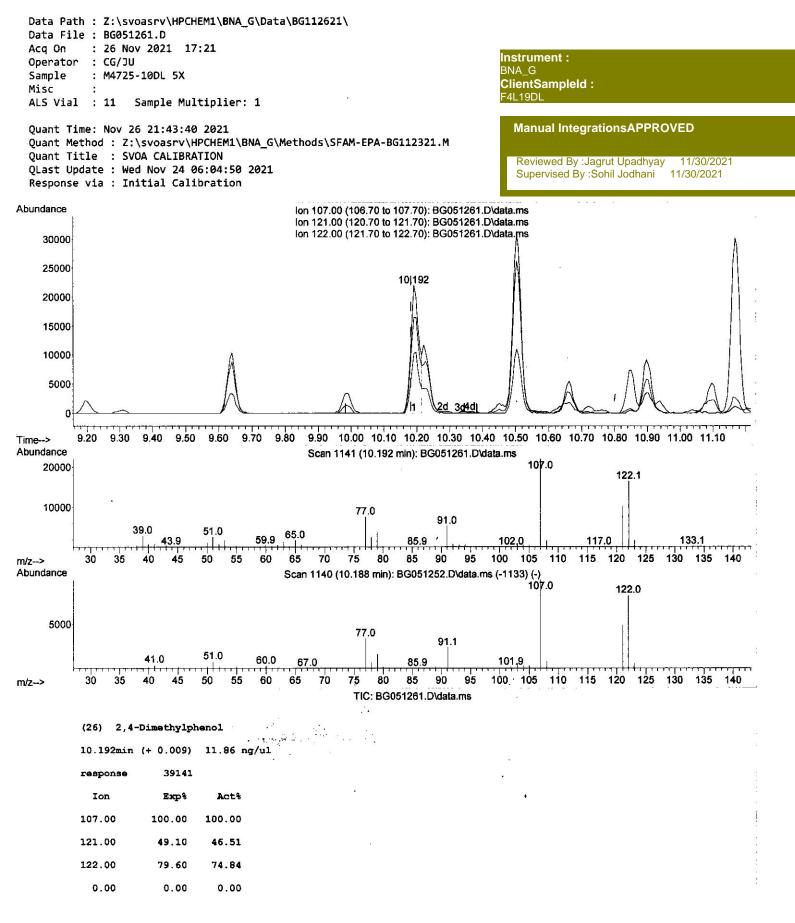
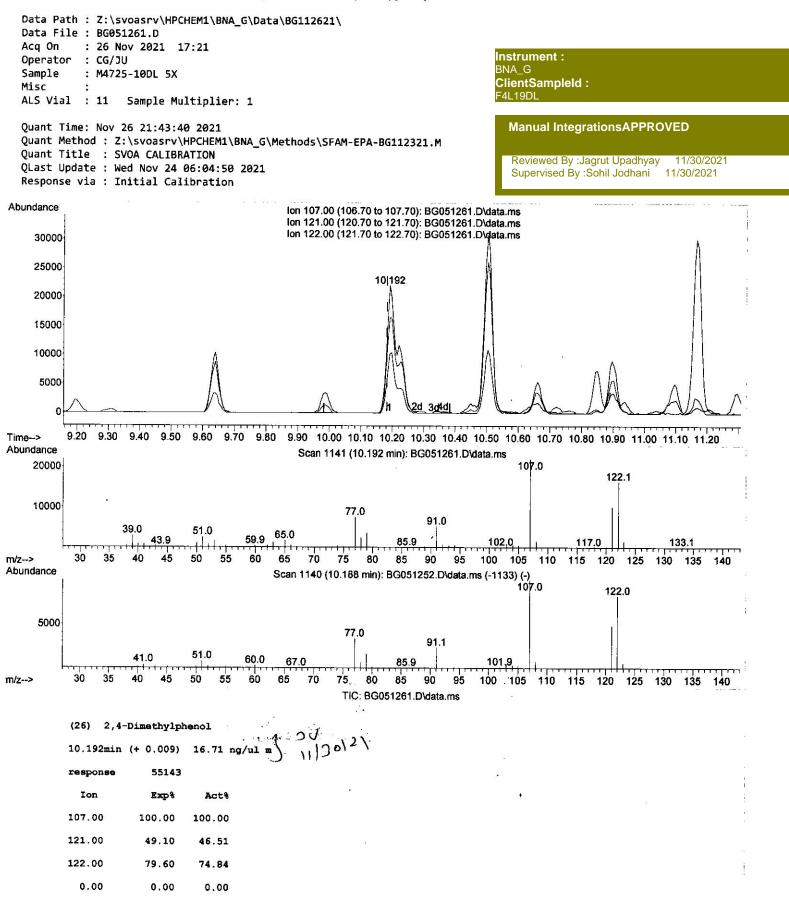


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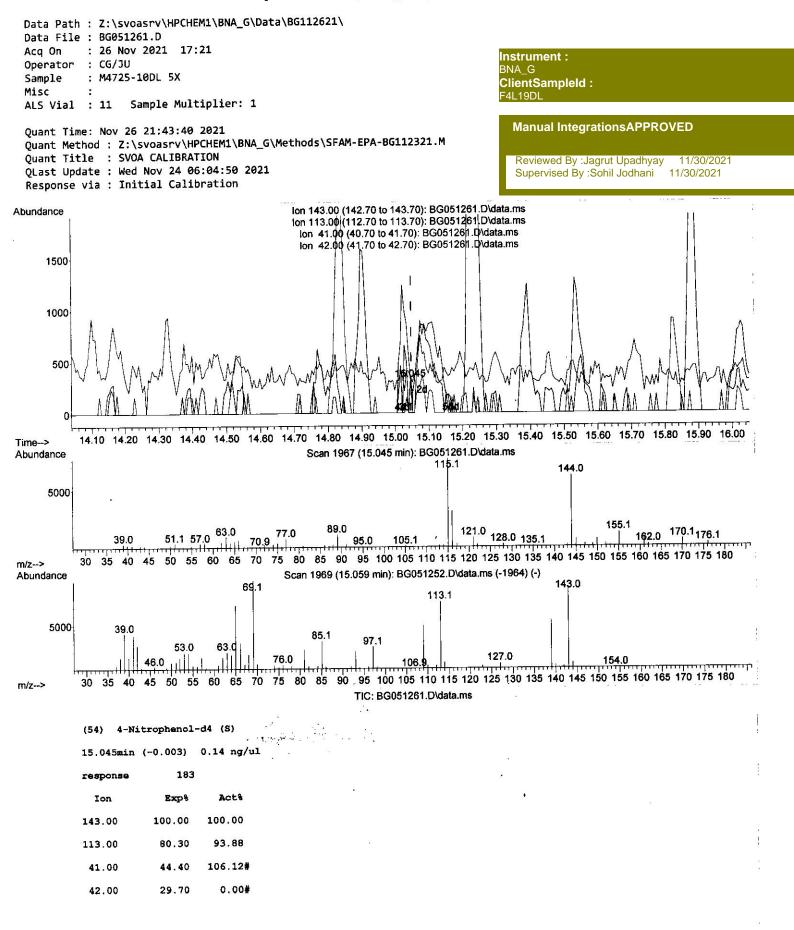
Quantitation Report (Qedit)



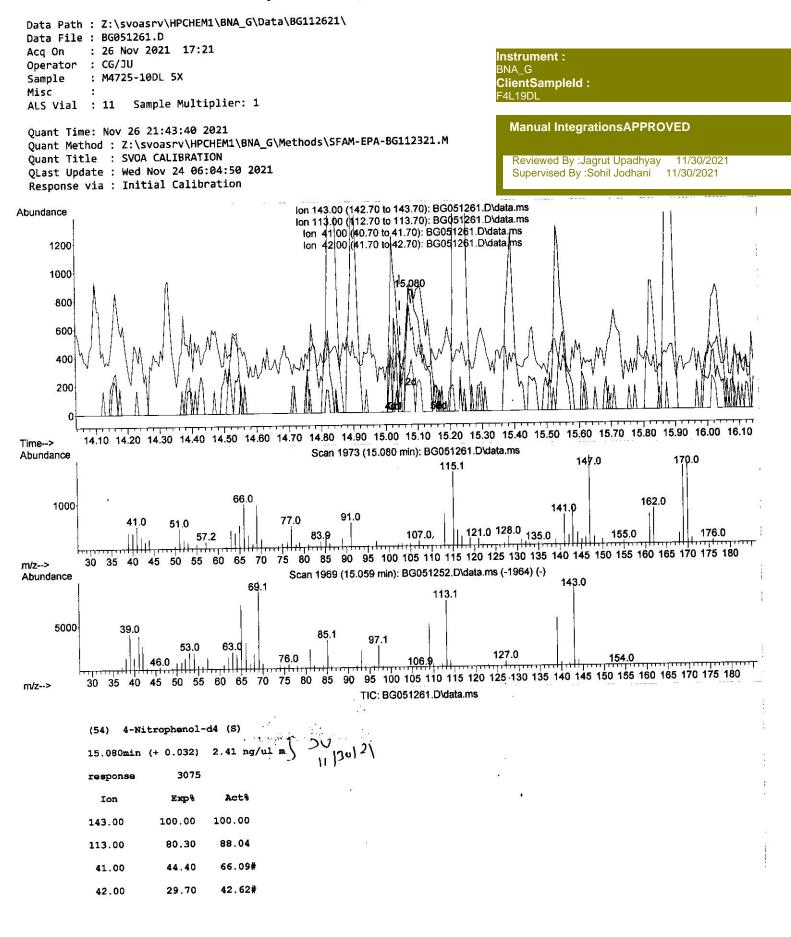
SFAM-EPA-BG112321.M Fri Nov 26 22:11:53 2021

Page: 1

Quantitation Report (Qedit)



Quantitation Report (Qedit)



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112621\						
Data File : BG051261.D						
Acg On : 26 Nov 2021 17:21						
Operator : CG/JU						strument :
Sample : M4725-10DL 5X						NA_G
Misc :						lientSampleId :
MISC : ALS Vial : 11 Sample Multiplier: 1						
Quant Time: Nov 26 21:43:40 2021						Manual IntegrationsAPPROVED
Quant Method : Z:\svoasrv\HPCHEM	1\BNA_G\I	Method	Is\SFAM-EPA	-BG112321.M		
Quant Title : SVOA CALIBRATION						Reviewed By :Jagrut Upadhyay 11/30/2021
QLast Update : Wed Nov 24 06:04:						Supervised By :Sohil Jodhani 11/30/2021
Response via : Initial Calibrati	.on					
Compound	R.T.	QIon	Response	Conc Units De	v(Min)	
Internal Standards					0.00	
 1,4-Dichlorobenzene-d4 	8.206		37091	20.000 ng/ul		
20) Naphthalene-d8	11.038		163668	20.000 ng/ul		
38) Acenaphthene-d10	14.839		102476	20.000 ng/ul		
64) Phenanthrene-d10	17.589		195889	20.000 ng/ul		
79) Chrysene-d12	21.890		170078	20.000 ng/ul		
88) Perylene-d12	25.303	264	170504	20.000 ng/ul	· 0.02	
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	Ød	0.000 ng/uL		
4) Pyridine-d5	3.987	84	3640	1.162 ng/ul		с.
7) Phenol-d5	7.372	99	5409	1.476 ng/ul		
<pre>9) Bis-(2-Chloroethyl)eth</pre>		67	13622	5.917 ng/ul	0.00	. 1
11) 2-Chlorophenol-d4	7.736	132	13015	4.930 ng/ul	0.00	
15) 4-Methylphenol-d8	8.917	113	9614	3.250 ng/ul	0.00	
21) Nitrobenzene-d5	9.387	128	8800	6.369 ng/ul		
24) 2-Nitrophenol-d4	10.109	143	9393	6.027 ng/ul		
28) 2,4-Dichlorophenol-d3	10.662	165	16019	6.058 ng/ul		
31) 4-Chloroaniline-d4	11.173		19647	5.078 ng/ul	20 CO. 10	
46) Dimethylphthalate-d6	14.228	166	53207	6.748 ng/ul		200121
49) Acenaphthylene-d8	14.534		68530 1	6.892 ng/ul		11/20121
54) 4-Nitrophenol-d4	15.080	143	3075m	2.409 ng/ul		\mathcal{M}^{1}
60) Fluorene-d10	15.826	176	47119 /	6.636 ng/ul		
65) 4,6-Dinitro-2-methylph		200	0d	0.000 ng/u]		
73) Anthracene-d10	17.683		70714	7.548 ng/ul 7.287 ng/ul		
81) Pyrene-d10	19.963		74987 65043	7.143 ng/ul		
92) Benzo(a)pyrene-d12	25.063	264	03043	7.145 HB/03	0.01	
Target Compounds				C	value	2)1
13) 2-Methylphenol	8.652	108	19743 ()	6.980 ng/u]	97	30
26) 2,4-Dimethylphenol	10.192	107	55143m	16.708 ng/u]		1/30/21
30) Naphthalene	11.097		2836914	318.557 ng/u	l# 92	
36) 2-Methylnaphthalene	12.671		25945	4.283 ng/u	L 92	
37) 1-Methylnaphthalene	12.894	142	70251	11.273 ng/u		
43) 1,1'-Biphenyl	13.664	154	14050	1.836 ng/u	L 97	
52) Acenaphthene	14.904	153	294241	45.419 ng/u		
56) Dibenzofuran	15.233		239458	25.626 ng/u		
61) Fluorene	15.885		187902	25.104 ng/u		
72) Phenanthrene	17.630		101783	9.411 ng/u		
74) Anthracene	17.724		29674	2.762 ng/u		
77) Carbazole	17.994		355545	37.709 ng/u		
80) Fluoranthene	19.634		53193	4.208 ng/u		
82) Pyrene	19.998	202	30973	2.505 ng/u	T4 20	

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

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