

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\  
 Data File : BN010391.D  
 Acq On : 02 Apr 2020 14:21  
 Operator : CG/JU  
 Sample : L2065-11DL 10X  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleID :  
 DBF07DL

Manual Integrations  
 APPROVED

mohammad  
 4/3/2020 2:19:15 PM

Quant Time: Apr 02 15:32:46 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN040120MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Apr 01 17:35:00 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.61  | 152  | 447243   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.38 | 136  | 1982096  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.23 | 164  | 1313466  | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.98 | 188  | 3044951  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.18 | 240  | 3531375  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.36 | 264  | 4031909  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |      |       |      |
|--------------------------------|-------|-----|--------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96  | 0d     | 0.00 | ng/uL |      |
| 5) Phenol-d5                   | 6.80  | 99  | 25665  | 0.69 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.96  | 67  | 57870  | 2.96 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.15  | 132 | 71255  | 2.41 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.31  | 113 | 50059  | 1.59 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.75  | 128 | 45428  | 3.09 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.46  | 143 | 41800  | 2.68 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.00 | 165 | 97638  | 3.02 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 0.00  | 131 | 0d     | 0.00 | ng/ul |      |
| 44) Dimethylphthalate-d6       | 13.65 | 166 | 378022 | 3.81 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.92 | 160 | 402904 | 3.48 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.44 | 143 | 10446  | 0.56 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.23 | 176 | 335809 | 3.89 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.35 | 200 | 29422  | 1.69 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.08 | 188 | 542251 | 3.88 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.38 | 212 | 646390 | 3.55 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.22 | 264 | 746945 | 3.72 | ng/ul | 0.00 |

## Target Compounds

|                         |       |     |           |         | Ovalue    |
|-------------------------|-------|-----|-----------|---------|-----------|
| 28) Naphthalene         | 10.42 | 128 | 33898724m | 317.913 | ng/ul     |
| 34) 2-Methylnaphthalene | 12.05 | 142 | 10169764  | 131.278 | ng/ul 99  |
| 41) 1,1'-Biphenyl       | 13.06 | 154 | 1940689   | 19.417  | ng/ul 100 |
| 50) Acenaphthene        | 14.30 | 153 | 6147849   | 71.653  | ng/ul 99  |
| 54) Dibenzofuran        | 14.64 | 168 | 5130772   | 42.455  | ng/ul 98  |
| 59) Fluorene            | 15.29 | 166 | 3051165   | 31.022  | ng/ul 99  |
| 70) Phenanthrene        | 17.02 | 178 | 4195198   | 24.864  | ng/ul 99  |
| 72) Anthracene          | 17.11 | 178 | 316989    | 1.878   | ng/ul 97  |
| 75) Carbazole           | 17.39 | 167 | 5550120   | 39.553  | ng/ul 99  |
| 77) Fluoranthene        | 19.05 | 202 | 470992    | 2.621   | ng/ul 99  |
| 80) Pyrene              | 19.40 | 202 | 286612    | 1.214   | ng/ul 99  |

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

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