

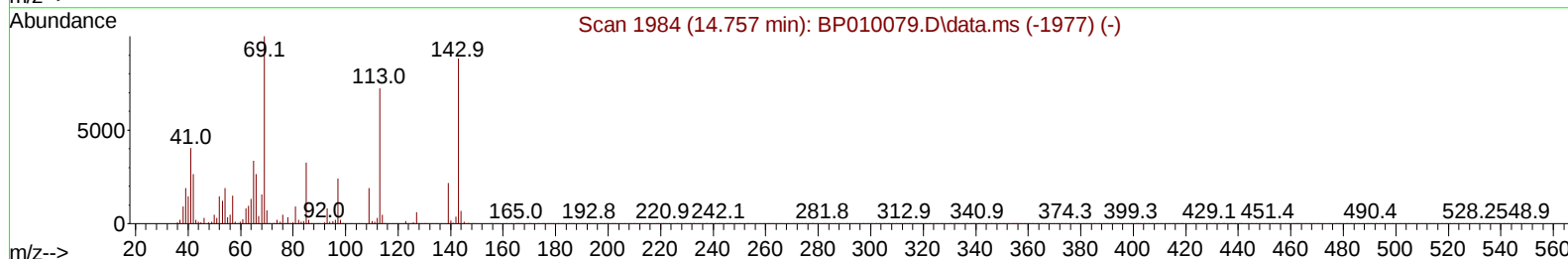
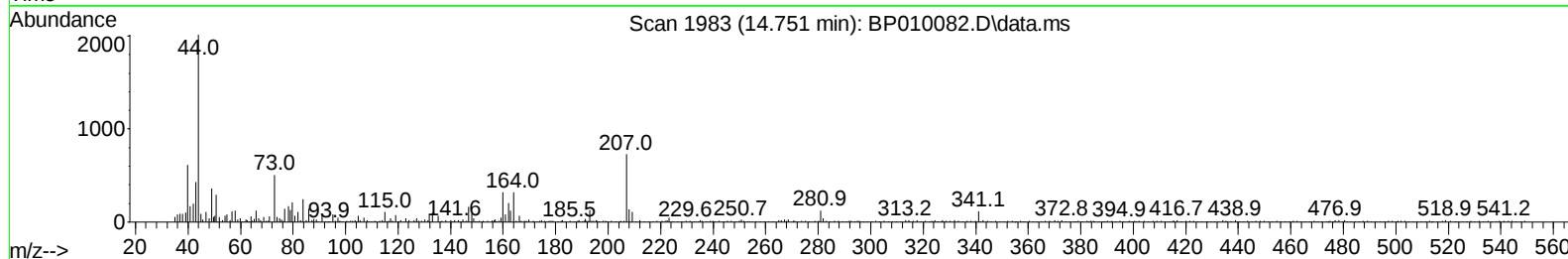
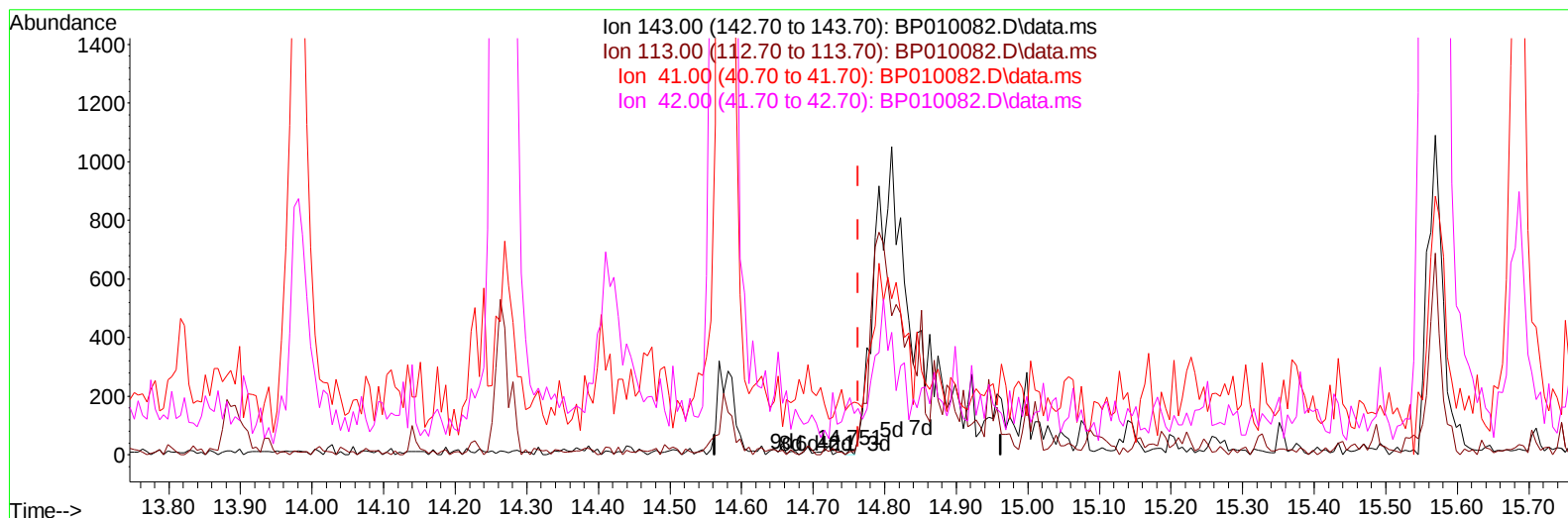
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
 Data File : BP010082.D
 Acq On : 25 Apr 2022 14:51
 Operator : CG/JU
 Sample : N2475-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A0BM3

Manual IntegrationsAPPROVED

Quant Time: Apr 26 01:00:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022
 Supervised By :Yogesh Patel 04/28/2022



TIC: BP010082.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.012) 0.00 ng/u1

response 7

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	57.89#
41.00	23.20	900.00#
42.00	18.00	1026.32#

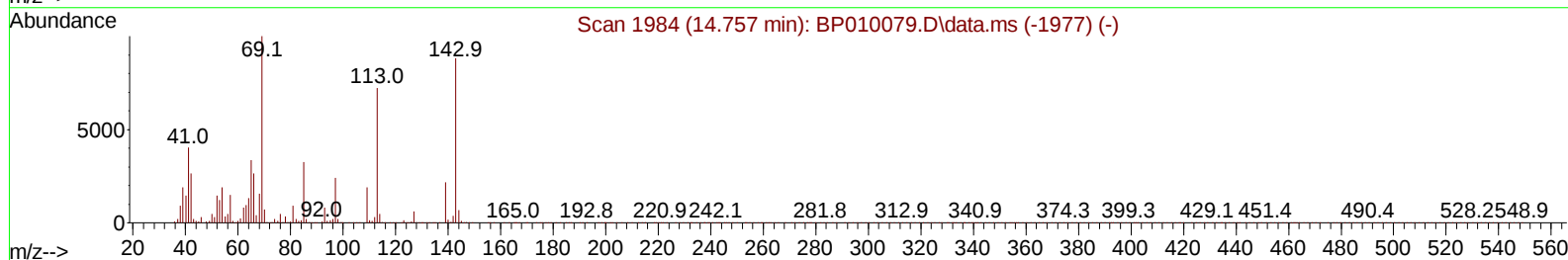
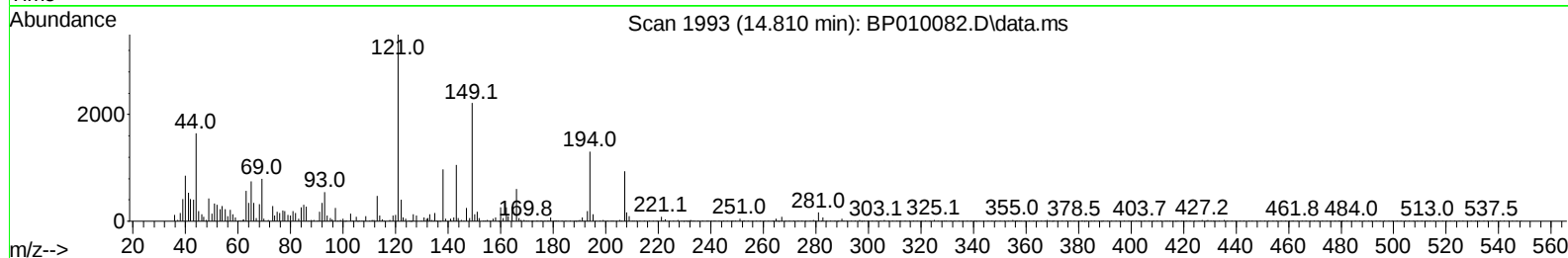
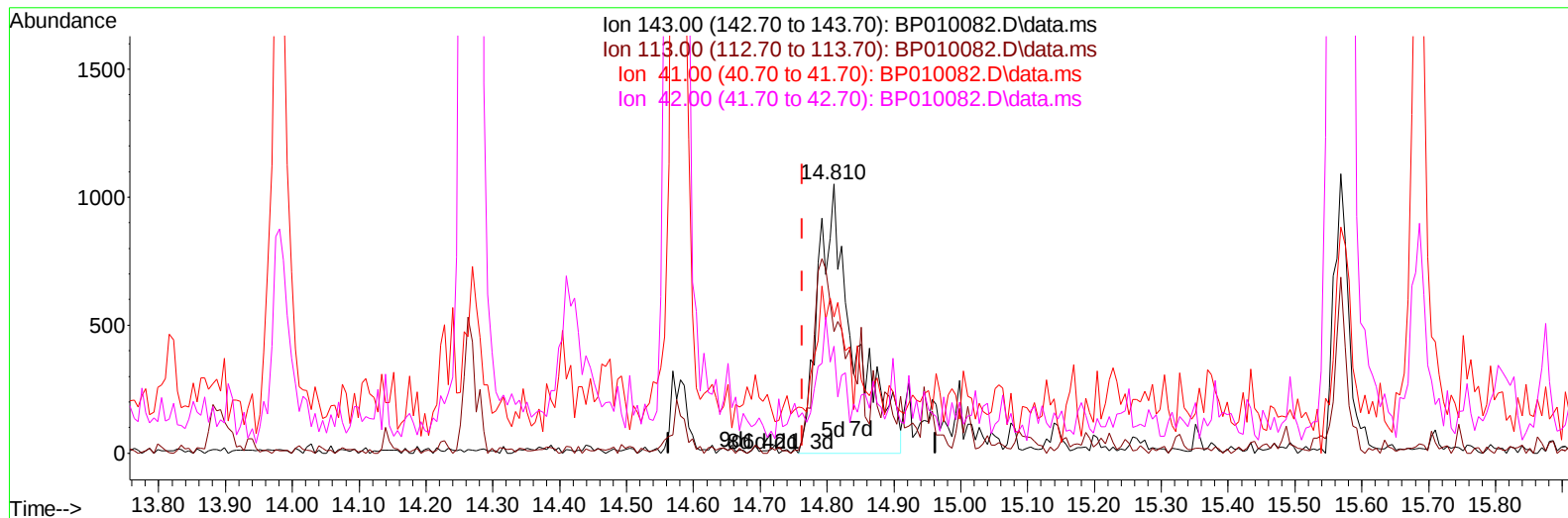
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TIC: BP010082.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.047) 0.68 ng/ul m

response 3935

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	45.10#
41.00	23.20	50.62#
42.00	18.00	39.77#

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 Misc :
 ALS Vial : 5 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	160706	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	634317	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	405595	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	812098	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.410	240	595545	20.000	ng/ul	# 0.00
88) Perylene-d12	23.874	264	486691	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	21993	5.968	ng/uL	0.00
4) Pyridine-d5	3.787	84	107801	10.398	ng/ul	0.00
7) Phenol-d5	7.105	99	42037	2.876	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	263168	30.003	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	118908	10.984	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	95564	7.891	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	159925	31.700	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	45541	8.265	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	121944	11.369	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	367557	23.899	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1035941	32.834	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1141590	29.896	ng/ul	0.00
54) 4-Nitrophenol-d4	14.810	143	3935m	0.677	ng/ul	0.05
60) Fluorene-d10	15.569	176	901342	33.211	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	33533	6.530	ng/ul	0.00
73) Anthracene-d10	17.428	188	1541004	39.223	ng/ul	0.00
81) Pyrene-d10	19.657	212	1643740	48.532	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	1056045	41.389	ng/ul	0.01

Target Compounds Qvalue

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

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