

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070319\  
 Data File : BE100338.D  
 Acq On : 3 Jul 2019 20:52  
 Operator : JU/SJ  
 Sample : K3561-06 5X  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_E  
 ClientSampleId :  
 EXT-016-SW098-062519

Manual Integrations  
 APPROVED

mohammad  
 7/5/2019 8:54:13 AM

Quant Time: Jul 04 01:35:06 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062819.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Jun 28 14:06:53 2019

Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.63  | 152  | 632      | 0.40 | ng    | -0.01    |
| 7) Naphthalene-d8         | 10.43 | 136  | 2512     | 0.40 | ng    | -0.01    |
| 13) Acenaphthene-d10      | 14.31 | 164  | 2057     | 0.40 | ng    | 0.00     |
| 19) Phenanthrene-d10      | 17.05 | 188  | 5731     | 0.40 | ng    | 0.00     |
| 27) Chrysene-d12          | 21.24 | 240  | 7854     | 0.40 | ng    | 0.00     |
| 34) Perylene-d12          | 23.67 | 264  | 9011     | 0.40 | ng    | 0.00     |

#### System Monitoring Compounds

|                             |       |     |      |      |    |       |
|-----------------------------|-------|-----|------|------|----|-------|
| 4) 2-Fluorophenol           | 5.20  | 112 | 74   | 0.05 | ng | -0.01 |
| 5) Phenol-d6                | 6.85  | 99  | 58m  | 0.03 | ng | 0.00  |
| 8) Nitrobenzene-d5          | 8.82  | 82  | 107  | 0.04 | ng | 0.00  |
| 11) 2-Methylnaphthalene-d10 | 12.04 | 152 | 235  | 0.05 | ng | -0.01 |
| 14) 2,4,6-Tribromophenol    | 15.81 | 330 | 67   | 0.05 | ng | -0.01 |
| 15) 2-Fluorobiphenyl        | 12.93 | 172 | 360  | 0.04 | ng | -0.01 |
| 25) Fluoranthene-d10        | 19.09 | 212 | 3368 | 0.04 | ng | 0.00  |
| 29) Terphenyl-d14           | 19.69 | 244 | 2103 | 0.13 | ng | -0.01 |

#### Target Compounds

|                                |       |     |        |        |    | Qvalue |
|--------------------------------|-------|-----|--------|--------|----|--------|
| 9) Naphthalene                 | 10.48 | 128 | 4932   | 0.178  | ng | # 97   |
| 12) 2-Methylnaphthalene        | 12.11 | 142 | 1062   | 0.218  | ng | 91     |
| 16) Acenaphthylene             | 14.02 | 152 | 7997   | 0.810  | ng | 99     |
| 17) Acenaphthene               | 14.37 | 154 | 791    | 0.141  | ng | 96     |
| 18) Fluorene                   | 15.35 | 166 | 2648   | 0.320  | ng | # 88   |
| 23) Phenanthrene               | 17.09 | 178 | 69300  | 4.989  | ng | 100    |
| 24) Anthracene                 | 17.19 | 178 | 7141   | 0.571  | ng | 96     |
| 26) Fluoranthene               | 19.12 | 202 | 135270 | 6.114  | ng | 99     |
| 28) Pyrene                     | 19.48 | 202 | 169329 | 8.143  | ng | 97     |
| 30) Benzo(a)anthracene         | 21.22 | 228 | 91377  | 3.538  | ng | 97     |
| 31) Chrysene                   | 21.28 | 228 | 108052 | 4.167  | ng | 98     |
| 32) Bis(2-ethylhexyl)phthalate | 21.15 | 149 | 402743 | 10.977 | ng | # 97   |
| 33) Indeno(1,2,3-cd)pyrene     | 26.22 | 276 | 59859  | 1.745  | ng | # 97   |
| 35) Benzo(b)fluoranthene       | 22.92 | 252 | 111871 | 4.579  | ng | # 96   |
| 36) Benzo(k)fluoranthene       | 22.96 | 252 | 37743m | 1.384  | ng |        |
| 37) Benzo(a)pyrene             | 23.55 | 252 | 79723m | 3.256  | ng |        |
| 38) Dibenzo(a,h)anthracene     | 26.22 | 278 | 15212  | 0.595  | ng | # 89   |
| 39) Benzo(a,h,i)perylene       | 27.01 | 276 | 67839  | 2.582  | ng | 98     |

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

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