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ERRATUM

Viral Escape at the Molecular Level Explained by Quantitative T-cell Receptor/Peptide/MHC Interactions and the Crystal Structure of a Peptide/MHC Complex

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It is regretted that due to a printing error Table 2 of the above paper contained the following errors:

Line 12:	$N^{\delta1}\rightarrow N^{\varepsilon2}$ Gln97	should read:	$O^{\delta1}\rightarrow N^{\varepsilon2}$ Gln97
Line 13:	$N^{\delta2}\rightarrow N^{\varepsilon1}$ Gln97	should read:	$N^{\delta2}\rightarrow O^{\varepsilon1}$ Gln97
Line 20:	N→1 O^{γ} Ser77	should read:	N→O^{γ} Ser77

The Table appears correctly below.

Table 2. Solvent accessibility of peptide residues and interaction partners

Residue	H-bonds[a]	Hydrophobic interactions[a,b]	Accesible surface: Side-chain atoms[c] (Å^2)	Accesible surface: Main-chain atoms[c] (Å^2)	Accesible surface: % of total[c,d]
P1Lys	N→OH Tyr171 O→OH Tyr159	Glu63 Trp167	30.9	0.0	10.6
P2Ala	N→$O^{\varepsilon1}$ Glu63 O→N^{ζ} Lys66	Tyr7 Lys66	2.3	1.0	1.1
P3Val	O→$O^{\varepsilon2}$ Glu9[e]	Tyr159	0.8	0.0	0.3
P4Tyr	O→$N^{\varepsilon2}$ His155	Lys66	112.0	2.4	39.3
P5Asn	N → $O^{\varepsilon1}$ Gln70 $O^{\delta1}\rightarrow N^{\varepsilon2}$ Gln97 $N^{\delta2}\rightarrow O^{\varepsilon1}$ Gln97	Trp73	0.0	12.8	4.4
P6Phe	N→OH Tyr156 O→$N^{\varepsilon1}$ Trp73	His155 Ala152	62.9	15.2	26.8
P7Ala	O→OH Tyr156[e]	Trp147	10.2	10.0	6.9
P8Thr	O→$N^{\varepsilon1}$ Trp147 O→N^{ζ} Lys146 $O^{\gamma1}\rightarrow N^{\zeta}$ Lys146	Trp73	26.5	3.9	10.5
P9Cys	N→O^{γ} Ser77 OT1→$O^{\gamma1}$ Thr143 OT1→OH Tyr84 OT2→$N^{\delta2}$ Asn80 OT2→N^{ζ} Lys146	Trp73 Ser77	0.0	0.0	0.0

[a] As identified by the program HBPLUS (McDonald & Thornton, 1994), using default parameters.
[b] Only for side-chains, and defined as carbon or sulfur atoms in contact with other carbon or sulfur atoms.
[c] Calculated using the program Surface (Collaborative Computational Project Number 4, 1994), using a 1.4 Å probe.
[d] Ratio between the accessible surface of a whole residue and the total accessible surface of the peptide.
[e] Hydrogen bonds mediated by an ordered water molecule.

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