

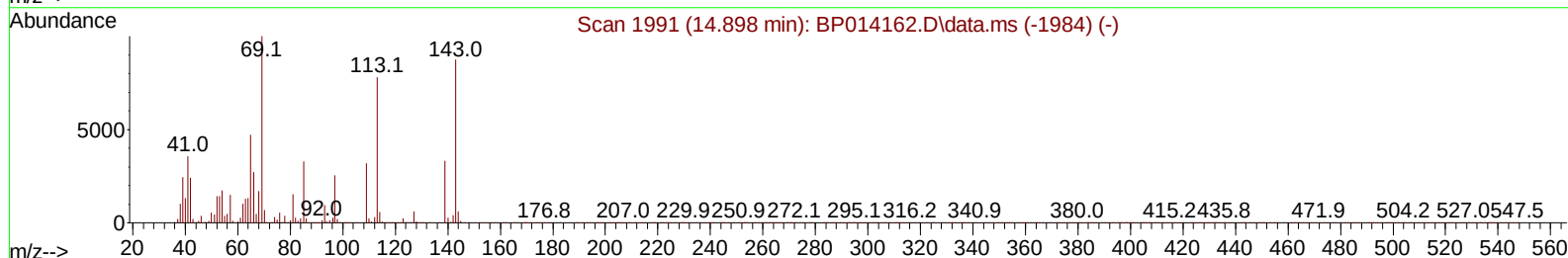
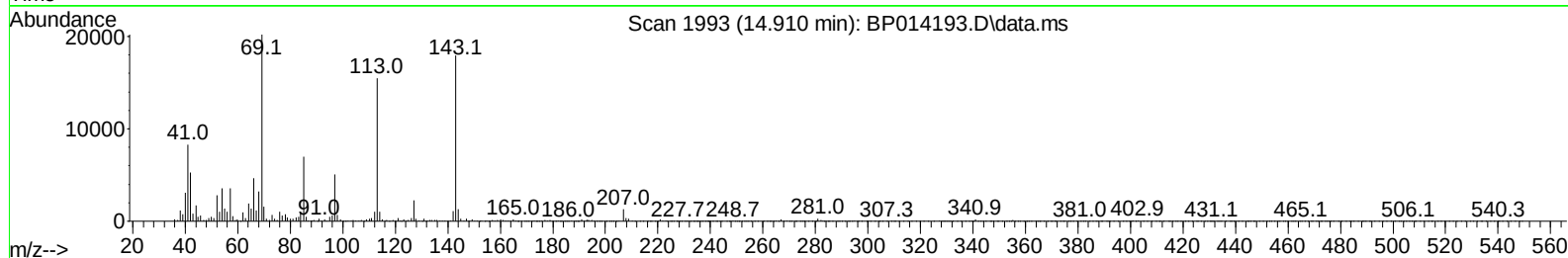
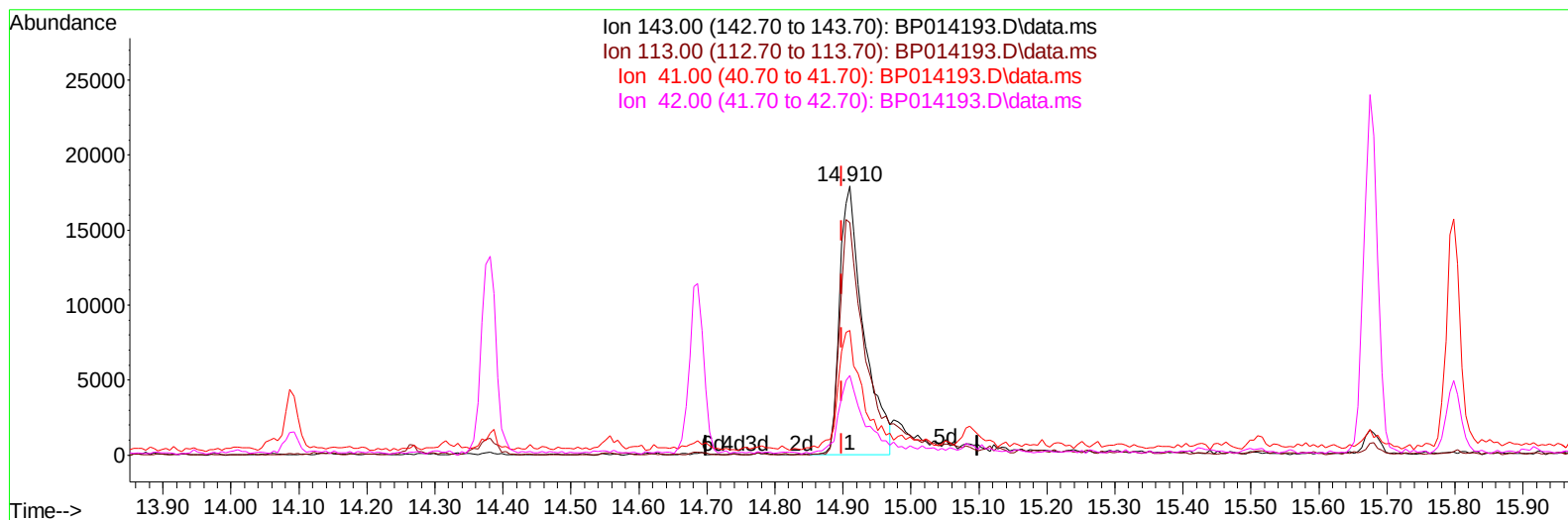
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032023\
 Data File : BP014193.D
 Acq On : 21 Mar 2023 22:19
 Operator : CG/JU
 Sample : 01949-03
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB933

Manual IntegrationsAPPROVED

Quant Time: Mar 22 01:22:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 21 04:47:07 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/22/2023
 Supervised By :mohammad ahmed 03/24/2023



TIC: BP014193.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.012) 5.16 ng/ul

response 43505

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	86.53
41.00	40.20	46.25
42.00	28.50	29.59

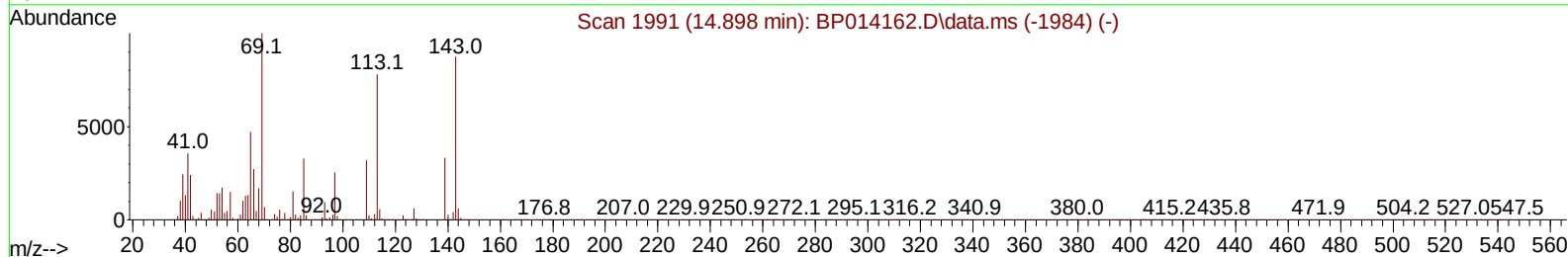
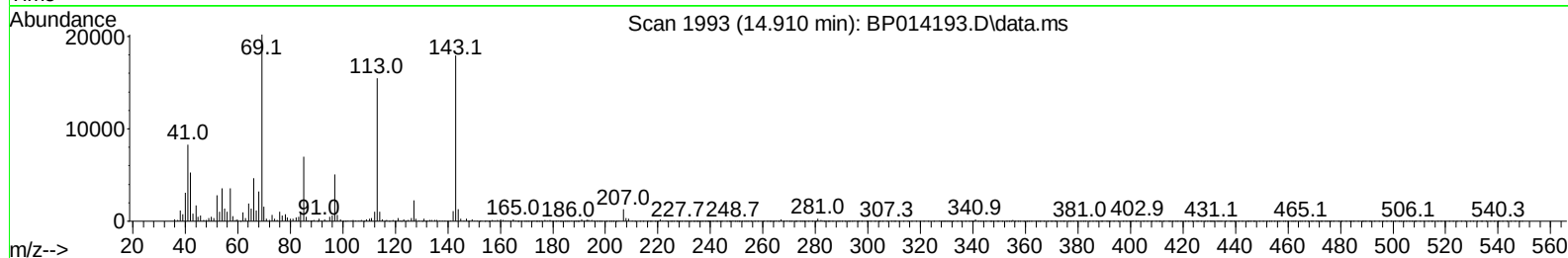
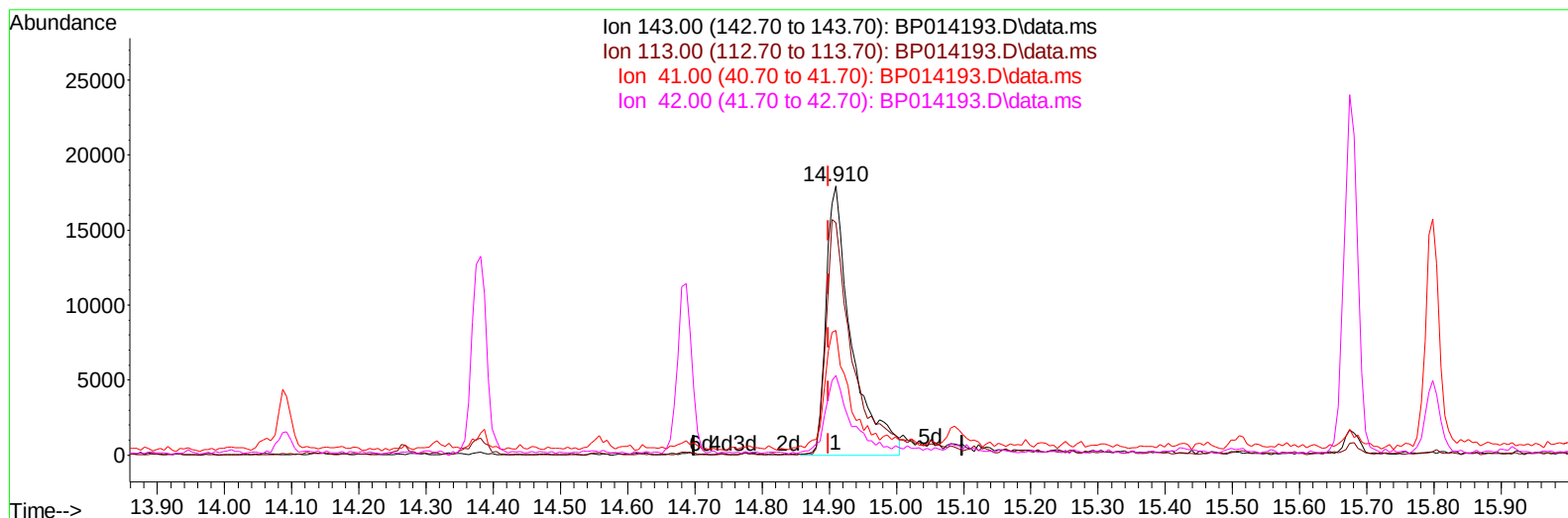
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(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.012) 5.64 ng/ul m

response 47521

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	86.53
41.00	40.20	46.25
42.00	28.50	29.59

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.046	152	191484	20.000	ng/ul	0.00
20) Naphthalene-d8	10.869	136	824846	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.687	164	563491	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	1291970	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.516	240	1146819	20.000	ng/ul	#-0.01
88) Perylene-d12	24.033	264	1029287	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	21372	4.217	ng/uL	0.00
4) Pyridine-d5	3.852	84	104199	7.426	ng/ul	0.00
7) Phenol-d5	7.205	99	126746	7.501	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.369	67	323963	32.873	ng/ul	0.00
11) 2-Chlorophenol-d4	7.575	132	321944	25.007	ng/ul	0.00
15) 4-Methylphenol-d8	8.757	113	240900	17.453	ng/ul	0.00
21) Nitrobenzene-d5	9.222	128	203069	30.855	ng/ul	0.00
24) 2-Nitrophenol-d4	9.946	143	239597	30.412	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.487	165	416819	28.118	ng/ul	0.00
31) 4-Chloroaniline-d4	11.010	131	466221	23.766	ng/ul	0.00
46) Dimethylphthalate-d6	14.092	166	1619905	33.874	ng/ul	0.00
49) Acenaphthylene-d8	14.381	160	1667876	32.923	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	47521m	5.639	ng/ul	0.01
60) Fluorene-d10	15.675	176	1364330	33.663	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	285043	29.191	ng/ul	0.00
73) Anthracene-d10	17.539	188	2298422	36.646	ng/ul	0.00
81) Pyrene-d10	19.763	212	2698220	42.659	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	2041464	37.287	ng/ul	0.00

Target Compounds Qvalue

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

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