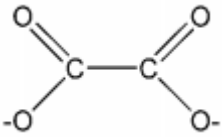


Question number	Answer	Marks	Guidance
1 (a)	$1s^2 2s^2 2p^6 3s^2 3p^6 3d^{10}$	1	allow [He] $2s^2$ , or [Ne] $3s^2$ , or [Ar] $3d^{10}$
	d-sub-shell / shell / orbitals / sub-level full (or not partially full)	1	can only score M2 if $d^{10}$ in M1 correct allow 'full d-orbital' if $d^{10}$ in M1 do not allow d-block
1 (b)	atom or ion or transition metal bonded to / surrounded by one or more ligands	1	Allow Lewis base instead of ligand
	by co-ordinate / dative (covalent) bonds / donation of an electron pair	1	can only score M2 if M1 correct
1 (c)	$H_2$ / hydrogen	1	do not allow H
	no lone / spare / non-bonded pair of electrons	1	only score M2 if M1 correct or give 'H' in M1
1 (d) (i)	+2 OR 2+ OR $Pd^{2+}$ OR II OR +II OR II+ OR two OR two plus	1	
1 (d) (ii)	Tetrahedral	1	these shapes can be in any order allow phonetic spelling, e.g., tetrahydal
	square planar	1	
2 (a) (i)	absorbs (certain frequencies of) (white) light / photons	1	not absorbs white / u.v. light
	d-electrons excited / promoted	1	or d-electrons move between levels / orbitals d-electrons can be implied elsewhere in answer
	the colour observed is the light not absorbed / light reflected / light transmitted	1	allow blue light transmitted penalise emission of light in M3

2 (a) (ii)	<p><math>E</math> is the energy gained by the (excited) electrons (of <math>\text{Cu}^{2+}</math>)</p> <p><math>h</math> (Planck's) constant</p> <p><math>\nu</math> frequency of light (absorbed by <math>\text{Cu}^{2+}(\text{aq})</math>)</p>	1  1  1	<p>allow:</p> <ul style="list-style-type: none"> <li>energy difference between orbitals / subshells</li> <li>energy of photon / light absorbed</li> <li>change in energy of the electrons</li> <li>energy lost by excited electrons</li> <li>energy of photon / light emitted</li> </ul> <p>do not allow wavelength</p> <p>If energy lost / photon lost / light emitted in M1 do not penalised light emitted</p>
2 (a) (iii)	<p><math>[\text{Cu}(\text{H}_2\text{O})_6]^{2+} + 4\text{Cl}^- \rightarrow [\text{CuCl}_4]^{2-} + 6\text{H}_2\text{O}</math></p> <p>Tetrahedral</p> <p><math>\text{Cl}^-</math> / Cl / chlorine too big (to fit more than 4 round Cu)</p>	1  1  1	<p>note that <math>[\text{CuCl}_4]^{2-}</math> is incorrect penalise charges shown separately on the ligand and overall penalise HCl</p> <p>allow water smaller than <math>\text{Cl}^-</math> explanation that change in shape is due to change in co-ordination number</p>
2 (b)	 <p>lone pair(s) on <math>\text{O}^-</math> / O</p>	1    1	<p>allow:</p> <ul style="list-style-type: none"> <li>ion drawn with any bond angles</li> <li>ion in square brackets with overall / 2- charge shown outside the brackets</li> <li>ion with delocalised <math>\text{O}=\text{C}-\text{O}</math> bonds in carboxylate group(s)</li> </ul> <p>allow position of lone pair(s) shown on O in the diagram even if the diagram is incorrect.</p>

