

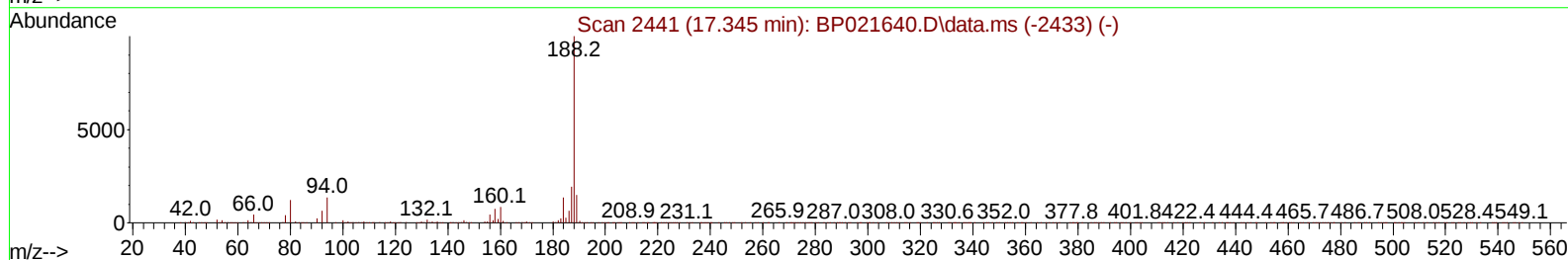
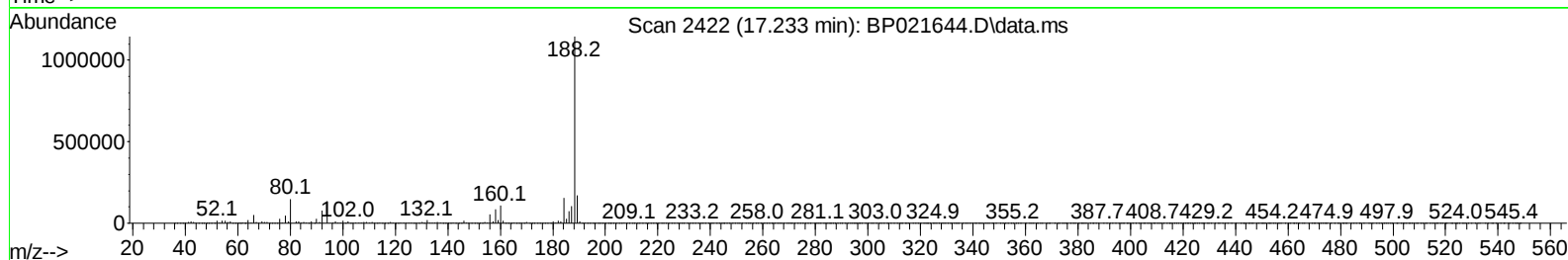
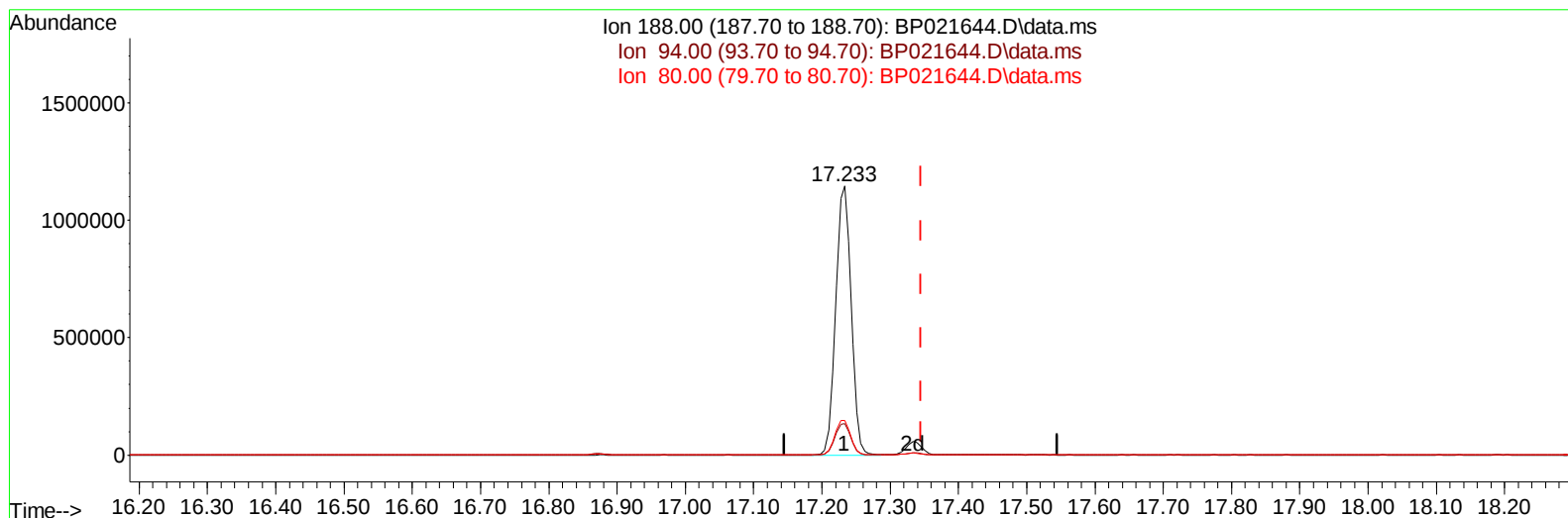
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021644.D
Acq On : 26 Aug 2024 18:48
Operator : RC/JU
Sample : P3695-09DL 10X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCNA4DL

Manual IntegrationsAPPROVED

Quant Time: Aug 26 23:21:35 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Aug 23 17:51:55 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/27/2024
Supervised By :mohammad ahmed 08/29/2024



TIC: BP021644.D\data.ms

(73) Anthracene-d10 (S)

17.233min (-0.112) 21.07 ng/u1

response 1792630

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	13.40	11.75
80.00	12.40	12.99
0.00	0.00	0.00

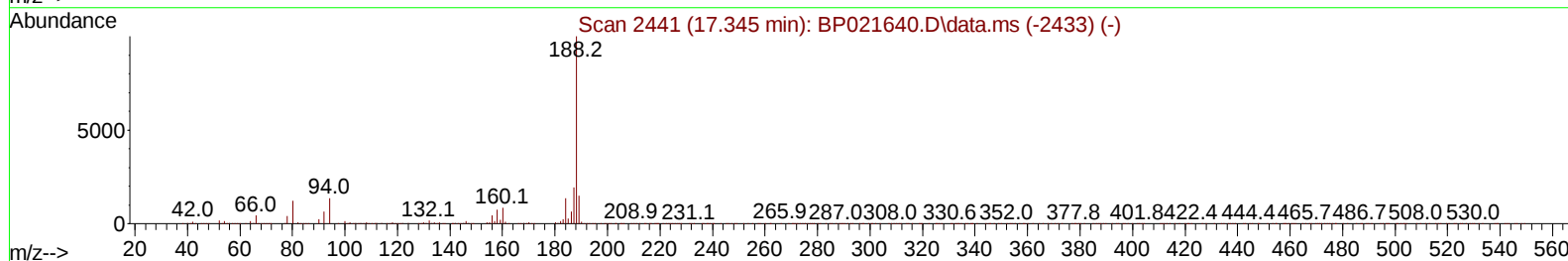
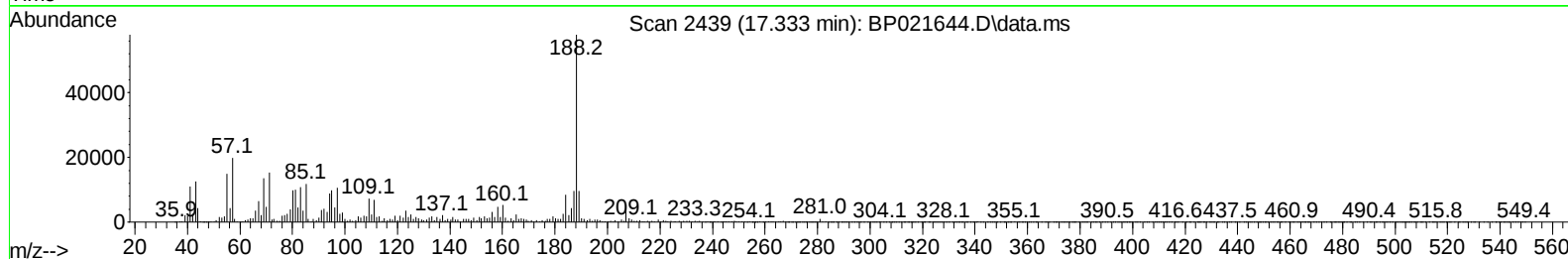
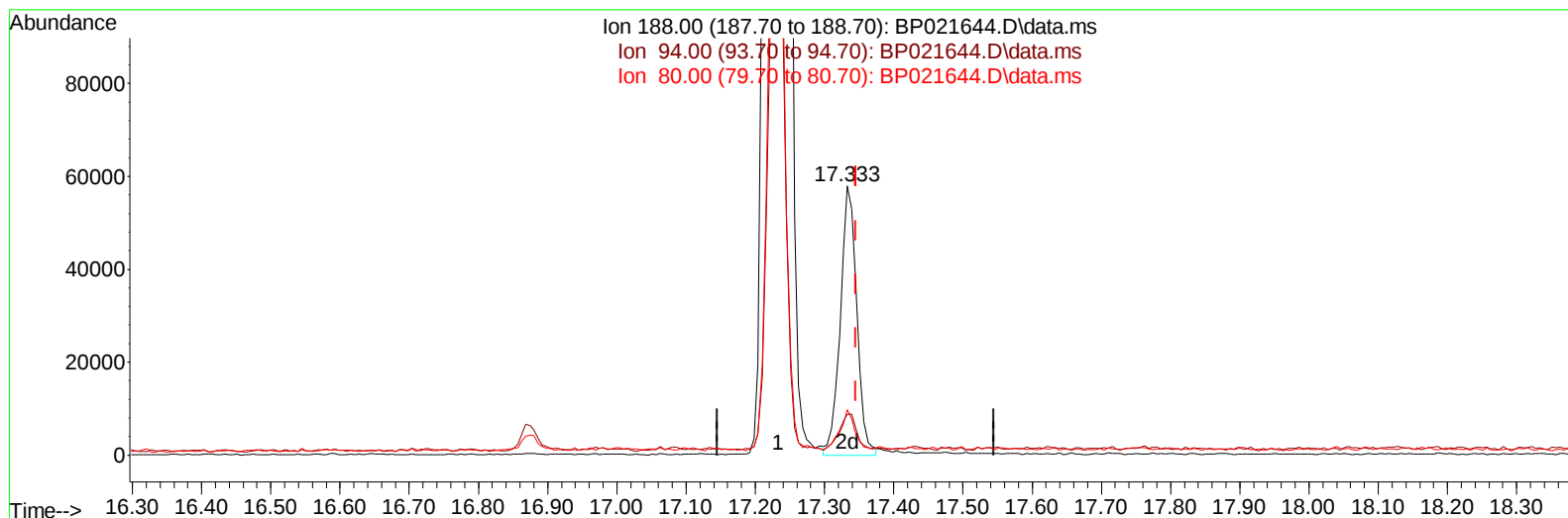
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TIC: BP021644.D\data.ms

(73) Anthracene-d10 (S)

17.333min (-0.012) 1.12 ng/ul m

response 95345

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	13.40	15.30
80.00	12.40	16.86#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/27/2024
 Supervised By :mohammad ahmed 08/29/2024

Quant Time: Aug 26 23:24:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 23 17:51:55 2024
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.798	152		285657	20.000	ng/ul	0.00
20) Naphthalene-d8	10.587	136		1212286	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164		826028	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188		1792630	20.000	ng/ul	0.00
79) Chrysene-d12	21.686	240		1794838	20.000	ng/ul	-0.03
88) Perylene-d12	25.092	264		2014370	20.000	ng/ul	-0.04
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.281	96		2266	0.263	ng/uL	0.00
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	6.963	99		24645	0.808	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67		15665	0.944	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132		16845	0.856	ng/ul	0.00
15) 4-Methylphenol-d8	8.493	113		16100	0.683	ng/ul	0.00
21) Nitrobenzene-d5	8.945	128		7604	0.780	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143		6234	0.630	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165		11388	0.584	ng/ul	0.00
31) 4-Chloroaniline-d4	10.716	131		16034	0.565	ng/ul	0.00
46) Dimethylphthalate-d6	13.839	166		60873	0.963	ng/ul	0.00
49) Acenaphthylene-d8	14.128	160		62263	0.877	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143		0d	0.000	ng/ul	
60) Fluorene-d10	15.445	176		58357	1.098	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200		0d	0.000	ng/ul	
73) Anthracene-d10	17.333	188		95345m	1.121	ng/ul	-0.01
81) Pyrene-d10	19.710	212		97332	0.949	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.856	264		68719	0.640	ng/ul	-0.05
Target Compounds							
						Qvalue	
71) Pentachlorophenol	16.874	266		418611	27.841	ng/ul	99

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

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