

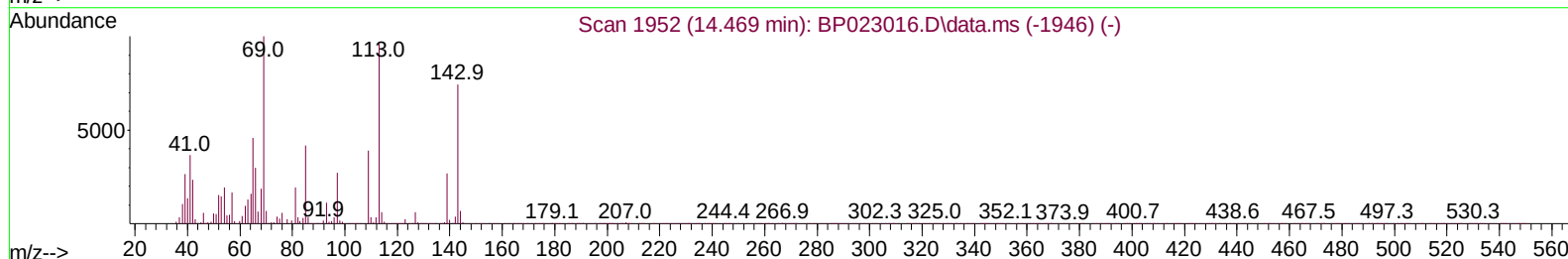
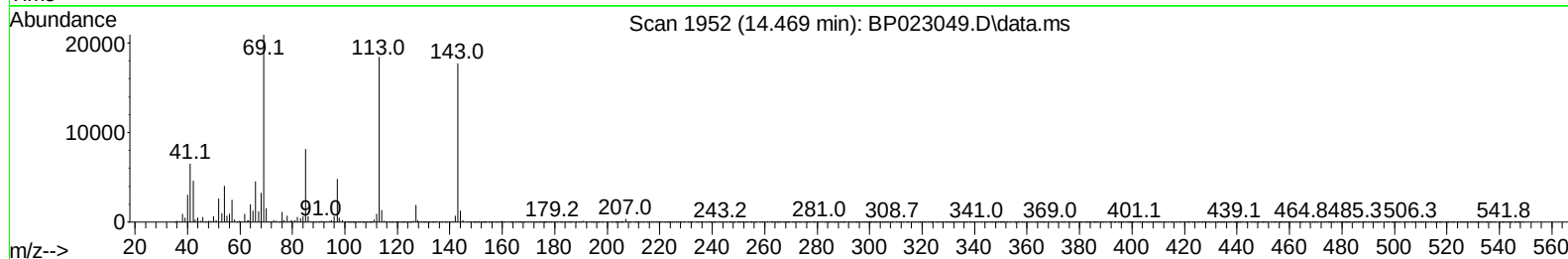
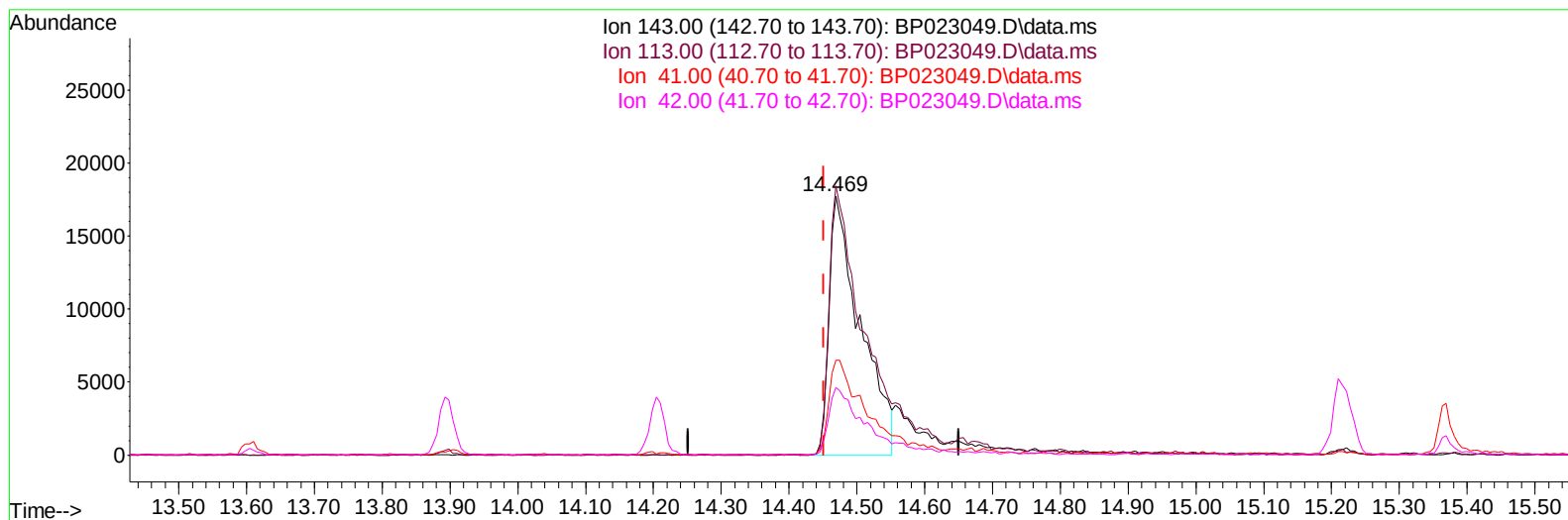
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111224\
Data File : BP023049.D
Acq On : 12 Nov 2024 16:35
Operator : RC/JU
Sample : PB164853BL
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK853

Manual IntegrationsAPPROVED

Quant Time: Nov 13 01:59:04 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 12 23:59:19 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2024
Supervised By :mohammad ahmed 11/14/2024



TIC: BP023049.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.018) 25.26 ng/ul

response 56638

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	103.79
41.00	39.20	36.62
42.00	25.90	26.10

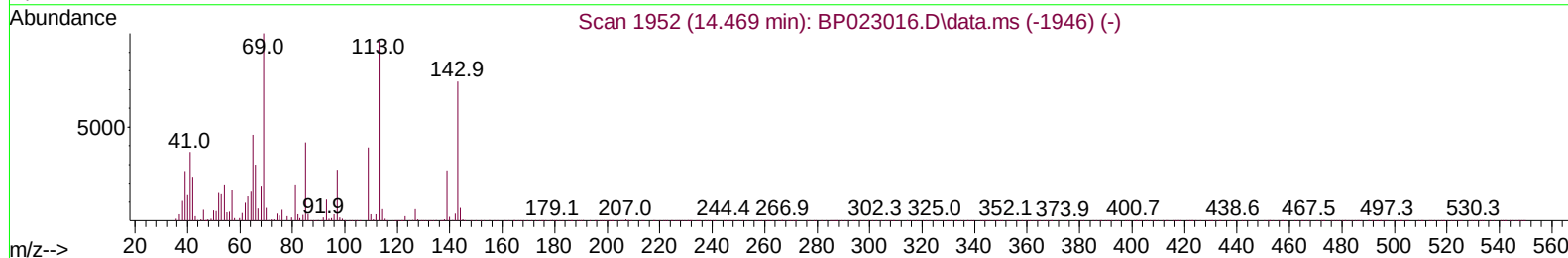
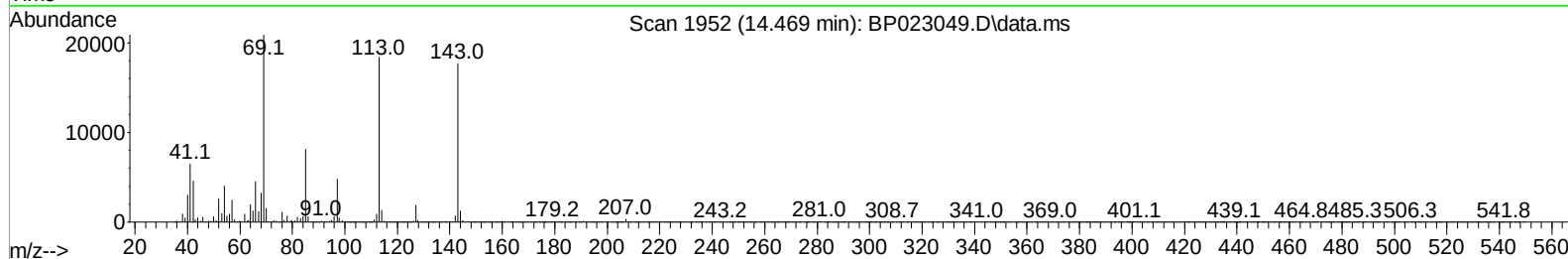
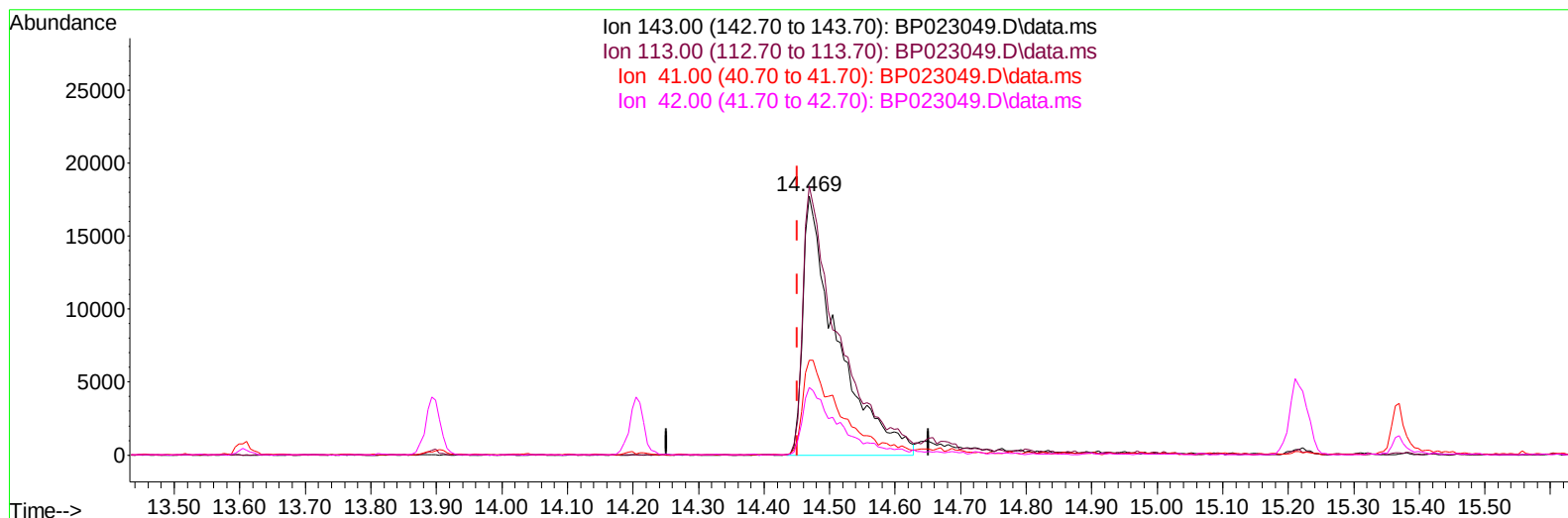
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(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.018) 28.99 ng/ul m

response 65000

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	103.79
41.00	39.20	36.62
42.00	25.90	26.10

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Quant Time: Nov 13 01:59:45 2024
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 QLast Update : Tue Nov 12 23:59:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	66748	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136	257536	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.210	164	206974	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.010	188	512984	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	582269	20.000	ng/ul	0.00
88) Perylene-d12	24.668	264	677037	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	9451	6.806	ng/uL	0.00
4) Pyridine-d5	3.552	84	124085	30.076	ng/ul	0.00
7) Phenol-d5	6.793	99	151937	30.789	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	94937	32.744	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	117893	32.620	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	128570	31.660	ng/ul	0.00
21) Nitrobenzene-d5	8.728	128	58841	32.819	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	70738	33.095	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	138865	31.023	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131	159766	29.918	ng/ul	0.00
46) Dimethylphthalate-d6	13.604	166	510802	34.242	ng/ul	-0.01
49) Acenaphthylene-d8	13.893	160	525837	33.783	ng/ul	0.00
54) 4-Nitrophenol-d4	14.469	143	65000m	28.995	ng/ul	0.02
60) Fluorene-d10	15.216	176	437262	33.433	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	83420	27.146	ng/ul	0.00
73) Anthracene-d10	17.122	188	741517	34.139	ng/ul	0.00
81) Pyrene-d10	19.492	212	959151	35.763	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.445	264	1153876	35.779	ng/ul	-0.02

Target Compounds Qvalue

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

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