

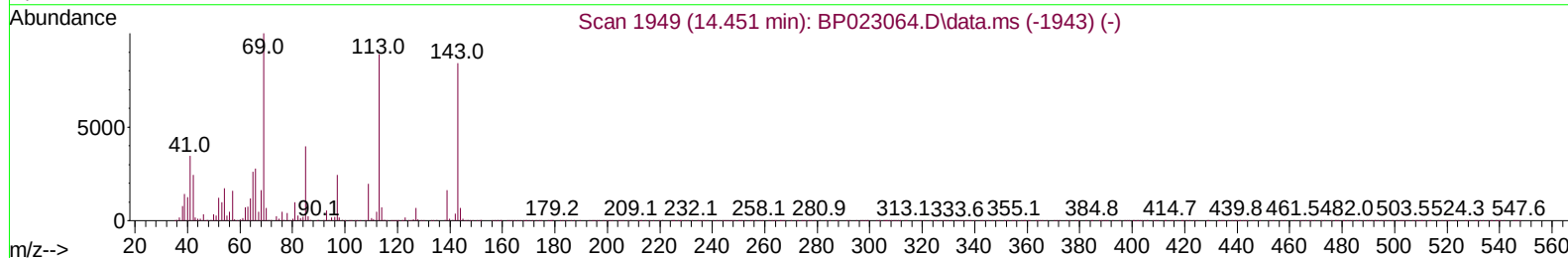
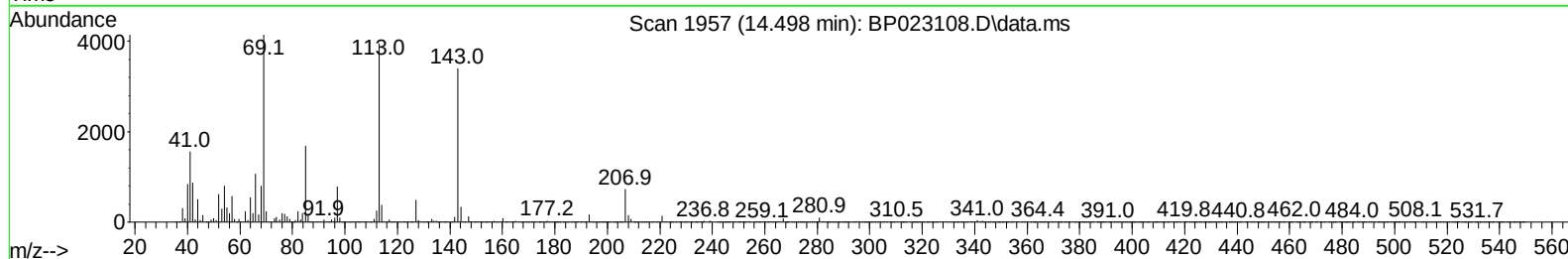
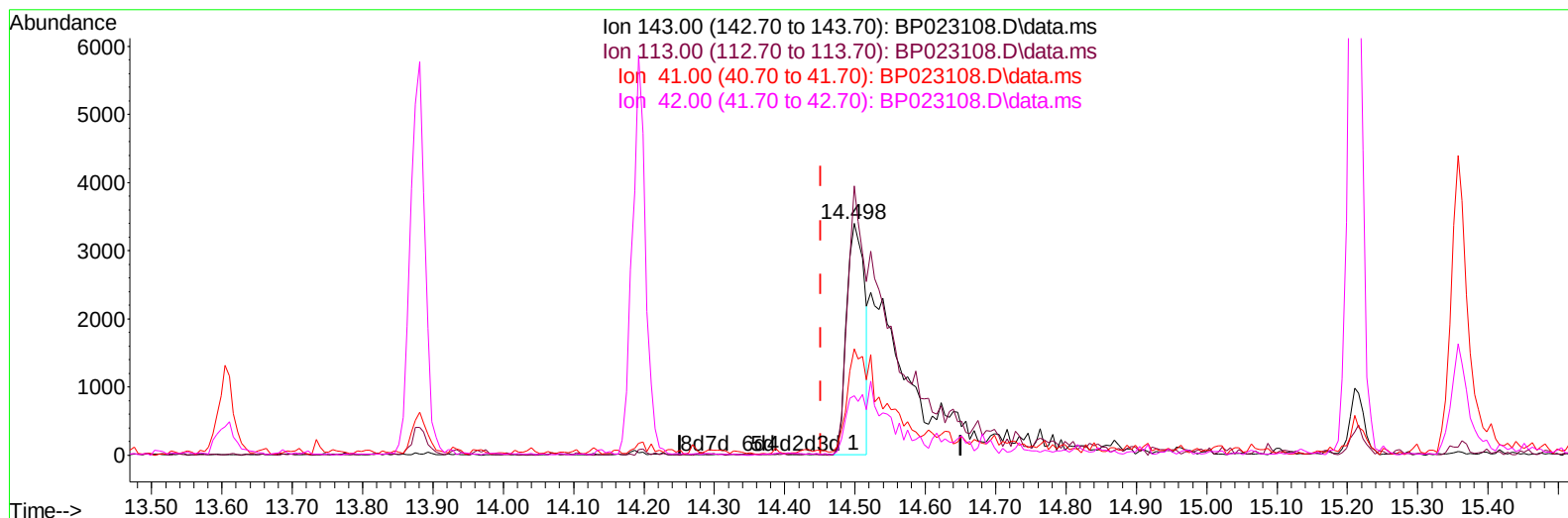
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023108.D
Acq On : 14 Nov 2024 15:03
Operator : RC/JU
Sample : P4777-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH7N7

Manual IntegrationsAPPROVED

Quant Time: Nov 14 23:37:15 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: BP023108.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.047) 2.34 ng/ul

response 5955

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	116.19
41.00	39.20	45.83
42.00	25.90	25.73

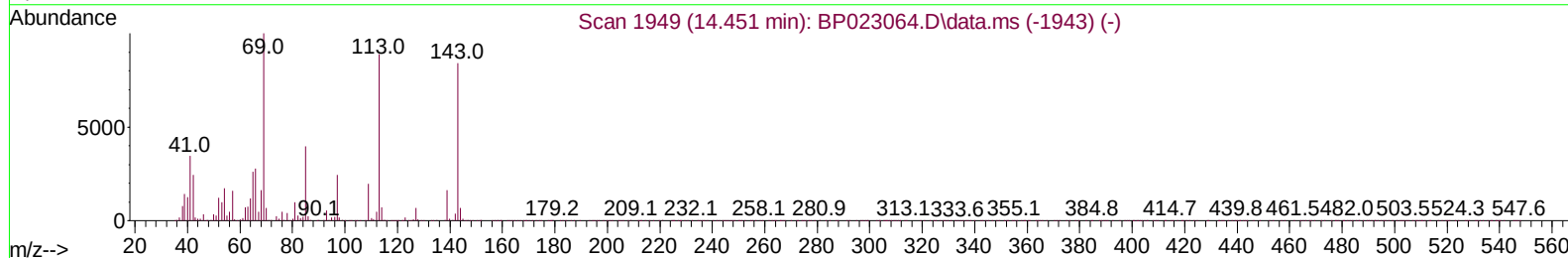
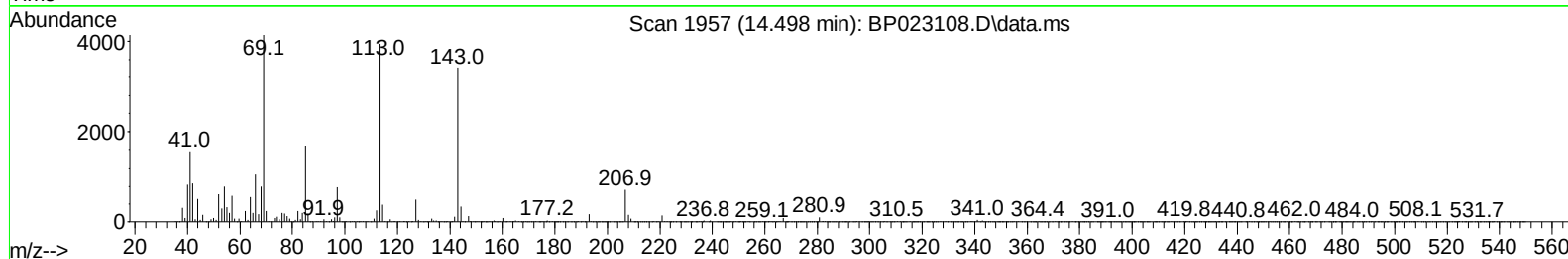
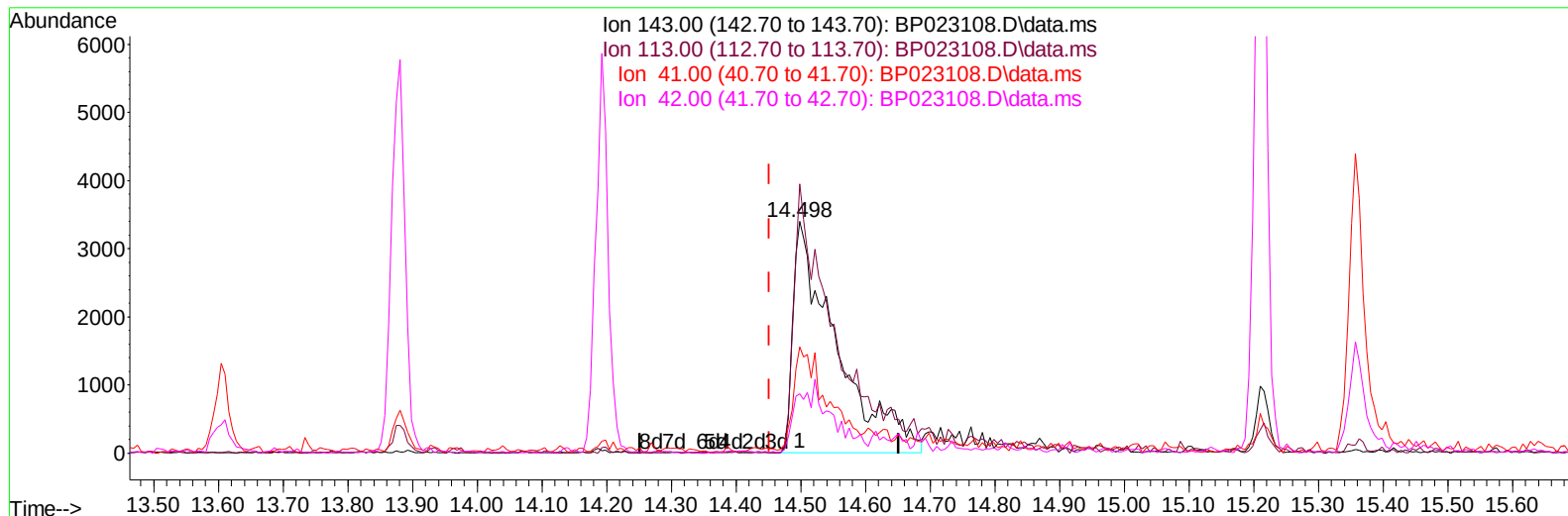
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TIC: BP023108.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.047) 6.22 ng/ul m

response 15809

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	116.19
41.00	39.20	45.83
42.00	25.90	25.73

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Quant Time: Nov 14 23:38:22 2024
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 QLast Update : Wed Nov 13 04:49:01 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	74015	20.000	ng/ul	0.00
20) Naphthalene-d8	10.328	136	282215	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.193	164	234778	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.004	188	585730	20.000	ng/ul	-0.01
79) Chrysene-d12	21.445	240	637318	20.000	ng/ul	0.00
88) Perylene-d12	24.663	264	746522	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	8452	5.489	ng/uL	0.00
4) Pyridine-d5	3.564	84	24356	5.324	ng/ul	0.02
7) Phenol-d5	6.787	99	41214	7.532	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	111842	34.788	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	118833	29.651	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	84288	18.718	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	70042	35.650	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	84735	36.176	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	160851	32.792	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131	131282	22.434	ng/ul	0.00
46) Dimethylphthalate-d6	13.604	166	623911	36.872	ng/ul	0.00
49) Acenaphthylene-d8	13.881	160	624881	35.392	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.498	143	15809m	6.217	ng/ul	0.05
60) Fluorene-d10	15.210	176	550792	37.126	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.357	200	110086	31.374	ng/ul	0.00
73) Anthracene-d10	17.098	188	954516	38.488	ng/ul	-0.02
81) Pyrene-d10	19.492	212	1218035	41.493	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.439	264	1423758	40.038	ng/ul	0.00

Target Compounds Qvalue

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

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