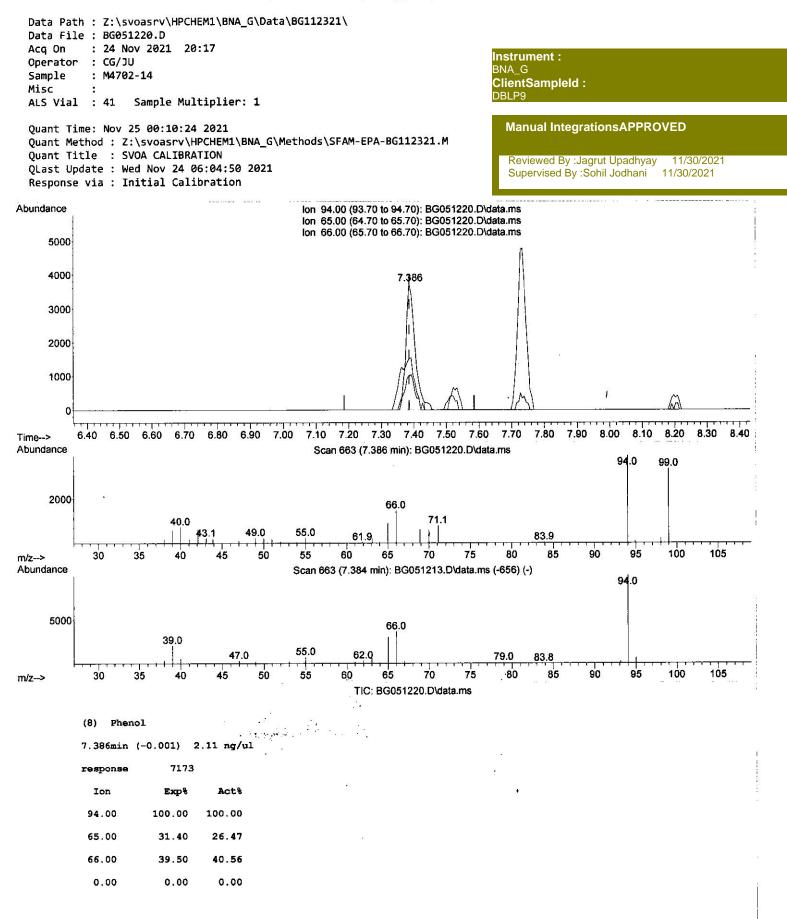


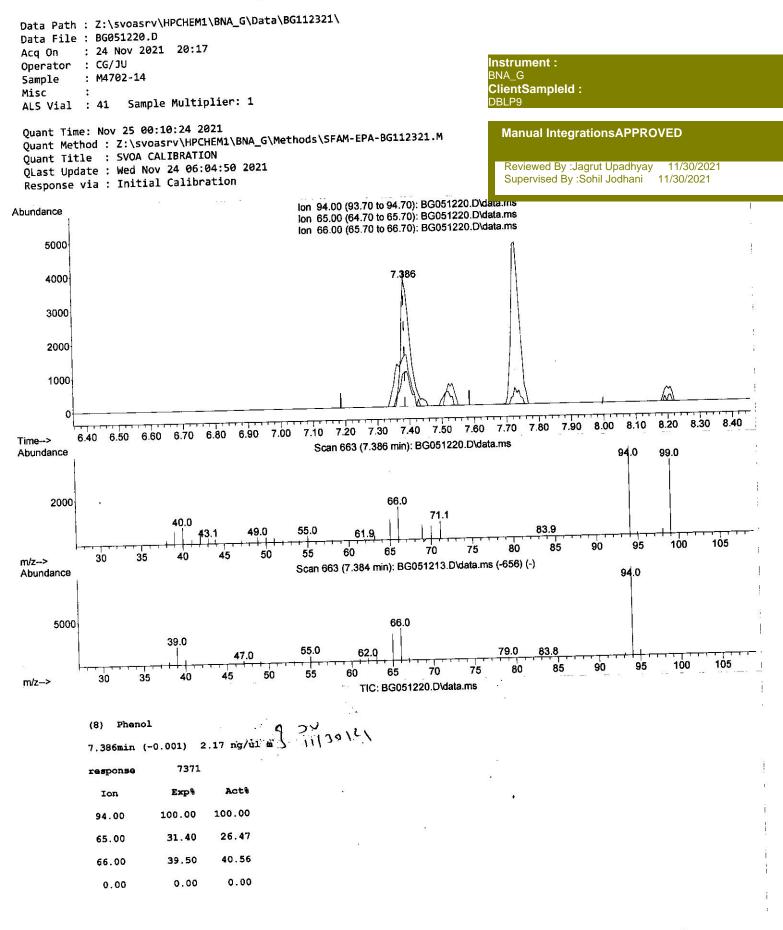
4

Quantitation Report (Qedit)

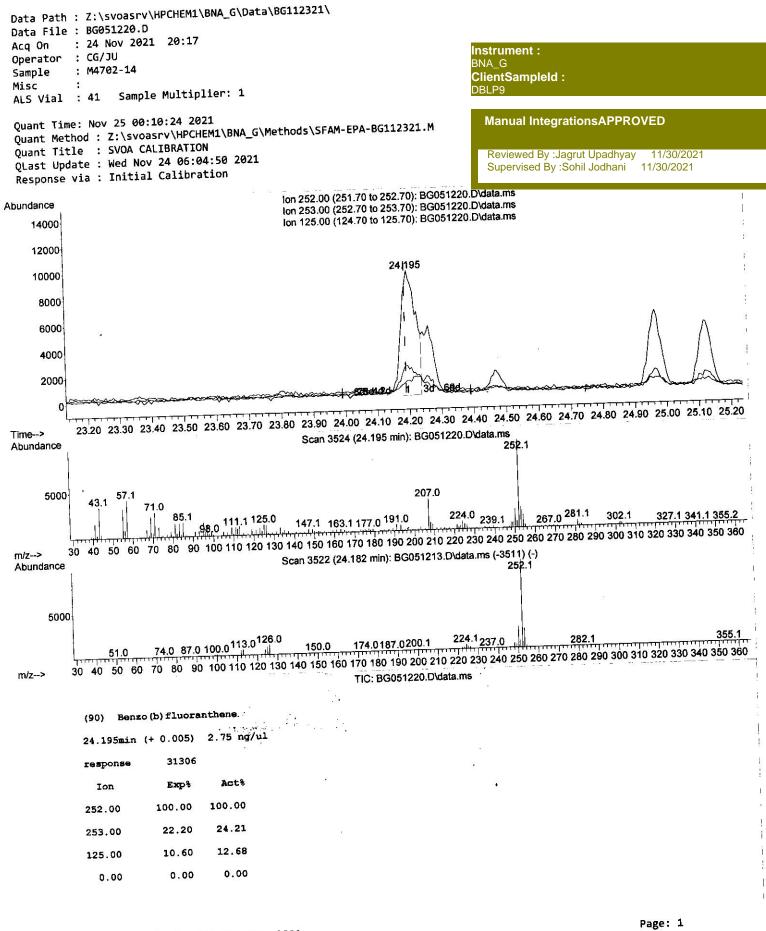


SFAM-EPA-BG112321.M Thu Nov 25 03:51:44 2021

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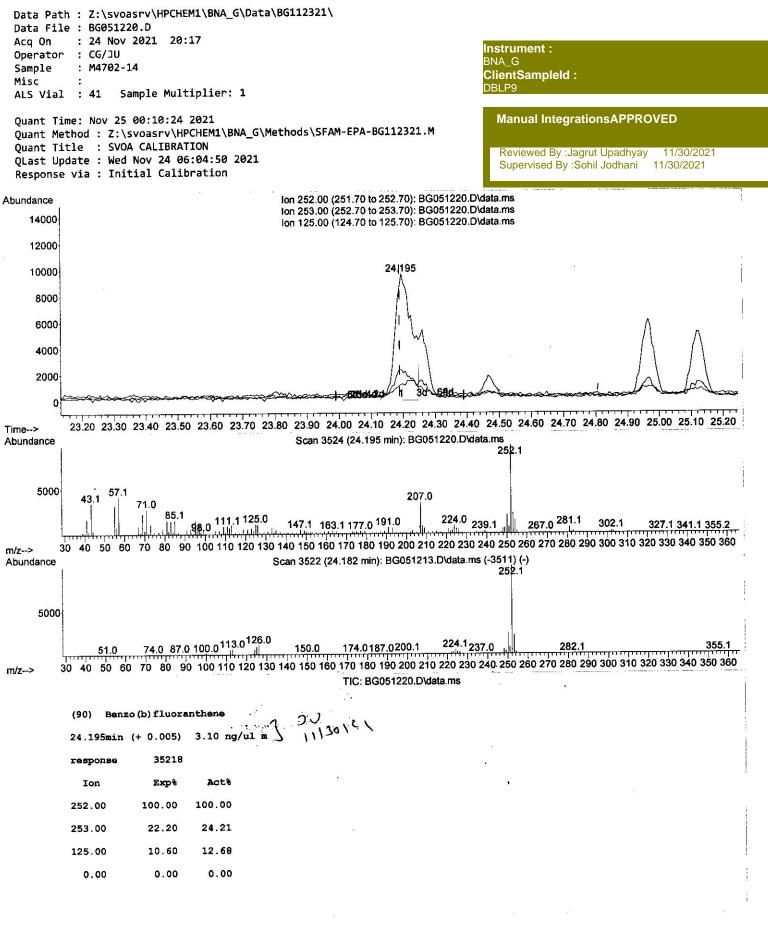


SFAM-EPA-BG112321.M Thu Nov 25 03:51:58 2021

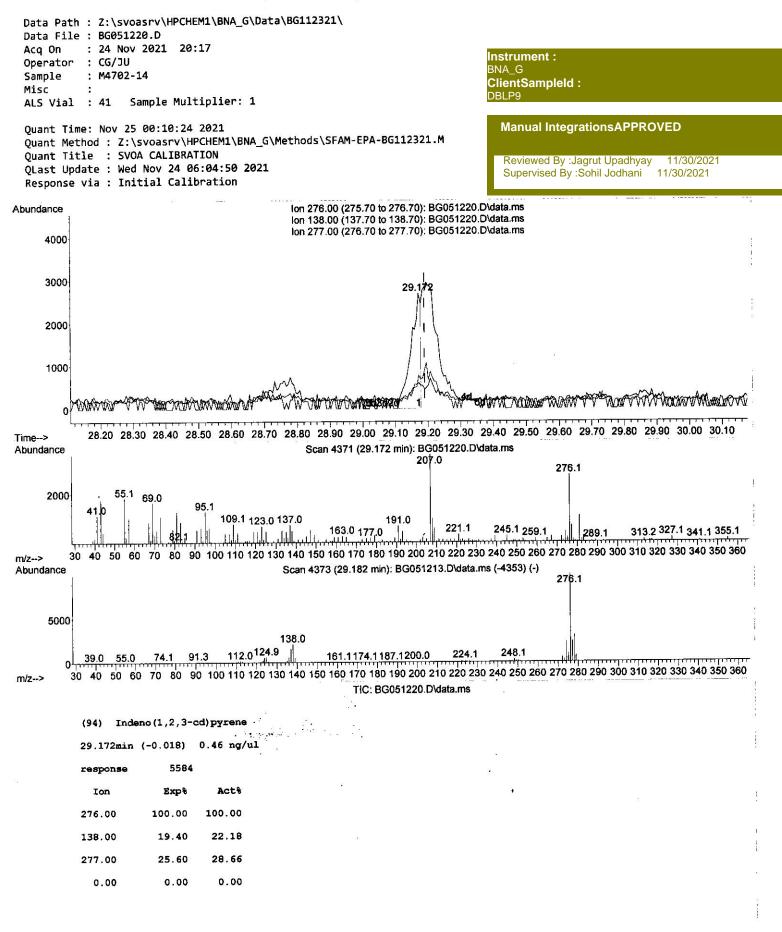


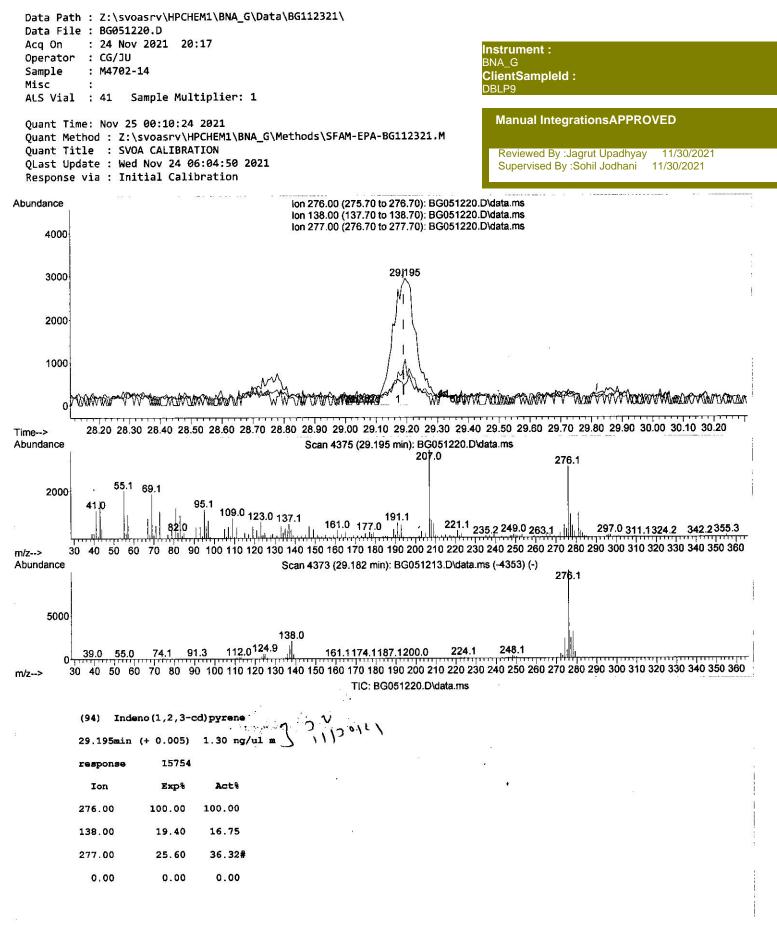
SFAM-EPA-BG112321.M Thu Nov 25 03:52:23 2021

Quantitation Report (Qedit)



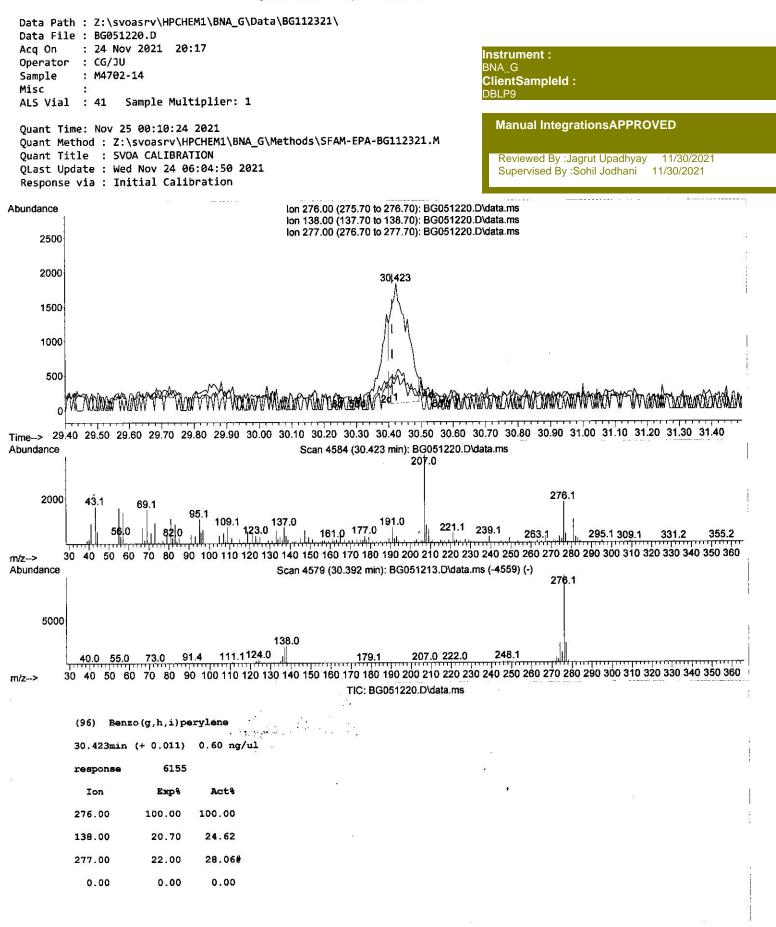
SFAM-EPA-BG112321.M Thu Nov 25 03:52:51 2021



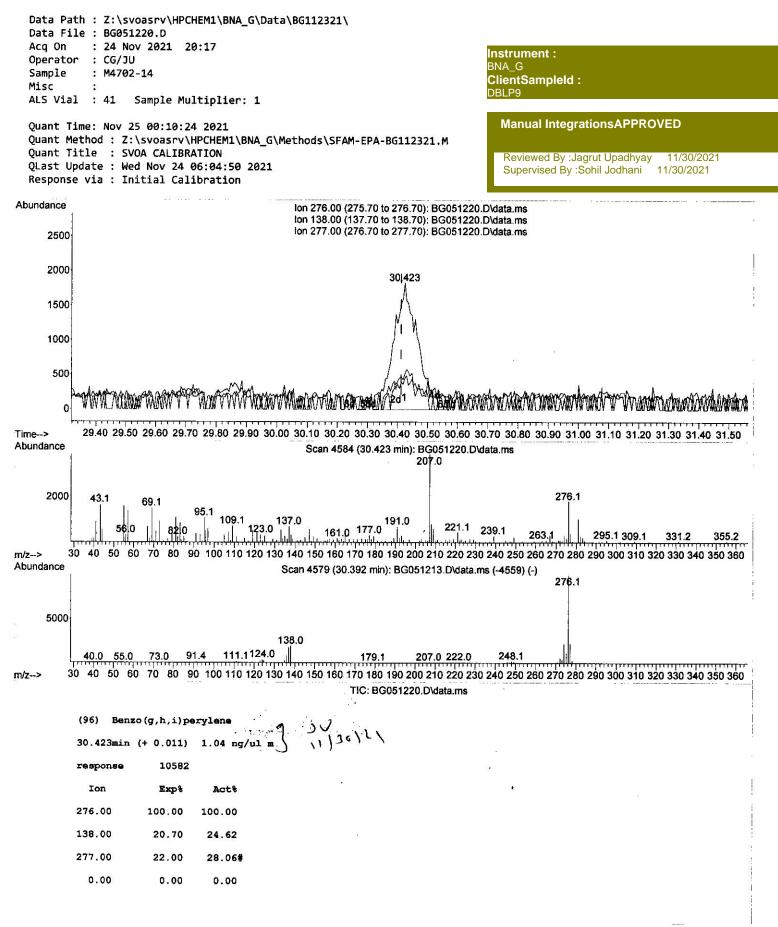


SFAM-EPA-BG112321.M Thu Nov 25 03:53:31 2021

Quantitation Report (Qedit)



SFAM-EPA-BG112321.M Thu Nov 25 03:53:55 2021



SFAM-EPA-BG112321.M Thu Nov 25 03:54:08 2021

| | ath : Z:\svoasrv\HPCHEM1\B ile : BG051220.D | NA_G\Data | a\BG1: | 12321\ | | | | |
|--------|--|------------------|------------|-----------------|------------------|--|--------------|---|
| Acq On | | | | | | | | |
| | or : CG/JU | | | | | | | trument : |
| Sample | | | | | | | | A_G |
| Misc | : | | | | | | | entSampleId : |
| | al : 41 Sample Multipli | er: 1 | | | | | DBI | LP9 |
| | Time: Nov 25 00:10:24 2021 | | | | | | Γ | Manual IntegrationsAPPROVED |
| | Method : Z:\svoasrv\HPCHEM Title : SVOA CALIBRATION | 1\BNA_G\I | Metho | ds\SFAM-EPA | -BG11232 | 1.M | | |
| | Update : Wed Nov 24 06:04: | 50 2021 | | | | | | Reviewed By :Jagrut Upadhyay 11/30/202 |
| | se via : Initial Calibrati | | | | | | Ľ | Supervised By :Sohil Jodhani 11/30/2021 |
| | Compound | | | Response | | | (Min) | |
| Inter | nal Standards | | | | | | | |
| | | 8.202 | 152 | 33129 | 20.000 | ng/ul | 0.00 | |
| 1000 | Naphthalene-d8 | 11.028 | 136 | 146727 | 20.000 | | 0.00 | |
| | Acenaphthene-d10 | 14.830 | 164 | 101025 | 20.000 | | 0.00 | |
| | Phenanthrene-d10 | 17.580 | 188 | 202938 | 20.000 | ng/ul | 0.00 | |
| | Chrysene-d12 | 21,880 | 240 | 168504 | 20.000 | ng/ul | 0.00 | |
| 88) | Perylene-d12 | 25.282 | 264 | 168414 | 20.000 | ng/ul | 0.00 | |
| Svste | m Monitoring Compounds | | | | | | | |
| | 1,4-Dioxane-d8 | 3.543 | 96 | 1657 | 1.738 | ng/uL | 0.00 | |
| | Pyridine-d5 | 3.978 | 84 | 7601 | 2.717 | | 0.00 | |
| 7) | Phenol-d5 | 7.362 | 99 | 35107 | 10.722 | ng/ul | 0.00 | |
| 9) | Bis-(2-Chloroethyl)eth | 7.515 | 67 | 22763 | 11.069 | ng/ul | 0.00 . | . 1 |
| 11) | 2-Chlorophenol-d4 | 7.732 | 132 | 26146 | 11,089 | | 0.00 | |
| 15) | 4-Methylphenol-d8 | 8.919 | 113 | 18968 | 7.179 | 1000 million 100 million 1 | 0.00 | |
| | Nitrobenzene-d5 | 9.377 | 128 | 14198 | 11.463 | - | 0.00 | |
| | 2-Nitrophenol-d4 | 10.106 | 143 | 16001 | 11.452 | | 0.00 | |
| | 2,4-Dichlorophenol-d3 | 10.652 | | 27653 | 11.665 | - | 0.00 | |
| | 4-Chloroaniline-d4 | 11.175 | | 13799 | 3.978 | | 0.01 | |
| | Dimethylphthalate-d6 | 14.225 | | 99429 | 12.791 | | 0.00 0.00 | |
| | Acenaphthylene-d8 | 14.530 | 160 | 125641 13632 | 12.818 10.834 | | 0.00 | |
| | 4-Nitrophenol-d4 | 15.053 15.823 | 143 176 | 90706 | 12.958 | 1000 | 0.00 | |
| | Fluorene-d10 4,6-Dinitro-2-methylph | | 200 | 9222 | 7.364 | | 0.00 | |
| | Anthracene-d10 | 17.680 | 188 | 132780 | 13.680 | 1928 | 0.00 | |
| | Pyrene-d10 | 19.959 | 212 | 160230 | 15.715 | | 0.00 | |
| | Benzo(a)pyrene-d12 | 25.047 | | 132575 | 14.740 | | 0.00 | |
| Tanga | t Compounds | | | | | 0v | alue | |
| | t Compounds Benzaldehyde | 7.333 | 77 | 2668 🦯 | 1.280 | | 95 | |
| • | Phenol | 7.386 | 94 | 7371m | | ng/ul | | \sim \sim |
| | Acetophenone | 9.031 | 105 | 21443 | | ng/ul | 92 | 50 |
| | Fluoranthene | 19.624 | 202 | 85995 | | ng/ul | 96 | 11/30/21 |
| | Pyrene | 19,989 | 202 | 89964 | | ng/ul | 96 | (I) |
| | Benzo(a)anthracene | 21.863 | 228 | 30416 | | ng/ul | 99 | , |
| | Chrysene | 21.927 | 228 | 33136 | | ng/ul | 99 | |
| | Benzo(b)fluoranthene | 24.195 | 252 | 35218m | | ng/ul | | |
| | Benzo(a)pyrene | 25.118 | | 14155 | 1.305 | ng/ul | 94 | |
| | Indeno(1,2,3-cd)pyrene | 29.195 | | 15754m | | ng/ul | | |
| | Benzo(g,h,i)perylene | 30.423 | 276 | 10582m | 4 037 | ng/ul | | |

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

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