

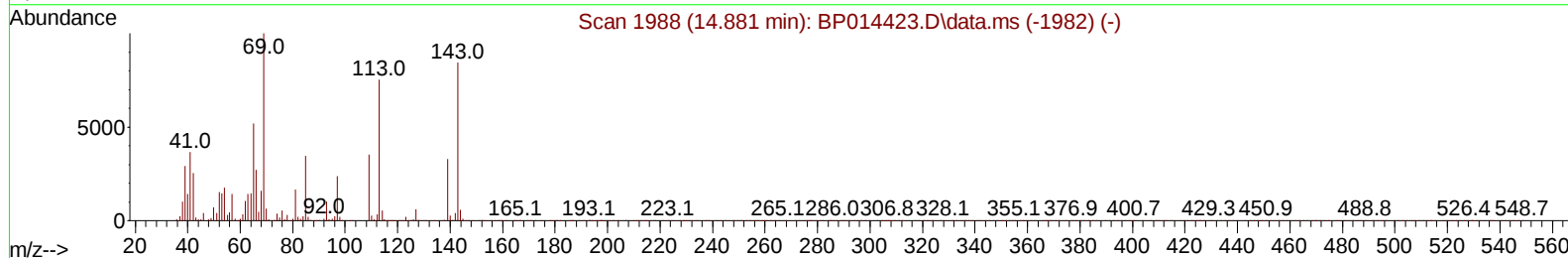
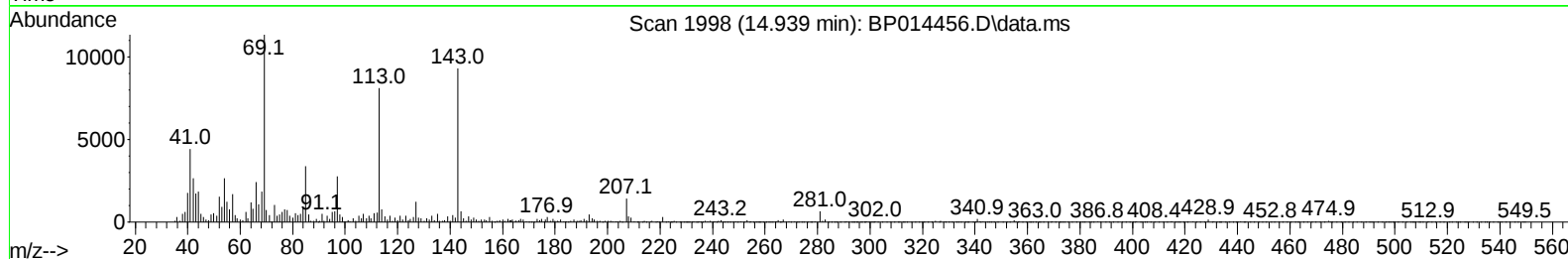
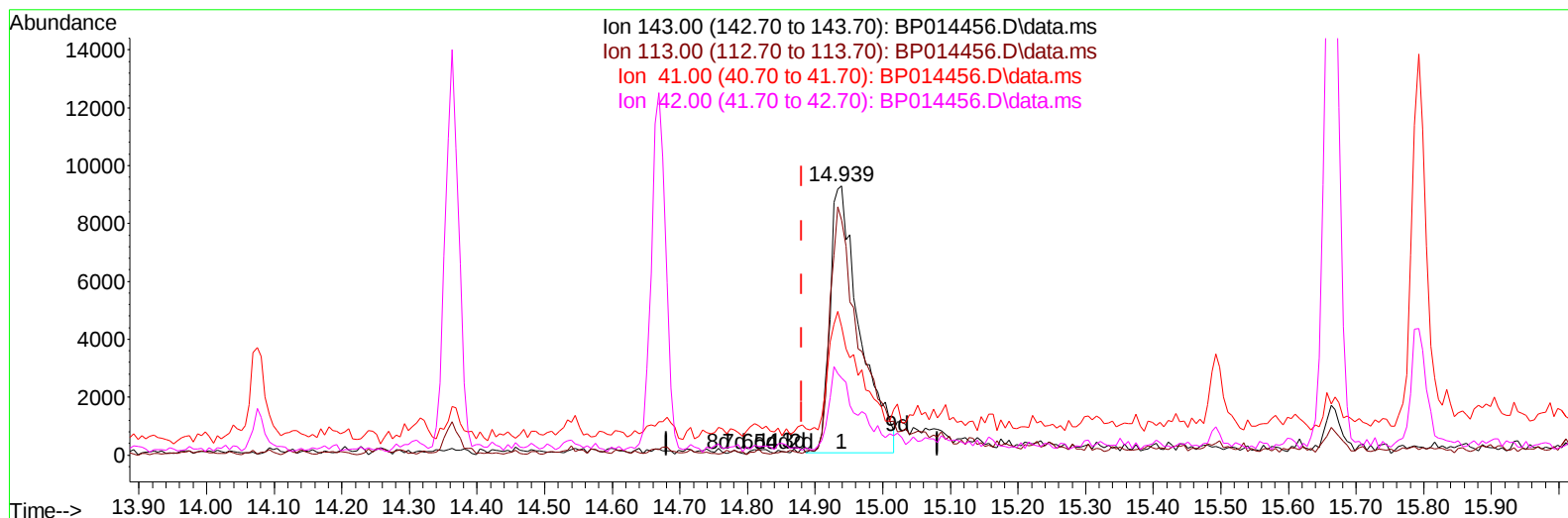
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014456.D
Acq On : 31 Mar 2023 22:10
Operator : CG/JU
Sample : 02092-10
Misc :
ALS Vial : 65 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB975

Manual IntegrationsAPPROVED

Quant Time: Mar 31 22:56:35 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: BP014456.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.939min (+ 0.059) 4.26 ng/u1

response 28088

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	87.35
41.00	44.20	47.70
42.00	31.50	28.73

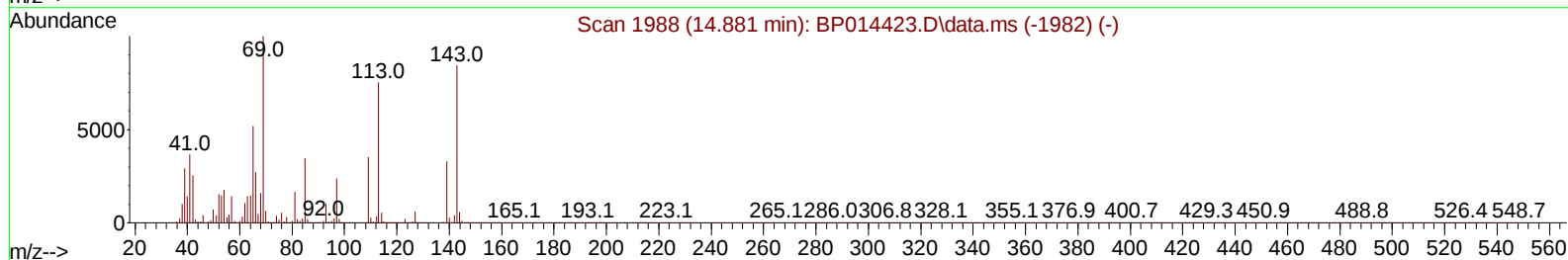
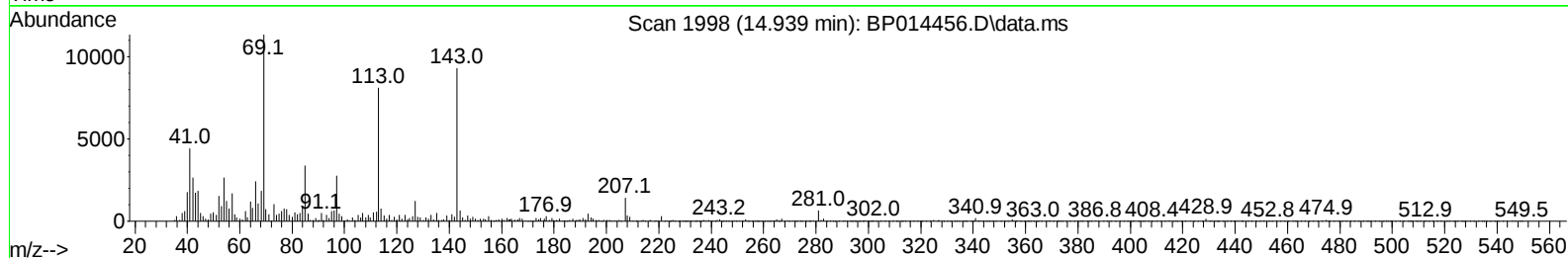
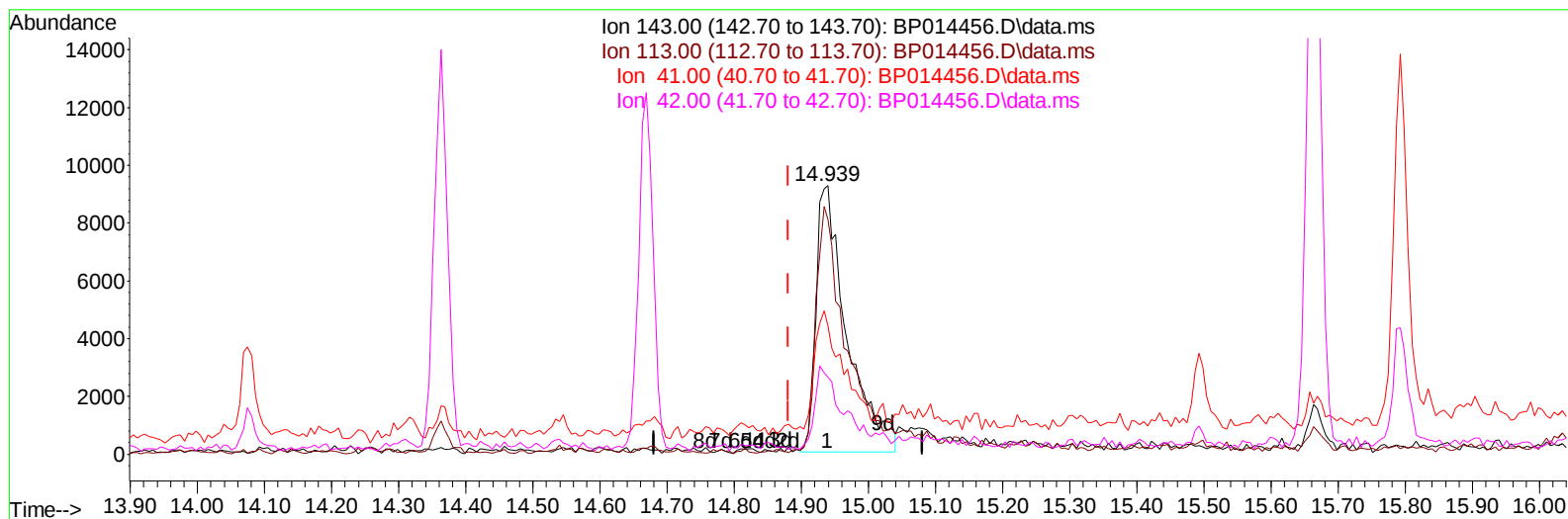
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TIC: BP014456.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.939min (+ 0.059) 4.44 ng/ul m

response 29281

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	87.35
41.00	44.20	47.70
42.00	31.50	28.73

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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	8.010	152	164867	20.000	ng/ul	0.00
20) Naphthalene-d8	10.840	136	772447	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	537916	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.433	188	1194633	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	945085	20.000	ng/ul	0.00
88) Perylene-d12	24.039	264	870936	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	16854	3.964	ng/uL	0.00
4) Pyridine-d5	3.840	84	35447	3.254	ng/ul	0.02
7) Phenol-d5	7.193	99	81537	5.363	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.340	67	259480	26.598	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	226978	20.654	ng/ul	0.00
15) 4-Methylphenol-d8	8.746	113	176689	14.406	ng/ul	0.02
21) Nitrobenzene-d5	9.193	128	161518	25.902	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	183296	25.670	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.469	165	329404	25.331	ng/ul	0.01
31) 4-Chloroaniline-d4	10.993	131	200741	11.719	ng/ul	0.01
46) Dimethylphthalate-d6	14.075	166	1361946	31.226	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1386836	28.656	ng/ul	0.00
54) 4-Nitrophenol-d4	14.939	143	29281m	4.444	ng/ul	0.06
60) Fluorene-d10	15.663	176	1109994	30.518	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	218577	24.866	ng/ul	0.00
73) Anthracene-d10	17.533	188	1795041	31.048	ng/ul	0.00
81) Pyrene-d10	19.763	212	2056941	32.402	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1459184	31.535	ng/ul	0.00

Target Compounds Qvalue

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

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