Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

Data File : BG050873.D

Acq On : 4 Nov 2021 8:34

Operator : CG/JU Sample : M4412-19

Misc

ALS Vial : 60 Sample Multiplier: 1

Quant Time: Nov 07 08:35:46 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

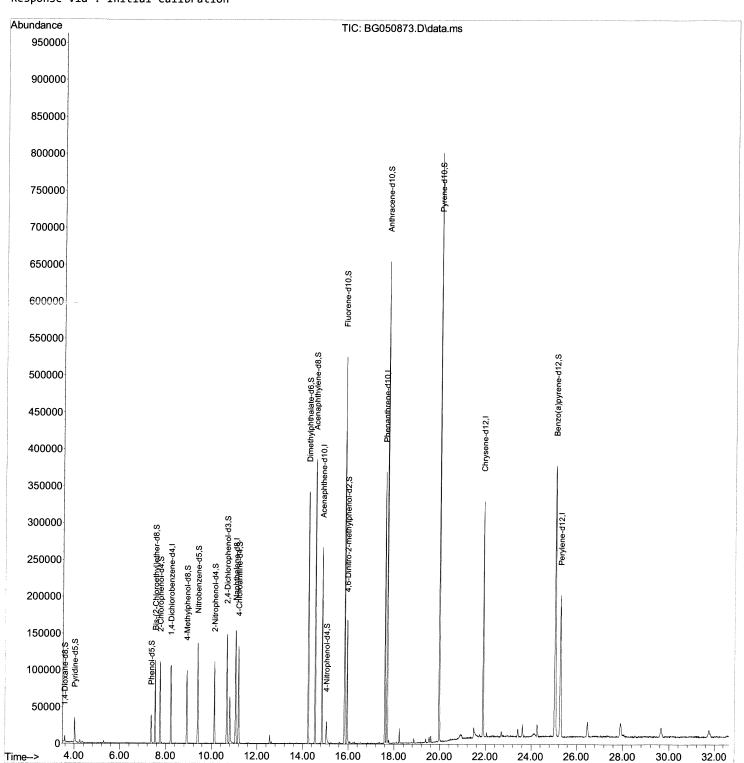
Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration

Instrument : BNA_G ClientSampleId : BG373

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/08/2021 Supervised By :mohammad ahmed 11/08/2021



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

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Acq On : 4 Nov 2021 8:34

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Quant Time: Nov 07 08:35:46 2021

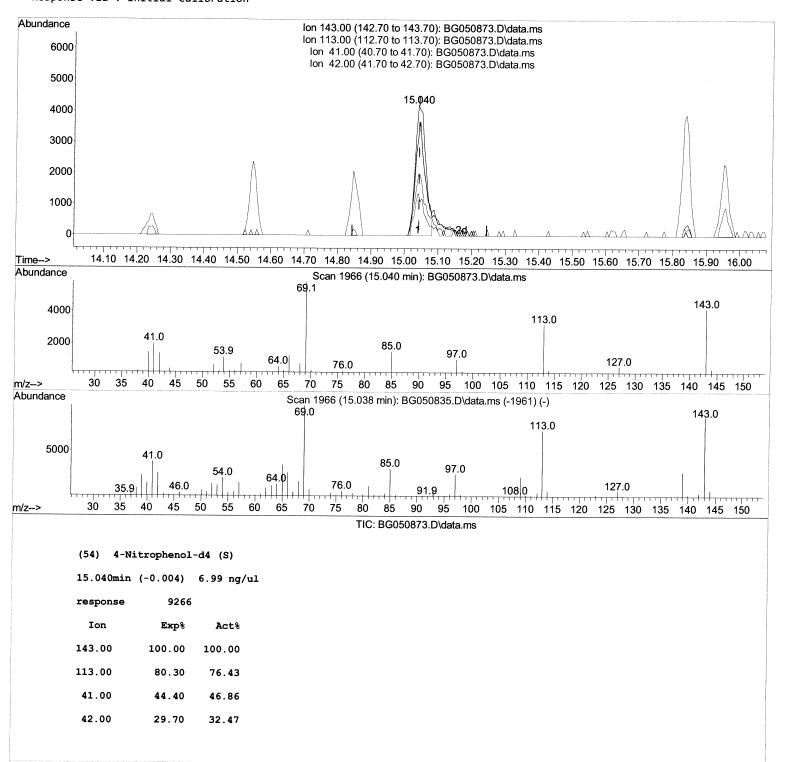
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration Instrument :
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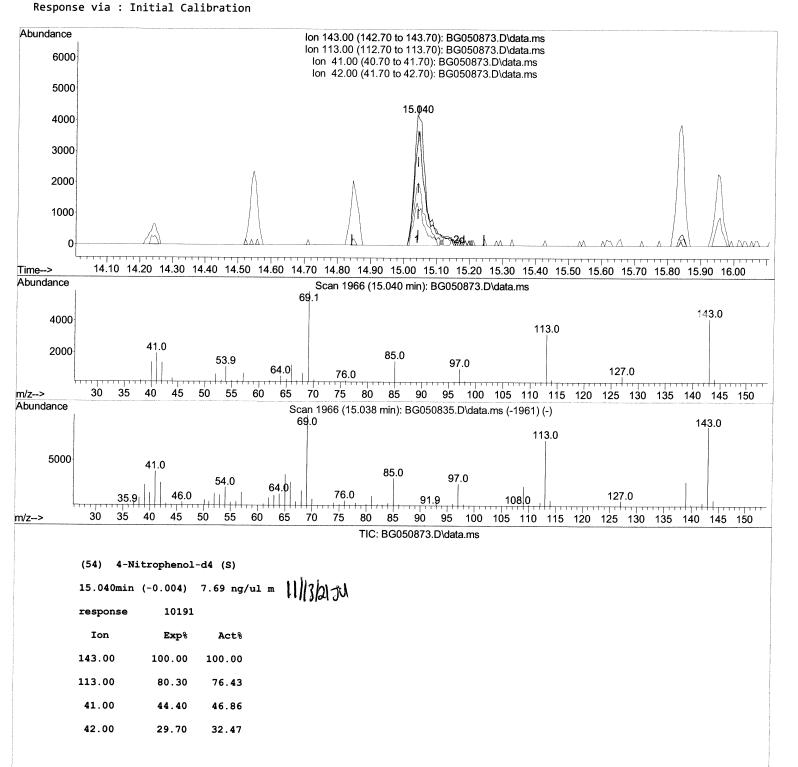
 $\label{lem:quant_method} \textbf{Quant Methods} : \textbf{Z:} \\ \textbf{SPAM-EPA-BG110321.M} \\ \\ \textbf{Quant Methods} : \textbf{Z:} \\ \textbf{SPAM-EPA-BG110321.M} \\ \\ \textbf{Quant Methods} : \textbf{Z:} \\ \textbf{Quant Methods} : \textbf{Quant Methods} : \textbf{Q:} \\ \textbf{Quant Methods} : \textbf{Q:} \\ \textbf{Quant Methods} : \textbf{Q:} \\ \textbf{Q:}$

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:49:05 2021

Instrument: BNA_G ClientSampleId :

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Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\

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Acq On : 4 Nov 2021 8:34 Operator : CG/JU

Sample : M4412-19

Misc

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Quant Title : SVOA CALIBRATION

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Compound	R.T.	QIon	Response	Conc Units Dev(M	lin)
Internal Standards					
 1,4-Dichlorobenzene-d4 	8.230	152	29193	20.000 ng/ul	0.00
20) Naphthalene-d8	11.050	136	131454	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.852	164	95501	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.596	188	220196	•	0.00
79) Chrysene-d12	21.891	240	191726	<u>o</u> ,	-0.02
88) Perylene-d12	25.293	264	188141	•	-0.02
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.583	96	4378	4.840 ng/uL	0.00
4) Pyridine-d5	4.018	84	23335	•	0.00
7) Phenol-d5	7.372	99	24446	•	0.00
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.543	67	57171	O,	0.00
11) 2-Chlorophenol-d4		132	50210	O.	0.00
15) 4-Methylphenol-d8	8.924	113	40055	•	0.00
21) Nitrobenzene-d5	9.399	128	33707	<u>.</u>	0.00
24) 2-Nitrophenol-d4	10.128	143	34480	Q .	0.00
28) 2,4-Dichlorophenol-d3	10.669	165	55761	O.	0.00
31) 4-Chloroaniline-d4	11.186	131	72624	<u>o</u> .	0.00
46) Dimethylphthalate-d6	14.247	166	230783	O .	0.00
49) Acenaphthylene-d8	14.546	160	277777	-	0.00
54) 4-Nitrophenol-d4	15.040	143	10191m>		t/98/111 00.0
60) Fluorene-d10	15.839	176	204410		9.00
65) 4,6-Dinitro-2-methylph	15.956	200	36956	<u>.</u>	9.00
73) Anthracene-d10	17.696	188	377899	Ο,	9.00
81) Pyrene-d10	19.969	212	438717	35.428 ng/ul (
92) Benzo(a)pyrene-d12	25.058	264	370724	_	9.02
arget Compounds				Ovalı	ıe

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