																					SV	M One	e vers	us One	e - Con	nfusion	Matri	X																		
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2	0	1	0	0	0	0	0	0	0 0	0	0	0	0	0	0	1	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0 0	0	0	0	0 0	0	0	0	0	0	0	0 0	) 0	0	0	0
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e Lab - 92	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0 2	0	0	0	0	0	0	0 0	0	0	0	0 0	0	0	0	0	0	0	0 0	0	0	0	0
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34 -	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0 2	0	0	0	0 0	0	0	0	0	0	0	0 0	0	0	0	0
35 -	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0 0	2	0	0	0 0	0	0	0	0	0	0	0 0	0	0	0	0
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42 -	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0 0	0	0	0	0 0	0	0	2	0	0	0	0 0	0	0	0	٥
43 -	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0 0	0	0	0	0 0	0	0	0	2	0	0	0 0	0	0	0	0
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45 -	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0 0	0	0	0	0 0	0	0	0	0	0	2	0 0	0	0	0	0
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48 -	0	0	0	0	0	0	0	0	υ 0	0	0	0	0	0	0	0	0							0 0			0	0	0	0	0 0	0	0	0	0 0	0	0	0	0	0	0	υ 0	2	0	0	J
49 -	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0 0	0	0	0	0 0	0	0	0	0	0	0	0 0	0	2	0	3
50 -	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0 0	0	0	0	0	0	0	0 0	0	0	0	0 0	0	0	0	0	0	0	0 0	0	0	2	0
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- 1.75 - 1.50 - 1.25 - 1.00 - 0.75 - 0.50 - 0.25