

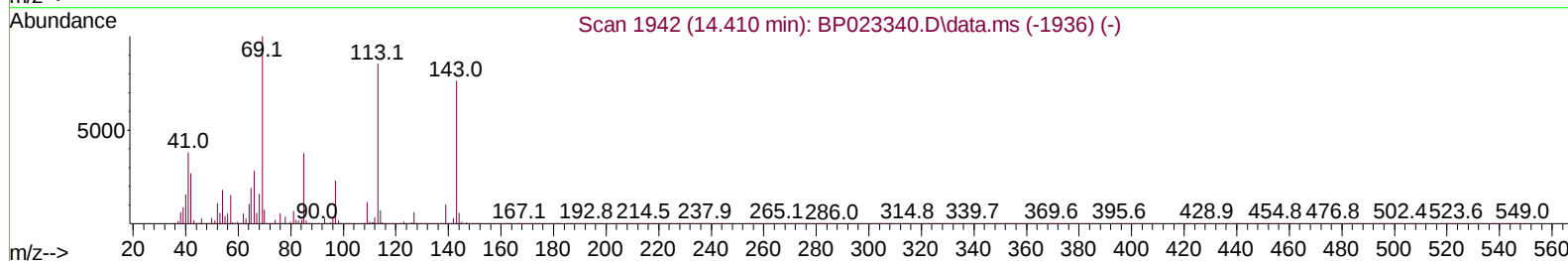
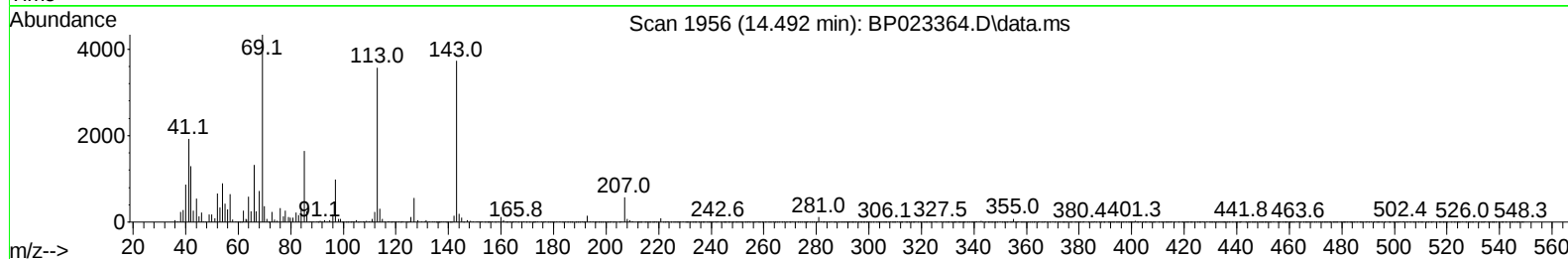
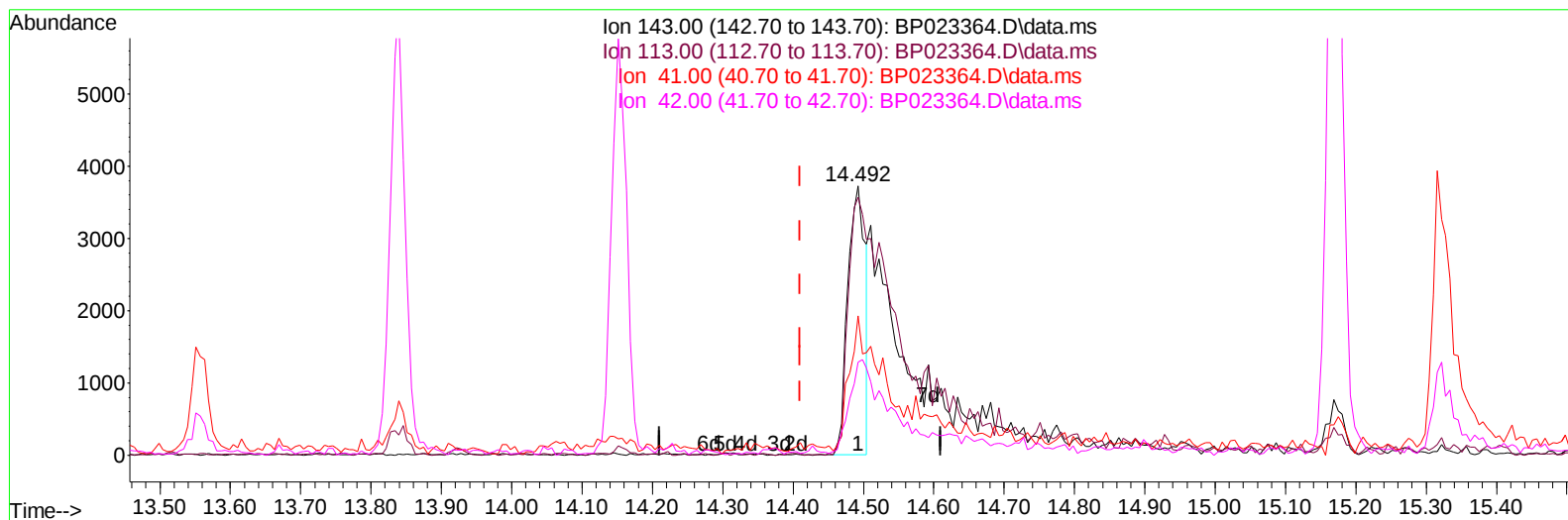
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121024\
Data File : BP023364.D
Acq On : 11 Dec 2024 03:21
Operator : RC/JU
Sample : P5111-11
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BG460

Manual IntegrationsAPPROVED

Quant Time: Dec 11 03:45:49 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 12/12/2024
Supervised By :mohammad ahmed 12/13/2024



TIC: BP023364.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.082) 2.83 ng/u1

response 6426

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	95.86
41.00	45.60	51.66
42.00	31.70	34.67

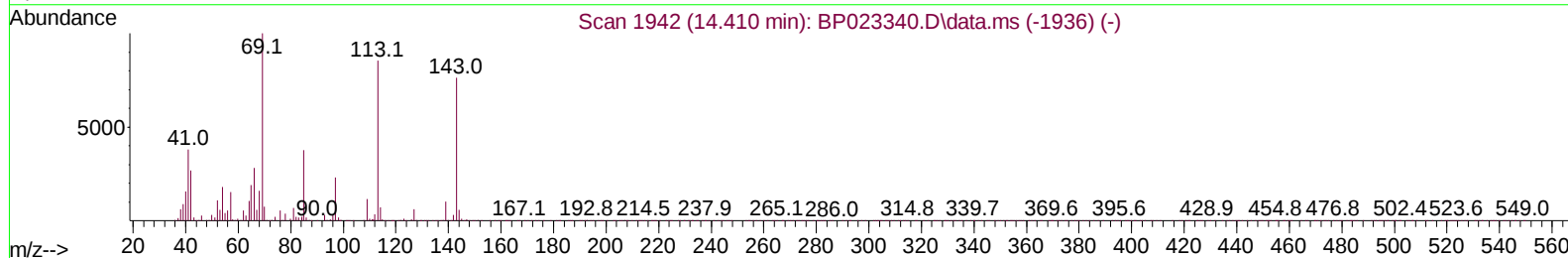
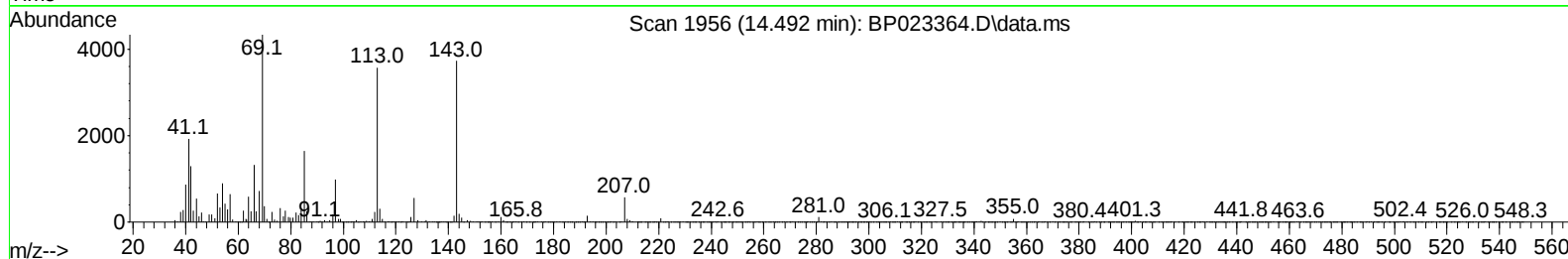
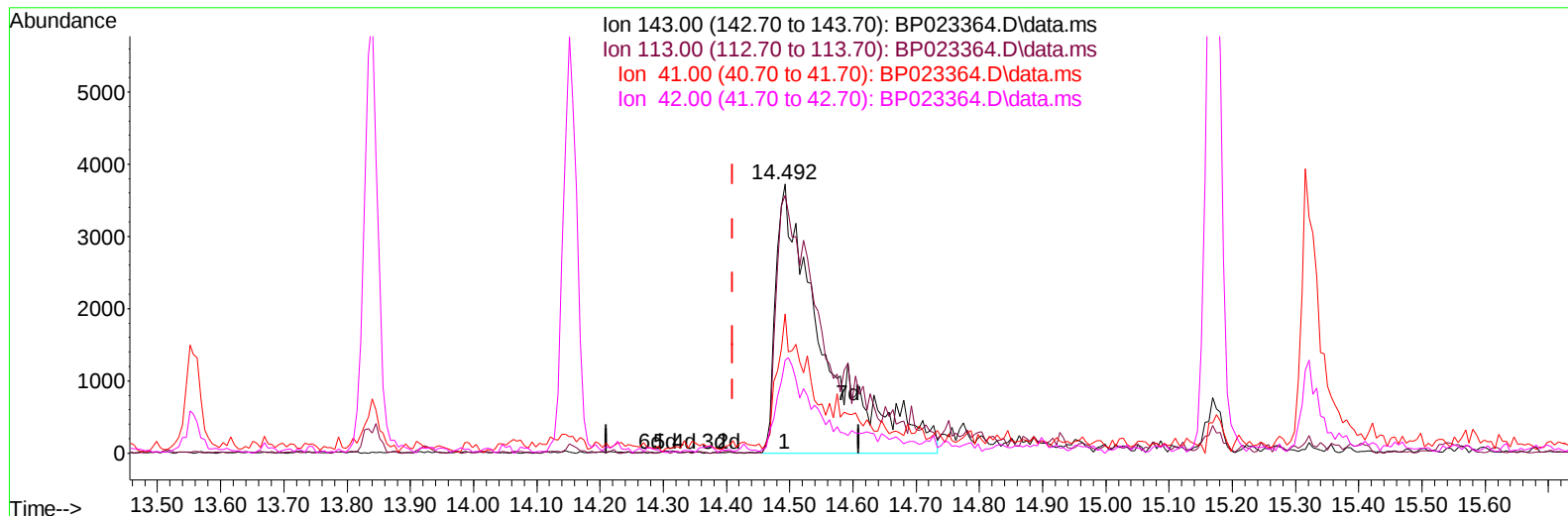
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(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.082) 8.73 ng/ul m

response 19815

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	95.86
41.00	45.60	51.66
42.00	31.70	34.67

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/12/2024
 Supervised By :mohammad ahmed 12/13/2024

Quant Time: Dec 11 03:46:35 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	7.540	152	85589	20.000	ng/ul	0.00
20) Naphthalene-d8	10.281	136	326249	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.157	164	238638	20.000	ng/ul	0.01
64) Phenanthrene-d10	16.957	188	563516	20.000	ng/ul	0.00
79) Chrysene-d12	21.386	240	623814	20.000	ng/ul	0.00
88) Perylene-d12	24.551	264	777676	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.099	96	10369	4.460	ng/uL	0.00
4) Pyridine-d5	3.534	84	21343	3.570	ng/ul	0.02
7) Phenol-d5	6.758	99	57287	8.531	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.893	67	141186	33.323	ng/ul	0.00
11) 2-Chlorophenol-d4	7.087	132	137786	27.799	ng/ul	0.00
15) 4-Methylphenol-d8	8.257	113	104677	19.313	ng/ul	0.00
21) Nitrobenzene-d5	8.681	128	75580	31.023	ng/ul	0.00
24) 2-Nitrophenol-d4	9.393	143	90709	32.306	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.940	165	172594	29.125	ng/ul	0.00
31) 4-Chloroaniline-d4	10.446	131	141193	21.268	ng/ul	0.00
46) Dimethylphthalate-d6	13.557	166	610186	32.937	ng/ul	0.00
49) Acenaphthylene-d8	13.839	160	611776	31.308	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	19815m	8.728	ng/ul	0.08
60) Fluorene-d10	15.169	176	514148	31.553	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.322	200	94208	27.662	ng/ul	0.01
73) Anthracene-d10	17.057	188	867940	33.748	ng/ul	0.00
81) Pyrene-d10	19.451	212	1101184	33.787	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.351	264	1363319	32.748	ng/ul	0.00

Target Compounds Qvalue

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

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