

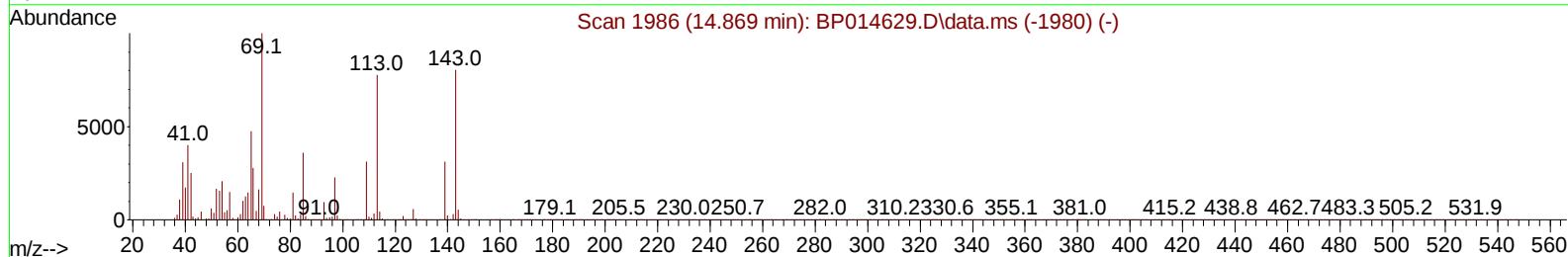
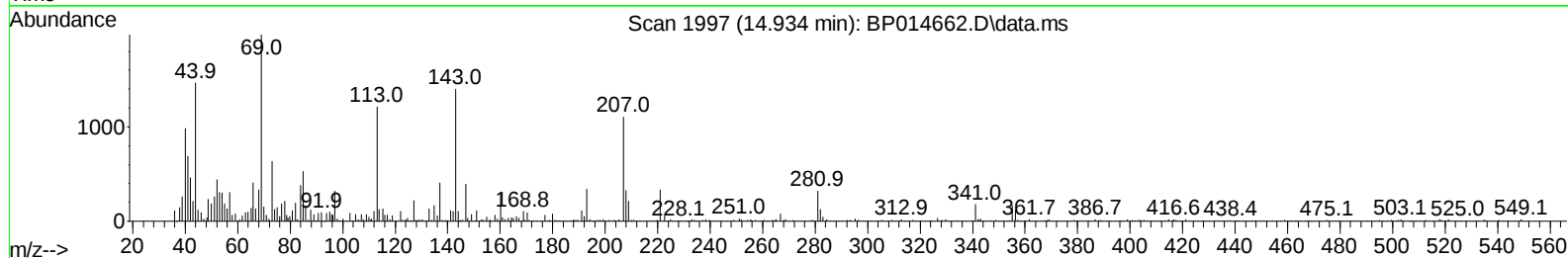
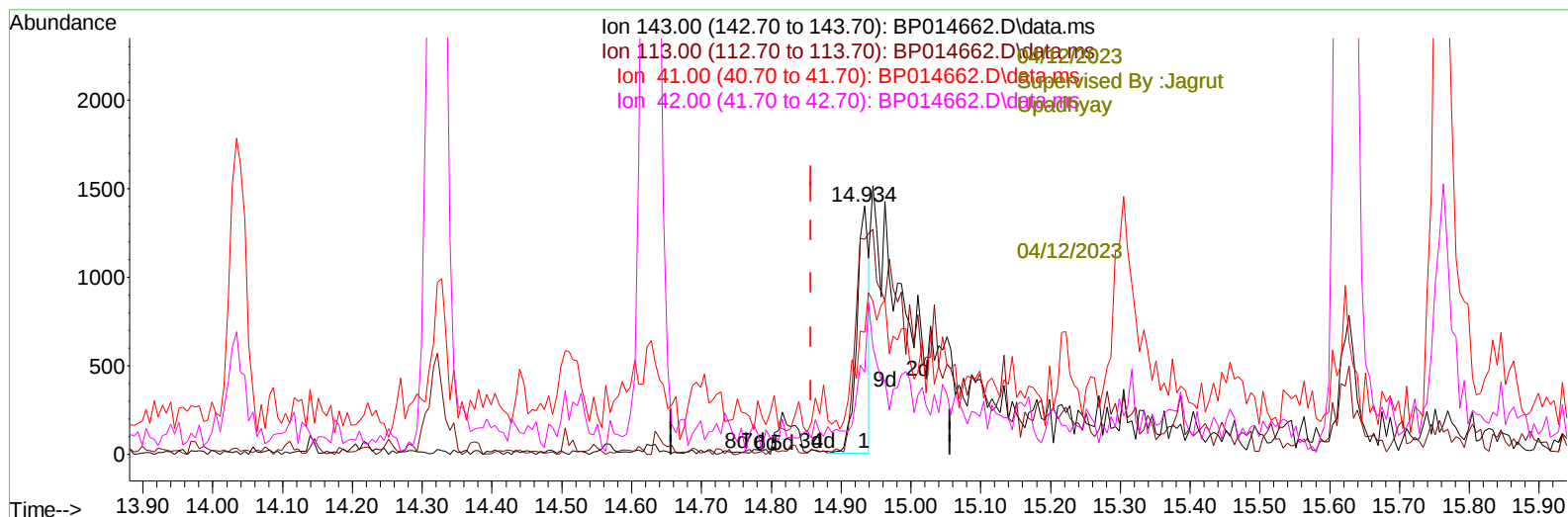
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041023\
Data File : BP014662.D
Acq On : 11 Apr 2023 18:35
Operator : CG/JU
Sample : 02251-11
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMC9

Manual IntegrationsAPPROVED

Quant Time: Apr 11 23:02:44 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



TIC: BP014662.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.934min (+ 0.077) 0.50 ng/u1

response 1679

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	86.59
41.00	44.20	49.14
42.00	31.50	32.95

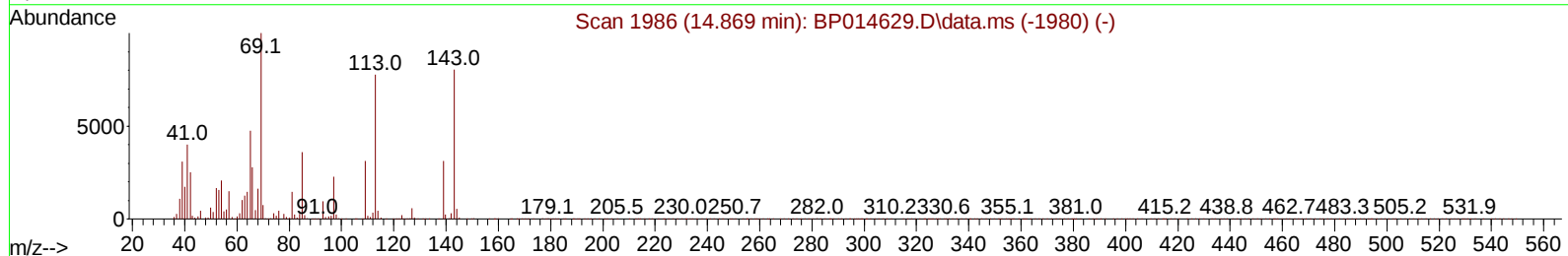
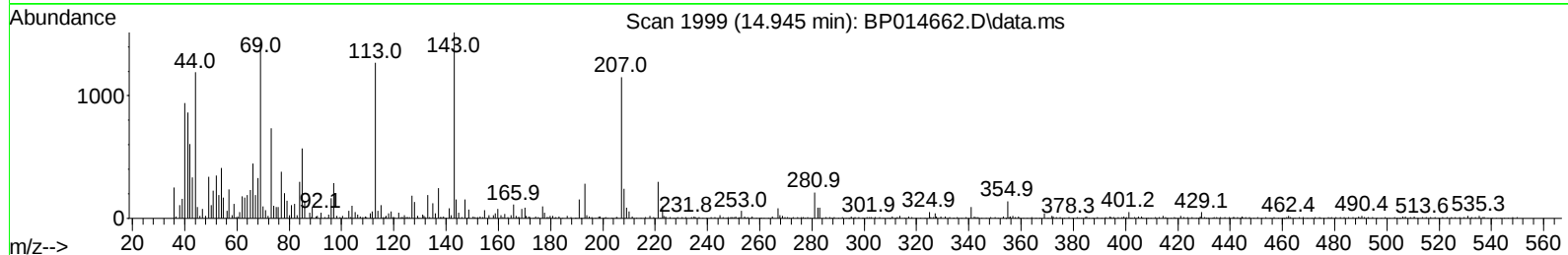
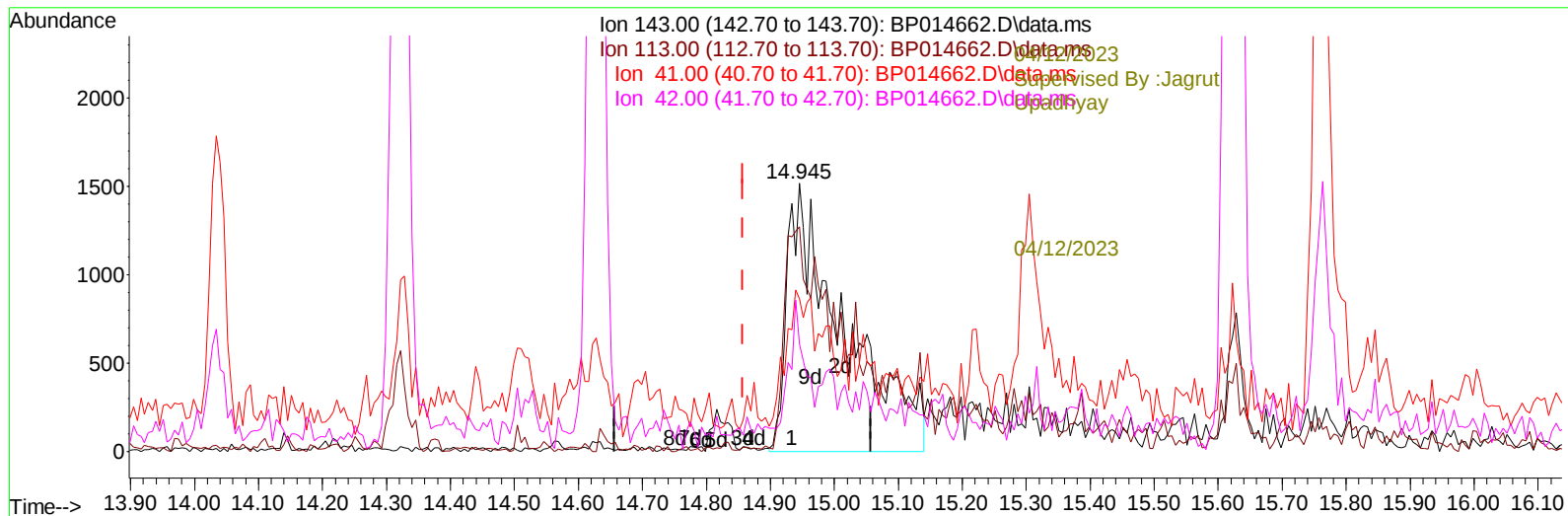
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(54) 4-Nitrophenol-d4 (S)

14.945min (+ 0.088) 2.74 ng/ul m

response 9189

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	83.77
41.00	44.20	56.99#
42.00	31.50	39.97#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----04/12/2023						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	109050	20.000	ng/ul	0.00
20) Naphthalene-d8	10.793	136	449029	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	273523	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	612783	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	633634	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	718005	20.000	ng/ul	0.00
-----04/12/2023						
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	6149	2.187	ng/uL	0.00
4) Pyridine-d5	3.817	84	34116	4.734	ng/ul	0.02
7) Phenol-d5	7.152	99	47664	4.739	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	184374	28.573	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	150587	20.716	ng/ul	0.00
15) 4-Methylphenol-d8	8.699	113	96979	11.954	ng/ul	0.00
21) Nitrobenzene-d5	9.152	128	93381	25.761	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	98362	23.697	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	166056	21.967	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	190654	19.147	ng/ul	0.00
46) Dimethylphthalate-d6	14.034	166	642948	28.991	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	681292	27.685	ng/ul	0.00
54) 4-Nitrophenol-d4	14.945	143	9189m	2.743	ng/ul	0.09
60) Fluorene-d10	15.622	176	550055	29.741	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	72950	16.179	ng/ul	0.00
73) Anthracene-d10	17.492	188	928157	31.297	ng/ul	0.00
81) Pyrene-d10	19.722	212	1186683	27.882	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1238254	32.460	ng/ul	0.00

Target Compounds Qvalue

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

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