

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121320\  
 Data File : BP004258.D  
 Acq On : 12 Dec 2020 14:35  
 Operator : CG/JU  
 Sample : L5000-01  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A54A1

Manual Integrations  
 APPROVED

mohammad  
 12/14/2020 2:46:03 PM

Quant Time: Dec 14 06:03:23 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 10 13:52:32 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.84  | 152  | 340319   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.63 | 136  | 1372242  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.48 | 164  | 852077   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.25 | 188  | 1856330  | 20.00 | ng/ul | 0.01     |
| 78) Chrysene-d12          | 21.33 | 240  | 1839489  | 20.00 | ng/ul | 0.01     |
| 86) Perylene-d12          | 23.70 | 264  | 2135200  | 20.00 | ng/ul | 0.02     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.33  | 96  | 16601   | 1.71  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.00  | 99  | 359062  | 12.75 | ng/ul | 0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.16  | 67  | 193840  | 11.03 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.37  | 132 | 295793  | 13.84 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.53  | 113 | 277656  | 13.05 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.99  | 128 | 140012  | 12.46 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.71  | 143 | 147293  | 15.89 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.25 | 165 | 294901  | 14.47 | ng/ul | 0.01 |
| 29) 4-Chloroaniline-d4         | 10.76 | 131 | 290267  | 11.49 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.89 | 166 | 862530  | 13.24 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.17 | 160 | 1124575 | 13.53 | ng/ul | 0.01 |
| 52) 4-Nitrophenol-d4           | 14.67 | 143 | 130642  | 11.88 | ng/ul | 0.02 |
| 58) Fluorene-d10               | 15.48 | 176 | 770632  | 13.07 | ng/ul | 0.01 |
| 63) 4,6-Dinitro-2-methylphenol | 15.60 | 200 | 16626   | 1.87  | ng/ul | 0.02 |
| 71) Anthracene-d10             | 17.34 | 188 | 1229579 | 13.59 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.58 | 212 | 1403978 | 14.65 | ng/ul | 0.01 |
| 90) Benzo(a)pyrene-d12         | 23.54 | 264 | 1585080 | 14.09 | ng/ul | 0.02 |

## Target Compounds

|                                |       |     |          |        | Ovalue |     |
|--------------------------------|-------|-----|----------|--------|--------|-----|
| 28) Naphthalene                | 10.68 | 128 | 401771   | 5.165  | ng/ul  | 99  |
| 34) 2-Methylnaphthalene        | 12.30 | 142 | 144868   | 2.768  | ng/ul  | 95  |
| 45) Dimethylphthalate          | 13.93 | 163 | 231742   | 3.494  | ng/ul  | 100 |
| 50) Acenaphthene               | 14.55 | 153 | 318313   | 5.450  | ng/ul  | 99  |
| 54) Dibenzofuran               | 14.88 | 168 | 497236   | 6.042  | ng/ul  | 99  |
| 59) Fluorene                   | 15.53 | 166 | 416392   | 6.326  | ng/ul  | 98  |
| 70) Phenanthrene               | 17.29 | 178 | 6373287  | 58.195 | ng/ul  | 98  |
| 72) Anthracene                 | 17.37 | 178 | 1360599  | 12.557 | ng/ul  | 100 |
| 75) Carbazole                  | 17.65 | 167 | 760871   | 8.629  | ng/ul  | 99  |
| 77) Fluoranthene               | 19.26 | 202 | 8227110  | 70.354 | ng/ul  | 97  |
| 80) Pyrene                     | 19.61 | 202 | 6966817  | 56.146 | ng/ul  | 97  |
| 83) Benzo(a)anthracene         | 21.32 | 228 | 3558354  | 29.347 | ng/ul  | 97  |
| 84) Bis(2-ethylhexyl)phthalate | 21.25 | 149 | 56014    | 1.010  | ng/ul# | 67  |
| 85) Chrysene                   | 21.37 | 228 | 3838533  | 32.270 | ng/ul  | 96  |
| 88) Benzo(b)fluoranthene       | 22.98 | 252 | 4840596  | 35.748 | ng/ul  | 98  |
| 89) Benzo(k)fluoranthene       | 23.02 | 252 | 1839707m | 13.229 | ng/ul  |     |
| 91) Benzo(a)pyrene             | 23.59 | 252 | 3412984  | 27.011 | ng/ul  | 97  |
| 92) Indeno(1,2,3-cd)pyrene     | 26.12 | 276 | 2509662  | 15.838 | ng/ul  | 94  |
| 93) Dibenzo(a,h)anthracene     | 26.13 | 278 | 608784   | 4.584  | ng/ul# | 85  |
| 94) Benzo(g,h,i)perylene       | 26.87 | 276 | 1892484  | 14.235 | ng/ul  | 94  |

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121320\  
Data File : BP004258.D  
Acq On : 12 Dec 2020 14:35  
Operator : CG/JU  
Sample : L5000-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A54A1

Manual Integrations  
APPROVED

mohammad  
12/14/2020 2:46:03 PM

Quant Time: Dec 14 06:03:23 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP121020MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Dec 10 13:52:32 2020  
Response via : Initial Calibration

