

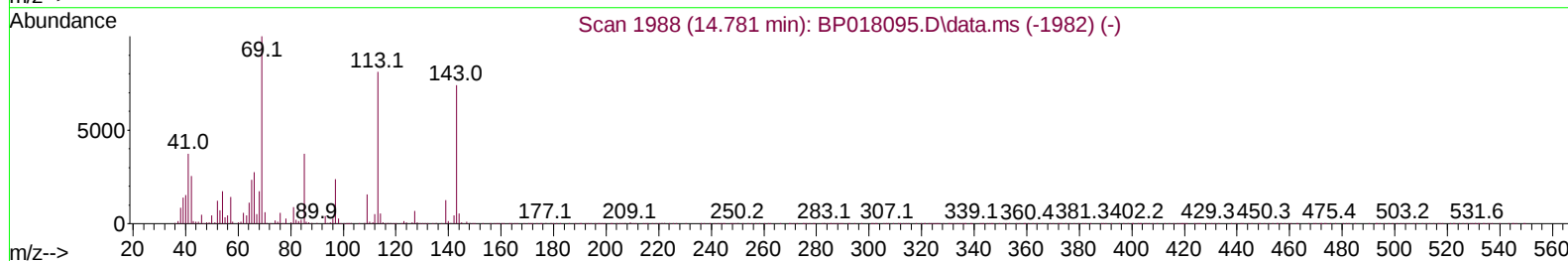
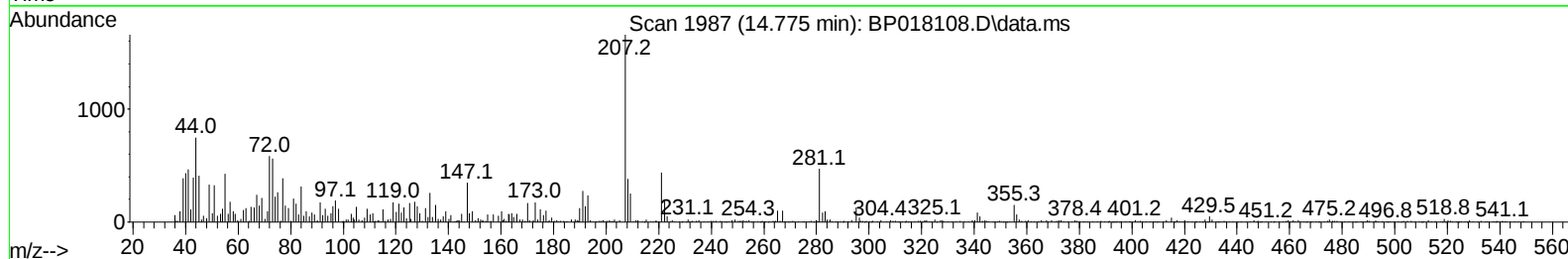
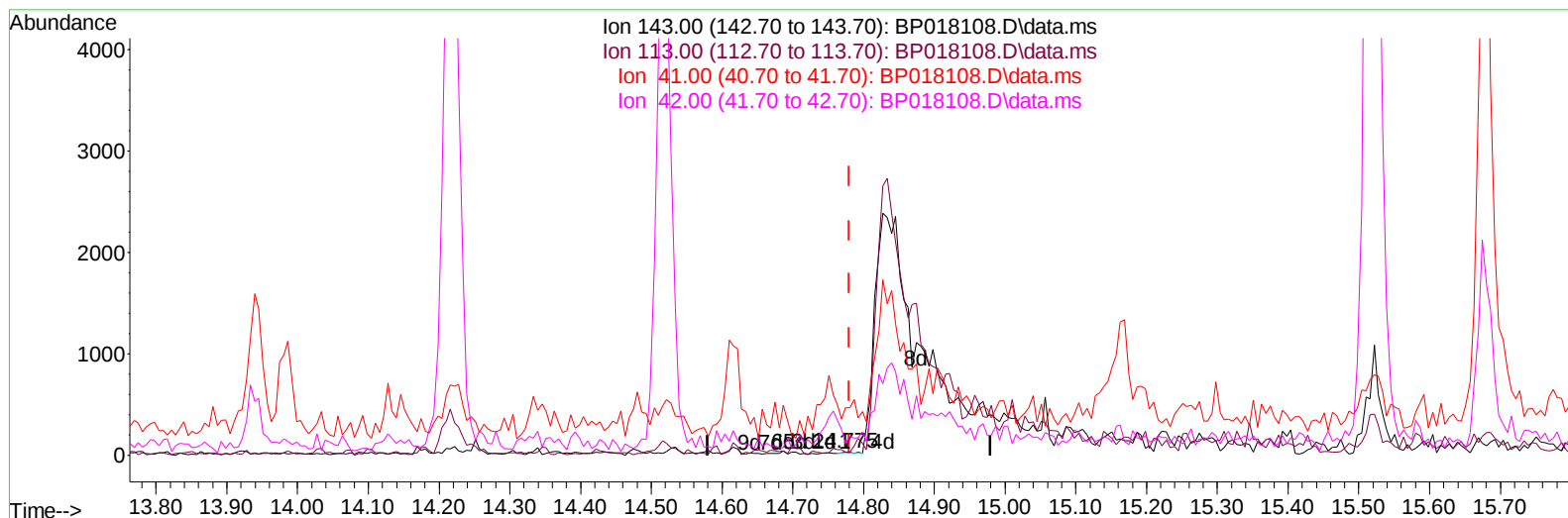
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018108.D
 Acq On : 12 Nov 2023 16:38
 Operator : MA/JU
 Sample : 05082-03
 Misc :
 ALS Vial : 88 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH345

Manual IntegrationsAPPROVED

Quant Time: Nov 12 22:05:55 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/13/2023
 Supervised By :mohammad ahmed 11/15/2023



TIC: BP018108.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.006) 0.01 ng/u1

response 17

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	96.43
41.00	47.70	1664.29#
42.00	30.10	407.14#

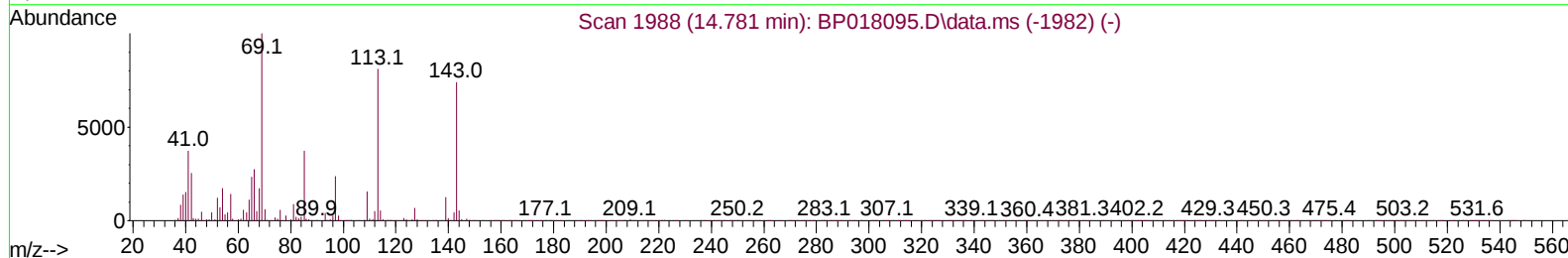
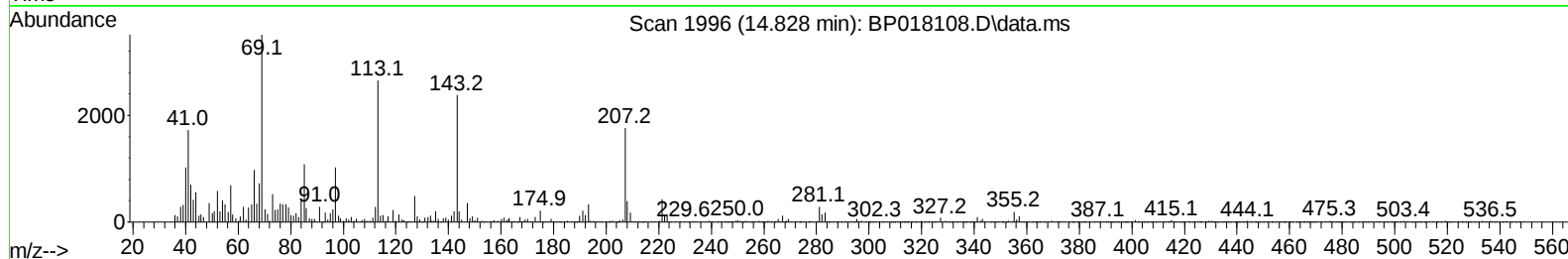
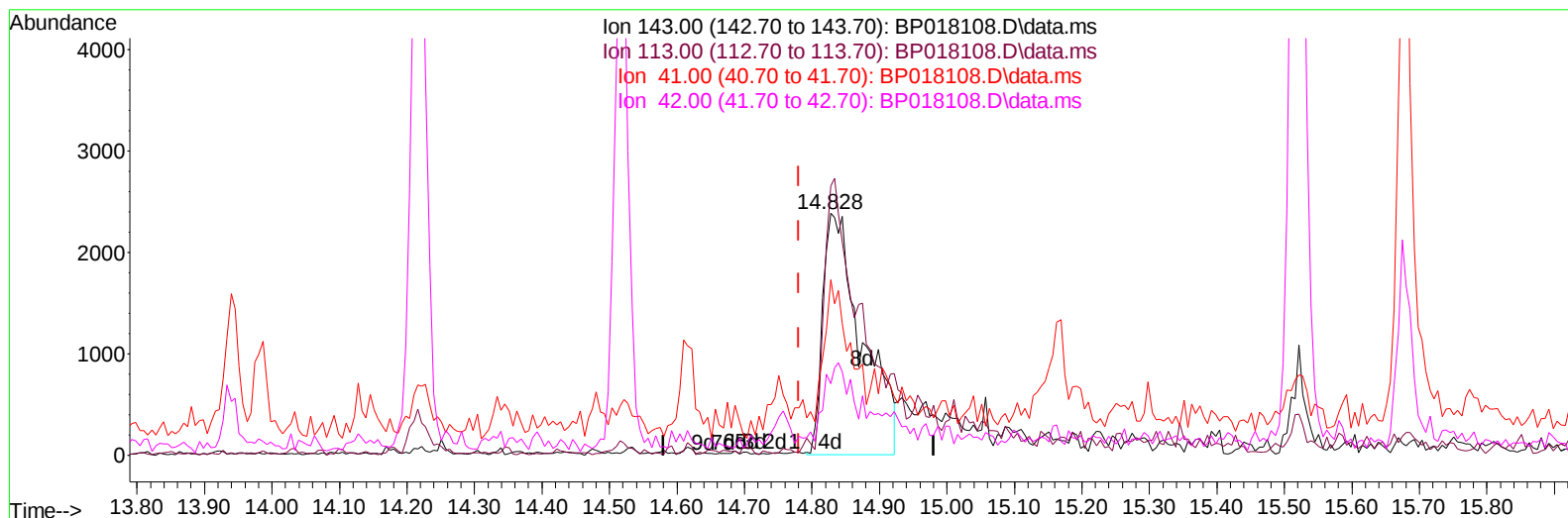
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TIC: BP018108.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.828min (+ 0.047) 3.94 ng/ul m

response 9593

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	111.27
41.00	47.70	72.51#
42.00	30.10	29.42

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 Misc :
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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/13/2023
 Supervised By :mohammad ahmed 11/15/2023

Quant Time: Nov 12 22:07:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	77924	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.663	136	303419	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.516	164	176587	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.298	188	379356	20.000	ng/ul	-0.01
79) Chrysene-d12	21.427	240	399793	20.000	ng/ul	0.00
88) Perylene-d12	23.927	264	463845	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	7801	3.532	ng/uL	0.00
4) Pyridine-d5	3.675	84	27790	4.825	ng/ul	0.01
7) Phenol-d5	7.022	99	44532	5.919	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	139046	31.451	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	131834	22.129	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	86361	14.009	ng/ul	0.00
21) Nitrobenzene-d5	9.051	128	85910	31.427	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	92879	29.990	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	143586	26.013	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	164442	21.724	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	532715	32.695	ng/ul	-0.01
49) Acenaphthylene-d8	14.216	160	561394	31.634	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.828	143	9593m	3.940	ng/ul	0.05
60) Fluorene-d10	15.522	176	442619	32.904	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	83567	30.472	ng/ul	0.00
73) Anthracene-d10	17.398	188	740715	36.102	ng/ul	-0.01
81) Pyrene-d10	19.657	212	939484	35.181	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.768	264	993646	36.841	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.264	88	49728	20.630	ng/uL	100

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

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