

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\  
 Data File : BP004306.D  
 Acq On : 15 Dec 2020 16:11  
 Operator : CG/JU  
 Sample : L5001-04  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A54C2

Manual Integrations  
 APPROVED

mohammad  
 12/17/2020 9:52:51 AM

Quant Time: Dec 15 17:36:06 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 15 09:00:06 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.85  | 152  | 349159   | 20.00 | ng/ul | 0.01     |
| 18) Naphthalene-d8        | 10.64 | 136  | 1458848  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.49 | 164  | 895710   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.26 | 188  | 1851474  | 20.00 | ng/ul | 0.01     |
| 78) Chrysene-d12          | 21.35 | 240  | 1712461  | 20.00 | ng/ul | 0.01     |
| 86) Perylene-d12          | 23.73 | 264  | 1933805  | 20.00 | ng/ul | 0.03     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.33  | 96  | 11851   | 1.19  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.01  | 99  | 291439  | 10.09 | ng/ul | 0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.17  | 67  | 157323  | 8.72  | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.38  | 132 | 234985  | 10.72 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.55  | 113 | 216616  | 9.93  | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.00  | 128 | 116816  | 9.78  | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.72  | 143 | 102864  | 10.44 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.26 | 165 | 243107  | 11.22 | ng/ul | 0.01 |
| 29) 4-Chloroaniline-d4         | 10.78 | 131 | 230385  | 8.58  | ng/ul | 0.01 |
| 44) Dimethylphthalate-d6       | 13.89 | 166 | 704270  | 10.29 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.18 | 160 | 952677  | 10.90 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.69 | 143 | 89107   | 7.71  | ng/ul | 0.01 |
| 58) Fluorene-d10               | 15.49 | 176 | 637336  | 10.28 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.62 | 200 | 2394    | 0.27  | ng/ul | 0.01 |
| 71) Anthracene-d10             | 17.35 | 188 | 966865  | 10.72 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.60 | 212 | 1001897 | 11.23 | ng/ul | 0.01 |
| 90) Benzo(a)pyrene-d12         | 23.57 | 264 | 1104229 | 10.84 | ng/ul | 0.03 |

## Target Compounds

|                            |       |     |         |       | Ovalue    |
|----------------------------|-------|-----|---------|-------|-----------|
| 6) Phenol                  | 7.03  | 94  | 58940   | 1.954 | ng/ul 95  |
| 45) Dimethylphthalate      | 13.94 | 163 | 133660  | 1.917 | ng/ul 99  |
| 70) Phenanthrene           | 17.30 | 178 | 173178  | 1.585 | ng/ul 97  |
| 77) Fluoranthene           | 19.27 | 202 | 674302  | 5.781 | ng/ul 96  |
| 80) Pyrene                 | 19.62 | 202 | 650216  | 5.629 | ng/ul 97  |
| 83) Benzo(a)anthracene     | 21.33 | 228 | 314383  | 2.785 | ng/ul# 87 |
| 85) Chrysene               | 21.39 | 228 | 429310  | 3.877 | ng/ul# 90 |
| 88) Benzo(b)fluoranthene   | 23.00 | 252 | 710844  | 5.796 | ng/ul# 80 |
| 89) Benzo(k)fluoranthene   | 23.05 | 252 | 201237m | 1.598 | ng/ul     |
| 91) Benzo(a)pyrene         | 23.62 | 252 | 424630  | 3.711 | ng/ul# 75 |
| 92) Indeno(1,2,3-cd)pyrene | 26.17 | 276 | 370227  | 2.580 | ng/ul# 92 |
| 94) Benzo(g,h,i)perylene   | 26.92 | 276 | 286197  | 2.377 | ng/ul# 84 |

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

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