

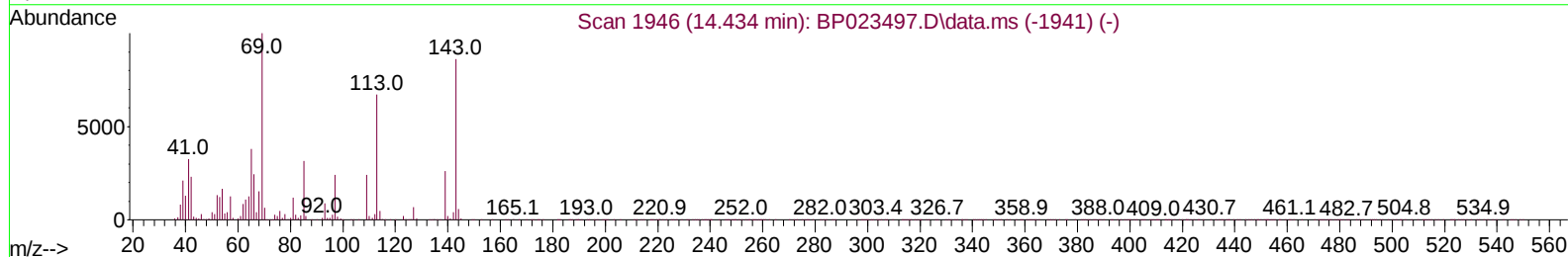
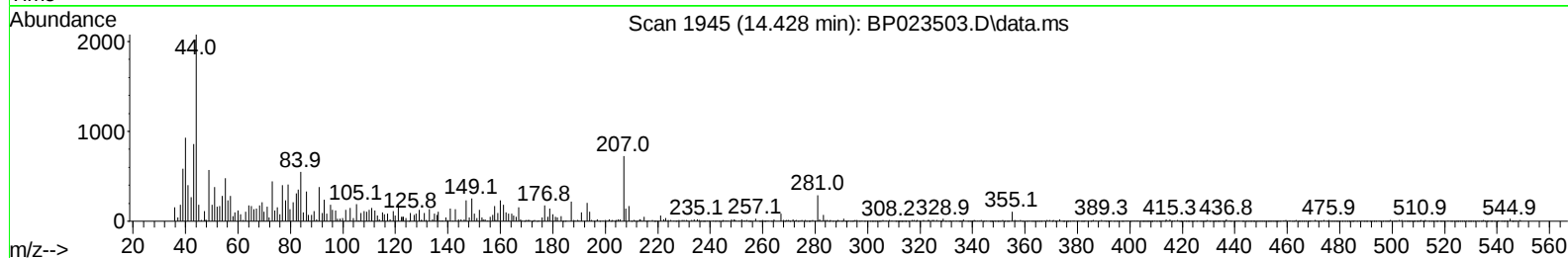
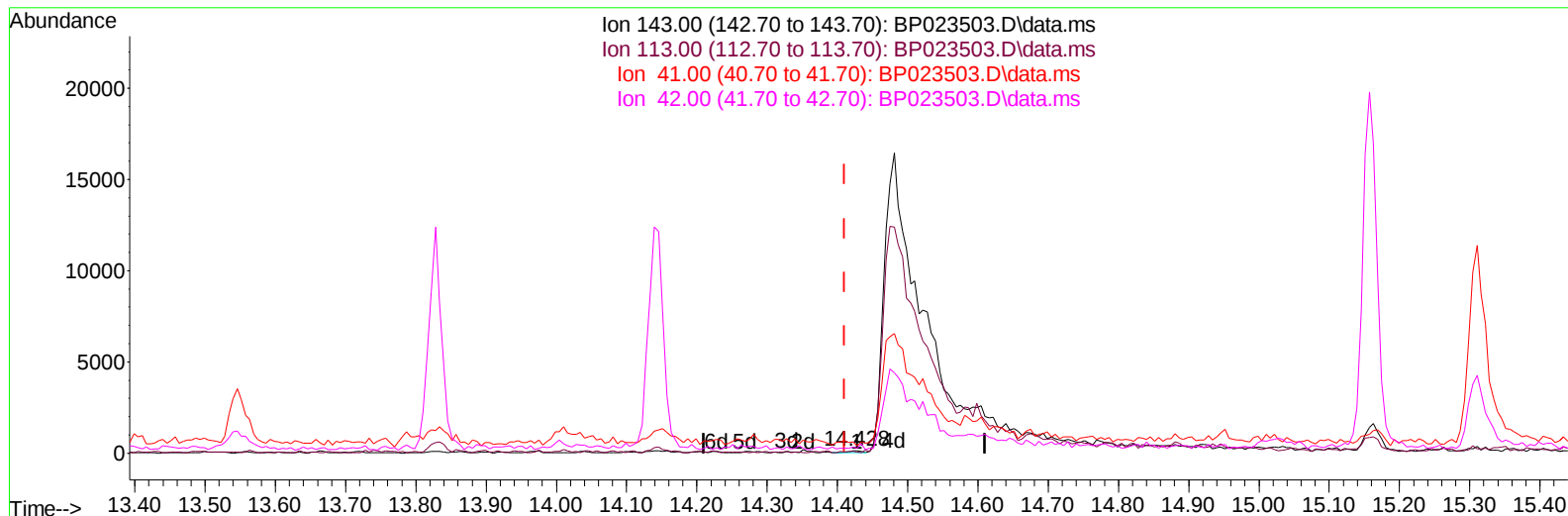
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023503.D
Acq On : 18 Dec 2024 18:22
Operator : RC/JU
Sample : P5235-04 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AT2

Manual IntegrationsAPPROVED

Quant Time: Dec 19 01:03:45 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 00:30:30 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/21/2024
Supervised By :mohammad ahmed 12/21/2024



TIC: BP023503.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.428min (+ 0.018) 0.02 ng/u1

response 150

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	48.89#
41.00	45.60	299.26#
42.00	31.70	200.74#

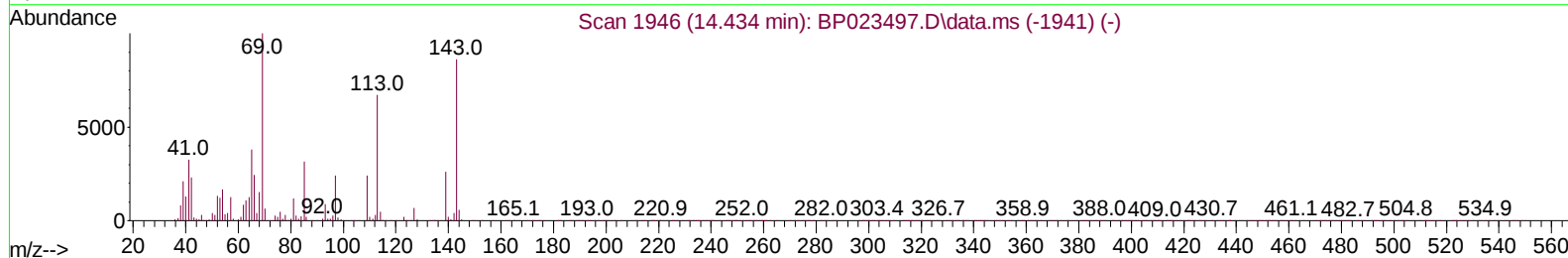
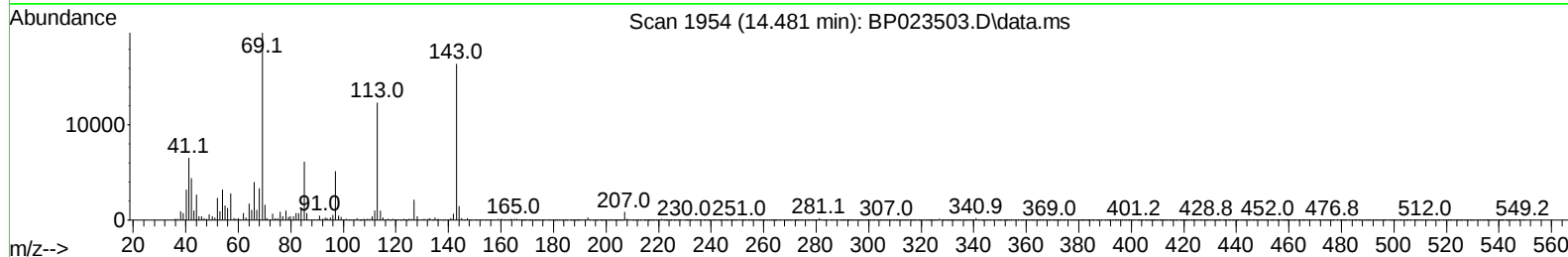
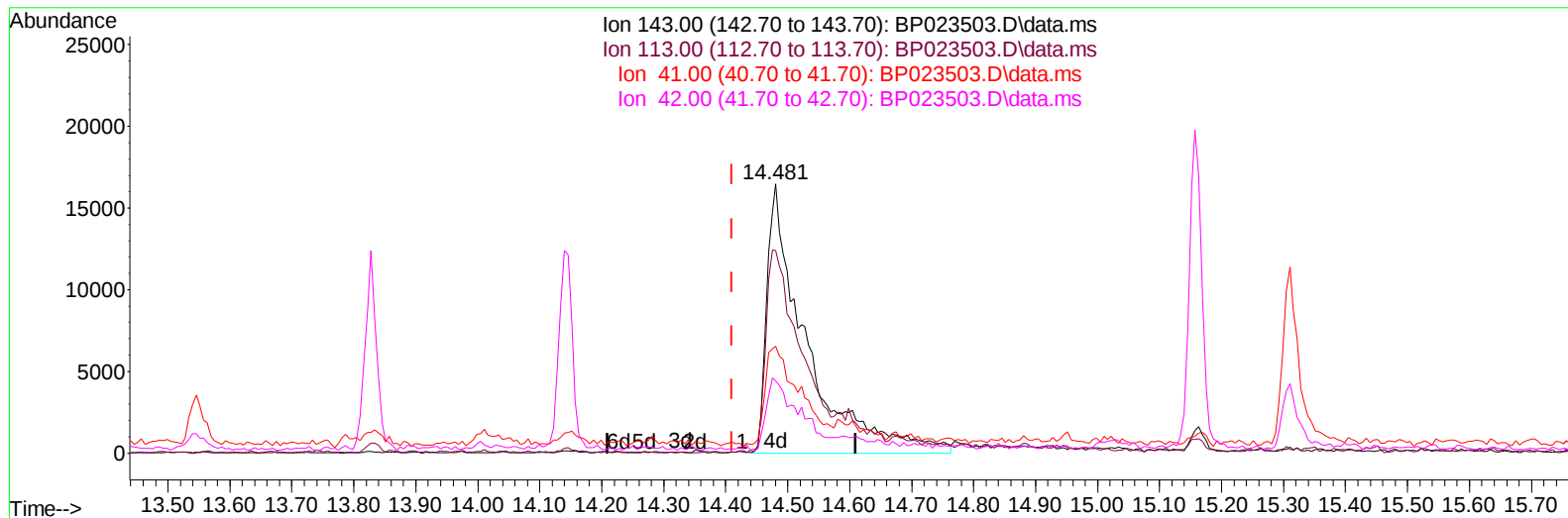
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(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.071) 11.24 ng/ul m

response 72435

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	75.23#
41.00	45.60	39.84
42.00	31.70	26.68

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 Sample : P5235-04 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AT2

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/21/2024
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Quant Time: Dec 19 03:53:11 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.522	152	245460	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.263	136	874532	20.000	ng/ul	#-0.02
38) Acenaphthene-d10	14.140	164	677272	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.951	188	1491365	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.380	240	1616145	20.000	ng/ul	# 0.00
88) Perylene-d12	24.551	264	1655270	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	27005	4.050	ng/uL	-0.01
4) Pyridine-d5	3.517	84	140513	8.196	ng/ul	0.00
7) Phenol-d5	6.752	99	164774	8.556	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.875	67	356148	29.310	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.081	132	354000	24.904	ng/ul	0.00
15) 4-Methylphenol-d8	8.246	113	310925	20.003	ng/ul	-0.01
21) Nitrobenzene-d5	8.669	128	187327	28.685	ng/ul	-0.02
24) 2-Nitrophenol-d4	9.375	143	226835	30.139	ng/ul	-0.02
28) 2,4-Dichlorophenol-d3	9.922	165	411287	25.892	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.428	131	413532	23.238	ng/ul	-0.01
46) Dimethylphthalate-d6	13.546	166	1615140	30.720	ng/ul	-0.01
49) Acenaphthylene-d8	13.828	160	1389283	25.052	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143	72435m	11.243	ng/ul	0.07
60) Fluorene-d10	15.157	176	1286922	27.828	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.310	200	278807	30.933	ng/ul	0.00
73) Anthracene-d10	17.045	188	2436968	35.804	ng/ul	0.00
81) Pyrene-d10	19.439	212	2990966	35.423	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.339	264	2928807	33.053	ng/ul	0.00

Target Compounds Qvalue

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

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