

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018256.D
 Acq On : 18 Nov 2023 12:07
 Operator : MA/JU
 Sample : 05370-08
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

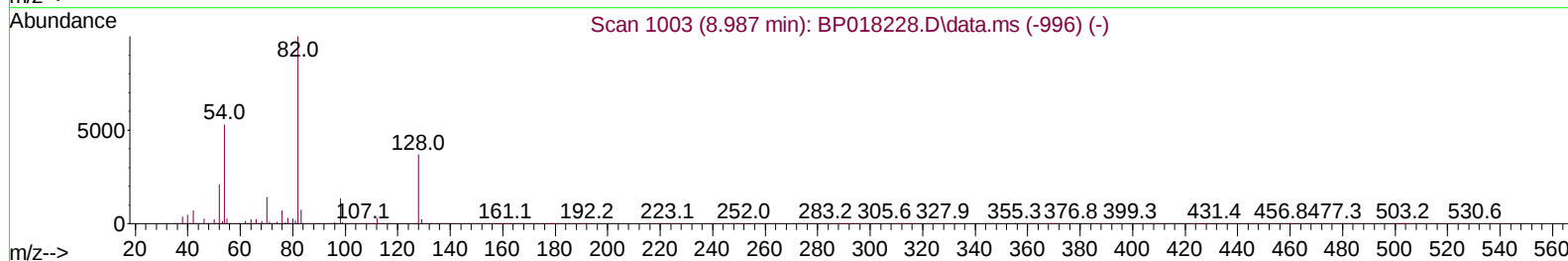
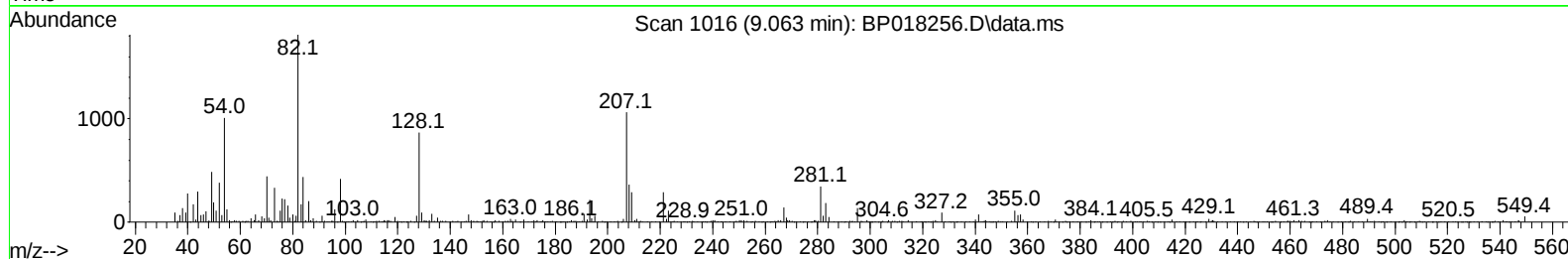
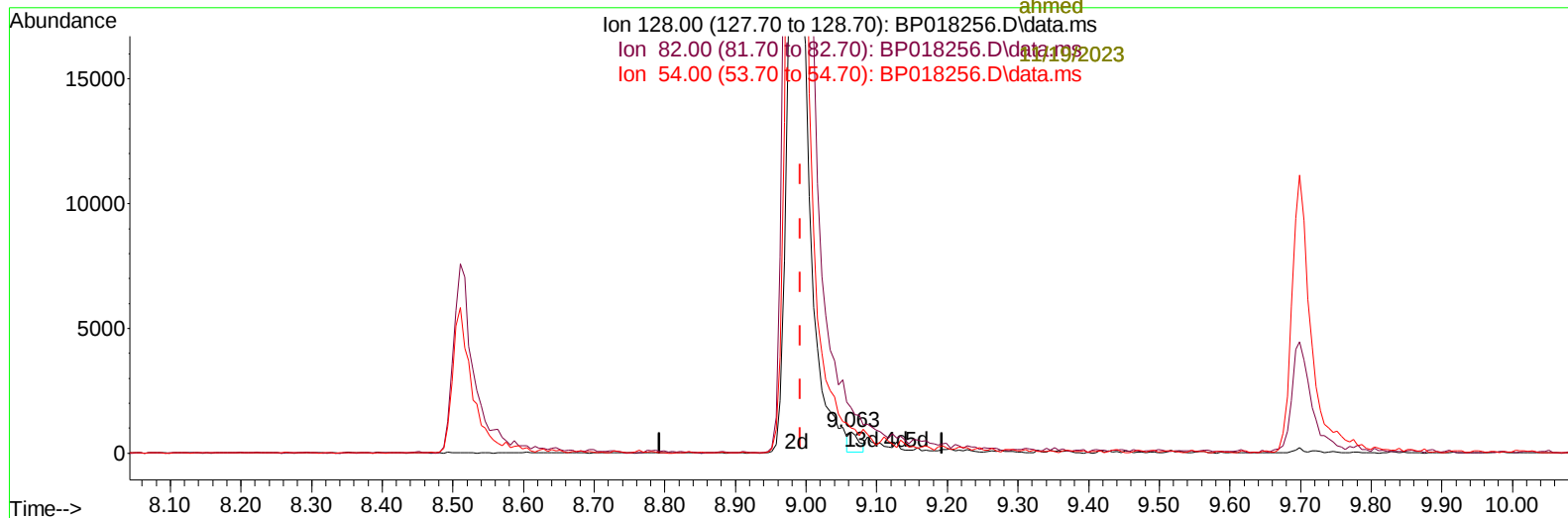
Instrument :
 BNA_P
 ClientSampleId :
 GCDQ4

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/19/2023
 Supervised By :mohammad ahmed 11/19/2023

11/19/2023
 Supervised By :mohammad
 ahmed

Quant Time: Nov 19 00:51:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration



TIC: BP018256.D\data.ms

(21) Nitrobenzene-d5 (S)

9.063min (+ 0.071) 0.45 ng/u1

response 774

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	256.70	209.58
54.00	125.00	116.28
0.00	0.00	0.00

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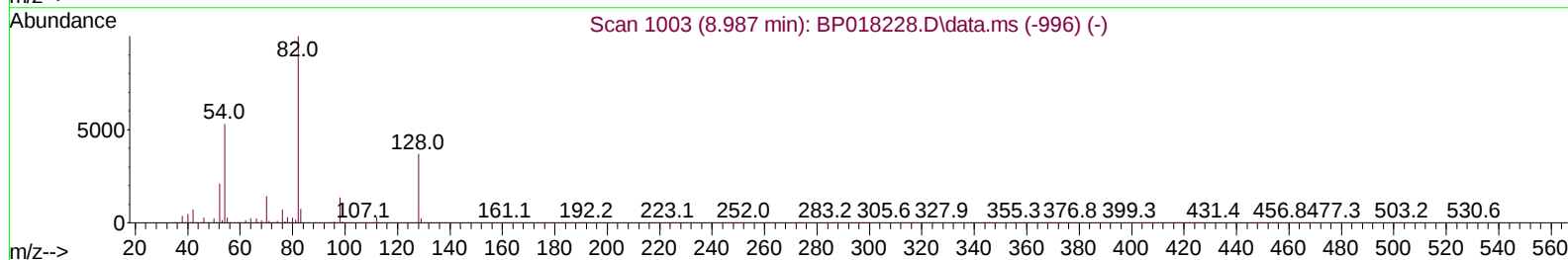
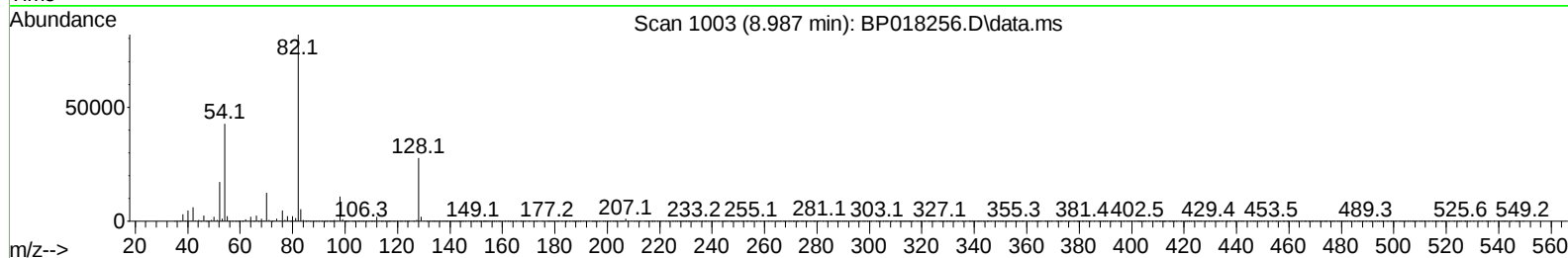
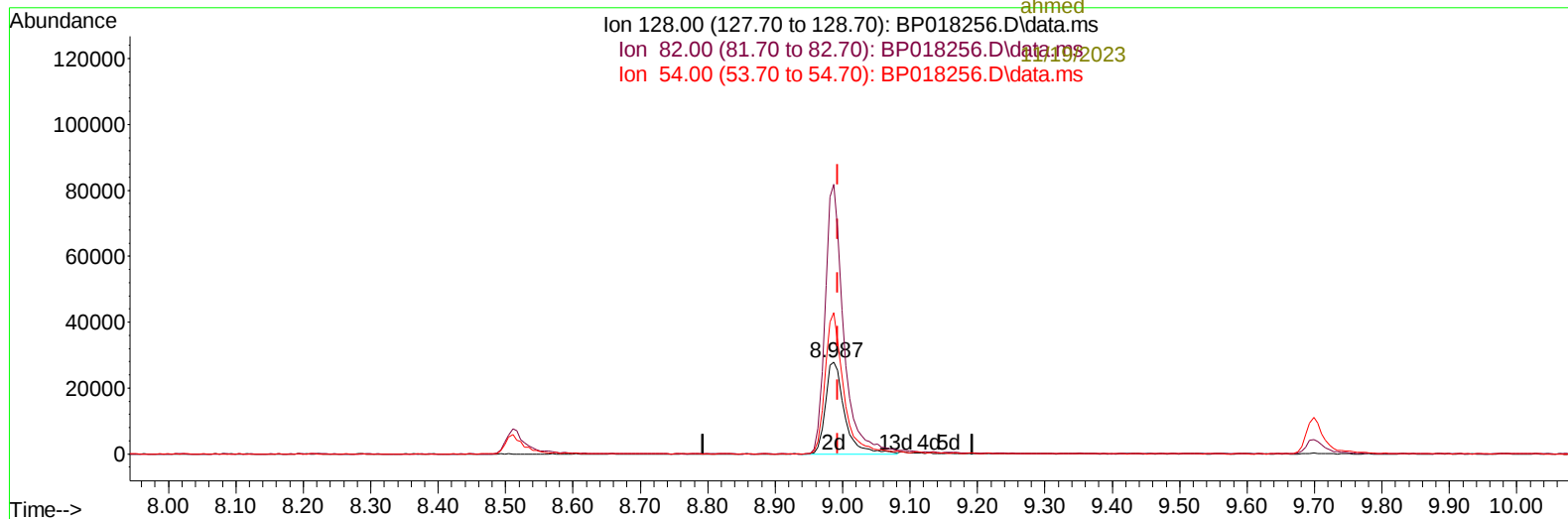
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Ion 128.00 (127.70 to 128.70): BP018256.D\data.ms
 Ion 82.00 (81.70 to 82.70): BP018256.D\data.ms
 Ion 54.00 (53.70 to 54.70): BP018256.D\data.ms



TIC: BP018256.D\data.ms

(21) Nitrobenzene-d5 (S)

8.987min (-0.006) 32.04 ng/u1 m

response 55184

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	256.70	294.13
54.00	125.00	154.06#
0.00	0.00	0.00

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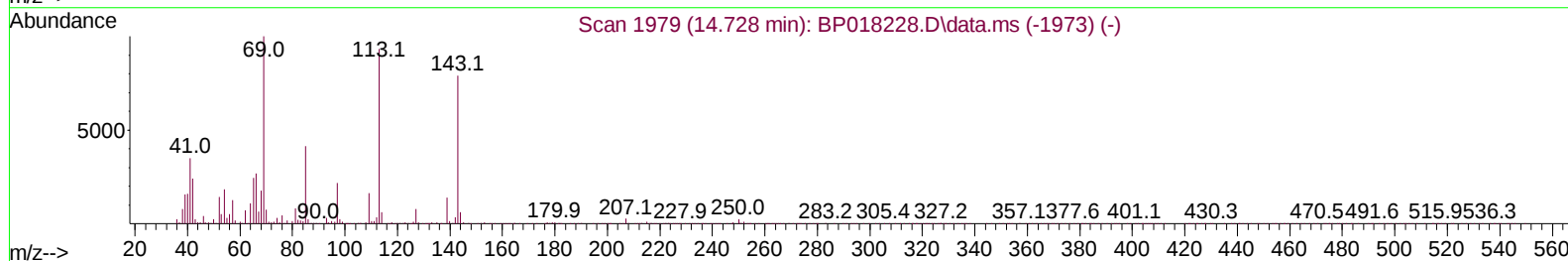
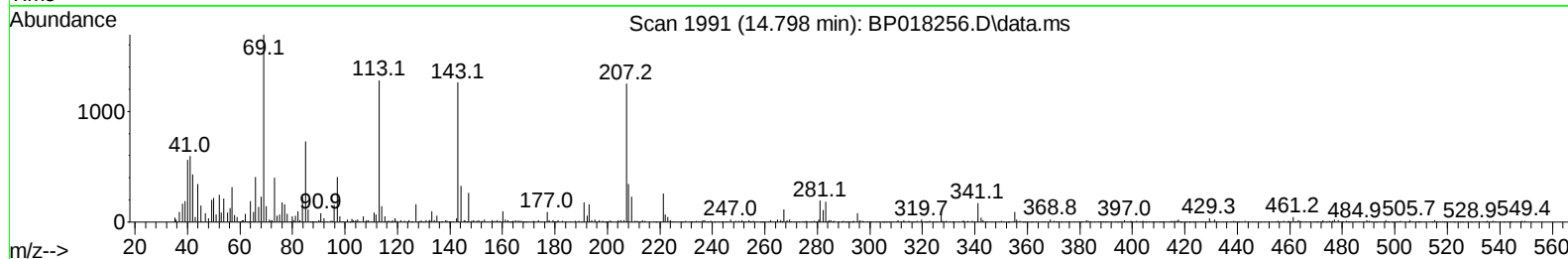
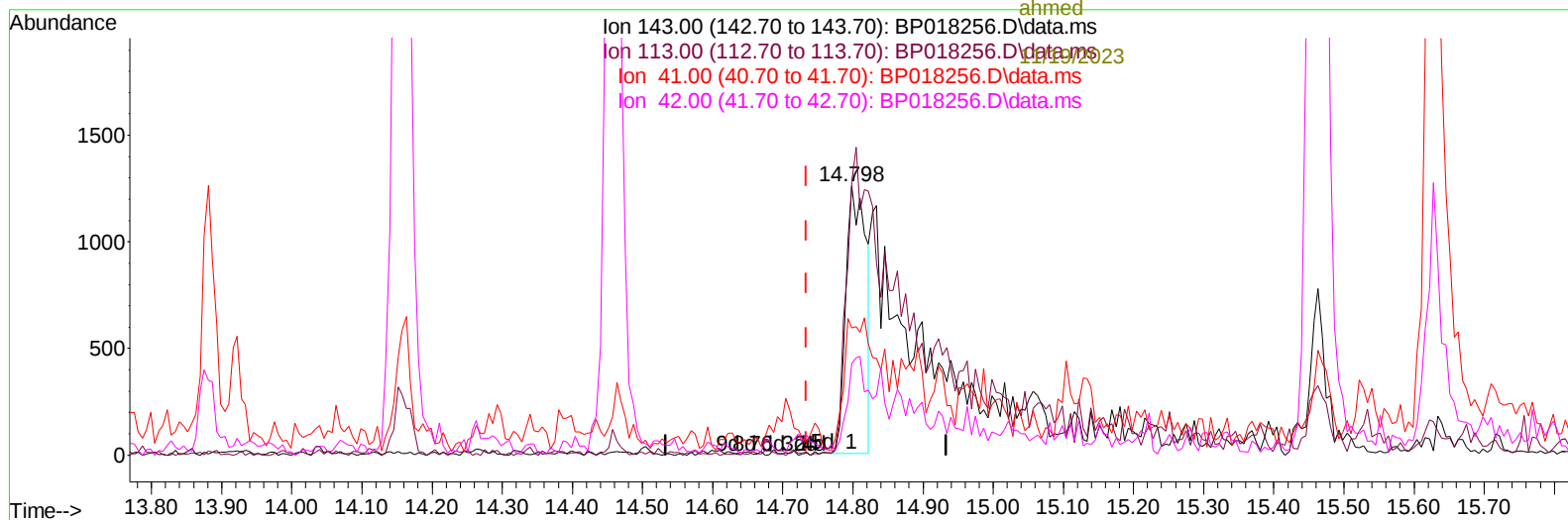
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(54) 4-Nitrophenol-d4 (S)**14.798min (+ 0.065) 1.52 ng/u1****response 2520**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	101.67
41.00	47.70	47.26
42.00	30.10	34.02

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Quant Time: Nov 19 02:26:50 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/19/2023						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	49002	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	191200	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	120420	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.240	188	290862	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	285099	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	323585	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	6266	4.511	ng/uL	0.00
4) Pyridine-d5	3.605	84	28738	7.935	ng/ul	0.01
7) Phenol-d5	6.964	99	30055	6.353	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.128	67	92577	33.299	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	95614	25.522	ng/ul	0.00
15) 4-Methylphenol-d8	8.511	113	58340	15.049	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	55184m	32.035	ng/ul	0.00
24) 2-Nitrophenol-d4	9.699	143	60537	31.020	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	94135	27.064	ng/ul	0.00
31) 4-Chloroaniline-d4	10.781	131	110954	23.261	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	390824	35.175	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	398890	32.961	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	7645m	4.605	ng/ul	0.06
60) Fluorene-d10	15.463	176	330231	35.999	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	56596	26.916	ng/ul	0.00
73) Anthracene-d10	17.345	188	626636	39.835	ng/ul	0.00
81) Pyrene-d10	19.604	212	786871	41.320	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	730361	38.817	ng/ul	0.00

Target Compounds Qvalue

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

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