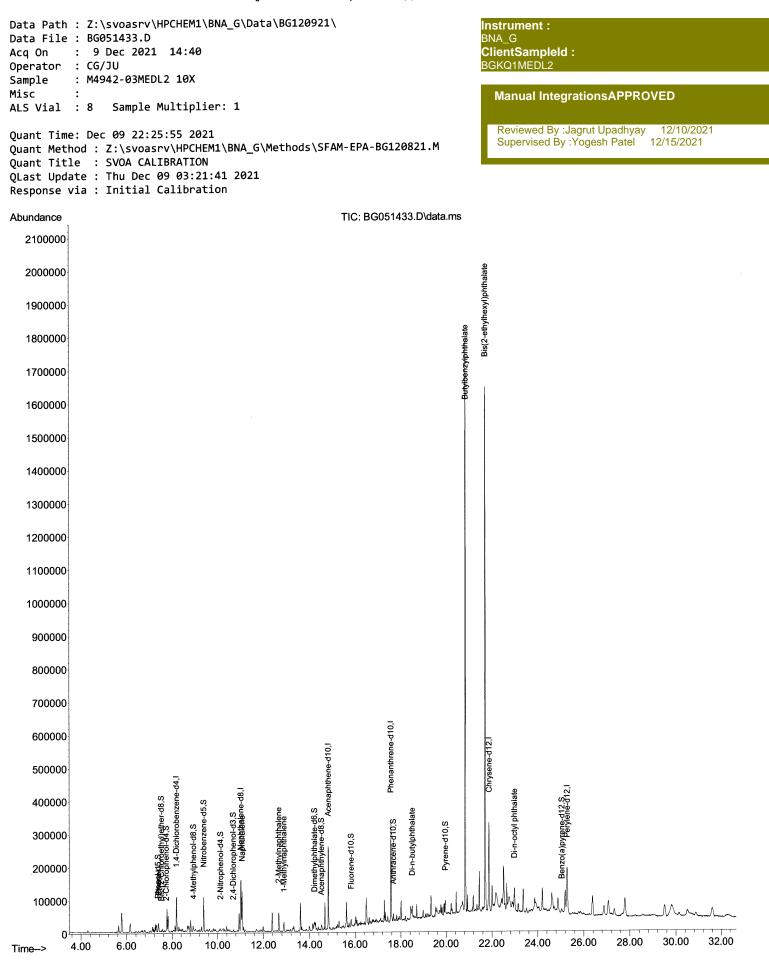
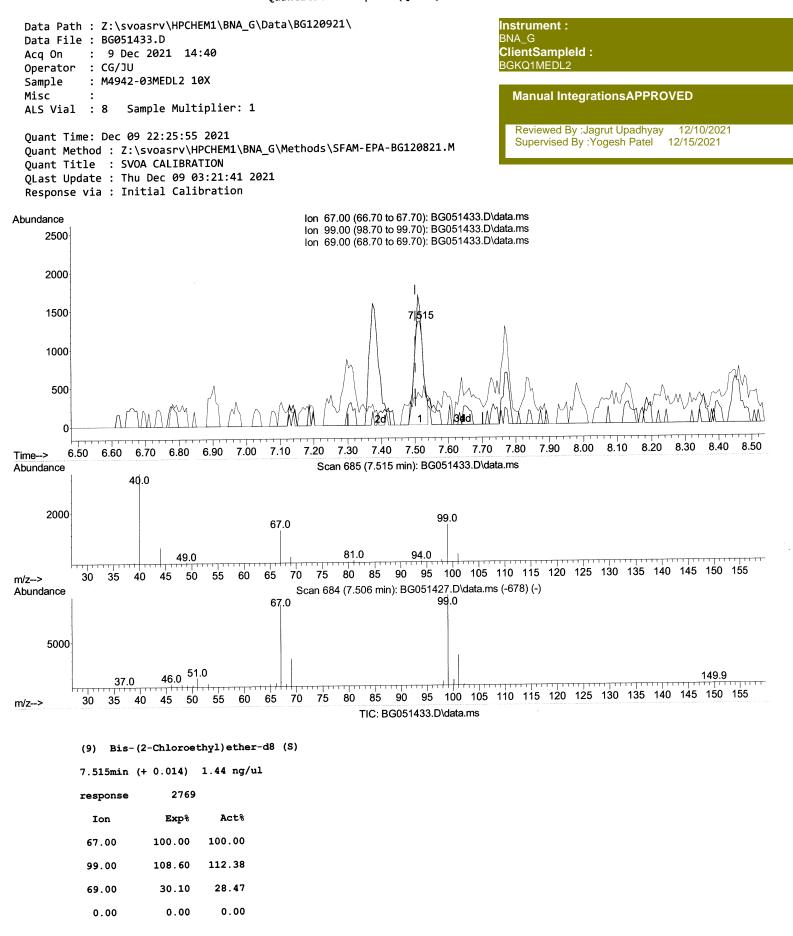
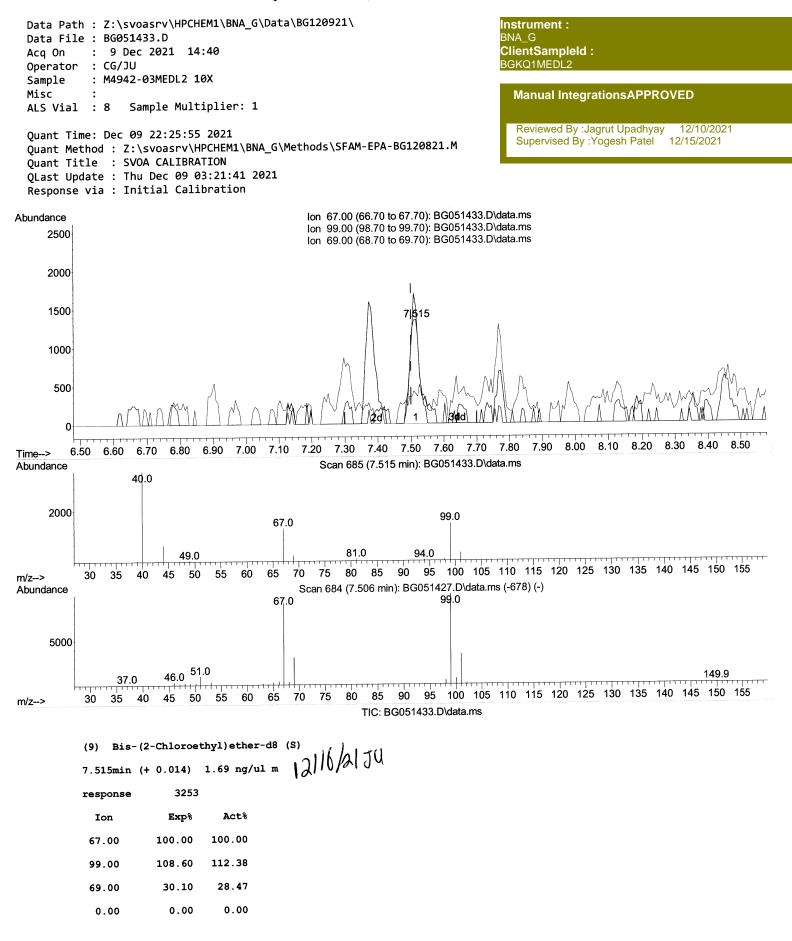
(QT Reviewed)

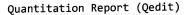


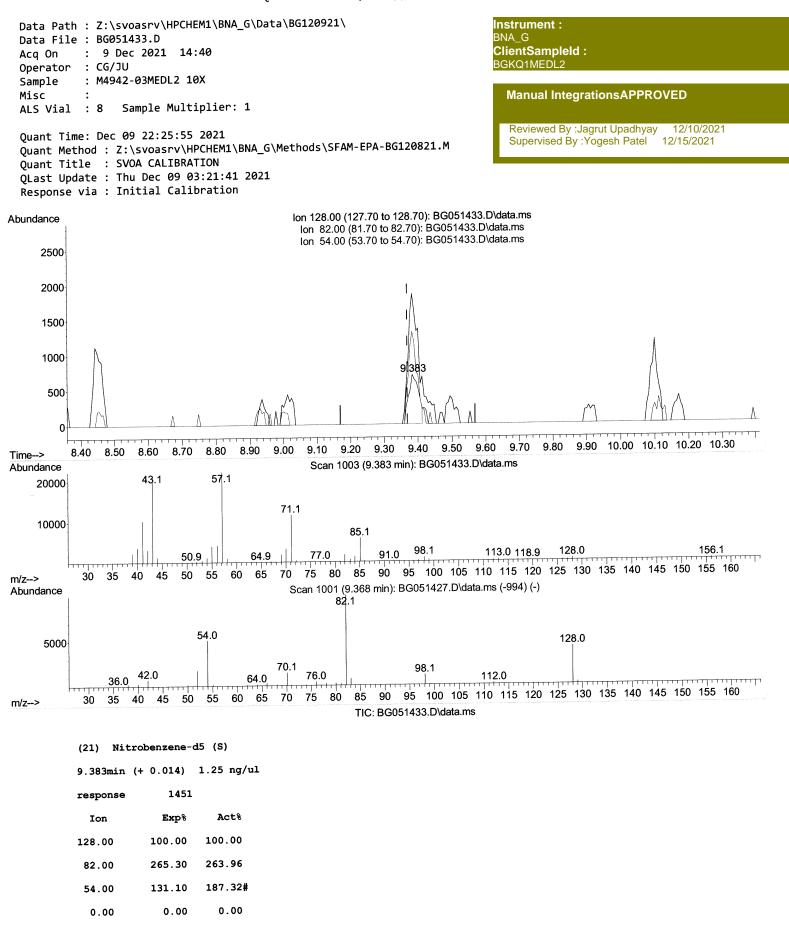




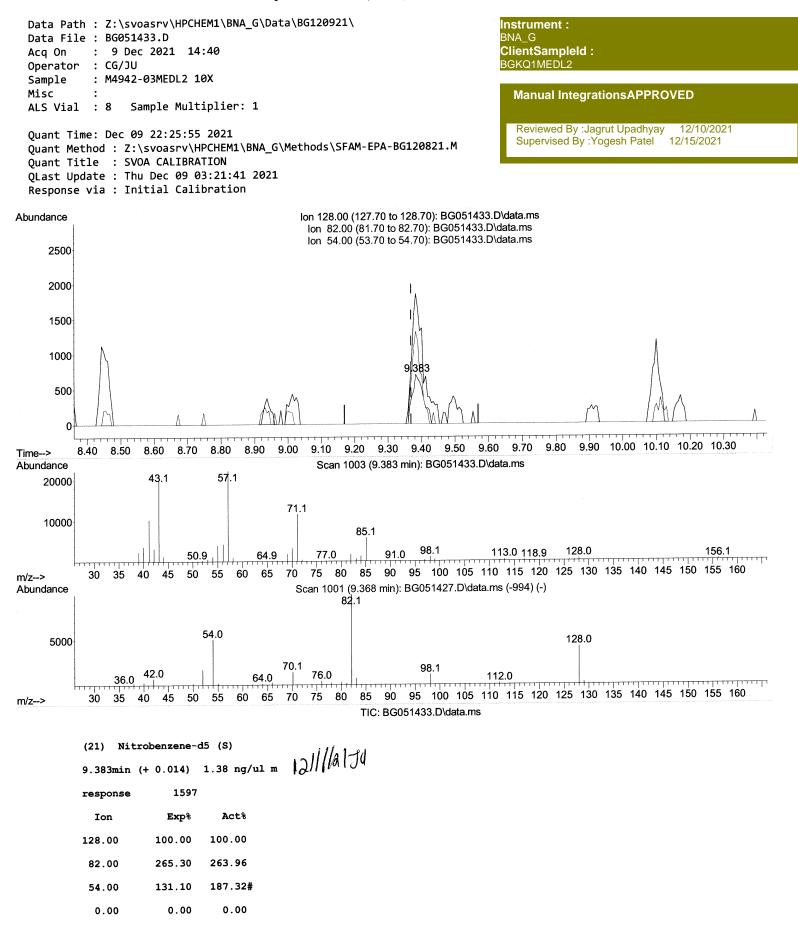


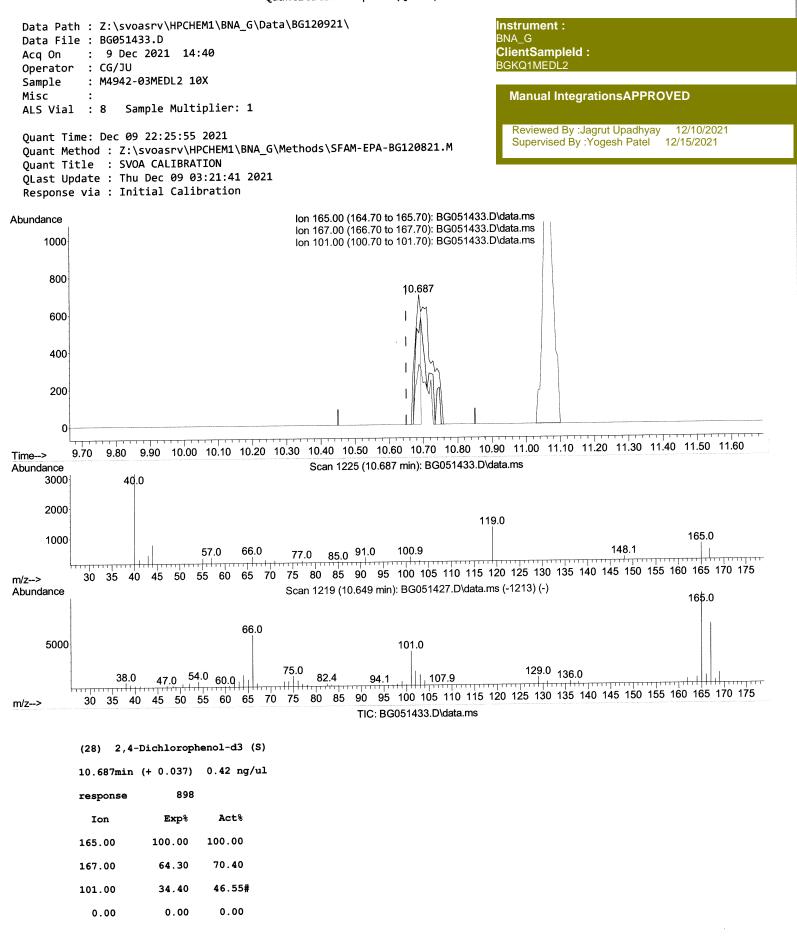




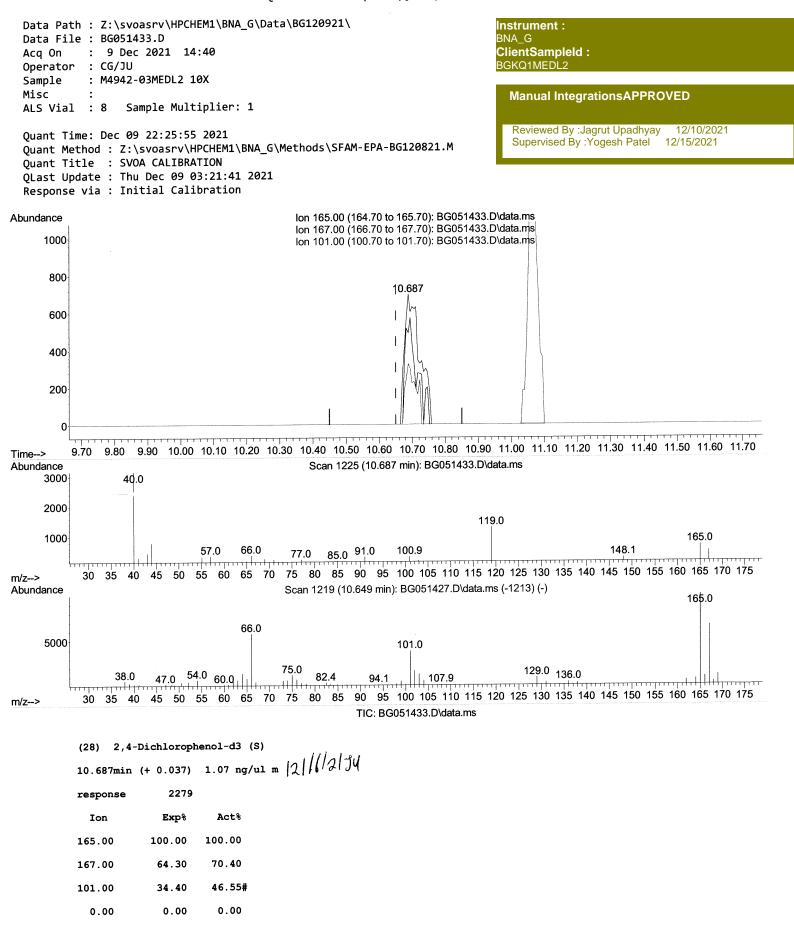




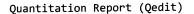


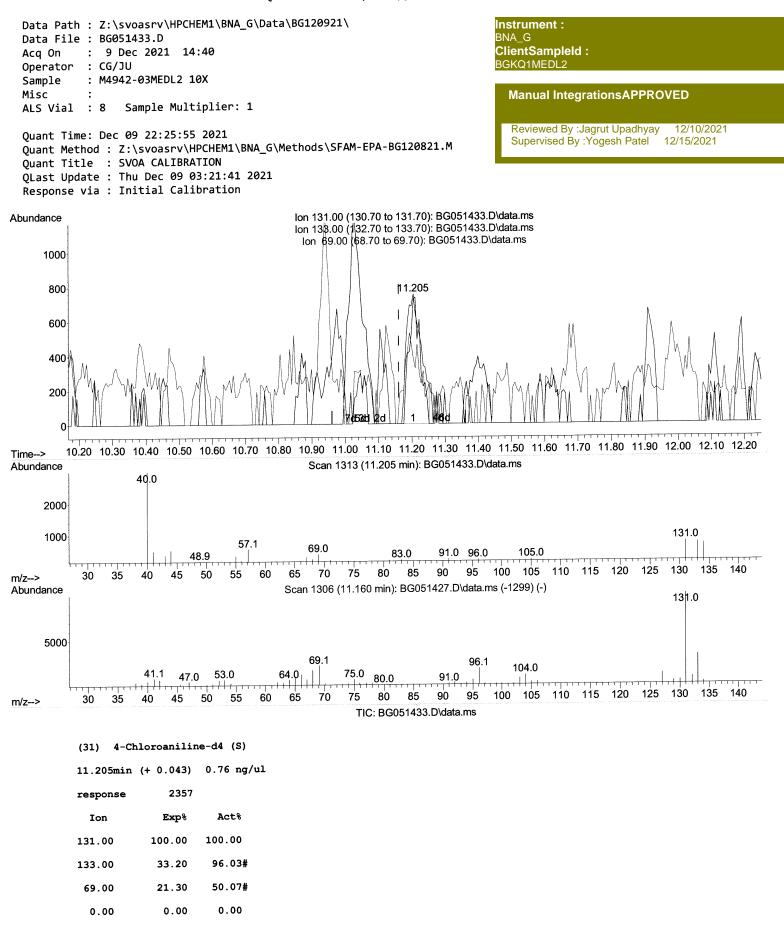


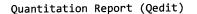


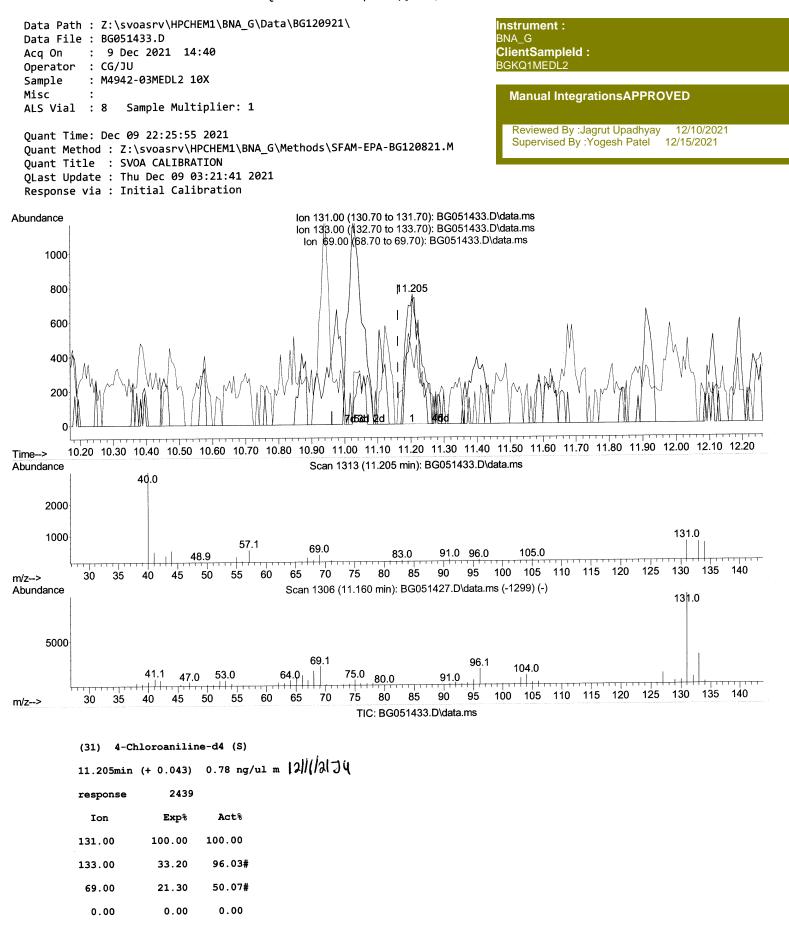


SFAM-EPA-BG120821.M Thu Dec 09 23:35:34 2021









| Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\ Data File : BG051433.D Acq On : 9 Dec 2021 14:40 Operator : CG/JU Sample : M4942-03MEDL2 10X | | | | | | Instrument : BNA_G ClientSampleId : BGKQ1MEDL2 | |
|--|---|---------------------|------------|--------------------|------------------|---|---|
| Misc : ALS Vial : 8 Sample Multiplier: 1 | | | | | | Manual IntegrationsAPPROVED | |
| Quant Quant QLast | Time: Dec 09 22:25:55 2021 Method : Z:\svoasrv\HPCHEM Title : SVOA CALIBRATION Update : Thu Dec 09 03:21: Ise via : Initial Calibrati | 1\BNA_G\ 41 2021 | \Metho | ds∖SFAM-EPA | -BG12082 | 21.M | Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :Yogesh Patel 12/15/2021 |
| | Compound | | | Response | | | |
| Inter | nal Standards | | | | | | |
| 1) | 1,4-Dichlorobenzene-d4 | 8.185 | | 29540 | 20.000 | - | 0.00 |
| 20) | Naphthalene-d8 | 11.017 | 136 | 133586 | 20.000 | | 0.00 |
| 38) | Acenaphthene-d10 | 14.818 | | 89914 | 20.000 | - | 0.00 |
| 64) | Phenanthrene-d10 | 17.574 | | 186173 | 20.000 | - | 0.00 |
| | Chrysene-d12 | 21.874 | | 163763 | 20.000 | - | 0.00 |
| 88) | Perylene-d12 | 25.270 | 264 | 162904 | 20.000 | ng/ul | 0.00 |
| Syste | em Monitoring Compounds | | | | | | |
| 3) | 1,4-Dioxane-d8 | 0.000 | 96 | 0 | | ng/uL | |
| 4) | Pyridine-d5 | 0.000 | 84 | 0d | | ng/ul | |
| • | Phenol-d5 | 7.374 | 99 | 3166 | 1.053 | ng/ul | 0.02 11/12/34 |
| 9) | Bis-(2-Chloroethyl)eth | 7.515 | 67 | 3253m > | 1.687 | ng/ul > | 0.02 0.01 12/16/21 34 |
| 11) | 2-Chlorophenol-d4 | 7.738 | 132 | 2688 | 1.250 | ng/ur | 0.01 |
| 15) | 4-Methylphenol-d8 | 8.937 | 113 | 2845 | | ng/ul | 0.03 0.01 12/16/21 JU |
| 21) | Nitrobenzene-d5 | 9.383 | 128 | 1597m > | | ng/ul> | |
| | 2-Nitrophenol-d4 | 10.118 | | 1364 | | ng/ul | 0.02 |
| | 2,4-Dichlorophenol-d3 | 10.687 | 165 | ^{2279m} > | | ng/ul | |
| 31) | 4-Chloroaniline-d4 | 11.205 | 131 | 2439m 🖍 | | ng/ul 1 | 0.04 |
| 46) | Dimethylphthalate-d6 | 14.219 | 166 | 10068 | | ng/ul | 0.00 |
| 49) | Acenaphthylene-d8 | 14.518 | | 13084 | | ng/ul | 0.00 |
| 54) | 4-Nitrophenol-d4 | 0.000 | | 0 | | ng/ul | |
| | Fluorene-d10 | 15.817 | | 8945 | | ng/ul | 0.00 |
| 65) | 4,6-Dinitro-2-methylph | | | 0d | | ng/ul | 0.00 |
| 73) | Anthracene-d10 | 17.673 | | 14897 | | ng/ul | 0.00 |
| | Pyrene-d10 | 19.953 | | 15382 | | ng/ul | 0.00 |
| 92) | Benzo(a)pyrene-d12 | 25.047 | 264 | 12742 | 1.516 | ng/ul | 0.01 |
| Target compounds | | | | | lue | | |
| 8) | Phenol | 7.403 | 94 | 13092 | | ng/ul | 96 |
| 30) | Naphthalene | 11.064 | 128 | 104580 | 14.256 | - | 98 |
| 36) | 2-Methylnaphthalene | 12.662 | 142 | 28393 | | ng/ul | 96 |
| 37) | 1-Methylnaphthalene | 12.879 | | 13763 | | ng/ul | 99 |
| 78) | Di-n-butylphthalate | 18.502 | | 21850 | | ng/ul | 99 |
| 83) | Butylbenzylphthalate | 20.840 | | 501300 | | - | |
| 86) | Bis(2-ethylhexyl)phtha | 21.710 | | 687466 | 95.417 | | 99 |
| | Di-n-octyl phthalate | 22.973 | 149 | 88568 | 7.393 | ng/ul | 100 |
| 83) 86) | Butylbenzylphthalate Bis(2-ethylhexyl)phtha | 20.840 21.710 | 149 149 | 501300 687466 | 96.642 95.417 | ng/ul | 96 99 |

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