# Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051324.D

Acq On : 3 Dec 2021 11:14

Operator : CG/JU Sample : M4833-21

Misc

ALS Vial : 29 Sample Multiplier: 1

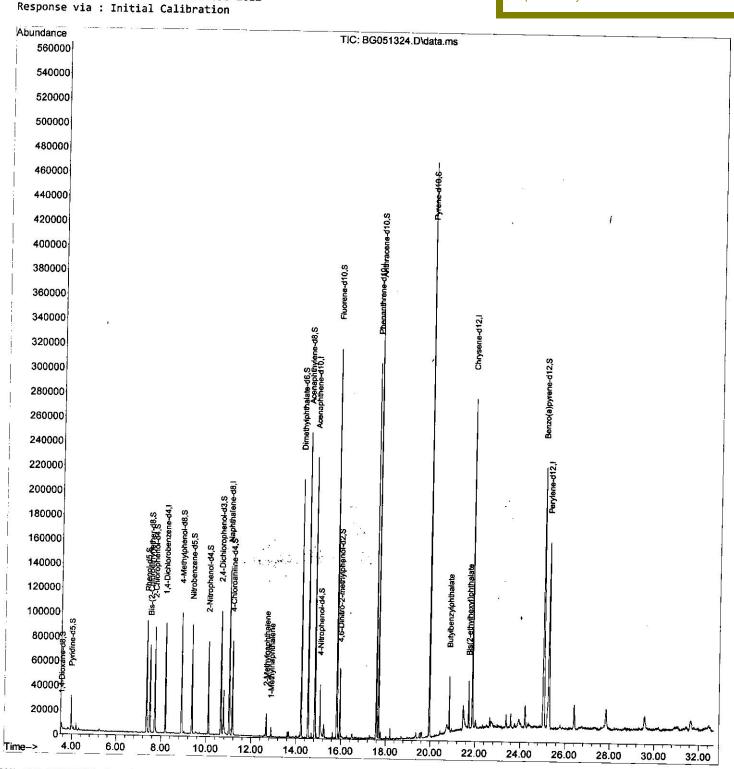
Quant Time: Dec 03 12:13:49 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Instrument :
BNA\_G
ClientSampleId :

## **Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



FAM-EPA-BG112321.M Fri Dec 03 12:16:09 2021

### Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120221\

Data File : BG051324.D

Acq On : 3 Dec 2021 11:14

Operator : CG/JU Sample : M4833-21

Misc

ALS Vial : 29 Sample Multiplier: 1

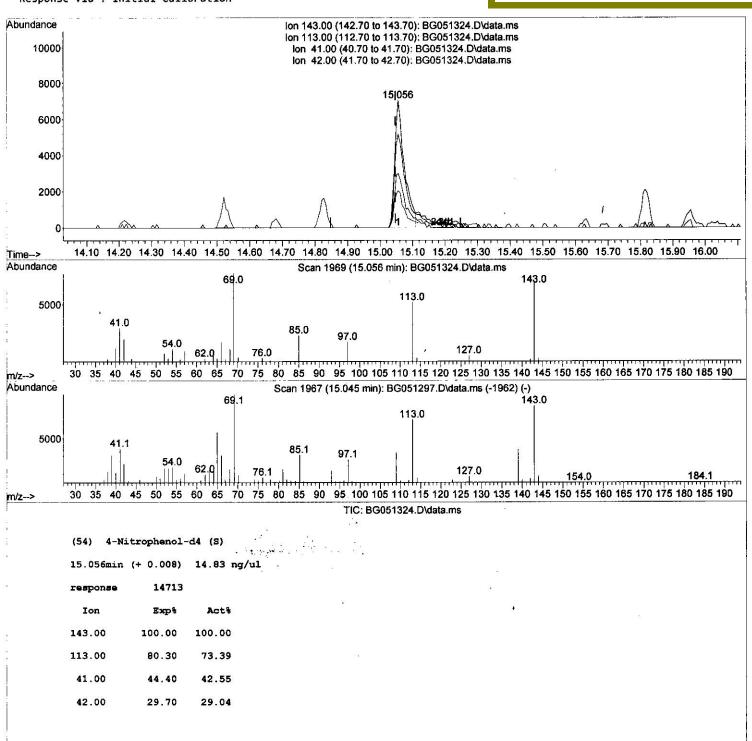
Quant Time: Dec 03 12:13:49 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument : BNA\_G ClientSampleId :

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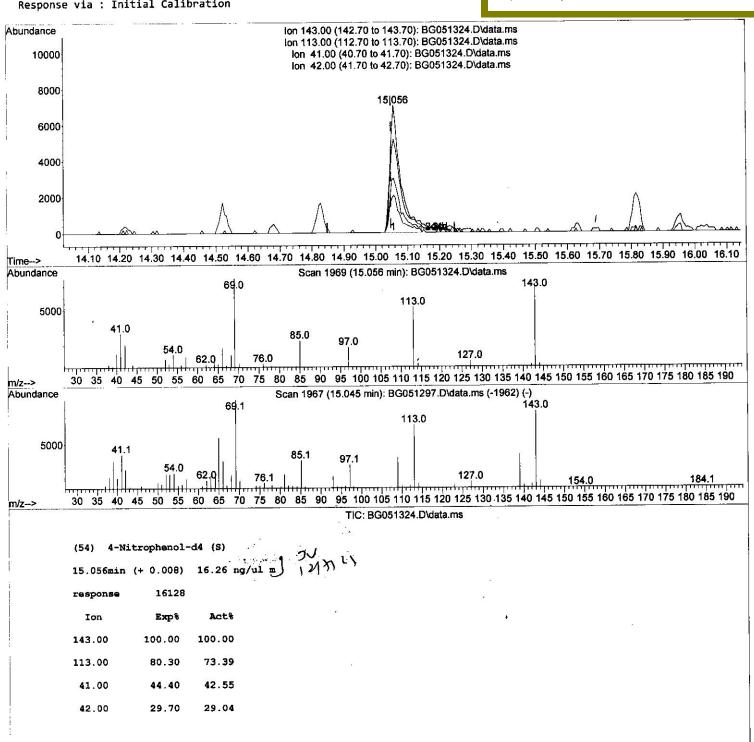
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#### Manual IntegrationsAPPROVED

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Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
Internal Standards					
<ol> <li>1,4-Dichlorobenzene-d4</li> </ol>	8.188	152	25871	20 000 ==/7	0.00
20) Naphthalene-d8	11.020		116348	20.000 ng/ul 20.000 ng/ul	-0.02
38) Acenaphthene-d10	14.827	164	79659	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.577	188	181355		0.00
79) Chrysene-d12	21.877	240	169450	20.000 ng/ul	0.00
88) Perylene-d12	25.279	264	166258	20.000 ng/ul 20.000 ng/ul	0.00 0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.528	96	2771	3 733 na/w	0.00
<ol><li>Pyridine-d5</li></ol>	3.963	84	18934	3.722 ng/uL	-0.02
7) Phenol-d5	7.353	99	56805	8.667 ng/ul	-0.02
9) Bis-(2-Chloroethyl)eth	7.506	67	37113	22.216 ng/ul	0.00
11) 2-Chlorophenol-d4	7.723	132	41893	23.111 ng/ul	0.00
15) 4-Methylphenol-d8	8.904	113	42398	22.753 ng/ul	0.00
21) Nitrobenzene-d5	9.374	128	22034	20.548 ng/ul	0.00
24) 2-Nitrophenol-d4	10.097	143	25404	22.435 ng/ul 22.930 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.644	165	41157	21.895 ng/ul	0.00
31) 4-Chloroaniline-d4	11.161	131	44464	16.166 ng/ul	0.00
46) Dimethylphthalate-d6	14.222	166	145304	23.706 ng/ul	0.00
49) Acenaphthylene-d8	14.521	160	184700	23.897 ng/ul	0.00
54) 4-Nitrophenol-d4	15.056	143	16128m	16.256 ng/ul	0.00
60) Fluorene-d10	15.814	176	128909	23.355 ng/ul	0.00
65) 4,6-Dinitro-2-methylph	15.955	200	13408	11.981 ng/ul	0.00
73) Anthracene-d10	17.676	188	213890	24.660 ng/ul	0.00
81) Pyrene-d10	19.956	212	258975	25.258 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.044	264	219610	24.733 ng/ul	0.00 0.00
Target Compounds				0	•
36) 2-Methylnaphthalene	12.665	142	9994	Qva	
37) 1-Methylnaphthalene	12.888	142	4647	2.321 ng/ul 1.049 ng/ul#	92
83) Butylbenzylphthalate	20.843	149	13317	2.600 ng/ul#	91
86) Bis(2-ethylhexyl)phtha	21.719	149	16953	2.300 ng/ul#	98 98
					70 

(#) = qualifier out of range (m) = manual integration (+) = signals summed . I final a suit of the

Instrument: BNA\_G ClientSampleId: ESQN8

## **Manual IntegrationsAPPROVED**

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12/1/21