

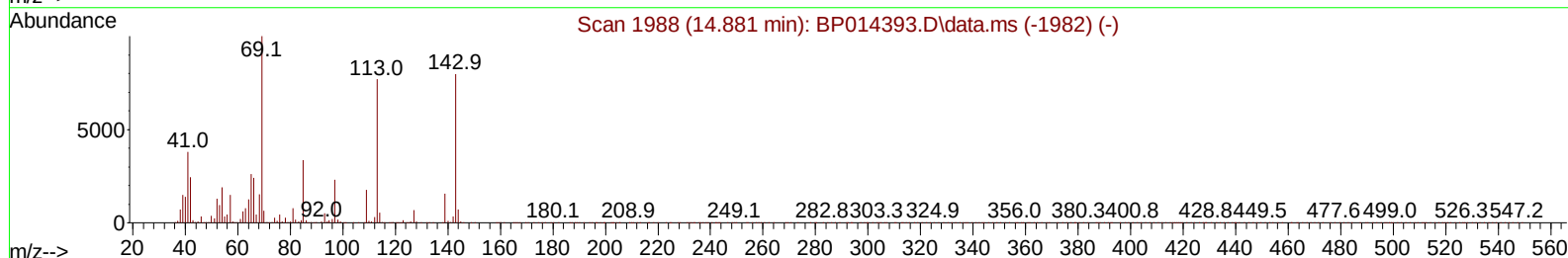
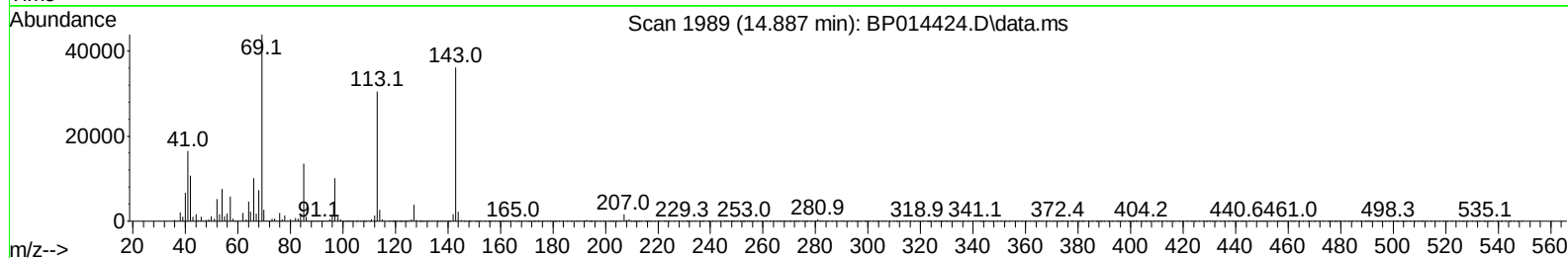
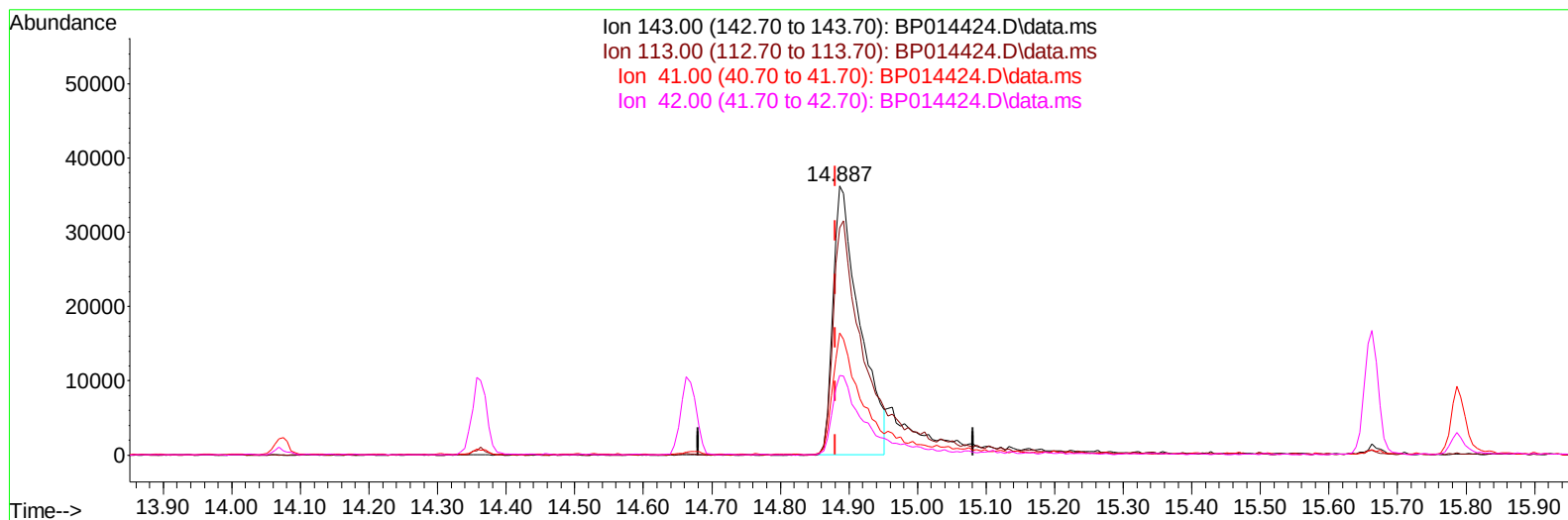
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014424.D
Acq On : 31 Mar 2023 04:11
Operator : CG/JU
Sample : PB151686BL
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK686

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/31/2023
Supervised By :Jagrut Upadhyay 03/31/2023

Quant Time: Mar 31 04:41:45 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



TIC: BP014424.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.887min (+ 0.006) 17.61 ng/ul

response 97265

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	84.42
41.00	44.20	45.49
42.00	31.50	29.60

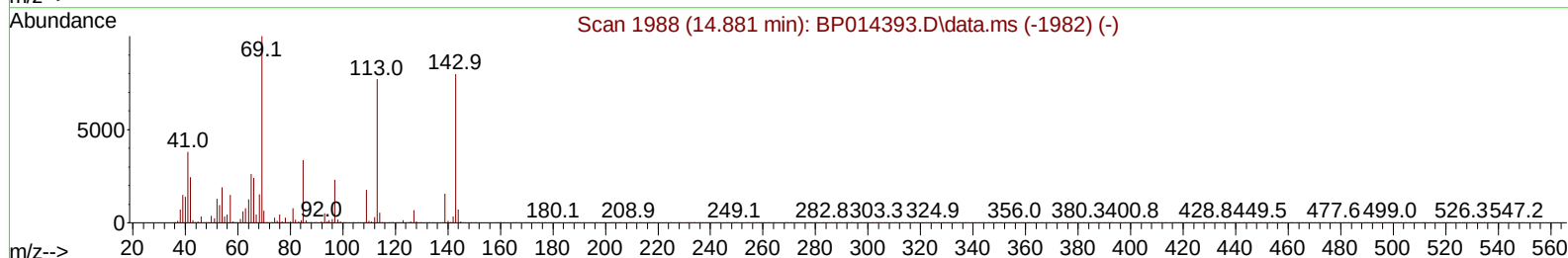
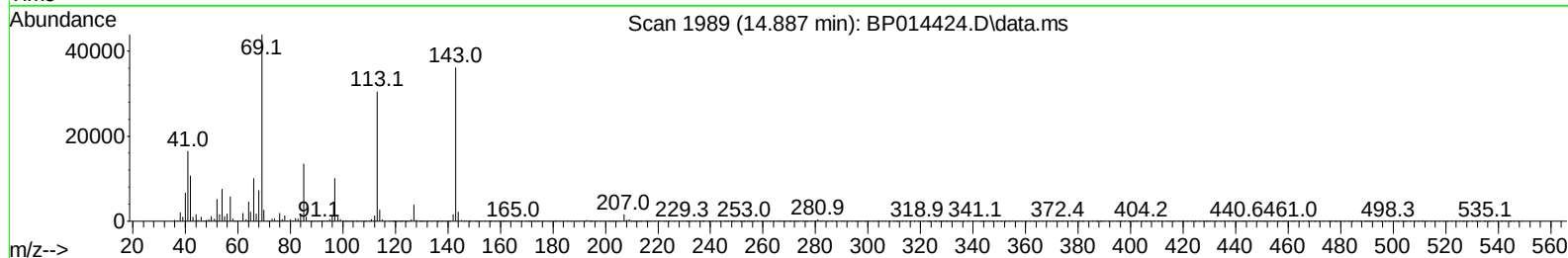
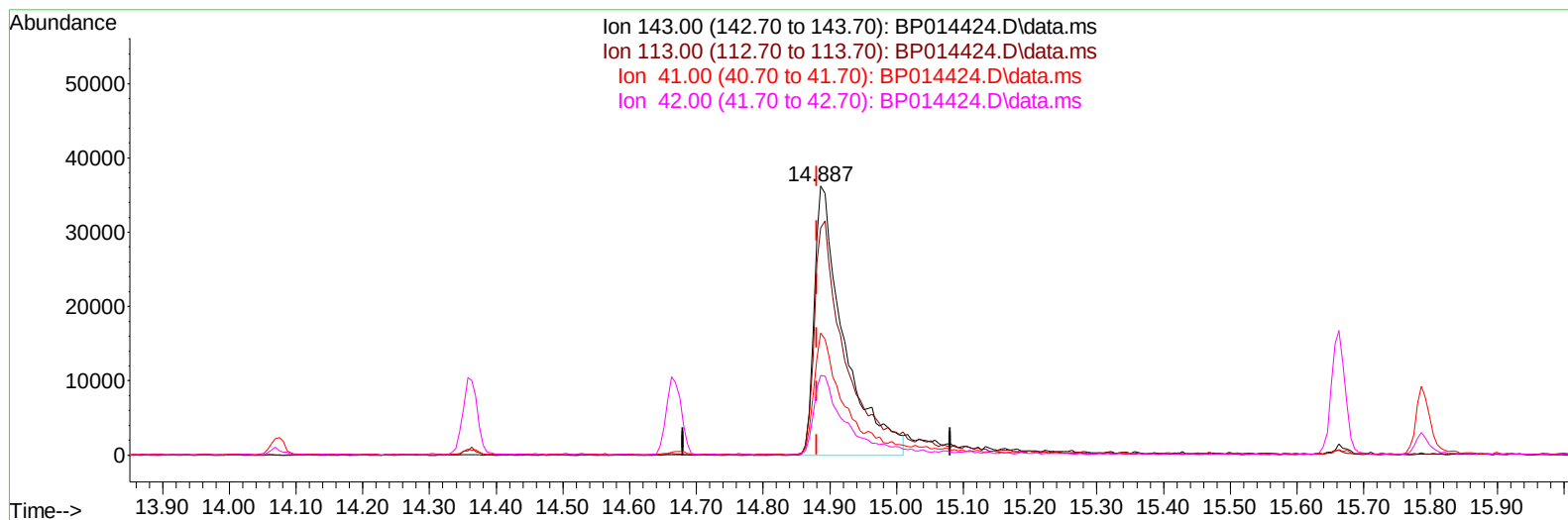
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(54) 4-Nitrophenol-d4 (S)

14.887min (+ 0.006) 20.20 ng/ul m

response 111618

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	84.42
41.00	44.20	45.49
42.00	31.50	29.60

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	172593	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	744018	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	451017	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	893839	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	806184	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	912879	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	24270	5.453	ng/uL	0.00
4) Pyridine-d5	3.817	84	273191	23.953	ng/ul	0.00
7) Phenol-d5	7.169	99	409193	25.707	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	281994	27.612	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	311092	27.041	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	328669	25.598	ng/ul	0.00
21) Nitrobenzene-d5	9.187	128	155966	25.967	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	178563	25.963	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	297039	23.715	ng/ul	0.00
31) 4-Chloroaniline-d4	10.987	131	276466	16.757	ng/ul	0.00
46) Dimethylphthalate-d6	14.075	166	987724	27.010	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1074384	26.477	ng/ul	0.00
54) 4-Nitrophenol-d4	14.887	143	111618m	20.204	ng/ul	0.00
60) Fluorene-d10	15.663	176	809559	26.546	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	155350	23.620	ng/ul	0.00
73) Anthracene-d10	17.528	188	1205485	27.867	ng/ul	0.00
81) Pyrene-d10	19.757	212	1362222	25.156	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	1400579	28.877	ng/ul	0.00

Target Compounds Qvalue

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

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