

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\  
 Data File : BM024533.D  
 Acq On : 18 Jan 2020 13:37  
 Operator : JU  
 Sample : L1109-05  
 Misc :  
 ALS Vial : 37 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 GASG4

Manual Integrations  
 APPROVED

mohammad  
 1/21/2020 7:59:48 AM

Quant Time: Jan 20 03:42:24 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jan 20 03:25:57 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.46  | 152  | 205981   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.22 | 136  | 905047   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.10 | 164  | 664368   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.85 | 188  | 1569989  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.06 | 240  | 1690888  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.17 | 264  | 1828707  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.06  | 96  | 11822   | 2.80  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.65  | 99  | 193743  | 12.69 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.80  | 67  | 124902  | 14.73 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 6.99  | 132 | 187398  | 14.37 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.16  | 113 | 169278  | 12.02 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.59  | 128 | 102126  | 16.06 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.31  | 143 | 110299  | 16.02 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.85  | 165 | 211139  | 13.62 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.35 | 131 | 242653  | 14.08 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.52 | 166 | 816271  | 15.68 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.79 | 160 | 889945  | 14.85 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.31 | 143 | 117764  | 13.55 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.10 | 176 | 731082  | 15.95 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.22 | 200 | 115170  | 12.85 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.95 | 188 | 1216990 | 16.67 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.25 | 212 | 1487334 | 16.46 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.04 | 264 | 1548078 | 16.32 | ng/ul | 0.00 |

## Target Compounds

|                         |       |     |         |        | Ovalue    |
|-------------------------|-------|-----|---------|--------|-----------|
| 28) Naphthalene         | 10.26 | 128 | 1261577 | 27.209 | ng/ul 100 |
| 34) 2-Methylnaphthalene | 11.89 | 142 | 780243  | 21.687 | ng/ul 100 |
| 41) 1,1'-Biphenyl       | 12.93 | 154 | 78665   | 1.631  | ng/ul 99  |
| 45) Dimethylphthalate   | 13.57 | 163 | 243697  | 4.605  | ng/ul 99  |
| 48) Acenaphthylene      | 13.82 | 152 | 100129  | 1.636  | ng/ul 96  |
| 54) Dibenzofuran        | 14.50 | 168 | 72120   | 1.189  | ng/ul 93  |
| 59) Fluorene            | 15.16 | 166 | 184627  | 3.599  | ng/ul 99  |
| 70) Phenanthrene        | 16.89 | 178 | 1182399 | 13.965 | ng/ul 100 |
| 72) Anthracene          | 16.99 | 178 | 221228  | 2.562  | ng/ul 99  |
| 77) Fluoranthene        | 18.92 | 202 | 382025  | 3.692  | ng/ul 100 |
| 80) Pyrene              | 19.28 | 202 | 478210  | 4.131  | ng/ul 99  |
| 83) Benzo(a)anthracene  | 21.04 | 228 | 183740m | 1.595  | ng/ul     |
| 85) Chrysene            | 21.09 | 228 | 143546  | 1.293  | ng/ul 97  |

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

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