		-			_		_	-		1
•	W	<b>'O</b>	RI	KS	SH	EI	ET	#2	1	

Name:		Date:	Period:	Seat #:
		te the correct Le	ewis electron-dot struct	ture for each. Note the shape of the
molecule (for compounds only)				
<b>F</b> # of valence e <sup>-</sup> 's =	O # of valence e <sup>-</sup> 's =	K # of val	ence e <sup></sup> 's =	Al # of valence e <sup>-</sup> 's =
<b>F</b> <sup></sup> # of valence e <sup></sup> 's =	<b>O</b> <sup>2</sup> — # of valence e <sup>-</sup> 's =	K <sup>+</sup> # of val	ence e <sup></sup> s =	<b>Al<sup>3+</sup></b> # of valence e <sup>-,</sup> s =
<b>F</b> <sub>2</sub> # of valence e <sup></sup> 's =	$H_2$ # of valence e <sup>-+</sup> s =	HF # of val	ence e <sup></sup> 's =	<b>NH</b> <sub>3</sub> # of valence e <sup>-</sup> 's =
<b>CH</b> <sub>4</sub> # of valence e <sup>-</sup> 's =	<b>NF</b> 3 # of valence e <sup>-</sup> 's =	SiF4 # of val	ence e <sup></sup> 's =	<b>C</b> <sub>2</sub> <b>H</b> <sub>6</sub> # of valence e <sup>-+</sup> s =
<b>MgH</b> <sub>2</sub> # of valence e <sup>-</sup> 's =	LiH # of valence e <sup></sup> 's =	AlH3 # of val	ence e <sup></sup> 's =	<b>BH</b> <sub>3</sub> # of valence e <sup>-</sup> 's =

		~~~	2
$C_2H_4 # of valence e^-'s =$	$C_2F_4 $ # of valence e <sup></sup> 's =	<b>CO</b> # of valence e <sup>-</sup> 's =	<b>O</b> <sub>2</sub> # of valence e <sup>-</sup> 's =
$CO_2$	$C_2H_2$	$N_2$	HCN
# of valence $e^{-s} = $	# of valence $e^{-s} = $	# of valence $e^{-s} = $	# of valence $e^{-s} = $
CN <sup></sup>	SO4 <sup>2—</sup>	PO <sub>4</sub> <sup>3</sup> —	ClO <sub>3</sub> <sup></sup>
# of valence $e^{-s} = $	# of valence $e^{-s} = $	# of valence $e^{-s} = $	# of valence $e^{-s} = $
CO3 <sup>2</sup>	NO <sub>3</sub> —	SO <sub>2</sub>	O <sub>3</sub>
# of valence $e^{-s} = $	# of valence e <sup>-</sup> 's =	# of valence $e^{-s} =$	# of valence $e^{-s} = $
SF <sub>6</sub>	XeF <sub>4</sub>	PCl <sub>5</sub>	SeF <sub>4</sub>
# of valence $e^{-s} = $	# of valence $e^{-s} = $	# of valence $e^{-s} = $	# of valence $e^{-s} = $
1		1	