

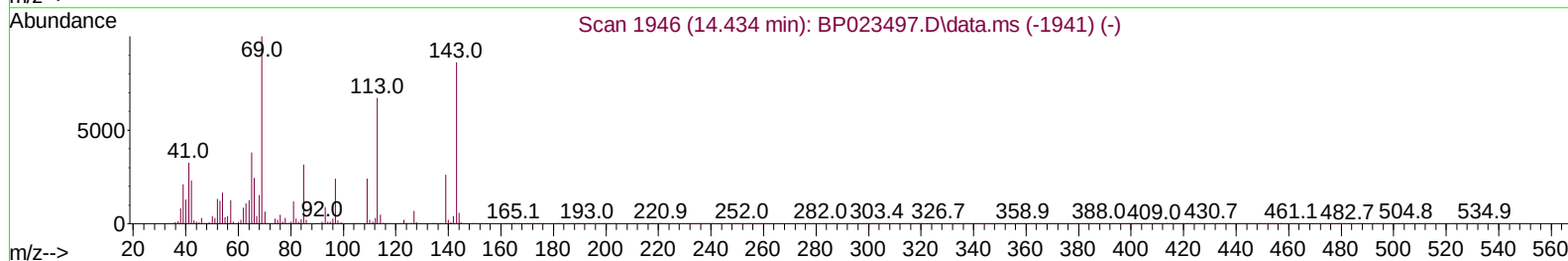
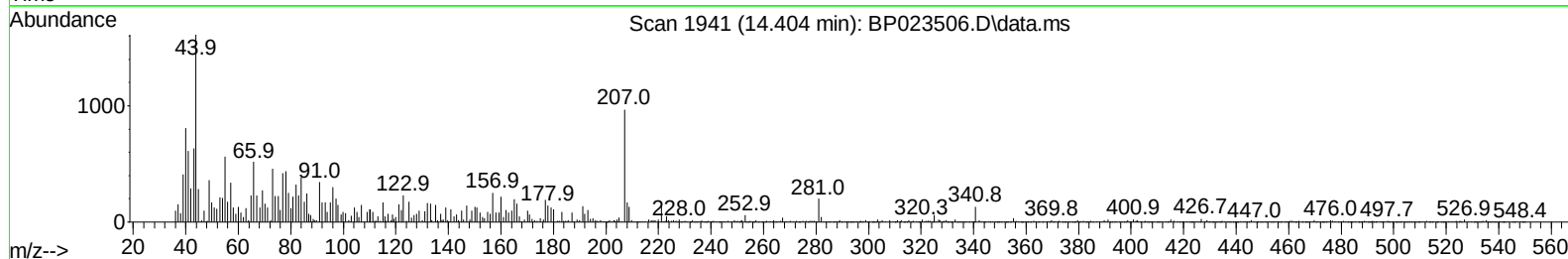
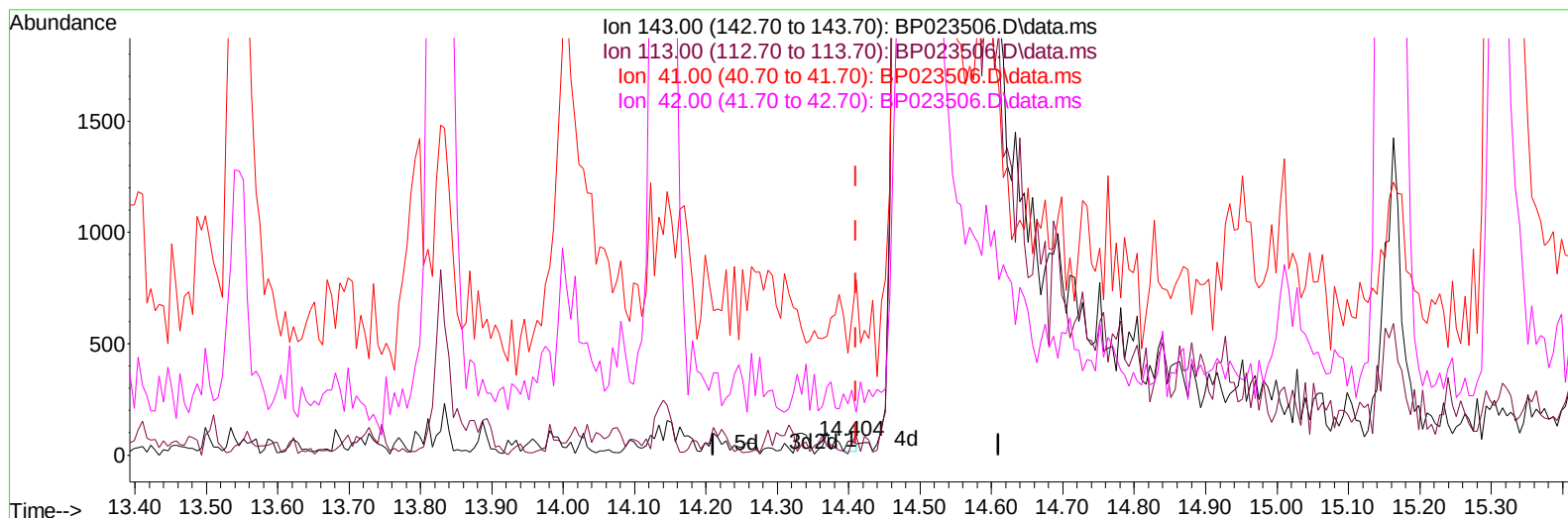
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023506.D
Acq On : 18 Dec 2024 20:23
Operator : RC/JU
Sample : P5235-07 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AT6

Manual IntegrationsAPPROVED

Quant Time: Dec 19 04:01:51 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: BP023506.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.404min (-0.006) 0.01 ng/uL

response 46

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	80.00#
41.00	45.60	941.54#
42.00	31.70	447.69#

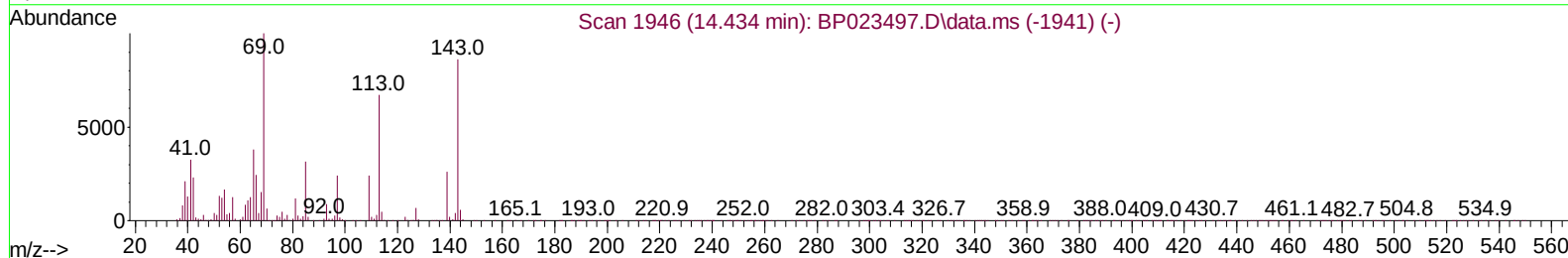
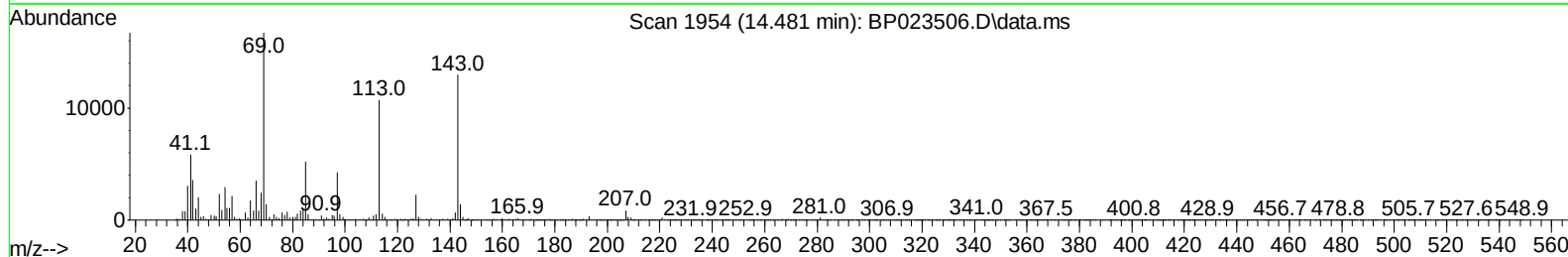
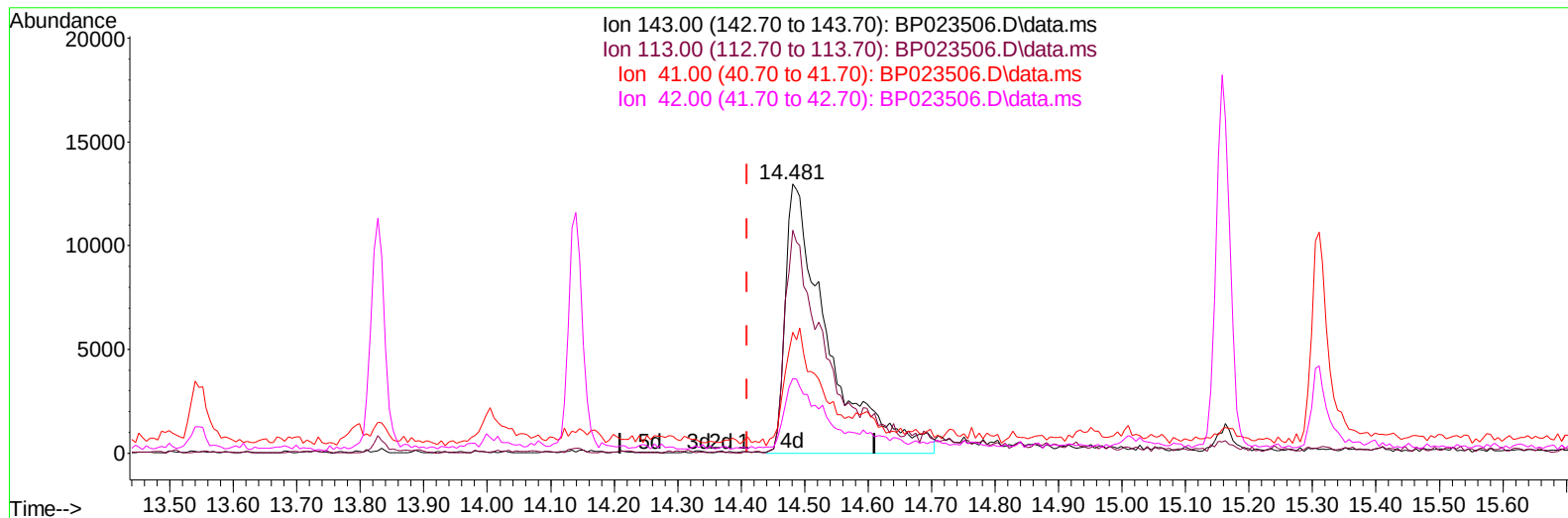
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(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.071) 9.98 ng/ul m

response 60489

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	82.81#
41.00	45.60	44.95
42.00	31.70	27.72

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 Sample : P5235-07 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/21/2024
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Quant Time: Dec 19 04:41:12 2024
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 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	236674	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.269	136	886893	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.140	164	637282	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.951	188	1409245	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.375	240	1479703	20.000	ng/ul	#-0.01
88) Perylene-d12	24.551	264	1531271	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.099	96	26659	4.147	ng/uL	0.00
4) Pyridine-d5	3.523	84	155624	9.414	ng/ul	0.00
7) Phenol-d5	6.752	99	168476	9.072	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.881	67	398034	33.973	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.081	132	384159	28.029	ng/ul	0.00
15) 4-Methylphenol-d8	8.246	113	301601	20.123	ng/ul	-0.01
21) Nitrobenzene-d5	8.669	128	204873	30.935	ng/ul	-0.02
24) 2-Nitrophenol-d4	9.381	143	240375	31.492	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.928	165	446415	27.711	ng/ul	0.00
31) 4-Chloroaniline-d4	10.428	131	459384	25.455	ng/ul	-0.01
46) Dimethylphthalate-d6	13.546	166	1590201	32.143	ng/ul	-0.01
49) Acenaphthylene-d8	13.828	160	1495347	28.656	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143	60489m	9.978	ng/ul	0.07
60) Fluorene-d10	15.163	176	1312614	30.165	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.310	200	264478	31.053	ng/ul	0.00
73) Anthracene-d10	17.045	188	2271499	35.318	ng/ul	0.00
81) Pyrene-d10	19.433	212	2819237	36.468	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.339	264	2817605	34.373	ng/ul	0.00

Target Compounds Qvalue

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

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