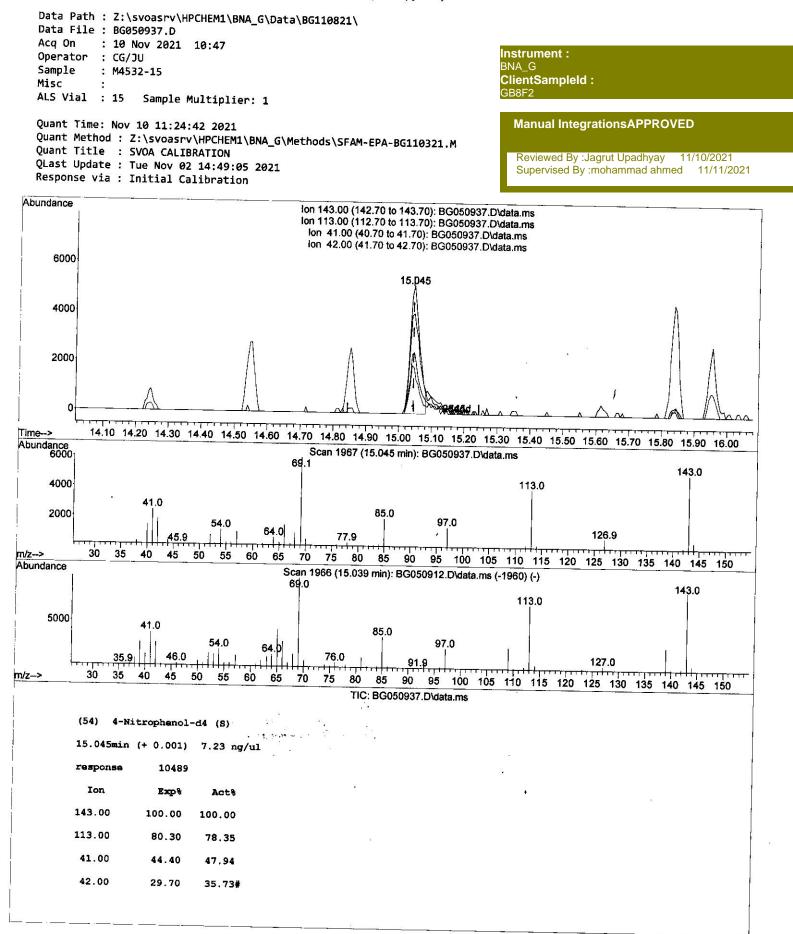


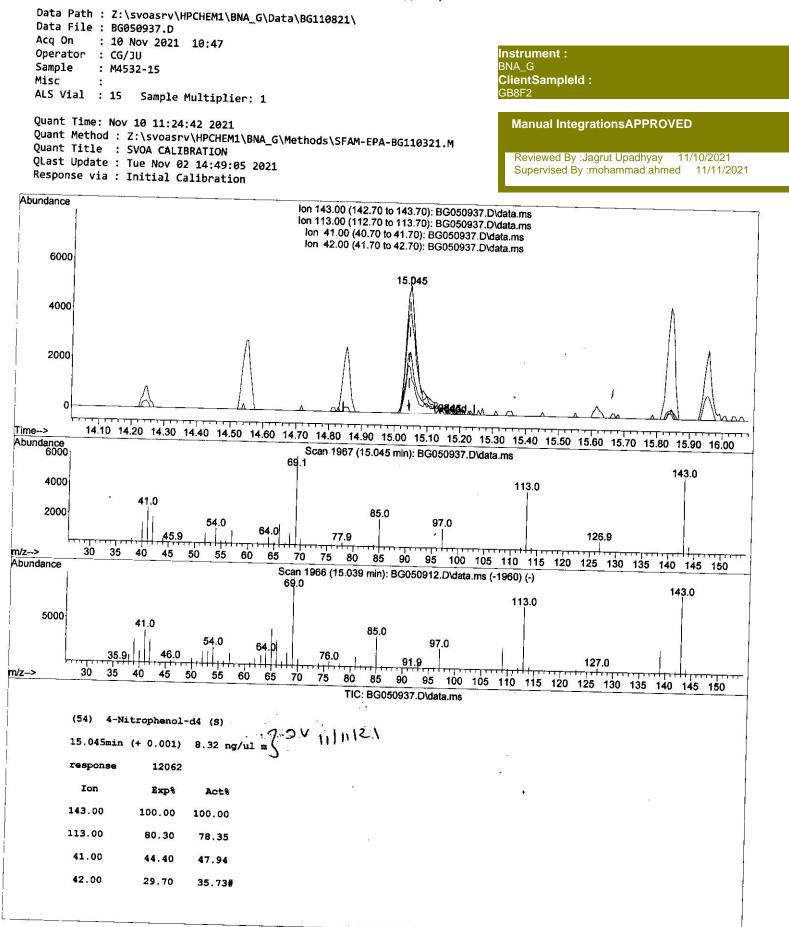
SFAM-EPA-BG110321.M Wed Nov 10 18:04:24 2021

Page: 2

٩



3



SFAM-EPA-BG110321.M Wed Nov 10 11:26:12 2021

Page: 1

Data Path : Z:\svoasrv\HPCHEM1 Data File : BG050937.D Acq On : 10 Nov 2021 10:47 Operator : CG/JU Sample : M4532-15 Misc : ALS Vial : 15 Sample Multip		ata\8G	110821\		Instrument : BNA_G ClientSampleId : GB8F2	
Quant Time: Nov 10 11:24:42 2021 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM~EPA-BG110321.M Quant Title : SVOA CALIBRATION					Manual IntegrationsAPPROVED Reviewed By :Jagrut Upadhyay 11/10/2021	
QLast Update : Tue Nov 02 14:49 Response via : Initial Calibrat	9:05 2021 tion				Supervised By :mohammad ahmed 11/11/2021	
Compound	R.T.	QIon	Response	Conc Units De	ev(Min)	
Internal Standards						
 1,4-Dichlorobenzene-d4 	8.229	152	35806	20.000 (1		
20) Naphthalene~d8	11.056		153764	20.000 ng/ul	0.00	
38) Acenaphthene-d10	14.851		104560	20.000 ng/ul 20.000 ng/ul	0.00	
64) Phenanthrene-d10	17.595		230838	20.000 ng/ul	-0.01	
79) Chrysene-d12	21.896		204736	20.000 ng/ul	-0.01	
88) Perylene-d12	25.292		199123	20.000 ng/ul	-0.01 -0.02	
Curta - N. Li			90000120000000-0 - 2		-0.02	
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	6524	5.881 ng/uL	0.00	
4) Pyridine-d5	4.017	84	18028	5.432 ng/ul	0.00	
7) Phenol-d5	7.372	99	30527	7.992 ng/ul	0.00	
9) Bis-(2-Chloroethyl)eth	7.542	67	75581	30.630 ng/ul	0.00	
11) 2-Chlorophenol-d4	7.754	132	65809	24.859 ng/ul	-0.01	
15) 4-Methylphenol-d8	8,929	113	51290	17.056 ng/ul	0.00	
21) Nitrobenzene-d5	9.399	128	43353	33.178 ng/ul		
24) 2-Nitrophenol-d4	10.127	143	45267	31.156 ng/ul	-0.01	
28) 2,4-Dichlorophenol-d3	10.668	165	70325	28.734 ng/ul	0.00	
31) 4-Chloroaniline-d4	11.185	131	92040	24.833 ng/ul	0.00	
46) Dimethylphthalate-d6	14.246	166	289146	36.146 ng/ul		
49) Acenaphthylene-d8	14.546	160	334608	33.573 ng/ul	-0.01 V (11/11 VC	
54) 4-Nitrophenol-d4	15.045	143	12062m \		0.00	
60) Fluorene-d10	15.838	176	250934	35.410 ng/ul		
65) 4,6-Dinitro-2-methylph	15.956	200	46489	33.218 ng/ul	0.00	
73) Anthracene-d10	17.695	188	465082	42.614 ng/ul		
81) Pyrene-d10	19.975	212	535159	40.470 ng/ul	0.00	
92) Benzo(a)pyrene-d12	25.057	264	450105	40.889 ng/ul	-0.02	
Target Compounds						

.

12

2

a ay yana a sa a sa

Target Compounds

Qvalue

÷

.

•

1

1

Aratar

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

FAM-EPA-BG110321.M Wed Nov 10 18:04:23 2021