

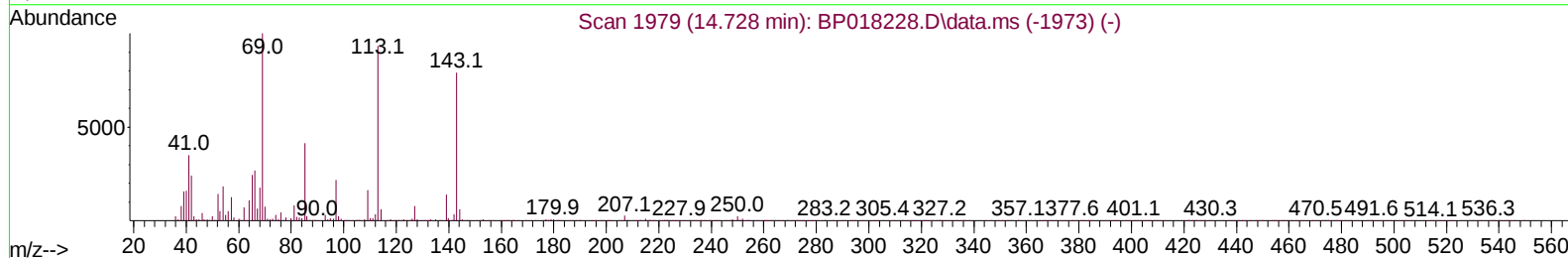
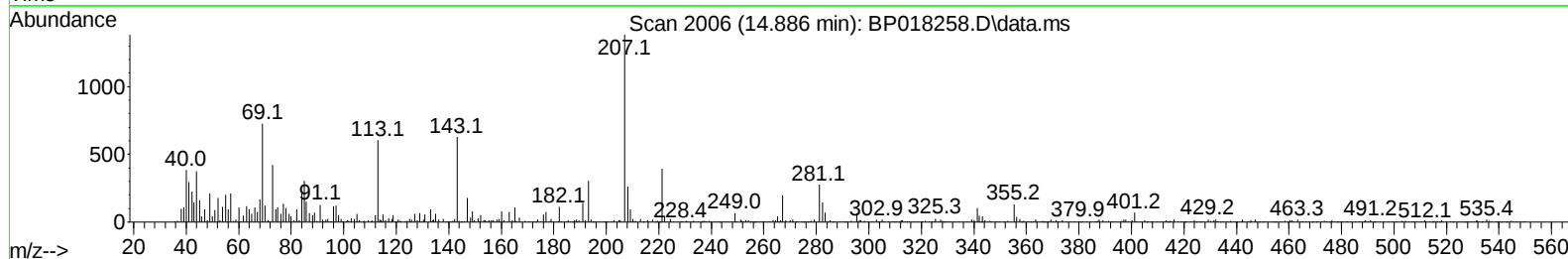
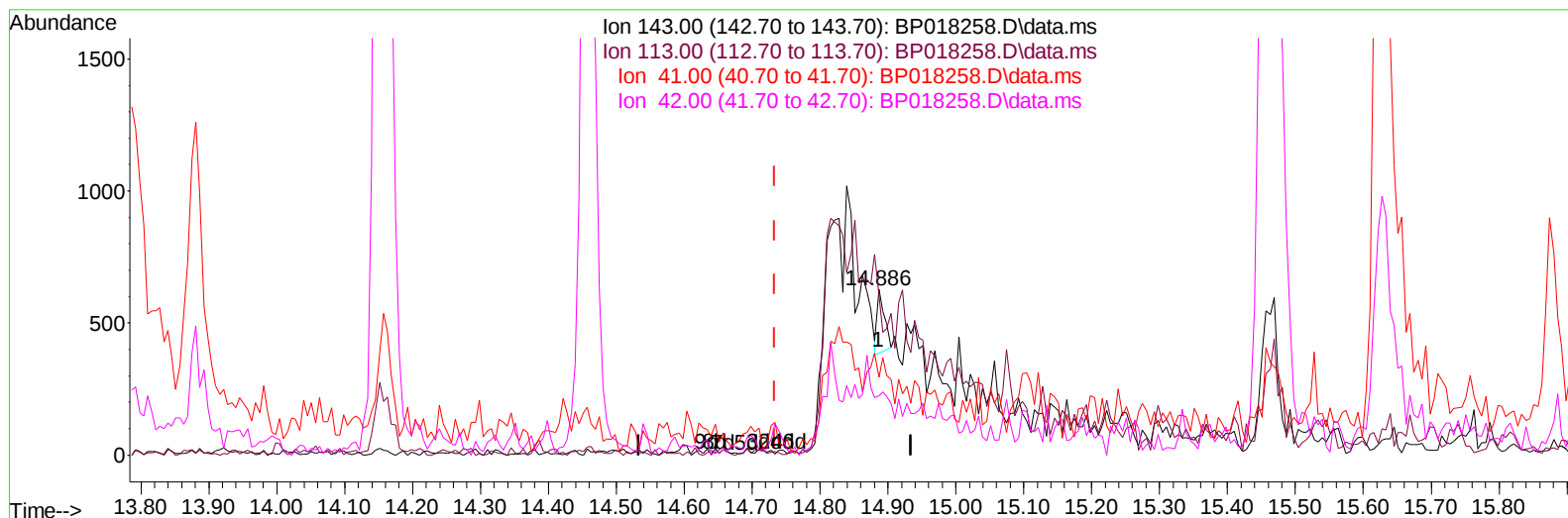
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018258.D
 Acq On : 18 Nov 2023 13:15
 Operator : MA/JU
 Sample : 05370-13
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDQ7

Manual IntegrationsAPPROVED

Quant Time: Nov 19 00:52:28 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

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



TIC: BP018258.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.886min (+ 0.153) 0.13 ng/u1

response 178

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	96.17
41.00	47.70	46.89
42.00	30.10	35.73

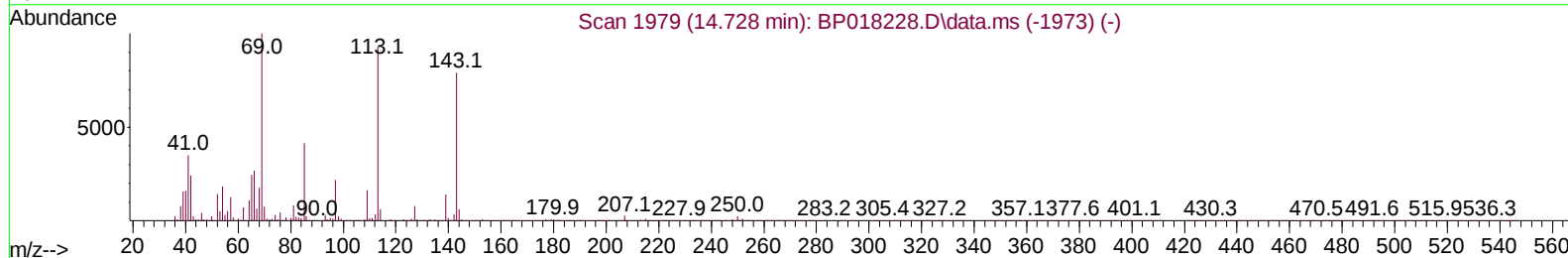
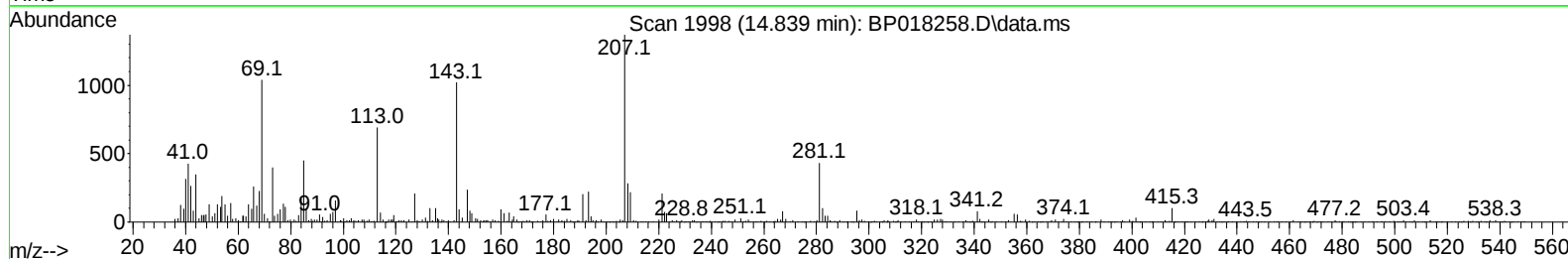
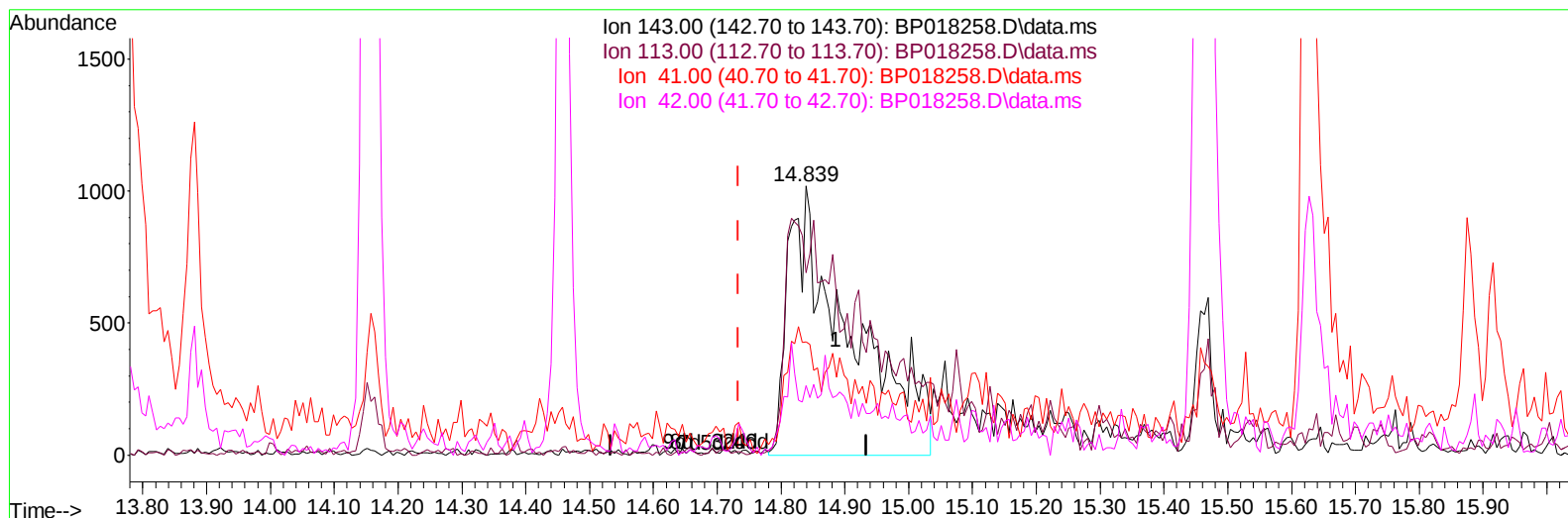
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(54) 4-Nitrophenol-d4 (S)

14.839min (+ 0.106) 4.95 ng/ul m

response 6885

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	67.71#
41.00	47.70	42.00
42.00	30.10	25.91

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Quant Time: Nov 19 02:30:54 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	43433	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	168622	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	100831	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	242160	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	224189	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	250881	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	6610	5.369	ng/uL	0.00
4) Pyridine-d5	3.605	84	33951	10.576	ng/ul	0.01
7) Phenol-d5	6.963	99	29769	7.099	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.128	67	90064	36.549	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	96182	28.966	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	59357	17.275	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	55226	36.352	ng/ul	0.00
24) 2-Nitrophenol-d4	9.699	143	60764	35.305	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	96362	31.414	ng/ul	0.00
31) 4-Chloroaniline-d4	10.781	131	115803	27.528	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	399792	42.972	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	416875	41.140	ng/ul	0.00
54) 4-Nitrophenol-d4	14.839	143	6885m	4.952	ng/ul	0.11
60) Fluorene-d10	15.463	176	338077	44.014	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	52599	30.046	ng/ul	0.00
73) Anthracene-d10	17.339	188	573141	43.761	ng/ul	0.00
81) Pyrene-d10	19.598	212	698858	46.669	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	668879	45.851	ng/ul	0.00

Target Compounds Qvalue

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

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