

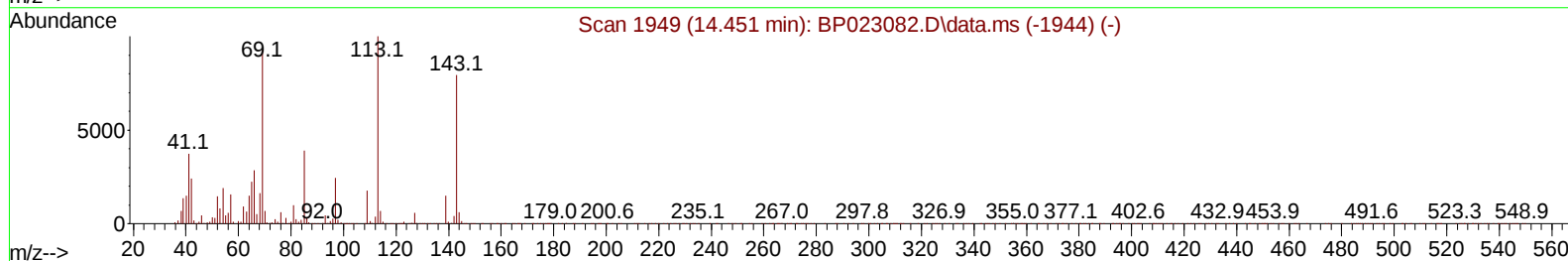
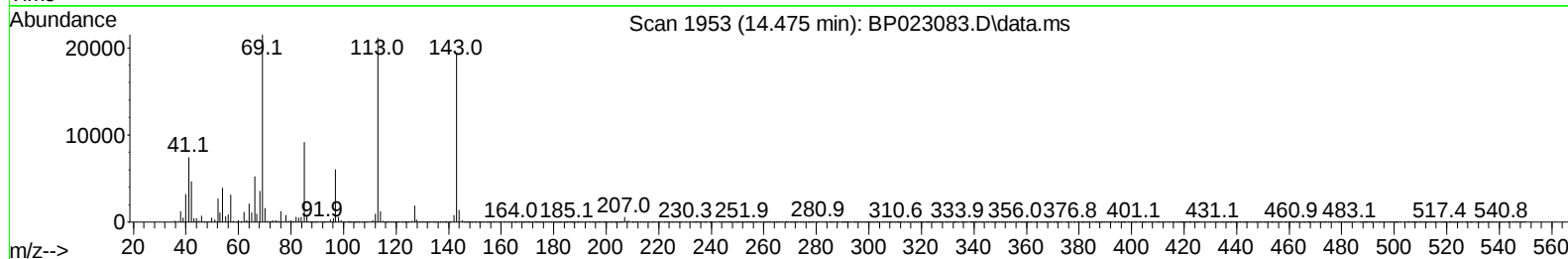
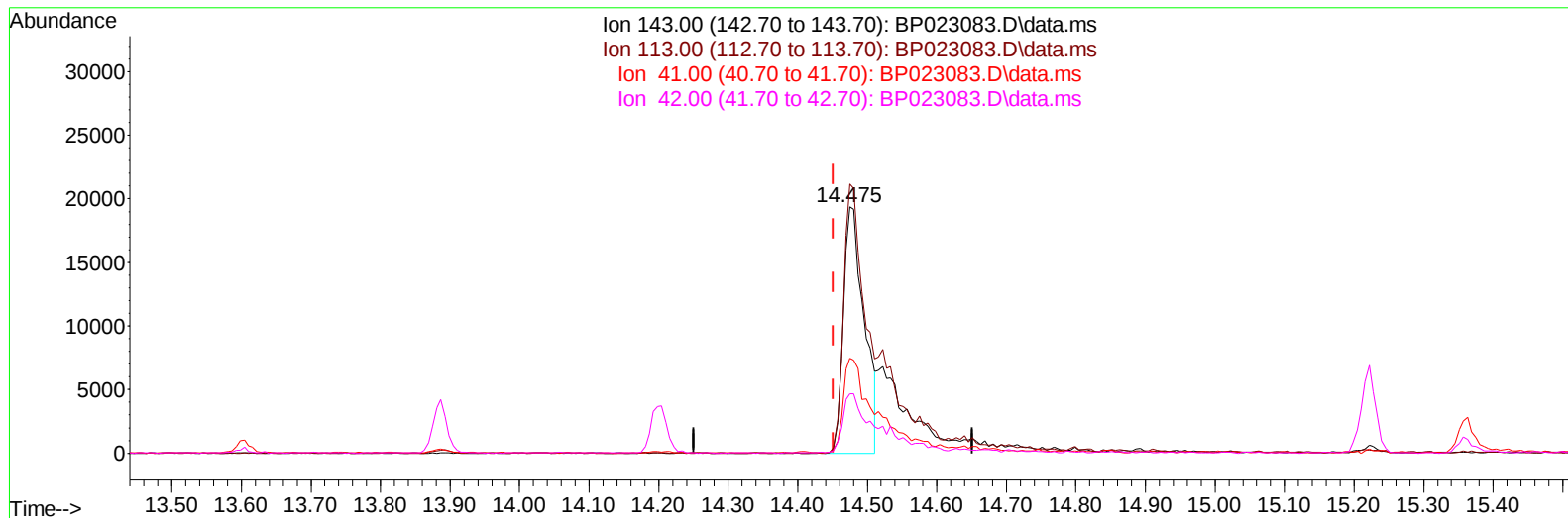
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023083.D
Acq On : 13 Nov 2024 16:11
Operator : RC/JU
Sample : PB164783BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK783

Manual IntegrationsAPPROVED

Quant Time: Nov 14 01:33:33 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 13 04:49:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/15/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BP023083.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.475min (+ 0.024) 18.40 ng/u1

response 40498

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	109.08
41.00	39.20	38.37
42.00	25.90	24.11

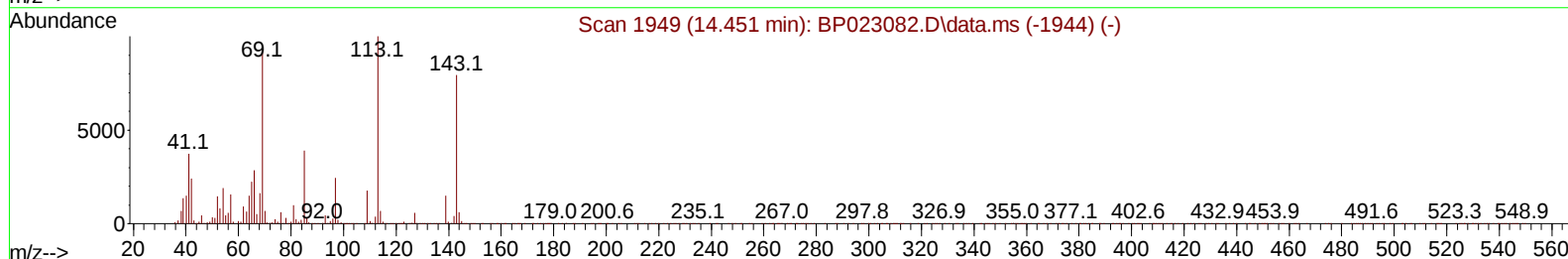
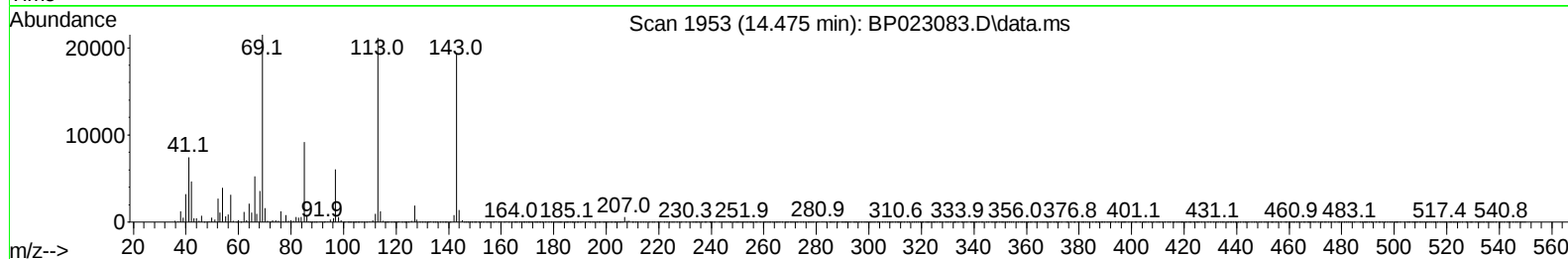
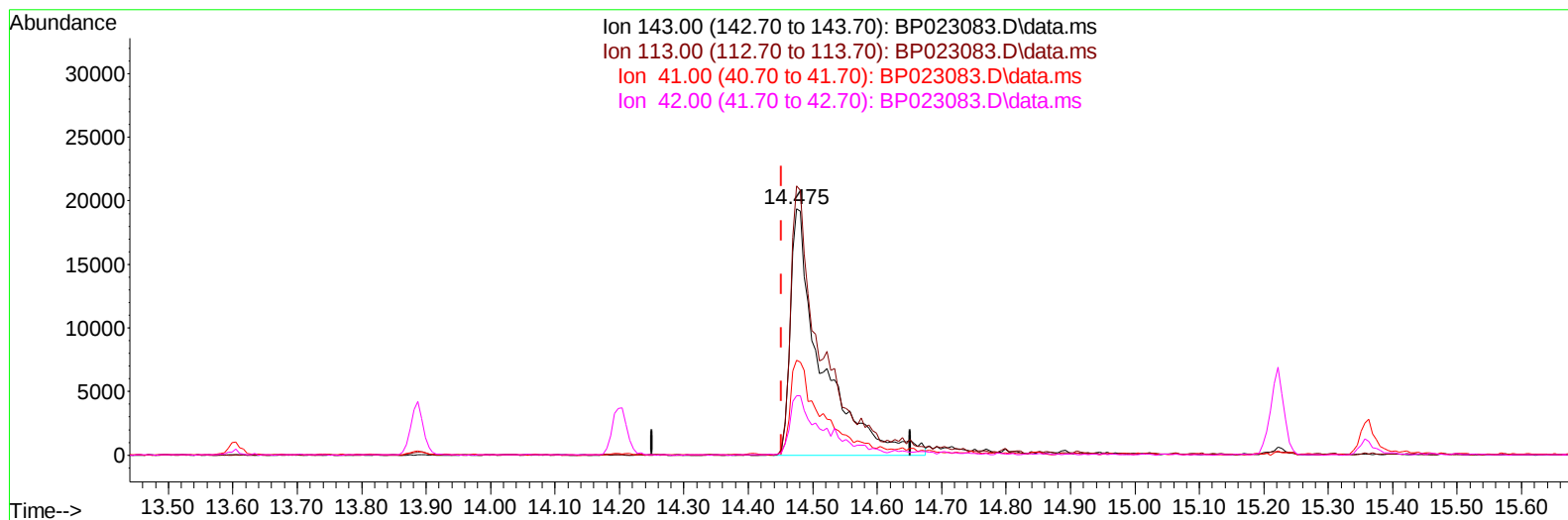
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TIC: BP023083.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.475min (+ 0.024) 29.39 ng/ul m

response 64676

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	109.08
41.00	39.20	38.37
42.00	25.90	24.11

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	67751	20.000	ng/ul	0.00
20) Naphthalene-d8	10.328	136	259259	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.204	164	203168	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	499221	20.000	ng/ul	-0.01
79) Chrysene-d12	21.439	240	582478	20.000	ng/ul	0.00
88) Perylene-d12	24.645	264	671763	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	9696	6.879	ng/uL	0.00
4) Pyridine-d5	3.552	84	124905	29.827	ng/ul	0.00
7) Phenol-d5	6.787	99	151222	30.190	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	95806	32.555	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	116747	31.824	ng/ul	0.00
15) 4-Methylphenol-d8	8.287	113	123740	30.020	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	57394	31.799	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	69259	32.187	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	131440	29.169	ng/ul	0.02
31) 4-Chloroaniline-d4	10.481	131	157592	29.315	ng/ul	0.00
46) Dimethylphthalate-d6	13.604	166	492679	33.646	ng/ul	0.00
49) Acenaphthylene-d8	13.887	160	511963	33.508	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.475	143	64676m	29.391	ng/ul	0.02
60) Fluorene-d10	15.222	176	425734	33.162	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.357	200	78781	26.343	ng/ul	0.00
73) Anthracene-d10	17.104	188	746208	35.303	ng/ul	-0.01
81) Pyrene-d10	19.486	212	968596	36.102	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.421	264	1142560	35.706	ng/ul	-0.02

Target Compounds Qvalue

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

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