

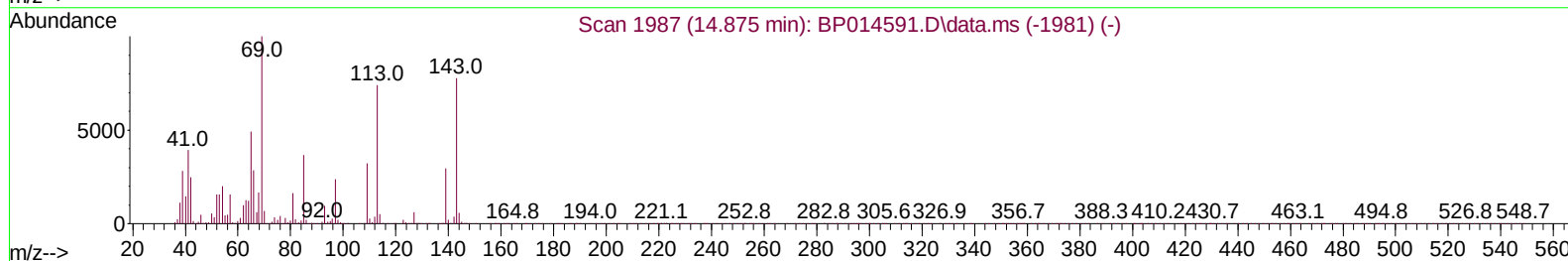
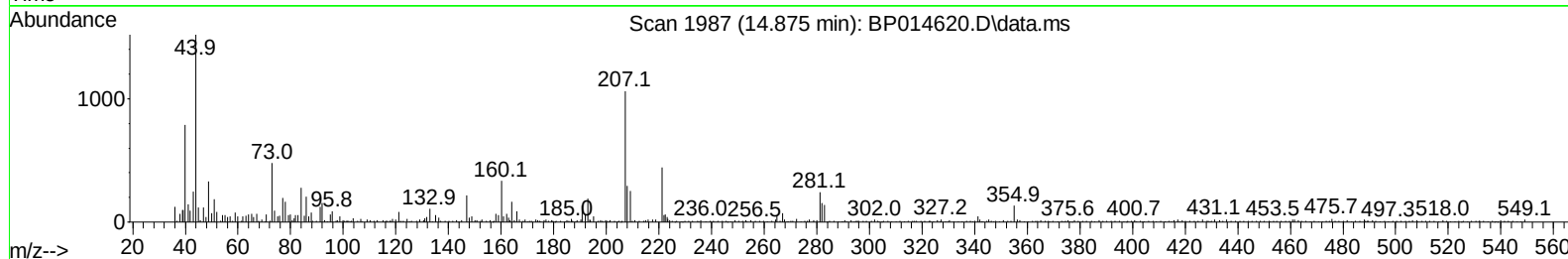
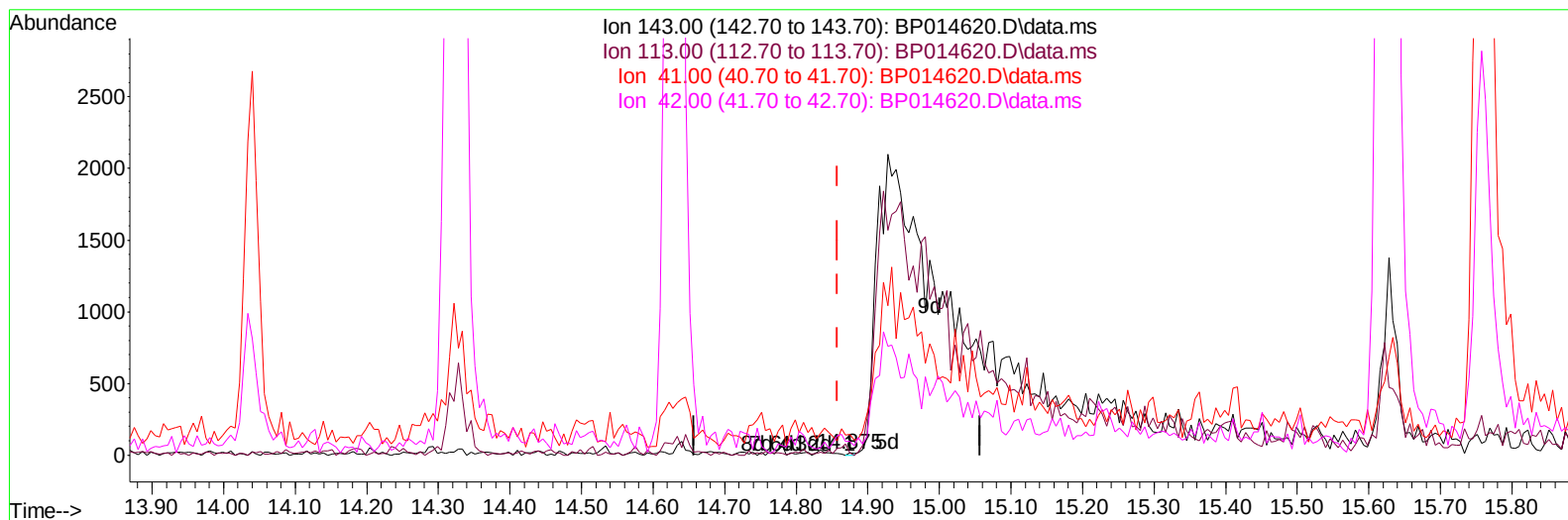
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040723\
Data File : BP014620.D
Acq On : 08 Apr 2023 00:02
Operator : CG/JU
Sample : 02225-14
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB2

Manual IntegrationsAPPROVED

Quant Time: Apr 08 01:52:08 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 07 00:45:34 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/10/2023
Supervised By :Jagrut Upadhyay 04/10/2023



TIC: BP014620.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (+ 0.018) 0.00 ng/ul

response 10

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	94.12
41.00	44.20	852.94#
42.00	31.50	552.94#

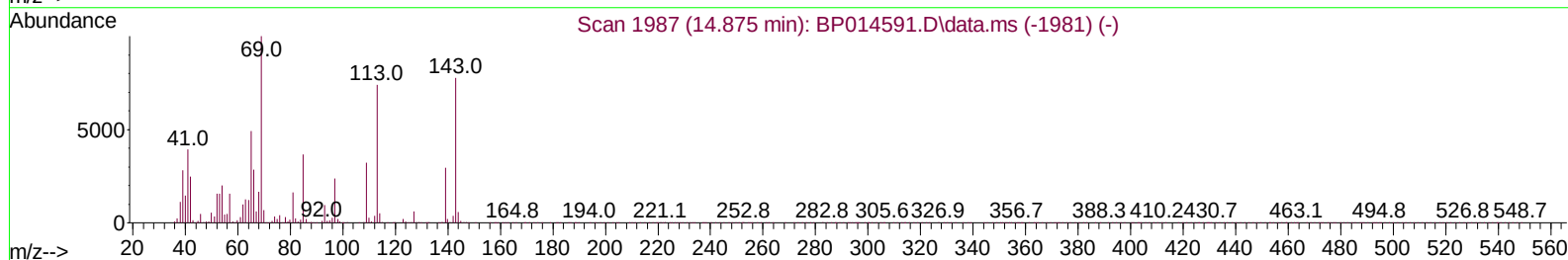
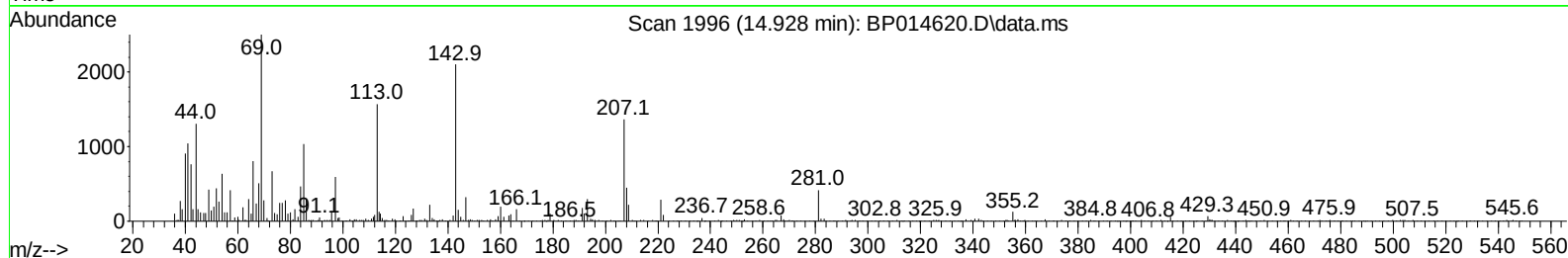
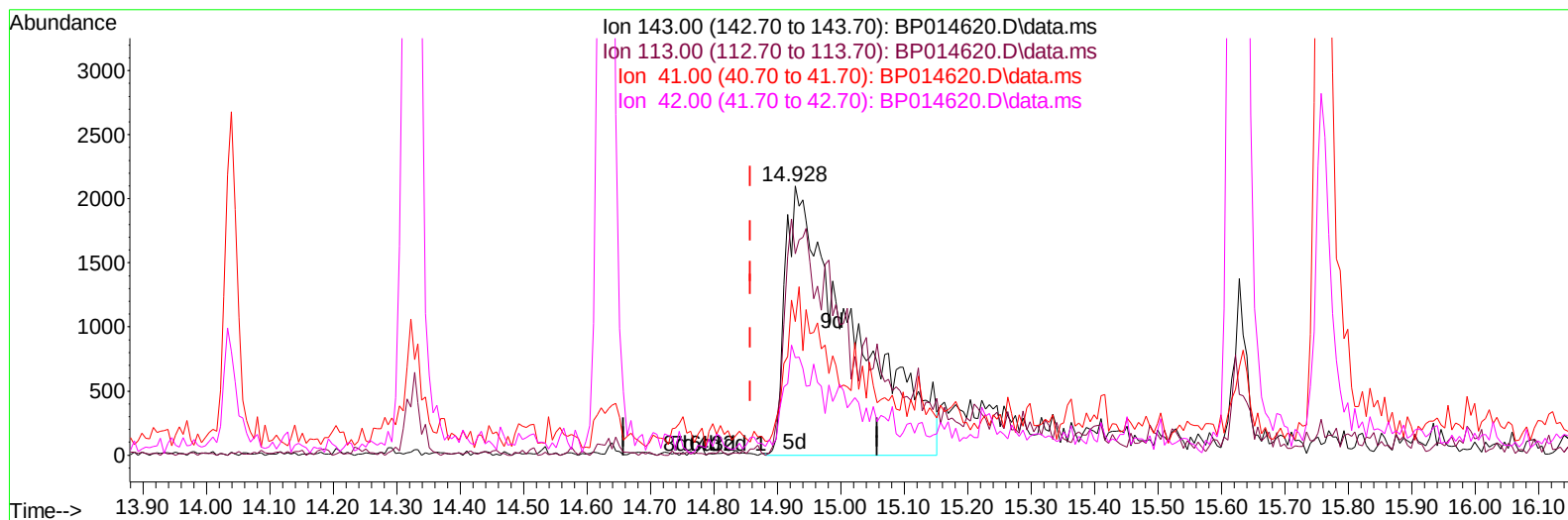
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(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.071) 3.72 ng/ul m

response 15407

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	74.80#
41.00	44.20	49.69
42.00	31.50	36.21

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	124535	20.000	ng/ul	0.00
20) Naphthalene-d8	10.799	136	533223	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.634	164	337789	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	741892	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	751176	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	855531	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	15813	4.924	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.152	99	63905	5.564	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	249756	33.893	ng/ul	0.00
11) 2-Chlorophenol-d4	7.511	132	211090	25.429	ng/ul	0.00
15) 4-Methylphenol-d8	8.699	113	128981	13.922	ng/ul	0.00
21) Nitrobenzene-d5	9.158	128	130799	30.386	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	143185	29.049	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	218237	24.311	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	68731	5.813	ng/ul	0.01
46) Dimethylphthalate-d6	14.040	166	876442	32.000	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	971736	31.975	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	15407m	3.724	ng/ul	0.07
60) Fluorene-d10	15.628	176	743586	32.556	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	135908	24.896	ng/ul	0.00
73) Anthracene-d10	17.492	188	1233676	34.360	ng/ul	0.00
81) Pyrene-d10	19.722	212	1523000	30.185	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.816	264	1626451	35.782	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.399	88	10090	2.859	ng/uL	97

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

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