

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG080522\  
 Data File : BG054535.D  
 Acq On : 5 Aug 2022 14:43  
 Operator : CG/JU  
 Sample : N3931-12  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 RVE-123

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 08/08/2022  
 Supervised By : Jagrut Upadhyay 08/08/2022

Quant Time: Aug 06 05:39:38 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jul 28 04:30:50 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.977  | 152  | 14996    | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 10.785 | 136  | 64896    | 20.000  | ng    | #-0.01   |
| 39) Acenaphthene-d10        | 14.616 | 164  | 55575    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.366 | 188  | 140189   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.631 | 240  | 150698   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 24.845 | 264  | 159118   | 20.000  | ng    | 0.01     |
|                             |        |      |          |         |       |          |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.562  | 112  | 98067    | 136.538 | ng    | 0.00     |
| 7) Phenol-d6                | 7.178  | 99   | 138808   | 141.041 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.140  | 82   | 92371    | 77.512  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 16.114 | 330  | 138296   | 118.519 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.235 | 172  | 305705   | 77.599  | ng    | -0.01    |
| 79) Terphenyl-d14           | 19.974 | 244  | 638729   | 84.083  | ng    | 0.00     |
|                             |        |      |          |         |       |          |
| Target Compounds            |        |      |          |         |       | Qvalue   |
| 49) Acenaphthylene          | 14.351 | 152  | 12092    | 2.558   | ng    | 95       |
| 71) Phenanthrene            | 17.407 | 178  | 21477    | 3.076   | ng    | 98       |
| 75) Fluoranthene            | 19.422 | 202  | 43926    | 4.616   | ng    | 96       |
| 78) Pyrene                  | 19.780 | 202  | 86885    | 10.048  | ng    | 99       |
| 81) Benzo(a)anthracene      | 21.614 | 228  | 42091    | 4.400   | ng    | 94       |
| 83) Chrysene                | 21.678 | 228  | 53707    | 5.942   | ng    | 95       |
| 87) Indeno(1,2,3-cd)pyrene  | 28.535 | 276  | 43657    | 3.607   | ng    | 94       |
| 88) Benzo(b)fluoranthene    | 23.823 | 252  | 62456m   | 6.637   | ng    |          |
| 89) Benzo(k)fluoranthene    | 23.870 | 252  | 23044m   | 2.461   | ng    |          |
| 90) Benzo(a)pyrene          | 24.686 | 252  | 63772    | 7.936   | ng    | 94       |
| 92) Benzo(g,h,i)perylene    | 29.681 | 276  | 59846    | 6.092   | ng    | # 91     |
| -----                       |        |      |          |         |       |          |

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

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