Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM111921\

Data File : BM033208.D

Acq On : 20 Nov 2021 10:58

Operator : CG/JU Sample : M4725-06

Misc

ALS Vial : 40 Sample Multiplier: 1

Quant Time: Nov 22 00:53:38 2021

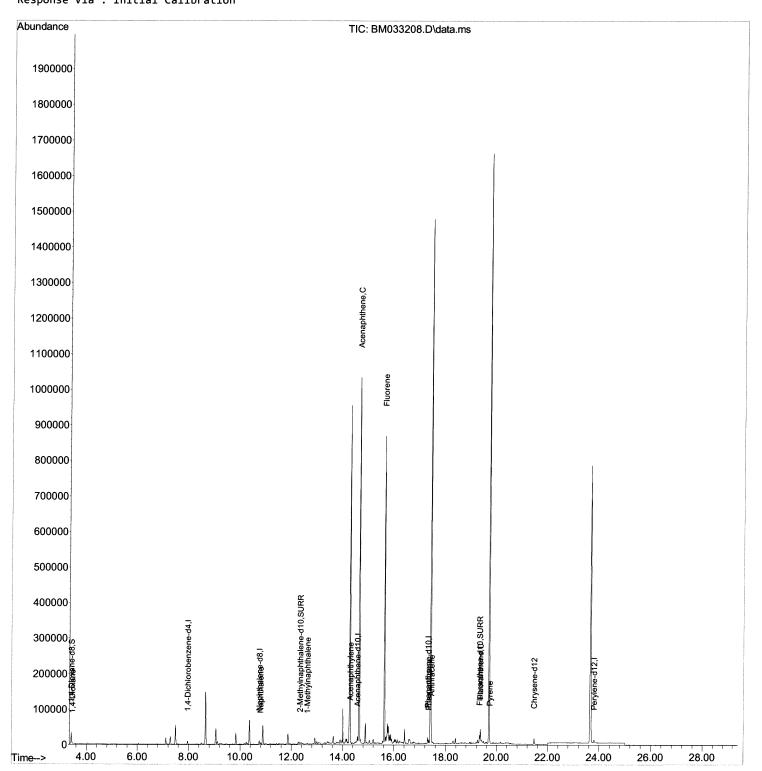
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM111921.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



# **Manual IntegrationsAPPROVED**



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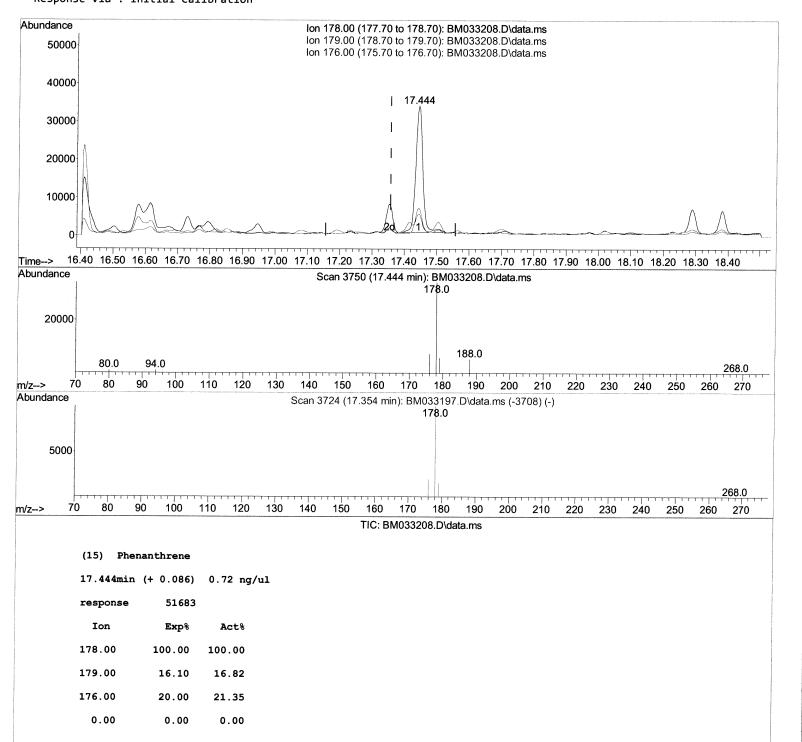
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM111921.M

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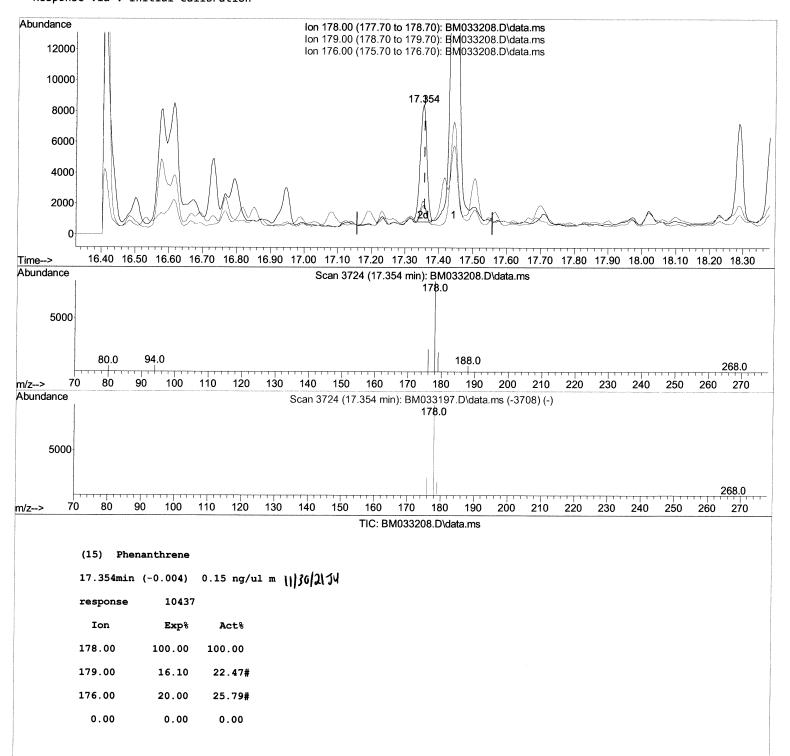
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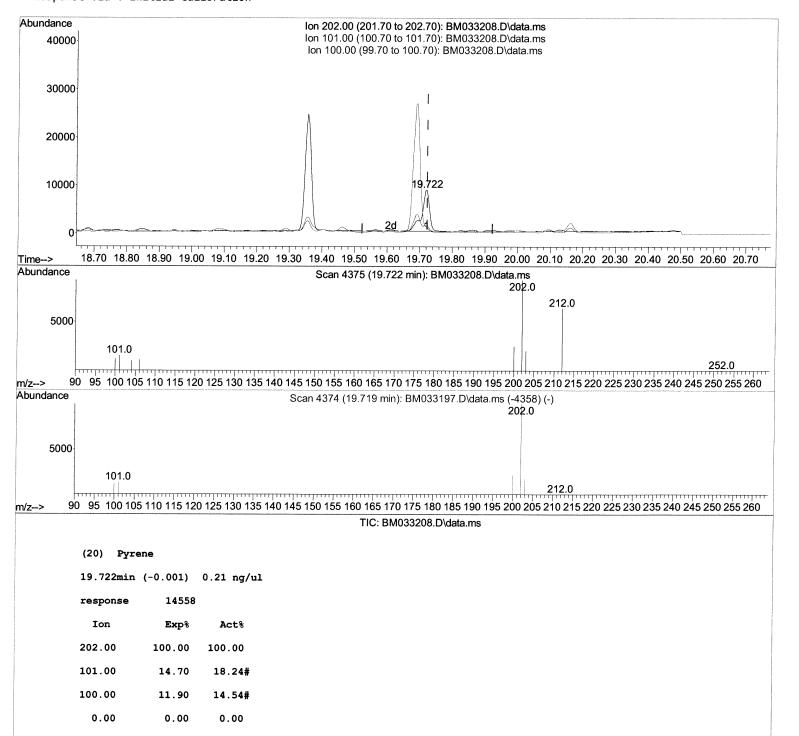
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM111921.M

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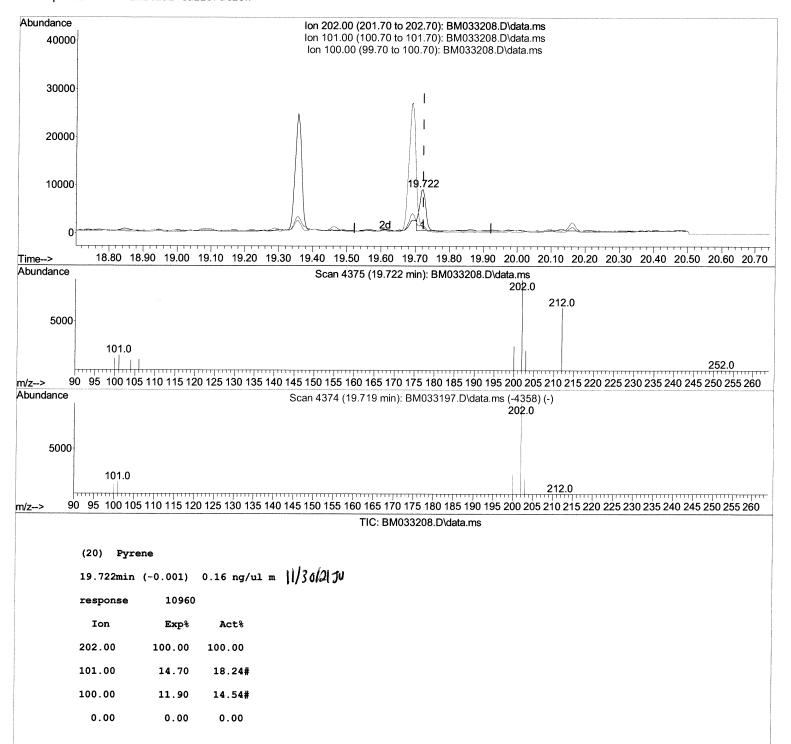
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM111921.M

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ClientSampleId:

F4L11

# Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc U	nits Dev	/(Min)
Internal Standards						
<ol> <li>1,4-Dichlorobenzene-d4</li> </ol>	7.960	152	4360	0.40	ng/ul	0.00
<ol><li>Naphthalene-d8</li></ol>	10.755	136	13556	0.40	ng/ul	# 0.00
<ol><li>9) Acenaphthene-d10</li></ol>	14.575	164	9534	0.40	ng/ul	# 0.00
13) Phenanthrene-d10	17.312	188	19831		ng/ul	0.00
17) Chrysene-d12	21.467	240	14748		ng/ul	
23) Perylene-d12	23.802	264	11526	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	19310	4.71	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.333	152	7590		ng/ul	0.00
18) Fluoranthene-d10	19.326	212	23472		ng/ul	
Target Compounds					Ov	alue
2) 1,4-Dioxane	3.448	88	348	0.08	ng/ul#	6
5) Naphthalene	10.804	128	2836		ng/ul#	73
<ol><li>8) 1-Methylnaphthalene</li></ol>	12.626	142	1827		ng/ul	
10) Acenaphthylene	14.298	152	11891		ng/ul#	
11) Acenaphthene	14.640	153	605696		ng/ul	
12) Fluorene	15.622	166	527980		ng/ul	
15) Phenanthrene	17.354	178	10437m >		•	1
16) Anthracene	17.444	178	50501		ng/ul	
19) Fluoranthene	19.356	202	32504		ng/ul	
20) Pyrene	19.722	202	10960m>		ng/ul>	

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