

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032719\  
 Data File : BF113342.D  
 Acq On : 27 Mar 2019 20:35  
 Operator : JU/SJ  
 Sample : K2097-01  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CS-20

Manual Integrations  
 APPROVED

Sohil  
 3/28/2019 8:41:23 AM

Quant Time: Mar 28 04:19:11 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 26 03:32:08 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min)  |
|-----------------------------|-------|------|----------|--------|-------|-----------|
| 1) 1,4-Dichlorobenzene-d4   | 6.70  | 152  | 173617   | 20.00  | ng    | 0.00      |
| 21) Naphthalene-d8          | 7.98  | 136  | 622890   | 20.00  | ng    | 0.00      |
| 39) Acenaphthene-d10        | 9.73  | 164  | 291795   | 20.00  | ng    | 0.00      |
| 64) Phenanthrene-d10        | 11.21 | 188  | 495398   | 20.00  | ng    | 0.00      |
| 76) Chrysene-d12            | 13.84 | 240  | 377011   | 20.00  | ng    | 0.00      |
| 87) Perylene-d12            | 15.26 | 264  | 320730   | 20.00  | ng    | 0.01      |
| System Monitoring Compounds |       |      |          |        |       |           |
| 5) 2-Fluorophenol           | 5.32  | 112  | 1269273  | 119.54 | ng    | 0.00      |
| 7) Phenol-d6                | 6.36  | 99   | 1654492  | 123.17 | ng    | 0.00      |
| 23) Nitrobenzene-d5         | 7.27  | 82   | 1029272  | 98.08  | ng    | 0.00      |
| 42) 2,4,6-Tribromophenol    | 10.52 | 330  | 393181   | 109.09 | ng    | 0.00      |
| 45) 2-Fluorobiphenyl        | 9.06  | 172  | 1610745  | 86.92  | ng    | 0.00      |
| 79) Terphenyl-d14           | 12.79 | 244  | 1328452  | 77.67  | ng    | 0.00      |
| Target Compounds            |       |      |          |        |       |           |
| 10) Phenol                  | 6.36  | 94   | 58538    | 3.817  | ng    | Qvalue 94 |
| 31) Naphthalene             | 8.00  | 128  | 364990   | 11.991 | ng    | 98        |
| 37) 2-Methylnaphthalene     | 8.70  | 142  | 123465   | 6.501  | ng    | 97        |
| 38) 1-Methylnaphthalene     | 8.79  | 142  | 50358    | 2.757  | ng    | 100       |
| 46) 1,1'-Biphenyl           | 9.16  | 154  | 60552    | 2.485  | ng    | 98        |
| 49) Acenaphthylene          | 9.59  | 152  | 318107   | 10.485 | ng    | 99        |
| 50) Dimethylphthalate       | 9.45  | 163  | 185442   | 8.576  | ng    | 99        |
| 55) Dibenzofuran            | 9.93  | 168  | 99373    | 3.869  | ng    | # 96      |
| 71) Phenanthrene            | 11.23 | 178  | 360522   | 14.654 | ng    | 98        |
| 72) Anthracene              | 11.29 | 178  | 312647   | 12.511 | ng    | 99        |
| 73) Carbazole               | 11.44 | 167  | 183463   | 7.694  | ng    | 99        |
| 75) Fluoranthene            | 12.42 | 202  | 1004914  | 36.133 | ng    | 99        |
| 78) Pvrene                  | 12.64 | 202  | 1054832  | 40.375 | ng    | 99        |
| 81) Benzo(a)anthracene      | 13.83 | 228  | 561958m  | 26.040 | ng    |           |
| 83) Chrysene                | 13.87 | 228  | 957541   | 43.683 | ng    | 98        |
| 86) Indeno(1,2,3-cd)pyrene  | 16.62 | 276  | 465638   | 24.752 | ng    | 97        |
| 88) Benzo(b)fluoranthene    | 14.86 | 252  | 1387311m | 69.631 | ng    |           |
| 89) Benzo(k)fluoranthene    | 14.89 | 252  | 370141m  | 21.155 | ng    |           |
| 90) Benzo(a)pvrene          | 15.20 | 252  | 472560   | 26.754 | ng    | 96        |
| 91) Dibenzo(a,h)anthracene  | 16.62 | 278  | 144333m  | 9.450  | ng    |           |
| 92) Benzo(g,h,i)perylene    | 17.02 | 276  | 430094   | 27.930 | ng    | 98        |

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

