

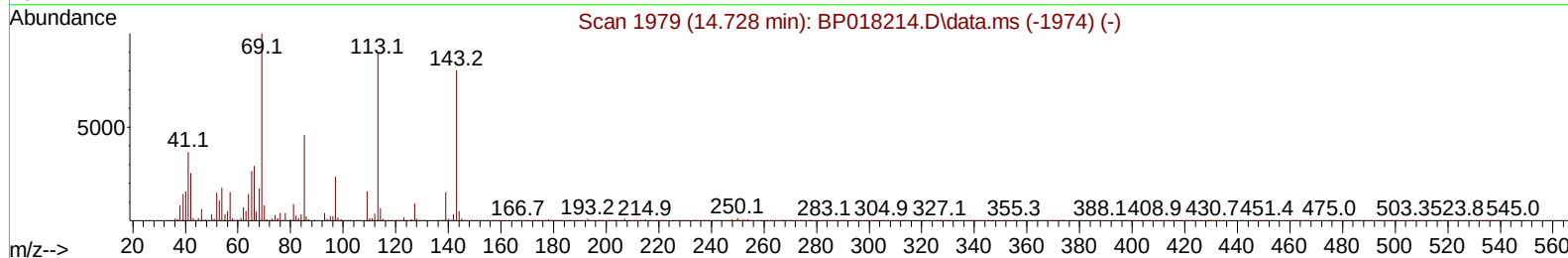
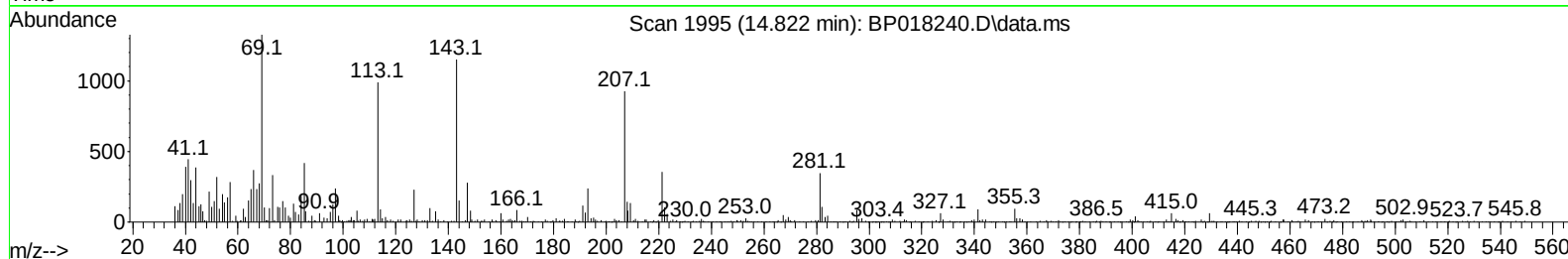
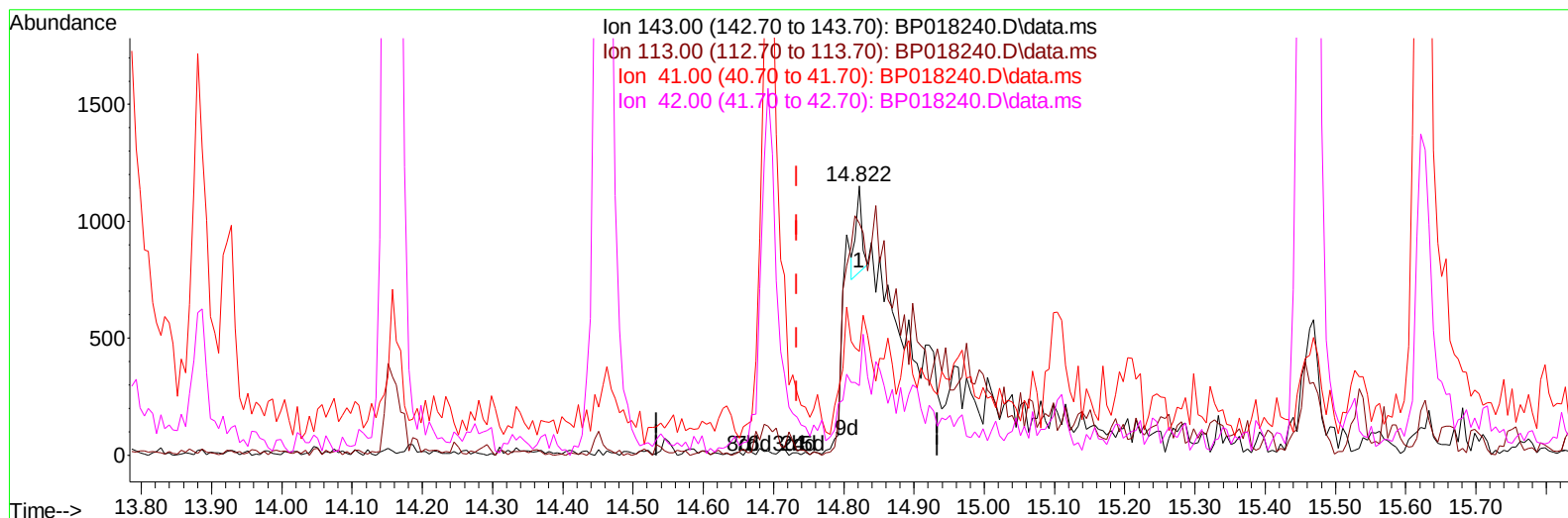
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018240.D
 Acq On : 18 Nov 2023 02:29
 Operator : MA/JU
 Sample : 05369-11
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDN8

Manual IntegrationsAPPROVED

Quant Time: Nov 18 02:57:38 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: BP018240.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (+ 0.088) 0.10 ng/ul

response 222

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	86.17
41.00	47.70	38.52
42.00	30.10	25.74

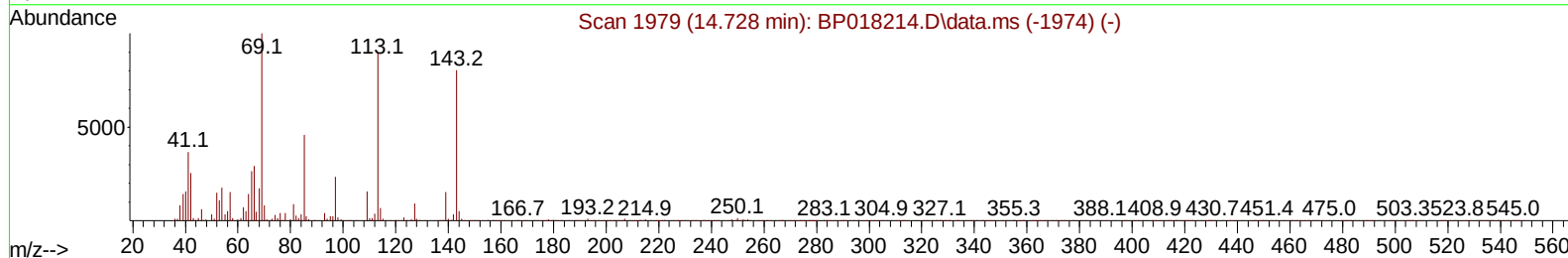
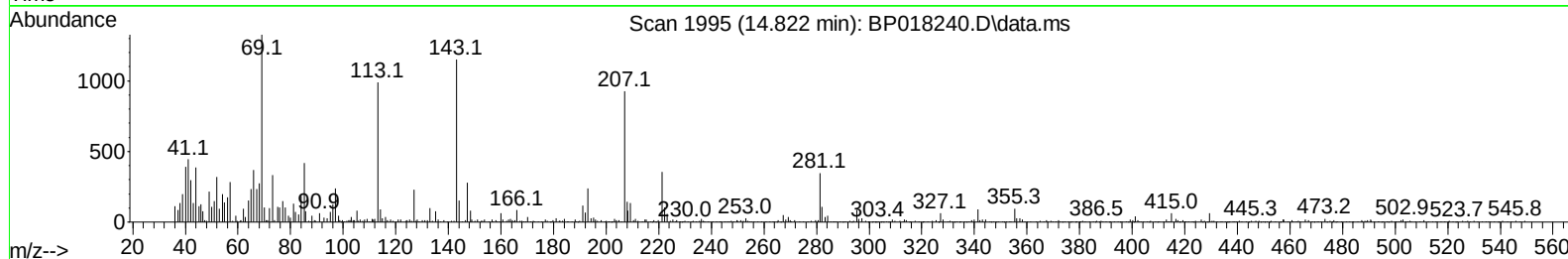
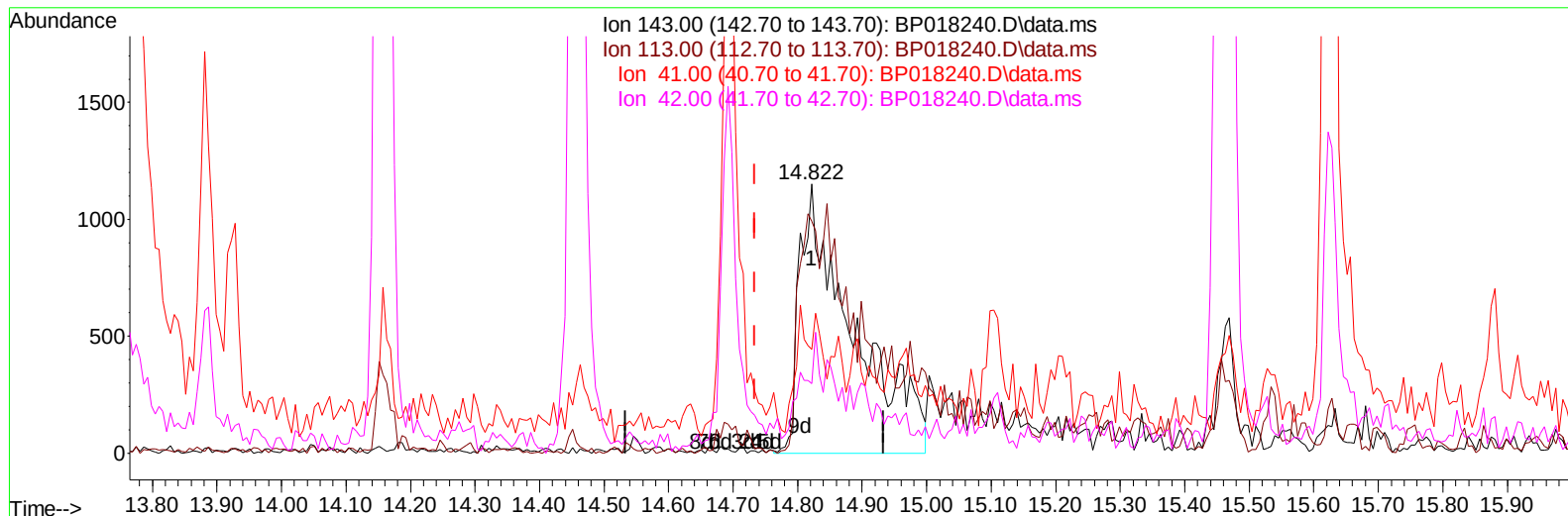
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TIC: BP018240.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (+ 0.088) 3.02 ng/ul m

response 6557

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	86.17
41.00	47.70	38.52
42.00	30.10	25.74

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Quant Time: Nov 18 02:58:15 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	65431	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	259983	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	157618	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	360346	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	364561	20.000	ng/ul	0.00
88) Perylene-d12	23.845	264	407227	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	8577	4.624	ng/uL	0.00
4) Pyridine-d5	3.611	84	20044	4.145	ng/ul	0.02
7) Phenol-d5	6.969	99	25649	4.060	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.128	67	104473	28.143	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	103067	20.604	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	58083	11.221	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	61930	26.440	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	73973	27.876	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	112806	23.851	ng/ul	0.00
31) 4-Chloroaniline-d4	10.787	131	107562	16.584	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	456021	31.357	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	450673	28.452	ng/ul	0.00
54) 4-Nitrophenol-d4	14.822	143	6557m	3.017	ng/ul	0.09
60) Fluorene-d10	15.463	176	377249	31.419	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	66916	25.687	ng/ul	0.00
73) Anthracene-d10	17.339	188	653841	33.549	ng/ul	0.00
81) Pyrene-d10	19.598	212	824212	33.847	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264	809464	34.185	ng/ul	0.00

Target Compounds Qvalue

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

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