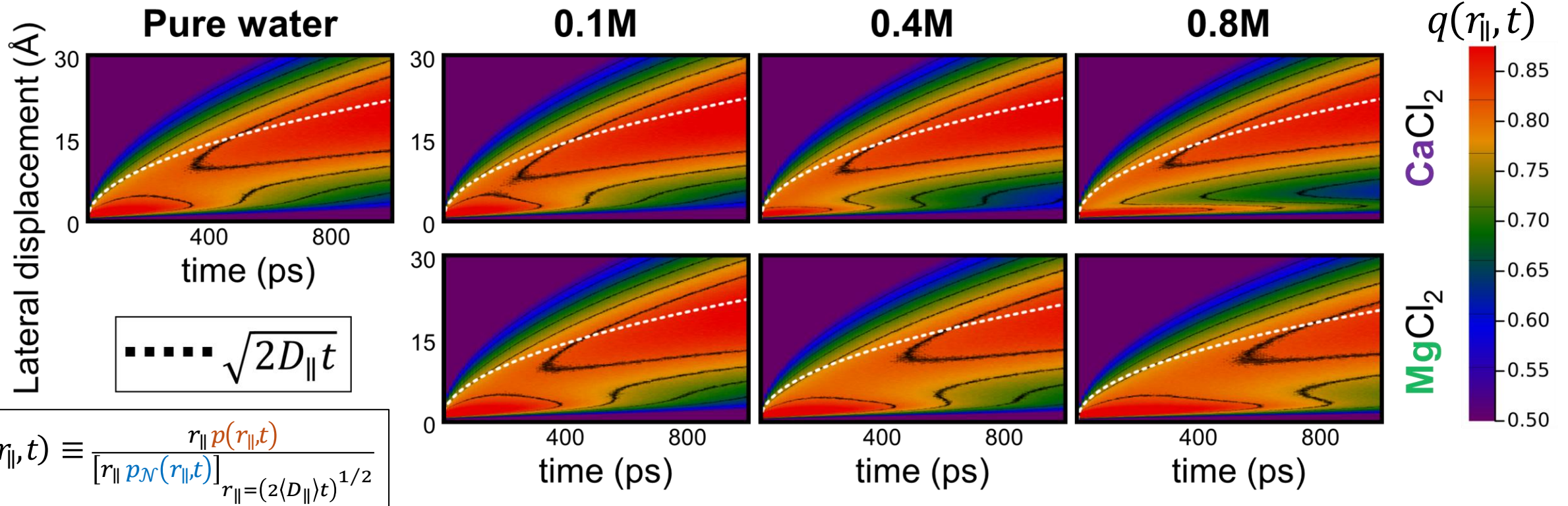


SI Result 1. Lateral Displacement Distribution

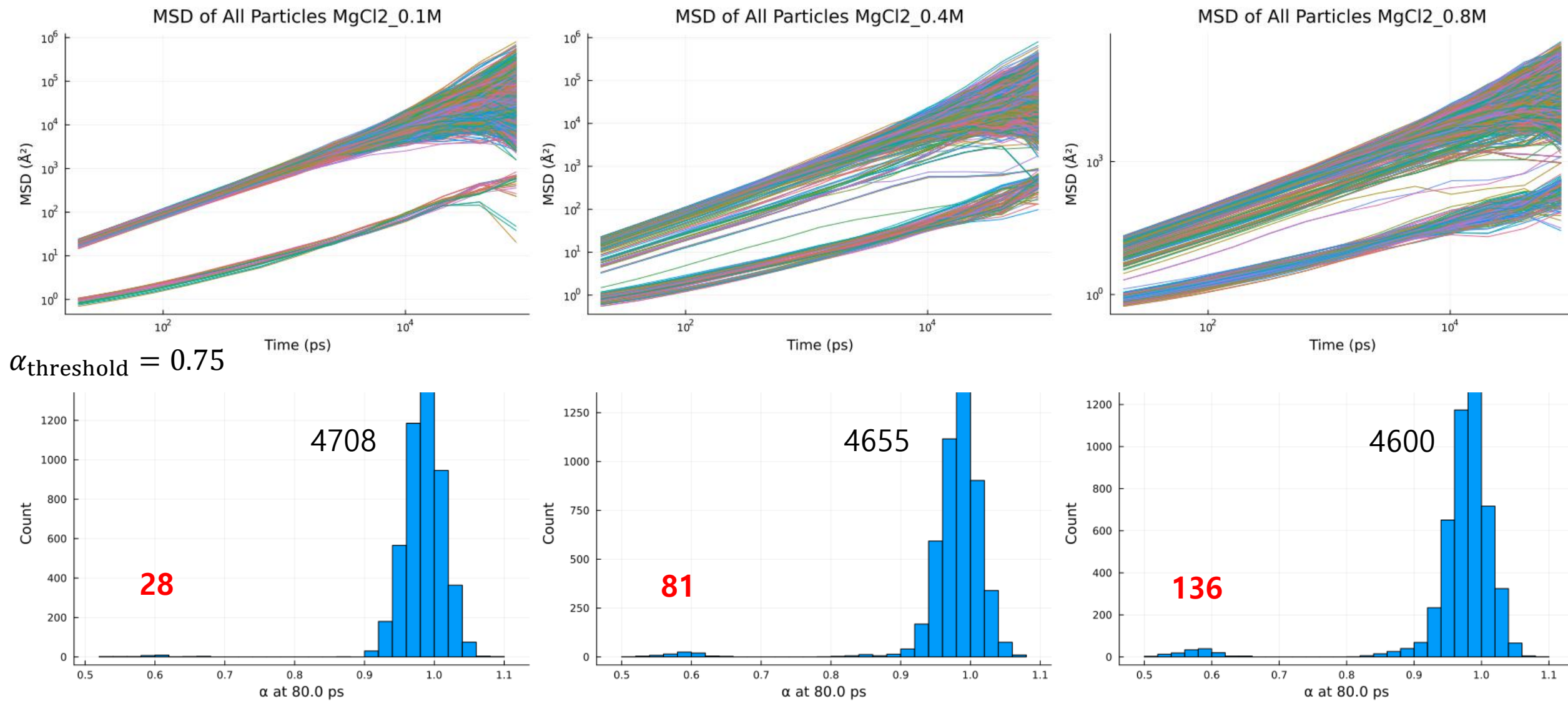


$$q(r_{\parallel}, t) \equiv \frac{r_{\parallel} p(r_{\parallel}, t)}{[r_{\parallel} p_{\mathcal{N}}(r_{\parallel}, t)]_{r_{\parallel} = (2\langle D_{\parallel} \rangle t)^{1/2}}}$$

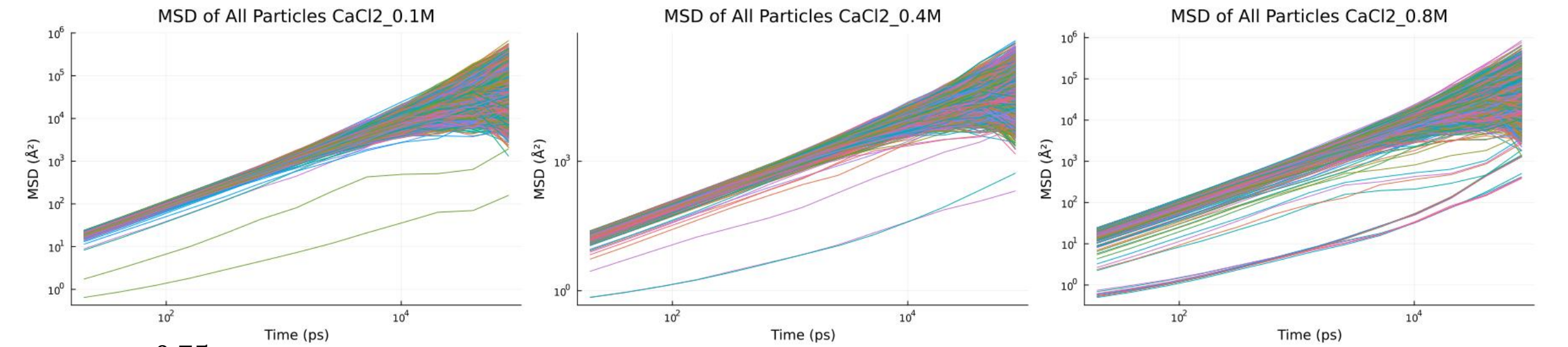
$p(r_{\parallel}, t)$: Radial distribution (simulation result) $p_{\mathcal{N}}(r_{\parallel}, t) \equiv (4\pi\langle D_{\parallel} \rangle t)^{-1} e^{-r_{\parallel}^2/4\langle D_{\parallel} \rangle t}$: Radial Gaussian distribution

- The peak position (**red**) = where the displacement (r_{\parallel}) is biased.
- At long times, peak position of $q(r_{\parallel}, t)$ converges to $\sqrt{2D_{\parallel}t}$. = At long times, $p(r_{\parallel}, t) \cong p_{\mathcal{N}}(r_{\parallel}, t)$.
- **CaCl₂**: As Conc. Increases, short r_{\parallel} region expands along the x-axis, while it shrinks along the y-axis.
- **MgCl₂**: As Conc. Increases, short r_{\parallel} region expands along the x-axis.
- **Both**: As Conc. Increases, the appearance of the peak at long r_{\parallel} is retarded in time.

SI Result 2. MSD of individual water molecules.



SI Result 2. MSD of individual water molecules.



$\alpha_{\text{threshold}} = 0.75$

