

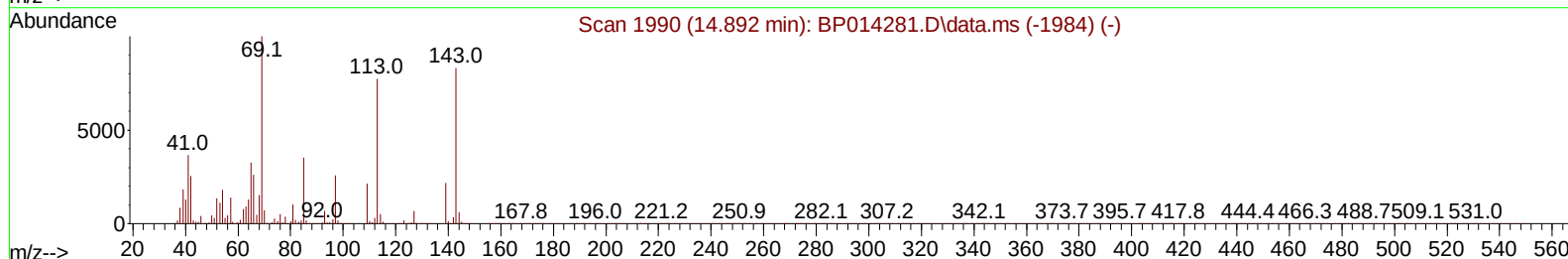
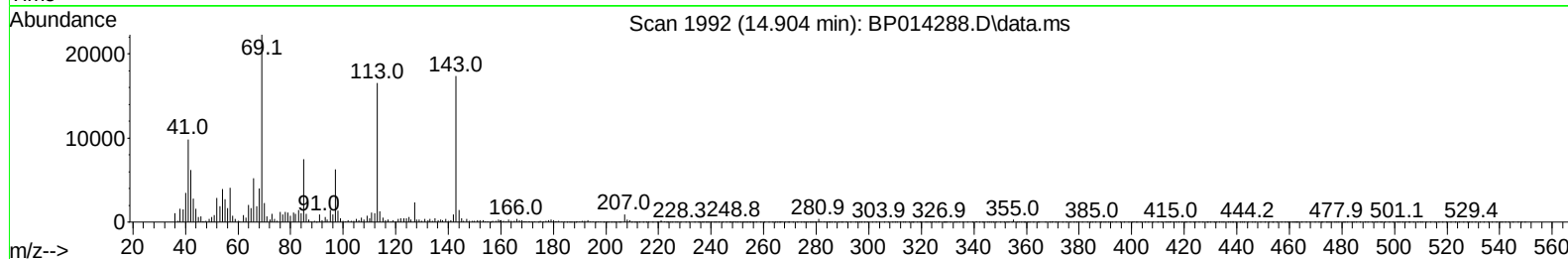
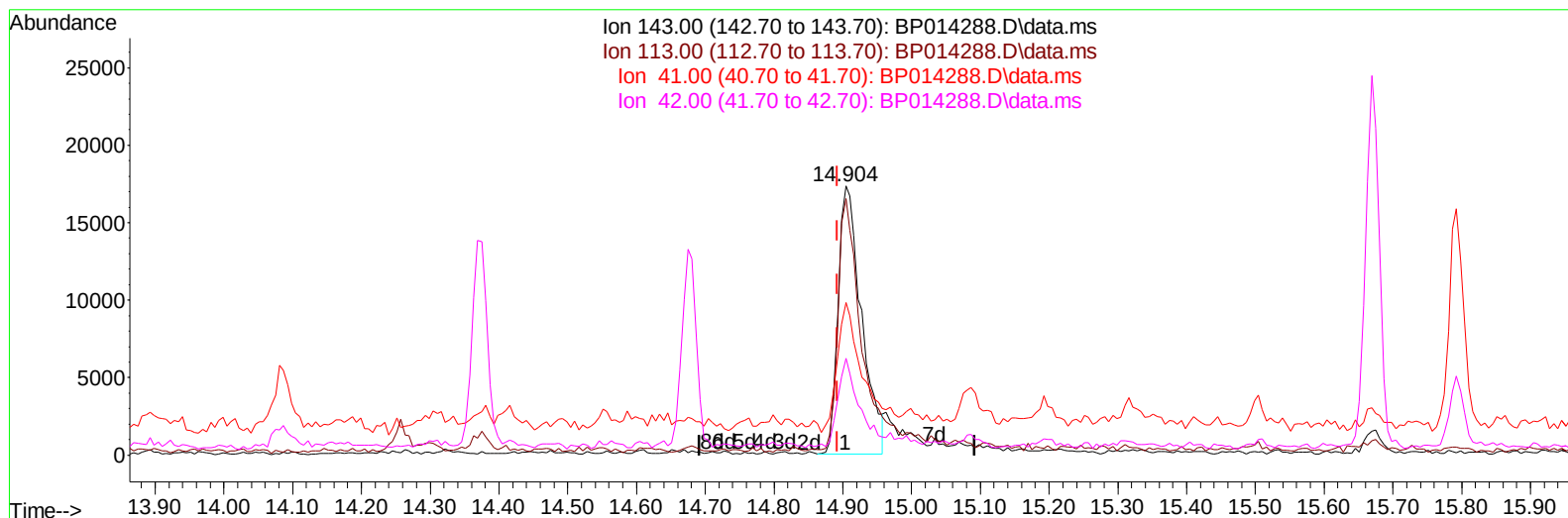
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032423\
 Data File : BP014288.D
 Acq On : 24 Mar 2023 13:11
 Operator : CG/JU
 Sample : 02037-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E45F2

Manual IntegrationsAPPROVED

Quant Time: Mar 24 22:35:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 24 22:30:41 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/27/2023
 Supervised By :Jagrut Upadhyay 03/27/2023



TIC: BP014288.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (+ 0.012) 4.76 ng/ul

response 41078

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	95.35
41.00	40.20	56.75#
42.00	28.50	35.89#

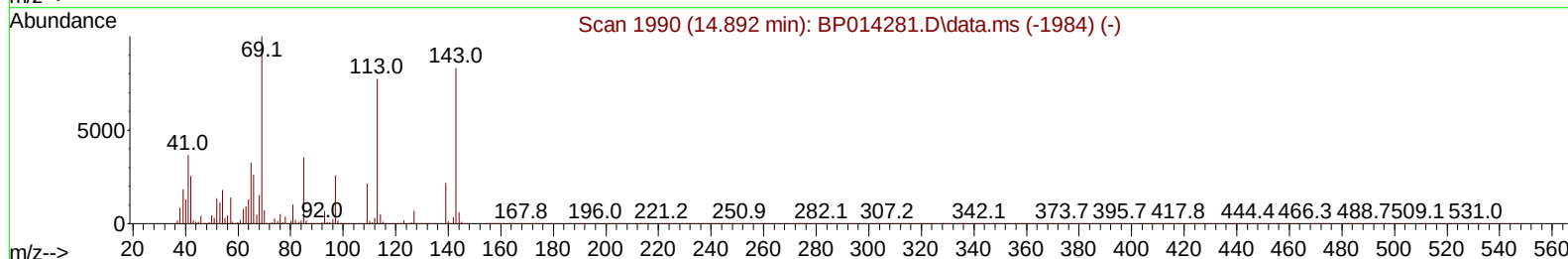
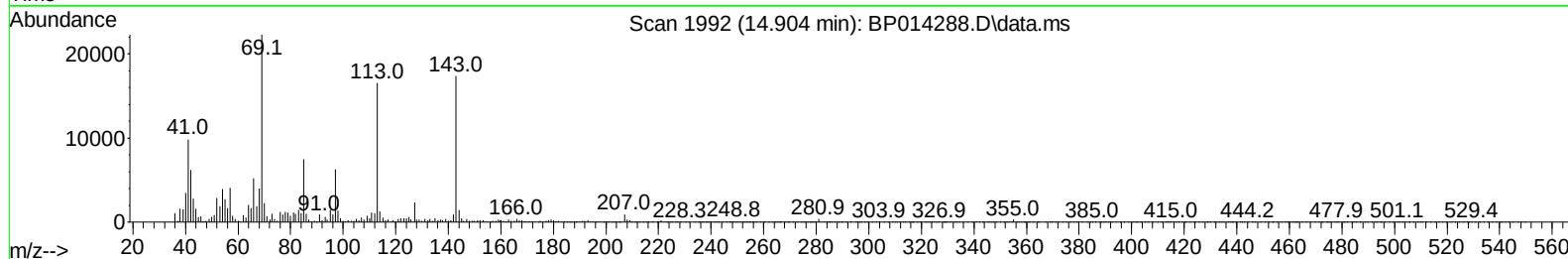
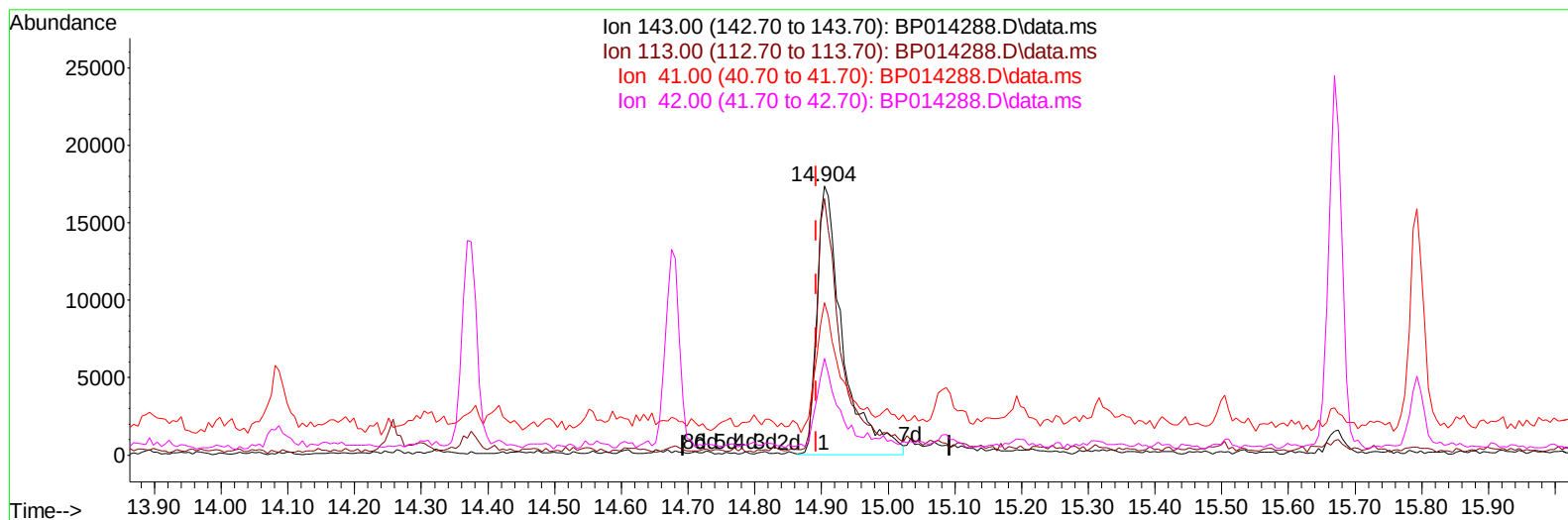
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TIC: BP014288.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (+ 0.012) 5.42 ng/ul m

response 46774

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	95.35
41.00	40.20	56.75#
42.00	28.50	35.89#

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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	8.034	152	182625	20.000	ng/ul	0.00
20) Naphthalene-d8	10.852	136	829744	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.675	164	576543	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.434	188	1332154	20.000	ng/ul	#-0.02
79) Chrysene-d12	21.516	240	1134187	20.000	ng/ul	#-0.02
88) Perylene-d12	24.039	264	1000629	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	25933	5.365	ng/uL	0.00
4) Pyridine-d5	3.840	84	110063	8.224	ng/ul	0.00
7) Phenol-d5	7.193	99	127284	7.898	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.358	67	312252	33.222	ng/ul	0.00
11) 2-Chlorophenol-d4	7.558	132	285349	23.240	ng/ul	-0.01
15) 4-Methylphenol-d8	8.746	113	251249	19.086	ng/ul	-0.01
21) Nitrobenzene-d5	9.205	128	194315	29.351	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.934	143	216553	27.325	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.475	165	389833	26.143	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.999	131	457419	23.180	ng/ul	-0.01
46) Dimethylphthalate-d6	14.081	166	1527335	31.215	ng/ul	-0.01
49) Acenaphthylene-d8	14.375	160	1632157	31.489	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	46774m	5.425	ng/ul	0.01
60) Fluorene-d10	15.669	176	1273273	30.705	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.793	200	240449	23.881	ng/ul	-0.01
73) Anthracene-d10	17.534	188	2105095	32.551	ng/ul	-0.01
81) Pyrene-d10	19.763	212	2426170	38.785	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.874	264	1748242	32.846	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.434	88	254855	48.841	ng/uL	91

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

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