

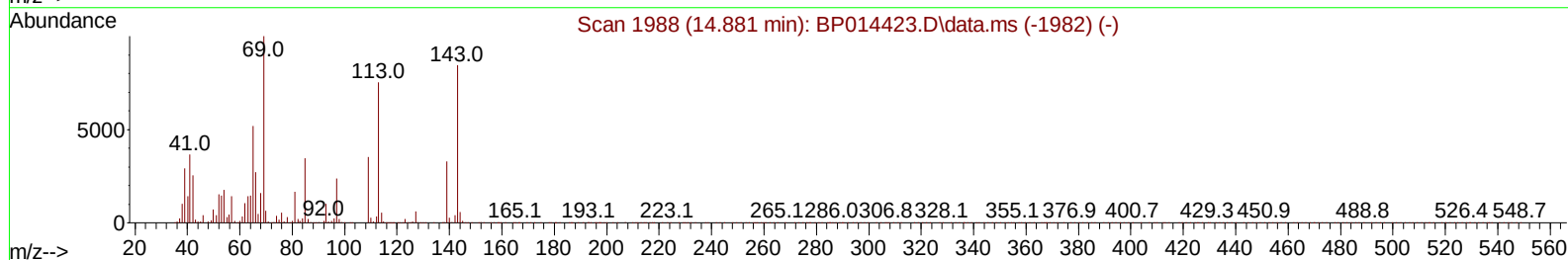
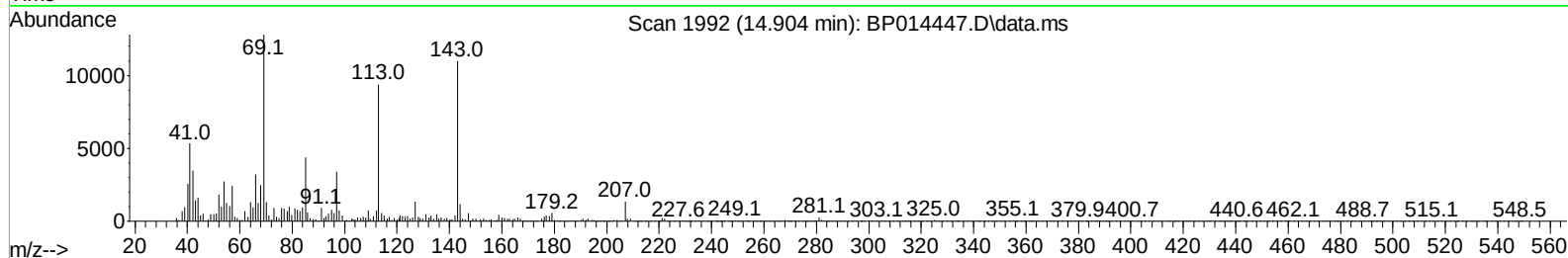
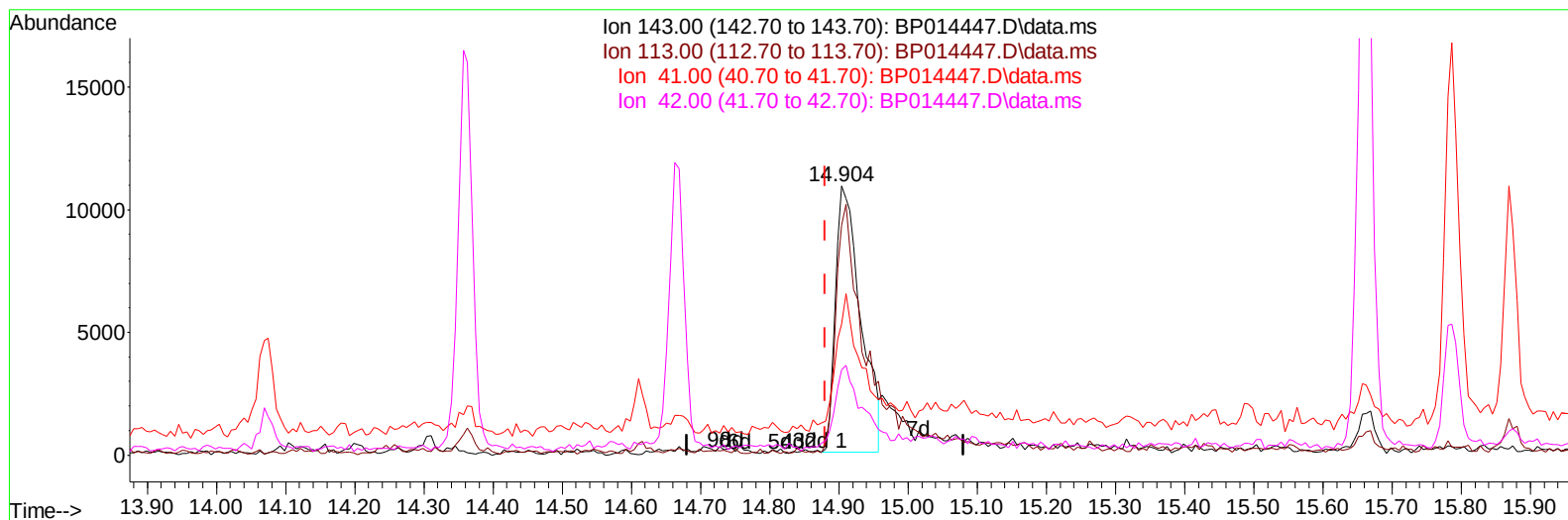
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014447.D
Acq On : 31 Mar 2023 17:06
Operator : CG/JU
Sample : 02093-01
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB964

Manual IntegrationsAPPROVED

Quant Time: Mar 31 22:54:55 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 31 04:41:29 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/03/2023
Supervised By :Jagrut Upadhyay 04/03/2023



TIC: BP014447.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (+ 0.023) 4.70 ng/ul

response 27472

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	85.54
41.00	44.20	48.65
42.00	31.50	31.84

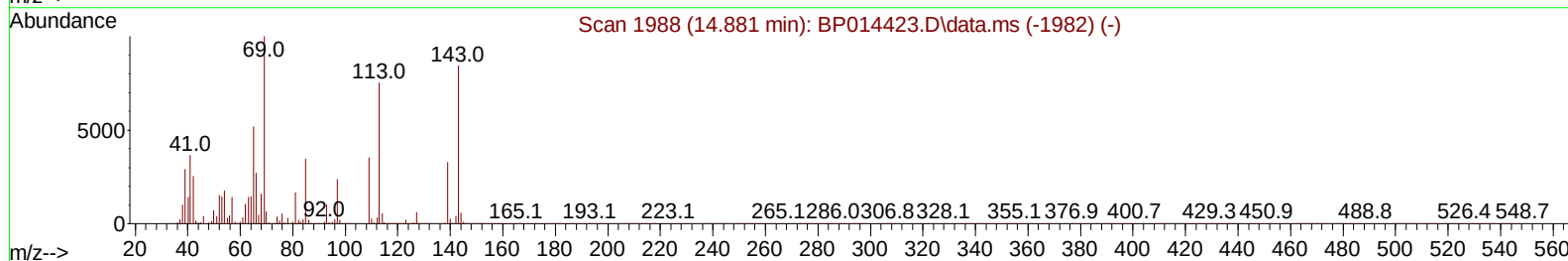
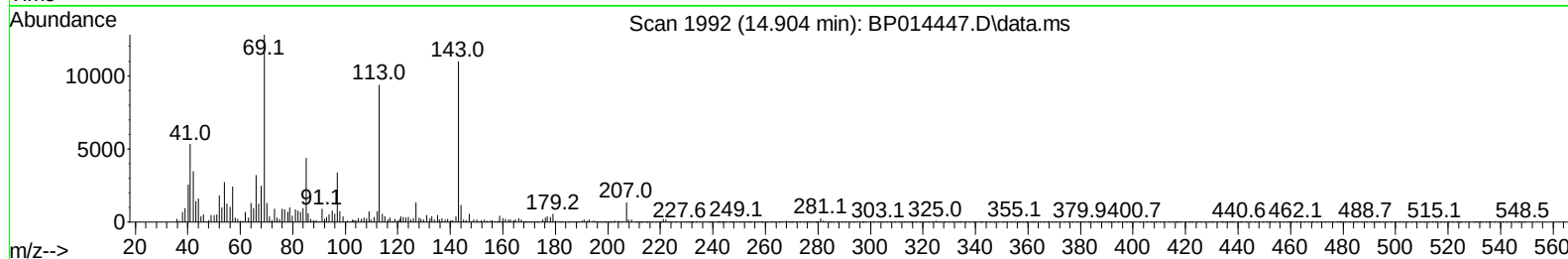
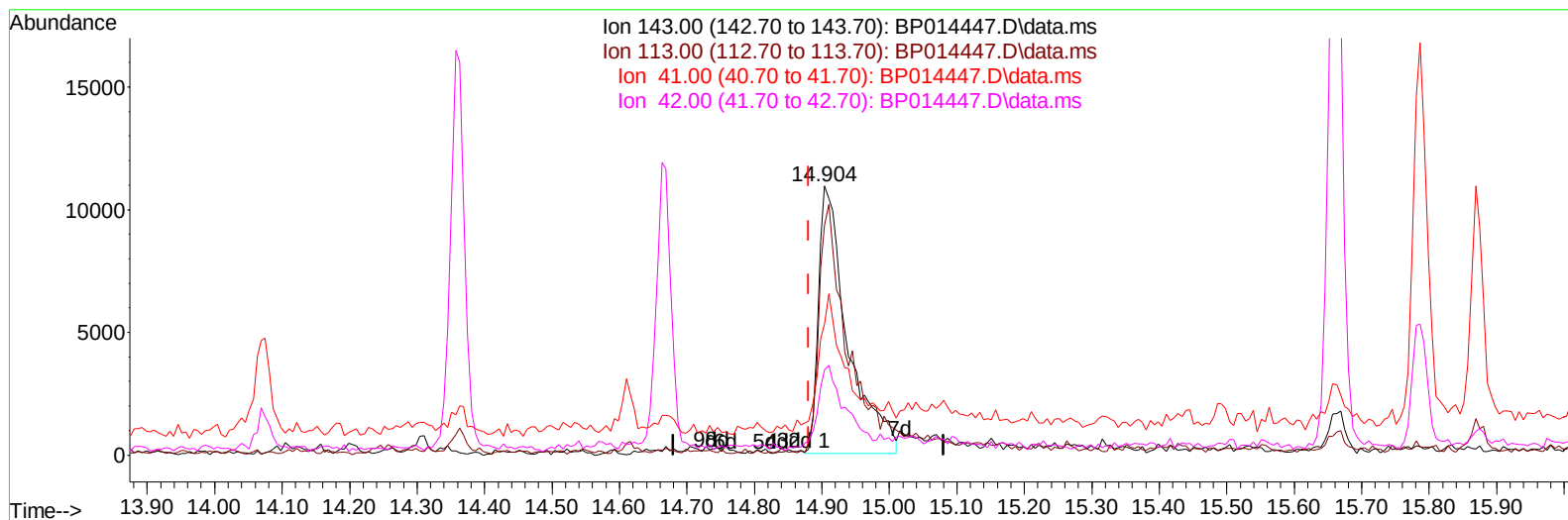
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TIC: BP014447.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (+ 0.023) 5.56 ng/ul m

response 32460

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	85.54
41.00	44.20	48.65
42.00	31.50	31.84

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	153912	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	710336	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	476731	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1021184	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	680505	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	642688	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	17038	4.293	ng/uL	0.00
4) Pyridine-d5	3.828	84	97604	9.596	ng/ul	0.01
7) Phenol-d5	7.175	99	108330	7.632	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	336284	36.925	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	290625	28.328	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	232176	20.278	ng/ul	0.00
21) Nitrobenzene-d5	9.187	128	203229	35.440	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	227636	34.667	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	403691	33.758	ng/ul	0.00
31) 4-Chloroaniline-d4	10.987	131	201484	12.791	ng/ul	0.00
46) Dimethylphthalate-d6	14.075	166	1600703	41.411	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1703223	39.711	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	32460m	5.559	ng/ul	0.02
60) Fluorene-d10	15.663	176	1315611	40.813	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	257308	34.244	ng/ul	0.00
73) Anthracene-d10	17.528	188	1998439	40.437	ng/ul	0.00
81) Pyrene-d10	19.757	212	2106362	46.082	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	1429240	41.857	ng/ul	0.00

Target Compounds Qvalue

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

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