

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP022420\  
 Data File : BP001874.D  
 Acq On : 24 Feb 2020 22:49  
 Operator : CG/JU  
 Sample : L1418-20DL 5X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleID :  
 C5AW4DL

Manual Integrations  
 APPROVED

mohammad  
 2/25/2020 3:42:35 PM

Quant Time: Feb 25 06:38:33 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP021820MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 25 05:55:24 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.65  | 152  | 369247   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.42 | 136  | 1968466  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.29 | 164  | 993916   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.04 | 188  | 2213999  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.13 | 240  | 2832107  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.34 | 264  | 2082819  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |      |       |      |
|--------------------------------|-------|-----|--------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.18  | 96  | 5261   | 0.59 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.83  | 99  | 79010  | 2.83 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.99  | 67  | 46484  | 2.87 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.19  | 132 | 72198  | 3.20 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.36  | 113 | 64406  | 2.91 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.79  | 128 | 41847  | 3.14 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.51  | 143 | 42547  | 3.53 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.05 | 165 | 96979  | 3.38 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.56 | 131 | 73230  | 2.16 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.70 | 166 | 240506 | 3.45 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.97 | 160 | 285309 | 3.32 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.49 | 143 | 22733  | 1.87 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.29 | 176 | 220733 | 3.52 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.40 | 200 | 13915  | 1.59 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.14 | 188 | 352616 | 3.62 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.38 | 212 | 555788 | 3.82 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.20 | 264 | 368900 | 3.67 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |        | Ovalue    |
|----------------------------|-------|-----|---------|--------|-----------|
| 48) Acenaphthylene         | 14.00 | 152 | 245772  | 2.638  | ng/ul 100 |
| 70) Phenanthrene           | 17.08 | 178 | 679979  | 5.569  | ng/ul 100 |
| 72) Anthracene             | 17.17 | 178 | 203438  | 1.687  | ng/ul 99  |
| 77) Fluoranthene           | 19.06 | 202 | 2963566 | 22.666 | ng/ul 99  |
| 80) Pyrene                 | 19.41 | 202 | 3418919 | 17.748 | ng/ul 100 |
| 83) Benzo(a)anthracene     | 21.12 | 228 | 1529818 | 8.203  | ng/ul 97  |
| 85) Chrysene               | 21.17 | 228 | 1403416 | 7.703  | ng/ul 98  |
| 88) Benzo(b)fluoranthene   | 22.69 | 252 | 1513759 | 12.351 | ng/ul 99  |
| 89) Benzo(k)fluoranthene   | 22.72 | 252 | 404216m | 3.182  | ng/ul     |
| 91) Benzo(a)pyrene         | 23.25 | 252 | 1116972 | 9.725  | ng/ul 98  |
| 92) Indeno(1,2,3-cd)pyrene | 25.57 | 276 | 730737  | 5.093  | ng/ul# 94 |
| 93) Dibenzo(a,h)anthracene | 25.59 | 278 | 152161  | 1.263  | ng/ul# 81 |
| 94) Benzo(g,h,i)perylene   | 26.26 | 276 | 727361  | 5.964  | ng/ul 97  |

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

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