Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112321\

Data File : BM033237.D

: 23 Nov 2021 15:03 Acq On

Operator : CG/JU

Sample : M4725-03DL 5X

Misc

ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 23 16:17:22 2021

Quant Method: Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Nov 19 15:41:12 2021

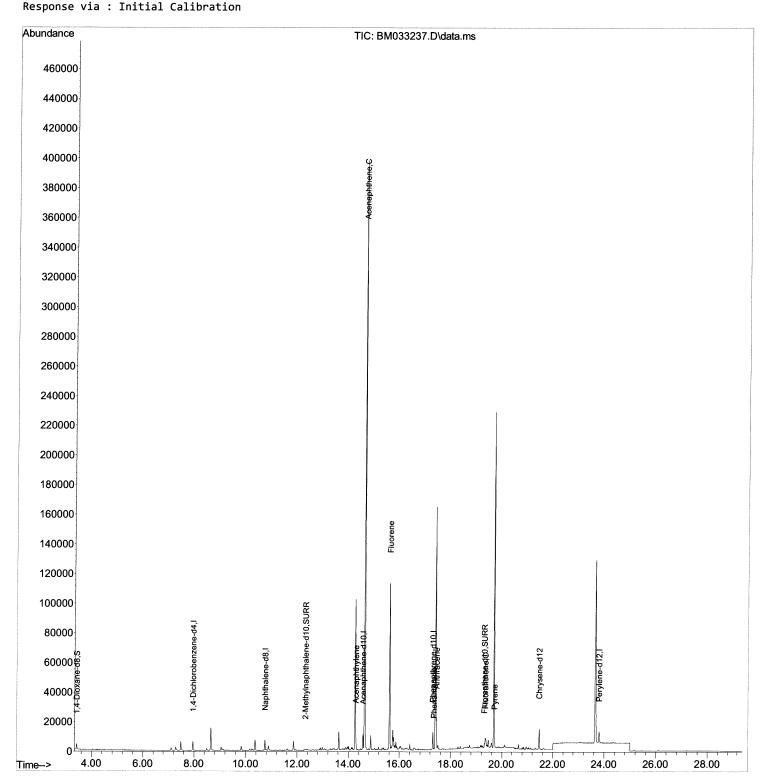
Instrument : BNA_M

ClientSampleId :

F4L06DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/23/2021 Supervised By :mohammad ahmed 11/26/2021



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112321\

Data File : BM033237.D

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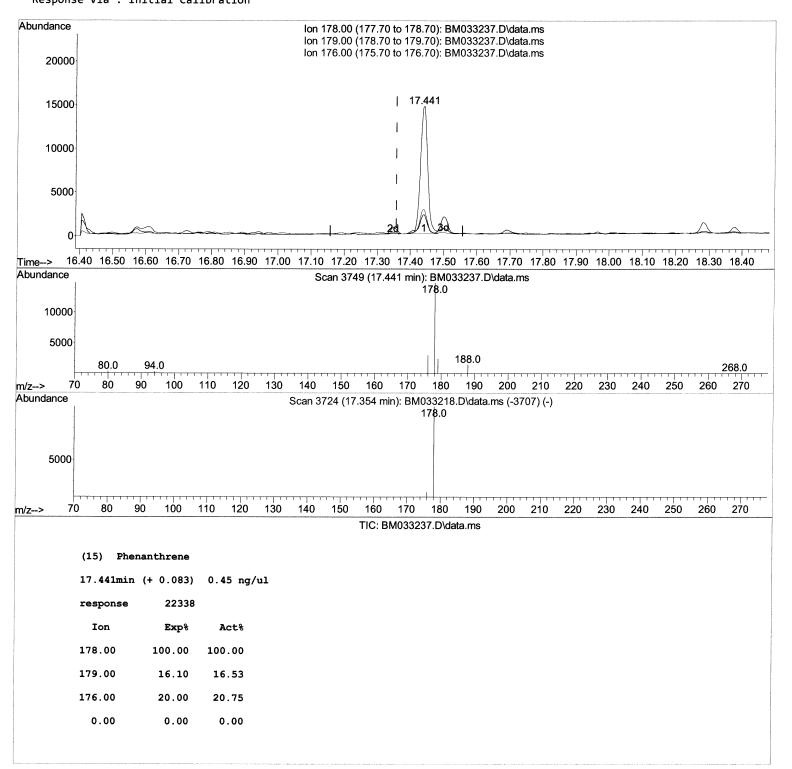
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Nov 19 15:41:12 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/23/2021 Supervised By :mohammad ahmed 11/26/2021



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112321\

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 $\label{thm:linear_quant_methods} \mbox{Quant Methods: $Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M} \label{thm:linear_methods}$

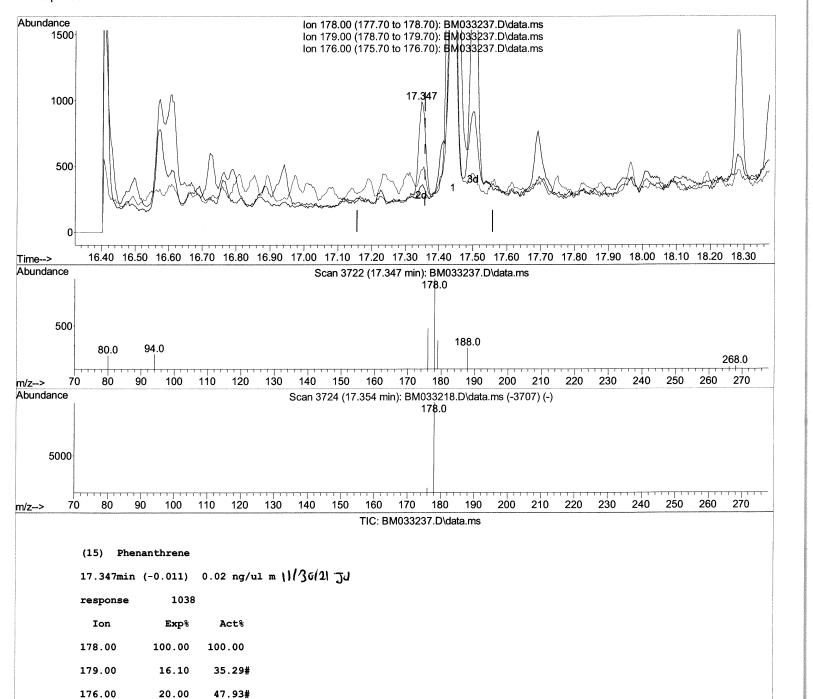
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Nov 19 15:41:12 2021 Response via : Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 11/23/2021 Supervised By: mohammad ahmed 11/26/2021



0.00

0.00

0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112321\

Data File : BM033237.D

Acq On : 23 Nov 2021 15:03

Operator : CG/JU Sample : M4725-: M4725-03DL 5X

Misc

ALS Vial : 8 Sample Multiplier: 1

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Nov 19 15:41:12 2021 Response via : Initial Calibration

Instrument : BNA_M ClientSampleId :

F4L06DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/23/2021 Supervised By:mohammad ahmed 11/26/2021

Compound	R.T.	QIon	Response	Conc Units Dev(Min)		
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.956	152	3215	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.750	136	9401	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.571	164	5864	0.400	ng/ul	#-0.01
13) Phenanthrene-d10	17.305	188	13572	0.400	ng/ul	-0.01
17) Chrysene-d12	21.465	240	13101	0.400	ng/ul	0.00
23) Perylene-d12	23.802	264	10505	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.413	96	2575	0.852	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.329	152	919	0.071	ng/ul	-0.01
18) Fluoranthene-d10	19.322	212	3869	0.102	ng/ul	-0.01
Target Compounds				Ovalue		
10) Acenaphthylene	14.291	152	5476	0.200	ng/ul#	90
11) Acenaphthene	14.632	153	232364	10.333	ng/ul	98
12) Fluorene	15.614	166	71081	2.801	ng/ul	98
15) Phenanthrene	17.347	178	1038m >	0.021	ng/ul≯	11/30/2174
16) Anthracene	17.441	178	22390	0.563	ng/ul	98
19) Fluoranthene	19.353	202	4318	0.068	ng/ul#	81
20) Pyrene	19.715	202	3827	0.061	ng/ul#	82

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