

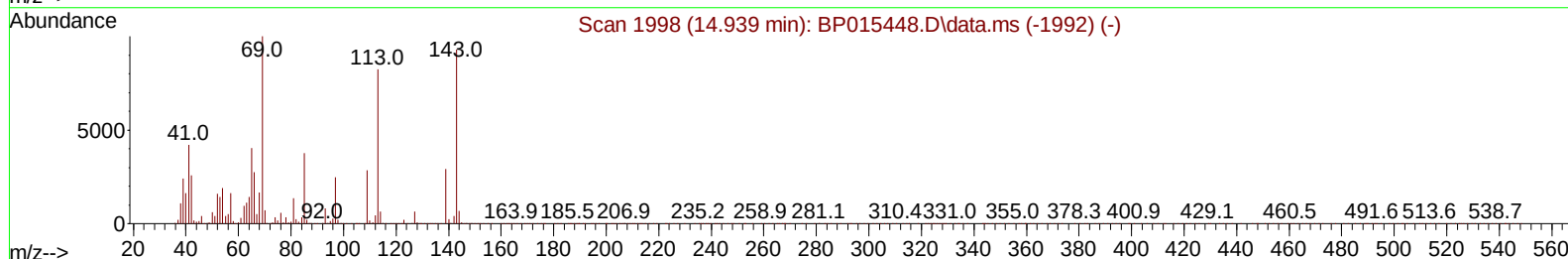
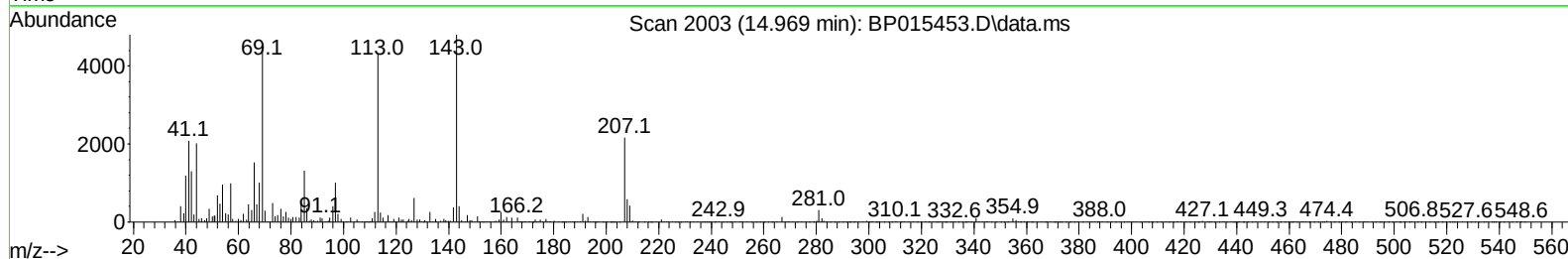
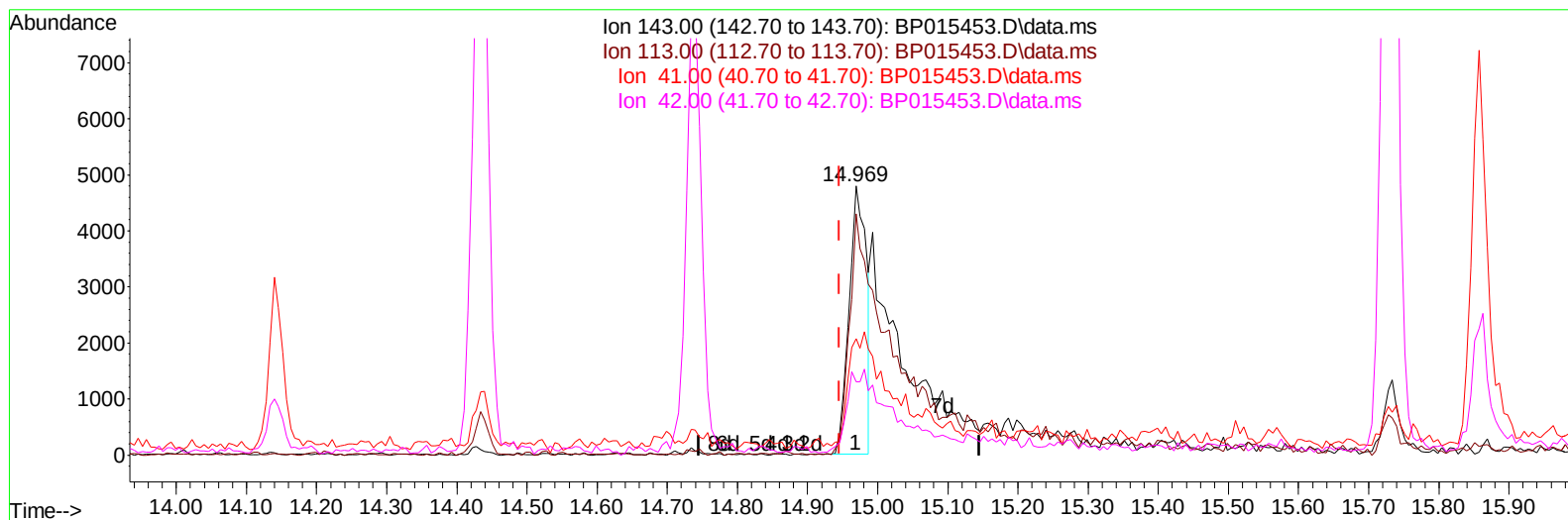
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060923\
 Data File : BP015453.D
 Acq On : 09 Jun 2023 15:45
 Operator : MA/JU
 Sample : 03097-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZEP5

Manual IntegrationsAPPROVED

Quant Time: Jun 09 22:58:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 07 23:46:15 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/12/2023
 Supervised By :mohammad ahmed 06/13/2023



TIC: BP015453.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.024) 1.91 ng/ul

response 8295

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	89.81
41.00	42.20	43.33
42.00	29.40	27.24

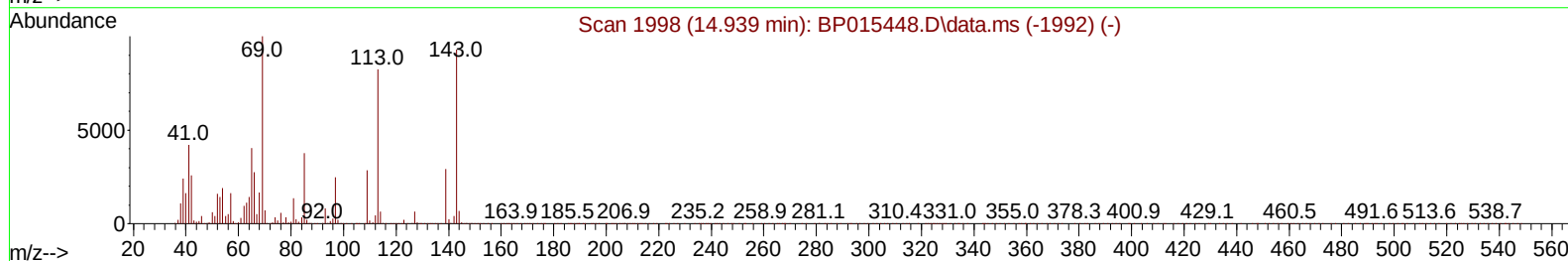
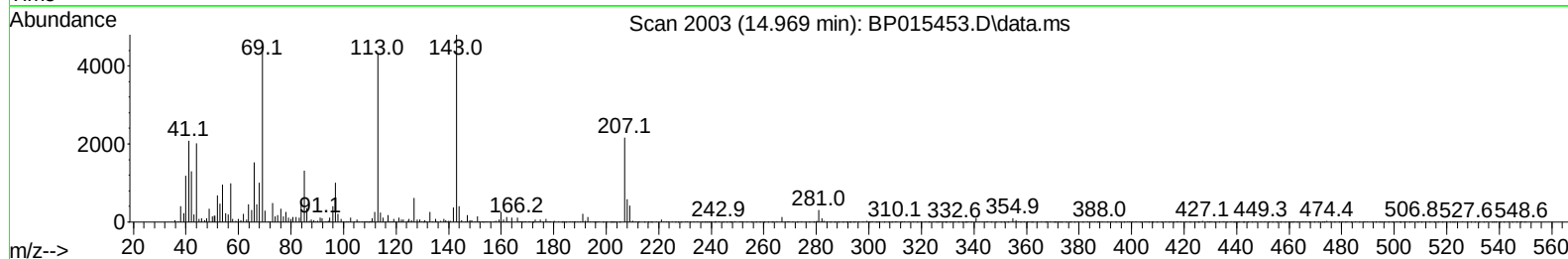
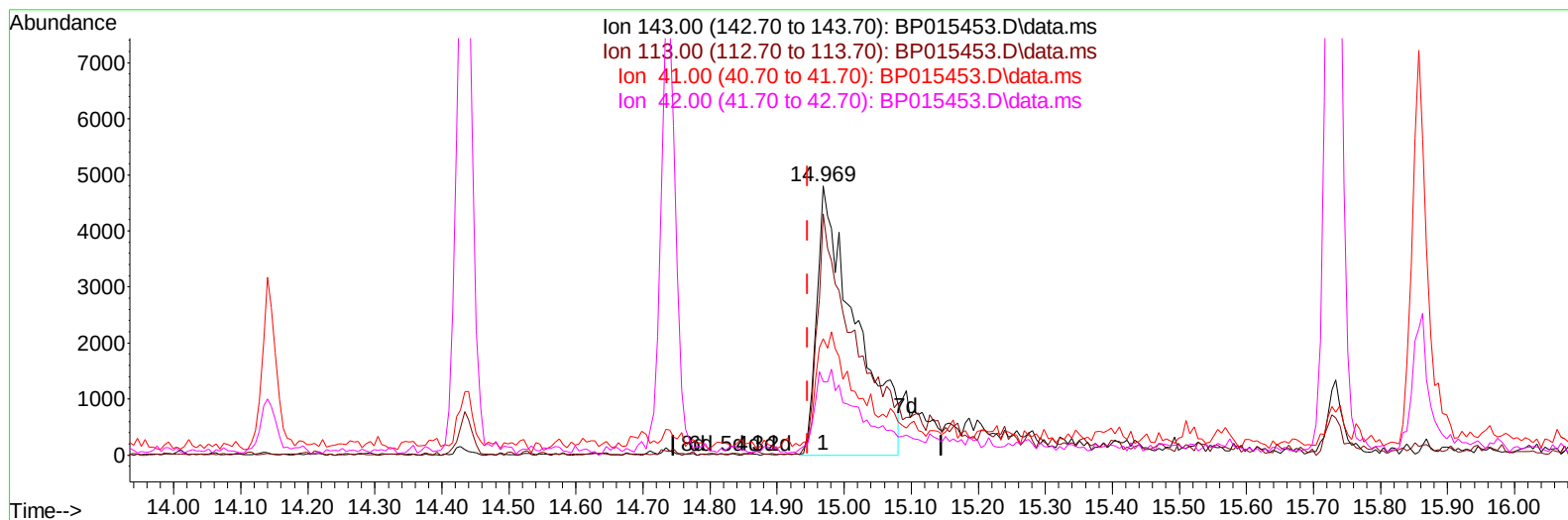
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(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.024) 4.41 ng/ul m

response 19137

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	89.81
41.00	42.20	43.33
42.00	29.40	27.24

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	92262	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.922	136	429434	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.740	164	307560	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.492	188	728094	20.000	ng/ul	-0.01
79) Chrysene-d12	21.575	240	740017	20.000	ng/ul	-0.01
88) Perylene-d12	24.127	264	760510	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	7867	3.157	ng/uL	0.00
4) Pyridine-d5	3.858	84	67461	10.424	ng/ul	0.00
7) Phenol-d5	7.252	99	71891	8.458	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	209482	41.267	ng/ul	0.00
11) 2-Chlorophenol-d4	7.622	132	196296	31.853	ng/ul	0.00
15) 4-Methylphenol-d8	8.810	113	146161	20.952	ng/ul	0.00
21) Nitrobenzene-d5	9.275	128	137846	40.886	ng/ul	0.00
24) 2-Nitrophenol-d4	9.999	143	140676	36.537	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	247058	35.176	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.063	131	337769	33.504	ng/ul	0.00
46) Dimethylphthalate-d6	14.146	166	1105376	45.577	ng/ul	0.00
49) Acenaphthylene-d8	14.434	160	1170846	44.150	ng/ul	0.00
54) 4-Nitrophenol-d4	14.969	143	19137m	4.406	ng/ul	0.02
60) Fluorene-d10	15.734	176	946758	46.292	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	126352	27.393	ng/ul	0.00
73) Anthracene-d10	17.592	188	1693122	51.110	ng/ul	0.00
81) Pyrene-d10	19.816	212	2096558	52.934	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.963	264	1987510	52.085	ng/ul	-0.01

Target Compounds Qvalue

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

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