## Cambridge International AS \& A Level

## CHEMISTRY

## Published

This mark scheme is published as an aid to teachers and candidates, to indicate the requirements of the examination. It shows the basis on which Examiners were instructed to award marks. It does not indicate the details of the discussions that took place at an Examiners' meeting before marking began, which would have considered the acceptability of alternative answers.

Mark schemes should be read in conjunction with the question paper and the Principal Examiner Report for Teachers.

Cambridge International will not enter into discussions about these mark schemes
Cambridge International is publishing the mark schemes for the October/November 2023 series for most Cambridge IGCSE, Cambridge International A and AS Level components, and some Cambridge O Level components.

## Generic Marking Principles

These general marking principles must be applied by all examiners when marking candidate answers. They should be applied alongside the specific content of the mark scheme or generic level descriptors for a question. Each question paper and mark scheme will also comply with these marking principles.

## GENERIC MARKING PRINCIPLE 1:

Marks must be awarded in line with:

- the specific content of the mark scheme or the generic level descriptors for the question
- the specific skills defined in the mark scheme or in the generic level descriptors for the question
- the standard of response required by a candidate as exemplified by the standardisation scripts.

GENERIC MARKING PRINCIPLE 2 :
Marks awarded are always whole marks (not half marks, or other fractions).

## GENERIC MARKING PRINCIPLE 3:

Marks must be awarded positively:

- marks are awarded for correct/valid answers, as defined in the mark scheme. However, credit is given for valid answers which go beyond the scope of the syllabus and mark scheme, referring to your Team Leader as appropriate
- marks are awarded when candidates clearly demonstrate what they know and can do
- marks are not deducted for errors
- marks are not deducted for omissions
- answers should only be judged on the quality of spelling, punctuation and grammar when these features are specifically assessed by the question as indicated by the mark scheme. The meaning, however, should be unambiguous.


## GENERIC MARKING PRINCIPLE 4:

Rules must be applied consistently, e.g. in situations where candidates have not followed instructions or in the application of generic level descriptors.

## GENERIC MARKING PRINCIPLE 5:

Marks should be awarded using the full range of marks defined in the mark scheme for the question (however; the use of the full mark range may be limited according to the quality of the candidate responses seen).

## GENERIC MARKING PRINCIPLE 6:

Marks awarded are based solely on the requirements as defined in the mark scheme. Marks should not be awarded with grade thresholds or grade descriptors in mind.

## Science-Specific Marking Principles

1 Examiners should consider the context and scientific use of any keywords when awarding marks. Although keywords may be present, marks should not be awarded if the keywords are used incorrectly.

2 The examiner should not choose between contradictory statements given in the same question part, and credit should not be awarded for any correct statement that is contradicted within the same question part. Wrong science that is irrelevant to the question should be ignored.

3 Although spellings do not have to be correct, spellings of syllabus terms must allow for clear and unambiguous separation from other syllabus terms with which they may be confused (e.g. ethane / ethene, glucagon / glycogen, refraction / reflection).

4 The error carried forward (ecf) principle should be applied, where appropriate. If an incorrect answer is subsequently used in a scientifically correct way, the candidate should be awarded these subsequent marking points. Further guidance will be included in the mark scheme where necessary and any exceptions to this general principle will be noted.

## 5 'List rule' guidance

For questions that require $\boldsymbol{n}$ responses (e.g. State two reasons ...):

- The response should be read as continuous prose, even when numbered answer spaces are provided.
- Any response marked ignore in the mark scheme should not count towards $\boldsymbol{n}$.
- Incorrect responses should not be awarded credit but will still count towards $\boldsymbol{n}$.
- Read the entire response to check for any responses that contradict those that would otherwise be credited. Credit should not be awarded for any responses that are contradicted within the rest of the response. Where two responses contradict one another, this should be treated as a single incorrect response.
- Non-contradictory responses after the first $\boldsymbol{n}$ responses may be ignored even if they include incorrect science.


## 6 Calculation specific guidance

Correct answers to calculations should be given full credit even if there is no working or incorrect working, unless the question states 'show your working'.

For questions in which the number of significant figures required is not stated, credit should be awarded for correct answers when rounded by the examiner to the number of significant figures given in the mark scheme. This may not apply to measured values.

For answers given in standard form (e.g. $a \times 10^{\eta}$ ) in which the convention of restricting the value of the coefficient (a) to a value between 1 and 10 is not followed, credit may still be awarded if the answer can be converted to the answer given in the mark scheme.

Unless a separate mark is given for a unit, a missing or incorrect unit will normally mean that the final calculation mark is not awarded. Exceptions to this general principle will be noted in the mark scheme.

7 Guidance for chemical equations
Multiples / fractions of coefficients used in chemical equations are acceptable unless stated otherwise in the mark scheme.
State symbols given in an equation should be ignored unless asked for in the question or stated otherwise in the mark scheme.

| Question | Answer | Marks |
| :---: | :---: | :---: |
| 1(a)(i) | $\begin{aligned} & \mathrm{M} 1: \mathrm{F}_{2}+\mathrm{ClO}_{2} \rightarrow \mathrm{FClO}_{2}+\mathrm{F}[1] \\ & \mathrm{M} 2: \mathrm{ClO}_{2}+\mathrm{F} \rightarrow \mathrm{FClO}_{2}[1] \end{aligned}$ <br> OR <br> M1: $\mathrm{F}_{2}+\mathrm{ClO}_{2} \rightarrow \mathrm{~F}_{2} \mathrm{ClO}_{2}$ <br> M2: $\mathrm{F}_{2} \mathrm{ClO}_{2}+\mathrm{ClO}_{2} \rightarrow 2 \mathrm{FClO}_{2}$ <br> two balanced equations MUST add to give the overall equation | 2 |
| 1(a)(ii) | first step AND <br> has one mole / molecule of $\mathrm{F}_{2}$ and $\mathrm{ClO}_{2}$ [1] <br> OR same moles of reactants as orders in rate equation | 1 |
| 1(b)(i) | $\begin{aligned} & \text { rate }=k\left[\mathrm{~F}_{2}\right]\left[\mathrm{ClO} \mathrm{O}_{2}\right] / \text { rate }=1.22\left[\mathrm{~F}_{2}\right][\mathrm{ClO} \\ & \text { AND second } / 2 \mathrm{BOTH}[1] \end{aligned}$ | 1 |
| 1(b)(ii) | $\begin{aligned} & \text { rate }=1.22 \times 2 \times 10^{-3} \times 2 \times 10^{-3}=4.88 \times 10^{-6}[1] \\ & \min 2 \operatorname{sfecf} 1(\text { b) (i) } \end{aligned}$ | 1 |
| 1(c)(i) | $k_{1}=(0.693 / 4)=0.173 \quad \text { OR } 1.73 \times 10^{-1}[1]$ <br> MUST BE 3SF $\mathrm{s}^{-1}[1]$ | 2 |
| 1(c)(ii) | at $4 \mathrm{~s}=0.001$ <br> at $8 \mathrm{~s}=0.0005$ <br> at $12 \mathrm{~s}=0.00025$ <br> smooth curve <br> ALL correct [1] | 1 |
| 1(c)(iii) | tangent drawn at $0.00100 \mathrm{~mol} \mathrm{dm}^{-3}$ <br> AND <br> gradient dependent on their rate $=\mathrm{Y} / \mathrm{X}$ in the range $=1.5-2.0 \times 10^{-4}[1]$ IGNORE sign | 1 |


| Question | Answer | Marks |
| :---: | :---: | :---: |
| 2(a) | $\left(K_{w}=\right)[\mathrm{H}+][\mathrm{OH}-] \mathrm{OR}\left(K_{w}=\right)[\mathrm{H} 3 \mathrm{O}+][\mathrm{OH}-][1]$ | 1 |
| 2(b)(i) | M1: pH values 1.5 AND 12.5 [1] <br> M2: conc of $\mathrm{HCl}=10-1.5=0.0316$ ecf [1] $\min 2 \mathrm{sf}$ | 2 |
| 2(b)(ii) | sodium chloride / NaCl AND 7 [1] | 1 |
| 2(b)(iii) | (mixture Y) 1 to 3 AND (mixture Z) 11 to 13 [1] | 1 |
| 2(c)(i) | $\mathrm{H}_{2} \mathrm{SO}_{4}>\mathrm{CH}_{3} \mathrm{CCl}_{2} \mathrm{COOH}>\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}[1] \mathrm{u} / \mathrm{c}$ <br> explanation <br> - $\mathrm{H}_{2} \mathrm{SO}_{4}$ is fully dissociated / strong acid <br> - $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH} / \mathrm{CH}_{3} \mathrm{CCl}_{2} \mathrm{COOH}$ are partly dissociated / weak acids <br> - $\mathrm{Cl} /$ chlorine is electron-withdrawing / electronegative <br> - alkyl group (in CH 3 CH 2 COOH ) is electron donating <br> - stabilises / destabilises anion OR weakens / strengthens O-H bond (linked correctly) <br> - correct reference to release of / donation of/ form $\mathrm{H}+$ / proton <br> any two [1] any four [2] | 3 |
| 2(c)(ii) | - $\mathrm{H}_{3} \mathrm{O}+$ and $\mathrm{H}_{2} \mathrm{O}$ <br> - $\mathrm{H}_{2} \mathrm{SO}_{4}$ and $\mathrm{HSO}_{4}^{-}$ <br> - $\mathrm{HSO}_{4}^{-}$and $\mathrm{SO}_{4}{ }^{2-}$ <br> any one correct pair [1] <br> all three correct pairs [2] | 2 |
| 2(d)(i) | (the) ratio of the concentrations (of a solute between) two solvents / two liquids (at) equilibrium [1] | 1 |
| 2(d)(ii) | $\begin{aligned} & \text { mass }=5-x / x=7.84 \text { OR mass }=x / 5-x=7.84 \\ & \text { mass }=0.5656 \mathrm{~g} \text { mass }=4.43 \mathrm{~g} \\ & \text { [1] min 2sf } \end{aligned}$ | 1 |


| Question | Answer | Marks |
| :---: | :--- | :---: |
| 2(d)(iii) | Any numbers in which volume of water is 7.84 times volume of hexane e.g. $78.4 \mathrm{~cm}^{3}$ water $10 \mathrm{~cm}^{3}$ hexane [1] u/c <br> e.g mass $=0.566 \mathrm{~g} \mathrm{177.4cm}^{3}$ water $22.6 \mathrm{~cm}^{3}$ hexane <br> ALLOW reverse ratio volume of hexane is 7.84 times volume of water when consistent with d(ii) | $\mathbf{1}$ |
| 2(d)(iv) | Q is $\mathrm{CH}_{3}\left(\mathrm{CH}_{2}\right)_{4} \mathrm{OH}$ <br> AND it is least polar <br> / contains a large non-polar hydrocarbon chain <br> /stronger id-id forces with hexane in Q OWTTE <br> $[1]$ | $\mathbf{1}$ |


| Question | Answer | Marks |
| :---: | :---: | :---: |
| 3(a)(i) | (number of possible) arrangements of particles / energy in a system OR measure / degree of disorder in / of a system [1] | 1 |
| 3(a)(ii) | positive / + <br> AND more gas molecules / particles in products OR more moles / molecules in products / RHS [1] | 1 |
| 3(b) | $\begin{aligned} & \Delta H=(2 \times 150)-(1 \times 496)(=-196)[1] \\ & \text { OR } \\ & \Delta H=(2 \times 150)+(4 \times 460)-(1 \times 496)-(4 \times 460)(=-196) \end{aligned}$ | 1 |
| 3(c) | $\begin{aligned} & \Delta G=\Delta H-\mathrm{T} \Delta S \text { seen or used with correct signs[1] } \\ & -238=-196-298 \Delta S \\ & \Delta S=42 / 298 \\ & \Delta S=(+) 0.141 / 0.1409(\mathrm{~kJ} \mathrm{~K}-1 \mathrm{~mol}-1) \\ & \quad \text { OR }(+) 141 / 140.9(\mathrm{~J} \mathrm{~K}-1 \mathrm{~mol}-1)[1] \text { ecf from } \Delta \mathrm{G}=\Delta \mathrm{H}+\mathrm{T} \Delta \mathrm{~S} \\ & 141=(2 \times 70)+\mathrm{S}(\mathrm{O} 2(\mathrm{~g}))-(2 \times 102) \\ & \mathrm{S}, \mathrm{O} 2(\mathrm{~g})=205 / 204.94\left(\mathrm{~J} \mathrm{~K}^{-1} \mathrm{~mol}^{-1}\right)[1] \text { ecf } \end{aligned}$ | 3 |


| Question | Answer | Marks |
| :---: | :---: | :---: |
| 3(d) | iron(III) chloride / $\mathrm{FeCl}^{3}$ <br> AND same state / phase as reactants / $\mathrm{H}_{2} \mathrm{O}_{2}$ [1] | 1 |
| 3(e)(i) | hydrogen peroxide / $\mathrm{H}_{2} \mathrm{O}_{2}$ AND +2.18 [1] | 1 |
| 3(e)(ii) | $\begin{aligned} & \Delta G=-\mathrm{n} E^{\ominus}{ }_{\text {cell }} F[1] \\ & \Delta G=(-2 \times 2.18 \times 96500)=-420.7(\mathrm{~kJ} \mathrm{~mol}-1)[1] \mathrm{ecf} \end{aligned}$ | 2 |
| 3(f)(i) | Nernst: $(E=) E^{9}+(0.059 / z) \log [0 x] /[r e d][1] u / c$ <br> $\mathrm{OR}(E=) E^{\ominus}+(\mathrm{RT} / \mathrm{zF}) \mathrm{ln}[\mathrm{ox}] /[\mathrm{red}]$ <br> $\mathrm{OR}(E=) 1.82+(0.059 / 1) \log (0.02 / 2)$ <br> $E=1.82+(0.059 / 1) \log (0.02 / 2)=(+) 1.702(\mathrm{~V})[1] \mathrm{min} 2 \mathrm{sf}$ <br> OR $E=1.82+[(8.314 \times 298 / 1) \times(96500)] \ln (0.02 / 2)=(+) 1.702$ | 2 |
| 3(f)(ii) | $\mathrm{H}_{2} \mathrm{O}_{2}+2 \mathrm{H}++2 \mathrm{Co}^{2+} \rightarrow 2 \mathrm{H}_{2} \mathrm{O}+2 \mathrm{Co}^{3+}[1]$ <br> ECF for reverse equation from (f)(i) if $E>1.77 \mathrm{~V}$ | 1 |
| 3(g)(i) | enthalpy change when one mole of gaseous ions forms an aqueous solution / dissolves in water [1] | 1 |


| Question | Answer | Marks |
| :---: | :---: | :---: |
| 3(g)(ii) | Substances with state symbols: <br> - AlF3(s) <br> - AlF3(aq) OR Al3+(aq) + (3)F-(aq) <br> - $\mathrm{A} l 3+(\mathrm{g})(3) \mathrm{F}^{-}(\mathrm{g})$ <br> [1] <br> changes identified and three correct arrow directions (one given) [1] | 2 |
| 3(g)(iii) | $\Delta H_{\text {latt }}=-4690+(3 \times-506)-(-209)=-5999(\mathrm{~kJ} \mathrm{~mol}-1)$ | 1 |


| Question | Answer | Marks |
| :---: | :---: | :---: |
| 4(a)(i) | - $\mathrm{Co}(\mathrm{OH})_{2} \mathrm{OR} \mathrm{Co}(\mathrm{OH})_{2}\left(\mathrm{H}_{2} \mathrm{O}\right)_{4}$ <br> - $\left[\mathrm{Co}\left(\mathrm{NH}_{3}\right)_{6}\right]^{2+} \mathrm{OR}\left[\mathrm{Co}\left(\mathrm{NH}_{3}\right)_{6}\right]^{3+}$ <br> - $\left[\mathrm{CoCl}_{4}\right]^{2-}$ <br> any two [1] all three [2] | 2 |
| 4(a)(ii) | pink to blue [1] | 1 |
| 4(b)(i) | $\begin{aligned} & 2 \mathrm{Ca}\left(\mathrm{NO}_{3}\right)_{2} \rightarrow 2 \mathrm{CaO}+4 \mathrm{NO}_{2}+\mathrm{O}_{2} \\ & \mathrm{OR} \mathrm{Ca}\left(\mathrm{NO}_{3}\right)_{2} \rightarrow \mathrm{CaO}+2 \mathrm{NO}_{2}+1 / 2 \mathrm{O}_{2}[1] \end{aligned}$ | 1 |


| Question | Answer | Marks |
| :---: | :---: | :---: |
| 4(b)(ii) | $\mathrm{M} 1: \mathrm{Mg}\left(\mathrm{NO}_{3}\right)_{2}$ below $480^{\circ} \mathrm{C}$, <br> $\mathrm{Ba}\left(\mathrm{NO}_{3}\right)_{2}$ above $520^{\circ} \mathrm{C}$ BOTH [1] u/c <br> M 2 : ionic radii of $\mathrm{M}^{2+}$ increases down the group <br> OR radii of $\mathrm{Ba}^{2+}$ is greater (than $\mathrm{Mg}^{2+}$ ) <br> OR charge density of $\mathrm{M}^{2+}$ decreases down the group <br> OR charge density of $\mathrm{Ba}^{2+}$ is smaller (than $\mathrm{Mg}^{2+}$ ) [1] ORA u/c <br> M3: (larger cations) polarise / distort anion / $\mathrm{NO}_{3}{ }^{-}$less OR <br> (larger ions) weaken $\mathrm{N}-\mathrm{O} / \mathrm{N}=\mathrm{O}$ (bond) less <br> [1] ORA | 3 |


| Question | Answer | Marks |
| :---: | :--- | :---: |
| $5(\mathrm{a})$ | more than one (stable) oxidation state [1] <br> empty / vacant (d) orbitals are energetically accessible <br> OR empty / vacant (d) orbitals can form dative bonds with ligands <br> [1] | $\mathbf{2}$ |
| $5(\mathrm{~b})(\mathrm{i})$ | oxid. no. $=+4$ AND <br> coord. no. $=8[1]$ | $\mathbf{1}$ |


| Question | Answer | Marks |
| :---: | :---: | :---: |
| 5(b)(ii) | bond angle must go from bond to bond OR CN group to CN group diagram <br> - for Ag: NC—Ag-CN <br> - angle $=$ labelled $180^{\circ}$ | 2 |
| 5(c)(i) | $K_{\text {stab }}=\left[\left[\mathrm{Cu}(\mathrm{CN})_{4}\right]^{3-}\right] /\left[\mathrm{Cu}^{+}\right]\left[\mathrm{CN}^{-}\right]^{4}[1]$ <br> ALLOW use of $\left[\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{+}\right]$ | 1 |
| 5(c)(ii) | $[\mathrm{Cu}+]=5.0 \times 10^{-19}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)[1] \mathrm{min} 1 \mathrm{sf}$ ECF from reversed ratio 5(c)(i) | 1 |
| 5(d)(i) | mol of $\mathrm{I}_{2}=0.5 \times 0.02 \times 20.1 / 1000=2.01 \times 10^{-4}$ [1] min 2 sf | 1 |
| 5(d)(ii) | mol of $\mathrm{Cu}=2.01 \times 10^{-4} \times 2 \times 4=1.608 / 1.61 \times 10^{-3}$ [1] ecf min 2 sf | 1 |
| 5(d)(iii) | $\%$ of $\mathrm{Cu}=100 \times\left(1.608 \times 10^{-3} \times 63.5\right) / 0.567=18.0$ [1] ecf | 1 |
| 5(d)(iv) | $\mathrm{Cu}^{2+}$ is $\mathrm{d}^{9} / \mathrm{d}$ shell / sub-shell / orbitals is / are not full AND Cu+ is $\mathrm{d}^{10} / \mathrm{d}$ shell / sub-shell / orbitals is / are full [1] d-d* transitions / d electron promotion not possible / possible (linked correctly) [1] | 2 |


| Question | Answer | Marks |
| :---: | :--- | :---: |
| 6(a)(i) | $\mathrm{NH}_{3}=$ monodentate <br> $\mathrm{EDTA}^{4-}=$ polydentate / hexadentate <br> $\mathrm{CN}^{-}=$monodentate <br> $\mathrm{C}_{2} \mathrm{O}_{4}^{2-}=$ bidentate <br> any two [1] all four [2] | $\mathbf{2}$ |
| 6(a)(ii) | (ligand that) donates 3 lone pairs to central metal atom / ion <br> OR <br> (ligand that) forms 3 dative bonds to central metal atom / ion [1] | $\mathbf{1}$ |
| 6(a)(iii) | ref to using the electrons / lone pair on (each) N / amine group [1] | $\mathbf{1}$ |
| 6(b)(i) | trans = non-polar <br> cis isomer 1 = polar <br> cis isomer 2 = polar ALL [1] | $\mathbf{1}$ |
| 6(b)(ii) | optical isomers / non-superimposable mirror images [1] <br> ALLOW enantiomers / they rotate polarised light differently | $\mathbf{1}$ |


| Question | Answer | Marks |
| :---: | :--- | :---: |
| $7(\mathrm{a})$ | $\mathrm{C}_{16} \mathrm{H}_{10} \mathrm{~N}_{2} \mathrm{O}_{7} \mathrm{~S}_{2}{ }^{2-} \mathrm{OR} \mathrm{C}_{16} \mathrm{H}_{10} \mathrm{~N}_{2} \mathrm{O}_{7} \mathrm{~S}_{2}[1]$ | $\mathbf{1}$ |


| Question | Answer | Marks |
| :---: | :---: | :---: |
| 7(b) | E: <br> [1] <br> F: <br> or <br> [1] | 3 |
| 7(c) | M1 step 1: $\mathrm{HNO}_{2} \mathrm{OR} \mathrm{NaNO}_{2}+\mathrm{HCl}[1]$ M2 step 1: $\mathrm{T} \leqslant 10^{\circ} \mathrm{C}[1]$ <br> M3 step 2: $\mathrm{NaOH}(\mathrm{aq}) /$ alkaline conditions [1] | 3 |
| 7(d) | 14/fourteen [1] | 1 |

Question

| Question |  |  |  | Answer | Marks |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 8(d)(ii) | environment | ठ | splitting pattern | explanation for SP | 3 |
|  | $\mathrm{CH}_{3}$ | $0.9-1.7$ | doublet | 1 H on neighbouring C <br> / next to CH <br> / one vicinal proton |  |
|  | $\mathrm{CH}$ | $2.2-3.0$ | multiplet / heptet / septet | 6 H on neighbouring C <br> / next to $2 \times\left(\mathrm{CH}_{3}\right)$ <br> $/$ six vicinal protons |  |
|  | COOH | 9.0-13.0 | singlet |  |  |
|  | [1] for $\delta$ values <br> [1] for splitting <br> [1] for explanations |  |  |  |  |
| 8(e)(i) |  <br> K |  |  |  | 2 |
| 8(e)(ii) | hydrolysis AND neutralisation / acid-base[1] |  |  |  | 1 |

Question

| Question | Answer | Marks |
| :---: | :---: | :---: |
| 9(a)(i) | M <br> [1] | 1 |
| 9(a)(ii) | hot (alkaline) $\mathrm{KMnO}_{4} / \mathrm{MnO}_{4}^{-}$[1] <br> $\mathrm{SOCl}_{2} \mathrm{OR} \mathrm{PCl} l_{5} \mathrm{OR} \mathrm{PCl}_{3}+$ heat[1] | 2 |
| 9(a)(iii) | oxidation [1] <br> (nucleophilic) substitution | 2 |

Question

| Question | Answer | Marks |
| :---: | :---: | :---: |
| 9(c) | chlorobenzene, chloroethane, benzoyl chloride [1] u / c <br> explanation linked to their order <br> - correct link to: strengthening (ArCl/RCl) C-Cl bond <br> OR weakening ( RCOCl ) $\mathrm{C}-\mathrm{Cl}$ bond <br> OR $\mathrm{C}-\mathrm{Cl}$ bond has partially double bond character ( ArCl ) <br> OR C-Clis more difficult to break (linked correctly) <br> chlorobenzene: <br> - Ione pair / p-orbital on Cloverlaps / delocalised / incorporated with ring benzoyl chloride: <br> - C of $\mathrm{C}-\mathrm{Cl}$ has most electron deficient <br> OR has an electronegative oxygen atom <br> / two electronegative atoms / electron withdrawing $\mathrm{C}=\mathrm{O}$ group <br> chloroethane <br> - electron donating effect / positive inductive effect of alkyl / R group <br> any two [1] any three [2] | 3 |

