

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM083119\  
 Data File : BM022469.D  
 Acq On : 31 Aug 2019 13:10  
 Operator : HP/JU  
 Sample : K4568-16  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SB-36-COMP

Manual Integrations  
 APPROVED

mohammad  
 9/5/2019 9:13:02 AM

Quant Time: Aug 31 16:02:32 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082919.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Aug 29 16:51:32 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.52  | 152  | 25208    | 20.00 | ng    | -0.01    |
| 21) Naphthalene-d8        | 10.29 | 136  | 90067    | 20.00 | ng    | 0.00     |
| 39) Acenaphthene-d10      | 14.18 | 164  | 57797    | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 16.95 | 188  | 132189   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 21.17 | 240  | 173203   | 20.00 | ng    | 0.00     |
| 87) Perylene-d12          | 23.36 | 264  | 201166   | 20.00 | ng    | 0.00     |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 5) 2-Fluorophenol           | 5.15  | 112  | 185147   | 128.92 | ng    | 0.00     |
| 7) Phenol-d6                | 6.72  | 99   | 233205   | 121.43 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.69  | 82   | 167432   | 81.42  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.69 | 330  | 143873   | 121.65 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.79 | 172  | 389037   | 76.25  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.60 | 244  | 951287   | 72.48  | ng    | 0.00     |

| Target Compounds           | R.T.  | QIon | Response | Conc  | Units | Qvalue |
|----------------------------|-------|------|----------|-------|-------|--------|
| 10) Phenol                 | 6.75  | 94   | 5123     | 2.655 | ng    | # 59   |
| 49) Acenaphthylene         | 13.90 | 152  | 46311    | 8.183 | ng    | 98     |
| 50) Dimethylphthalate      | 13.66 | 163  | 43712    | 8.696 | ng    | 99     |
| 72) Anthracene             | 17.09 | 178  | 22750    | 3.001 | ng    | 99     |
| 75) Fluoranthene           | 19.02 | 202  | 26987    | 2.630 | ng    | 97     |
| 78) Pyrene                 | 19.39 | 202  | 80563    | 6.825 | ng    | 97     |
| 81) Benzo(a)anthracene     | 21.15 | 228  | 39285    | 3.160 | ng    | # 89   |
| 83) Chrysene               | 21.20 | 228  | 44283    | 3.742 | ng    | 93     |
| 86) Indeno(1,2,3-cd)pyrene | 25.53 | 276  | 51880    | 4.359 | ng    | 97     |
| 88) Benzo(b)fluoranthene   | 22.70 | 252  | 98319m   | 7.035 | ng    |        |
| 90) Benzo(a)pyrene         | 23.26 | 252  | 81673    | 6.584 | ng    | 95     |
| 92) Benzo(a,h,i)perylene   | 26.20 | 276  | 50327    | 4.668 | ng    | 98     |

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

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