

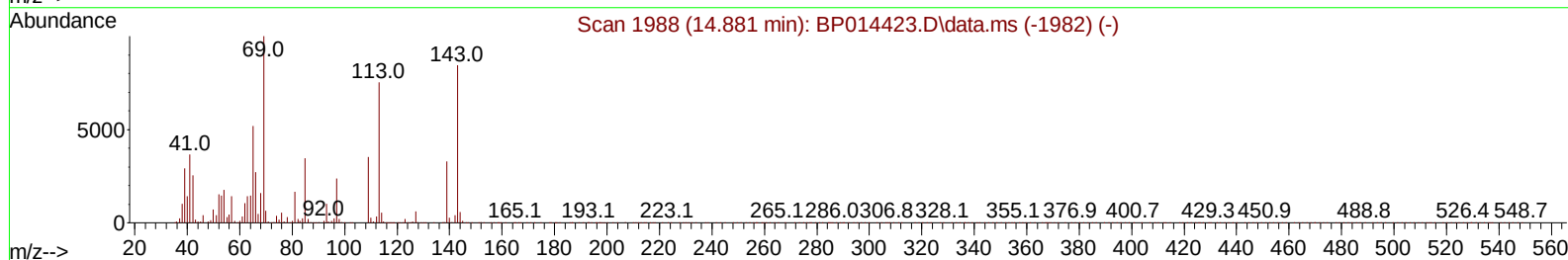
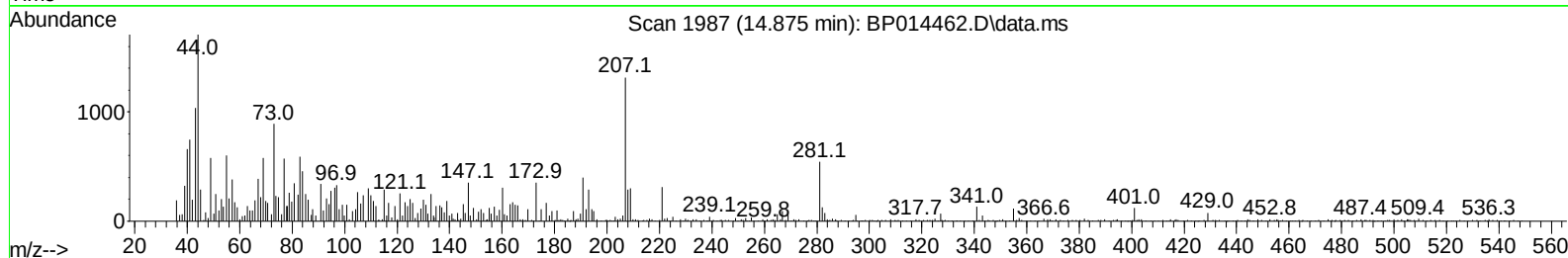
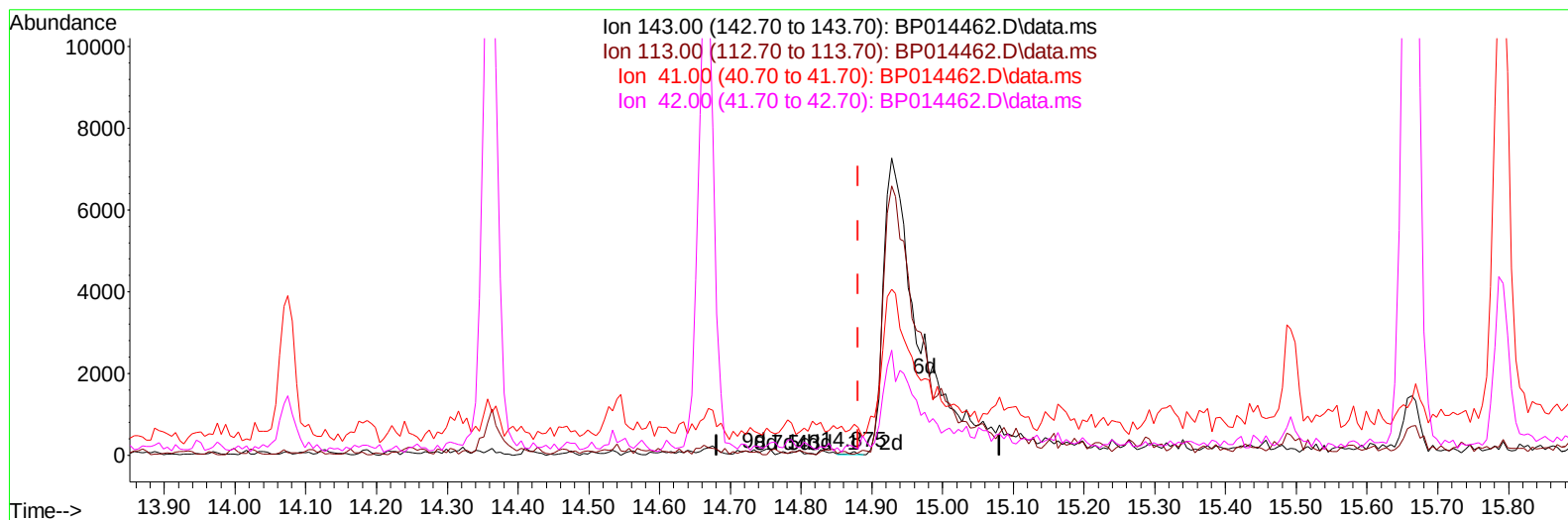
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014462.D
Acq On : 01 Apr 2023 02:06
Operator : CG/JU
Sample : 02093-06
Misc :
ALS Vial : 72 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB975

Manual IntegrationsAPPROVED

Quant Time: Apr 01 02:39:39 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: BP014462.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (-0.006) 0.01 ng/u1

response 48

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	25.30#
41.00	44.20	901.20#
42.00	31.50	239.76#

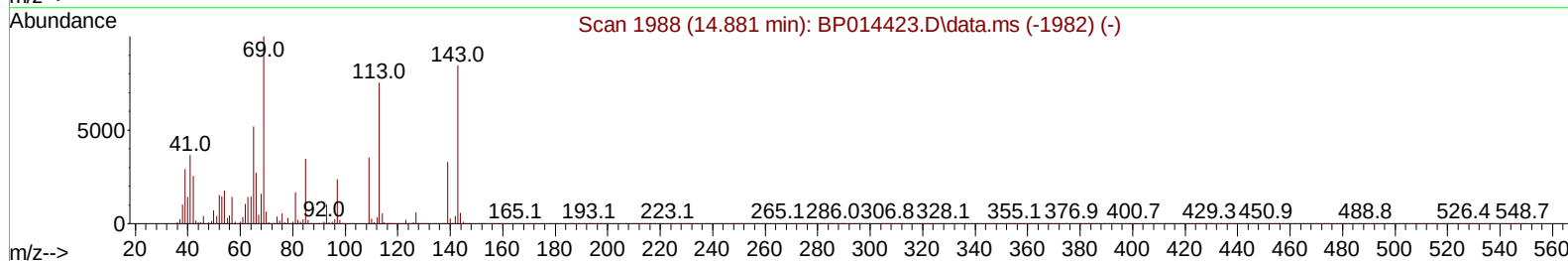
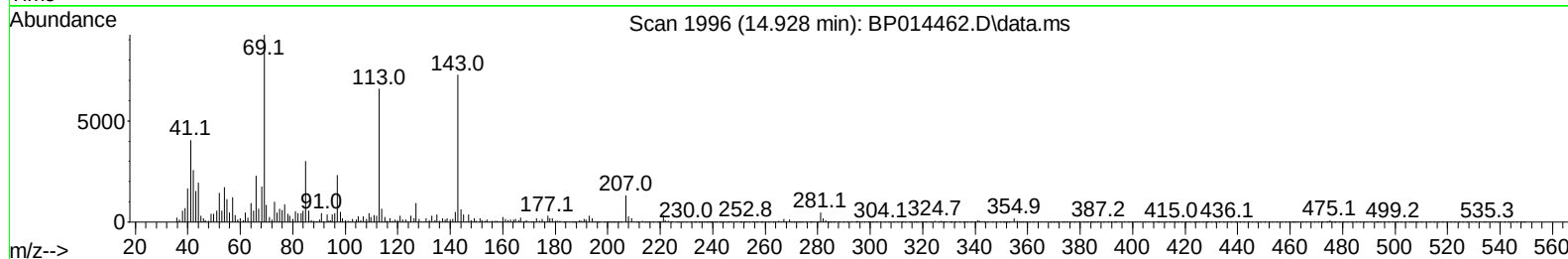
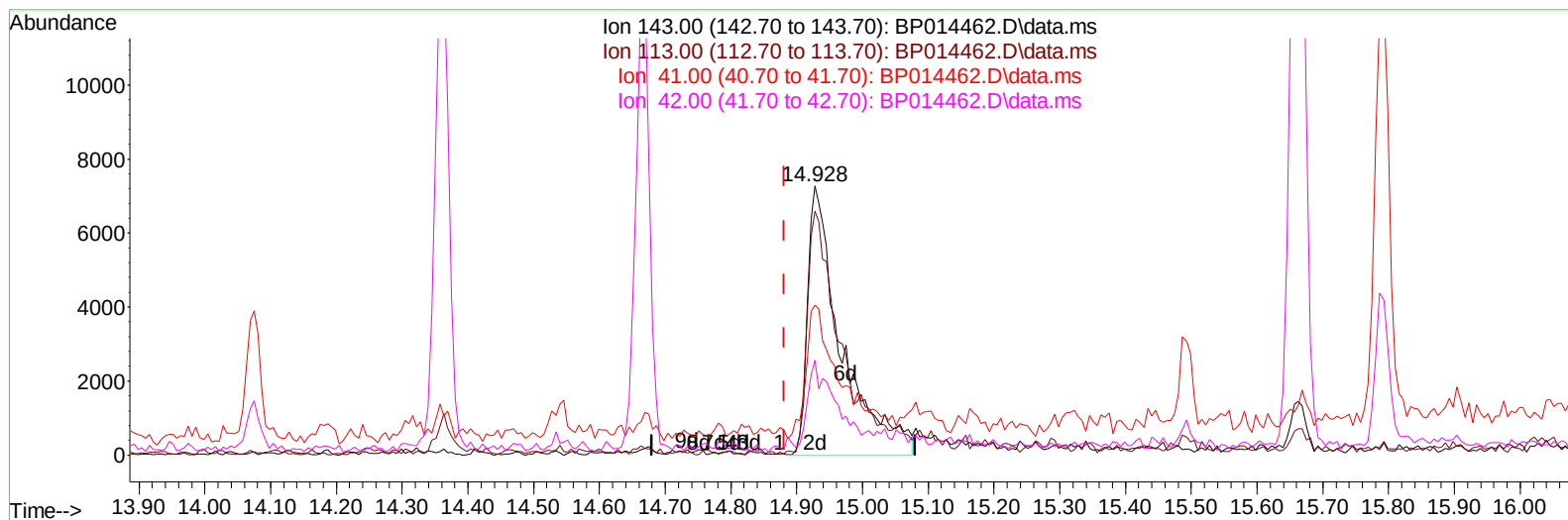
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(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.047) 4.04 ng/ul m

response 25980

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	90.60
41.00	44.20	55.83#
42.00	31.50	35.38

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	172626	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	780926	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	525481	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1112012	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	767768	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	818616	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	17703	3.977	ng/uL	0.00
4) Pyridine-d5	3.840	84	43416	3.806	ng/ul	0.02
7) Phenol-d5	7.187	99	89304	5.609	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.340	67	286914	28.089	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	252045	21.904	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	186626	14.533	ng/ul	0.00
21) Nitrobenzene-d5	9.193	128	172459	27.356	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	194184	26.900	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	351489	26.735	ng/ul	0.00
31) 4-Chloroaniline-d4	10.987	131	196757	11.362	ng/ul	0.00
46) Dimethylphthalate-d6	14.075	166	1416176	33.238	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1445853	30.583	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	25980m	4.036	ng/ul	0.05
60) Fluorene-d10	15.663	176	1139503	32.071	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	200202	24.468	ng/ul	0.00
73) Anthracene-d10	17.528	188	1754605	32.603	ng/ul	0.00
81) Pyrene-d10	19.757	212	1841416	35.707	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1456529	33.489	ng/ul	0.00

Target Compounds Qvalue

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

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