

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082719\  
 Data File : BG042517.D  
 Acq On : 27 Aug 2019 23:21  
 Operator : HP/JU  
 Sample : K4526-19  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 T9-12-13-L-(0-1.9)

Manual Integrations  
 APPROVED

mohammad  
 8/29/2019 7:58:25 AM

Quant Time: Aug 28 03:35:06 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Aug 22 14:29:15 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 8.00  | 152  | 33422    | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.82 | 136  | 127972   | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.66 | 164  | 97806    | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.41 | 188  | 249389   | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.68 | 240  | 297448   | 20.00  | ng    | -0.01    |
| 87) Perylene-d12            | 24.92 | 264  | 352483   | 20.00  | ng    | -0.02    |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.55  | 112  | 149574   | 90.64  | ng    | 0.00     |
| 7) Phenol-d6                | 7.16  | 99   | 200058   | 89.75  | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 9.16  | 82   | 109494   | 59.13  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 16.15 | 330  | 123848   | 89.09  | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 13.28 | 172  | 323549   | 55.95  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.02 | 244  | 692983   | 50.37  | ng    | 0.00     |
| Target Compounds            |       |      |          |        |       |          |
| 10) Phenol                  | 7.19  | 94   | 5962     | 2.631  | ng    | # 75     |
| 50) Dimethylphthalate       | 14.11 | 163  | 20698    | 2.933  | ng    | # 96     |
| 71) Phenanthrene            | 17.45 | 178  | 77964    | 6.294  | ng    | 97       |
| 75) Fluoranthene            | 19.46 | 202  | 160282   | 10.602 | ng    | 97       |
| 78) Pyrene                  | 19.83 | 202  | 141248   | 7.745  | ng    | 97       |
| 81) Benzo(a)anthracene      | 21.67 | 228  | 73049    | 4.037  | ng    | 99       |
| 83) Chrysene                | 21.73 | 228  | 80334    | 4.604  | ng    | 98       |
| 86) Indeno(1,2,3-cd)pyrene  | 28.62 | 276  | 57232    | 2.413  | ng    | # 94     |
| 88) Benzo(b)fluoranthene    | 23.89 | 252  | 112238   | 5.792  | ng    | 99       |
| 89) Benzo(k)fluoranthene    | 23.96 | 252  | 38716m   | 2.079  | ng    |          |
| 90) Benzo(a)pyrene          | 24.77 | 252  | 75985    | 4.132  | ng    | 97       |
| 92) Benzo(a,h,i)perylene    | 29.76 | 276  | 55817    | 2.998  | ng    | 94       |

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

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