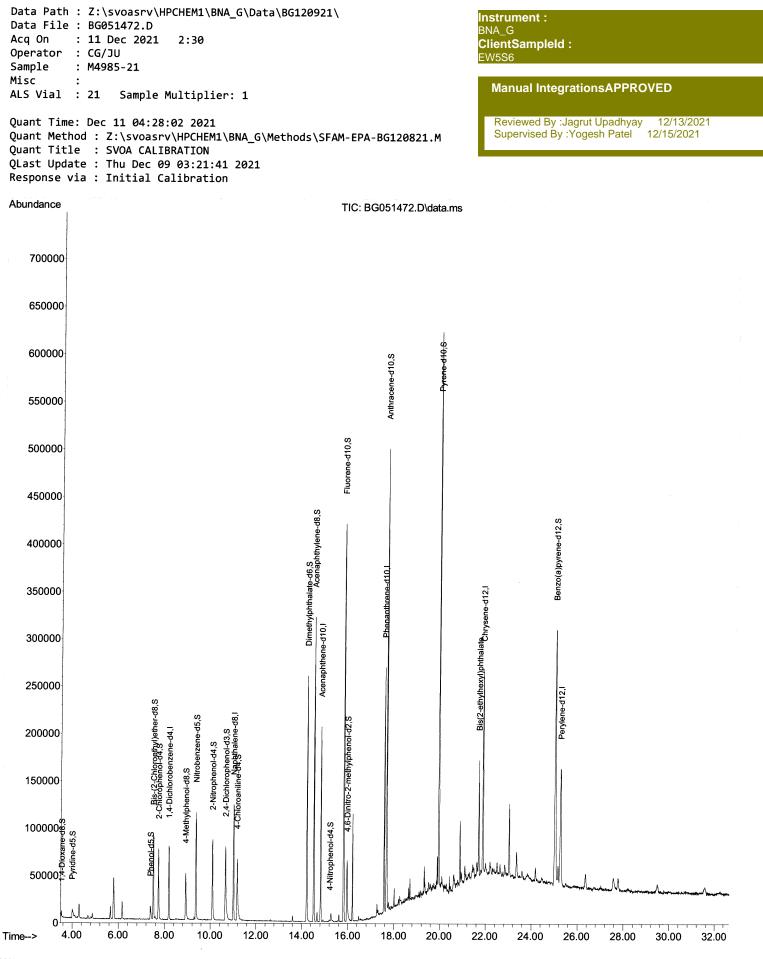
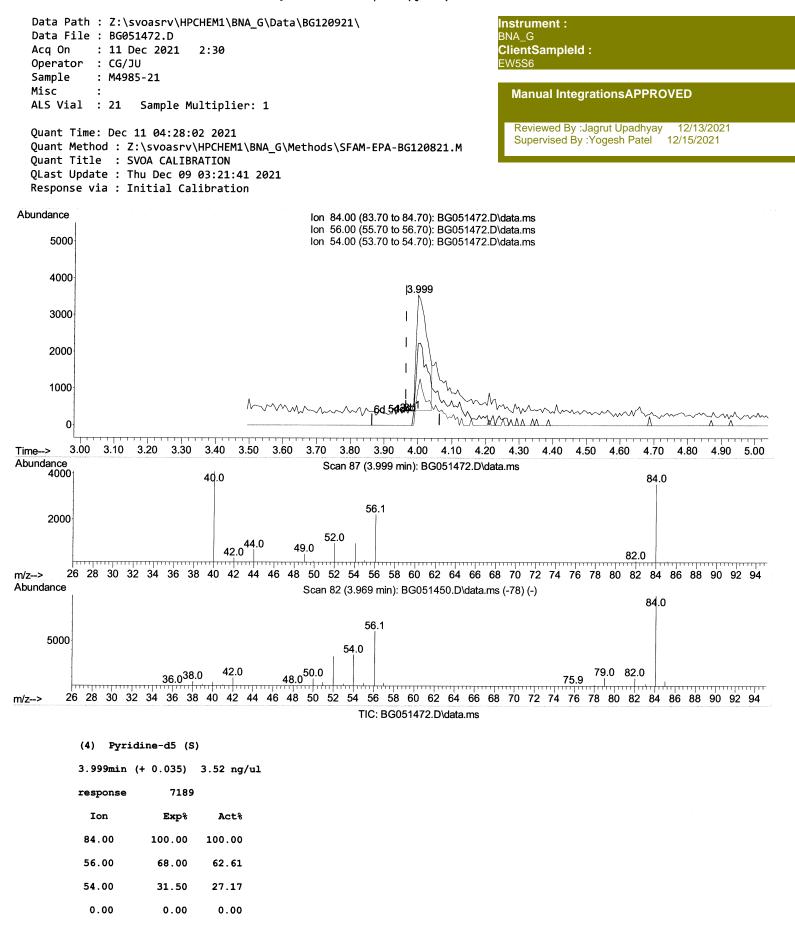
(QT Reviewed)

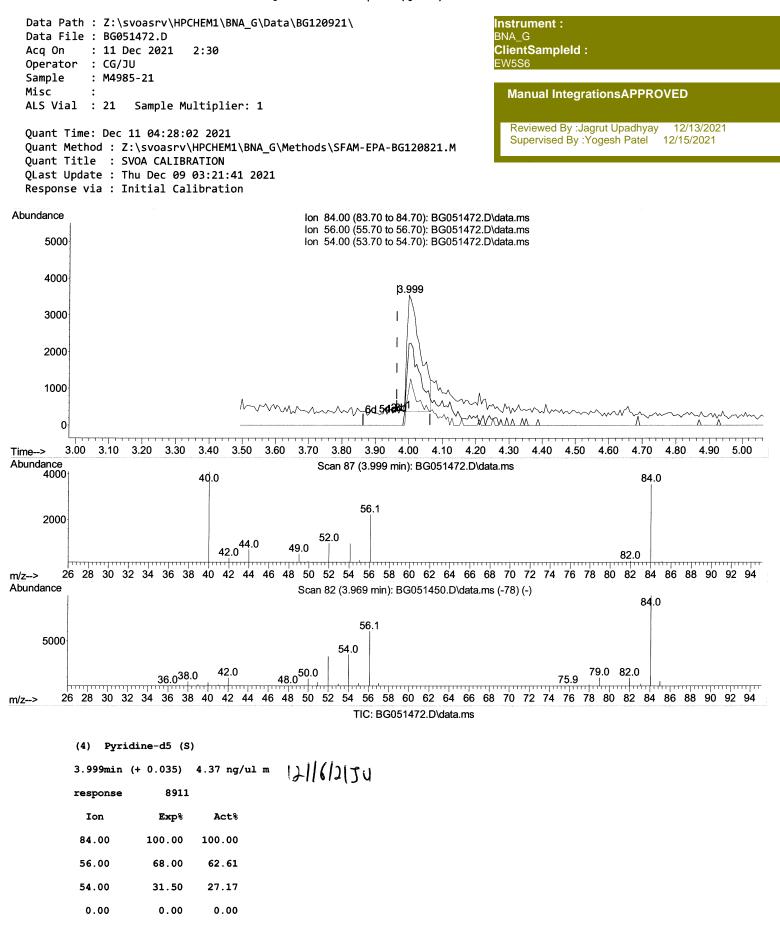


SFAM-EPA-BG120821.M Sat Dec 11 05:13:36 2021

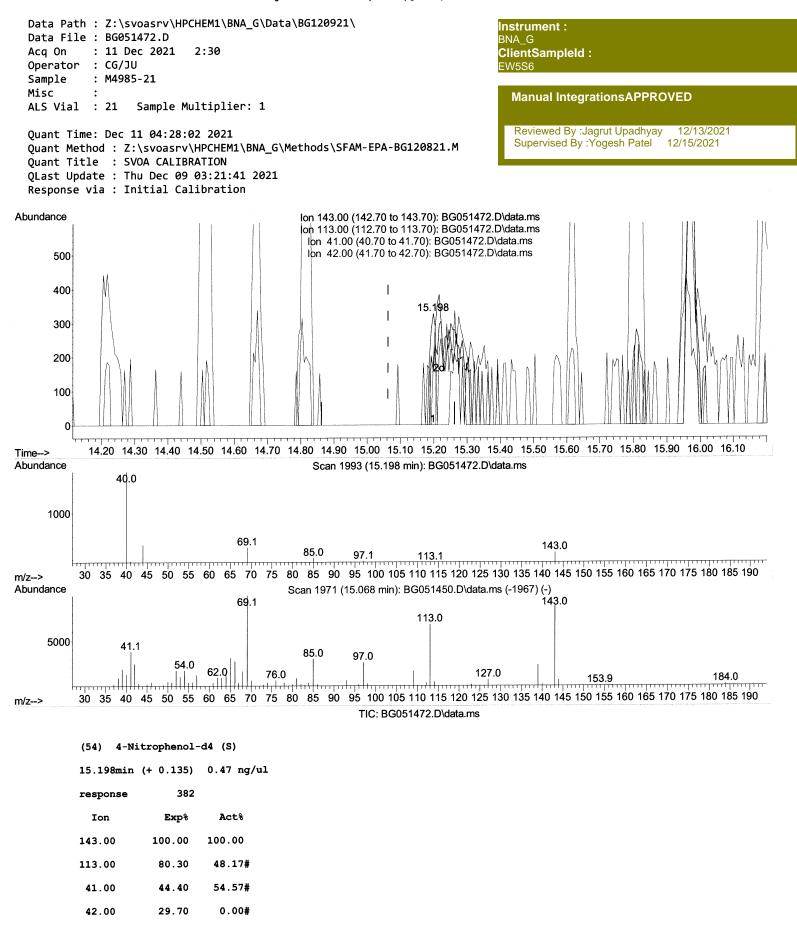
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Page: 2
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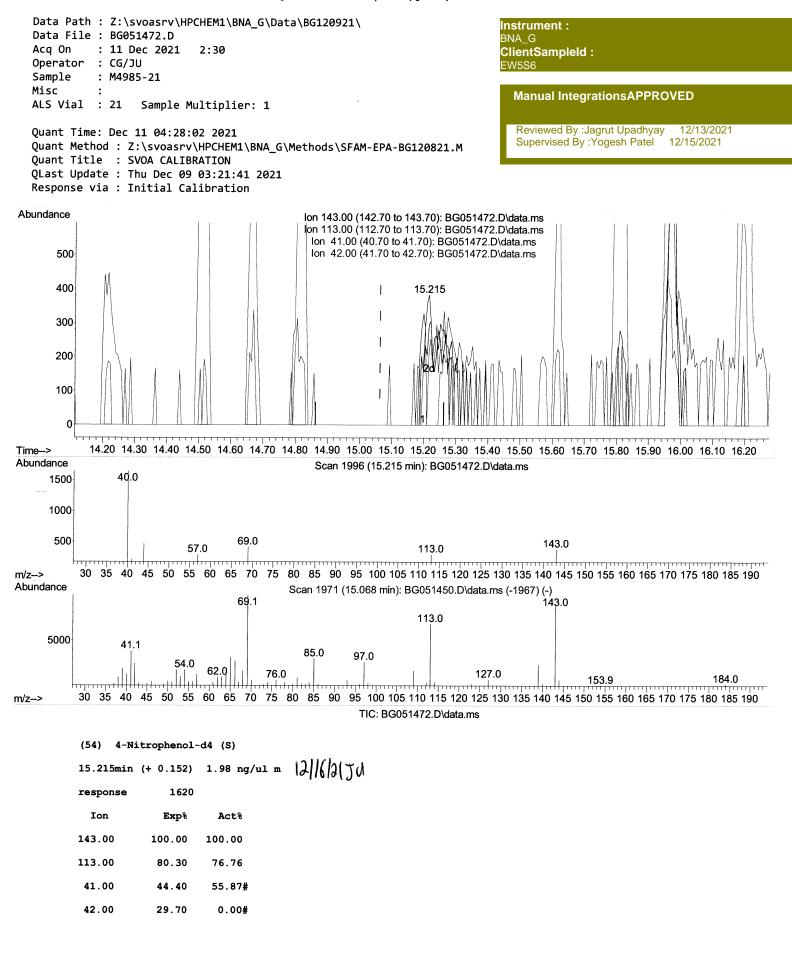












(QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120921\ Data File : BG051472.D Acq On : 11 Dec 2021 2:30 Operator : CG/JU Sample : M4985-21 Misc : ALS Vial : 21 Sample Multiplier: 1 Quant Time: Dec 11 04:28:02 2021 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M Quant Title : SVOA CALIBRATION QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration						BN CI EV	strument : IA_G ientSampleId : V5S6 Manual IntegrationsAPPROVED Reviewed By :Jagrut Upadhyay 12/13/2021 Supervised By :Yogesh Patel 12/15/2021
Compound	R.T. 0)Ion R	esponse	Conc Un	its Dev(Min)	
		-	•		•		
Internal Standards							
 1,4-Dichlorobenzene-d4 	8.188	152	23333	20.000	ng/ul	0.0	0
20) Naphthalene-d8	11.008	136	104288	20.000		0.0	0
38) Acenaphthene-d10	14.816	164	70327	20.000	ng/ul	0.0	0
64) Phenanthrene-d10	17.565	188 :	160200	20.000	ng/ul	0.0	0
79) Chrysene-d12	21.872	240 :	147617	20.000	ng/ul	0.0	0
88) Perylene-d12	25.268	264 :	141274	20.000	ng/ul	0.0	0
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.535	96	3461	4.871	ng/ul	0.00	
4) Pyridine-d5	3.999	84	8911m \	4.367		0.00	12/1(12134
7) Phenol-d5	7.389	99	12758	5.371	ng/ul	0.03	
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.501	67	49010	32.174		0.00	
11) 2-Chlorophenol-d4		132	40543	23.990		0.00	
15) 4-Methylphenol-d8		113	26853	14.390	.	0.01	
21) Nitrobenzene-d5		128	28781	31.814		0.00	
24) 2-Nitrophenol-d4		143	30467	29.762		0.00	
28) 2,4-Dichlorophenol-d3		165	44195	26.538		0.01	
31) 4-Chloroaniline-d4	11.173		53664	22.031		0.01	
46) Dimethylphthalate-d6			L88522	34.643	•	0.00	
49) Acenaphthylene-d8			240165	34 846	ng/ul	a aa	
54) 4-Nitrophenol-d4	15.215	143	1620m >	1.977	ng/ul>	0.15	12/661 JJ
60) Fluorene-d10	15.809	176 1	174520	36.026	ng/ul	0.00	
65) 4,6-Dinitro-2-methylph	15.961	200	16671	17.512		0.01	
73) Anthracene-d10	17.665	188 2	299716	39.985		0.00	
81) Pyrene-d10	19.945	212 3	849213	39.358		0.00	
92) Benzo(a)pyrene-d12	25.027		804082	41.731	0	0.00	
Target Compounds					0.75	luo.	
86) Bis(2-ethylhexyl)phtha	21 696 1	1/0	47135	7 250	Qva]		
		149 	4/133	7.258		97	

1

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