

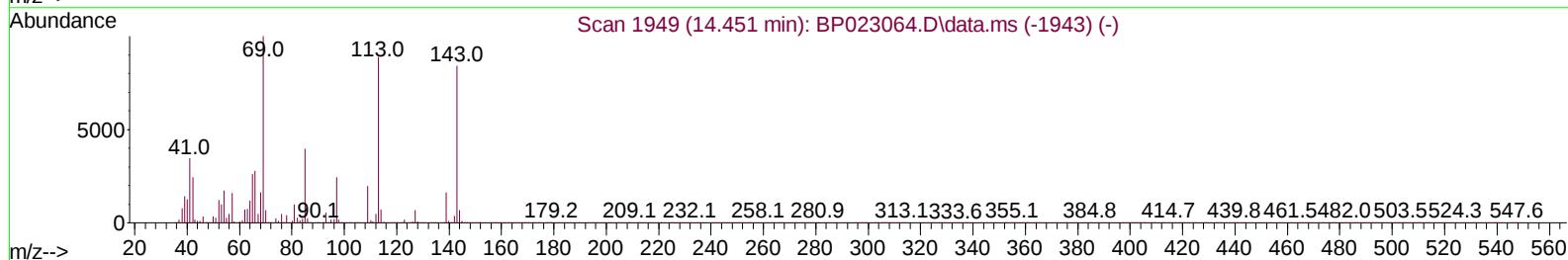
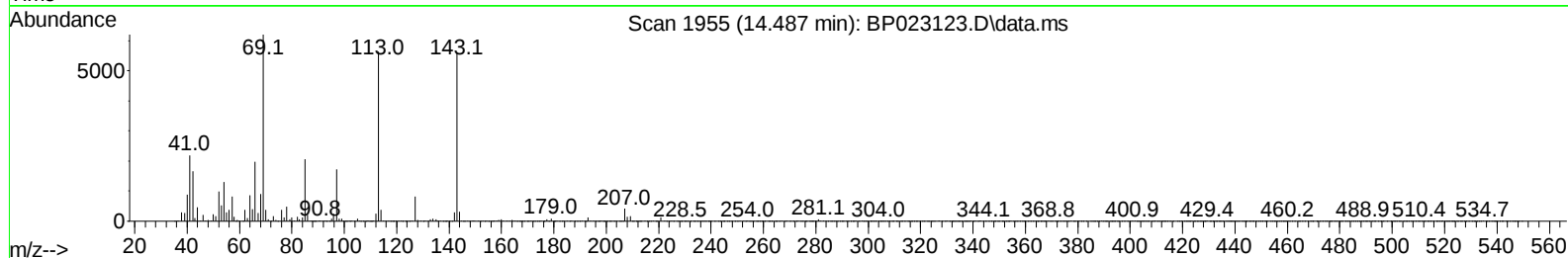
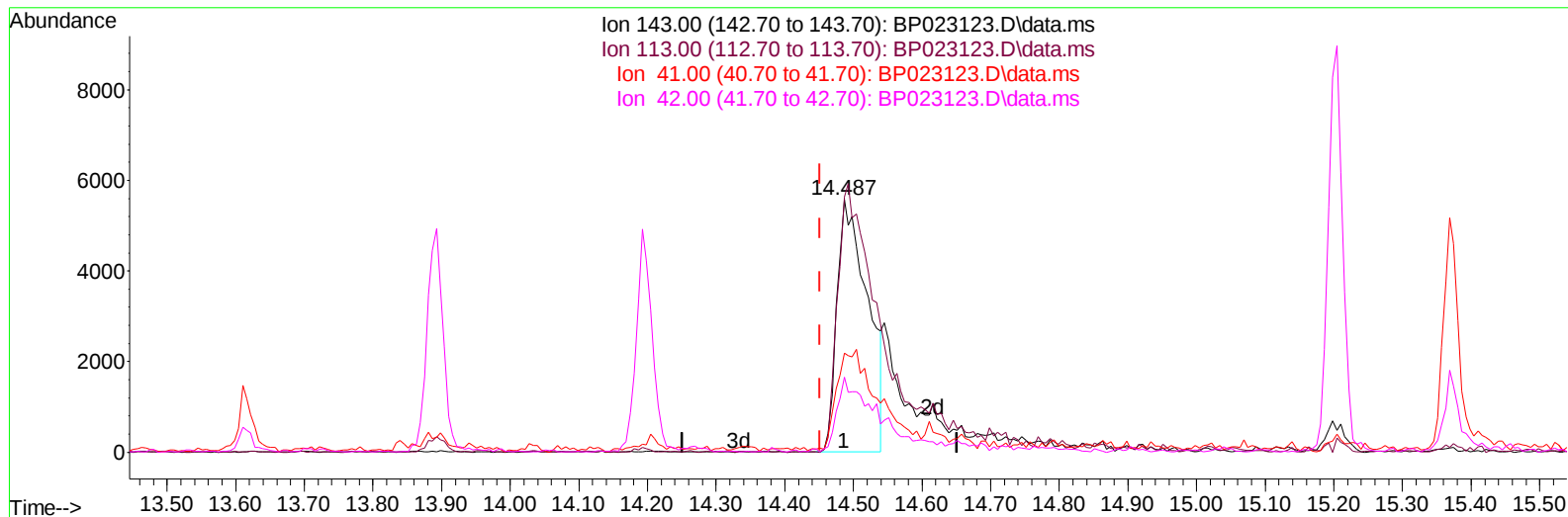
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023123.D
Acq On : 15 Nov 2024 03:55
Operator : RC/JU
Sample : P4801-09
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCPB0

Manual IntegrationsAPPROVED

Quant Time: Nov 15 04:26:53 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: BP023123.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.035) 6.79 ng/u1

response 17273

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	100.79
41.00	39.20	39.18
42.00	25.90	29.57

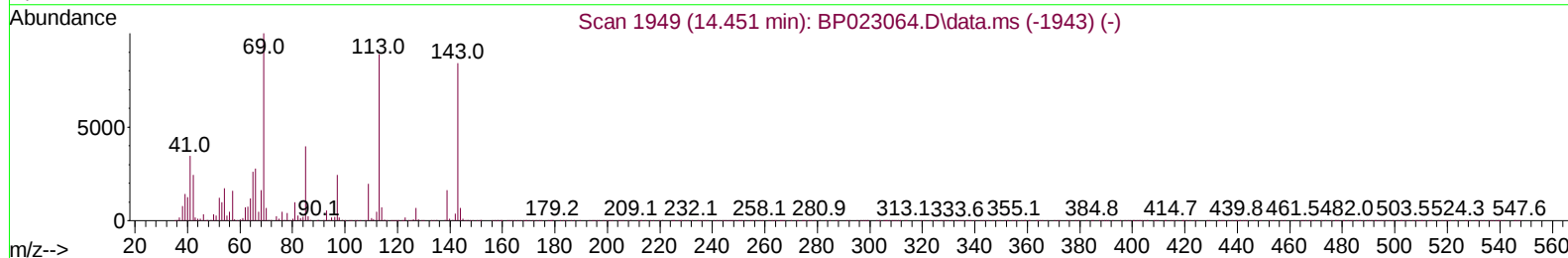
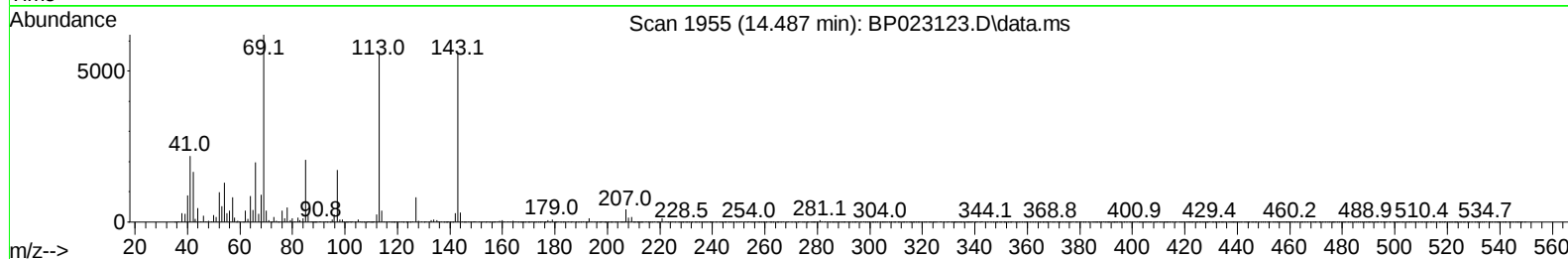
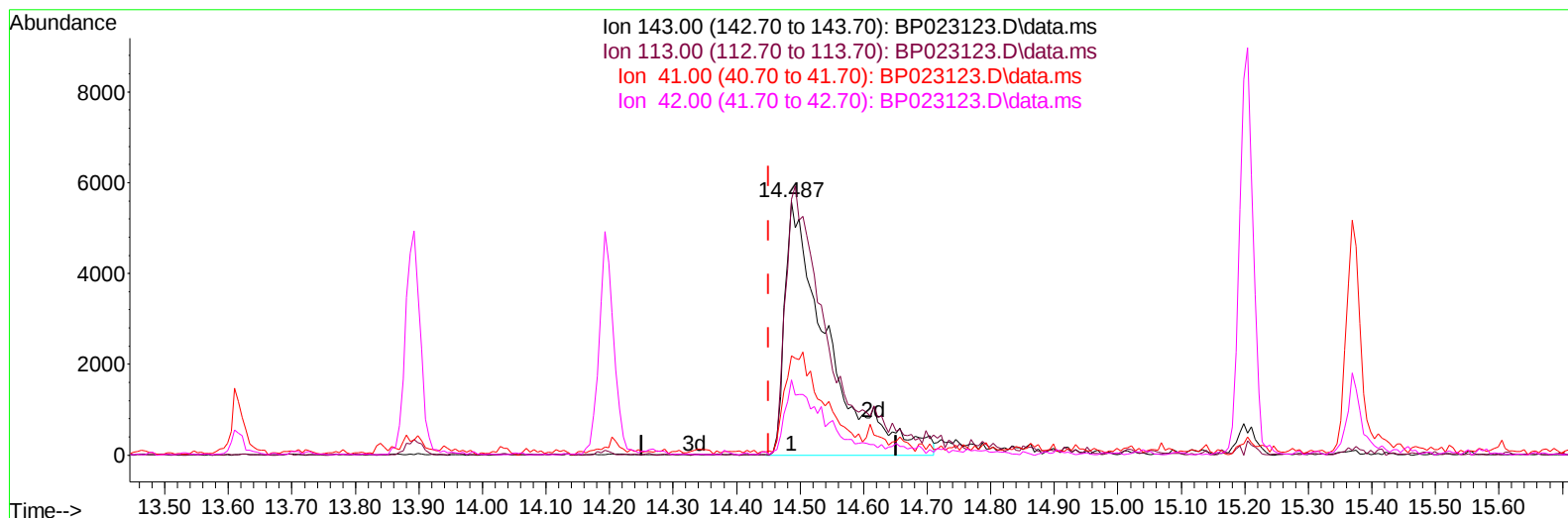
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TIC: BP023123.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.035) 10.20 ng/ul m

response 25940

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	100.79
41.00	39.20	39.18
42.00	25.90	29.57

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Quant Time: Nov 15 04:27:37 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	77282	20.000	ng/ul	0.00
20) Naphthalene-d8	10.322	136	291391	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.198	164	234730	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	611118	20.000	ng/ul	-0.02
79) Chrysene-d12	21.445	240	730391	20.000	ng/ul	0.00
88) Perylene-d12	24.668	264	875230	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	7771	4.833	ng/uL	0.00
4) Pyridine-d5	3.564	84	24243	5.075	ng/ul	0.02
7) Phenol-d5	6.787	99	48325	8.458	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	97551	29.060	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	116009	27.723	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	92192	19.608	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	62104	30.614	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	75948	31.404	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	150849	29.785	ng/ul	0.00
31) 4-Chloroaniline-d4	10.487	131	121752	20.150	ng/ul	0.00
46) Dimethylphthalate-d6	13.616	166	633939	37.472	ng/ul	0.00
49) Acenaphthylene-d8	13.892	160	589288	33.383	ng/ul	0.00
54) 4-Nitrophenol-d4	14.487	143	25940m	10.203	ng/ul	0.04
60) Fluorene-d10	15.204	176	539433	36.368	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.375	200	123996	33.870	ng/ul	0.01
73) Anthracene-d10	17.110	188	978338	37.810	ng/ul	0.00
81) Pyrene-d10	19.492	212	1365185	40.579	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.445	264	1673335	40.137	ng/ul	0.00

Target Compounds Qvalue

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

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