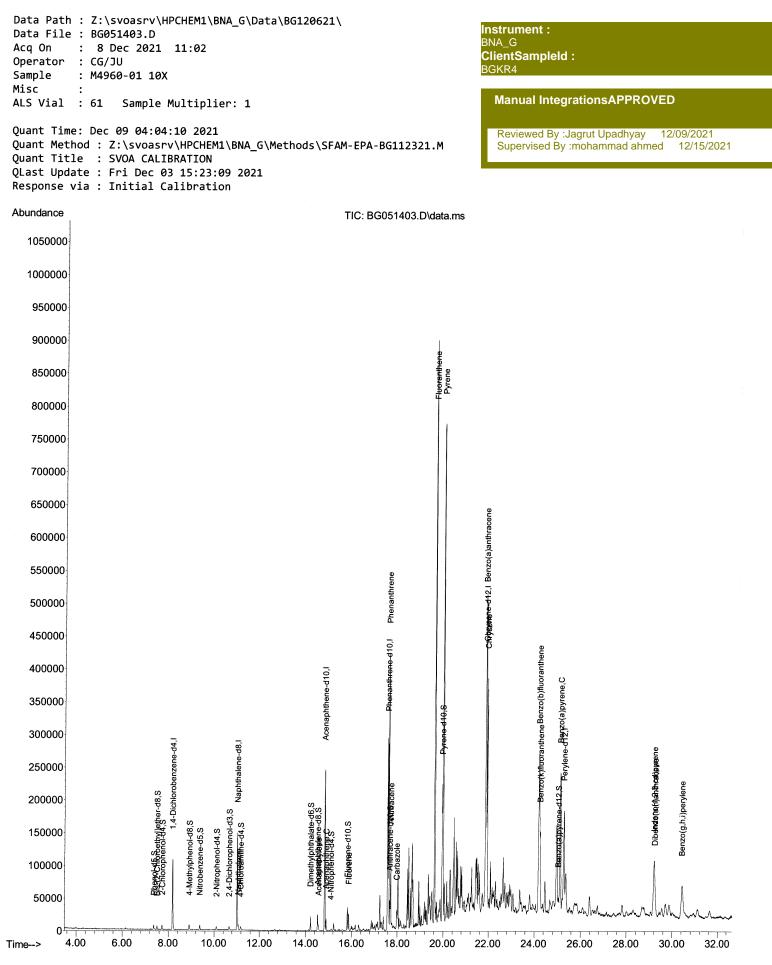
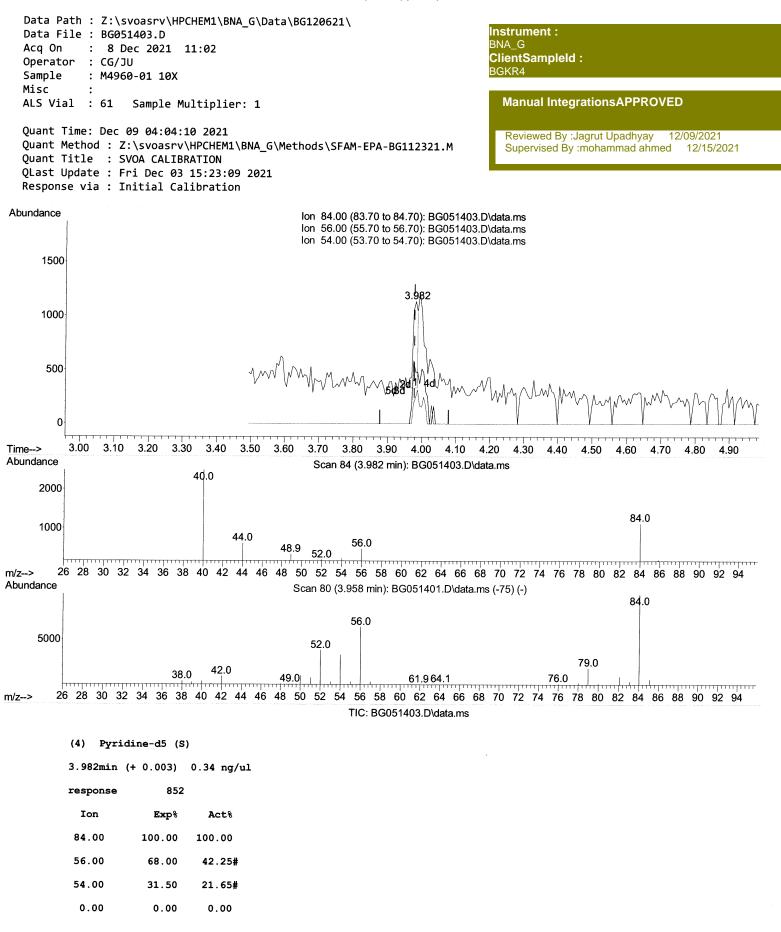
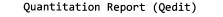
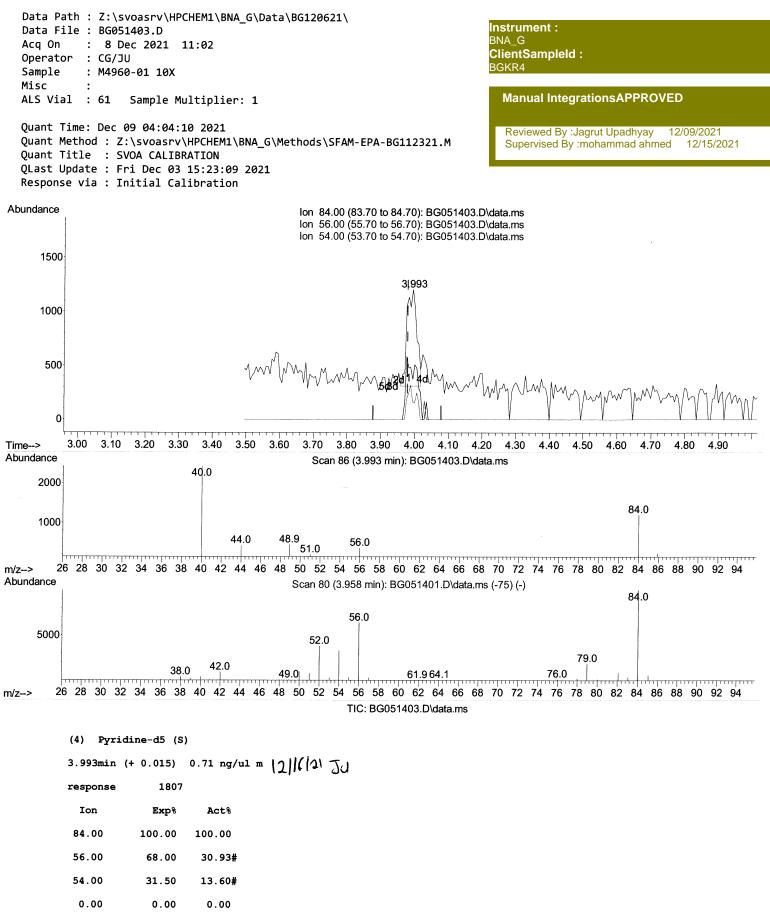
(QT Reviewed)



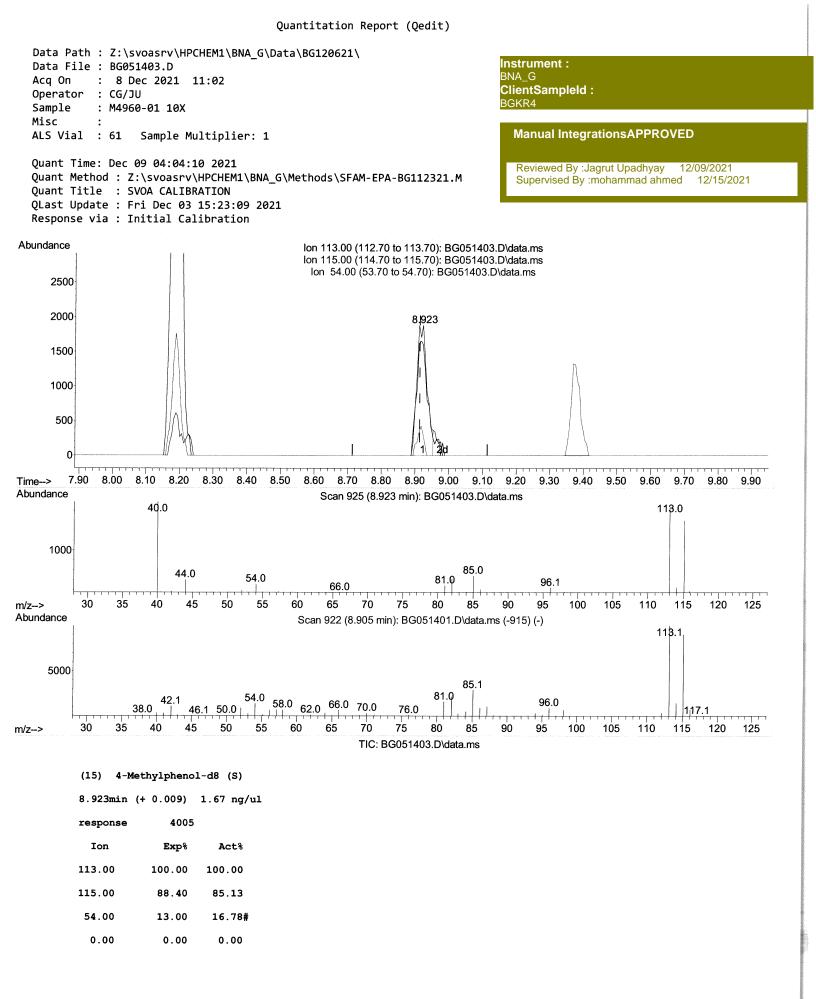


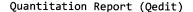


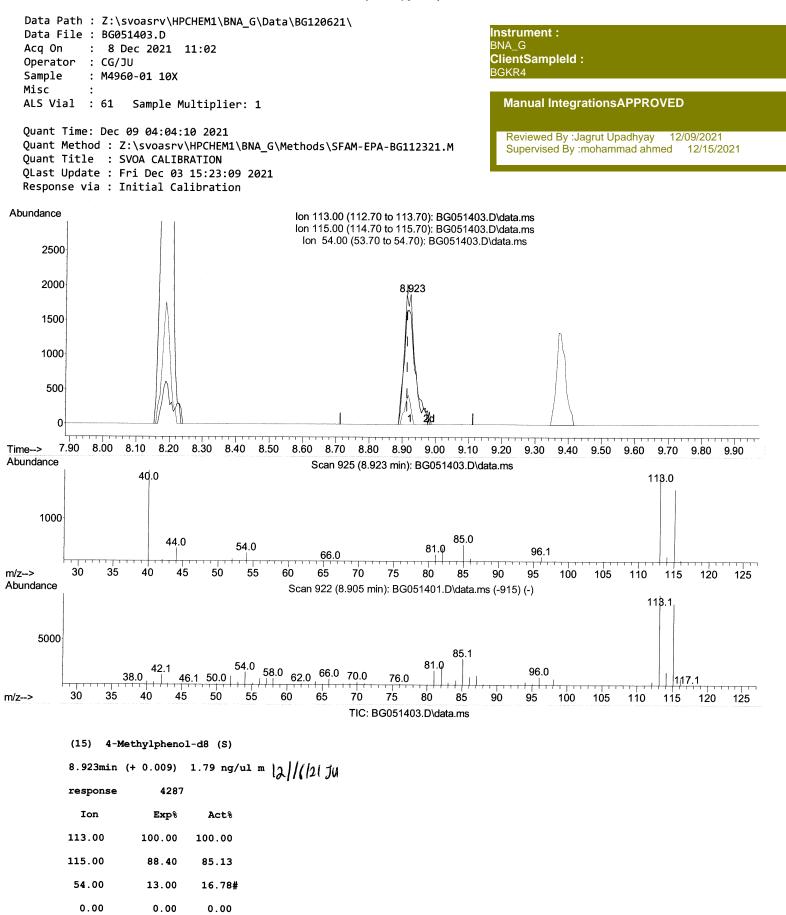


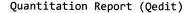


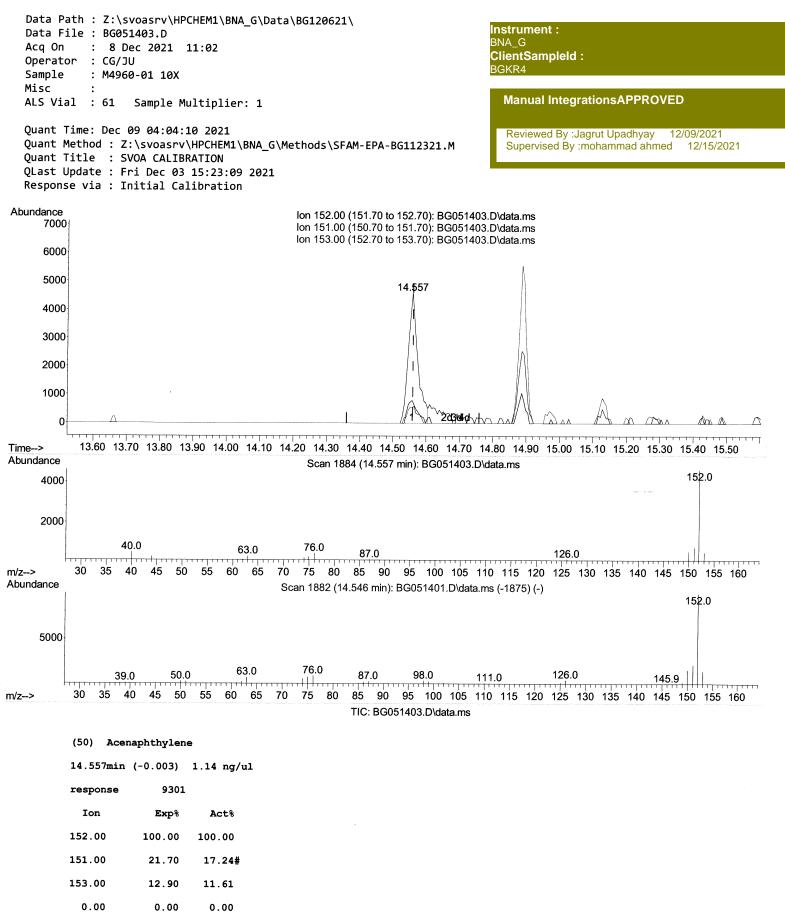
SFAM-EPA-BG112321.M Thu Dec 09 04:21:42 2021

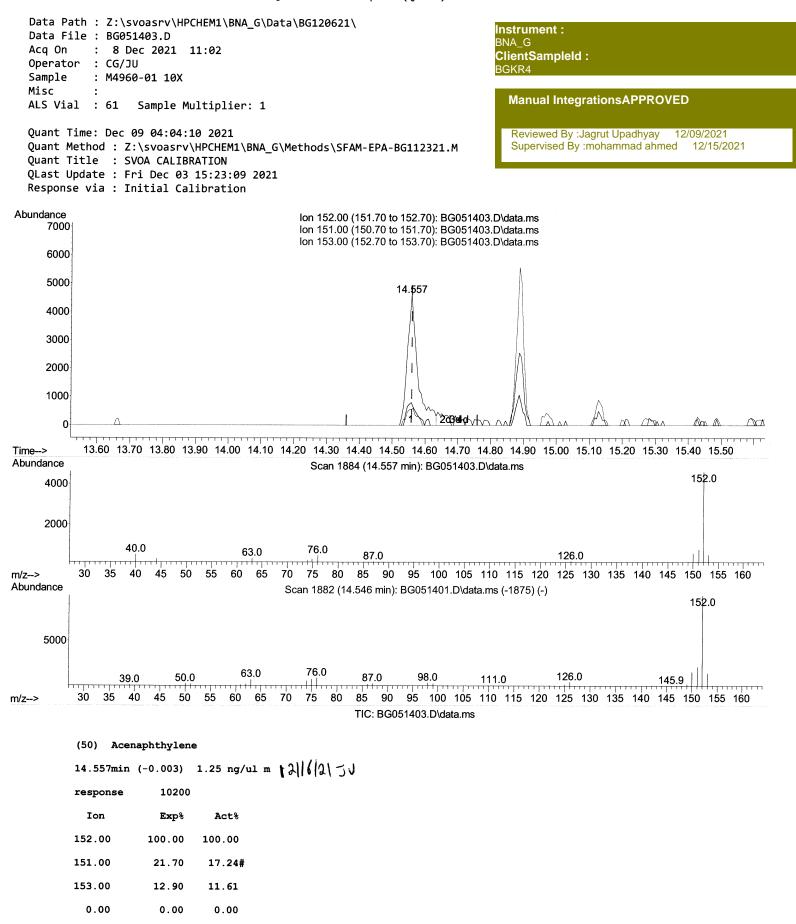




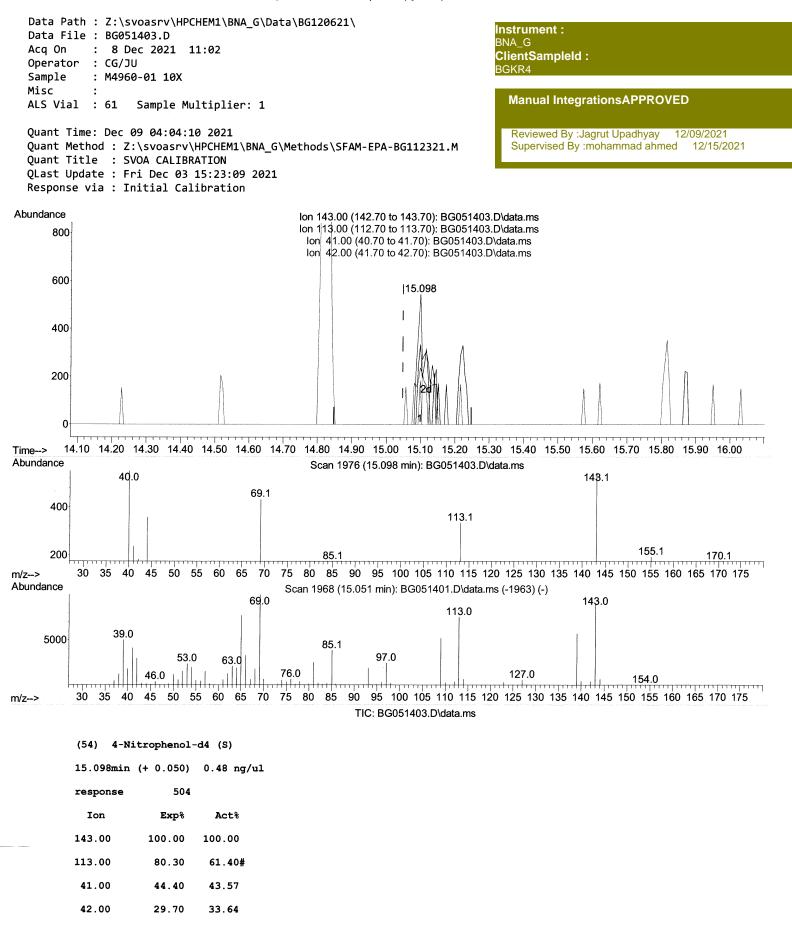




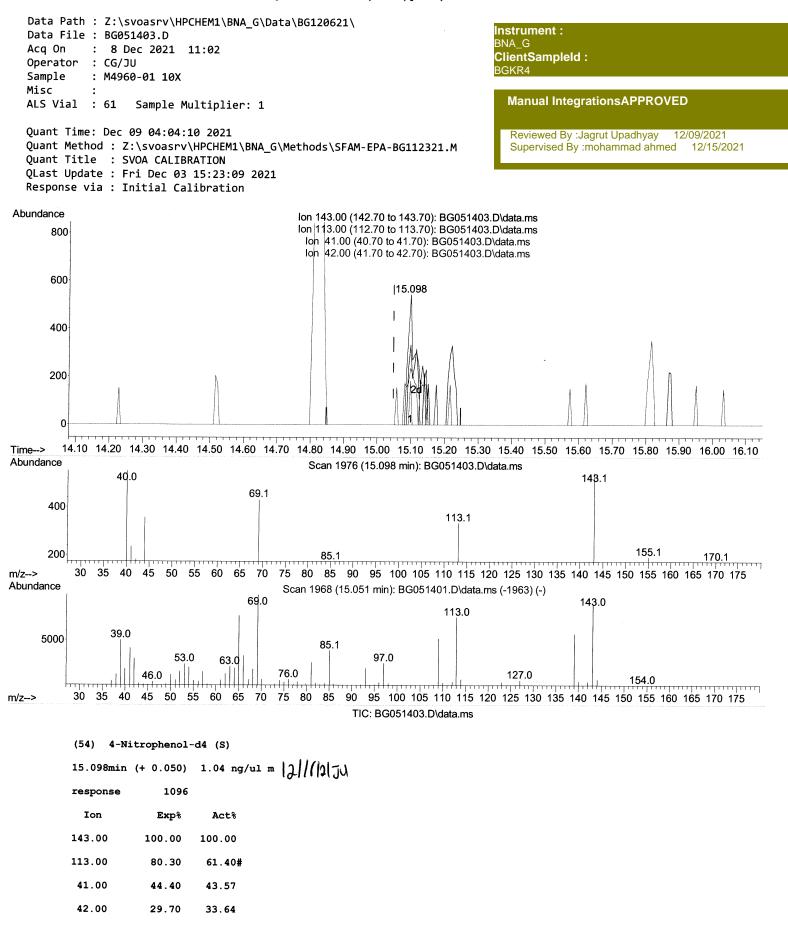




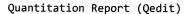


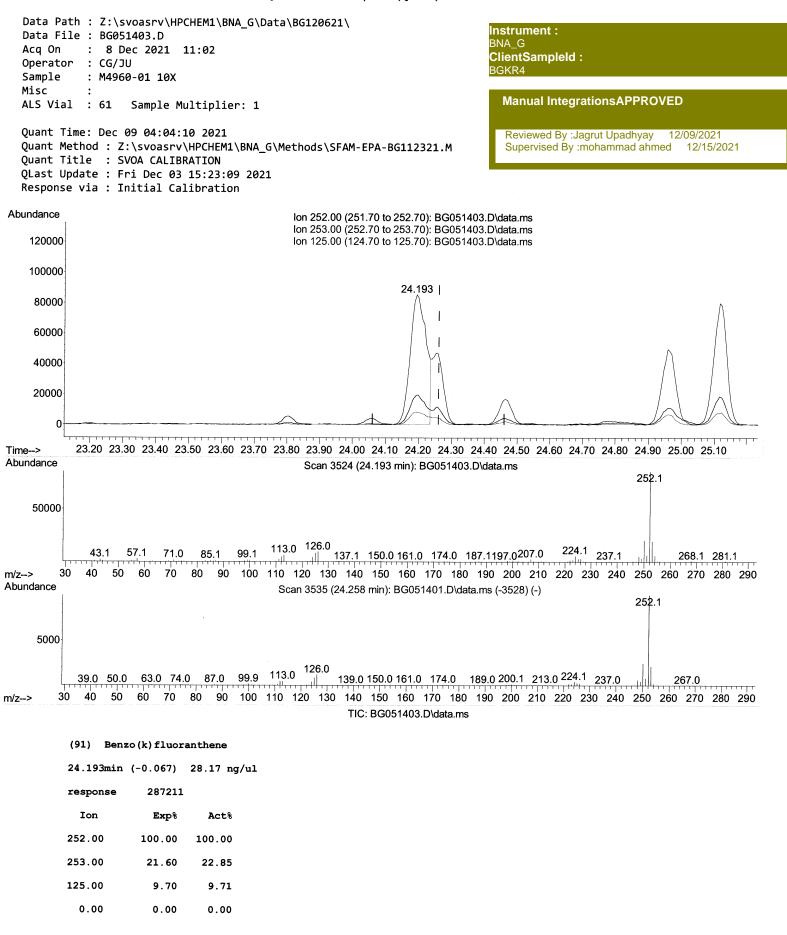




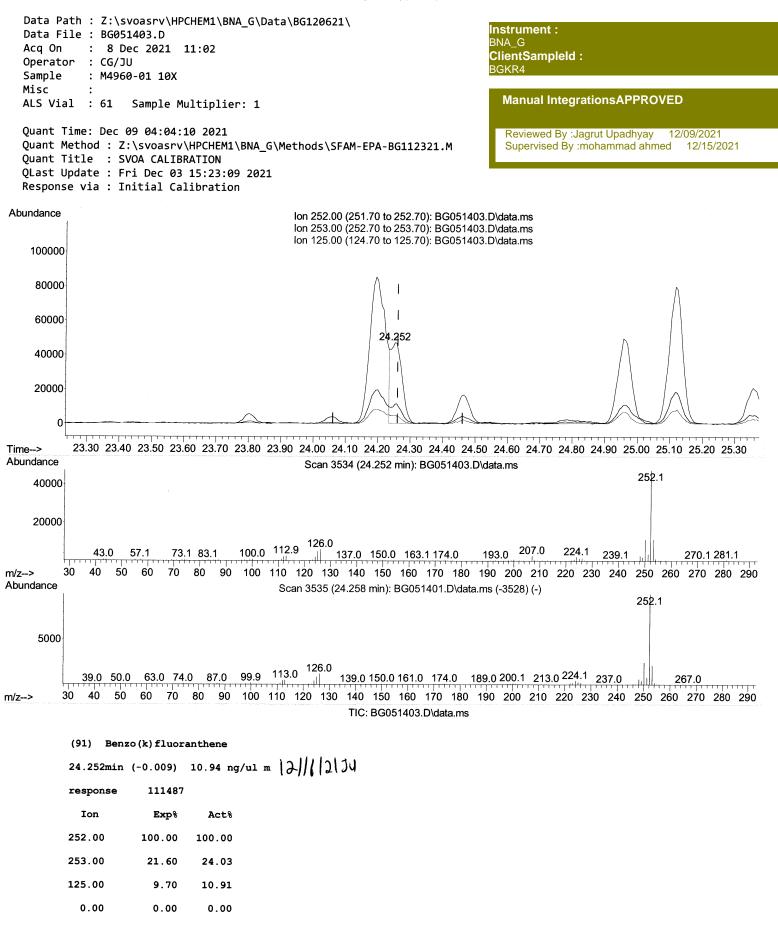


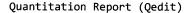
SFAM-EPA-BG112321.M Thu Dec 09 04:23:14 2021

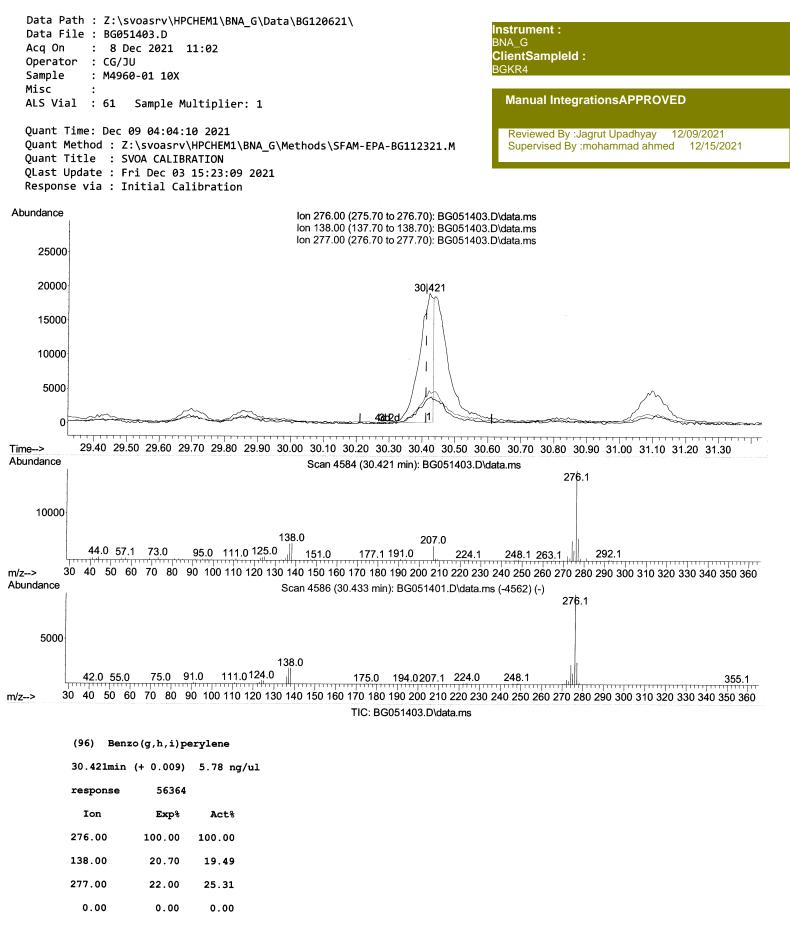




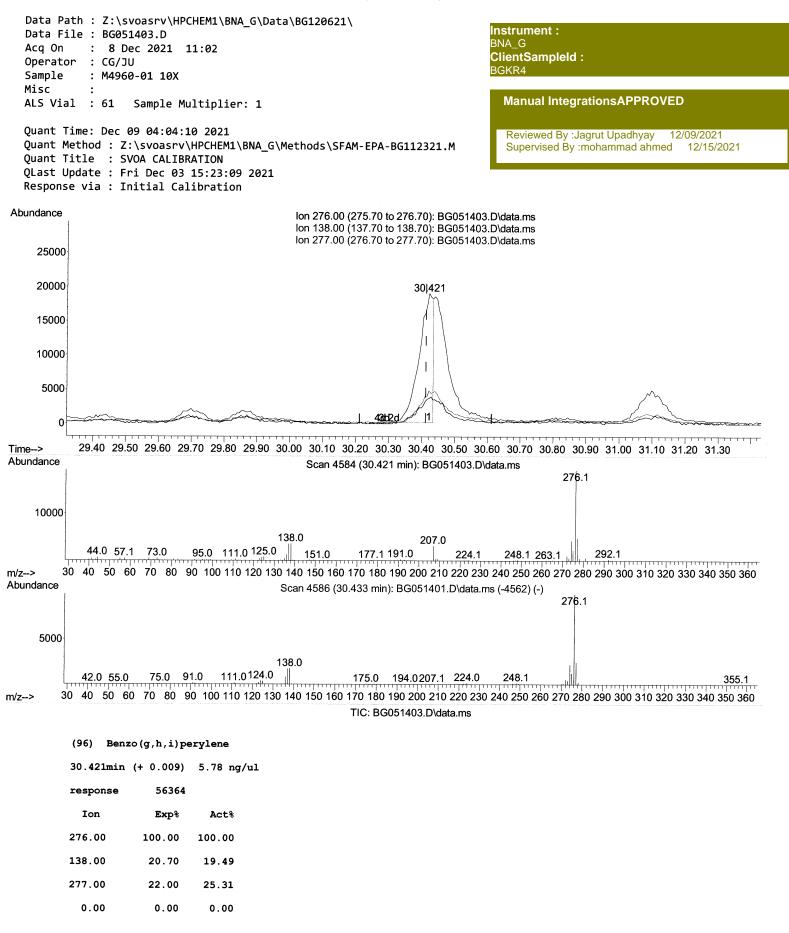


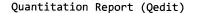


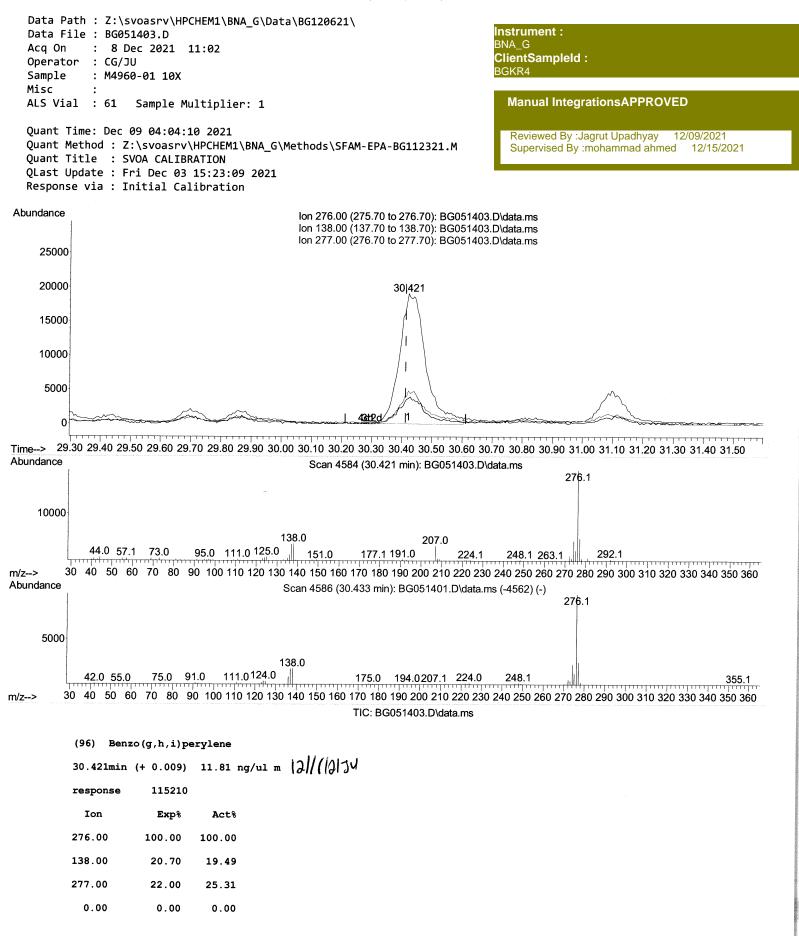












|   | <b>C</b>                 |  |  |
|---|--------------------------|--|--|
| Data Path : Z:\svoasrv\HPCHEM1                                    | BNA_G\Data\B             | G120621\                                   | lu of rum on f                           |
| Data File : BG051403.D  |                          |  | Instrument :<br>BNA G                    |
| Acq On : 8 Dec 2021 11:02   |                          |  | ClientSampleId :                         |
| Operator : CG/JU<br>Sample : M4960-01 10X                         |                          |  | BGKR4                                    |
| Sample : M4960-01 10X<br>Misc :                                   |                          |  |  |
| ALS Vial : 61 Sample Multipl                                      | ier: 1                   |  | Manual IntegrationsAPPROVED              |
|   |                          |  |  |
| Quant Time: Dec 09 04:04:10 202                                   |                          |  | Reviewed By :Jagrut Upadhyay 12/09/2021  |
| Quant Method : Z:\svoasrv\HPCHE<br>Quant Title : SVOA CALIBRATION | MI\BNA_G\Met             | nods\SFAM-EPA-BG112321.M                   | Supervised By :mohammad ahmed 12/15/2021 |
| QLast Update : Fri Dec 03 15:23                                   |                          |  |  |
| Response via : Initial Calibrat                                   |                          |  |  |
|   |                          |  |  |
| Compound  | R.T. QIO                 | n Response Conc Units Dev                  | (Min)                                    |
| Internal Standards  |                          |  |  |
| 1) 1,4-Dichlorobenzene-d4   | 8.188 15                 | 2 30042 20.000 ng/ul                       | -0.01                                    |
| 20) Naphthalene-d8  | 11.020 13                | 0.   | 0.00                                     |
| 38) Acenaphthene-d10  | 14.822 16                | •  | 0.00                                     |
| 64) Phenanthrene-d10  | 17.577 18                | -  | 0.00                                     |
| 79) Chrysene-d12  | 21.878 24                | 0 158319 20.000 ng/ul                      | 0.00                                     |
| 88) Perylene-d12  | 25.280 26                | 160964 20.000 ng/ul                        | 0.00                                     |
| System Monitoring Compounds                                       |                          |  |  |
| 3) 1,4-Dioxane-d8   | 0.000 9                  | 5 0 0.000 ng/uL                            |  |
| 4) Pyridine-d5  | 3.993 84                 | 1807m ≥ 0.712 ng/ul>                       | 0.01 ja/16/2174                          |
| 7) Phenol-d5  | 7.366 99                 | 9 5563 1.874 ng/ul                         | 0.00                                     |
| <pre>9) Bis-(2-Chloroethyl)eth</pre>                              | 7.507 6                  | -  | 0.00                                     |
| <pre>11) 2-Chlorophenol-d4</pre>                                  | 7.730 132                | 2 3937 1.841 ng/ul                         | 0.00                                     |
| <pre>15) 4-Methylphenol-d8</pre>                                  | 8.923 113                | 8 4287m> 1.789 ng/ul>                      | 0.001211/12134                           |
| 21) Nitrobenzene-d5   | 9.375 128                | 8 1795 1.620 ng/ul                         | 0.00                                     |
| 24) 2-Nitrophenol-d4  | 10.104 143               | 0,   | 0.00                                     |
| 28) 2,4-Dichlorophenol-d3   | 10.668 165               | 0,   | 0.01                                     |
| 31) 4-Chloroaniline-d4<br>46) Dimethylphthalate-d6                | 11.173 131<br>14.217 166 | 0,   | 0.00                                     |
| 49) Acenaphthylene-d8   | 14.522 166               | 0,   | 0.00<br>0.00                             |
| 54) 4-Nitrophenol-d4  | 15.098 143               | 1096m > 1.035  mg/ul                       | - 0.05 (2111)1)                          |
| 60) Fluorene-d10  | 15.815 176               |  | 0.00                                     |
| 65) 4,6-Dinitro-2-methylph  |                          |  |  |
| 73) Anthracene-d10  | 17.671 188               | -  | 0.00                                     |
| 81) Pyrene-d10  | 19.951 212               | 0,   | 0.00                                     |
| 92) Benzo(a)pyrene-d12  | 25.045 264               | 22368 2.602 ng/ul                          | 0.00                                     |
| Target Compounds  |                          | 0.97                                       | alue                                     |
| 30) Naphthalene   | 11.067 128               |  | 94                                       |
| 50) Acenaphthylene  | 14.557 152               |  |  |
| 52) Acenaphthene  | 14.886 153               | -  | 97                                       |
| 61) Fluorene  | 15.874 166               | -  | 95                                       |
| 72) Phenanthrene  | 17.619 178               | 261358 26.042 ng/ul                        | 100                                      |
| 74) Anthracene  | 17.707 178               | 0,   | 98                                       |
| 77) Carbazole   | 17.983 167               | 0.   | 96                                       |
| 80) Fluoranthene  | 19.622 202               | 0,   | 97                                       |
| 82) Pyrene<br>85) Benzo(a)anthracene                              | 19.986 202<br>21.855 228 | 0,   | 97                                       |
| 87) Chrysene  | 21.925 228               | 277392 25.833 ng/ul<br>249390 24.176 ng/ul | 99<br>98                                 |
| 90) Benzo(b)fluoranthene  | 24.193 252               | 287211 26.440 ng/ul                        | 98                                       |
| 91) Benzo(k)fluoranthene  | 24.252 252               | 111487m ≻ 10.937 ng/ul >                   |  |
| 93) Benzo(a)pyrene  | 25.116 252               | 225764 21.785 ng/ul                        | 96                                       |
| 94) Indeno(1,2,3-cd)pyrene  | 29.199 276               | 142388 12.278 ng/ul                        | 97                                       |
| 95) Dibenzo(a,h)anthracene  | 29.228 278               | 35767 3.635 ng/ul                          | 93                                       |
| 96) Benzo(g,h,i)perylene  | 30.421 276               | 115210m > 11.808 ng/ul =                   | > [0][(1012)                             |
|   |                          |  |  |
|   |                          |  |  |

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

1