

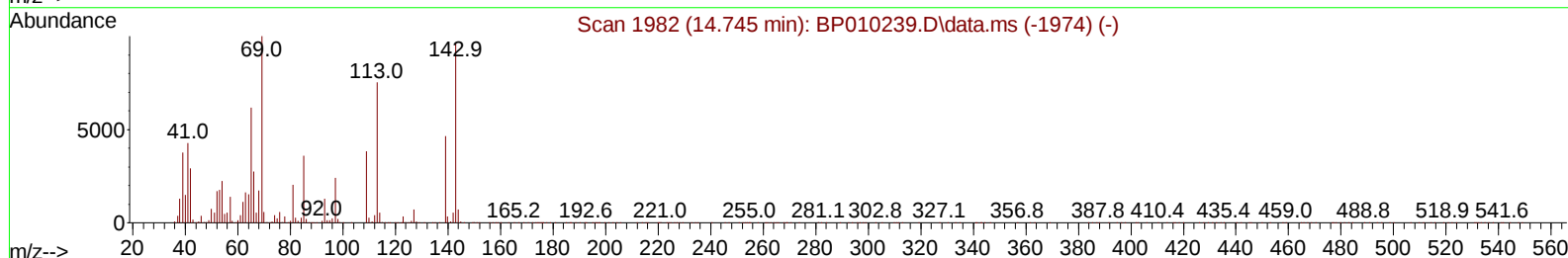
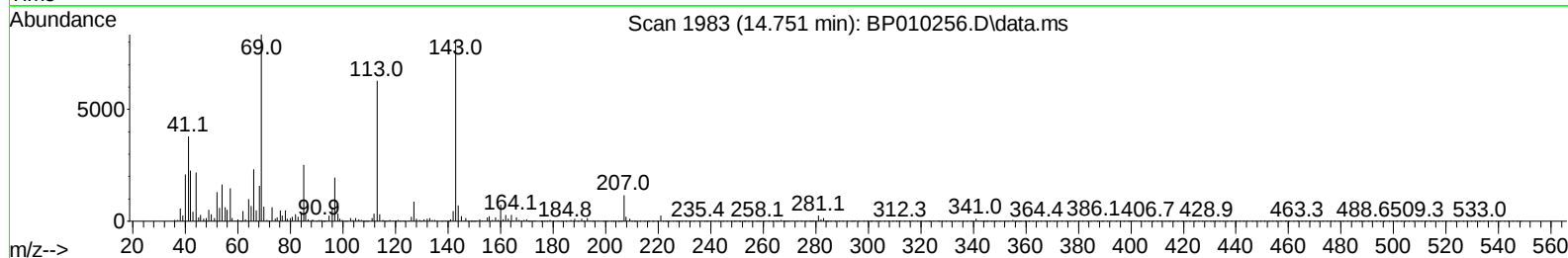
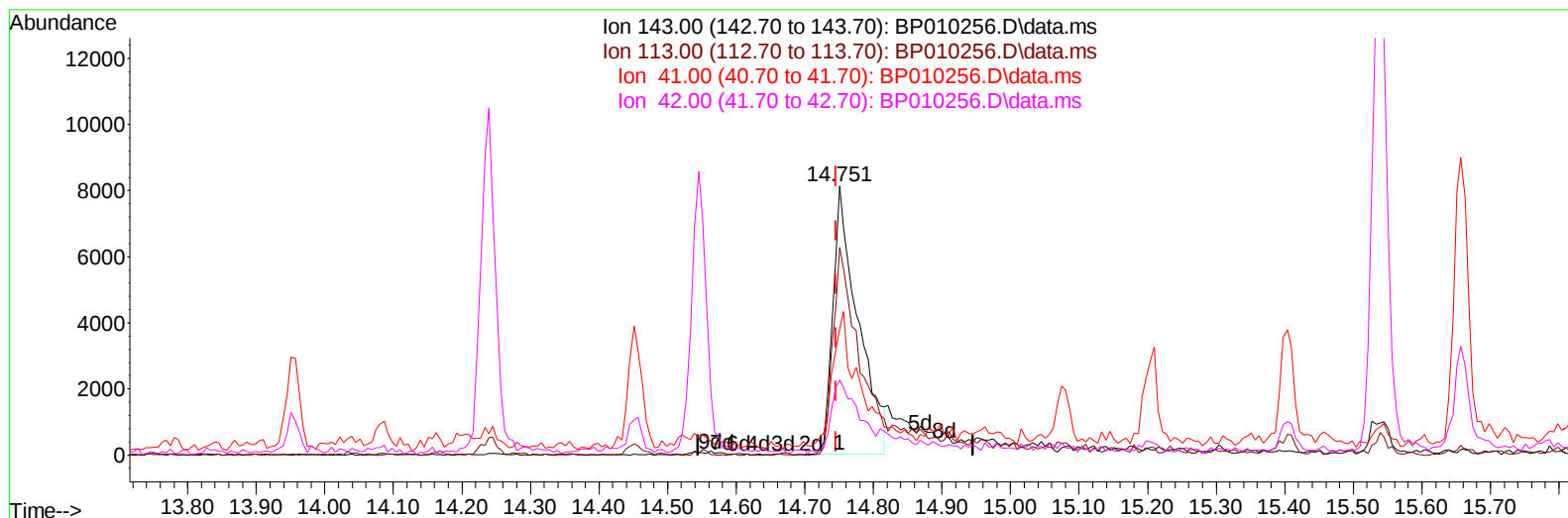
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
 Data File : BP010256.D
 Acq On : 05 May 2022 20:14
 Operator : CG/JU
 Sample : N2678-05
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH0C0

Manual IntegrationsAPPROVED

Quant Time: May 06 02:08:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 06 01:05:07 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/10/2022
 Supervised By :mohammad ahmed 05/10/2022



TIC: BP010256.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 3.69 ng/u1

response 20637

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.32#
41.00	23.20	46.74#
42.00	18.00	27.88#

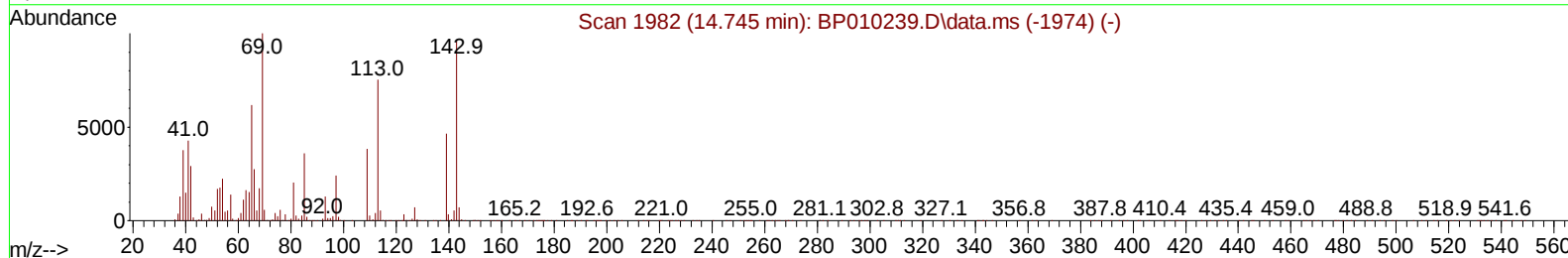
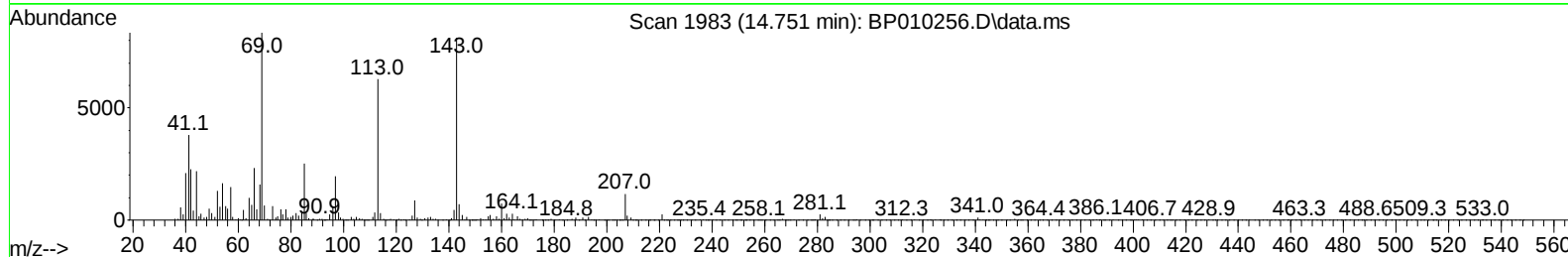
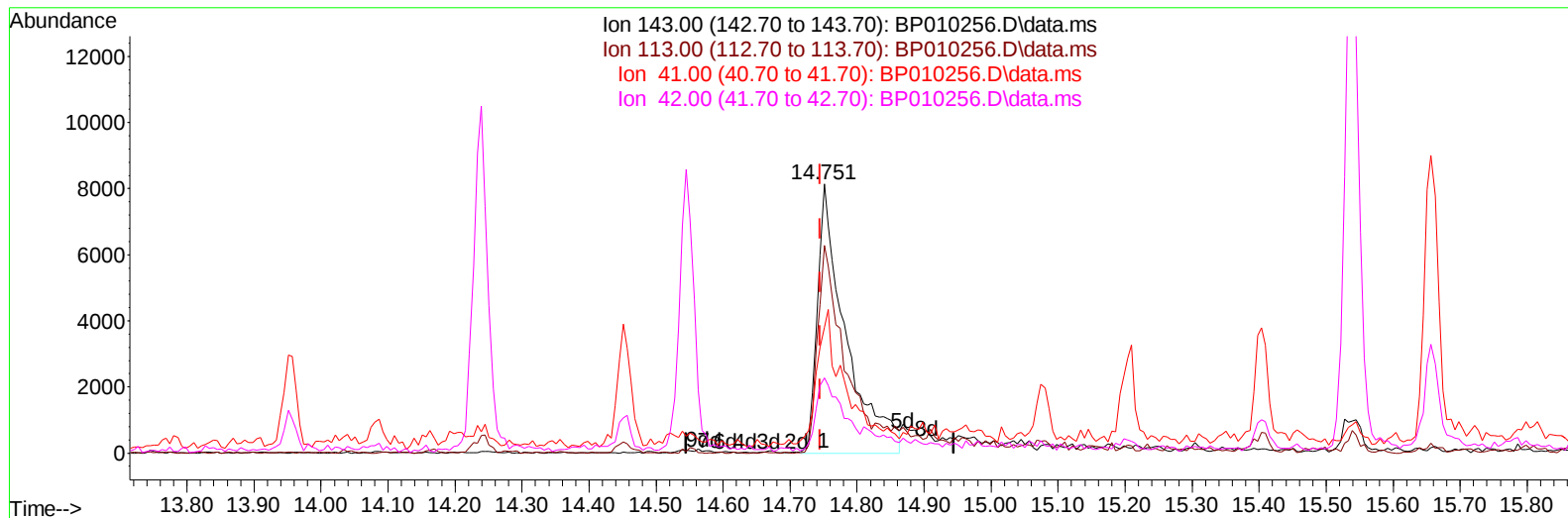
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TIC: BP010256.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 4.24 ng/ul m

response 23707

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.32#
41.00	23.20	46.74#
42.00	18.00	27.88#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	159761	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.716	136	662360	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.545	164	390349	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.298	188	706532	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.380	240	659380	20.000	ng/ul	# 0.00
88) Perylene-d12	23.821	264	664819	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	20442	5.580	ng/uL	0.00
4) Pyridine-d5	3.781	84	67400	6.539	ng/ul	0.01
7) Phenol-d5	7.081	99	99273	6.831	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	315780	36.215	ng/ul	0.00
11) 2-Chlorophenol-d4	7.445	132	288680	26.824	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	190592	15.830	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	190422	36.147	ng/ul	0.00
24) 2-Nitrophenol-d4	9.798	143	200131	34.785	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.339	165	307622	27.465	ng/ul	0.00
31) 4-Chloroaniline-d4	10.857	131	255446	15.906	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1091480	35.946	ng/ul	0.00
49) Acenaphthylene-d8	14.239	160	1239568	33.730	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	23707m	4.241	ng/ul	0.00
60) Fluorene-d10	15.539	176	903046	34.573	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.657	200	142231	31.837	ng/ul	0.00
73) Anthracene-d10	17.398	188	1395248	40.819	ng/ul	0.00
81) Pyrene-d10	19.627	212	1598979	42.640	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.662	264	1489429	42.734	ng/ul	0.00

Target Compounds Qvalue

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

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