

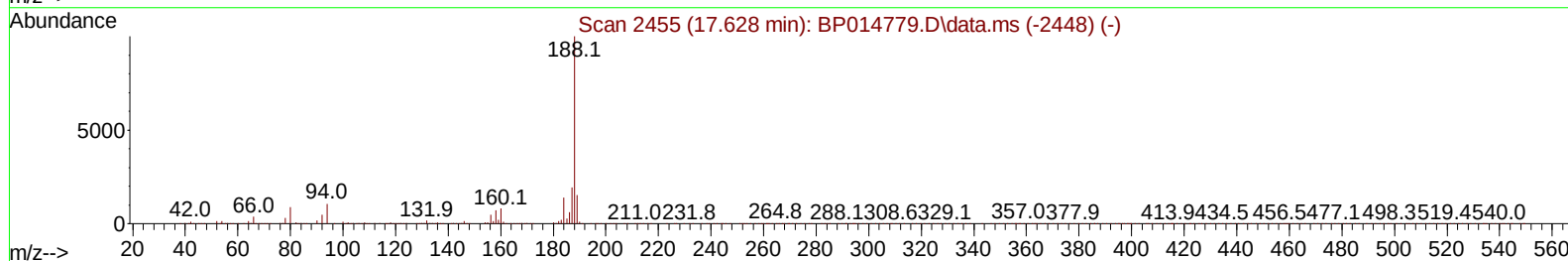
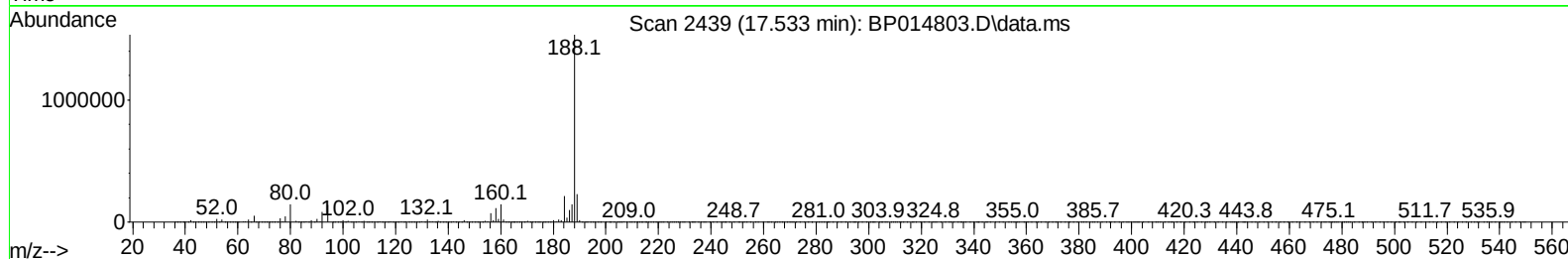
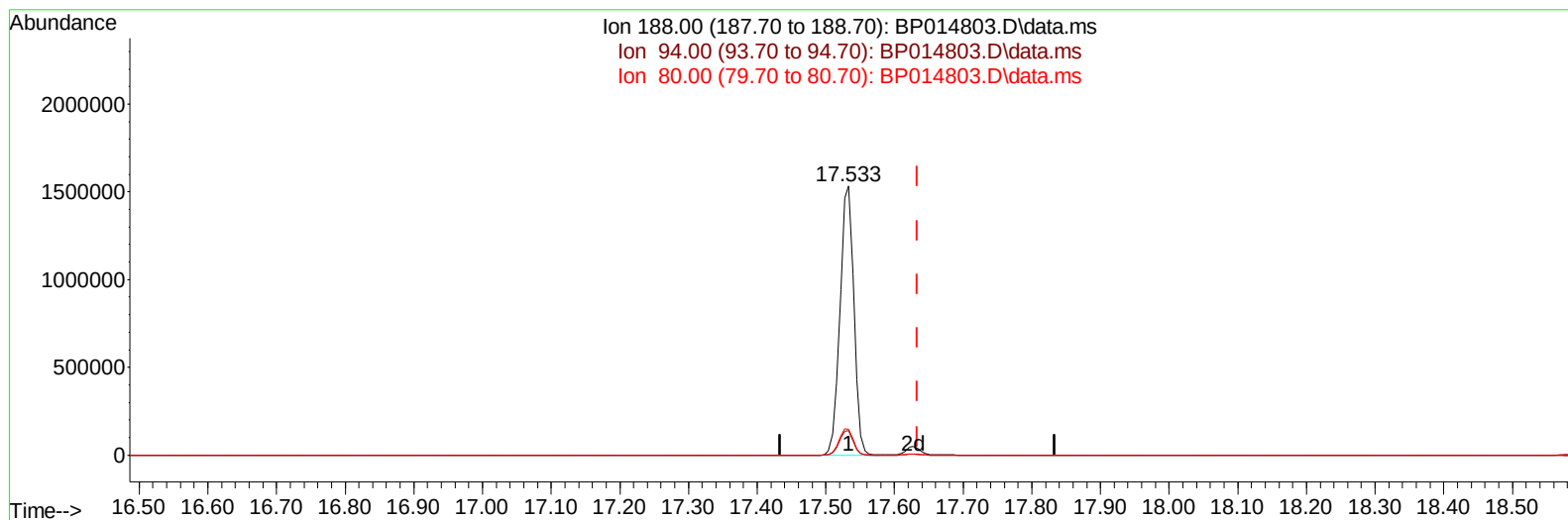
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
Data File : BP014803.D
Acq On : 21 Apr 2023 22:30
Operator : CG/JU
Sample : 02296-19DL2 50X
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCBM8DL2

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/24/2023
Supervised By :Jagrut Upadhyay 04/24/2023

Quant Time: Apr 21 23:56:55 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 20 02:32:50 2023
Response via : Initial Calibration



TIC: BP014803.D\data.ms

(73) Anthracene-d10 (S)

17.533min (-0.100) 20.70 ng/u1

response 2154549

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	10.50	9.10
80.00	9.20	9.58
0.00	0.00	0.00

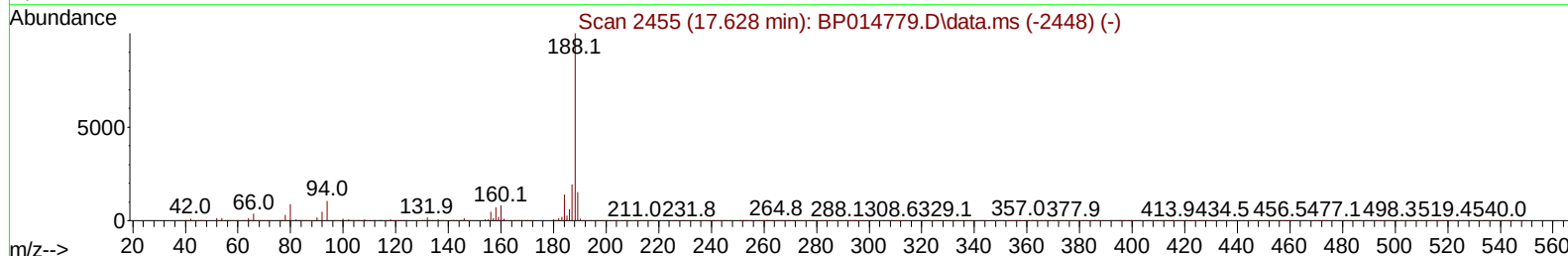
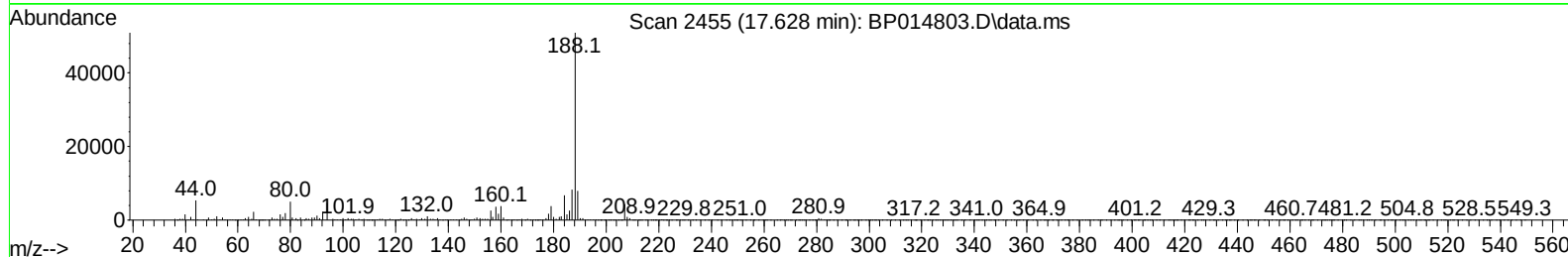
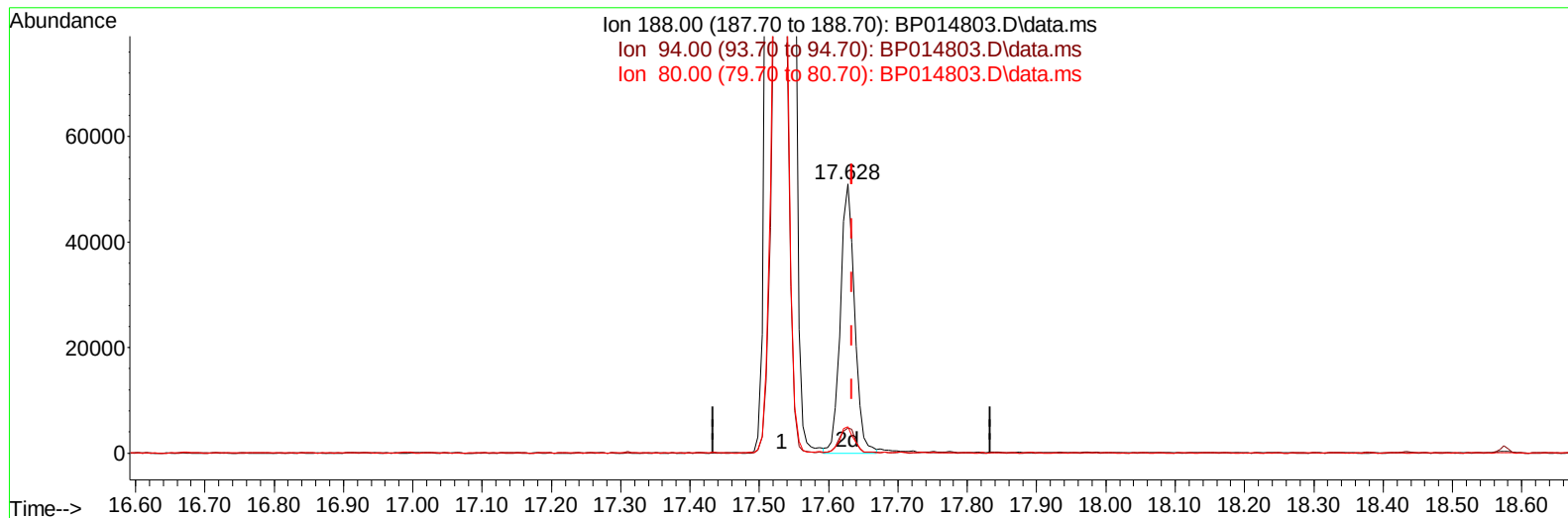
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(73) Anthracene-d10 (S)

17.628min (-0.006) 0.69 ng/ul m

response 72018

Ion	Exp%	Act%
188.00	100.00	100.00
94.00	10.50	9.30
80.00	9.20	9.91
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.152	152	379839	20.000	ng/ul	0.00
20) Naphthalene-d8	10.981	136	1605176	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.781	164	1028980	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.533	188	2154549	20.000	ng/ul	0.00
79) Chrysene-d12	21.598	240	1434326	20.000	ng/ul	0.00
88) Perylene-d12	24.174	264	1441106	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.287	99	4501	0.137	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.469	67	14281	0.705	ng/ul	0.00
11) 2-Chlorophenol-d4	7.675	132	10688	0.438	ng/ul	0.00
15) 4-Methylphenol-d8	8.851	113	7652	0.299	ng/ul	0.00
21) Nitrobenzene-d5	9.322	128	5845	0.581	ng/ul	0.00
24) 2-Nitrophenol-d4	10.046	143	4926	0.542	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.587	165	11680	0.474	ng/ul	0.00
31) 4-Chloroaniline-d4	11.110	131	6376	0.169	ng/ul	0.00
46) Dimethylphthalate-d6	14.181	166	54507	0.669	ng/ul	0.00
49) Acenaphthylene-d8	14.475	160	50288	0.533	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.769	176	46133	0.660	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.628	188	72018m	0.692	ng/ul	0.00
81) Pyrene-d10	19.839	212	64985	0.791	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.004	264	40624	0.544	ng/ul	-0.01
Target Compounds						
					Qvalue	
30) Naphthalene	11.034	128	4753959	52.845	ng/ul	100
36) 2-Methylnaphthalene	12.622	142	296347	4.934	ng/ul	99
37) 1-Methylnaphthalene	12.840	142	201580	3.386	ng/ul	100
52) Acenaphthene	14.845	153	253995	3.640	ng/ul	99
56) Dibenzofuran	15.175	168	248284	2.528	ng/ul	99
61) Fluorene	15.822	166	82681	1.031	ng/ul	99
72) Phenanthrene	17.575	178	177474	1.462	ng/ul	99
77) Carbazole	17.922	167	109817	1.016	ng/ul	99

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

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