
1:A:193:PRO:CD	1:A:230:ARG:HD2	2.42	0.50
1:B:84:ASP:N	1:B:84:ASP:OD2	2.45	0.50
1:B:144:LEU:HB2	1:B:167:MET:SD	2.52	0.49
1:B:87:TYR:HE1	1:B:105:THR:HG22	1.77	0.49
1:A:2:ILE:HD11	1:A:122:THR:HG23	1.93	0.49
1:B:53:LYS:HG3	1:B:54:PRO:CA	2.35	0.49
1:B:111:GLN:NE2	1:B:148:ASN:O	2.45	0.49
1:B:134:LYS:HD2	1:B:134:LYS:N	2.26	0.49
1:B:263:ARG:HB3	1:B:263:ARG:HH11	1.77	0.49
1:B:107:GLU:O	1:B:109:LYS:HG2	2.12	0.49
1:A:111:GLN:NE2	1:A:147:TYR:CE1	2.76	0.49
1:B:93:LYS:HB3	1:B:94:PHE:CE1	2.48	0.48
1:B:90:LEU:N	1:B:90:LEU:CD1	2.77	0.48
1:B:134:LYS:HG3	1:B:139:THR:CG2	2.44	0.48
1:B:152:HIS:CE1	1:B:246:THR:HG23	2.49	0.48
1:A:270:PRO:N	1:A:271:PRO:HD2	2.29	0.48
1:B:111:GLN:HG3	1:B:111:GLN:O	2.14	0.47
1:A:115:LEU:HD12	1:A:224:TYR:CE1	2.48	0.47
1:A:168:PRO:HG2	1:A:171:PHE:HB3	1.95	0.47
1:A:152:HIS:HB2	1:A:248:SER:HB2	1.95	0.47
1:A:146:PHE:HA	1:A:251:VAL:O	2.15	0.47
1:B:216:ASP:OD1	1:B:216:ASP:N	2.43	0.47
1:B:112:ARG:O	1:B:113:SER:C	2.53	0.47
1:B:23:LEU:HA	1:B:93:LYS:HG3	1.96	0.47
1:A:10:ILE:HD11	1:A:89:LEU:HD21	1.97	0.47
1:B:19:LYS:O	1:B:19:LYS:HD3	2.14	0.47
1:A:193:PRO:HD2	1:A:230:ARG:HD2	1.96	0.47
1:A:219:LEU:HD23	1:A:219:LEU:C	2.35	0.47
1:B:144:LEU:HD12	1:B:145:VAL:N	2.30	0.47
1:B:116:THR:O	1:B:119:ASP:HB2	2.15	0.46
1:A:209:PRO:HG2	1:A:215:VAL:CG2	2.43	0.46
1:B:271:PRO:CG	1:B:309:GLN:HE22	2.28	0.46
1:B:117:PRO:HB3	1:B:221:ALA:HB2	1.96	0.46
1:B:214:GLN:O	1:B:214:GLN:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:LYS:HE3	1:B:74:GLU:C	2.37	0.46
1:A:255:LYS:HE2	1:A:255:LYS:HB3	1.70	0.46
1:B:73:GLU:OE2	1:B:73:GLU:N	2.48	0.46
1:B:68:PHE:CD2	1:B:159:HIS:HB3	2.50	0.46
1:A:85:GLY:O	1:A:106:SER:OG	2.14	0.46
1:A:204:PHE:O	1:A:259:CYS:HA	2.16	0.46
1:A:152:HIS:HB2	1:A:248:SER:CB	2.47	0.45
1:A:84:ASP:CG	1:A:86:GLU:HB2	2.36	0.45
1:A:232:LEU:HA	1:A:232:LEU:HD12	1.76	0.45
1:A:134:LYS:O	1:A:135:ASP:CB	2.64	0.45
1:B:108:PHE:O	1:B:109:LYS:HD3	2.17	0.45
1:A:26:THR:HG21	1:A:50:LEU:CD2	2.44	0.45
1:A:23:LEU:HD12	1:A:23:LEU:C	2.36	0.45
1:B:121:MET:HA	1:B:220:LEU:HD12	1.98	0.45
1:B:116:THR:N	1:B:119:ASP:OD2	2.36	0.45
1:B:126:VAL:HG11	1:B:164:ILE:HD13	1.99	0.45
1:A:216:ASP:OD1	1:A:216:ASP:C	2.56	0.45
1:B:122:THR:OG1	1:B:123:ALA:N	2.50	0.44
1:B:263:ARG:HB3	1:B:263:ARG:NH1	2.33	0.44
1:B:267:LYS:HE2	1:B:273:MET:CE	2.47	0.44
1:B:78:THR:CG2	1:B:89:LEU:HB3	2.44	0.44
1:A:108:PHE:O	1:A:109:LYS:HG3	2.16	0.44
1:B:247:LYS:HD3	1:B:247:LYS:HA	1.77	0.44
1:A:110:ASP:O	1:A:113:SER:OG	2.20	0.44
1:B:67:PRO:HG2	1:B:68:PHE:CD1	2.53	0.44
1:B:135:ASP:O	1:B:136:ASP:C	2.51	0.44
1:A:269:GLY:O	1:A:273:MET:HG3	2.18	0.44
1:A:277:ASN:OD1	1:A:277:ASN:C	2.55	0.44
1:B:186:LEU:HA	1:B:187:PRO:HD3	1.73	0.44
1:A:13:LYS:HE3	1:A:76:VAL:O	2.17	0.44
1:A:80:ASP:OD2	1:A:81:LEU:N	2.50	0.44
1:A:271:PRO:HG2	3:A:511:HOH:O	2.18	0.43
1:B:136:ASP:C	1:B:138:VAL:H	2.21	0.43
1:A:246:THR:OG1	1:A:247:LYS:N	2.49	0.43
1:B:13:LYS:HE2	1:B:76:VAL:O	2.18	0.43
1:A:261:VAL:HG11	1:A:281:TYR:CD1	2.53	0.43
1:B:178:LEU:HD23	1:B:205:VAL:HG21	1.99	0.43
1:A:90:LEU:O	1:A:102:LEU:N	2.50	0.43
1:B:23:LEU:HD23	1:B:75:LEU:CD2	2.48	0.43
1:B:19:LYS:HD3	1:B:19:LYS:C	2.39	0.43
1:B:36:PRO:HG2	1:B:301:HIS:HB2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:HIS:HD2	1:B:160:LYS:N	2.16	0.43
1:B:146:PHE:HB2	1:B:278:SER:CB	2.49	0.43
1:A:234:PHE:CE2	1:A:262:PRO:HG3	2.54	0.43
1:B:215:VAL:HA	1:B:219:LEU:HD13	2.00	0.43
1:B:148:ASN:OD1	1:B:277:ASN:HB2	2.18	0.43
1:B:4:GLU:HG2	1:B:5:ASN:N	2.20	0.43
1:A:13:LYS:HA	1:A:13:LYS:HD3	1.84	0.43
1:B:115:LEU:CB	1:B:120:LEU:HD11	2.48	0.42
1:B:296:LEU:HD22	1:B:303:VAL:HG22	2.00	0.42
1:A:292:LYS:O	1:A:296:LEU:HG	2.18	0.42
1:A:117:PRO:HB3	1:A:221:ALA:HB2	2.00	0.42
1:B:8:GLN:NE2	1:B:12:ASP:OD1	2.52	0.42
1:A:6:LEU:O	1:A:9:LYS:HB2	2.19	0.42
1:A:263:ARG:HD2	1:A:310:CYS:O	2.19	0.42
1:B:204:PHE:O	1:B:259:CYS:HA	2.19	0.42
1:B:66:ASP:HA	1:B:67:PRO:HD3	1.88	0.42
1:B:308:LEU:HD11	1:B:313:PRO:HA	2.00	0.42
1:A:162:LEU:HD23	1:A:162:LEU:N	2.35	0.42
1:B:267:LYS:HE2	1:B:273:MET:HE3	2.02	0.42
1:A:111:GLN:HB2	1:A:158:ASP:C	2.40	0.42
1:A:145:VAL:HG22	1:A:164:ILE:HG13	2.00	0.42
1:A:307:LEU:HA	1:A:307:LEU:HD23	1.91	0.42
1:B:136:ASP:O	1:B:138:VAL:C	2.57	0.42
1:A:133:ASP:C	1:A:135:ASP:H	2.23	0.42
1:A:169:GLU:OE1	1:B:140:CYS:N	2.38	0.42
1:B:272:LEU:HD23	1:B:272:LEU:HA	1.93	0.41
1:B:68:PHE:HD2	1:B:159:HIS:CG	2.38	0.41
1:A:11:HIS:CE1	1:A:15:VAL:CG1	3.03	0.41
1:A:179:CYS:O	1:A:182:LYS:HB2	2.20	0.41
1:A:100:HIS:HD2	1:A:101:SER:N	2.18	0.41
1:B:78:THR:HG23	1:B:89:LEU:H	1.85	0.41
1:B:178:LEU:O	1:B:194:LEU:HD22	2.21	0.41
1:A:48:PRO:HA	1:A:288:LYS:O	2.20	0.41
1:B:100:HIS:C	1:B:100:HIS:ND1	2.74	0.41
1:B:118:SER:O	1:B:122:THR:HG23	2.21	0.41
1:A:234:PHE:CD2	1:A:262:PRO:HG3	2.56	0.41
1:A:90:LEU:HB2	1:A:102:LEU:HB2	2.03	0.41
1:A:145:VAL:HB	1:A:253:LEU:HB3	2.03	0.41
1:A:315:THR:O	1:A:316:ALA:C	2.59	0.41
1:B:97:VAL:HA	1:B:98:PRO:HD2	1.79	0.41
1:A:193:PRO:HA	1:A:206:LEU:CD2	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:GLU:HB2	1:B:193:PRO:CD	2.50	0.41
1:B:3:GLU:HG2	1:B:81:LEU:HD12	2.03	0.41
1:B:272:LEU:HD13	1:B:296:LEU:CD2	2.50	0.41
1:B:246:THR:OG1	1:B:246:THR:O	2.34	0.41
1:B:177:VAL:HA	1:B:180:ASN:HB2	2.03	0.41
1:B:15:VAL:HG23	1:B:16:ALA:N	2.36	0.40
1:A:5:ASN:O	1:A:9:LYS:HD3	2.22	0.40
1:A:145:VAL:HA	1:A:163:GLN:O	2.20	0.40
1:A:131:GLN:C	1:A:133:ASP:H	2.25	0.40
1:B:23:LEU:C	1:B:23:LEU:HD12	2.41	0.40
1:B:14:PHE:CD1	1:B:23:LEU:HD21	2.56	0.40
1:A:187:PRO:HB3	1:A:192:GLU:HA	2.03	0.40
1:B:187:PRO:HB3	1:B:193:PRO:HD3	2.04	0.40
1:B:144:LEU:HD12	1:B:145:VAL:H	1.87	0.40
1:B:132:GLY:C	1:B:134:LYS:N	2.74	0.40
1:A:208:LEU:HB3	1:A:209:PRO:HD2	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:LYS:NZ	1:B:218:ASP:OD1[1_554]	2.17	0.03

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	290/333 (87%)	272 (94%)	18 (6%)	0	100	100
1	B	289/333 (87%)	266 (92%)	20 (7%)	3 (1%)	19	45
All	All	579/666 (87%)	538 (93%)	38 (7%)	3 (0%)	34	63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	135	ASP
1	B	137	ASP
1	B	134	LYS

### 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	270/303 (89%)	244 (90%)	26 (10%)	10	24
1	B	269/303 (89%)	246 (91%)	23 (9%)	13	29
All	All	539/606 (89%)	490 (91%)	49 (9%)	12	26

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	HIS
1	A	31	LYS
1	A	49	SER
1	A	50	LEU
1	A	79	GLU
1	A	82	ASN
1	A	104	VAL
1	A	131	GLN
1	A	135	ASP
1	A	155	SER
1	A	158	ASP
1	A	162	LEU
1	A	170	LYS
1	A	180	ASN
1	A	190	ASN
1	A	192	GLU
1	A	198	LYS
1	A	213	ASP
1	A	217	GLU
1	A	226	SER

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Mol	Chain	Res	Type
1	A	255	LYS
1	A	263	ARG
1	A	273	MET
1	A	278	SER
1	A	287	VAL
1	A	294	GLU
1	B	49	SER
1	B	50	LEU
1	B	72	ASP
1	B	84	ASP
1	B	90	LEU
1	B	100	HIS
1	B	103	LEU
1	B	108	PHE
1	B	112	ARG
1	B	134	LYS
1	B	135	ASP
1	B	136	ASP
1	B	170	LYS
1	B	177	VAL
1	B	182	LYS
1	B	184	HIS
1	B	197	ASP
1	B	232	LEU
1	B	246	THR
1	B	247	LYS
1	B	287	VAL
1	B	298	GLU
1	B	315	THR

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

Mol	Chain	Res	Type
1	A	82	ASN
1	A	157	GLN

### 5.3.3 RNA ⓘ

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](#)

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	B4P	A	401	-	42,58,58	1.61	9 (21%)	51,91,91	3.08	11 (21%)
2	B4P	B	401	-	42,58,58	1.64	9 (21%)	51,91,91	3.14	12 (23%)

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	B4P	A	401	-	-	0/30/70/70	0/6/6/6
2	B4P	B	401	-	-	0/30/70/70	0/6/6/6

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	B4P	C2F-C3F	-3.72	1.43	1.53
2	A	401	B4P	O2F-C2F	-3.39	1.34	1.43
2	A	401	B4P	C2F-C3F	-3.31	1.44	1.53
2	B	401	B4P	O2F-C2F	-3.24	1.35	1.43
2	A	401	B4P	C2E-C3E	-3.07	1.45	1.53
2	B	401	B4P	C2E-C3E	-3.05	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	B4P	O3E-C3E	-2.99	1.35	1.43
2	B	401	B4P	O3E-C3E	-2.93	1.35	1.43
2	A	401	B4P	O4E-C4E	-2.83	1.38	1.45
2	B	401	B4P	O4E-C4E	-2.46	1.39	1.45
2	A	401	B4P	O5F-C5F	-2.04	1.36	1.44
2	B	401	B4P	O5F-C5F	-2.01	1.36	1.44
2	A	401	B4P	C6A-N6A	2.58	1.42	1.34
2	A	401	B4P	C6B-N6B	2.58	1.42	1.34
2	B	401	B4P	C6B-N6B	2.60	1.43	1.34
2	B	401	B4P	C6A-N6A	2.63	1.43	1.34
2	A	401	B4P	O4E-C1E	3.93	1.46	1.41
2	B	401	B4P	O4E-C1E	4.42	1.46	1.41

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	B4P	N3B-C2B-N1B	-12.58	119.27	128.89
2	B	401	B4P	N3B-C2B-N1B	-11.57	120.03	128.89
2	B	401	B4P	N3A-C2A-N1A	-10.81	120.61	128.89
2	A	401	B4P	C1F-N9B-C4B	-10.66	110.86	126.94
2	B	401	B4P	C1F-N9B-C4B	-10.18	111.58	126.94
2	A	401	B4P	N3A-C2A-N1A	-9.34	121.74	128.89
2	B	401	B4P	PG-O3B-PB	-6.15	115.47	132.73
2	A	401	B4P	PG-O3B-PB	-4.86	119.09	132.73
2	B	401	B4P	PB-O3A-PA	-4.32	120.60	132.73
2	A	401	B4P	C4A-C5A-N7A	-3.69	106.09	109.48
2	B	401	B4P	C4A-C5A-N7A	-3.53	106.24	109.48
2	A	401	B4P	PB-O3A-PA	-2.86	124.69	132.73
2	B	401	B4P	PD-O3G-PG	-2.79	124.89	132.73
2	A	401	B4P	C1E-N9A-C4A	-2.13	123.72	126.94
2	B	401	B4P	C4F-O4F-C1F	-2.12	107.39	109.72
2	B	401	B4P	O2G-PG-O3G	2.00	114.17	105.09
2	B	401	B4P	O5E-C5E-C4E	2.04	116.64	109.12
2	A	401	B4P	O4F-C1F-N9B	2.19	112.69	108.10
2	A	401	B4P	O3A-PA-O5E	2.45	109.44	102.94
2	B	401	B4P	O4F-C1F-N9B	3.07	114.53	108.10
2	A	401	B4P	O4E-C1E-N9A	3.40	115.22	108.10
2	B	401	B4P	O3G-PD-O5F	3.61	112.52	102.94
2	A	401	B4P	O3G-PD-O5F	4.61	115.17	102.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	B4P	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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	296/333 (88%)	0.17	19 (6%) 23 21	43, 75, 114, 141	0
1	B	295/333 (88%)	0.32	18 (6%) 25 23	58, 90, 128, 150	0
All	All	591/666 (88%)	0.25	37 (6%) 23 22	43, 84, 123, 150	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	MET	5.7
1	A	190	ASN	5.6
1	A	1	MET	5.2
1	B	191	ALA	4.3
1	A	246	THR	4.0
1	B	104	VAL	3.8
1	B	272	LEU	3.5
1	B	108	PHE	3.5
1	B	157	GLN	3.5
1	A	75	LEU	3.3
1	B	110	ASP	3.1
1	B	270	PRO	2.9
1	A	81	LEU	2.8
1	B	188	THR	2.7
1	A	247	LYS	2.7
1	B	54	PRO	2.7
1	B	158	ASP	2.7
1	B	182	LYS	2.7
1	A	4	GLU	2.6
1	B	106	SER	2.6
1	A	189	PHE	2.6
1	B	190	ASN	2.5
1	A	88	LYS	2.5
1	A	134	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	316	ALA	2.4
1	B	69	ALA	2.3
1	B	107	GLU	2.3
1	A	73	GLU	2.3
1	A	236	GLN	2.3
1	A	5	ASN	2.2
1	B	269	GLY	2.2
1	A	90	LEU	2.2
1	A	77	VAL	2.2
1	B	246	THR	2.1
1	A	78	THR	2.1
1	A	10	ILE	2.1
1	A	132	GLY	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 [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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	B4P	A	401	53/53	0.98	0.14	-0.65	50,63,89,91	0
2	B4P	B	401	53/53	0.97	0.18	-0.67	54,80,101,105	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.