Quantitation Report (Qedit)

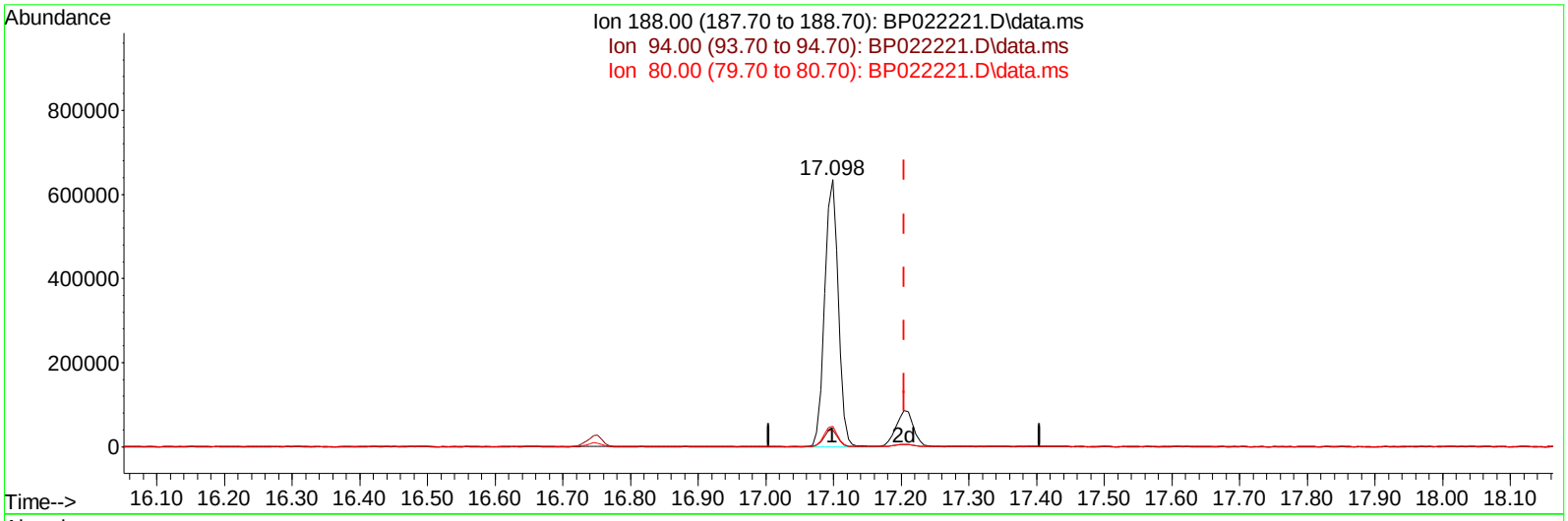
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022221.D
Acq On : 04 Oct 2024 09:25
Operator : RC/JU
Sample : P4018-03MEDL 10X
Misc :
ALS Vial : 27 Sample Multiplier: 1

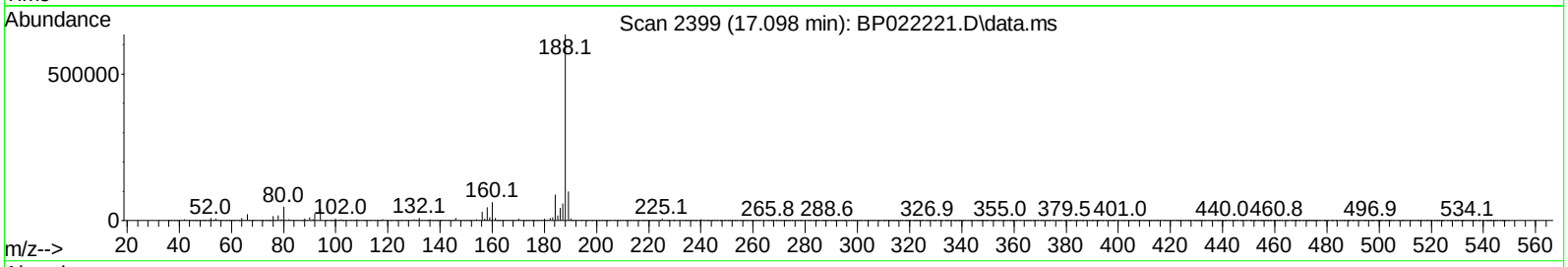
Quant Time: Oct 04 12:47:43 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 04 05:25:28 2024
Response via : Initial Calibration

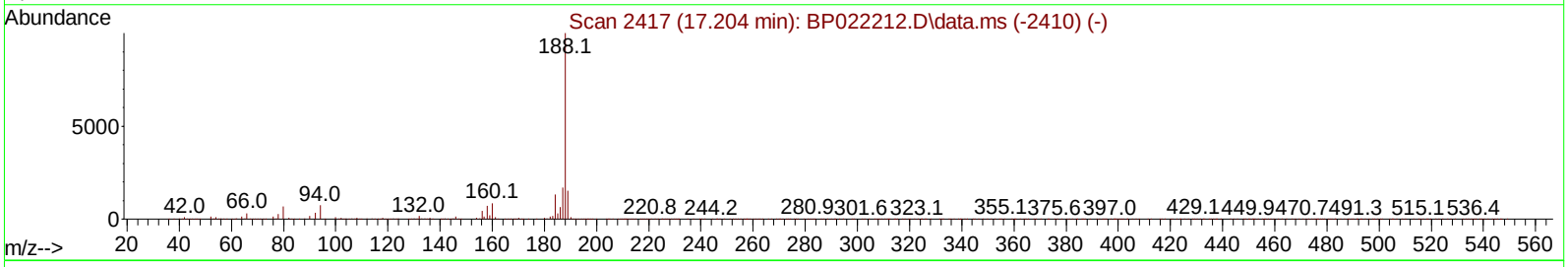
Instrument :
BNA_P
ClientSampleId :
DDCA1MEDL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/04/2024
Supervised By :mohammad ahmed 10/05/2024







TIC: BP022221.D\data.ms

(73) Anthracene-d10 (S)

17.098min (-0.106) 21.42 ng/ul

response 894421

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	7.50	6.35
80.00	8.00	7.57
0.00	0.00	0.00

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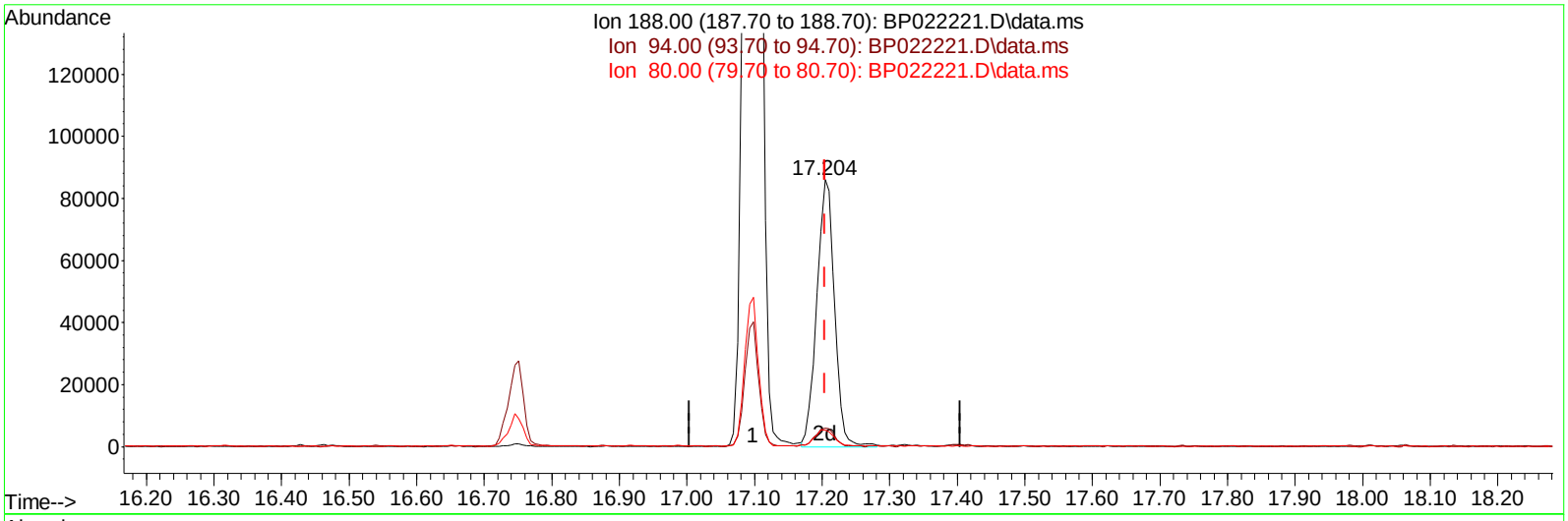

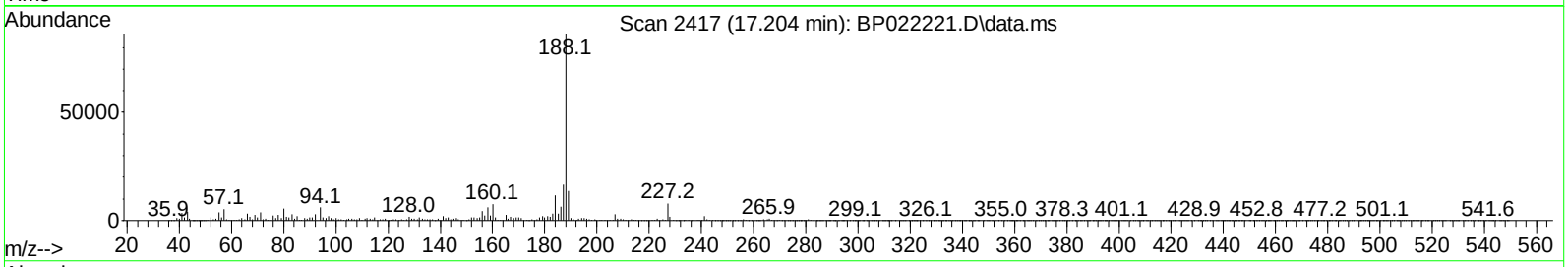

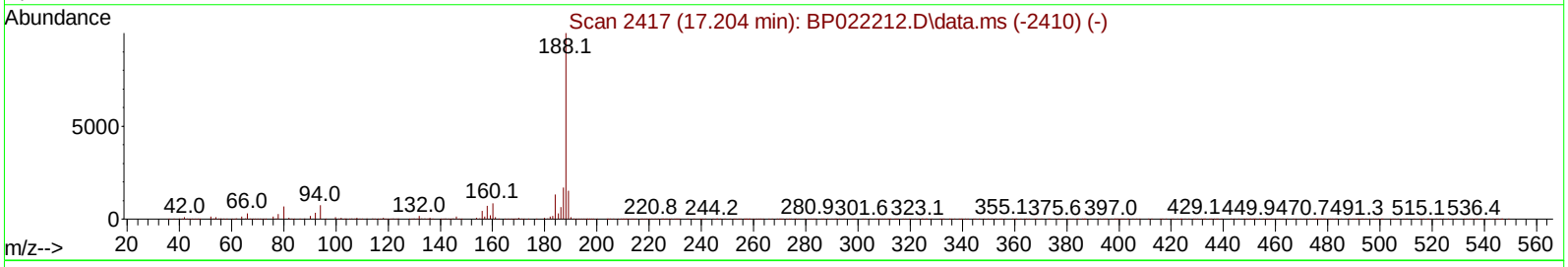

TIC: BP022221.D\data.ms

(73) Anthracene-d10 (S)

17.204min (+ 0.000) 3.74 ng/ul m

response 156367

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	7.50	7.04
80.00	8.00	6.29#
0.00	0.00	0.00

Quantitation Report (LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022221.D
Acq On : 04 Oct 2024 09:25
Operator : RC/JU
Sample : P4018-03MEDL 10X
Misc :
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Oct 05 02:23:25 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 04 05:25:28 2024
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
DDCA1MEDL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/04/2024
Supervised By :mohammad ahmed 10/05/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	127233	20.000	ng/ul	0.00
20) Naphthalene-d8	10.452	136	490477	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.298	164	377495	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	894421	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	994309	20.000	ng/ul	0.00
88) Perylene-d12	24.810	264	1174785	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	2057	0.633	ng/uL	0.00
4) Pyridine-d5	3.640	84	22825	2.458	ng/ul	0.01
7) Phenol-d5	6.881	99	30294	2.911	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	20637	3.220	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	24696	3.280	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	23887	2.896	ng/ul	0.00
21) Nitrobenzene-d5	8.828	128	11141	3.035	ng/ul	0.00
24) 2-Nitrophenol-d4	9.546	143	12384	2.975	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	26078	2.885	ng/ul	0.00
31) 4-Chloroaniline-d4	10.587	131	32410	3.005	ng/ul	0.00
46) Dimethylphthalate-d6	13.704	166	98762	3.384	ng/ul	0.00
49) Acenaphthylene-d8	13.987	160	103603	3.353	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.310	176	84564	3.276	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.434	200	8553	1.540	ng/ul	0.00
73) Anthracene-d10	17.204	188	156367m	3.745	ng/ul	0.00
81) Pyrene-d10	19.575	212	184798	3.587	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.580	264	206188	3.337	ng/ul	0.00
Target Compounds						Qvalue
36) 2-Methylnaphthalene	12.110	142	45512	2.373	ng/ul	96
37) 1-Methylnaphthalene	12.322	142	27494	1.421	ng/ul#	95
58) 2,3,4,6-Tetrachlorophenol	14.945	232	192015	22.017	ng/ul	99
71) Pentachlorophenol	16.751	266	2838452	326.200	ng/ul	100

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

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