

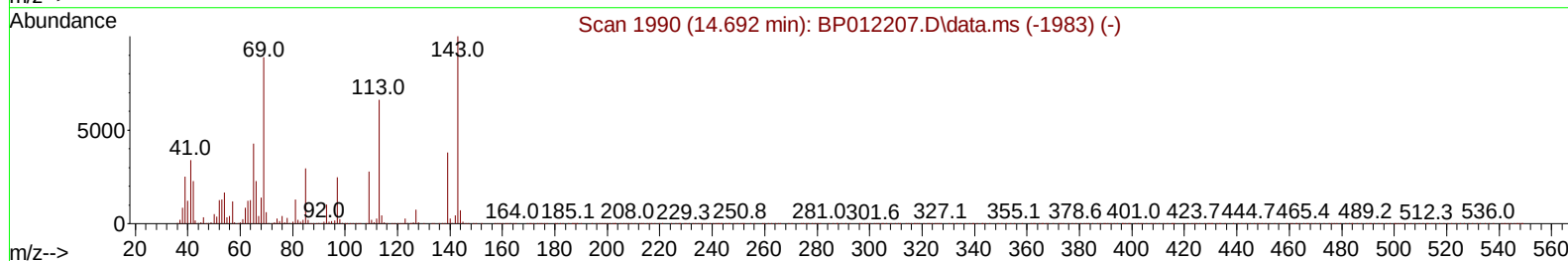
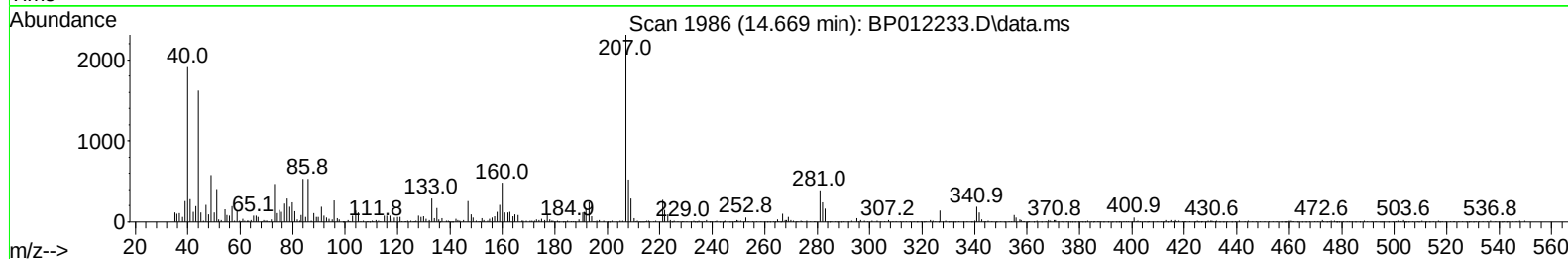
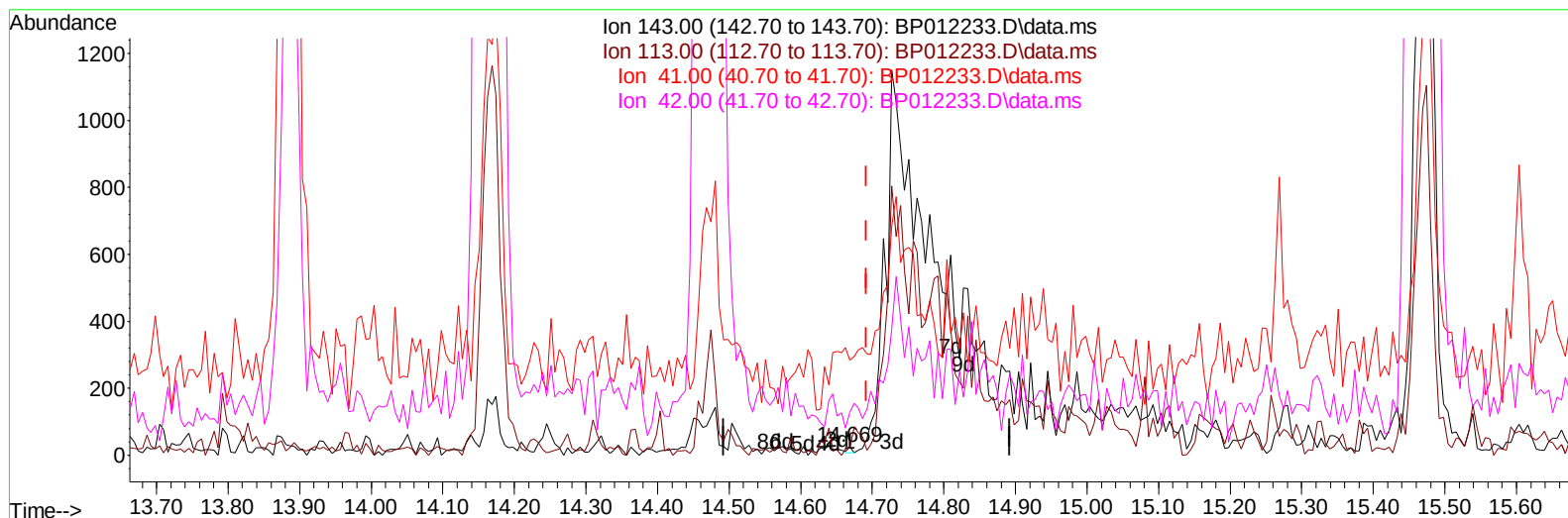
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012233.D
 Acq On : 21 Oct 2022 02:13
 Operator : CG/JU
 Sample : N5144-09
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COBT0

Manual IntegrationsAPPROVED

Quant Time: Oct 21 02:54:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/21/2022
 Supervised By :mohammad ahmed 10/24/2022



TIC: BP012233.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.024) 0.00 ng/u1

response 6

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.20	88.00#
41.00	35.10	1132.00#
42.00	23.90	516.00#

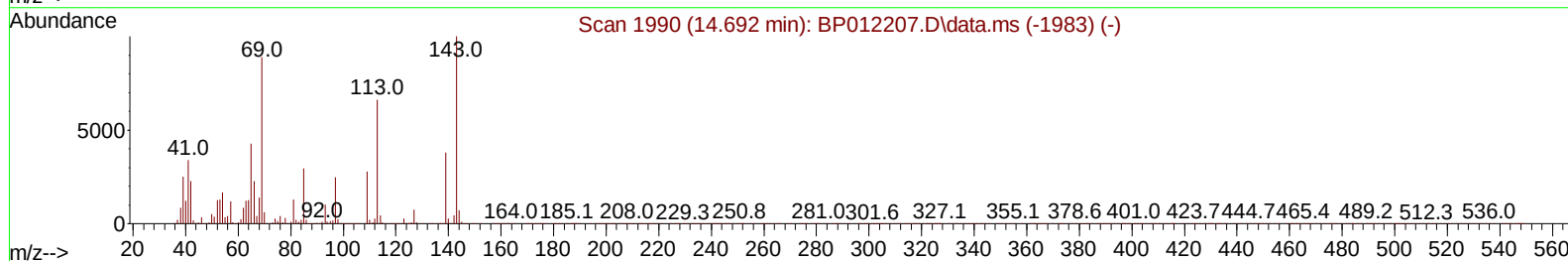
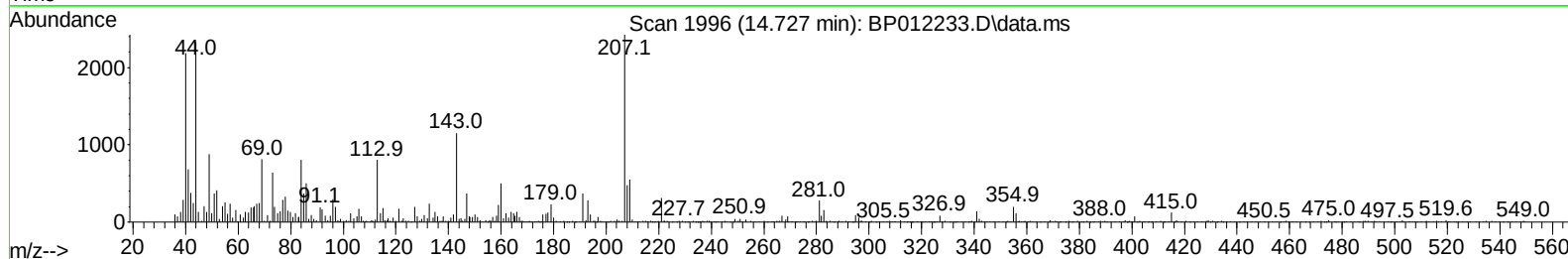
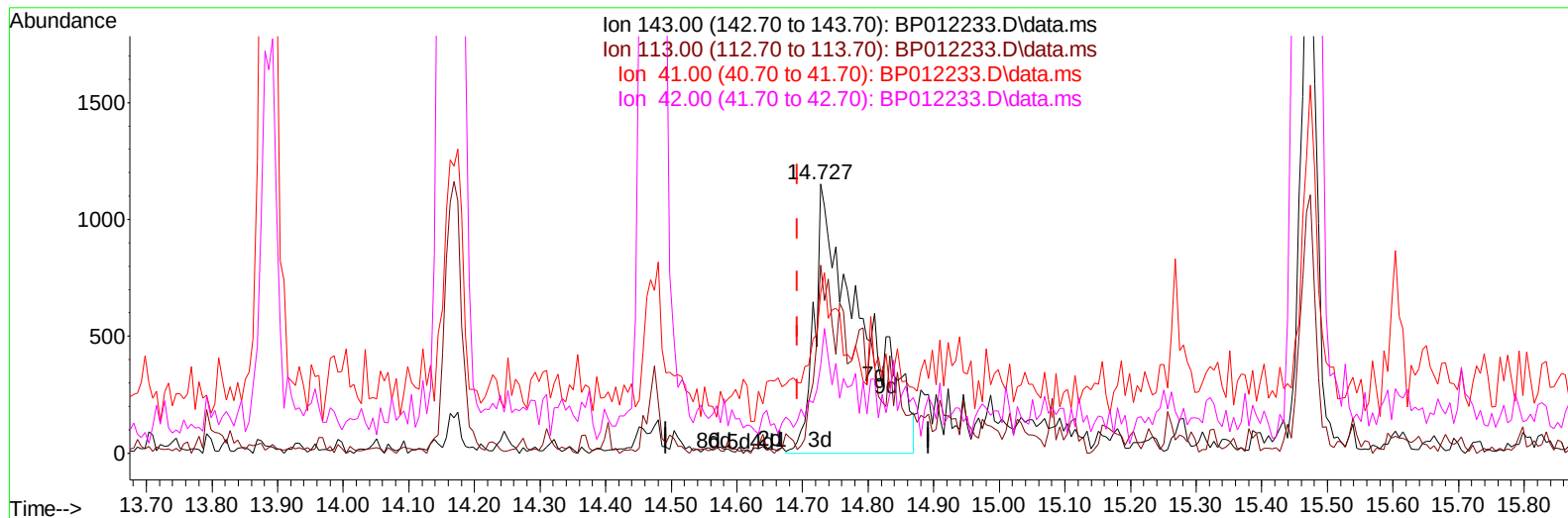
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(54) 4-Nitrophenol-d4 (S)

14.727min (+ 0.035) 0.37 ng/ul m

response 5621

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.20	69.85
41.00	35.10	59.51#
42.00	23.90	32.75#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	428543	20.000	ng/ul	0.00
20) Naphthalene-d8	10.628	136	1876478	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.474	164	1165444	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188	2343215	20.000	ng/ul	0.00
79) Chrysene-d12	21.321	240	1752464	20.000	ng/ul	-0.01
88) Perylene-d12	23.727	264	1705754	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.257	96	40690	3.843	ng/uL	0.00
4) Pyridine-d5	3.681	84	393976	12.912	ng/ul	0.00
7) Phenol-d5	6.993	99	204597	5.530	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	675281	30.097	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	640608	22.748	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	405434	14.100	ng/ul	-0.01
21) Nitrobenzene-d5	8.992	128	457004	32.902	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	338398	22.581	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	715808	25.465	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	972593	24.352	ng/ul	-0.01
46) Dimethylphthalate-d6	13.886	166	2632802	32.476	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	2984143	32.248	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.727	143	5621m	0.370	ng/ul	0.04
60) Fluorene-d10	15.469	176	2303570	33.028	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	8269	0.620	ng/ul	0.00
73) Anthracene-d10	17.333	188	3850391	37.488	ng/ul	0.00
81) Pyrene-d10	19.574	212	3880906	38.896	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.574	264	3228229	37.739	ng/ul	-0.01

Target Compounds Qvalue

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

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