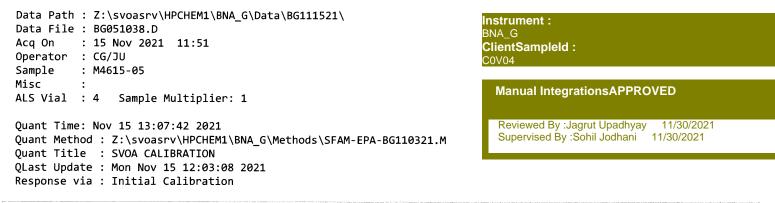
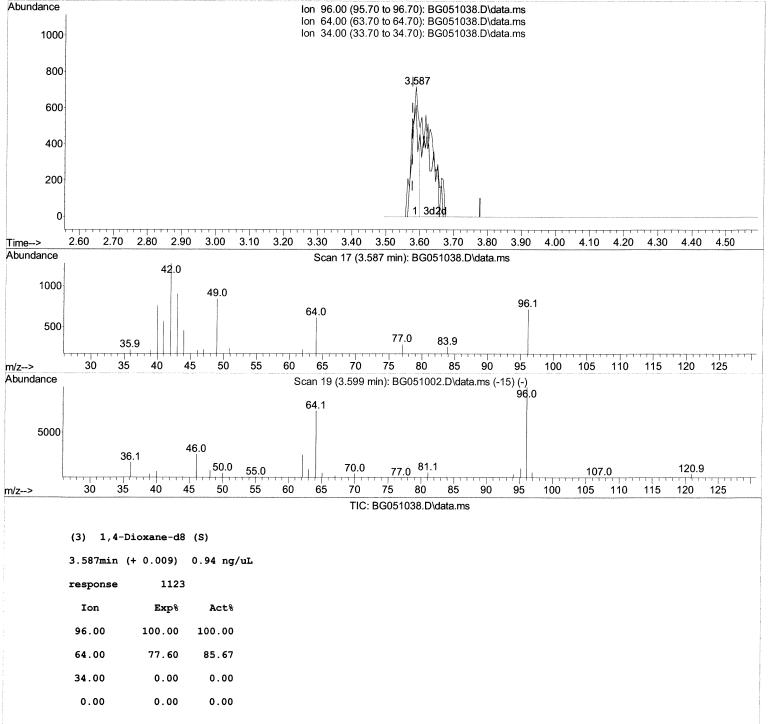
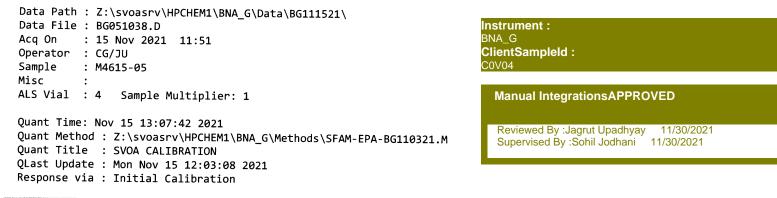
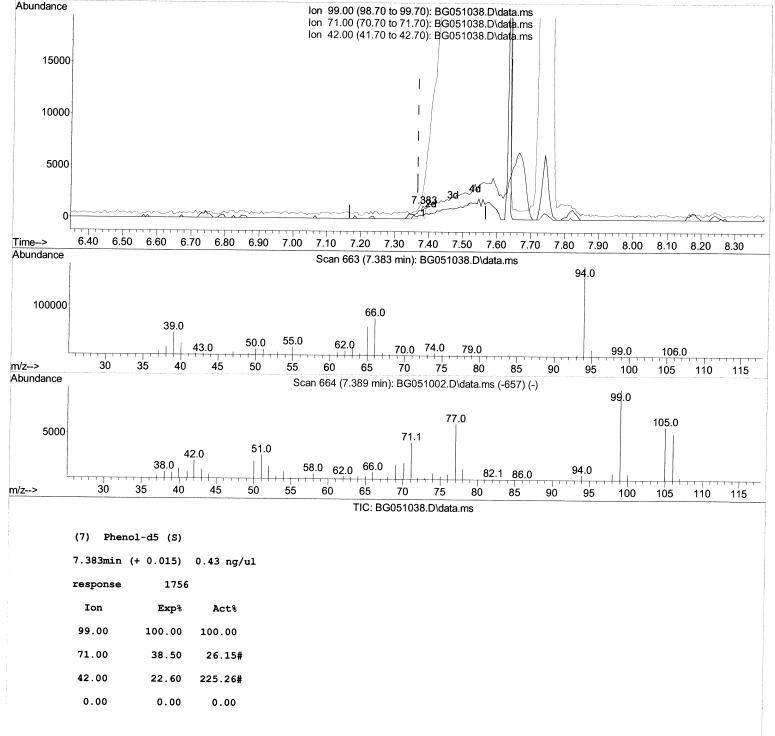
Data File Acq On Operator Sample Misc	: M4615-05 :		Instrument : BNA_G ClientSampleId : C0V04 Manual IntegrationsAPPROVED
Quant Time: Quant Title QLast Updat	: 4 Sample Multiplier: 1 : Nov 15 12:35:13 2021 e : te : Mon Nov 15 12:03:08 2021 ia : Initial Calibration		Reviewed By :Jagrut Upadhyay 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021
Abundance		ГІС: BG051038.D\data.ms	
1.7e+07			
1.6e+07			
1.5e+07			
1.4e+07	To a contract of the contract		
1.3e+07	<u>د</u>		
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2000000 සි ස	Pics Actifing and Lighter and S.S. 1.4-Dichlorobentzene-d4,1 Methylipanol, J.P. (1998) Mitrobenzene-d5, S 2.4-Dichlorophenol-d4, S 2.4-Dichlorophenol-d4, S 2.4-Dichlorophenol-d6, S Naphthalate-d6, S An Mitrophralate-d6, S An Mitrophralat	4.6-Dinitro-2 ⁻ / ///////////////////////////////////	Chrysene-d12,1 Beryddiabyfgyp-d12,S
10000000	Pitching 1,4-Dich 1,4-Dich 1,4-Dich 1,4-Dich 2,4-Dich Naphi Acenat	4,6-Dini Arthr	Benzel
0 Time> 4.0		6.00 18.00 20.00	22.00 24.00 26.00 28.00 30.00 32.00

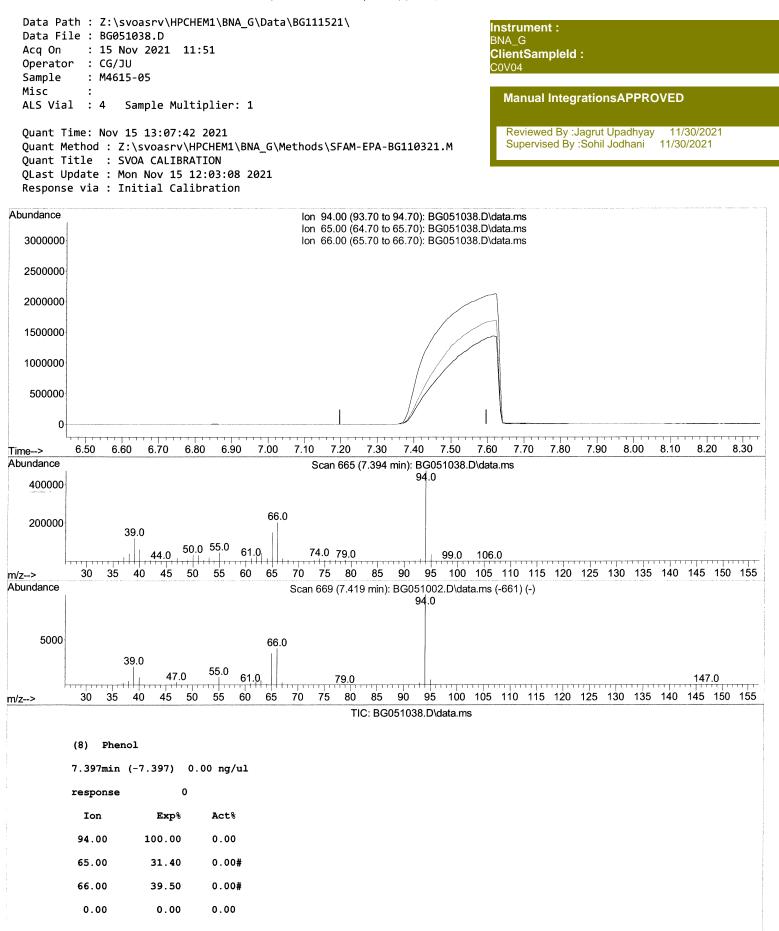




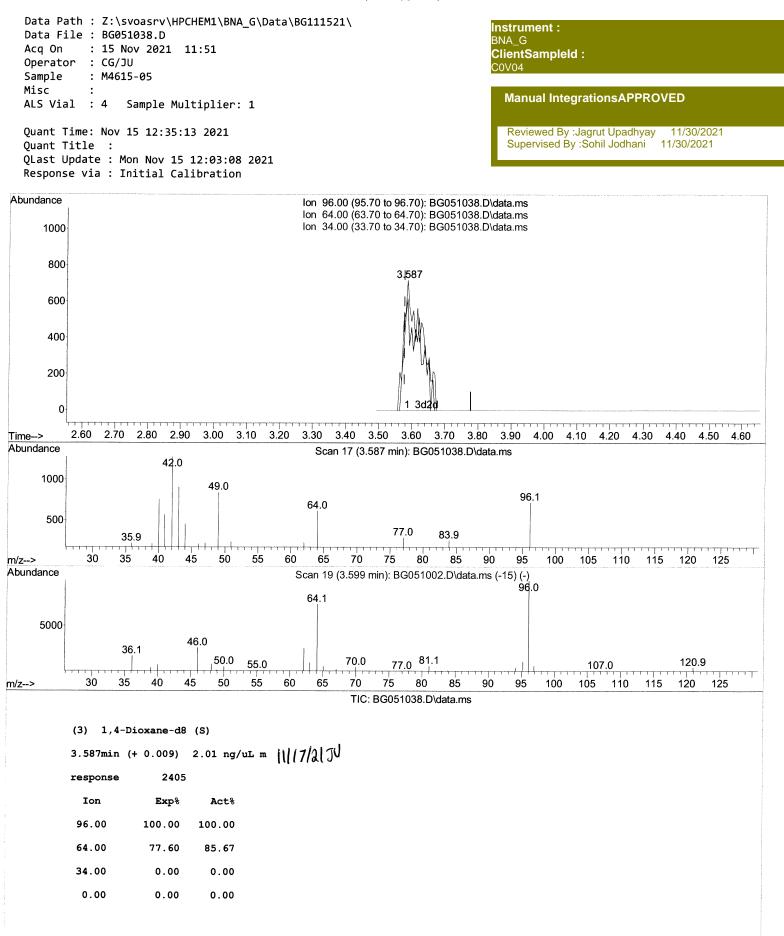
Quantitation Report (Qedit)

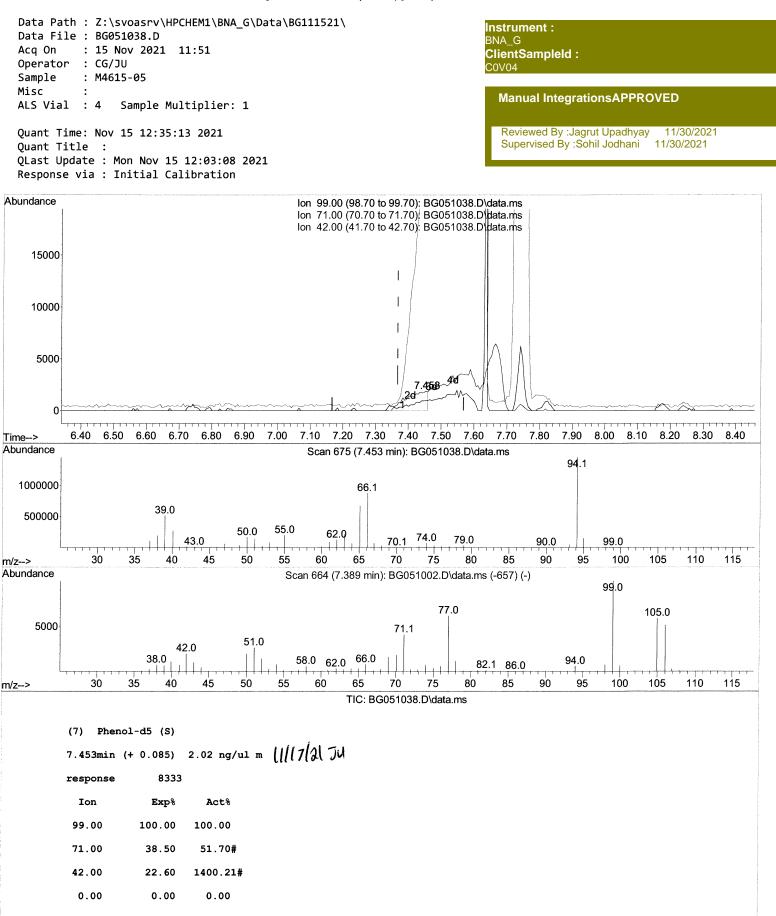




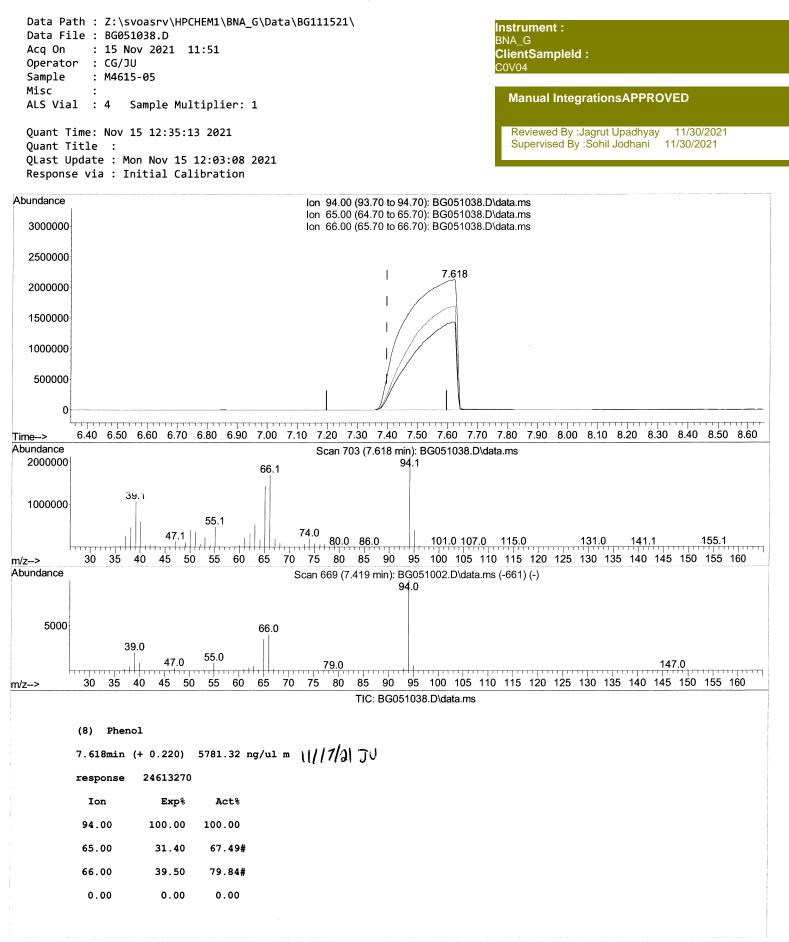


Quantitation Report (Qedit)









Data Path : Z:\svoasrv\HPCHEM1\ Data File : BG051038.D Acq On : 15 Nov 2021 11:51 Operator : CG/JU Sample : M4615-05 Misc : ALS Vial : 4 Sample Multipli	er: 1	Instrument : BNA_G ClientSampleId : C0V04 Manual IntegrationsAPPROVED							
Quant Time: Nov 15 12:35:13 2021 Quant Title : QLast Update : Mon Nov 15 12:03:08 2021						Reviewed By :Jagrut Upadhyay 11/30/2021 Supervised By :Sohil Jodhani 11/30/2021			
Response via : Initial Calibration Compound R.T. QIon Response Conc Units Dev(Min)									
Internal Standards									
 1,4-Dichlorobenzene-d4 	8.240	152	38578	20.000	ng/ul	0.01			
20) Naphthalene-d8	11.055			20.000 i	ng/ul	0.00			
38) Acenaphthene-d10	14.850 17.594	164	121019	20.000 i	ng/ul	0.00			
64) Phenanthrene-d10	17.594	188	250296	20.000 i	ng/ul	0.00			
79) Chrysene-d12	21.895			20.000 r	ng/ul	0.00			
88) Perylene-d12	25.309	264	203356	20.000 r	ng/ul	0.00			
System Monitoring Compounds									
3) 1,4-Dioxane-d8	3.587	96	2405m 🕆	2 012	aα/uL >	0.00 11/17/21JU			
4) Pyridine-d5	0.000	84		0 000 -	~~/]				
7) Phenol-d5	7.453	99		> 2 025 r	ng/u⊥ ng/u1 ≯	0.09 11/17/21 7			
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.671	67		4.980 r	ng/ui	0.13			
11) 2-Chlorophenol-d4	7.800	132	26272	9.211 r		0.05			
15) 4-Methylphenol-d8	8.945	113	31900	9.211 1 9.846 r	•	0.02			
21) Nitrobenzene-d5	9.416	128	15635	10.083 r		0.01			
24) 2-Nitrophenol-d4	10.138	143	18026		-	0.01			
28) 2,4-Dichlorophenol-d3	10.138	165	29422	10.455 r 10.130 r		0.02			
31) 4-Chloroaniline-d4	11.202	131	1560	0.355 r	-	0.01			
46) Dimethylphthalate-d6	14.245	166	95317	10.295 r	0	0.00			
49) Acenaphthylene-d8	14.551	160	124663	10.295 r 10.807 r	0	0.00			
					-				
60) Fluorene-d10	15.062	143	14189	8.452 n	0	0.02			
-	15.837		81205	9.901 n	•	0.00			
65) 4,6-Dinitro-2-methylph	17 (04	200	12800	8.435 n	-	0.01			
 73) Anthracene-d10 81) Pyrene-d10 92) Benzo(a) pyrene-d12 	10.000	188	131324	11.097 n	•	0.00			
81) Pyrene-d10	19.980	212	139647	10.149 n		0.00			
92) Benzo(a)pyrene-d12	25.062	264	112612	10.017 n	ig/uI	0.00			
Target Compounds				Qva	lue unalo au				
8) Phenol	7.618	94	24613270m>5			1100 (117/21 Ju			
<pre>13) 2-Methylphenol</pre>	8.687	108	8579	2.726 n	g/ul	99			
16) Acetophenone	9.075		90546	17.990 n		98			
18) 4-Methylphenol	9.004	108	68647	20.488 n	g/ul	91			

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