

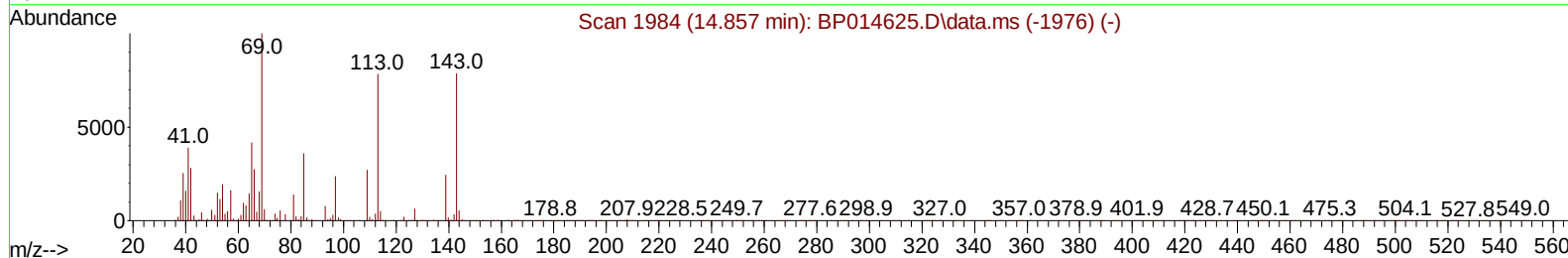
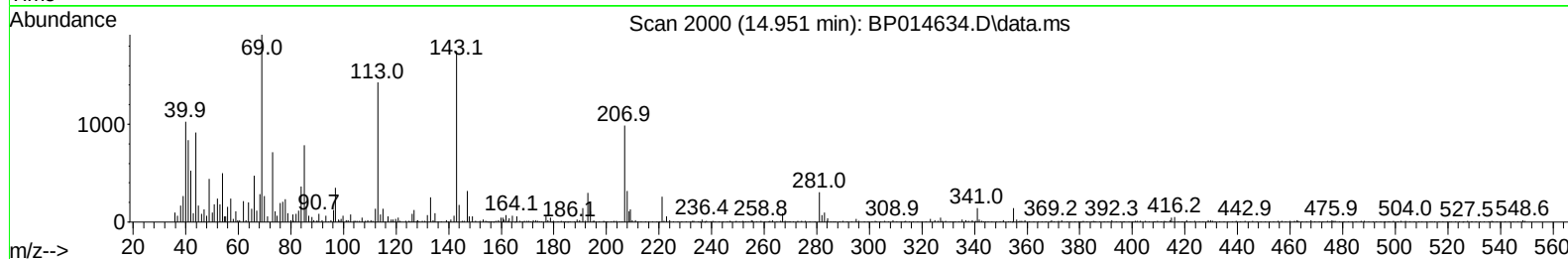
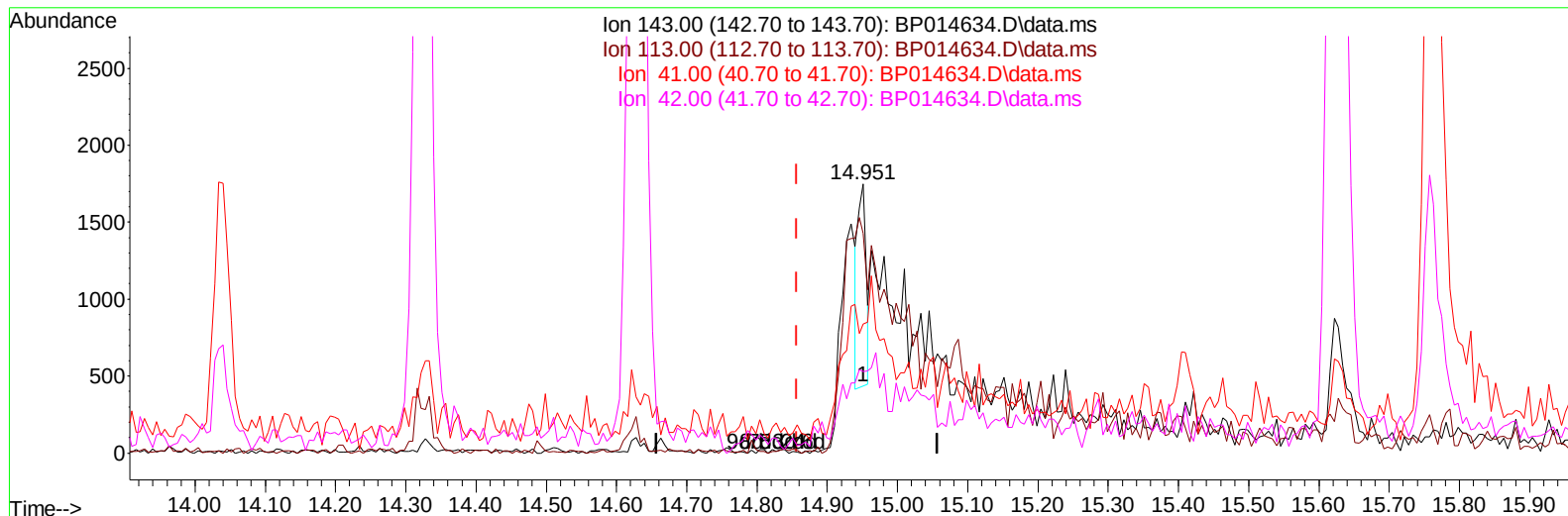
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
Data File : BP014634.D
Acq On : 10 Apr 2023 19:22
Operator : CG/JU
Sample : 02250-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHDx3

Manual IntegrationsAPPROVED

Quant Time: Apr 11 01:24:07 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 11 01:21:43 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/11/2023
Supervised By :Jagrut Upadhyay 04/11/2023



TIC: BP014634.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.094) 0.30 ng/ul

response 1060

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	81.68
41.00	44.20	48.08
42.00	31.50	30.22

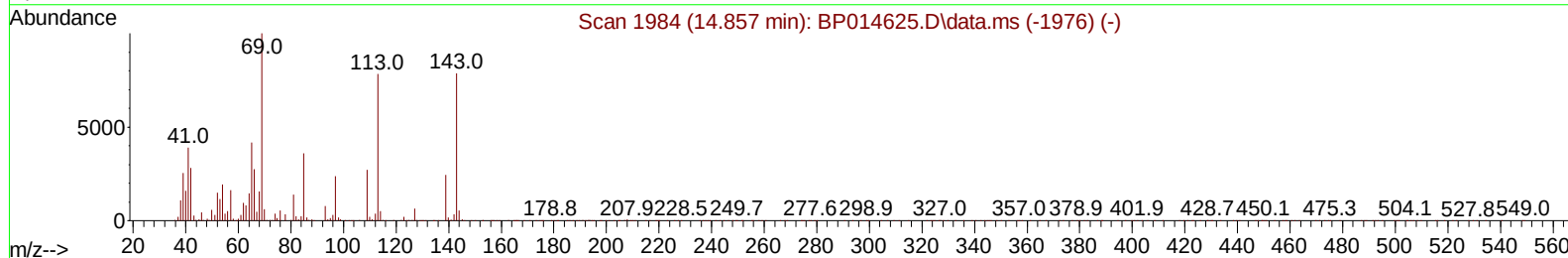
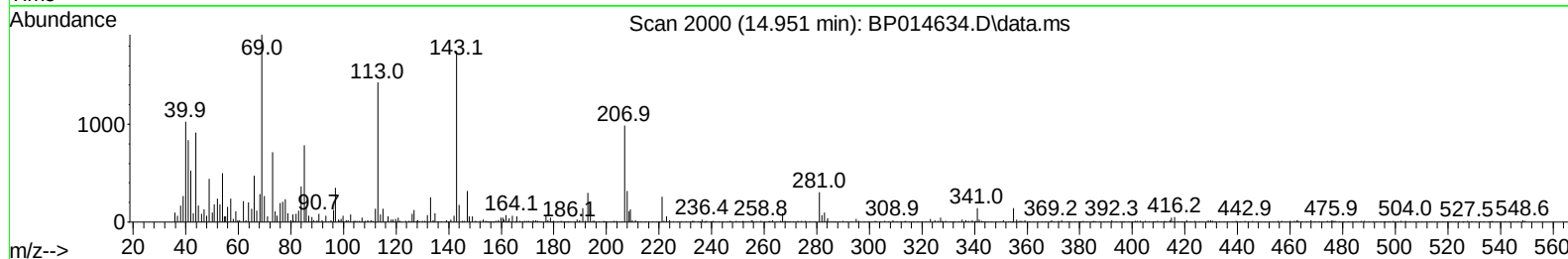
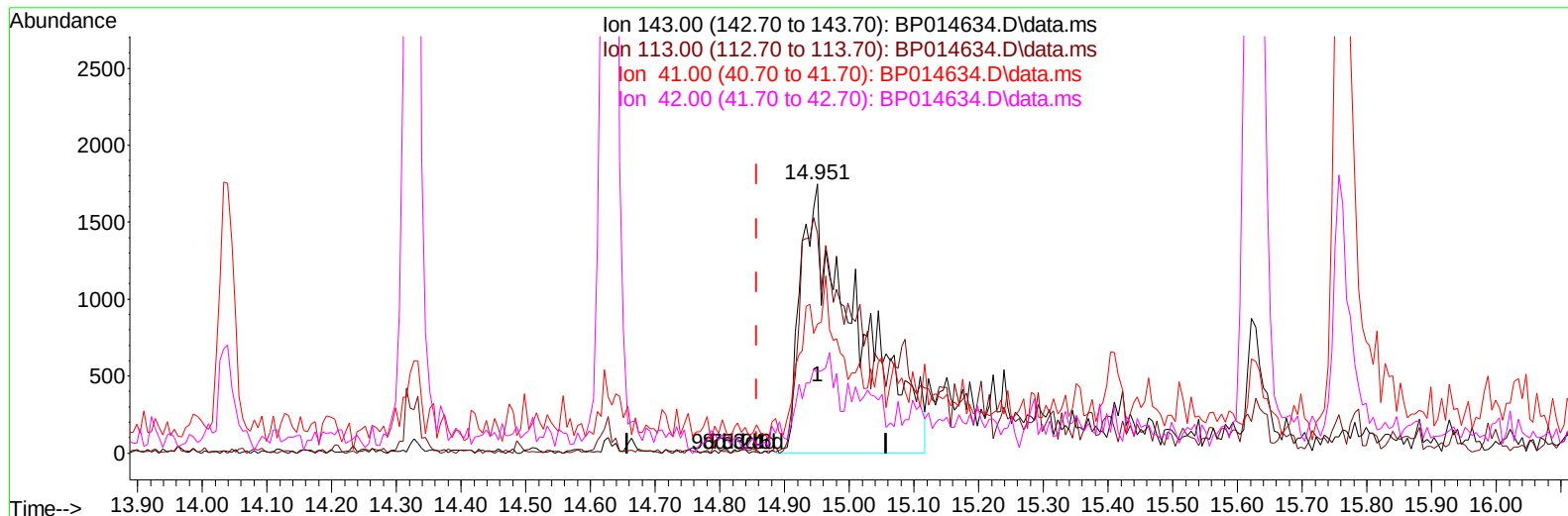
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(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.094) 3.02 ng/ul m

response 10720

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	81.68
41.00	44.20	48.08
42.00	31.50	30.22

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	117192	20.000	ng/ul	0.00
20) Naphthalene-d8	10.799	136	492526	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	289661	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	603123	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	581434	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	691343	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	13500	4.467	ng/uL	0.00
4) Pyridine-d5	3.817	84	28384	3.665	ng/ul	0.02
7) Phenol-d5	7.152	99	57422	5.313	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	177826	25.644	ng/ul	0.00
11) 2-Chlorophenol-d4	7.510	132	157266	20.132	ng/ul	0.00
15) 4-Methylphenol-d8	8.705	113	100581	11.537	ng/ul	0.00
21) Nitrobenzene-d5	9.157	128	89266	22.451	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	101286	22.246	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	159759	19.267	ng/ul	0.00
31) 4-Chloroaniline-d4	10.951	131	174383	15.966	ng/ul	0.00
46) Dimethylphthalate-d6	14.040	166	600724	25.578	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	636545	24.426	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	10720m	3.021	ng/ul	0.09
60) Fluorene-d10	15.628	176	511867	26.135	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	85055	19.166	ng/ul	0.00
73) Anthracene-d10	17.492	188	887700	30.413	ng/ul	0.00
81) Pyrene-d10	19.722	212	1061850	27.189	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1130038	30.766	ng/ul	0.00

Target Compounds Qvalue

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

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