

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020921\  
 Data File : BP004701.D  
 Acq On : 09 Feb 2021 13:07  
 Operator : CG/JU  
 Sample : SSTD16004  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD160004

Manual Integrations  
 APPROVED

mohammad  
 2/10/2021 12:16:24 PM

Quant Time: Feb 09 13:36:39 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP020921.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 09 13:12:53 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.775  | 152  | 50957    | 20.00  | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.569 | 136  | 205342   | 20.00  | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.422 | 164  | 127793   | 20.00  | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.175 | 188  | 272457   | 20.00  | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.263 | 240  | 285042   | 20.00  | ng/ul | 0.01     |
| 88) Perylene-d12              | 23.562 | 264  | 268801   | 20.00  | ng/ul | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.00   | ng/uL |          |
| 4) Pyridine-d5                | 3.658  | 84   | 629694   | 195.42 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.952  | 99   | 751103   | 180.17 | ng/ul | 0.02     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.116  | 67   | 383079   | 148.40 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 0.000  | 132  | 0d       | 0.00   | ng/ul |          |
| 15) 4-Methylphenol-d8         | 8.493  | 113  | 588869   | 182.76 | ng/ul | 0.02     |
| 21) Nitrobenzene-d5           | 0.000  | 128  | 0d       | 0.00   | ng/ul |          |
| 24) 2-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.00   | ng/ul |          |
| 28) 2,4-Dichlorophenol-d3     | 0.000  | 165  | 0d       | 0.00   | ng/ul |          |
| 31) 4-Chloroaniline-d4        | 10.710 | 131  | 786740   | 170.26 | ng/ul | 0.01     |
| 46) Dimethylphthalate-d6      | 0.000  | 166  | 0d       | 0.00   | ng/ul |          |
| 49) Acenaphthylene-d8         | 0.000  | 160  | 0d       | 0.00   | ng/ul |          |
| 54) 4-Nitrophenol-d4          | 14.645 | 143  | 309026   | 196.05 | ng/ul | 0.04     |
| 60) Fluorene-d10              | 0.000  | 176  | 0d       | 0.00   | ng/ul |          |
| 65) 4,6-Dinitro-2-methylph... | 15.551 | 200  | 328739   | 177.85 | ng/ul | 0.02     |
| 73) Anthracene-d10            | 0.000  | 188  | 0d       | 0.00   | ng/ul |          |
| 81) Pyrene-d10                | 0.000  | 212  | 0d       | 0.00   | ng/ul |          |
| 92) Benzo(a)pyrene-d12        | 0.000  | 264  | 0d       | 0.00   | ng/ul |          |
|                               |        |      |          |        |       |          |
| Target Compounds              |        |      |          |        |       | Qvalue   |
| 5) Pyridine                   | 3.675  | 79   | 625510   | 188.44 | ng/ul | 95       |
| 6) Benzaldehyde               | 6.916  | 77   | 171799   | 69.09  | ng/ul | 96       |
| 8) Phenol                     | 6.981  | 94   | 792736   | 177.84 | ng/ul | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.210  | 93   | 571854   | 172.50 | ng/ul | 96       |
| 13) 2-Methylphenol            | 8.222  | 108  | 590144   | 180.08 | ng/ul | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.310  | 45   | 444487   | 121.25 | ng/ul | 96       |
| 16) Acetophenone              | 8.604  | 105  | 979697   | 170.61 | ng/ul | 99       |
| 18) 4-Methylphenol            | 8.569  | 108  | 621279   | 177.47 | ng/ul | 95       |
| 32) 4-Chloroaniline           | 10.740 | 127  | 807234   | 168.96 | ng/ul | 100      |
| 34) Caprolactam               | 11.545 | 113  | 186412m  | 165.98 | ng/ul |          |
| 40) Hexachlorocyclopentadiene | 12.587 | 237  | 531827   | 138.96 | ng/ul | 99       |
| 51) 3-Nitroaniline            | 14.339 | 138  | 265720   | 161.34 | ng/ul | 95       |
| 53) 2,4-Dinitrophenol         | 14.539 | 184  | 249360   | 170.93 | ng/ul | 98       |
| 55) 4-Nitrophenol             | 14.657 | 109  | 333203   | 155.65 | ng/ul | 96       |
| 63) 4-Nitroaniline            | 15.516 | 138  | 241325   | 145.47 | ng/ul | 98       |
| 66) 4,6-Dinitro-2-methylph... | 15.563 | 198  | 331790   | 176.02 | ng/ul | 97       |
| 70) Atrazine                  | 16.651 | 200  | 572605   | 157.10 | ng/ul | 98       |
| 71) Pentachlorophenol         | 16.822 | 266  | 396250   | 151.28 | ng/ul | 97       |
| 77) Carbazole                 | 17.586 | 167  | 2345146  | 177.19 | ng/ul | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.180 | 252  | 1066160  | 151.60 | ng/ul | 100      |
| 89) Di-n-octyl phthalate      | 22.068 | 149  | 3365107  | 230.20 | ng/ul | 100      |
| -----                         |        |      |          |        |       |          |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SSTD160004

**Manual Integrations**

**APPROVED**

mohammad

2/10/2021 12:16:24 PM

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020921\  
Data File : BP004701.D  
Acq On : 09 Feb 2021 13:07  
Operator : CG/JU  
Sample : SSTD16004  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTD160004

Manual Integrations  
APPROVED

mohammad  
2/10/2021 12:16:24 PM

Quant Time: Feb 09 13:36:39 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP020921.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Feb 09 13:12:53 2021  
Response via : Initial Calibration

