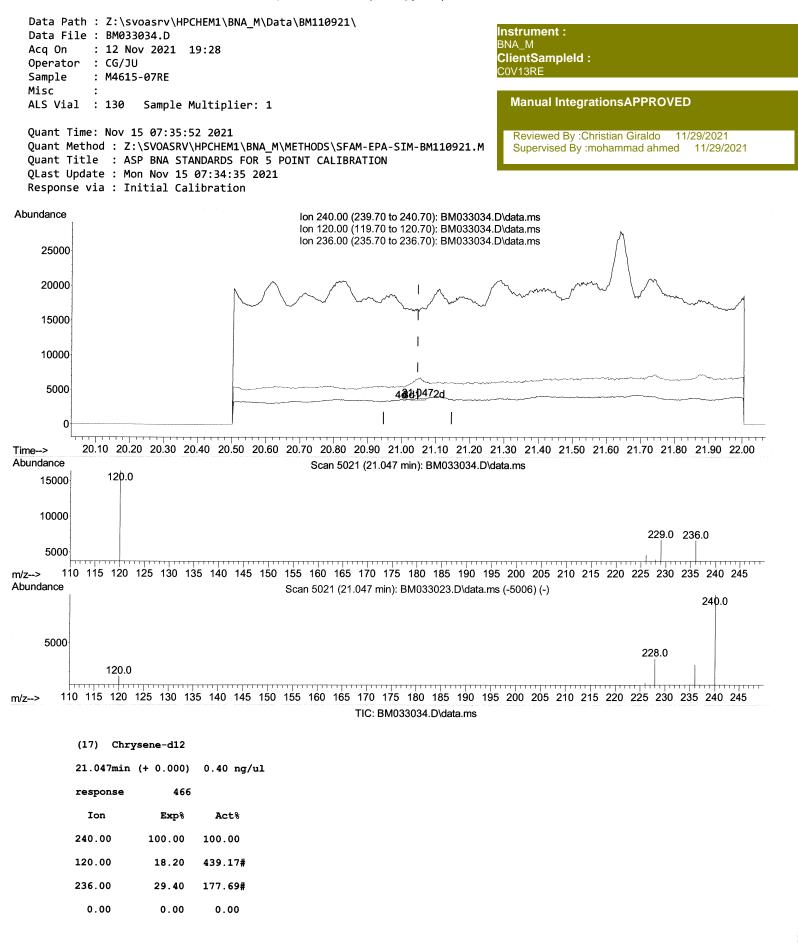
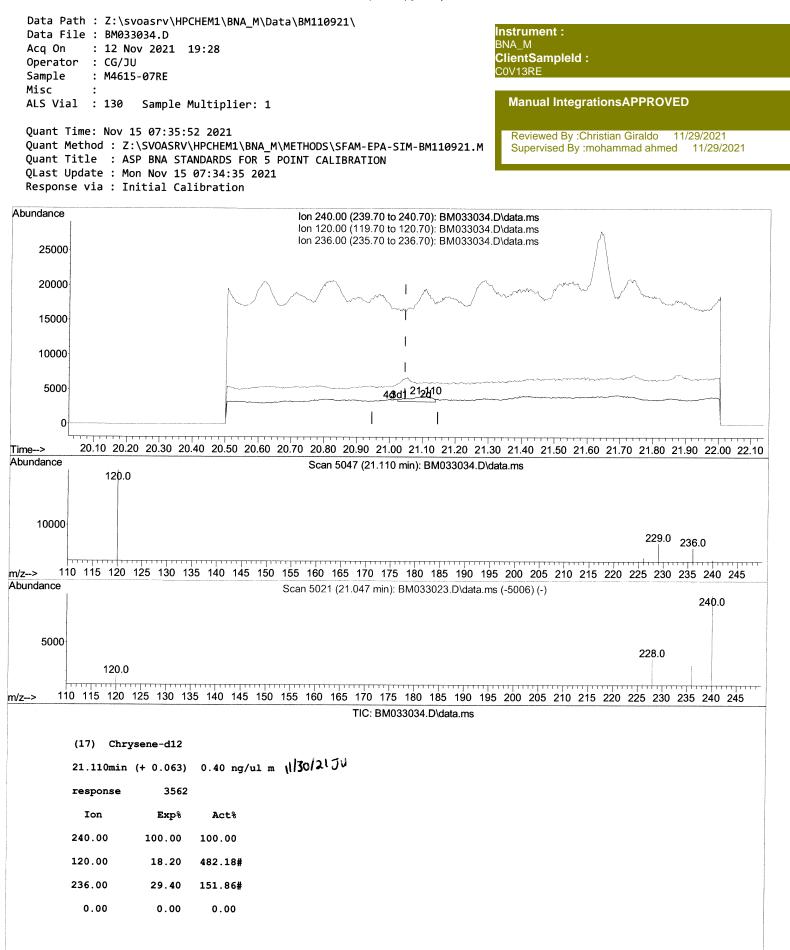
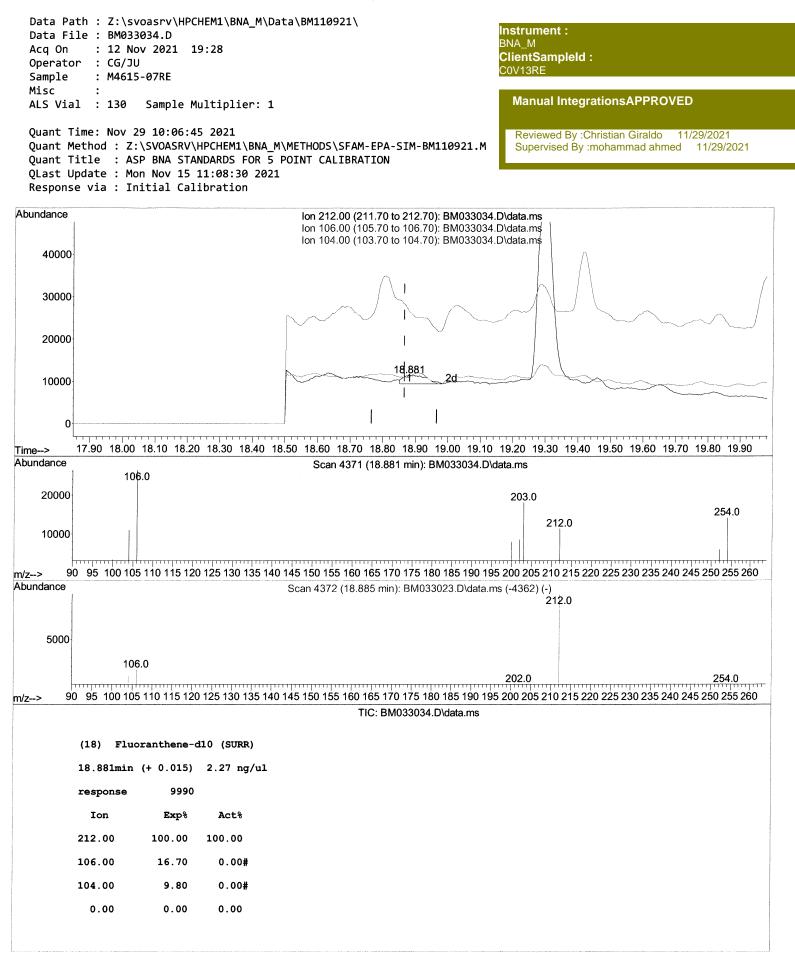


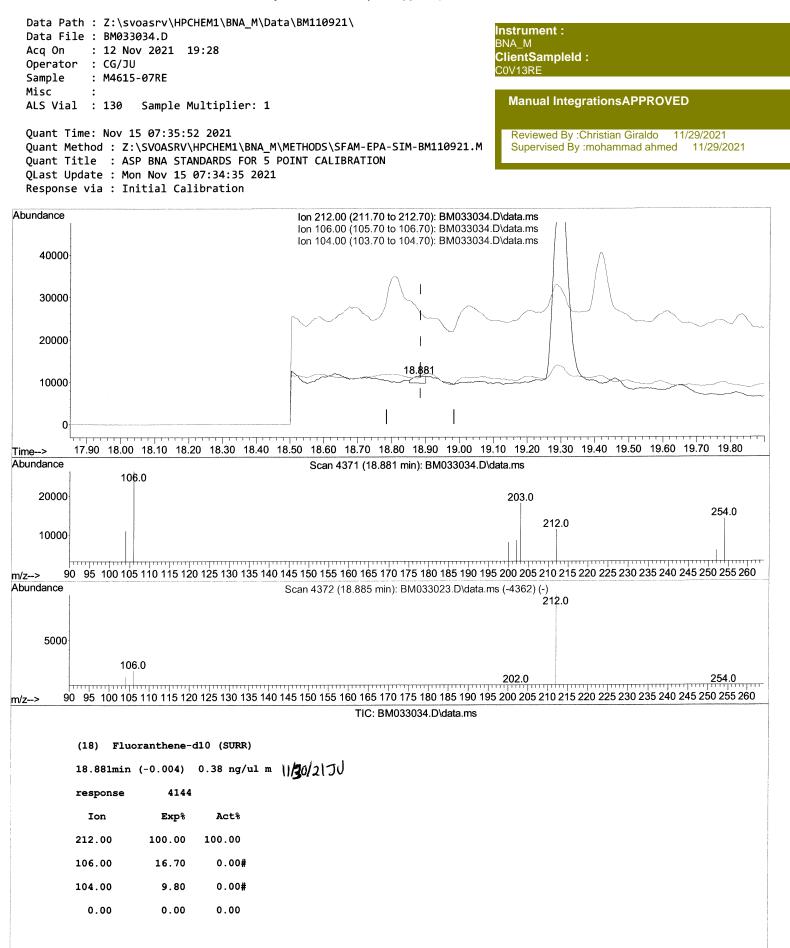
Time--> 4.00 6.00 8.00 10.00 12.00 14.00 16.00 18.00 20.00 22.00 24.00 26.00 28.00











Data Path : Z:\svoasrv\HPCHEM1\ Data File : BM033034.D Acq On : 12 Nov 2021 19:28 Operator : CG/JU Sample : M4615-07RE Misc : ALS Vial : 130 Sample Multip	_	ta\BM:	110921\	Instrument : BNA_M ClientSampleId : C0V13RE Manual IntegrationsAPPROVED
Quant Time: Nov 15 07:35:52 2021 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110921.M Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION QLast Update : Mon Nov 15 07:34:35 2021 Response via : Initial Calibration				
Compound		-		Conc Units Dev(Min)
Internal Standards				
 1,4-Dichlorobenzene-d4 	7.452	152	2013	0.40 ng/ul 0.00
A) Naphthalene-d8	10.230	136	6804	0.40 ng/ul # 0.00
9) Acenaphthene-d10	14.121	164	6187	0.40 ng/ul # 0.01
13) Phenanthrene-d10	16.872	188	6568	0.40 ng/ul # 0.02
17) Chrysene-d12	21.110	240	3562m>	0.40 ng/ul > 0.06 /30/2 JU
23) Perylene-d12	23.156		53900	0.40 ng/ul # 0.00
System Monitoring Compounds				
3) 1,4-Dioxane-d8	2.860	96	2795	1.09 ng/ul 0.00
6) 2-Methylnaphthalene-d10	11.827	152	1916	0.20 ng/ul 0.00
18) Fluoranthene-d10	18.881	212	4144m>	
Target Compounds				Qvalue
5) Naphthalene	10.280	128	1257	0.06 ng/ul# 77
7) 2-Methylnaphthalene	11.903	142	3055	0.22 ng/ul 95
8) 1-Methylnaphthalene	12.128	142	2145	0.16 ng/ul 90
12) Fluorene	15.156	166	9937	0.36 ng/ul# 16
15) Phenanthrene	16.900		13229	0.61 ng/ul# 1

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