

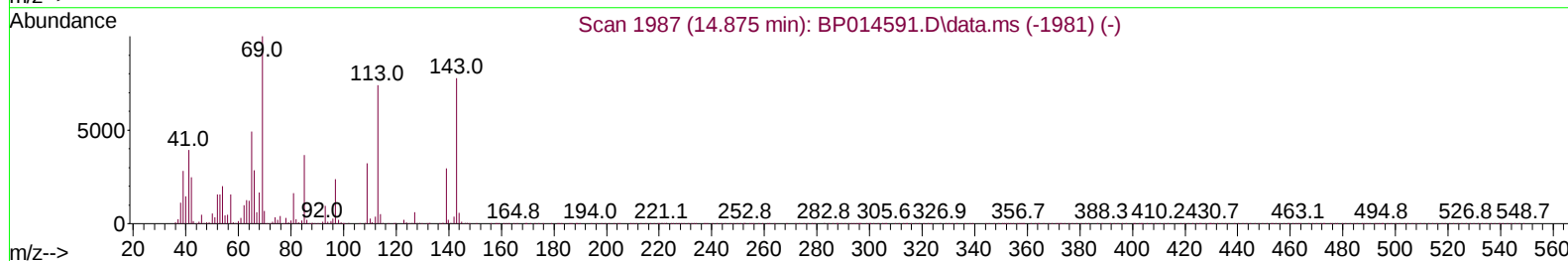
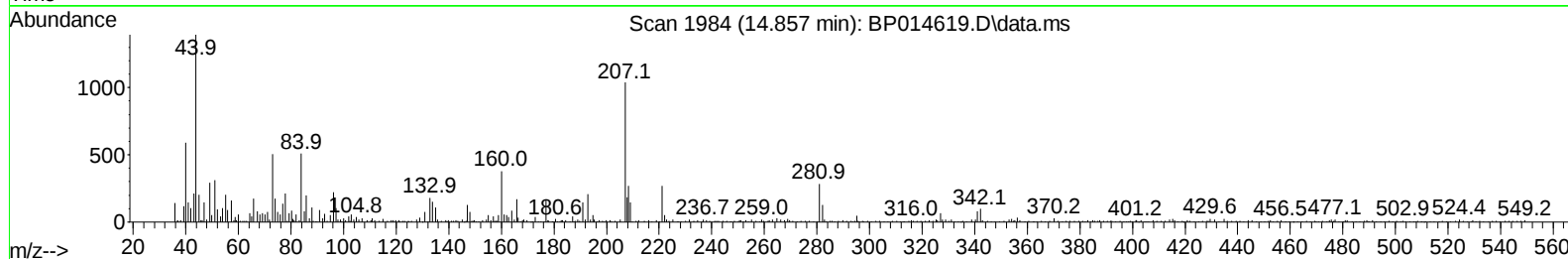
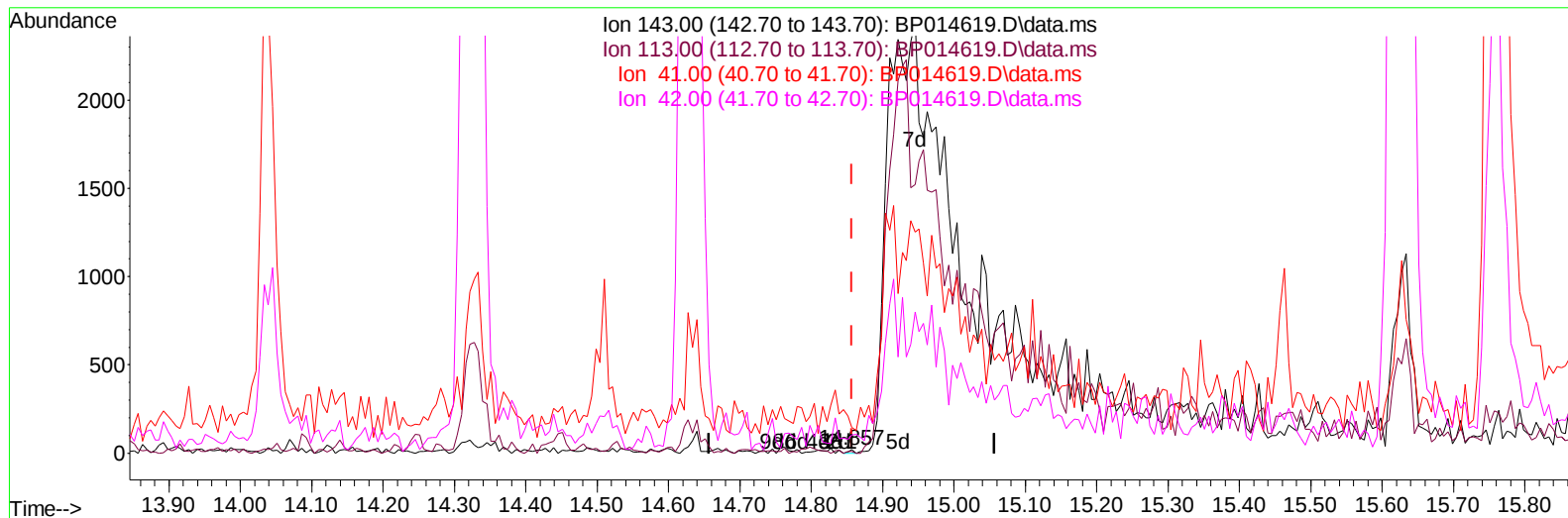
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040723\
Data File : BP014619.D
Acq On : 07 Apr 2023 23:28
Operator : CG/JU
Sample : 02225-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB1

Manual IntegrationsAPPROVED

Quant Time: Apr 08 01:51:56 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: BP014619.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (-0.000) 0.00 ng/u1

response 10

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	36.36#
41.00	44.20	659.09#
42.00	31.50	472.73#

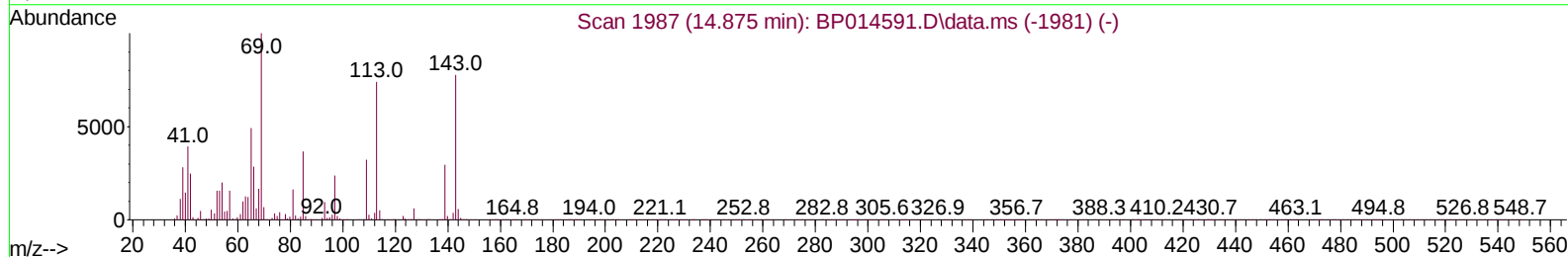
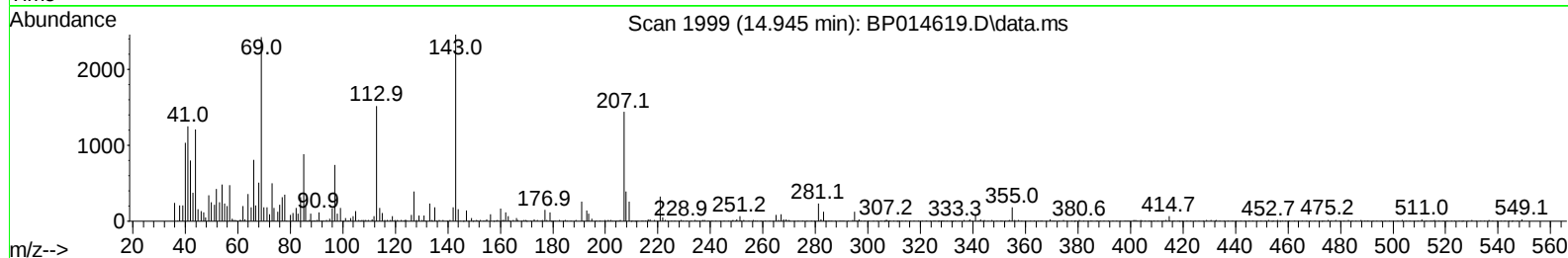
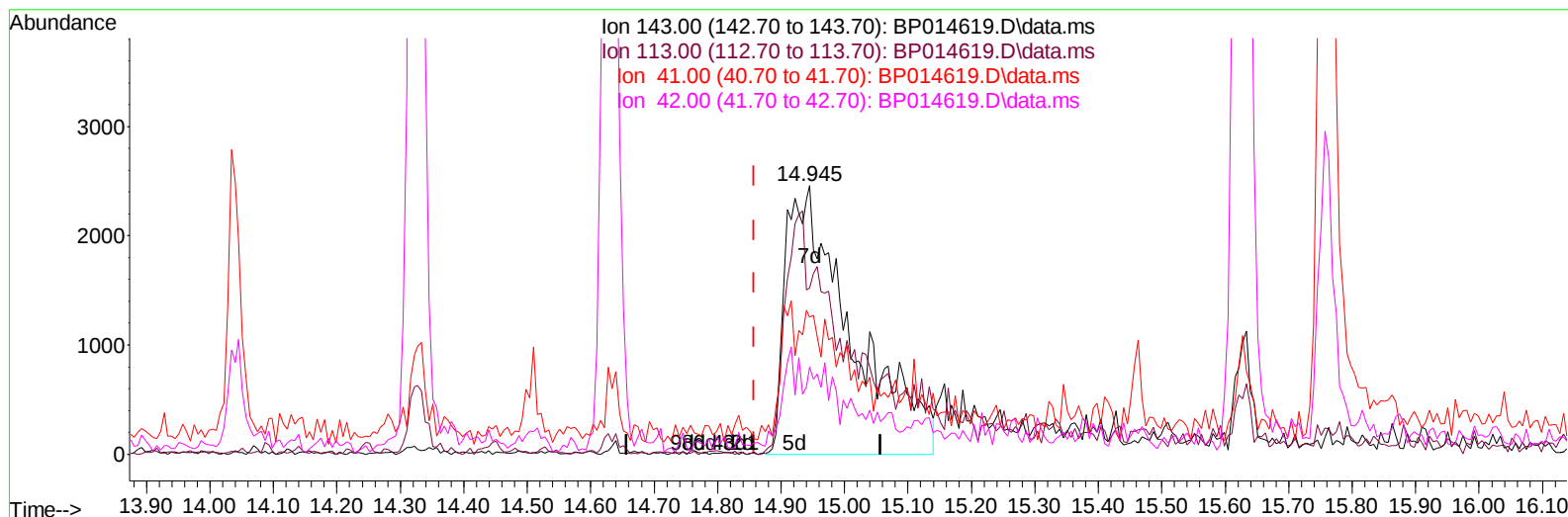
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(54) 4-Nitrophenol-d4 (S)

14.945min (+ 0.088) 4.15 ng/ul m

response 17683

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	61.84#
41.00	44.20	50.98
42.00	31.50	32.55

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	127479	20.000	ng/ul	0.00
20) Naphthalene-d8	10.798	136	545611	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.634	164	347479	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	761822	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	715198	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	869765	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	15349	4.669	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.152	99	66109	5.623	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	256257	33.972	ng/ul	0.00
11) 2-Chlorophenol-d4	7.510	132	206863	24.344	ng/ul	0.00
15) 4-Methylphenol-d8	8.699	113	129485	13.654	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	135563	30.777	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	148847	29.512	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	224719	24.465	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	111663	9.229	ng/ul	0.00
46) Dimethylphthalate-d6	14.039	166	939513	33.347	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	1028327	32.894	ng/ul	0.00
54) 4-Nitrophenol-d4	14.945	143	17683m	4.154	ng/ul	0.09
60) Fluorene-d10	15.628	176	811951	34.558	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.757	200	153477	27.379	ng/ul	0.00
73) Anthracene-d10	17.492	188	1365206	37.029	ng/ul	0.00
81) Pyrene-d10	19.722	212	1654165	34.433	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1735281	37.552	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.399	88	26182	7.248	ng/uL	Qvalue 98

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

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