

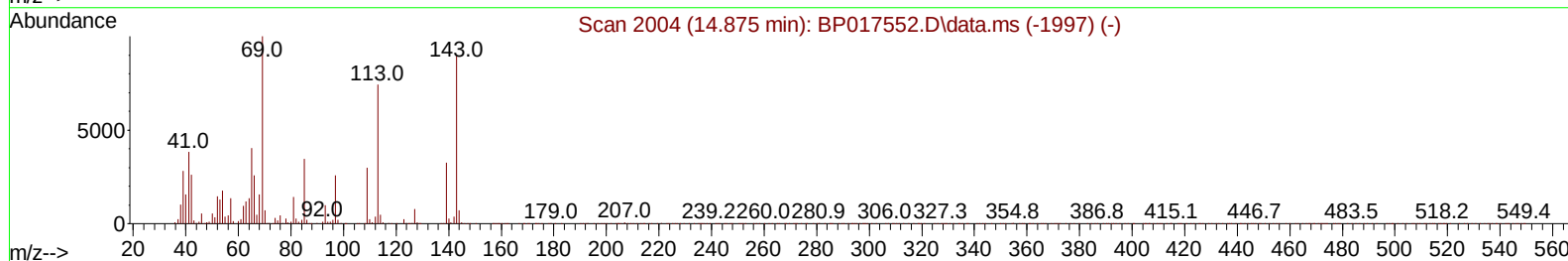
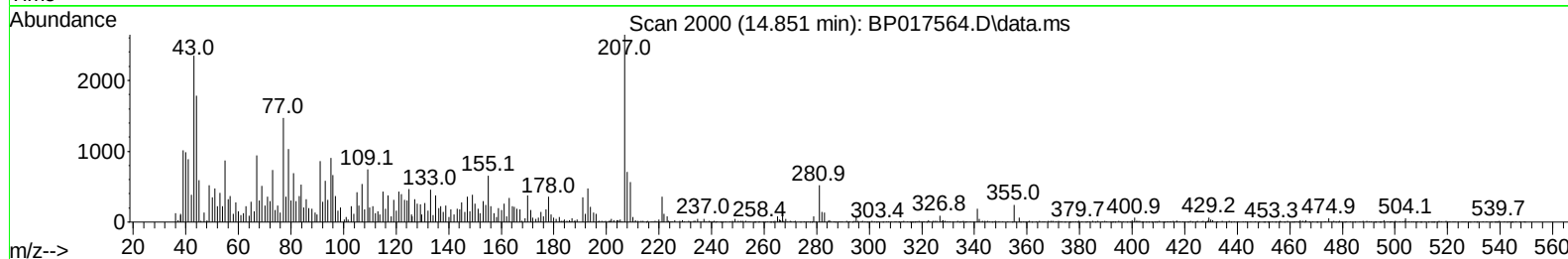
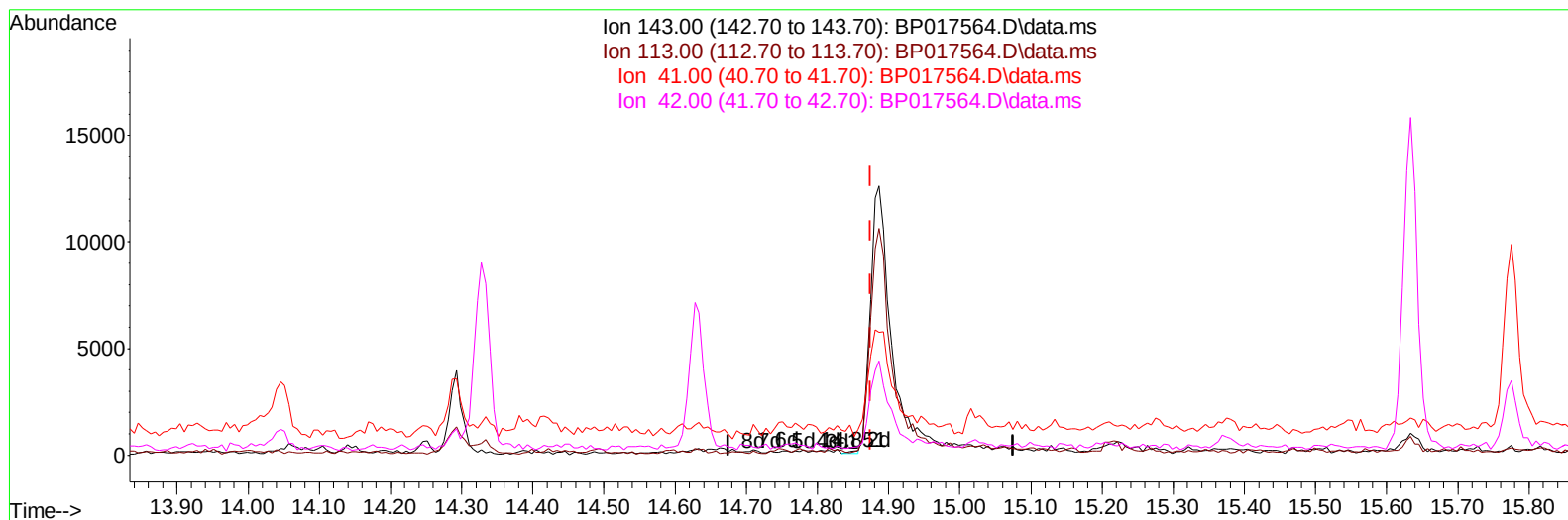
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017564.D
 Acq On : 05 Oct 2023 14:14
 Operator : MA/JU
 Sample : 04679-09
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PMW-3(2030927)

Manual IntegrationsAPPROVED

Quant Time: Oct 06 01:59:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 04 16:33:47 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/06/2023
 Supervised By :mohammad ahmed 10/07/2023



TIC: BP017564.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.851min (-0.023) 0.03 ng/ul

response 107

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.00	84.86
41.00	43.80	480.00#
42.00	29.00	207.57#

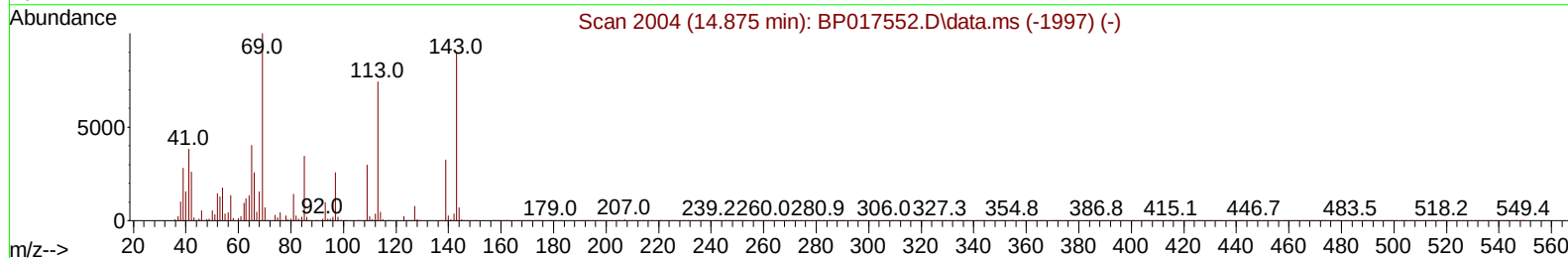
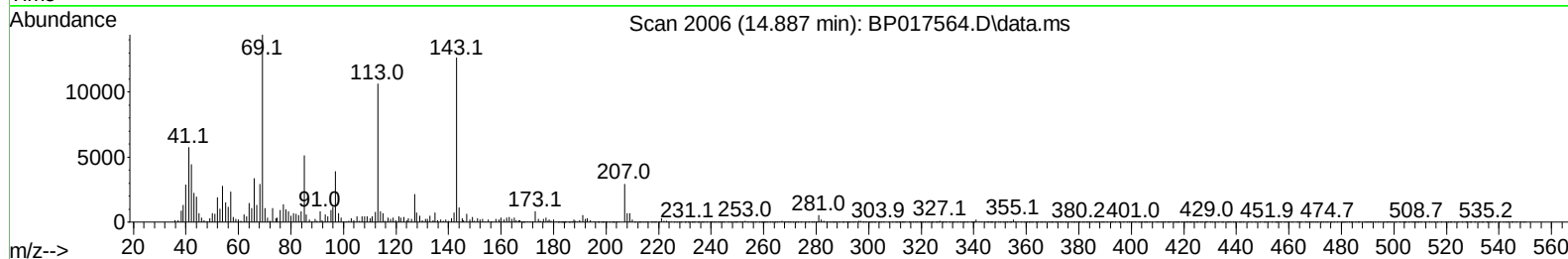
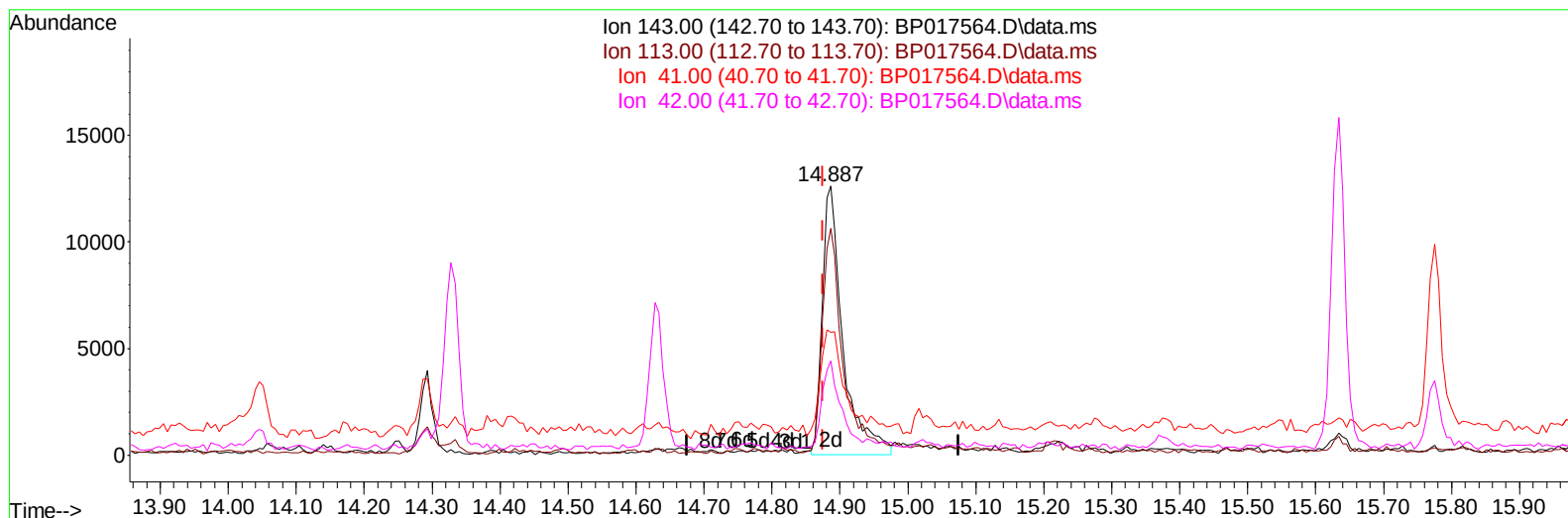
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TIC: BP017564.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.887min (+ 0.012) 7.34 ng/ul m

response 26538

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.00	84.22
41.00	43.80	45.54
42.00	29.00	35.03#

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Quant Time: Oct 06 02:36:35 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	96870	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	398153	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.634	164	264721	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.422	188	624094	20.000	ng/ul	0.00
79) Chrysene-d12	21.557	240	631082	20.000	ng/ul	0.00
88) Perylene-d12	24.133	264	684438	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	9963	4.112	ng/uL	0.00
4) Pyridine-d5	3.728	84	78214	12.332	ng/ul	0.00
7) Phenol-d5	7.099	99	74039	9.084	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	192583	38.723	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132	197696	30.806	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	151676	23.246	ng/ul	-0.01
21) Nitrobenzene-d5	9.146	128	121404	39.634	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	133423	38.620	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	239287	37.360	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	304570	33.381	ng/ul	0.00
46) Dimethylphthalate-d6	14.046	166	914491	44.287	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	918303	40.331	ng/ul	0.00
54) 4-Nitrophenol-d4	14.887	143	26538m	7.339	ng/ul	0.01
60) Fluorene-d10	15.634	176	763999	43.621	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	137077	35.231	ng/ul	0.00
73) Anthracene-d10	17.522	188	1328415	45.984	ng/ul	0.00
81) Pyrene-d10	19.775	212	1674932	47.763	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.963	264	1650552	47.785	ng/ul	-0.01

Target Compounds Qvalue

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

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