

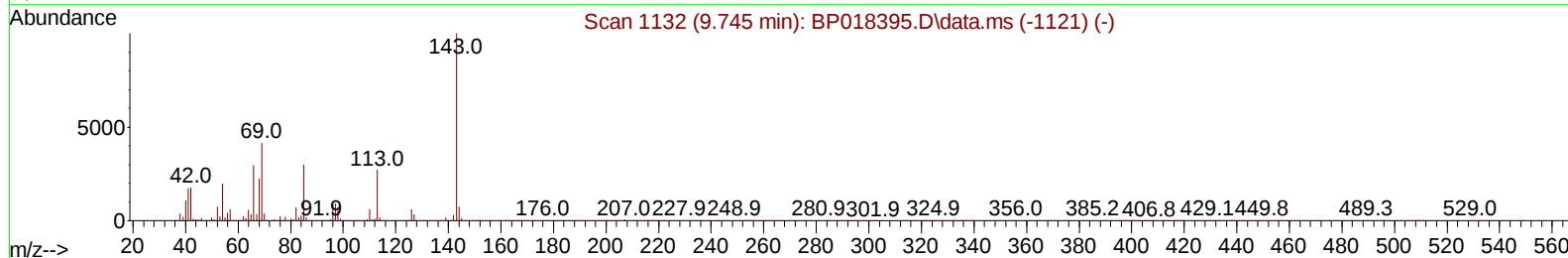
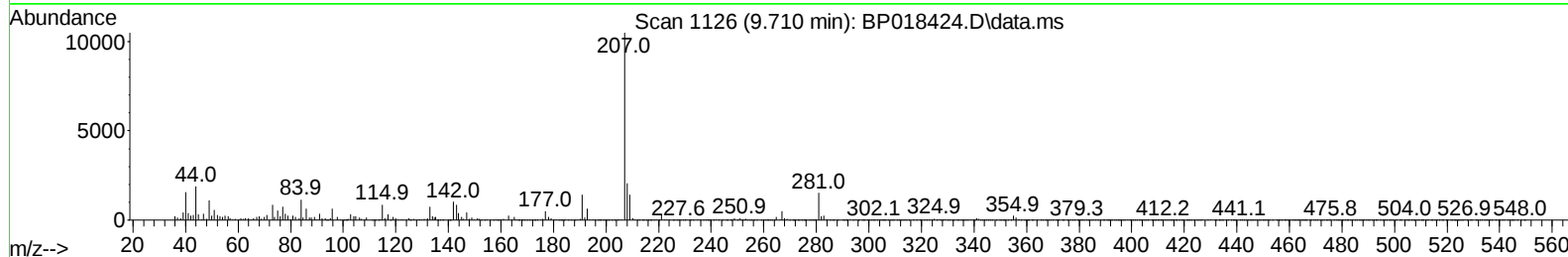
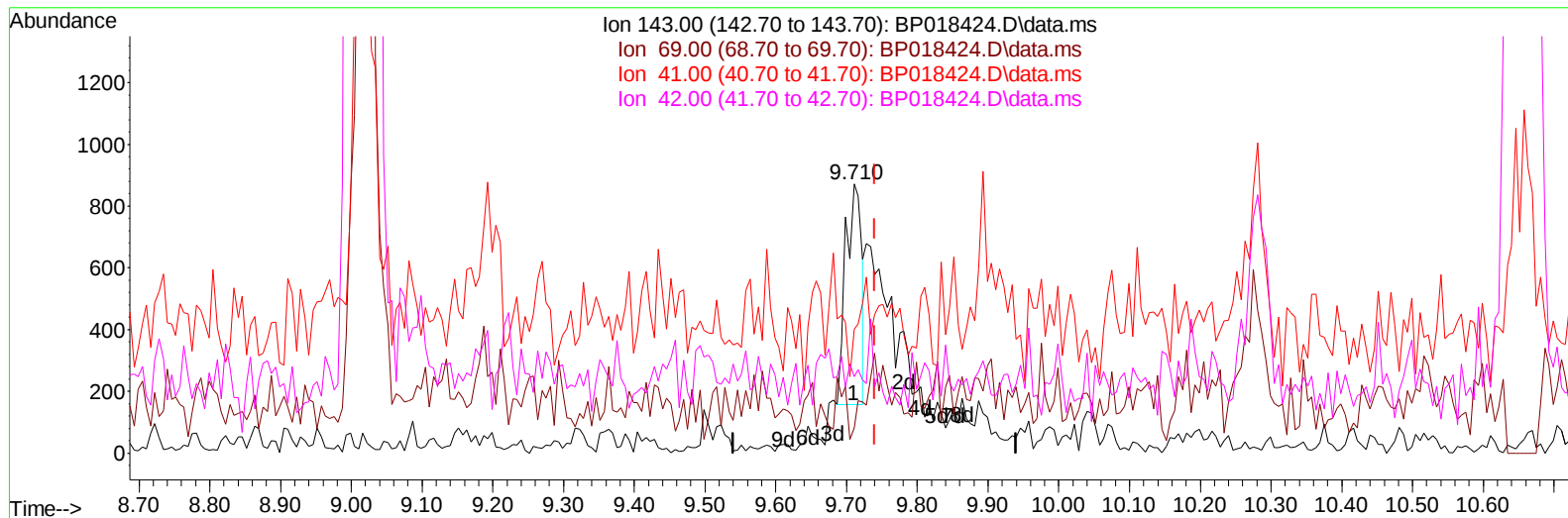
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018424.D
 Acq On : 28 Nov 2023 05:05
 Operator : MA/JU
 Sample : 05391-01
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4A06

Manual IntegrationsAPPROVED

Quant Time: Nov 28 05:57:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 28 02:41:10 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/29/2023
 Supervised By :mohammad ahmed 12/01/2023



TIC: BP018424.D\data.ms

(24) 2-Nitrophenol-d4 (S)

9.710min (-0.029) 0.07 ng/u1

response 1107

Ion	Exp%	Act%
143.00	100.00	100.00
69.00	44.40	9.41#
41.00	21.30	46.04#
42.00	20.90	28.36#

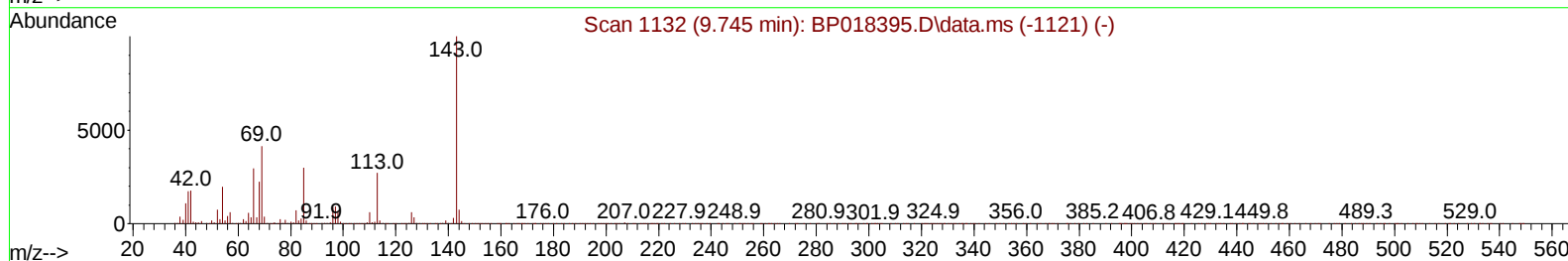
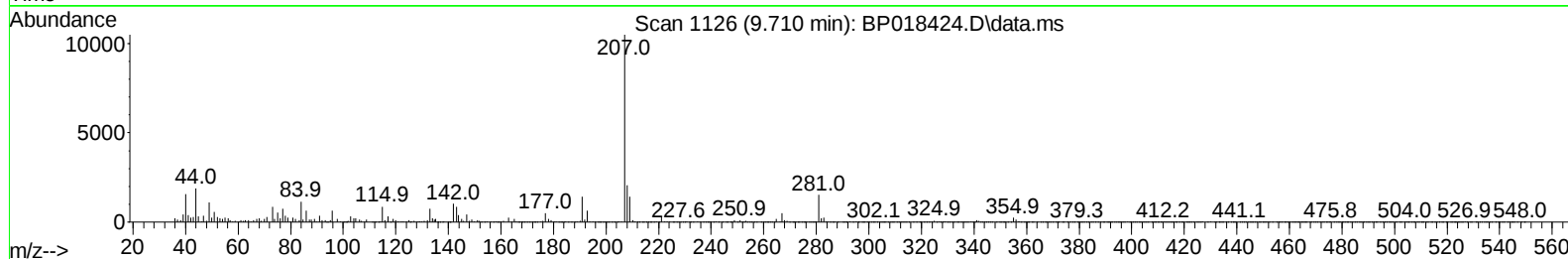
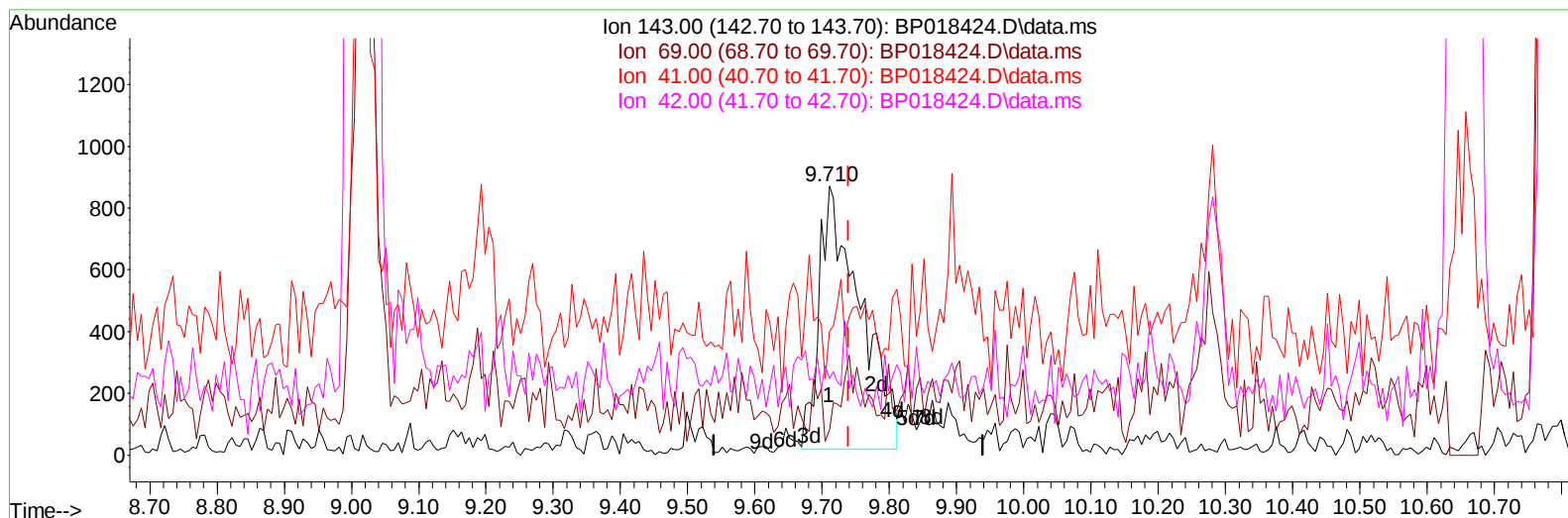
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(24) 2-Nitrophenol-d4 (S)

9.710min (-0.029) 0.23 ng/ul m

response 3605

Ion	Exp%	Act%
143.00	100.00	100.00
69.00	44.40	9.41#
41.00	21.30	46.04#
42.00	20.90	28.36#

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.858	152		510647	20.000	ng/ul	0.00
20) Naphthalene-d8	10.657	136		2088006	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.493	164		1258887	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188		2260699	20.000	ng/ul	0.00
79) Chrysene-d12	21.392	240		1743576	20.000	ng/ul	0.00
88) Perylene-d12	23.939	264		2081871	20.000	ng/ul	0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.264	96		53780	3.945	ng/uL	0.00
4) Pyridine-d5	3.687	84		500674	13.363	ng/ul	0.00
7) Phenol-d5	7.022	99		11357	0.243	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67		664373	23.807	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132		3271	0.083	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113		59190	1.572	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128		466373	27.499	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143		3605m	0.230	ng/ul	-0.03
28) 2,4-Dichlorophenol-d3	10.275	165		2012	0.053	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131		1279596	25.107	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166		1969201	17.580	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160		3260066	25.542	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143		58	0.004	ng/ul	0.00
60) Fluorene-d10	15.487	176		2344471	25.023	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200		42	0.004	ng/ul	0.00
73) Anthracene-d10	17.345	188		3144213	26.021	ng/ul	0.00
81) Pyrene-d10	19.592	212		3174895	22.879	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.769	264		3241686	26.858	ng/ul	0.03

Target Compounds Qvalue

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

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