

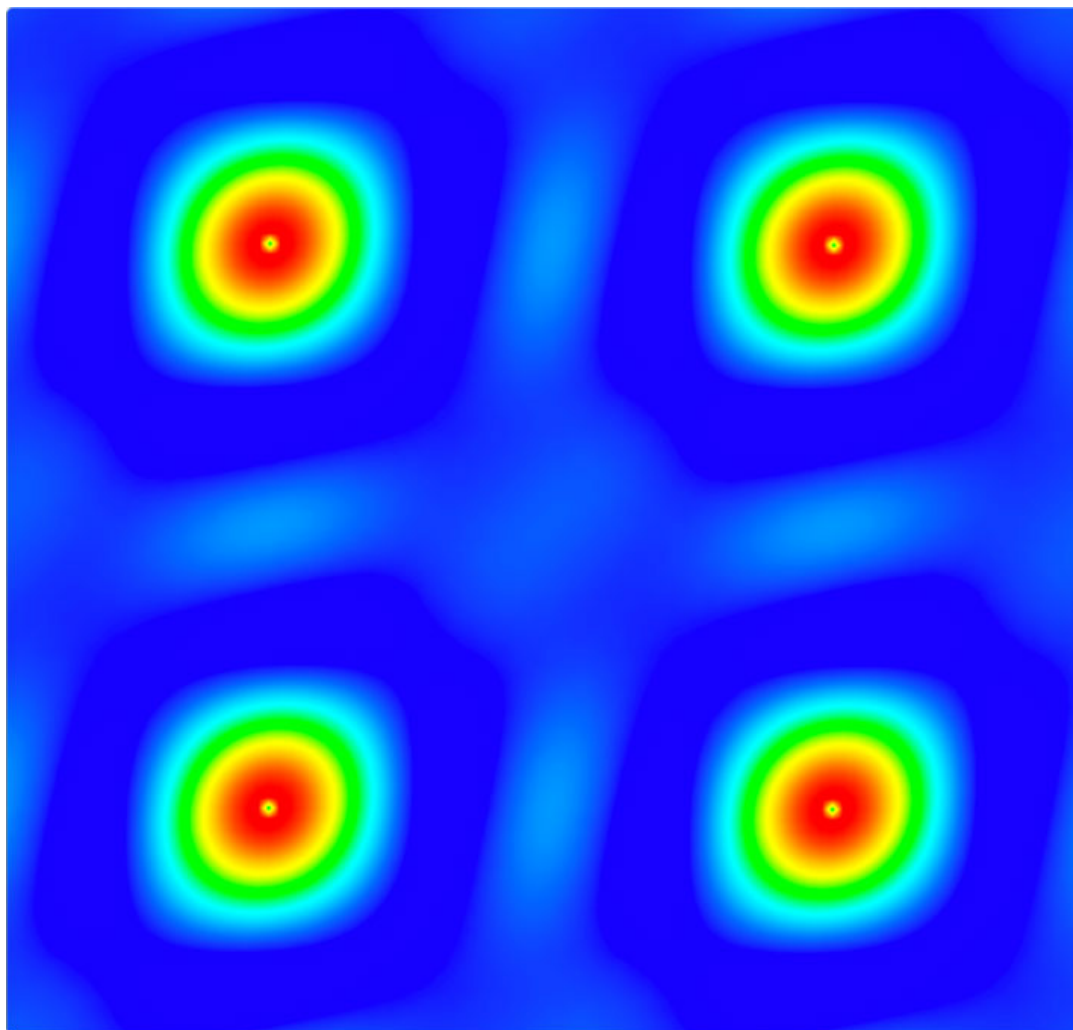


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Simulated electron density for a simple cubic crystal with cell parameter 3 Å.

This sequence shows the effect of increasing the resolution from 3 Å (separate atoms barely resolvable) down to a fraction of an Ångstrom. Note the prominent fringes which become weaker as the resolution increases.

N.B., at 0.1 Å we start to see moiré effects due to limited sampling ("aliasing").

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