

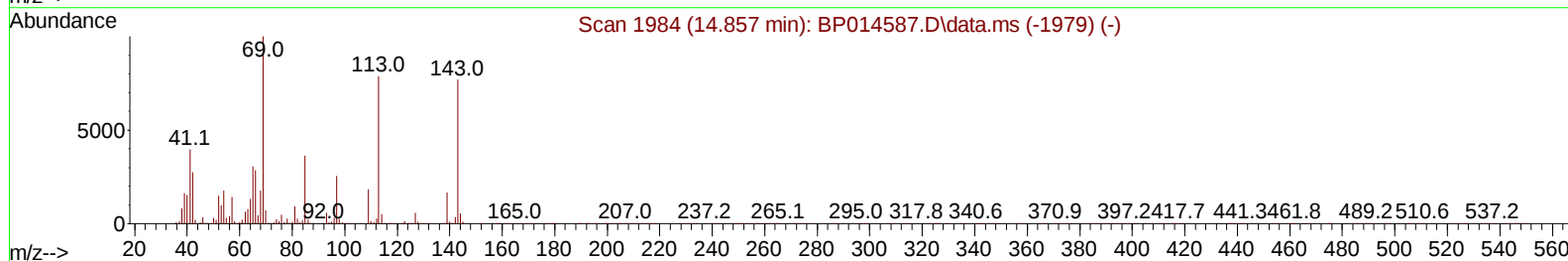
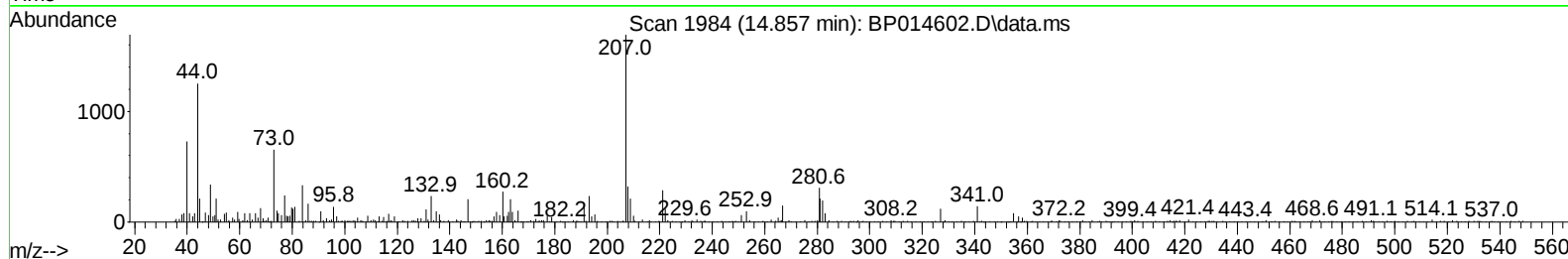
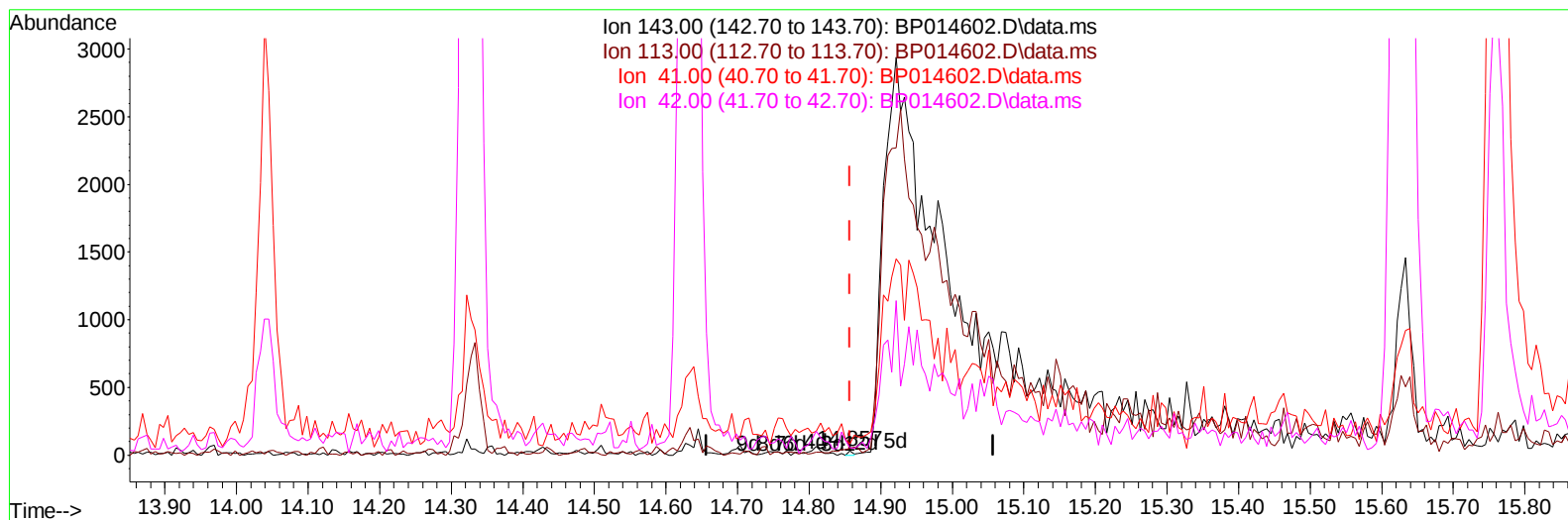
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
 Data File : BP014602.D
 Acq On : 06 Apr 2023 23:29
 Operator : CG/JU
 Sample : 02225-04
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AC9

Manual IntegrationsAPPROVED

Quant Time: Apr 07 00:48:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 07 00:45:34 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/07/2023
 Supervised By :Jagrut Upadhyay 04/07/2023



TIC: BP014602.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (-0.000) 0.00 ng/u1

response 14

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	159.38#
41.00	44.20	253.13#
42.00	31.50	168.75#

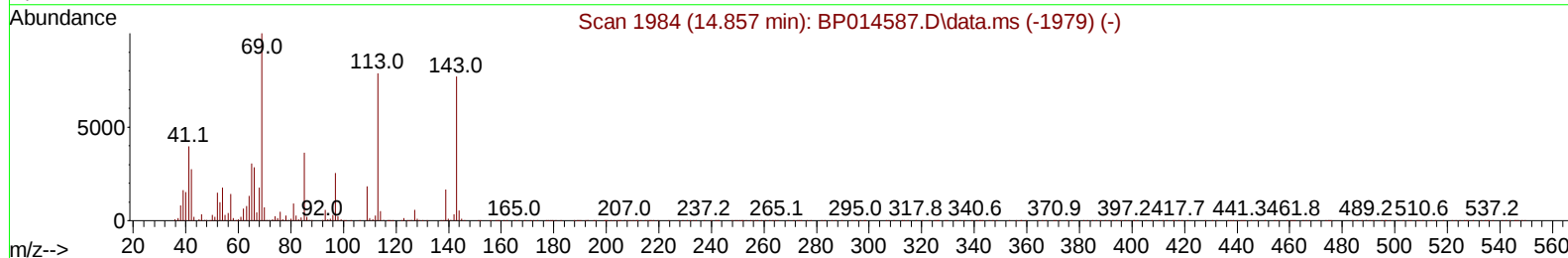
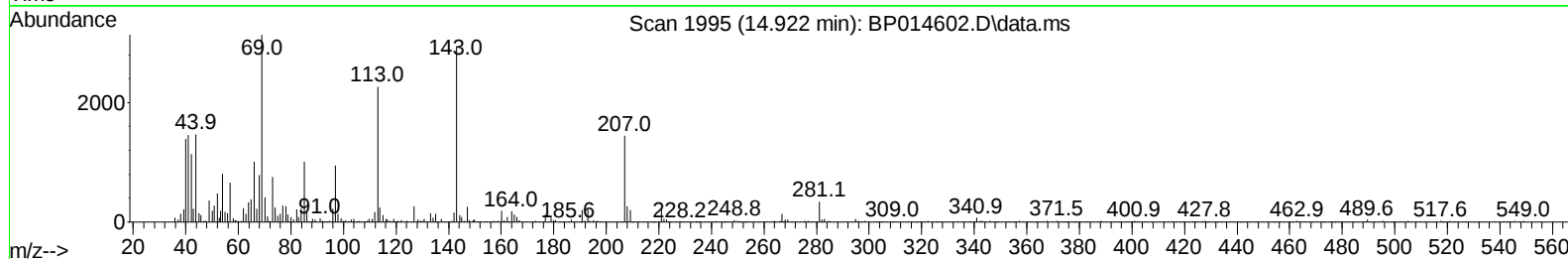
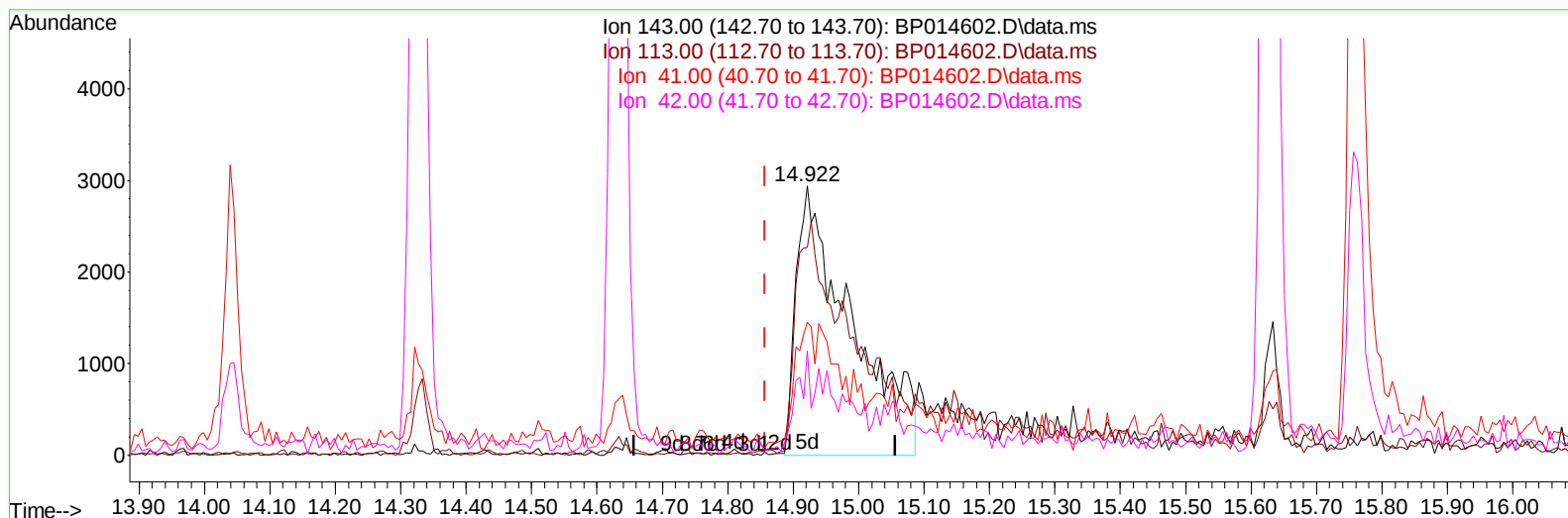
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(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.064) 2.86 ng/ul m

response 17272

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	77.17#
41.00	44.20	49.37
42.00	31.50	38.75#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	163493	20.000	ng/ul	0.00
20) Naphthalene-d8	10.798	136	733520	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.633	164	493652	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	1073864	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	811315	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	829460	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	16878	4.003	ng/uL	0.00
4) Pyridine-d5	3.810	84	70050	6.484	ng/ul	0.02
7) Phenol-d5	7.157	99	78431	5.202	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67	273951	28.318	ng/ul	0.00
11) 2-Chlorophenol-d4	7.516	132	225475	20.690	ng/ul	0.00
15) 4-Methylphenol-d8	8.704	113	155223	12.762	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	150113	25.350	ng/ul	0.00
24) 2-Nitrophenol-d4	9.887	143	172673	25.466	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	263046	21.301	ng/ul	0.00
31) 4-Chloroaniline-d4	10.951	131	295793	18.185	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	1122006	28.032	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	1209395	27.231	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	17272m	2.856	ng/ul	0.06
60) Fluorene-d10	15.627	176	958875	28.727	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	163665	20.713	ng/ul	0.00
73) Anthracene-d10	17.492	188	1556296	29.946	ng/ul	0.00
81) Pyrene-d10	19.727	212	1722678	31.611	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1306368	29.644	ng/ul	0.00

Target Compounds Qvalue

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

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