

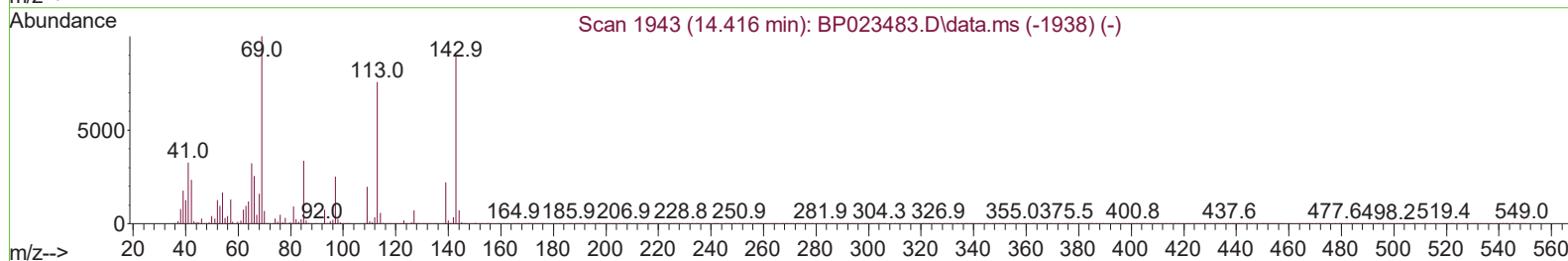
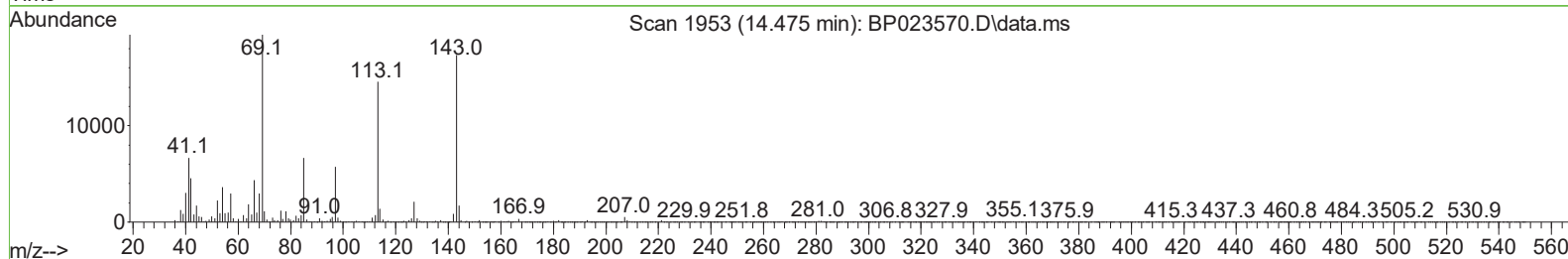
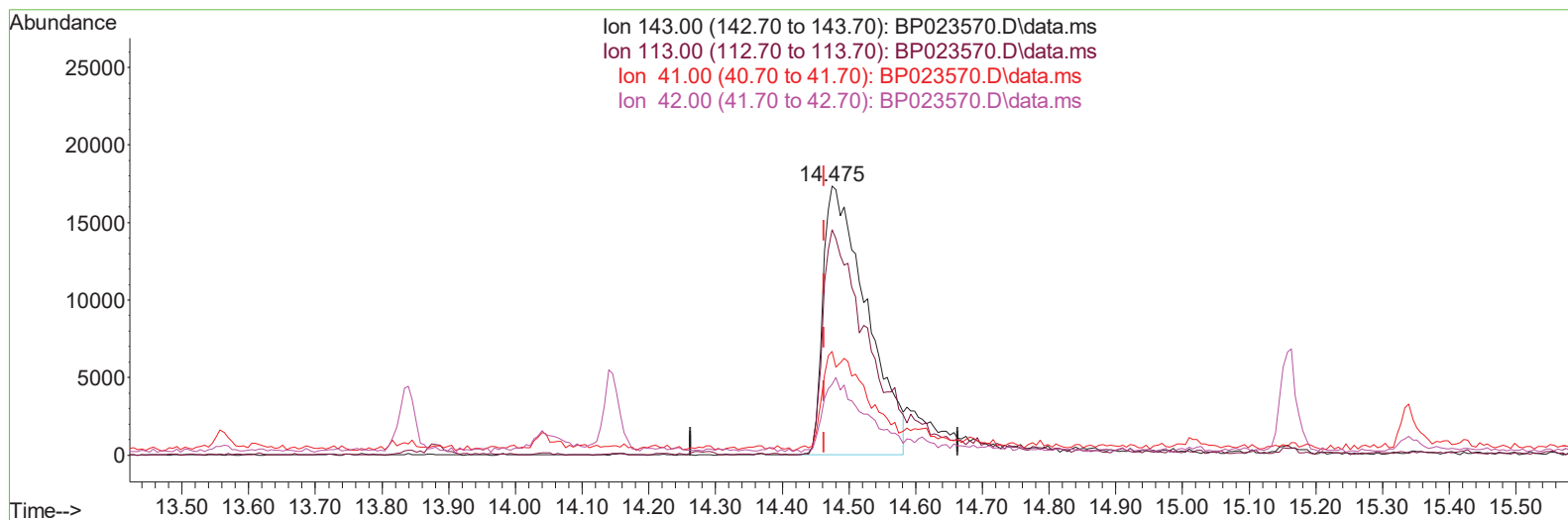
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122324\
Data File : BP023570.D
Acq On : 23 Dec 2024 19:45
Operator : RC/JU
Sample : P5302-17
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E2AQ0

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/30/2024
Supervised By :mohammad ahmed 12/30/2024

Quant Time: Dec 24 04:48:24 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 24 02:14:57 2024
Response via : Initial Calibration



TIC: BP023570.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.475min (+ 0.012) 25.52 ng/ul

response 78260

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	83.80#
41.00	45.60	38.51
42.00	31.70	26.36

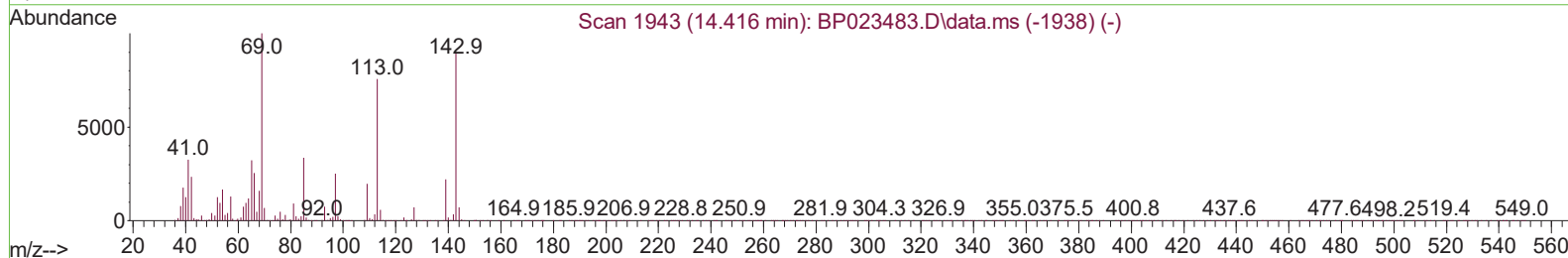
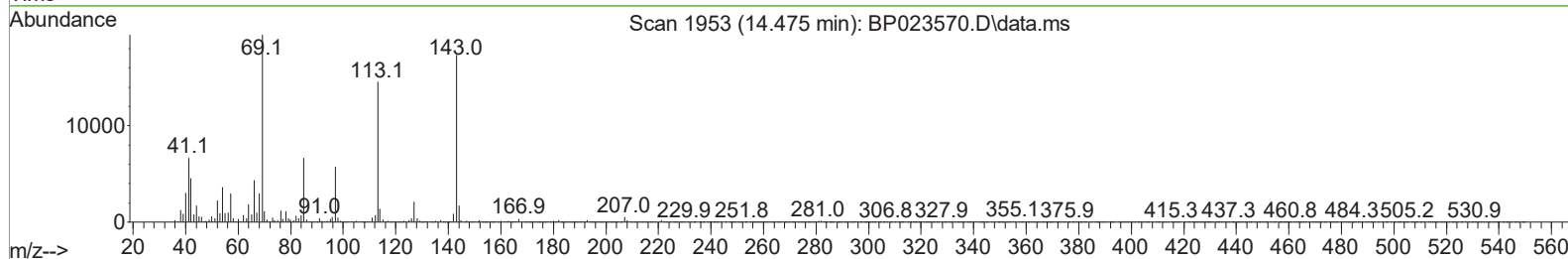
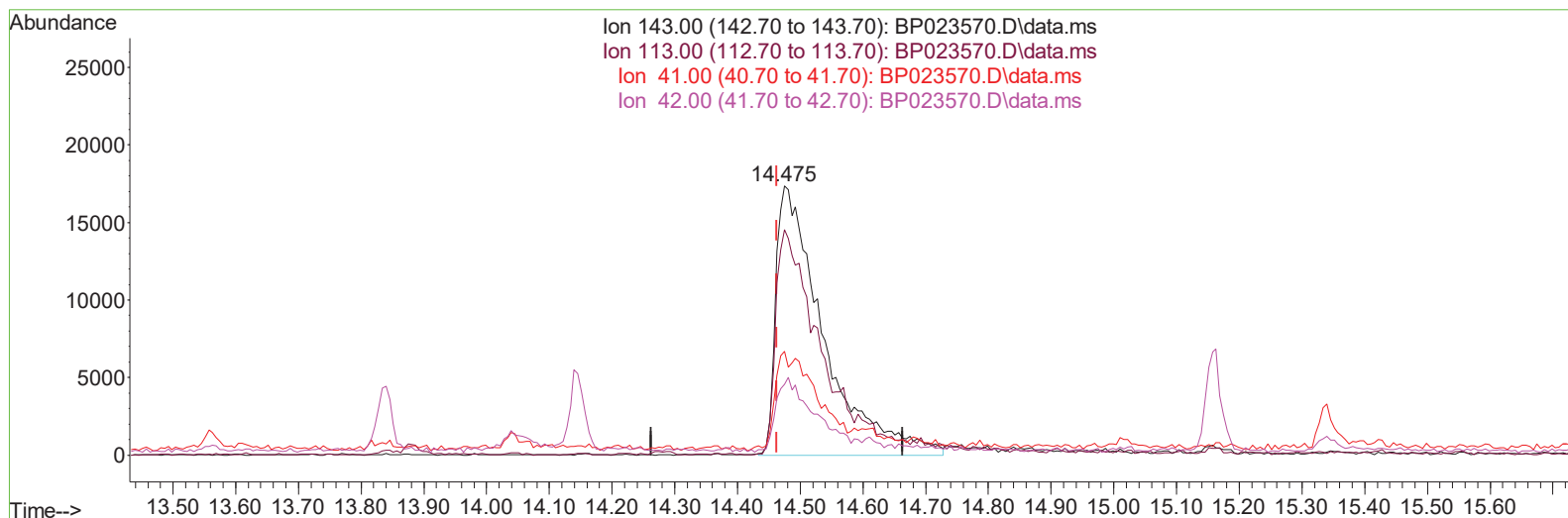
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TIC: BP023570.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.475min (+ 0.012) 29.79 ng/ul m

response 91351

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.10	83.80#
41.00	45.60	38.51
42.00	31.70	26.36

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Quant Time: Dec 24 04:50:30 2024
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 QLast Update : Tue Dec 24 02:14:57 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.522	152	126637	20.000	ng/u1	0.02
20) Naphthalene-d8	10.275	136	455256	20.000	ng/u1	# 0.02
38) Acenaphthene-d10	14.145	164	322299	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.963	188	686425	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.392	240	824497	20.000	ng/u1	# 0.00
88) Perylene-d12	24.574	264	963951	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	14189	4.125	ng/uL	0.00
4) Pyridine-d5	3.505	84	153664	17.373	ng/u1	0.02
7) Phenol-d5	6.746	99	249318	25.092	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	6.875	67	155378	24.786	ng/u1	0.02
11) 2-Chlorophenol-d4	7.075	132	175839	23.977	ng/u1	0.02
15) 4-Methylphenol-d8	8.252	113	196070	24.449	ng/u1	0.02
21) Nitrobenzene-d5	8.675	128	85967	25.288	ng/u1	0.02
24) 2-Nitrophenol-d4	9.387	143	103382	26.386	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	9.940	165	191623	23.173	ng/u1	0.02
31) 4-Chloroaniline-d4	10.434	131	233855	25.244	ng/u1	0.01
46) Dimethylphthalate-d6	13.563	166	646295	25.831	ng/u1	0.00
49) Acenaphthylene-d8	13.840	160	685543	25.977	ng/u1	0.00
54) 4-Nitrophenol-d4	14.475	143	91351m	29.794	ng/u1	0.01
60) Fluorene-d10	15.163	176	547262	24.867	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.339	200	74753	18.019	ng/u1	0.00
73) Anthracene-d10	17.063	188	893609	28.525	ng/u1	0.00
81) Pyrene-d10	19.451	212	1102875	25.603	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.351	264	1290743	25.013	ng/u1	-0.02

Target Compounds Qvalue

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

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