

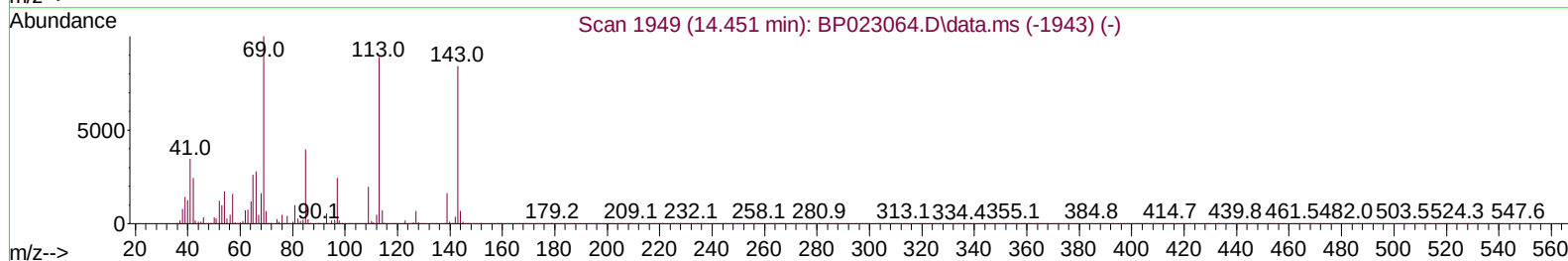
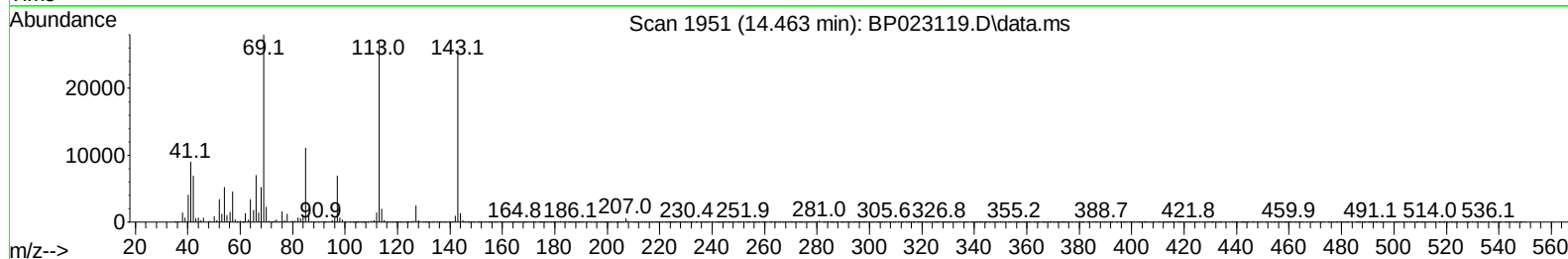
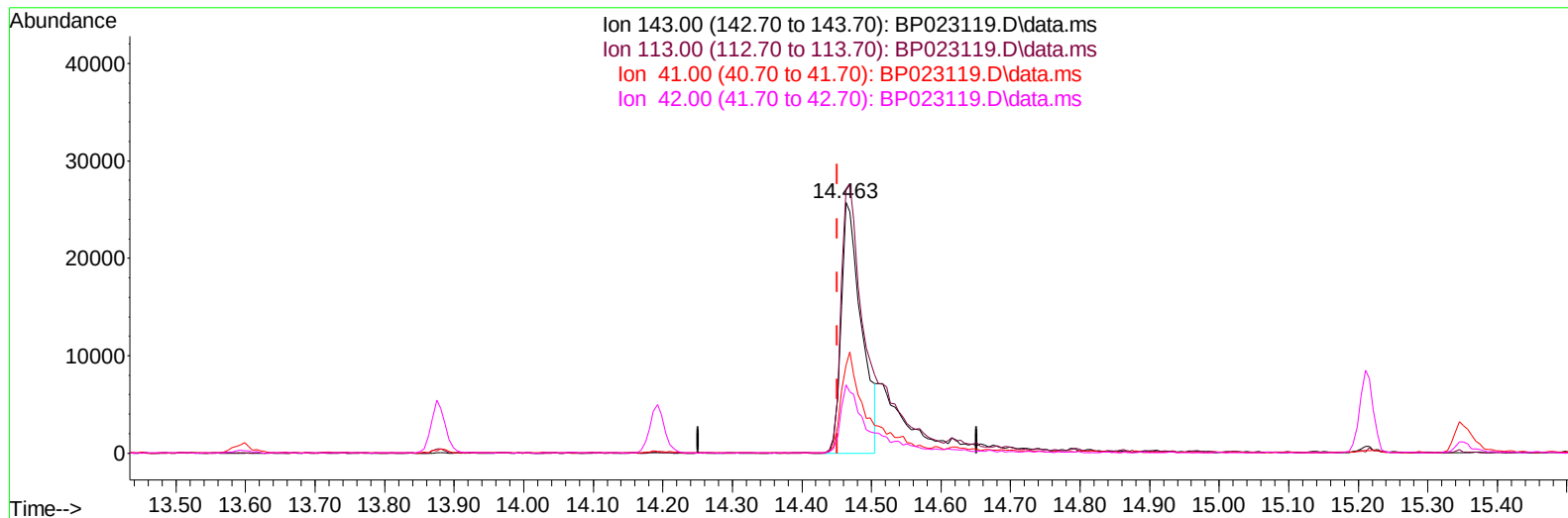
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023119.D
Acq On : 15 Nov 2024 01:12
Operator : RC/JU
Sample : PB164889BL
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK889

Manual IntegrationsAPPROVED

Quant Time: Nov 15 01:36:52 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: BP023119.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.463min (+ 0.012) 18.86 ng/ul

response 51744

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	105.00
41.00	39.20	35.01
42.00	25.90	27.09

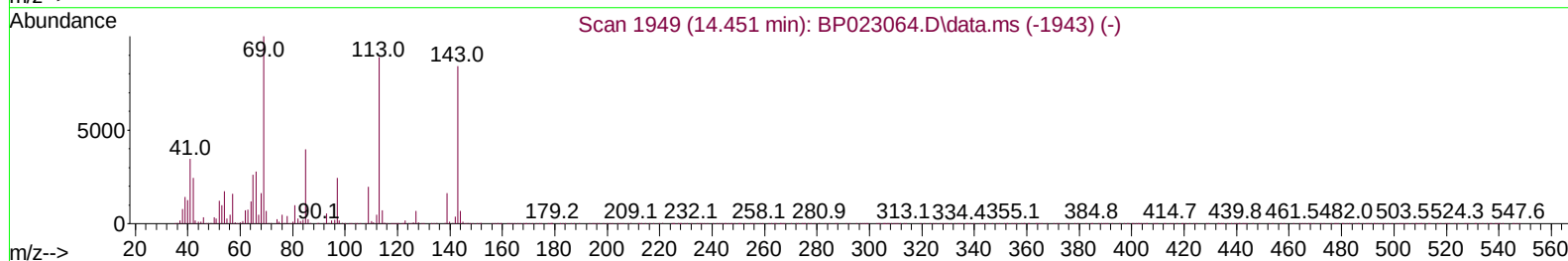
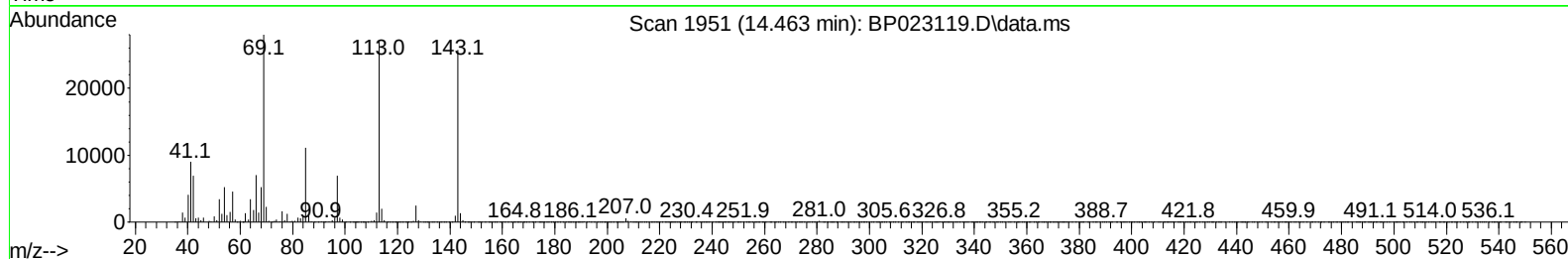
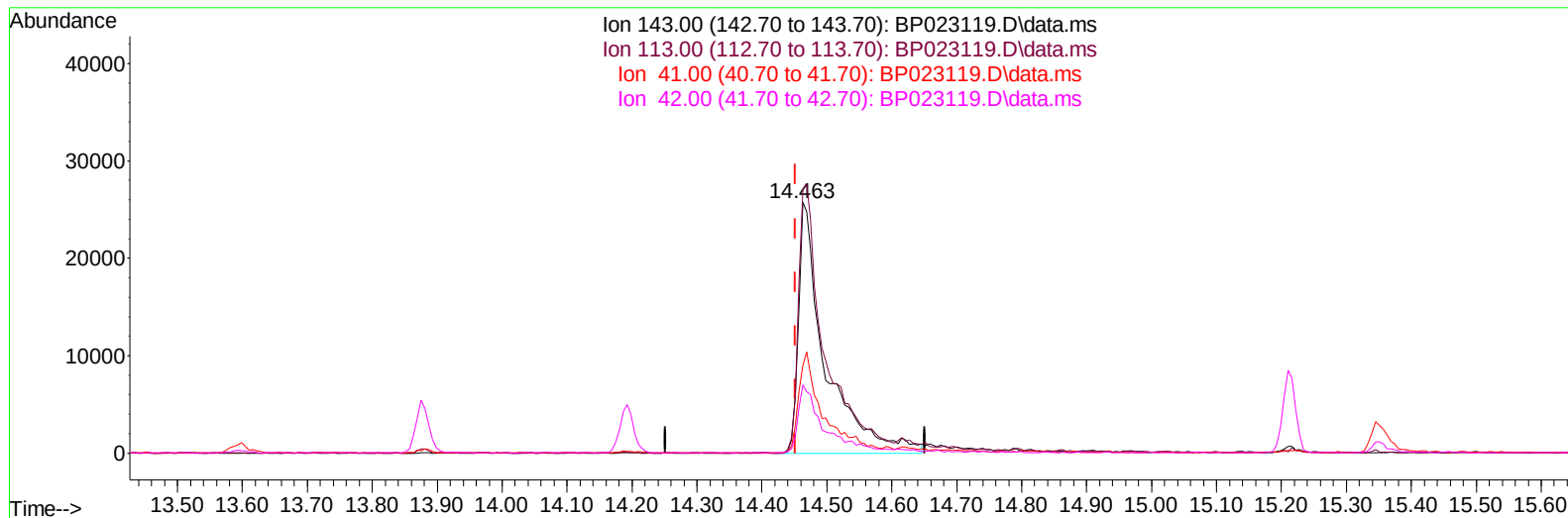
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TIC: BP023119.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.463min (+ 0.012) 27.02 ng/ul m

response 74150

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	105.00
41.00	39.20	35.01
42.00	25.90	27.09

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 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK889

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/15/2024
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Quant Time: Nov 15 01:44:37 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	78586	20.000	ng/ul	0.00
20) Naphthalene-d8	10.328	136	299333	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.192	164	253335	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.004	188	643406	20.000	ng/ul	-0.01
79) Chrysene-d12	21.439	240	721226	20.000	ng/ul	0.00
88) Perylene-d12	24.662	264	789186	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	9742	5.959	ng/uL	0.00
4) Pyridine-d5	3.546	84	129689	26.699	ng/ul	0.00
7) Phenol-d5	6.787	99	159523	27.456	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	99541	29.161	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	125443	29.480	ng/ul	0.00
15) 4-Methylphenol-d8	8.287	113	135240	28.286	ng/ul	0.00
21) Nitrobenzene-d5	8.728	128	65214	31.295	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	79192	31.876	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	153590	29.521	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131	175602	28.292	ng/ul	0.00
46) Dimethylphthalate-d6	13.592	166	607430	33.268	ng/ul	-0.02
49) Acenaphthylene-d8	13.875	160	598973	31.440	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.463	143	74150m	27.023	ng/ul	0.01
60) Fluorene-d10	15.216	176	522319	32.628	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.351	200	103210	26.778	ng/ul	-0.01
73) Anthracene-d10	17.098	188	924145	33.923	ng/ul	-0.02
81) Pyrene-d10	19.486	212	1144395	34.449	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.439	264	1253798	33.353	ng/ul	0.00

Target Compounds Qvalue

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

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