

Data File : BM012277.D

Acq On : 18 Oct 2017 08:20

Operator : SJ/JU

Sample : I5747-02

Misc :

ALS Vial : 28 Sample Multiplier: 1

Instrument :

BNA\_M

ClientSampleId :

B-23(1-3)-100517

Manual Integrations

APPROVED

Sohil

10/18/2017 7:33:53 PM

Quant Time: Oct 30 08:21:38 2017

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Oct 18 08:12:28 2017

Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.78  | 152  | 211813   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.59 | 136  | 778803   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.44 | 164  | 511141   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.19 | 188  | 1202718  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.37 | 240  | 1232564  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.66 | 264  | 937661   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.20  | 96  | 17945   | 4.03  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.96  | 99  | 486732  | 28.32 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.12  | 67  | 296845  | 28.80 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.32  | 132 | 461793  | 31.68 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.50  | 113 | 304788  | 20.60 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.95  | 128 | 175538  | 29.54 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.67  | 143 | 213724  | 30.78 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.22 | 165 | 406887  | 30.69 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.73 | 131 | 424151  | 34.18 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.84 | 166 | 1413148 | 31.28 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.13 | 160 | 1693544 | 30.97 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.67 | 143 | 131537  | 19.37 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.43 | 176 | 1192824 | 32.17 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.58 | 200 | 158140  | 19.16 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.29 | 188 | 1836111 | 30.87 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.58 | 212 | 2404402 | 38.43 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.52 | 264 | 1426056 | 31.55 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |      |        | Qvalue |
|----------------------------|-------|-----|---------|------|--------|--------|
| 4) Benzaldehyde            | 6.93  | 77  | 62438   | 5.44 | ng/ul  | 96     |
| 6) Phenol                  | 6.98  | 94  | 113833  | 6.41 | ng/ul  | 97     |
| 69) Phenanthrene           | 17.23 | 178 | 208432  | 3.05 | ng/ul  | 98     |
| 74) Fluoranthene           | 19.24 | 202 | 523948  | 6.35 | ng/ul# | 96     |
| 77) Pyrene                 | 19.61 | 202 | 592706  | 7.28 | ng/ul# | 95     |
| 80) Benzo(a)anthracene     | 21.35 | 228 | 299598  | 3.74 | ng/ul  | 97     |
| 82) Chrysene               | 21.40 | 228 | 330067  | 4.38 | ng/ul  | 96     |
| 85) Benzo(b)fluoranthene   | 22.97 | 252 | 446582  | 7.30 | ng/ul# | 97     |
| 86) Benzo(k)fluoranthene   | 23.01 | 252 | 140024m | 2.37 | ng/ul  |        |
| 88) Benzo(a)pyrene         | 23.56 | 252 | 267604  | 4.64 | ng/ul# | 95     |
| 89) Indeno(1,2,3-cd)pyrene | 26.00 | 276 | 210948  | 3.44 | ng/ul  | 97     |
| 91) Benzo(g,h,i)perylene   | 26.73 | 276 | 189464  | 3.78 | ng/ul  | 95     |

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

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