

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061220\  
 Data File : BG045754.D  
 Acq On : 13 Jun 2020 00:05  
 Operator : CG/JU  
 Sample : L2979-02  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 15-SP

Manual Integrations  
 APPROVED

mohammad  
 6/15/2020 8:45:21 AM

Quant Time: Jun 13 01:01:40 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG051520.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 01 15:07:00 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.95  | 152  | 62354    | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 10.76 | 136  | 254065   | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 14.59 | 164  | 169034   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 17.34 | 188  | 355050   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 21.60 | 240  | 293284   | 20.00 | ng    | 0.00     |
| 87) Perylene-d12          | 24.75 | 264  | 321394   | 20.00 | ng    | 0.00     |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 5) 2-Fluorophenol           | 5.58  | 112  | 369980   | 115.96 | ng    | 0.04     |
| 7) Phenol-d6                | 7.21  | 99   | 560634   | 123.46 | ng    | 0.07     |
| 23) Nitrobenzene-d5         | 9.11  | 82   | 364035   | 78.14  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.10 | 330  | 239265   | 110.68 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.21 | 172  | 831591   | 83.45  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.96 | 244  | 1041342  | 82.81  | ng    | 0.00     |

| Target Compounds           | R.T.  | QIon | Response | Conc  | Units | Qvalue |
|----------------------------|-------|------|----------|-------|-------|--------|
| 9) Benzaldehyde            | 7.09  | 77   | 15208    | 5.140 | ng    | 91     |
| 50) Dimethylphthalate      | 14.05 | 163  | 52914    | 4.563 | ng    | 100    |
| 71) Phenanthrene           | 17.38 | 178  | 67713    | 4.369 | ng    | 97     |
| 75) Fluoranthene           | 19.40 | 202  | 89644    | 4.547 | ng    | 91     |
| 78) Pyrene                 | 19.76 | 202  | 79234    | 4.739 | ng    | 95     |
| 81) Benzo(a)anthracene     | 21.58 | 228  | 35343m   | 2.163 | ng    |        |
| 83) Chrysene               | 21.65 | 228  | 44671    | 2.844 | ng    | 98     |
| 86) Indeno(1,2,3-cd)pyrene | 28.32 | 276  | 41854    | 2.037 | ng    | # 70   |
| 88) Benzo(b)fluoranthene   | 23.76 | 252  | 61554    | 3.571 | ng    | # 93   |
| 90) Benzo(a)pyrene         | 24.60 | 252  | 40400    | 2.517 | ng    | # 96   |
| 92) Benzo(g,h,i)perylene   | 29.43 | 276  | 38658    | 2.396 | ng    | 97     |

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

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