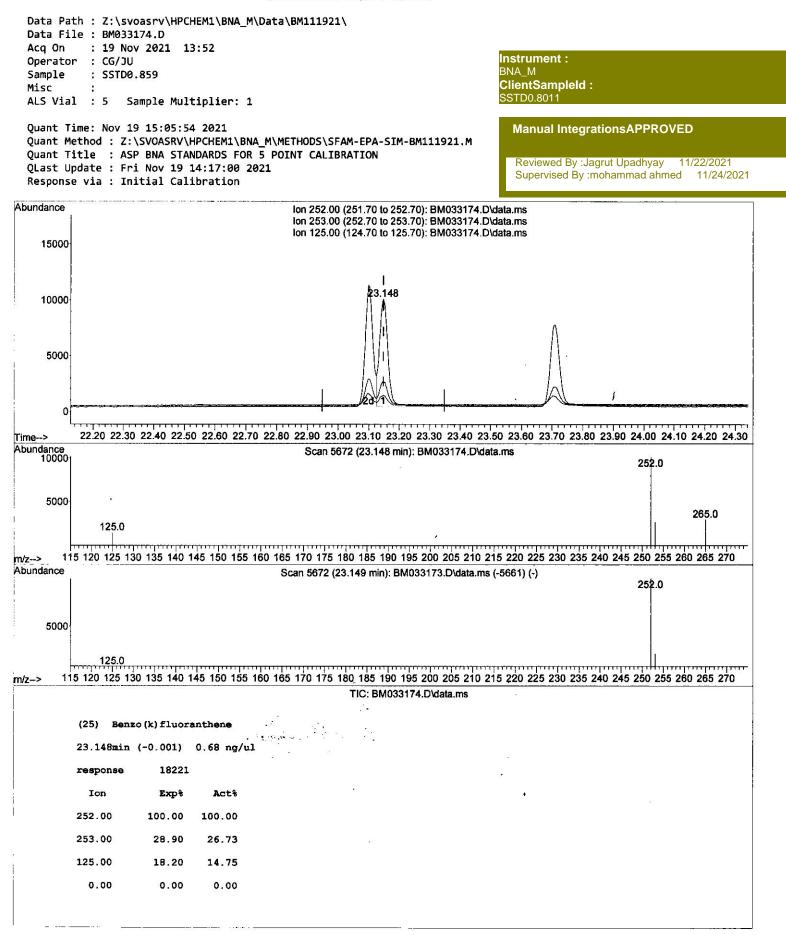
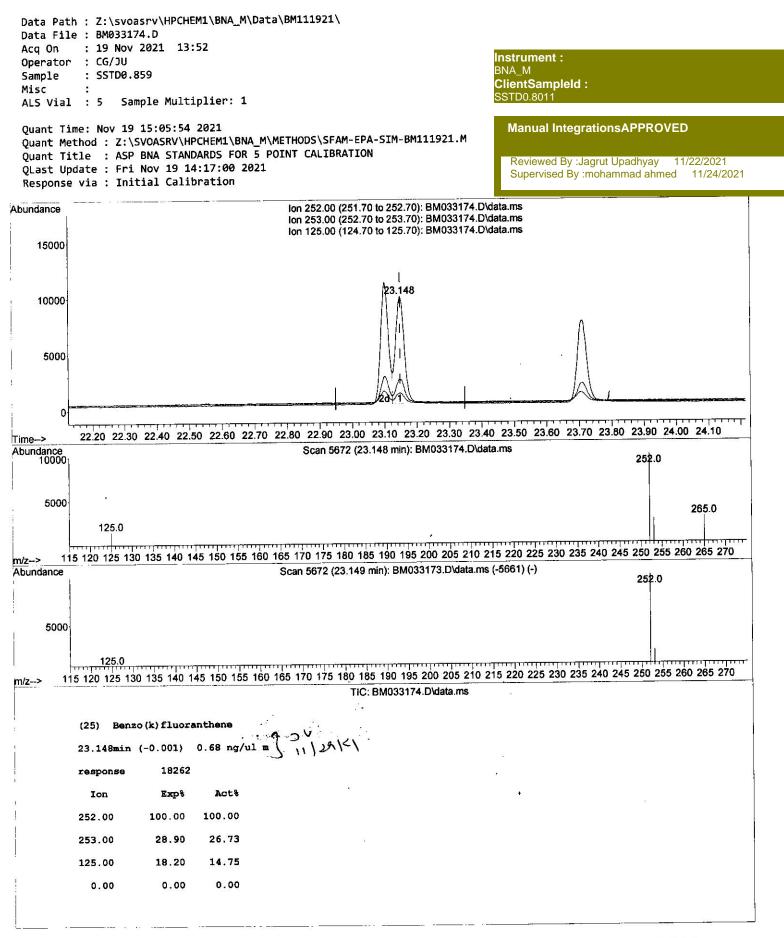


SFAM-EPA-SIM-BM111921.M Fri Nov 19 15:13:01 2021



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SFAM-EPA-SIM-BM111921.M Fri Nov 19 15:12:26 2021

Page: 1

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111921\ Data File : BM033174.D Acq On : 19 Nov 2021 13:52 Operator : CG/JU BNA_M : SSTD0.859 Sample Misc ALS Vial : 5 Sample Multiplier: 1 Quant Time: Nov 19 15:05:54 2021 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION QLast Update : Fri Nov 19 14:17:00 2021 Response via : Initial Calibration R.T. QIon Response Conc Units Dev(Min) Compound -----Internal Standards 0.00 0.400 ng/ul 7.963 152 2404 1) 1,4-Dichlorobenzene-d4 0.400 ng/ul # 0.00 6505 10.760 136 4) Naphthalene-d8 0.400 ng/ul 0.00 3744 14.583 164 9) Acenaphthene-d10 0.00 17.316 188 7980 0.400 ng/ul 13) Phenanthrene-d10 0.00 0.400 ng/ul 6670 21.472 240 17) Chrysene-d12 0.400 ng/ul # 0.00 5487 23.812 264 23) Perylene-d12 System Monitoring Compounds 0.00 0.584 ng/ul 1793 3.420 96 3) 1,4-Dioxane-d8 0.00 0.761 ng/ul 6) 2-Methylnaphthalene-d10 12.339 152 7028 0.00 0.744 ng/ul 19.334 212 15035 18) Fluoranthene-d10 Qvalue Target Compounds 91 0.678 ng/ul 88 2137 3.455 2) 1,4-Dioxane 98 0.829 ng/ul 10.810 128 15664 5) Naphthalene 98 0.782 ng/ul 10238 7) 2-Methylnaphthalene 12.415 142 0.791 ng/ul 100 10140 12.631 142 8) 1-Methylnaphthalene 0.846 ng/ul# 99 13771 14.303 152 10) Acenaphthylene 0.791 ng/ul 99 14.644 153 11237 11) Acenaphthene 98 0.762 ng/ul 15.626 166 12771 12) Fluorene 96 16.965 266 1.009 ng/ul 2097 14) Pentachlorophenol 0.761 ng/ul 99 17.357 178 20188 15) Phenanthrene 0.775 ng/ul 99 18051 17.451 178 16) Anthracene 97 0.719 ng/ul 19.360 202 22490 19) Fluoranthene 96 0.691 ng/ul 22479 19.722 202 20) Pyrene 99 21.455 228 0.717 ng/ul 17651 21) Benzo(a)anthracene 99 0.684 ng/ul 18994 21.508 228 22) Chrysene 18529 18262m } J1/29121 0.736 ng/ul 94 23.099 252 24) Benzo(b)fluoranthene 0.677 ng/ul 23.148 252 25) Benzo(k)fluoranthene 91 0.760 ng/ul 15780 23.708 252 26) Benzo(a)pyrene 88 26.208 276 0.839 ng/ul# 20037 27) Indeno(1,2,3-cd)pyrene 0.850 ng/ul# 92 16068 26.223 278 28) Dibenzo(a,h)anthracene 92 . 17637 0.851 ng/ul# 26.942 276 29) Benzo(g,h,i)perylene -------------

(#) = qualifier out of range (m) = manual integration (+) = signals summed a de service de la destrucción de la de

Instrument :

ClientSampleId :

SSTD0.8011

Manual IntegrationsAPPROVED

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Reviewed By : Jagrut Upadhyay 11/22/2021 Supervised By :mohammad ahmed 11/24/2021

SFAM-EPA-SIM-BM111921.M Fri Nov 19 15:12:59 2021