

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070623\  
 Data File : BP016072.D  
 Acq On : 07 Jul 2023 15:52  
 Operator : MA/JU  
 Sample : 03283-11  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EXYX9

Quant Time: Jul 07 20:28:03 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062623.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jul 06 23:18:26 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 8.034  | 152  | 210999   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.857 | 136  | 968351   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.681 | 164  | 635712   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.445 | 188  | 1388230  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12          | 21.557 | 240  | 1170450  | 20.000 | ng/u1 | 0.02     |
| 88) Perylene-d12          | 24.092 | 264  | 1094558  | 20.000 | ng/u1 | 0.01     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.352  | 96  | 27998   | 4.774  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.817  | 84  | 89658   | 6.065  | ng/u1 | 0.01 |
| 7) Phenol-d5                  | 7.269  | 99  | 82607   | 4.576  | ng/u1 | 0.05 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.363  | 67  | 357577  | 32.015 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.587  | 132 | 304985  | 22.379 | ng/u1 | 0.02 |
| 15) 4-Methylphenol-d8         | 8.799  | 113 | 206622  | 14.488 | ng/u1 | 0.03 |
| 21) Nitrobenzene-d5           | 9.234  | 128 | 199930  | 27.016 | ng/u1 | 0.02 |
| 24) 2-Nitrophenol-d4          | 9.969  | 143 | 218660  | 25.316 | ng/u1 | 0.02 |
| 28) 2,4-Dichlorophenol-d3     | 10.557 | 165 | 240273  | 15.611 | ng/u1 | 0.05 |
| 31) 4-Chloroaniline-d4        | 11.057 | 131 | 474933  | 24.020 | ng/u1 | 0.04 |
| 46) Dimethylphthalate-d6      | 14.098 | 166 | 1597149 | 32.505 | ng/u1 | 0.01 |
| 49) Acenaphthylene-d8         | 14.375 | 160 | 1662352 | 30.096 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.975 | 143 | 109     | 0.017  | ng/u1 | 0.01 |
| 60) Fluorene-d10              | 15.687 | 176 | 1317195 | 32.557 | ng/u1 | 0.01 |
| 65) 4,6-Dinitro-2-methylph... | 15.881 | 200 | 192126  | 22.504 | ng/u1 | 0.03 |
| 73) Anthracene-d10            | 17.551 | 188 | 2176597 | 34.325 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.780 | 212 | 2600211 | 34.022 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.916 | 264 | 1877031 | 33.995 | ng/u1 | 0.00 |

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

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

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