

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081720\  
 Data File : BP003007.D  
 Acq On : 17 Aug 2020 17:44  
 Operator : CG/JU  
 Sample : L3691-09  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 S711-N-SO-0-0.5-081320

Manual Integrations  
 APPROVED

mohammad  
 8/19/2020 2:21:02 PM

Quant Time: Aug 18 04:13:50 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 17 14:07:35 2020  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4      | 7.69  | 152  | 368369   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8             | 10.46 | 136  | 1322662  | 20.00  | ng    | 0.00     |
| 39) Acenaphthene-d10           | 14.32 | 164  | 878208   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10           | 17.07 | 188  | 1659721  | 20.00  | ng    | 0.00     |
| 76) Chrysene-d12               | 21.17 | 240  | 1990162  | 20.00  | ng    | 0.00     |
| 86) Perylene-d12               | 23.42 | 264  | 1903514  | 20.00  | ng    | 0.00     |
| System Monitoring Compounds    |       |      |          |        |       |          |
| 5) 2-Fluorophenol              | 5.29  | 112  | 1787146  | 74.25  | ng    | 0.00     |
| 7) Phenol-d6                   | 6.86  | 99   | 2143465  | 64.14  | ng    | 0.00     |
| 23) Nitrobenzene-d5            | 8.83  | 82   | 1274130  | 52.18  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol       | 15.82 | 330  | 841879   | 77.13  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl           | 12.95 | 172  | 3153400  | 49.06  | ng    | 0.00     |
| 79) Terphenyl-d14              | 19.65 | 244  | 4206380  | 42.31  | ng    | 0.00     |
| Target Compounds               |       |      |          |        |       |          |
| 10) Phenol                     | 6.89  | 94   | 91520    | 2.719  | ng    | 95       |
| 32) Benzoic acid               | 9.73  | 122  | 15978m   | 8.778  | ng    |          |
| 50) Dimethylphthalate          | 13.78 | 163  | 852757   | 12.415 | ng    | 100      |
| 71) Phenanthrene               | 17.12 | 178  | 412169   | 4.235  | ng    | 99       |
| 75) Fluoranthene               | 19.10 | 202  | 1558068  | 13.895 | ng    | 98       |
| 78) Pyrene                     | 19.45 | 202  | 1305936  | 9.387  | ng    | 99       |
| 81) Benzo(a)anthracene         | 21.16 | 228  | 753618   | 5.559  | ng    | 100      |
| 83) Chrysene                   | 21.20 | 228  | 965261   | 7.548  | ng    | 98       |
| 84) Bis(2-ethylhexyl)phthalate | 21.10 | 149  | 224555   | 2.916  | ng    | 97       |
| 87) Indeno(1,2,3-cd)pyrene     | 25.70 | 276  | 796724   | 5.226  | ng    | 95       |
| 88) Benzo(b)fluoranthene       | 22.74 | 252  | 1604641m | 11.946 | ng    |          |
| 89) Benzo(k)fluoranthene       | 22.78 | 252  | 413395m  | 3.256  | ng    |          |
| 90) Benzo(a)pyrene             | 23.32 | 252  | 868366   | 7.154  | ng    | 100      |
| 92) Benzo(a,h,i)perylene       | 26.40 | 276  | 805560   | 6.421  | ng    | 98       |

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

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