

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062724\  
 Data File : BP020697.D  
 Acq On : 28 Jun 2024 18:21  
 Operator : MA/JU  
 Sample : PB161630BL  
 Misc :  
 ALS Vial : 55 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK630

Quant Time: Jun 28 22:51:01 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062524.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jun 25 18:30:06 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.746  | 152  | 207473   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.516 | 136  | 878547   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.375 | 164  | 584876   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.192 | 188  | 1184961  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.651 | 240  | 817423   | 20.000 | ng/u1 | 0.01     |
| 88) Perylene-d12          | 25.016 | 264  | 945259   | 20.000 | ng/u1 | 0.02     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.270  | 96  | 26703   | 5.527  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.681  | 84  | 349268  | 24.804 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 6.922  | 99  | 494938  | 29.086 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.087  | 67  | 301459  | 27.944 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.281  | 132 | 409308  | 29.431 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.446  | 113 | 410204  | 29.681 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 8.893  | 128 | 207484  | 31.988 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.605  | 143 | 234203  | 36.081 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.140 | 165 | 419677  | 31.388 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.657 | 131 | 469551  | 25.059 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.787 | 166 | 1325528 | 32.596 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.063 | 160 | 1480066 | 30.501 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.604 | 143 | 182364  | 29.350 | ng/u1 | 0.01 |
| 60) Fluorene-d10              | 15.393 | 176 | 1122080 | 32.790 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.522 | 200 | 155284  | 26.004 | ng/u1 | 0.02 |
| 73) Anthracene-d10            | 17.292 | 188 | 1659318 | 31.201 | ng/u1 | 0.01 |
| 81) Pyrene-d10                | 19.681 | 212 | 1643420 | 33.466 | ng/u1 | 0.01 |
| 92) Benzo(a)pyrene-d12        | 24.786 | 264 | 1480563 | 32.199 | ng/u1 | 0.02 |

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

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

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