

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\  
 Data File : BN010061.D  
 Acq On : 03 Mar 2020 14:39  
 Operator : CG/JU  
 Sample : L1507-17DL 30X  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleID :  
 DBD15DL

Manual Integrations  
 APPROVED

mohammad  
 3/4/2020 7:51:27 AM

Quant Time: Mar 03 15:27:30 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 03 12:28:17 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.37  | 152  | 61565    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.11 | 136  | 224335   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.00 | 164  | 146100   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.77 | 188  | 329756   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 20.99 | 240  | 314873   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.09 | 264  | 318921   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |      |     |    |      |       |  |
|--------------------------------|------|-----|----|------|-------|--|
| 3) 1,4-Dioxane-d8              | 0.00 | 96  | 0  | 0.00 | ng/uL |  |
| 5) Phenol-d5                   | 0.00 | 99  | 0  | 0.00 | ng/ul |  |
| 7) Bis-(2-Chloroethyl)ether-d  | 0.00 | 67  | 0  | 0.00 | ng/ul |  |
| 9) 2-Chlorophenol-d4           | 0.00 | 132 | 0  | 0.00 | ng/ul |  |
| 13) 4-Methylphenol-d8          | 0.00 | 113 | 0  | 0.00 | ng/ul |  |
| 19) Nitrobenzene-d5            | 0.00 | 128 | 0  | 0.00 | ng/ul |  |
| 22) 2-Nitrophenol-d4           | 0.00 | 143 | 0  | 0.00 | ng/ul |  |
| 26) 2,4-Dichlorophenol-d3      | 0.00 | 165 | 0  | 0.00 | ng/ul |  |
| 29) 4-Chloroaniline-d4         | 0.00 | 131 | 0d | 0.00 | ng/ul |  |
| 44) Dimethylphthalate-d6       | 0.00 | 166 | 0d | 0.00 | ng/ul |  |
| 47) Acenaphthylene-d8          | 0.00 | 160 | 0d | 0.00 | ng/ul |  |
| 52) 4-Nitrophenol-d4           | 0.00 | 143 | 0  | 0.00 | ng/ul |  |
| 58) Fluorene-d10               | 0.00 | 176 | 0d | 0.00 | ng/ul |  |
| 63) 4,6-Dinitro-2-methylphenol | 0.00 | 200 | 0  | 0.00 | ng/ul |  |
| 71) Anthracene-d10             | 0.00 | 188 | 0d | 0.00 | ng/ul |  |
| 79) Pyrene-d10                 | 0.00 | 212 | 0d | 0.00 | ng/ul |  |
| 90) Benzo(a)pyrene-d12         | 0.00 | 264 | 0d | 0.00 | ng/ul |  |

## Target Compounds

|                            |       |     |         |         | Ovalue |    |
|----------------------------|-------|-----|---------|---------|--------|----|
| 28) Naphthalene            | 10.16 | 128 | 510361  | 42.188  | ng/ul  | 99 |
| 34) 2-Methylnaphthalene    | 11.78 | 142 | 395620  | 47.097  | ng/ul  | 97 |
| 41) 1,1'-Biphenyl          | 12.83 | 154 | 114679  | 10.220  | ng/ul  | 98 |
| 50) Acenaphthene           | 14.07 | 153 | 703698  | 73.791  | ng/ul  | 99 |
| 54) Dibenzofuran           | 14.42 | 168 | 632167  | 47.229  | ng/ul  | 98 |
| 59) Fluorene               | 15.07 | 166 | 552924  | 49.222  | ng/ul  | 98 |
| 70) Phenanthrene           | 16.82 | 178 | 3114909 | 165.517 | ng/ul  | 98 |
| 72) Anthracene             | 16.90 | 178 | 324100  | 16.847  | ng/ul  | 99 |
| 75) Carbazole              | 17.18 | 167 | 131010  | 7.733   | ng/ul  | 99 |
| 77) Fluoranthene           | 18.85 | 202 | 2439174 | 110.650 | ng/ul  | 96 |
| 80) Pyrene                 | 19.21 | 202 | 1696920 | 74.751  | ng/ul# | 93 |
| 83) Benzo(a)anthracene     | 20.97 | 228 | 441257  | 20.392  | ng/ul  | 98 |
| 85) Chrysene               | 21.03 | 228 | 253673  | 11.908  | ng/ul  | 97 |
| 88) Benzo(b)fluoranthene   | 22.47 | 252 | 238497  | 11.534  | ng/ul  | 99 |
| 89) Benzo(k)fluoranthene   | 22.52 | 252 | 70997m  | 3.626   | ng/ul  |    |
| 91) Benzo(a)pyrene         | 23.00 | 252 | 136579  | 7.344   | ng/ul  | 98 |
| 92) Indeno(1,2,3-cd)pyrene | 25.13 | 276 | 37041   | 1.678   | ng/ul# | 89 |
| 94) Benzo(g,h,i)perylene   | 25.75 | 276 | 24229   | 1.309   | ng/ul  | 98 |

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

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