

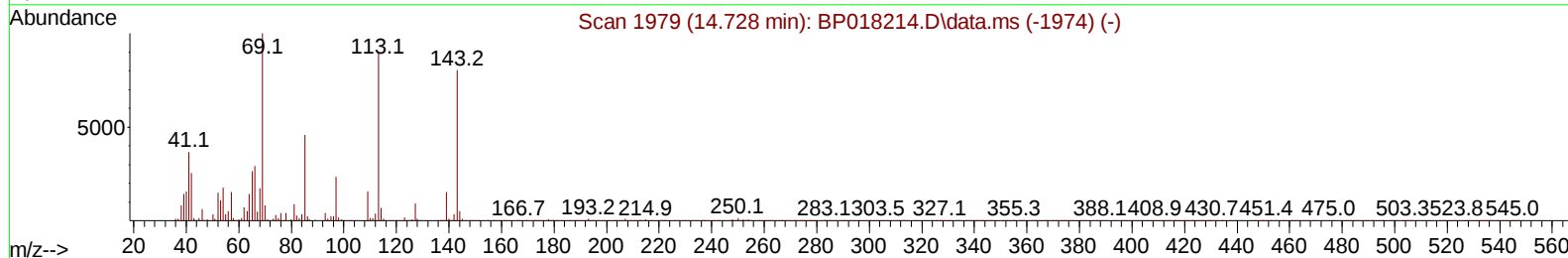
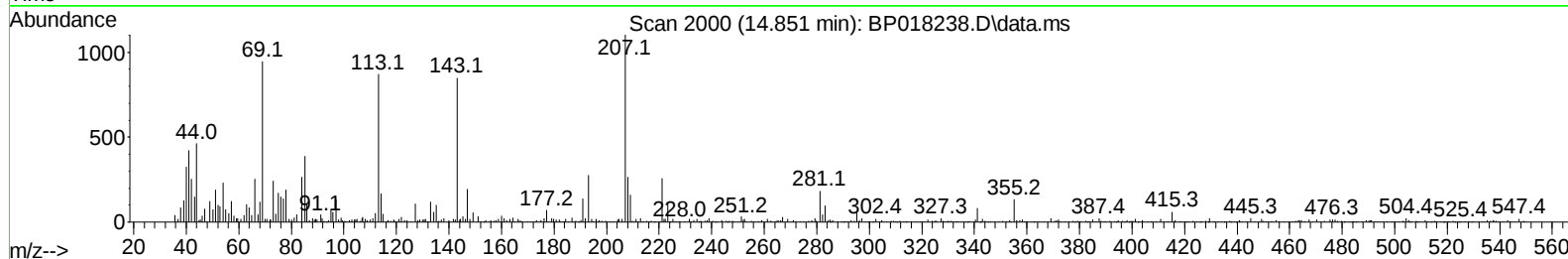
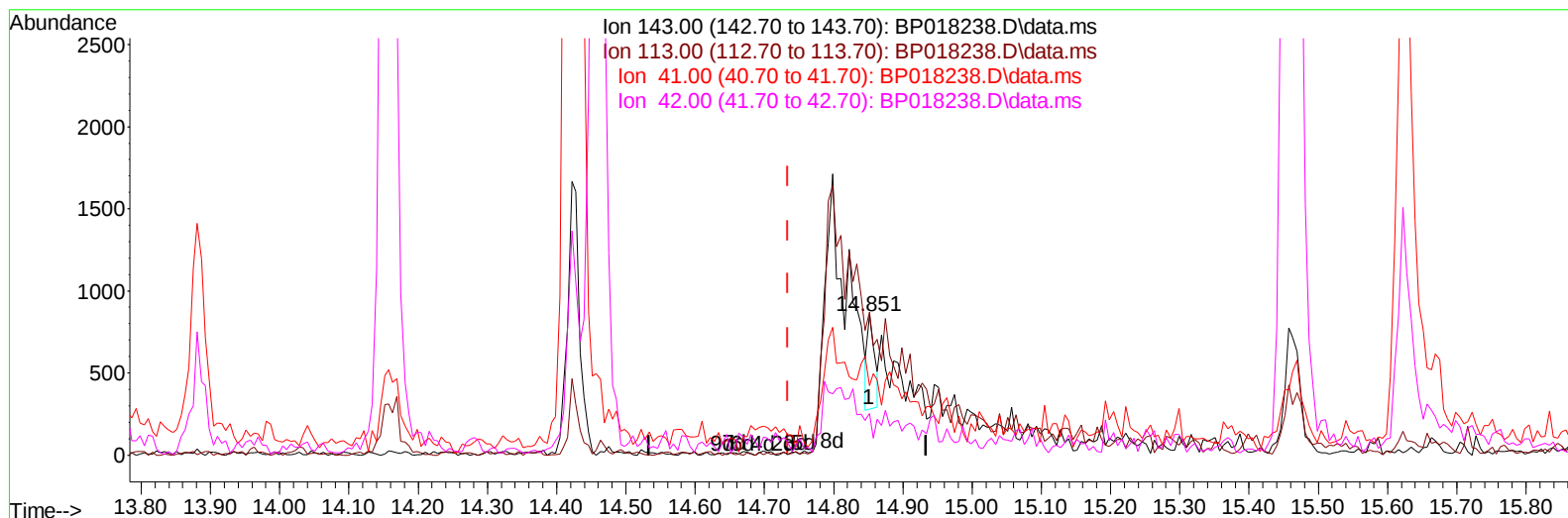
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018238.D
 Acq On : 18 Nov 2023 01:21
 Operator : MA/JU
 Sample : 05369-22
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDR7

Manual IntegrationsAPPROVED

Quant Time: Nov 18 01:51:17 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: BP018238.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.851min (+ 0.118) 0.20 ng/u1

response 413

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	102.47
41.00	47.70	49.76
42.00	30.10	29.88

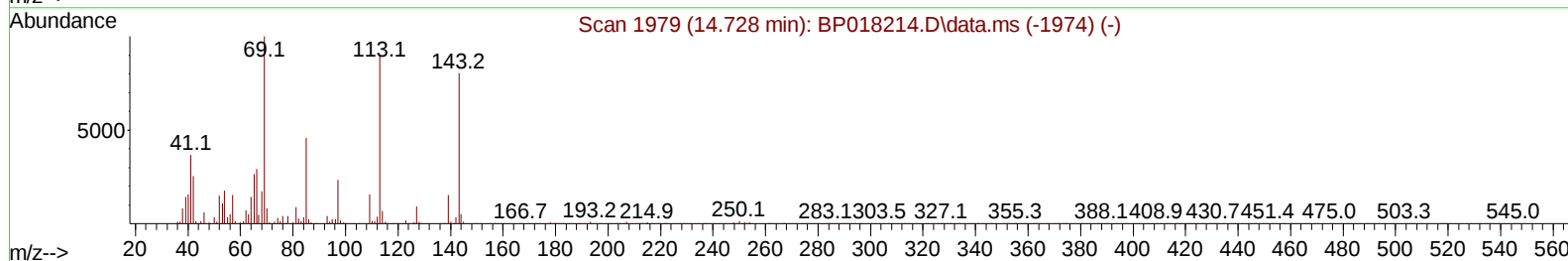
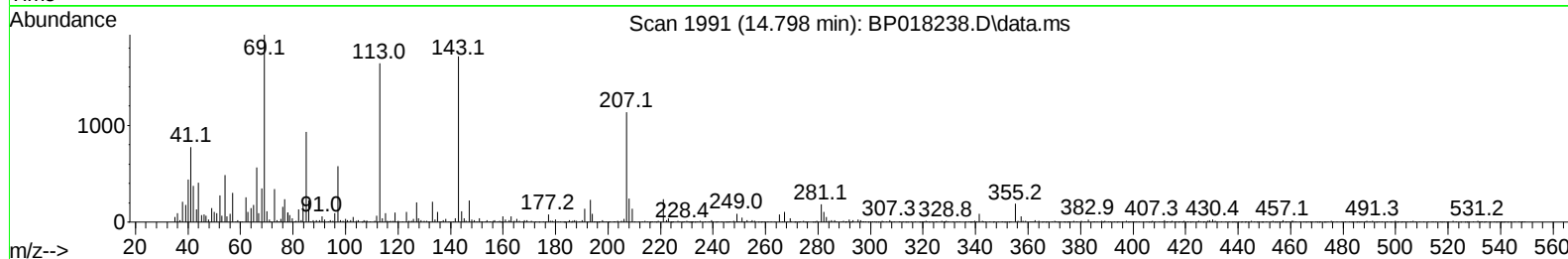
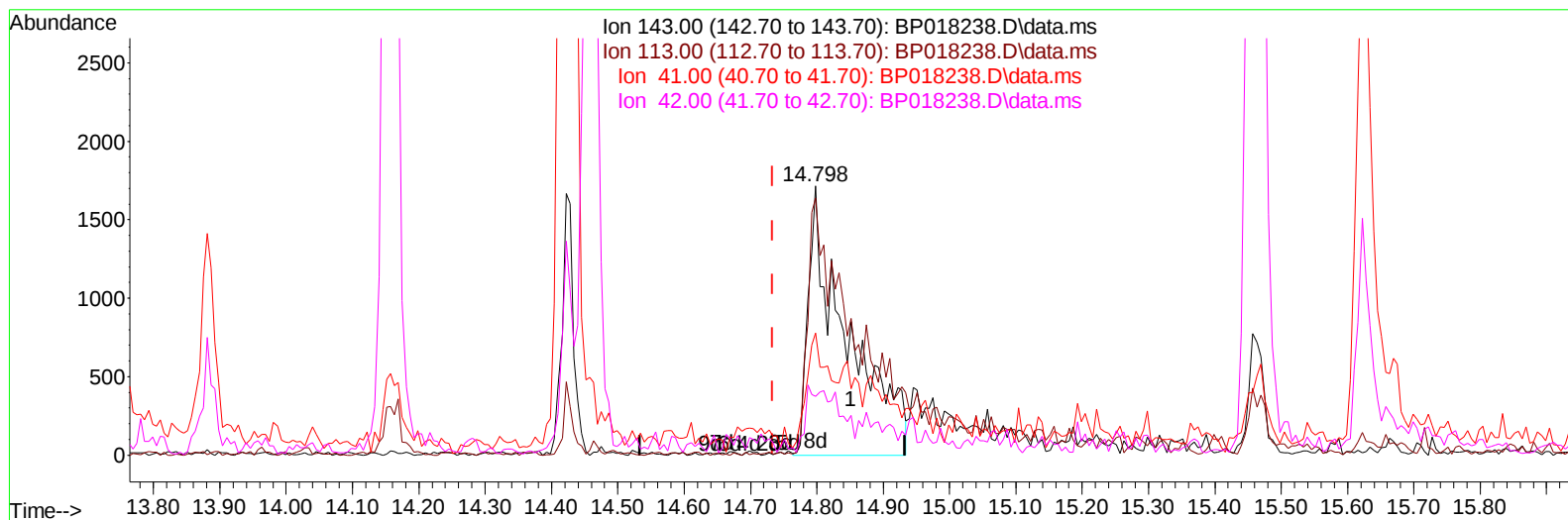
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(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.065) 3.30 ng/ul m

response 6874

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	95.57
41.00	47.70	45.33
42.00	30.10	21.76#

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Manual IntegrationsAPPROVED

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

Quant Time: Nov 18 01:51:52 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	54064	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	221422	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	150906	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.240	188	348506	20.000	ng/ul	0.00
79) Chrysene-d12	21.375	240	337187	20.000	ng/ul	0.00
88) Perylene-d12	23.845	264	378491	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	7637	4.983	ng/uL	0.00
4) Pyridine-d5	3.599	84	47949	11.999	ng/ul	0.00
7) Phenol-d5	6.964	99	30203	5.787	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.128	67	103188	33.641	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	96849	23.431	ng/ul	0.00
15) 4-Methylphenol-d8	8.511	113	60428	14.128	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	64991	32.579	ng/ul	0.00
24) 2-Nitrophenol-d4	9.699	143	67034	29.661	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	109612	27.212	ng/ul	0.00
31) 4-Chloroaniline-d4	10.775	131	148249	26.838	ng/ul	-0.01
46) Dimethylphthalate-d6	13.881	166	474934	34.110	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	475821	31.375	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	6874m	3.304	ng/ul	0.06
60) Fluorene-d10	15.463	176	424338	36.913	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	68526	27.199	ng/ul	0.00
73) Anthracene-d10	17.340	188	665128	35.288	ng/ul	0.00
81) Pyrene-d10	19.598	212	820505	36.430	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264	800846	36.389	ng/ul	0.00

Target Compounds Qvalue

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

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