Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\

Data File : BG050882.D

Acq On : 8 Nov 2021 16:12

Operator : CG/JU Sample : M4492-02

Misc

ALS Vial : 8 Sample Multiplier: 1

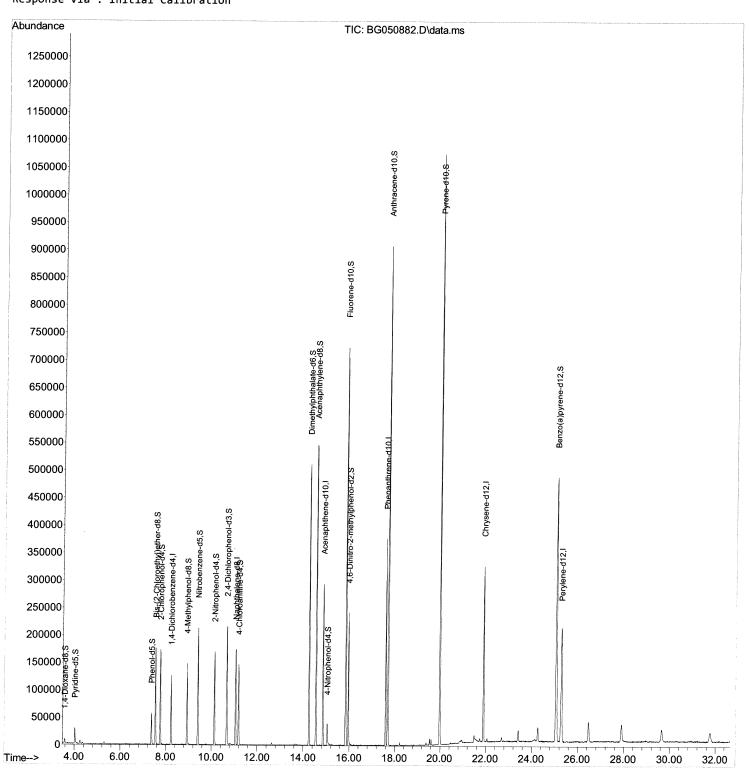
Quant Time: Nov 08 16:01:48 2021

Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : BG1R9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/08/2021 Supervised By :mohammad ahmed 11/09/2021



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\

Data File : BG050882.D

Acq On : 8 Nov 2021 16:12

Operator : CG/JU Sample : M4492-02

Misc

ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 08 16:01:48 2021

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

Quant Title : SVOA CALIBRATION

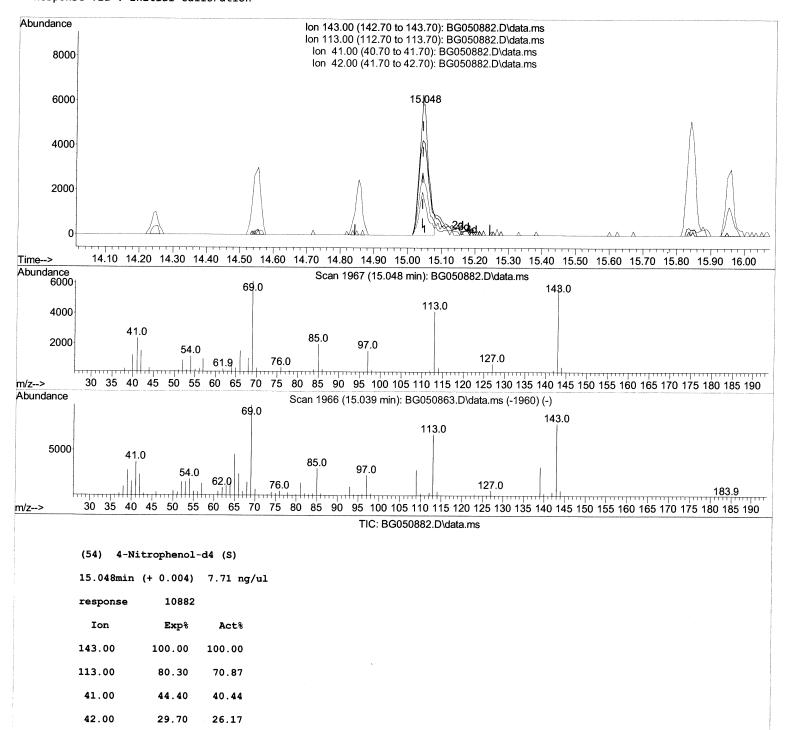
QLast Update : Tue Nov 02 14:49:05 2021 Response via: Initial Calibration

Instrument: BNA_G ClientSampleId :

BG1R9

Manual Integrations APPROVED

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Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG110821\

Data File : BG050882.D

Acq On : 8 Nov 2021 16:12

Operator : CG/JU Sample : M4492-02

Misc

ALS Vial : 8 Sample Multiplier: 1

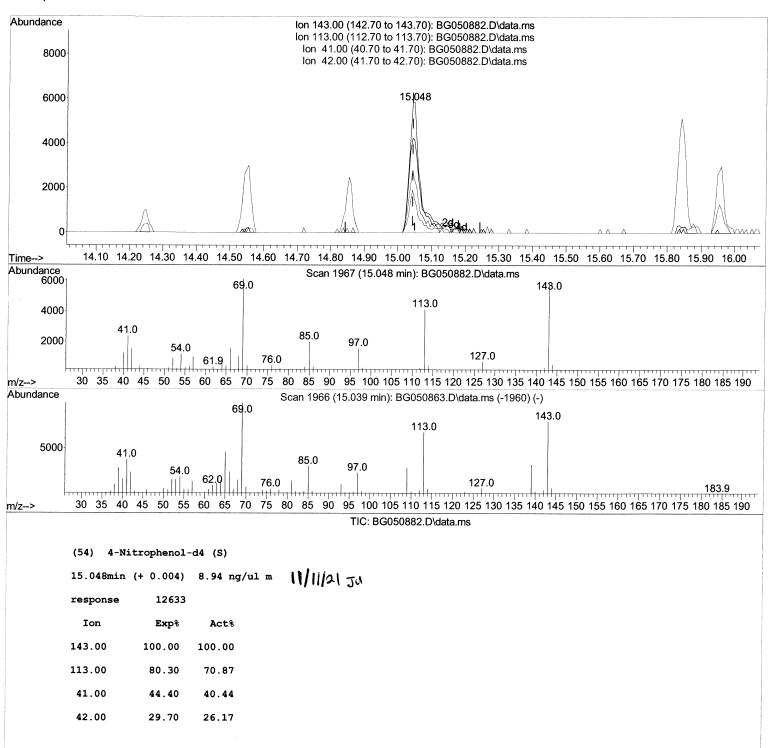
Quant Time: Nov 08 16:01:48 2021

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 ClientSampleld: BG1R9

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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc Units Dev(Min)
Internal Standards					-
 1,4-Dichlorobenzene-d4 	8.233	152	33891	20.000 ng/ul 0	.00
20) Naphthalene-d8		136	149456	20.000 ng/ul 0	.00
38) Acenaphthene-d10	14.854	164	101811	_	.00
64) Phenanthrene-d10	17.598	188	226803	20.000 ng/ul 0	.00
79) Chrysene-d12	21.899	240	195955	20.000 ng/ul 0	.00
88) Perylene-d12	25.295	264	195411	20.000 ng/ul -0	.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.585	96	4637	4.416 ng/uL 0.	00
4) Pyridine-d5	4.020	84	17961	5.718 ng/ul 0.	00
7) Phenol-d5	7.375	99	35304	9.764 ng/ul 0.	00
9) Bis-(2-Chloroethyl)eth	7.545	67	90902	38.921 ng/ul 0.	00
	7.757		77009	30.734 ng/ul 0.	0 0
<pre>15) 4-Methylphenol-d8</pre>	8.926	113	57536	20.214 ng/ul 0.	0 0
21) Nitrobenzene-d5	9.402	128	50680	39.903 ng/ul 0.	0 0
24) 2-Nitrophenol-d4	10.130	143	54475	38.574 ng/ul 0.	0 0
28) 2,4-Dichlorophenol-d3	10.671	165	80519	33.847 ng/ul 0.	00
31) 4-Chloroaniline-d4	11.188	131	79860	22.168 ng/ul 0.	00
46) Dimethylphthalate-d6	14.249	166	330102	42.380 ng/ul 0.	0 0
	14.555	160	400940	41.315 ng/ul 0.	00
54) 4-Nitrophenol-d4	15.048	143	12633m	> 8.945 ng/ul > 0.0	00 IIIIAI JU
60) Fluorene-d10	15.841	176	291648	42.267 ng/ul 0.0	
65) 4,6-Dinitro-2-methylph	15.959	200	53913	39.208 ng/ul 0.0	9 0
73) Anthracene-d10	17.698	188	517121	48.226 ng/ul 0.0	20
81) Pyrene-d10	19.972	212	595221	47.029 ng/ul 0.0	30
92) Benzo(a)pyrene-d12	25.066	264	498490	46.145 ng/ul 0.0	90
Target Compounds Qvalue					

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