Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120921\

Data File : BG051452.D

Acq On : 10 Dec 2021 9:52

Operator : CG/JU

Sample : M4870-08DL2 50X

Misc

ALS Vial : 29 Sample Multiplier: 1

Quant Time: Dec 10 10:37:21 2021

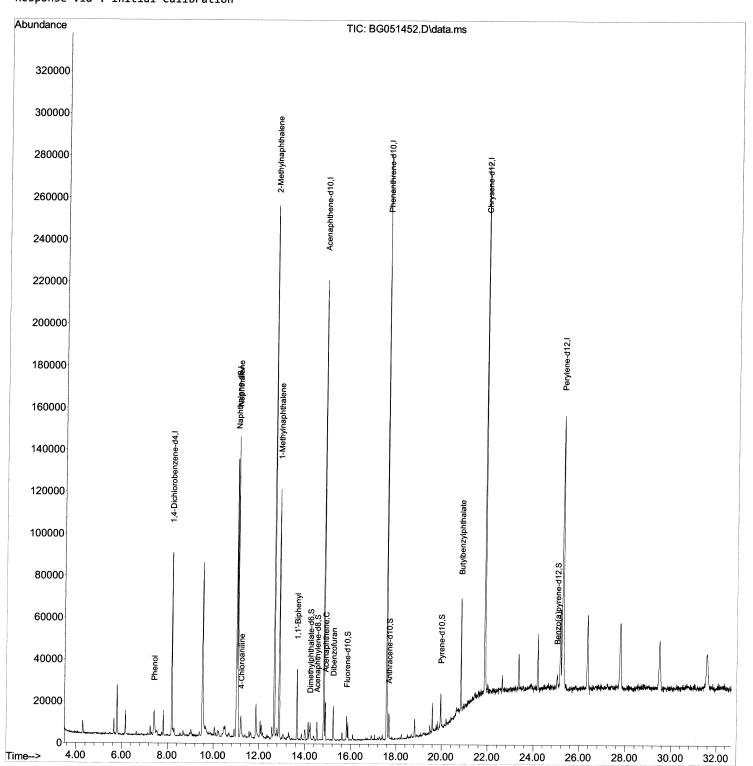
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration



## **Manual IntegrationsAPPROVED**



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120921\

Data File : BG051452.D

Acq On : 10 Dec 2021 9:52

Operator : CG/JU

Sample : M4870-08DL2 50X

Misc

ALS Vial : 29 Sample Multiplier: 1

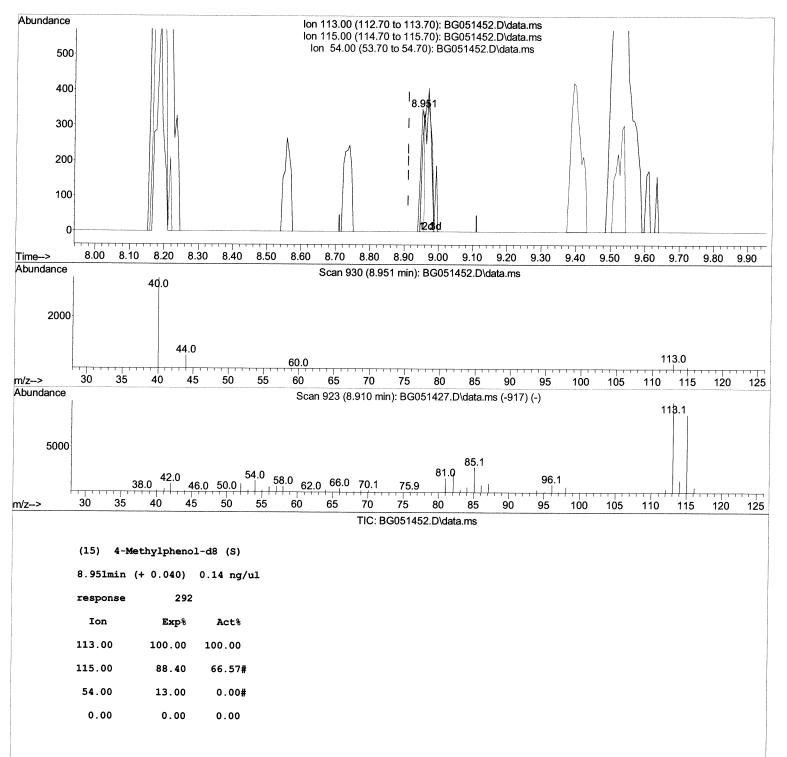
Quant Time: Dec 10 12:41:42 2021

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration



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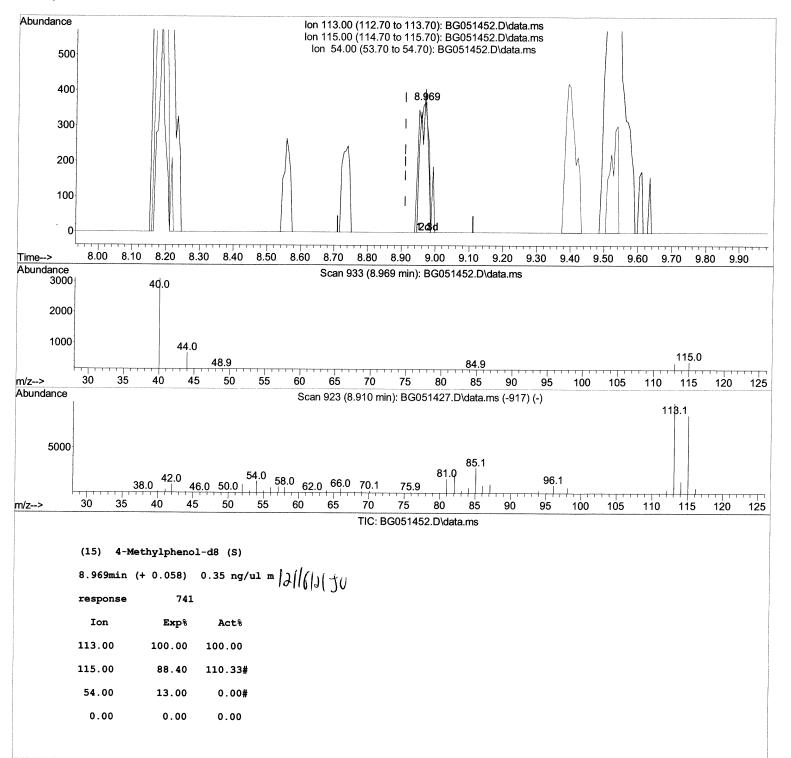
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG120821.M

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Misc

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ALS Vial : 29 Sample Multiplier: 1

Quant Time: Dec 10 12:41:42 2021

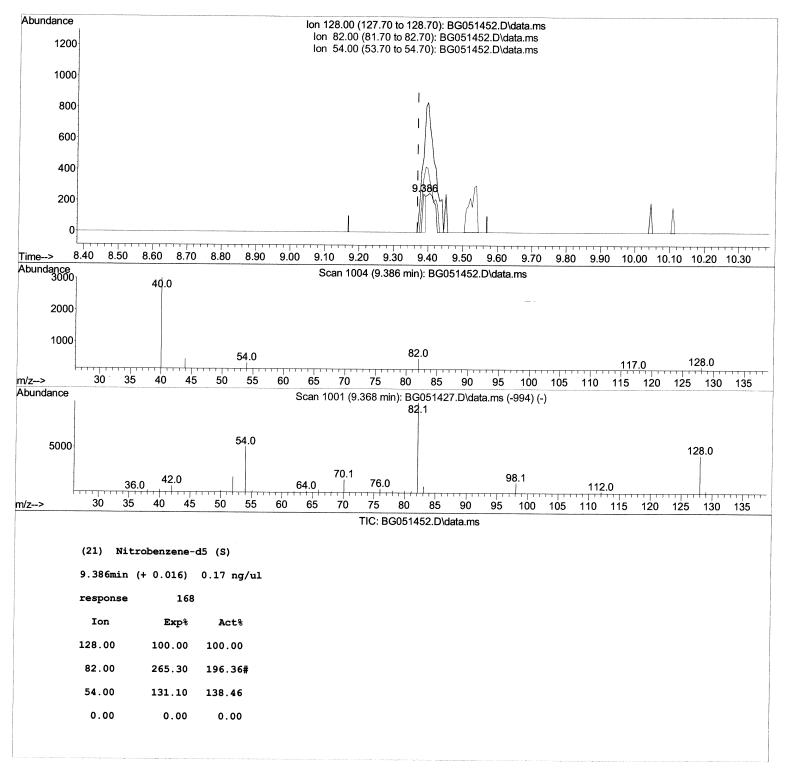
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration



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Quant Time: Dec 10 10:37:21 2021

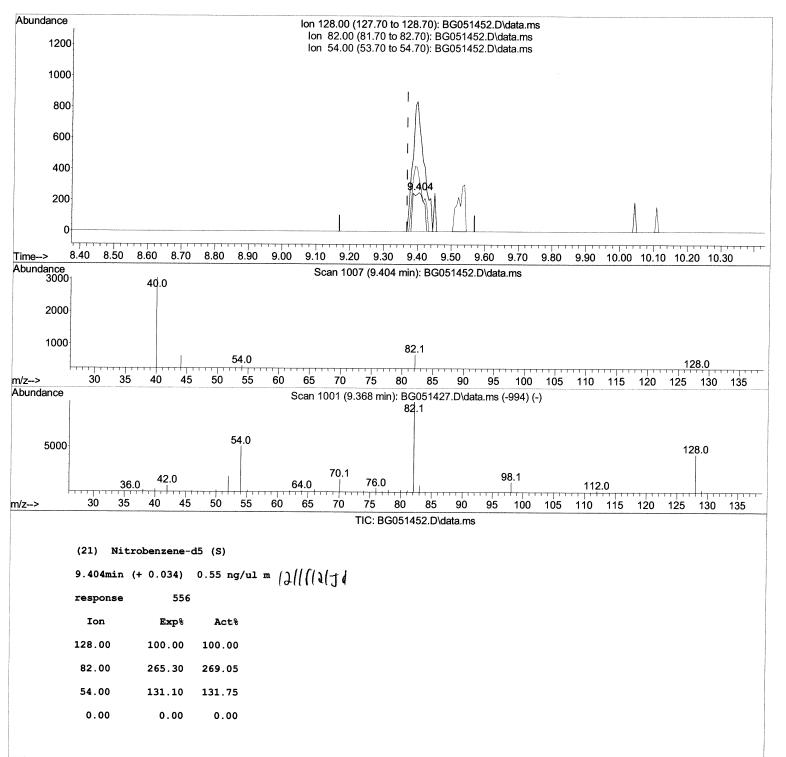
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG120821.M

Quant Title : SVOA CALIBRATION

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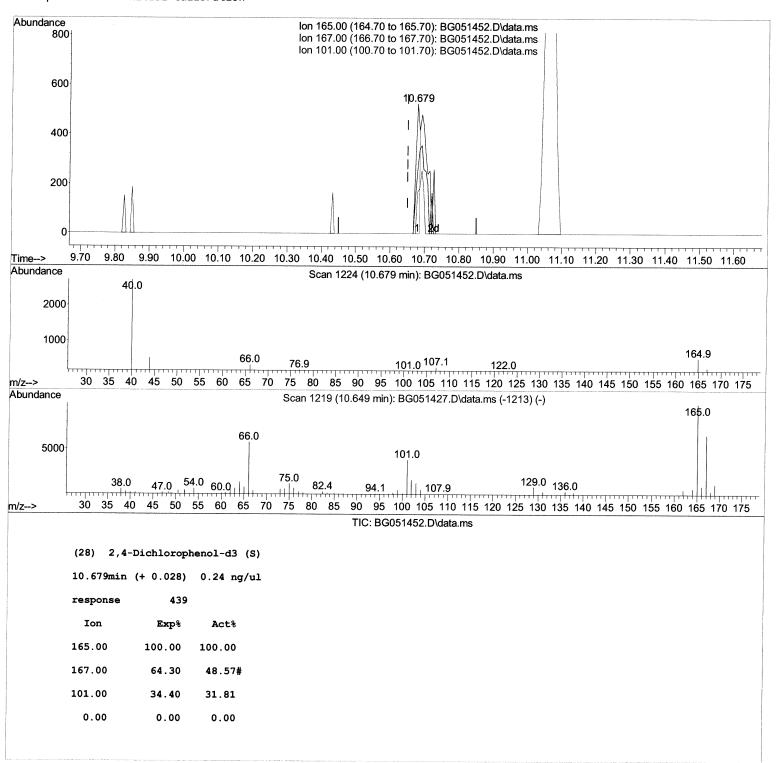
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG120821.M

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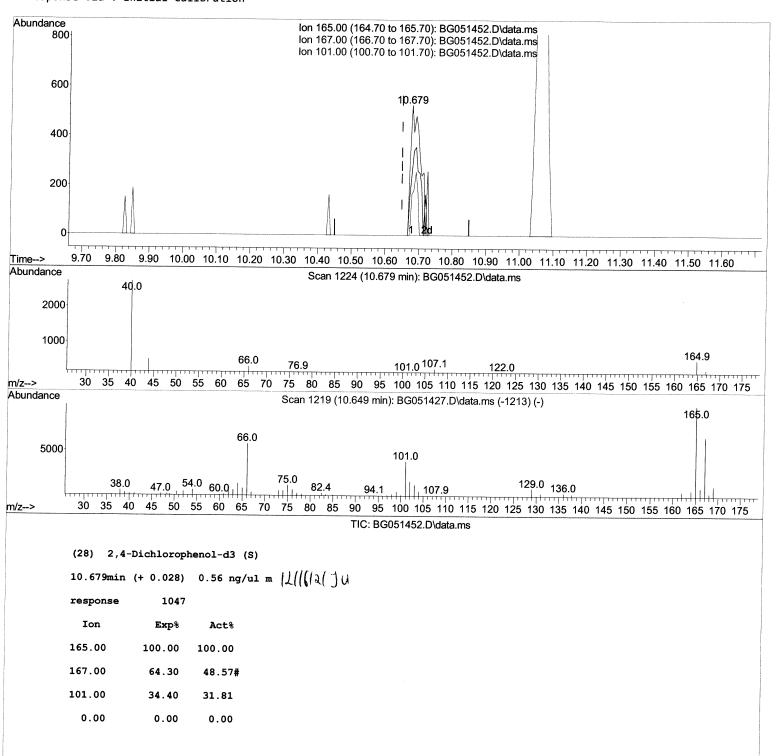
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Quant Time: Dec 10 10:37:21 2021

 $\label{thm:linear_Quant_Methods} \textbf{Quant Methods: Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG120821.M}$ 

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 03:21:41 2021 Response via : Initial Calibration Instrument: BNA\_G ClientSampleld: BGKP5DL2

# **Manual IntegrationsAPPROVED**

	Compound	R.T.	QIon	Response	Conc Ur	its Dev	(Min)
Inte	rnal Standards						
	1,4-Dichlorobenzene-d4	8.187	152	26178	20.000	ng/ul	0.00
	Naphthalene-d8	11.014	136	116618		ng/ul	0.00
•	Acenaphthene-d10	14.815		79754		ng/ul	0.00
-	Phenanthrene-d10	17.571		182122		ng/ul	0.00
	Chrysene-d12	21.871	240	167397		ng/ul	0.00
	Perylene-d12	25.267		159718		ng/ul	0.00
Syst	em Monitoring Compounds						
	1,4-Dioxane-d8	0.000	96	0	0.000	ng/uL	
	Pyridine-d5	0.000	84	Ød		ng/ul	
	Phenol-d5	0.000	99	0d		ng/ul	
9)	Bis-(2-Chloroethyl)eth	7.523	67	1021		ng/ul	0.02
	2-Chlorophenol-d4	7.747	132	1066		ng/ul	0.02
	4-Methylphenol-d8	8.969	113	741m >		ng/ul>	
	Nitrobenzene-d5	9.404	128	556m <	0.550	ng/ul	
24)	2-Nitrophenol-d4	10.120	143	480	a 110	ng/ul	0 02
	2,4-Dichlorophenol-d3		165	1047m	> 0.562	ng/ul >	= 0.03 12/11/12/J
	4-Chloroaniline-d4	11.196	131	1052	0.386	ng/ul	0.03
	Dimethylphthalate-d6	14.227	166	6342		ng/ul	0.01
	Acenaphthylene-d8	14.521	160	8039		ng/ul	0.00
	4-Nitrophenol-d4	0.000	143	0		ng/ul	
	Fluorene-d10	15.826	176	5844		ng/ul	0.02
	4,6-Dinitro-2-methylph	0.000	200	0		ng/ul	0.02
	Anthracene-d10	17.682	188	10797		ng/ul	0.01
81)	Pyrene-d10	19.962	212	11269		ng/ul	0.01
92)	Benzo(a)pyrene-d12	25.038	264	8686		ng/ul	0.00
Target Compounds						Qvalue	
8)	Phenol	7.418	94	11967	4.387	-	93
30)	Naphthalene	11.061	128	124634	19.461		97
	4-Chloroaniline	11.207	127	5305		ng/ul#	75
36)	2-Methylnaphthalene	12.659	142	121137	28.365	•	99
	1-Methylnaphthalene	12.876	142	55813	12.697		90
43)	1,1'-Biphenyl	13.657	154	19472	3.266		97
	Acenaphthene				1.713		97
	Dibenzofuran	15.226	168	8669 12942	1.804		93
83)	Butylbenzylphthalate	20.837	149	16833	3.175		89

<sup>(#) =</sup> qualifier out of range (m) = manual integration (+) = signals summed