

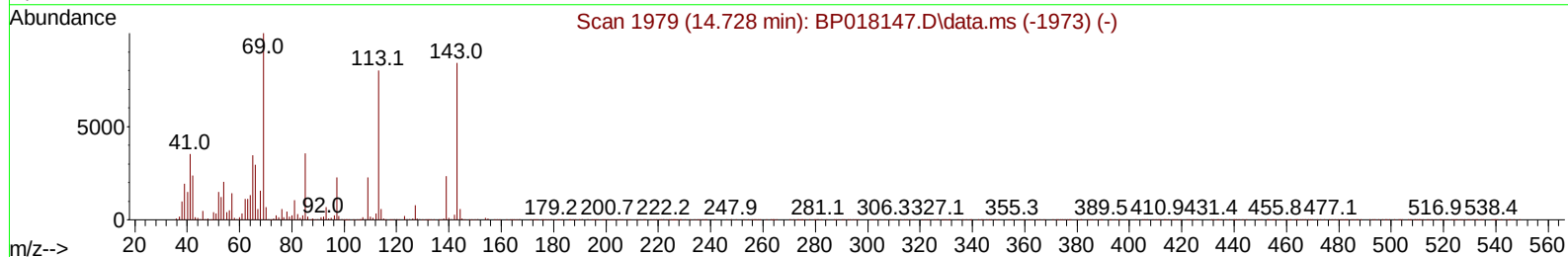
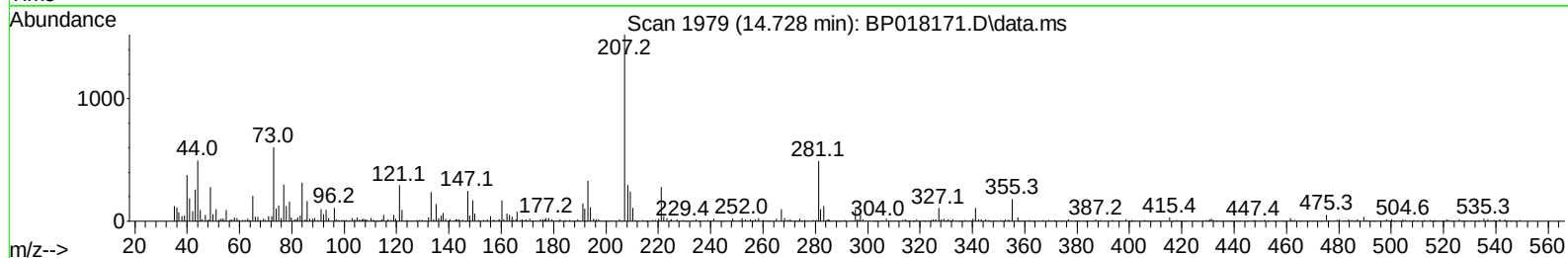
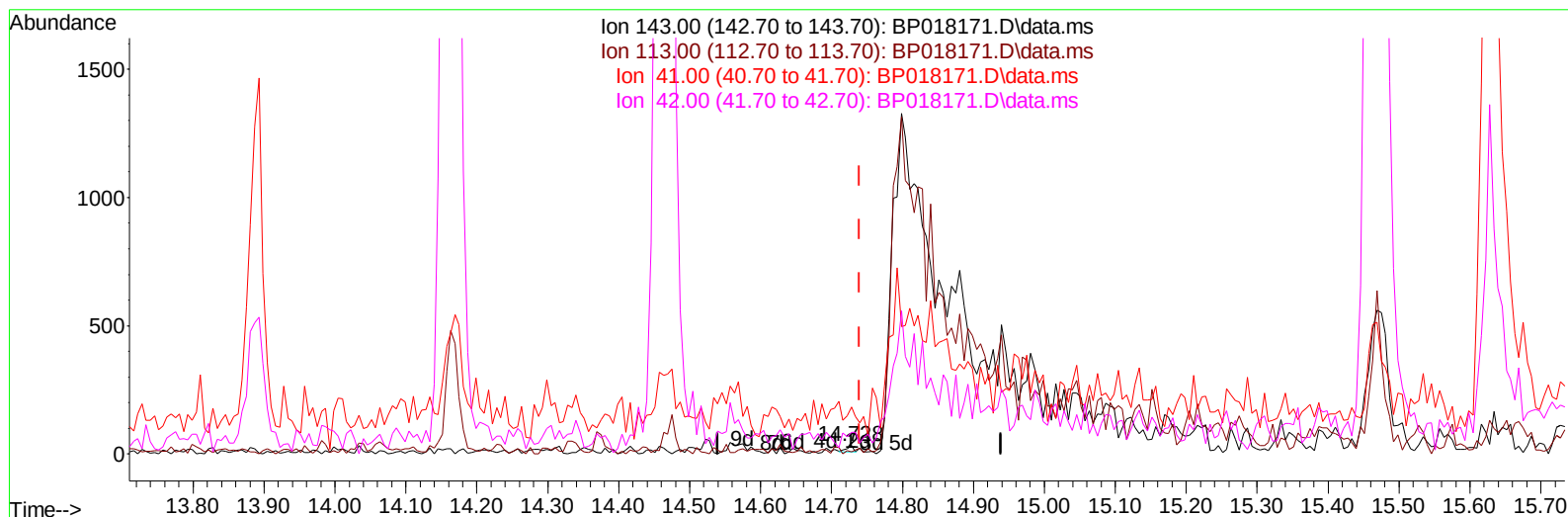
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
Data File : BP018171.D
Acq On : 15 Nov 2023 00:03
Operator : MA/JU
Sample : 05203-01 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AN8

Manual IntegrationsAPPROVED

Quant Time: Nov 15 00:42:04 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 13 15:50:18 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/16/2023
Supervised By :mohammad ahmed 11/16/2023



TIC: BP018171.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (-0.012) 0.01 ng/u1

response 17

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	23.53#
41.00	47.70	541.18#
42.00	30.10	250.00#

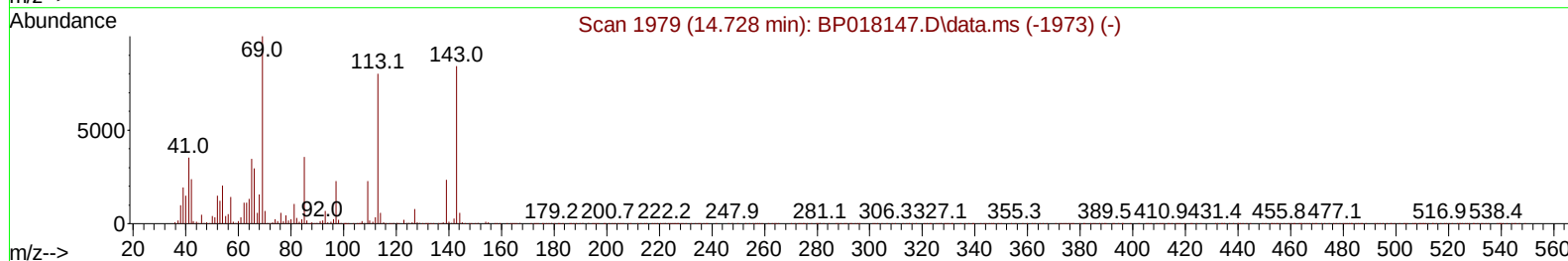
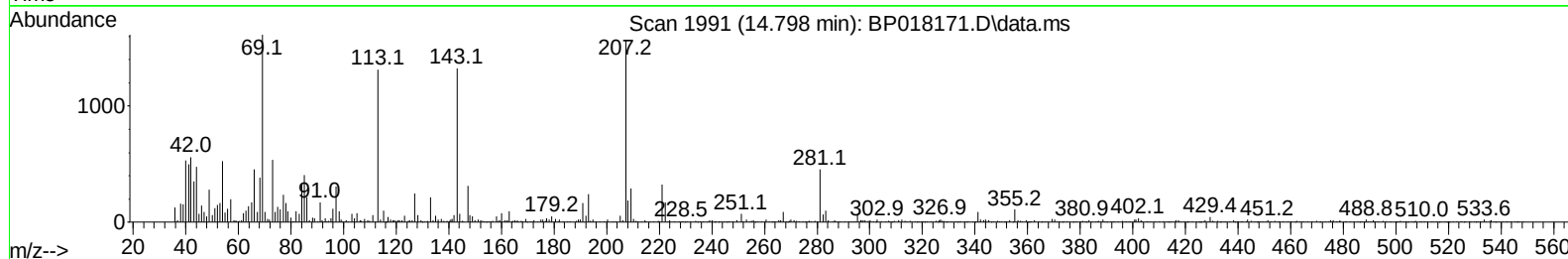
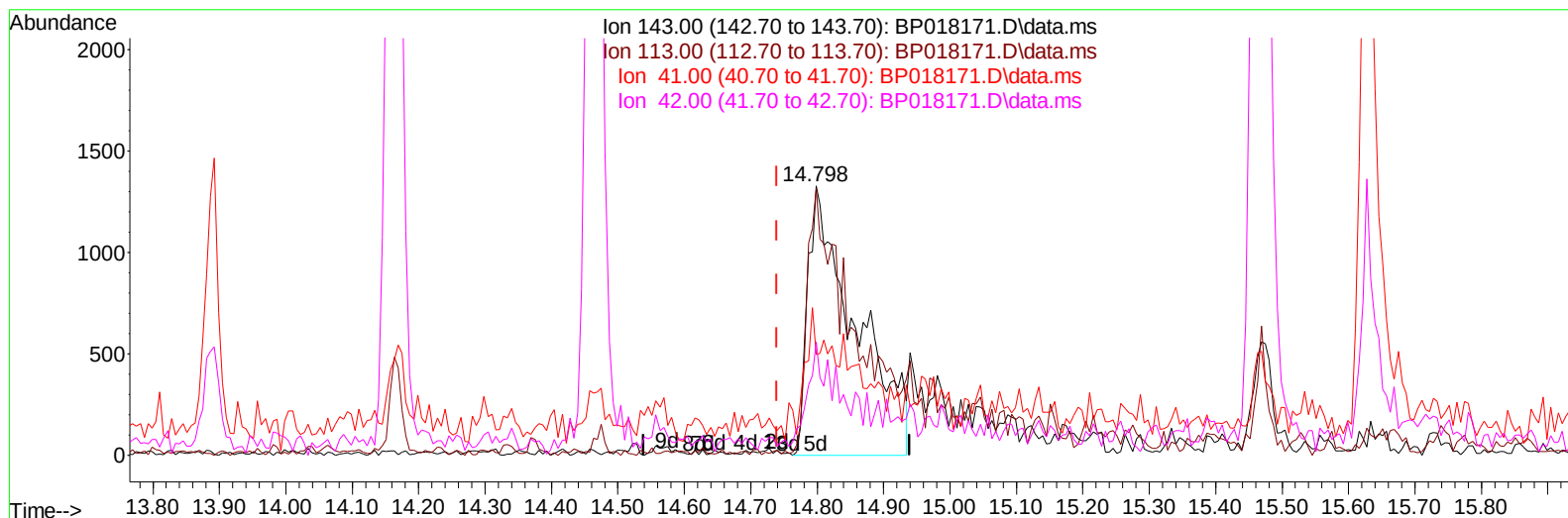
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TIC: BP018171.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.059) 2.30 ng/ul m

response 6591

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	99.10
41.00	47.70	37.45#
42.00	30.10	42.13#

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Instrument :
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Quant Time: Nov 15 01:53:17 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	76509	20.000	ng/ul	0.00
20) Naphthalene-d8	10.605	136	316848	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164	207568	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.245	188	479783	20.000	ng/ul	-0.01
79) Chrysene-d12	21.375	240	468958	20.000	ng/ul	-0.02
88) Perylene-d12	23.845	264	520569	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	5514	2.542	ng/uL	0.00
4) Pyridine-d5	3.611	84	24481	4.329	ng/ul	0.00
7) Phenol-d5	6.964	99	26093	3.533	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	105955	24.409	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	91873	15.707	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	57513	9.502	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	62706	21.966	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	66842	20.668	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	105638	18.327	ng/ul	0.00
31) 4-Chloroaniline-d4	10.787	131	133360	16.871	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	486311	25.392	ng/ul	-0.01
49) Acenaphthylene-d8	14.163	160	482989	23.154	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	6591m	2.303	ng/ul	0.06
60) Fluorene-d10	15.469	176	400844	25.351	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	55348	15.957	ng/ul	-0.04
73) Anthracene-d10	17.345	188	690028	26.592	ng/ul	-0.01
81) Pyrene-d10	19.604	212	853109	27.235	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.680	264	827726	27.345	ng/ul	-0.02

Target Compounds Qvalue

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

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