

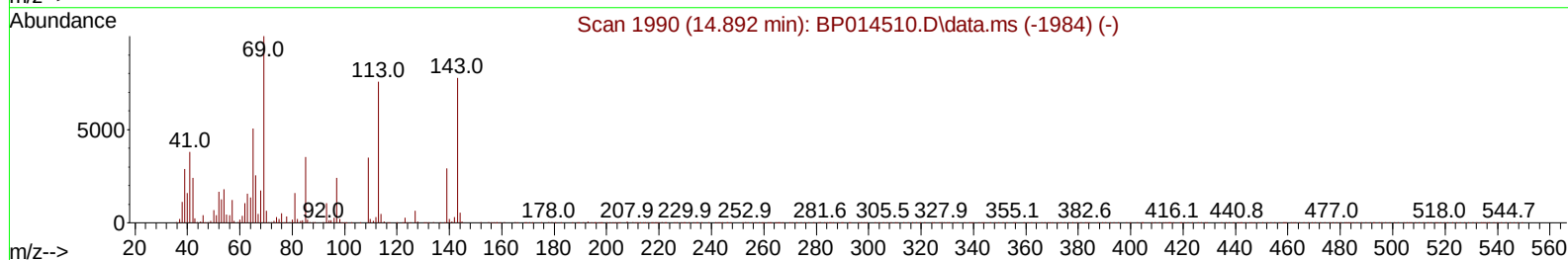
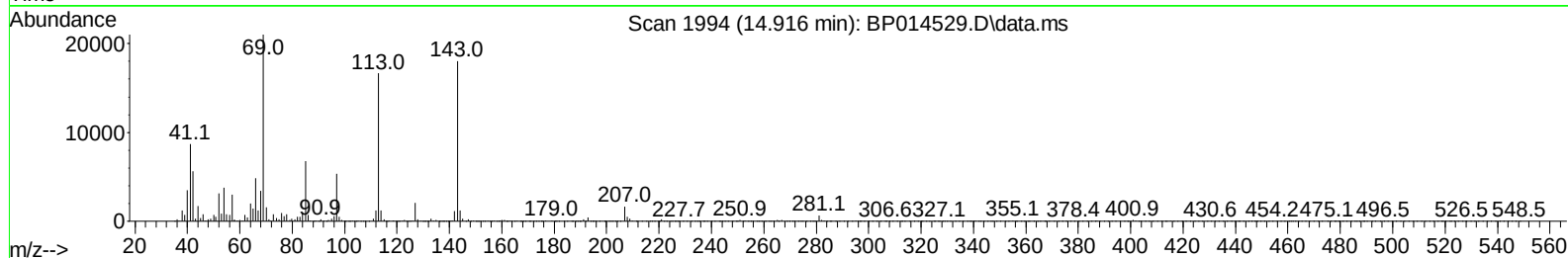
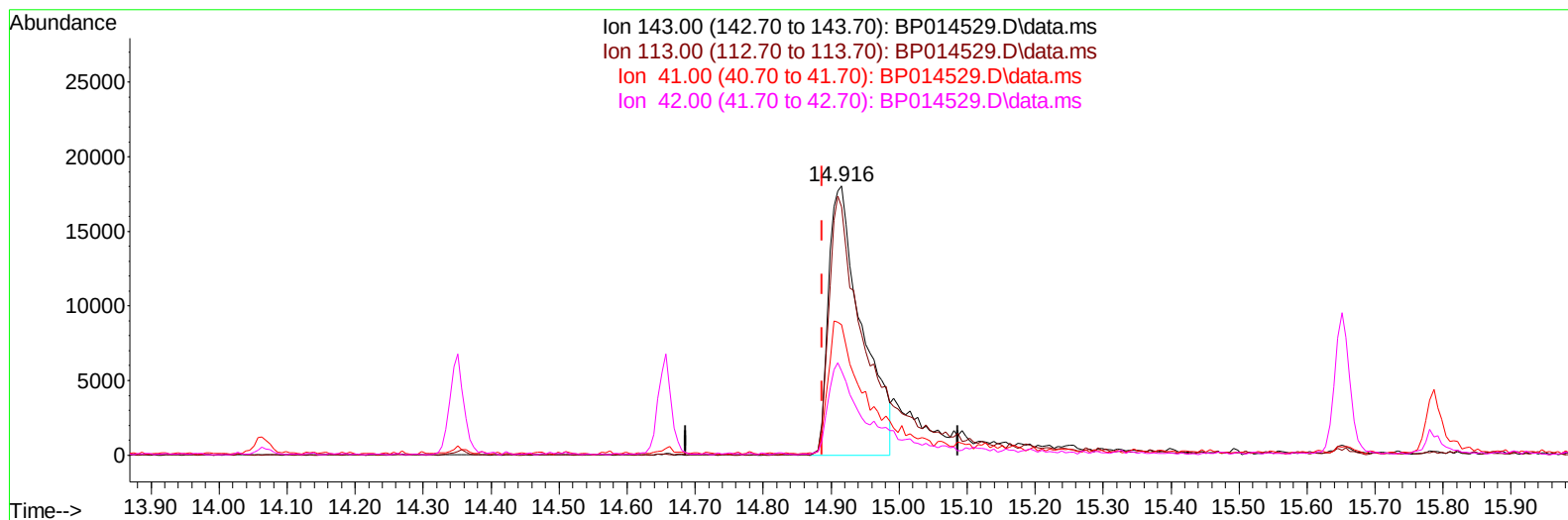
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014529.D
Acq On : 04 Apr 2023 15:44
Operator : CG/JU
Sample : PB151869BL
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK869

Manual IntegrationsAPPROVED

Quant Time: Apr 04 23:45:45 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 04 23:43:56 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/05/2023
Supervised By :Jagrut Upadhyay 04/05/2023



TIC: BP014529.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.916min (+ 0.029) 21.14 ng/ul

response 60568

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	92.32
41.00	44.20	48.36
42.00	31.50	31.34

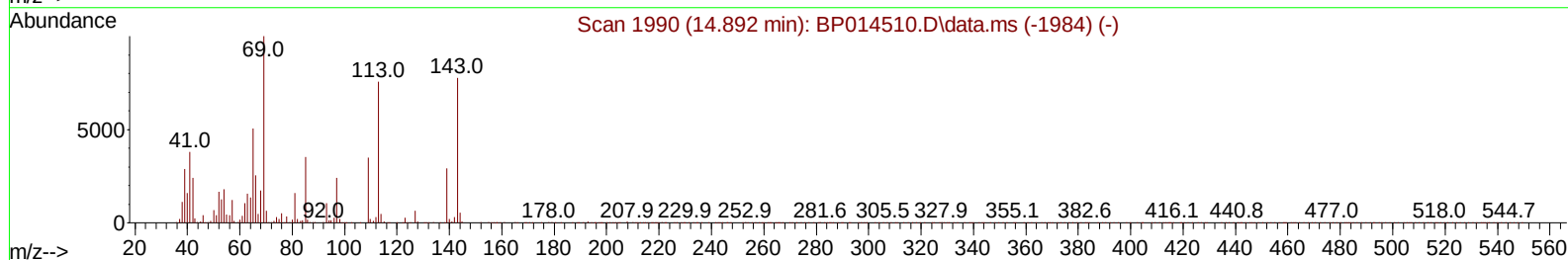
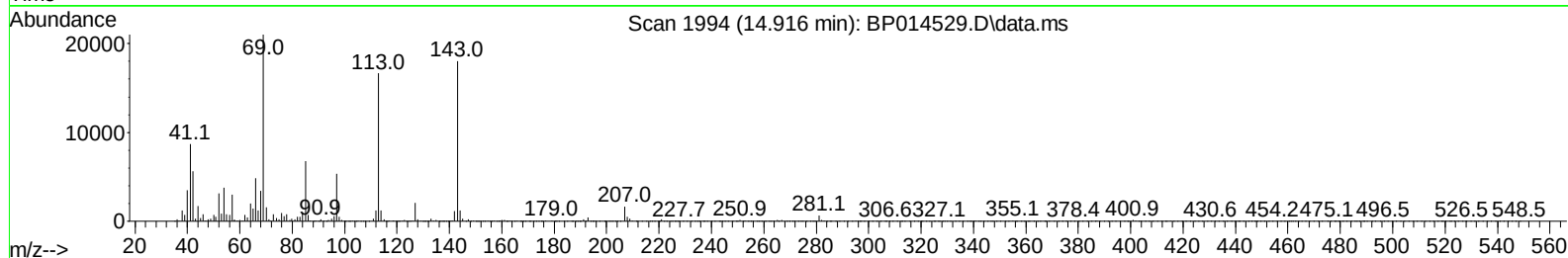
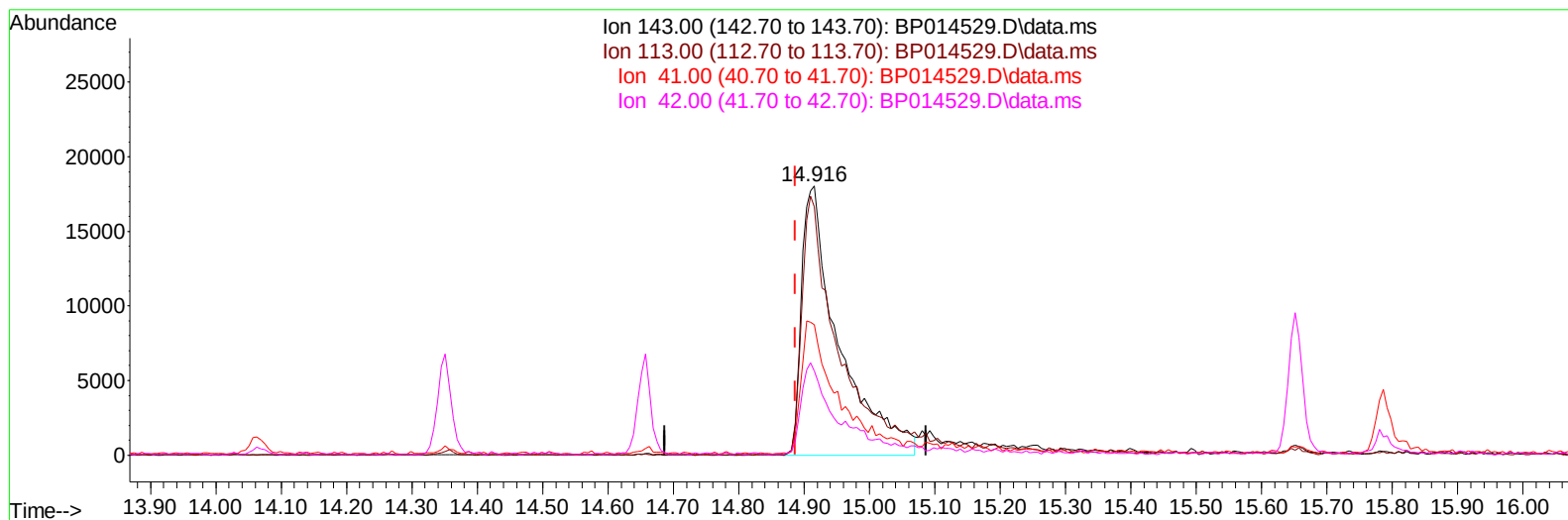
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(54) 4-Nitrophenol-d4 (S)

14.916min (+ 0.029) 25.05 ng/ul m

response 71767

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	92.32
41.00	44.20	48.36
42.00	31.50	31.34

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.004	152	87021	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	365851	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	233903	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	553925	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	569744	20.000	ng/ul	0.00
88) Perylene-d12	24.021	264	644469	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	14668	6.537	ng/uL	0.00
4) Pyridine-d5	3.816	84	173685	30.203	ng/ul	0.00
7) Phenol-d5	7.169	99	233650	29.113	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	151745	29.470	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	176357	30.403	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	178747	27.612	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	84340	28.556	ng/ul	0.00
24) 2-Nitrophenol-d4	9.910	143	94550	27.958	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	156621	25.429	ng/ul	0.00
31) 4-Chloroaniline-d4	10.975	131	208496	25.700	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	559184	29.485	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	604852	28.742	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	71767m	25.048	ng/ul	0.03
60) Fluorene-d10	15.651	176	488370	30.879	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	75985	18.643	ng/ul	0.00
73) Anthracene-d10	17.516	188	814979	30.401	ng/ul	0.00
81) Pyrene-d10	19.745	212	1012858	26.466	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.851	264	1071276	31.287	ng/ul	0.00

Target Compounds Qvalue

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

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