

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042321\  
 Data File : BN014383.D  
 Acq On : 23 Apr 2021 18:39  
 Operator : CG/JU  
 Sample : M2065-08  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BG561

Manual Integrations  
 APPROVED

mohammad  
 4/26/2021 11:37:43 AM

Quant Time: Apr 24 01:19:56 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN041521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 23 23:19:59 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.69  | 152  | 122397   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.46 | 136  | 504137   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.32 | 164  | 286329   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.07 | 188  | 593871   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.26 | 240  | 608597   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.49 | 264  | 604613   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.21  | 96  | 5430   | 1.69  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.63  | 84  | 18838m | 2.21  | ng/ul | 0.02 |
| 7) Phenol-d5                   | 6.85  | 99  | 105693 | 10.33 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.02  | 67  | 75072  | 12.25 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.22  | 132 | 91373  | 11.35 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.38  | 113 | 84541  | 10.90 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 8.83  | 128 | 47544  | 13.05 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.55  | 143 | 35864  | 10.05 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.08 | 165 | 75828  | 9.87  | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.59 | 131 | 86506  | 7.59  | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 13.73 | 166 | 268435 | 12.94 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 14.01 | 160 | 342794 | 13.74 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.52 | 143 | 10696  | 2.66  | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.32 | 176 | 242384 | 13.22 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.43 | 200 | 3835   | 1.21  | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.16 | 188 | 353267 | 12.74 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.46 | 212 | 385723 | 13.06 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.34 | 264 | 393358 | 12.26 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |              | Ovalue |
|----------------------------|-------|-----|---------|--------------|--------|
| 50) Acenaphthylene         | 14.04 | 152 | 37691   | 1.420 ng/ul  | 96     |
| 80) Fluoranthene           | 19.13 | 202 | 98381   | 2.635 ng/ul  | 99     |
| 82) Pyrene                 | 19.49 | 202 | 111056  | 2.794 ng/ul  | 99     |
| 85) Benzo(a)anthracene     | 21.25 | 228 | 148032m | 3.764 ng/ul  |        |
| 87) Chrysene               | 21.30 | 228 | 135524  | 3.556 ng/ul  | 97     |
| 90) Benzo(b)fluoranthene   | 22.82 | 252 | 159643  | 3.947 ng/ul# | 95     |
| 91) Benzo(k)fluoranthene   | 22.86 | 252 | 75504m  | 1.875 ng/ul  |        |
| 93) Benzo(a)pyrene         | 23.39 | 252 | 132535  | 3.703 ng/ul# | 97     |
| 94) Indeno(1,2,3-cd)pyrene | 25.72 | 276 | 76130   | 1.605 ng/ul  | 98     |
| 96) Benzo(g,h,i)perylene   | 26.40 | 276 | 53637   | 1.399 ng/ul  | 99     |

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

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