

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP010224\  
 Data File : BP019229.D  
 Acq On : 03 Jan 2024 03:07  
 Operator : MA/JU  
 Sample : 125 PPM TOX STD  
 Misc : GCMS CONFIRMATION  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 125 PPM TOX STD

Quant Time: Jan 03 04:40:27 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP122723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 29 03:48:11 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.799  | 152  | 248008   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.593 | 136  | 876549   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.445 | 164  | 495243   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.198 | 188  | 1015635  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.304 | 240  | 977423   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.669 | 264  | 1224400  | 20.000 | ng/ul | 0.00     |

|                               |       |     |    |       |       |  |
|-------------------------------|-------|-----|----|-------|-------|--|
| System Monitoring Compounds   |       |     |    |       |       |  |
| 3) 1,4-Dioxane-d8             | 0.000 | 96  | 0d | 0.000 | ng/uL |  |
| 4) Pyridine-d5                | 0.000 | 84  | 0d | 0.000 | ng/ul |  |
| 7) Phenol-d5                  | 0.000 | 99  | 0d | 0.000 | ng/ul |  |
| 9) Bis-(2-Chloroethyl)eth...  | 0.000 | 67  | 0d | 0.000 | ng/ul |  |
| 11) 2-Chlorophenol-d4         | 0.000 | 132 | 0d | 0.000 | ng/ul |  |
| 15) 4-Methylphenol-d8         | 0.000 | 113 | 0d | 0.000 | ng/ul |  |
| 21) Nitrobenzene-d5           | 0.000 | 128 | 0d | 0.000 | ng/ul |  |
| 24) 2-Nitrophenol-d4          | 0.000 | 143 | 0d | 0.000 | ng/ul |  |
| 28) 2,4-Dichlorophenol-d3     | 0.000 | 165 | 0d | 0.000 | ng/ul |  |
| 31) 4-Chloroaniline-d4        | 0.000 | 131 | 0d | 0.000 | ng/ul |  |
| 46) Dimethylphthalate-d6      | 0.000 | 166 | 0d | 0.000 | ng/ul |  |
| 49) Acenaphthylene-d8         | 0.000 | 160 | 0d | 0.000 | ng/ul |  |
| 54) 4-Nitrophenol-d4          | 0.000 | 143 | 0d | 0.000 | ng/ul |  |
| 60) Fluorene-d10              | 0.000 | 176 | 0d | 0.000 | ng/ul |  |
| 65) 4,6-Dinitro-2-methylph... | 0.000 | 200 | 0d | 0.000 | ng/ul |  |
| 73) Anthracene-d10            | 0.000 | 188 | 0d | 0.000 | ng/ul |  |
| 81) Pyrene-d10                | 0.000 | 212 | 0d | 0.000 | ng/ul |  |
| 92) Benzo(a)pyrene-d12        | 0.000 | 264 | 0d | 0.000 | ng/ul |  |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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