## Transition Metal Basics

## Mark Scheme

| Level | International A Level |
| :--- | :--- |
| Subject | Chemistry |
| Exam Board | Edexcel |
| Topic | Transition Metals \& Organic Nitrogen Chemistry |
| Sub Topic | Transition Metal Basics |
| Booklet | Mark Scheme |


| Time Allowed: | 86 minutes |
| :--- | :---: |
| Score: | $/ 71$ |
| Percentage: | $/ 100$ |

Grade Boundaries:

| A* | A | B | C | D | E | U |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| $>85 \%$ | $77.5 \%$ | $70 \%$ | $62.5 \%$ | $57.5 \%$ | $45 \%$ | $<45 \%$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1}$ | B |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2}$ | A |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{3}$ | D |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{4}$ | A |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{5}$ | D |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{6}$ | C |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{7}$ | B |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{8}$ | C |  | 1 |
| Question <br> Number Correct Answer Reject Mark <br> $\mathbf{9}$ B  1 |  |  |  |


| Question <br> Number | Correct Answer | Reject | Mark |
| ---: | :--- | :--- | :--- |
| $\mathbf{1 0}$ | B |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 1}$ B |  | $\mathbf{1}$ |  |


| Question <br> Number | Correct Answer | Reject | Mark |
| ---: | :--- | :--- | :---: |
| $\mathbf{1 2}$ | A |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| ---: | :--- | :--- | :---: |
| $\mathbf{1 3}$ | C |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| ---: | :--- | :--- | :---: |
| $\mathbf{1 4}$ | D |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| ---: | :--- | :--- | :---: |
| $\mathbf{1 5}$ | A |  | $\mathbf{1}$ |

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( a ) ( i i )}$ | If name and formula are given, both <br> must be correct |  |  |
|  | $\mathbf{X}=$ (aqueous) ammonia / $\mathrm{NH}_{3}(\mathrm{aq})$ <br> ALLOW <br> $\mathrm{NH}_{3} /$ ammonium hydroxide (1) <br> $\mathbf{Y =}$ potassium iodide / KI <br> ALLOW <br> other soluble iodides <br> IGNORE references to concentration | $\mathbf{X}=\mathrm{NaOH}$ | iodide / I- <br> KI and acid <br> HI |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( a ) ( i i i )}$ | (Product is) ethanoic acid $/ \mathrm{CH}_{3} \mathrm{COOH} /$ <br> ethanoate( ions) $/ \mathrm{CH}_{3} \mathrm{COO}^{-} \quad(1)$ <br> IGNORE carboxylic <br> Ethanal is a reducing agent / reduces <br> $\mathrm{Cu}^{2+}$ <br> Stand alone marks <br> IGNORE <br> references to oxidation of ethanol <br> products of reduction (e.g. Cu) |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( a ) ( i v )}$ | (Iodine is formed quantitatively and is <br> determined by) titration against sodium <br> thiosulfate solution (of known <br> concentration) | Colorimetry | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 6 ( b ) ( i )}$ | (3)d orbitals / (3)d subshell split (by the (1) <br> attached ligands) <br> Electrons are promoted (from lower to <br> higher energy d orbital(s) / levels) <br> OR <br> Electrons move from lower to higher split <br> energy d orbital(s) / levels) <br> ALLOW <br> d-d transitions occur |  |  |
|  | Absorbing energy /photons of a certain <br> frequency (in the visible region) <br> ALLOW <br> Absorbing light <br> Reflected / transmitted / remaining light is <br> coloured / yellow / in the visible region <br> ALLOW (1) <br> Complementary colour seen <br> Reflected / transmitted / remaining light / <br> frequency is seen <br> Penalise omission of (3)d once only. <br> Ignore reference to electrons relaxing / <br> dropping to the ground state |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( b ) ( i i )}$ | Colour depends on the frequency <br> /wavelength /energy of the absorbed (1) <br> light |  |  |
| Different ligands split the d orbitals to <br> a different extent |  | $\mathbf{2}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :---: |
| $\mathbf{1 6 ( c ) ( i )}$ | $2 \mathrm{Cu}^{+}(\mathrm{aq}) \rightarrow \mathrm{Cu}(\mathrm{s})+\mathrm{Cu}^{2+}(\mathrm{aq})$ <br> ALLOW <br> reversible arrows <br> Electrons |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( c ) ( i i )}$ | The copper(I) is oxidized to <br> copper(II) and (in the same reaction) <br> reduced to copper((0)) |  |  |
| OR <br> Copper changes from +1 to 0 and +2 | IGNORE <br> Reference to a Cu atom |  | $\mathbf{1}$ |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 16(c)(iii) | Relevant reduction potentials are $\begin{aligned} & \mathrm{Cu}^{2+}+\mathrm{e}^{-} \rightleftharpoons \mathrm{Cu}^{+} \mathrm{E}^{\ominus}=+0.15(\mathrm{~V}) \\ & \mathrm{Cu}^{+}+\mathrm{e}^{-} \rightleftharpoons \mathrm{Cu} \mathrm{E}^{\ominus}=+0.52(\mathrm{~V}) \end{aligned}$ <br> ALLOW single arrows $\begin{equation*} \mathrm{E}_{\text {cell }}^{\ominus}=0.52-0.15=(+) 0.37(\mathrm{~V}) \tag{1} \end{equation*}$ <br> TE on incorrect $\mathrm{E}^{\ominus}$ values providing $\mathrm{E}^{\ominus}$ cell is positive <br> ( $\mathrm{E}^{\circ}{ }_{\text {cell }}$ positive so reaction thermodynamically favourable) |  | 2 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( a ) ( i )}$ | $3 d^{5} 4 s^{1}$ |  |  |
|  | $/ 4 s^{1} 3 d^{5}$ |  | 1 |
|  | ALLOW |  |  |
|  | Complete configuration <br> $1 s^{2} 2 s^{2} 2 p^{6} 3 s^{2} 3 p^{6} 4 s^{1} 3 d^{5}$ |  |  |
|  | ALLOW |  |  |
|  | Capitals and subscripts |  |  |


| Question Number | Correct Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{aligned} & 17 \\ & \text { (a) (ii) } \end{aligned}$ | It is $4 s^{1}$ rather than $4 s^{2}$ because with two of the reasons below <br> $3 d^{5} /$ half- filled $3 d$ sub shell is particularly stable <br> The paired electrons repel <br> All six electrons are in separate orbitals (minimizing repulsion) <br> ALLOW <br> The energy required to promote/ transfer 4s to 3 d is small OR <br> The energy difference between 4 s and $3 d$ is small |  | 2 |

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| Question Number | Correct Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(b)(i) | $\begin{aligned} & \left(\mathrm{E}^{\ominus} \mathrm{Zn}^{2+}(\mathrm{aq}) \mid \mathrm{Zn}(\mathrm{~s})=-0.76 \mathrm{~V}\right. \\ & \mathrm{E}^{\ominus} \mathrm{Cr}^{3+}(\mathrm{aq}), \mathrm{Cr}^{2+}(\mathrm{aq}) \mid \mathrm{Pt}=-0.41 \mathrm{~V} \\ & \mathrm{E}^{\ominus}\left[\mathrm{Cr}_{2} \mathrm{O}_{7}^{2-}(\mathrm{aq})+7 \mathrm{H}^{+}(\mathrm{aq})\right], \\ & \left.\left[2 \mathrm{Cr}^{3+}(\mathrm{aq})+7 \mathrm{H}_{2} \mathrm{O}(\mathrm{I})\right] \mid \mathrm{Pt}=+1.33 \mathrm{~V}\right) \end{aligned}$ <br> If no other mark is scored, data scores (1) however shown <br> Calculation of $E^{\ominus}$ cell values: <br> $\mathrm{E}^{\ominus}{ }_{\text {cell }}$ for first step $=$ $\begin{equation*} 1.33--0.76=(+) 2.09(\mathrm{~V}) \tag{1} \end{equation*}$ <br> $\mathrm{E}^{\ominus}{ }_{\text {cell }}$ for second step $=$ $\begin{equation*} -0.41--0.76=(+) 0.35(\mathrm{~V}) \tag{1} \end{equation*}$ <br> As (both) values are positive, (both) reactions are spontaneous/feasible <br> Third mark is independent |  | 3 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( b ) ( i i )}$ | Orange to green to blue <br> IGNORE qualifying words eg pale blue |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7}$ | The small amount of hydrogen <br> (b)(iii) <br> produced (does not present a serious <br> risk) |  | 1 |
| ALLOW <br> "Less" for small amount <br> Indication of ventilation |  |  |  |

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| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( c ) ( i )}$ | It is bridging/ bidentate ligand | Polydentate | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( c ) ( i i )}$ | Dative (covalent) (bonds)/ <br> co-ordinate (bonds) | 1 |  |


| Question Number | Correct Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{aligned} & 17 \\ & \text { (c) (iii) } \end{aligned}$ | Any two from: <br> Chromium atoms/ ions are covalently bonded/bonded to each other <br> OR <br> Two (chromium) ions/ chromium atoms in the complex <br> Each ethanoate ligand forms bonds to two different atoms/ ions <br> Ethanoate ions are not normally bidentate ligands <br> ALLOW <br> Contains both monodentate and bidentate ligands <br> Allow six ligands and complex not octahedral | Just "two different ligands" | 2 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7}$ <br> $\mathbf{( c ) ( i v ) ~}$ | The energies of the d electron levels <br> are split to different extents (by <br> different ligands) |  | 2 |
|  | ALLOW <br> d-d (orbitals) splitting is different <br> OR <br> d-d transitions are different | (1) |  |
| So different energy/ frequency/ <br> wavelength light absorbed | (1) | (...(just) <br> transmitted |  |

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| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( c ) ( v )}$ | There are two peaks as two different <br> hydrogen environments (1) |  | 2 |
|  | ElTHER <br> The areas due to hydrogen in water <br> molecules compared to hydrogen in <br> ethanoate ions is in the ratio 1 to 3/ <br> 4 to 12 <br> OR <br> As there are 4 hydrogen atoms in <br> water and 12 hydrogen atoms in <br> ethanoate ions |  |  |


| Question Number | Correct Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(d) | First mark <br> Dilution factor: <br> moles of chromium (II) ethanoate in $25.0 \mathrm{~cm}^{3}$ $\begin{equation*} =\frac{2.66 \times 10^{-3}}{10}=2.66 \times 10^{-4} \tag{1} \end{equation*}$ <br> Second mark <br> Ratio of manganate(VII) to chromium <br> 4 mol manganate(VII) react with 5 mol of chromium (II) <br> OR <br> 8 mol mangante(VII) react with 5 mol of chromium(II) ethanoate <br> Third mark <br> moles of manganate(VII) ion $\begin{align*} & =\frac{4 \times 5.32 \times 10^{-4}}{5} \text { OR } \frac{8 \times 2.66 \times 10^{-4}}{5} \\ & =4.256 \times 10^{-4} \tag{1} \end{align*}$ <br> Fourth mark <br> Volume of manganate(VII) solution $\begin{align*} & =\frac{4.256 \times 10^{-4}}{0.00750} \times 1000 \\ & =56.75 \mathrm{~cm}^{3} \tag{1} \end{align*}$ <br> Correct answer no working (4) <br> $28.375 \mathrm{~cm}^{3}$ gets (3) <br> Fifth mark <br> This is unsuitable/ inaccurate because it requires refilling the burette hence increasing burette error <br> OR <br> Better to use more concentrated potassium manganate(VII) OR less chromium ethanoate |  | 5 |

(Total for Question 17 = 21 marks)

| Question <br> Number | Acceptable Answer | Reject | Mark |  |
| :---: | :--- | :---: | :---: | :---: |
| $\mathbf{1 8 ( a ) ( i )}$ | $\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}$ <br> $\mathrm{ALLOW}\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{4}\right]^{2+}$ <br> $\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{4}(\mathrm{OH})_{2}$ <br> $\mathrm{ALLOW} \mathrm{Cu}(\mathrm{OH})_{2}$ | (1) | $\mathrm{Cu}^{2+}(\mathrm{aq)}$ | 3 |
|  | $\left[\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{4}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2}\right]^{2+}$ <br> ALLOW $\left[\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{4}\right]^{2+}$ | (1) |  |  |
| ALLOW <br> Ligand in any order <br> Omission of square brackets | $\left[\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{6}\right]^{2+}$ |  |  |  |


| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(a)(ii) | (3)d orbitals / (3)d subshell split <br> (by the attached ligands) <br> Electrons are promoted (from lower to higher energy d orbital(s) / levels) <br> OR <br> Electrons move from lower to higher energy (d orbital(s) / levels) ALLOW <br> d-d transitions occur /electrons are excited <br> (1) <br> Absorbing energy /photons of a certain frequency (in the visible region) <br> ALLOW <br> Absorbing light <br> Reflected / transmitted / remaining light is coloured / in the visible region <br> ALLOW <br> Complementary colour seen <br> Reflected / transmitted / remaining <br> light / frequency is seen <br> Penalise omission of (3)d once only. Ignore reference to electrons relaxing / dropping to the ground state | Orbital / shell / subshells split d-d splitting <br> Emitted <br> 'Reverse' for 'complementary' | 4 |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :---: | :--- | :--- | :---: |
| 18(a)(iii) | The (different) ligands split the (3)d <br> orbitals / subshell to a different <br> extent | Orbital / shell / <br> subshells unless <br> penalised in <br> 22(a)(ii) | 2 |
|  | (So) the energy absorbed / reflected <br> ltransmitted is different <br> OR | Emitted unless <br> Radiation (ALLOW light) is at a <br> penalised in <br> different frequency | (1) |



Total for Question 18 = 14 marks

