Quantitation Report (Qedit)

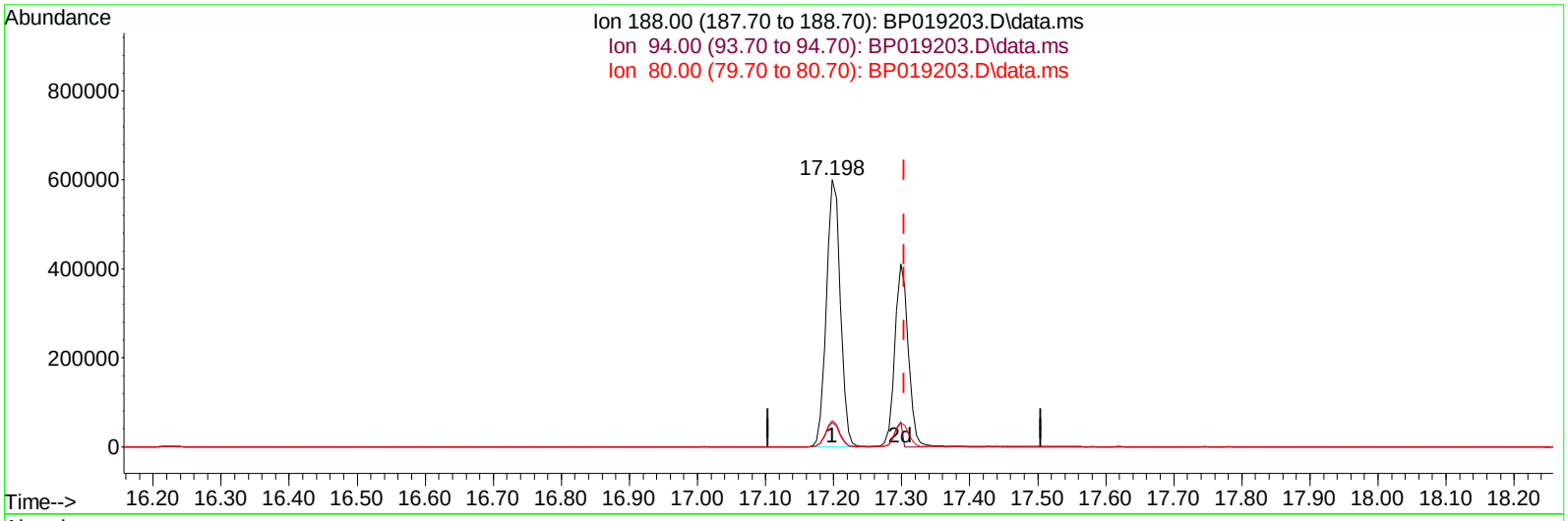
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019203.D
Acq On : 02 Jan 2024 12:18
Operator : MA/JU
Sample : 06020-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

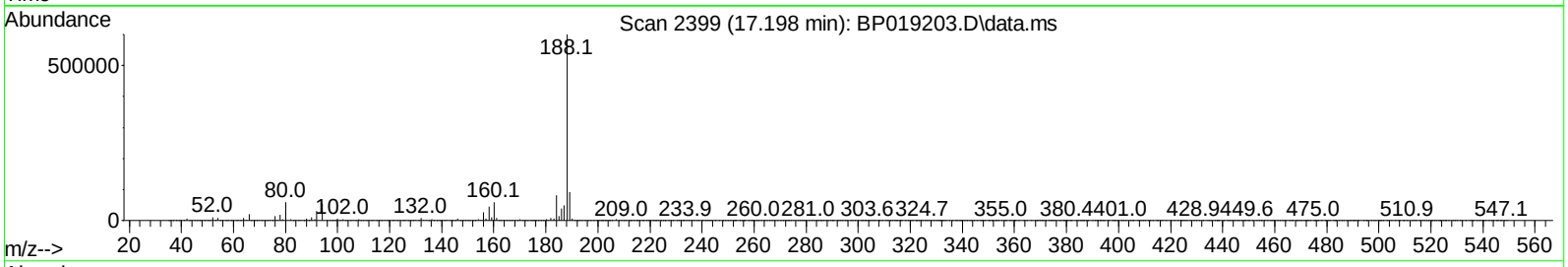
Quant Time: Jan 02 12:45:01 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 29 03:48:11 2023
Response via : Initial Calibration

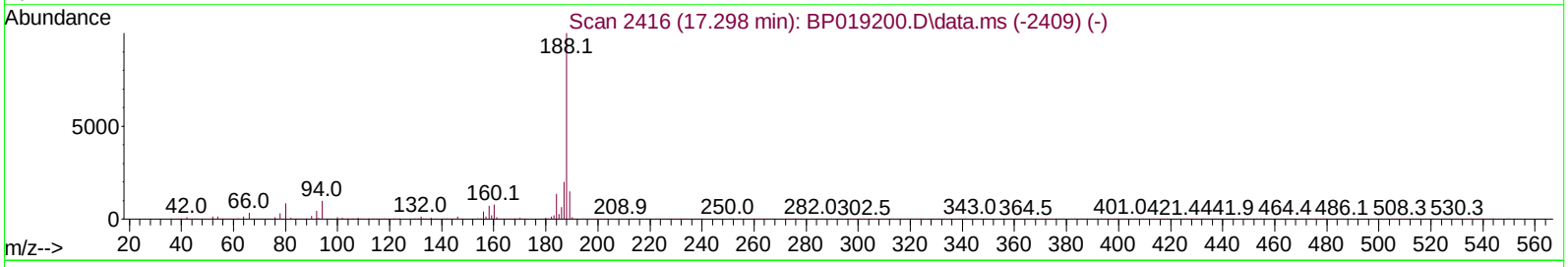
Instrument :
BNA_P
ClientSampleId :
BH3P1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/04/2024
Supervised By :mohammad ahmed 01/05/2024







TIC: BP019203.D\data.ms

(73) Anthracene-d10 (S)

17.198min (-0.106) 19.15 ng/ul

response 849778

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	10.50	9.08
80.00	9.10	9.87
0.00	0.00	0.00

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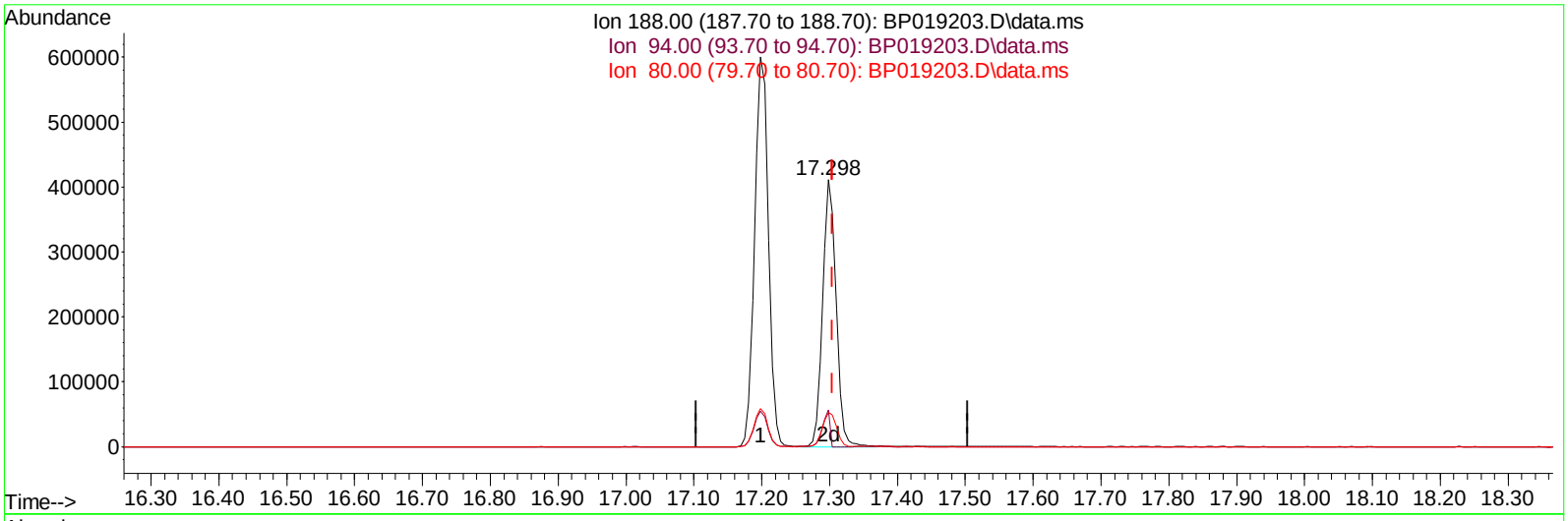

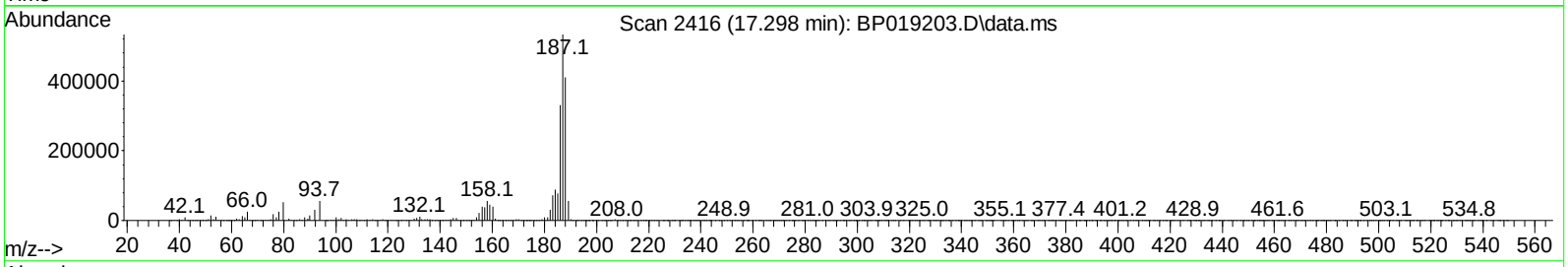

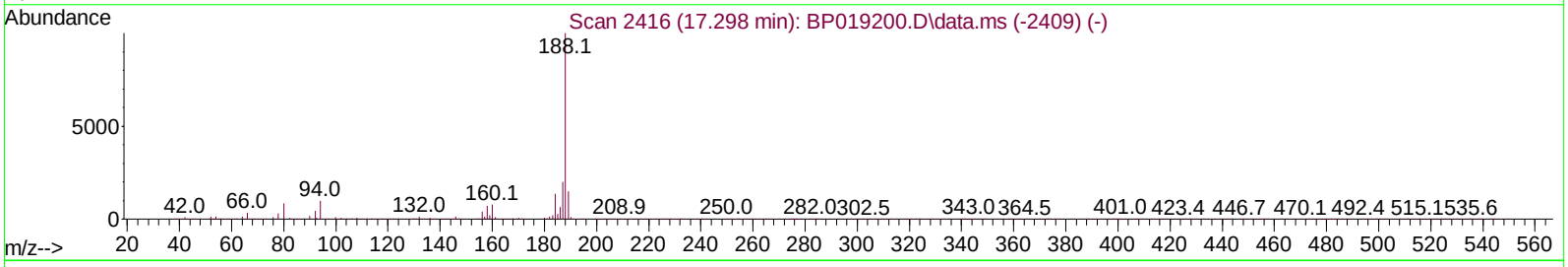

TIC: BP019203.D\data.ms

(73) Anthracene-d10 (S)

17.298min (-0.006) 12.78 ng/ul m

response 567312

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	10.50	13.78#
80.00	9.10	12.95#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
Data File : BP019203.D
Acq On : 02 Jan 2024 12:18
Operator : MA/JU
Sample : 06020-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Jan 02 12:46:28 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 29 03:48:11 2023
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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	169675	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	627483	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.446	164	326454	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	849778	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	920112	20.000	ng/ul	0.00
88) Perylene-d12	23.669	264	252575	20.000	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	24840	5.000	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.975	99	105550	6.937	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	380348	43.208	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	382904	31.215	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	152641	12.885	ng/ul	0.00
21) Nitrobenzene-d5	8.958	128	218372	40.259	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.681	143	225467	38.461	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	418089	35.319	ng/ul	0.00
31) 4-Chloroaniline-d4	10.722	131	57	0.004	ng/ul	-0.02
46) Dimethylphthalate-d6	13.863	166	1208393	41.907	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	62460	1.996	ng/ul	0.00
54) 4-Nitrophenol-d4	14.663	143	47649	9.786	ng/ul	0.00
60) Fluorene-d10	15.440	176	971075	38.982	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	181201	32.526	ng/ul	0.00
73) Anthracene-d10	17.298	188	567312m	12.784	ng/ul	0.00
81) Pyrene-d10	19.545	212	968483	17.271	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.516	264	219837	15.142	ng/ul	0.00

Target Compounds Qvalue

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

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