

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG020224\  
 Data File : BG060548.D  
 Acq On : 2 Feb 2024 1:04  
 Operator : MA/JU  
 Sample : P1313-02  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 COMP-2

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 02/02/2024  
 Supervised By :Sohil Jodhani 02/02/2024

Quant Time: Feb 02 01:42:51 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG010824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jan 09 04:07:12 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.920  | 152  | 146262   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.723 | 136  | 594888   | 20.000 | ng    | # 0.00   |
| 39) Acenaphthene-d10          | 14.559 | 164  | 443894   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.309 | 188  | 1099091  | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 21.575 | 240  | 1040701  | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 24.736 | 264  | 1195874  | 20.000 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.494  | 112  | 699219   | 78.631 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.086  | 99   | 886285   | 67.308 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 9.083  | 82   | 650658   | 51.647 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 16.052 | 330  | 577977   | 88.498 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.179 | 172  | 1712670  | 53.642 | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.930 | 244  | 2763982  | 59.888 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| Target Compounds              |        |      |          |        |       | Qvalue   |
| 71) Phenanthrene              | 17.350 | 178  | 657255   | 12.372 | ng    | 93       |
| 72) Anthracene                | 17.444 | 178  | 127594   | 2.406  | ng    | 94       |
| 75) Fluoranthene              | 19.366 | 202  | 1214063  | 19.027 | ng    | 88       |
| 78) Pyrene                    | 19.730 | 202  | 1067681  | 16.065 | ng    | # 86     |
| 81) Benzo(a)anthracene        | 21.551 | 228  | 553746   | 8.595  | ng    | 92       |
| 83) Chrysene                  | 21.616 | 228  | 502398   | 7.929  | ng    | 93       |
| 84) Bis(2-ethylhexyl)phtha... | 21.457 | 149  | 267258   | 7.735  | ng    | 97       |
| 87) Indeno(1,2,3-cd)pyrene    | 28.337 | 276  | 277053   | 3.332  | ng    | 95       |
| 88) Benzo(b)fluoranthene      | 23.725 | 252  | 711774m  | 10.075 | ng    |          |
| 89) Benzo(k)fluoranthene      | 23.790 | 252  | 148501m  | 2.130  | ng    |          |
| 90) Benzo(a)pyrene            | 24.589 | 252  | 502213   | 7.869  | ng    | 96       |
| 92) Benzo(g,h,i)perylene      | 29.460 | 276  | 283675   | 4.076  | ng    | 99       |
| -----                         |        |      |          |        |       |          |

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

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