

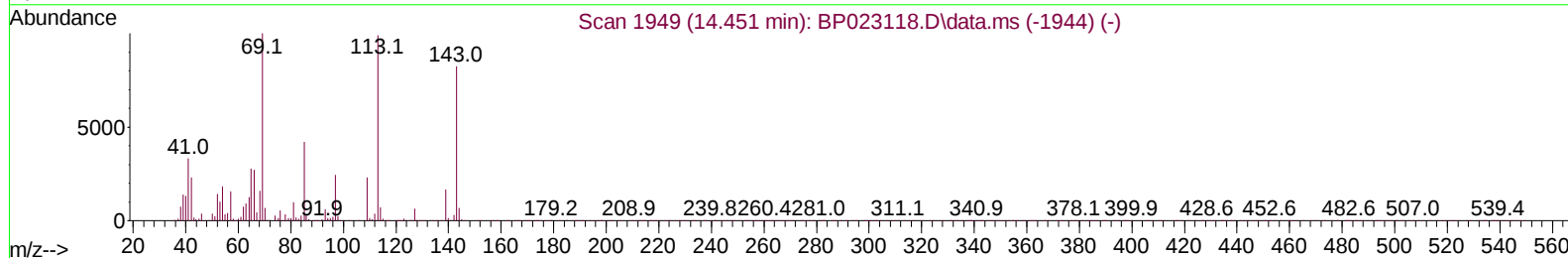
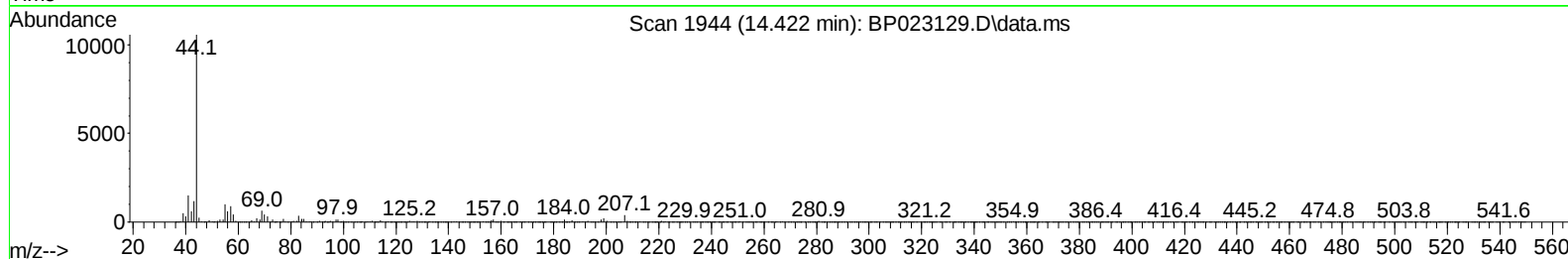
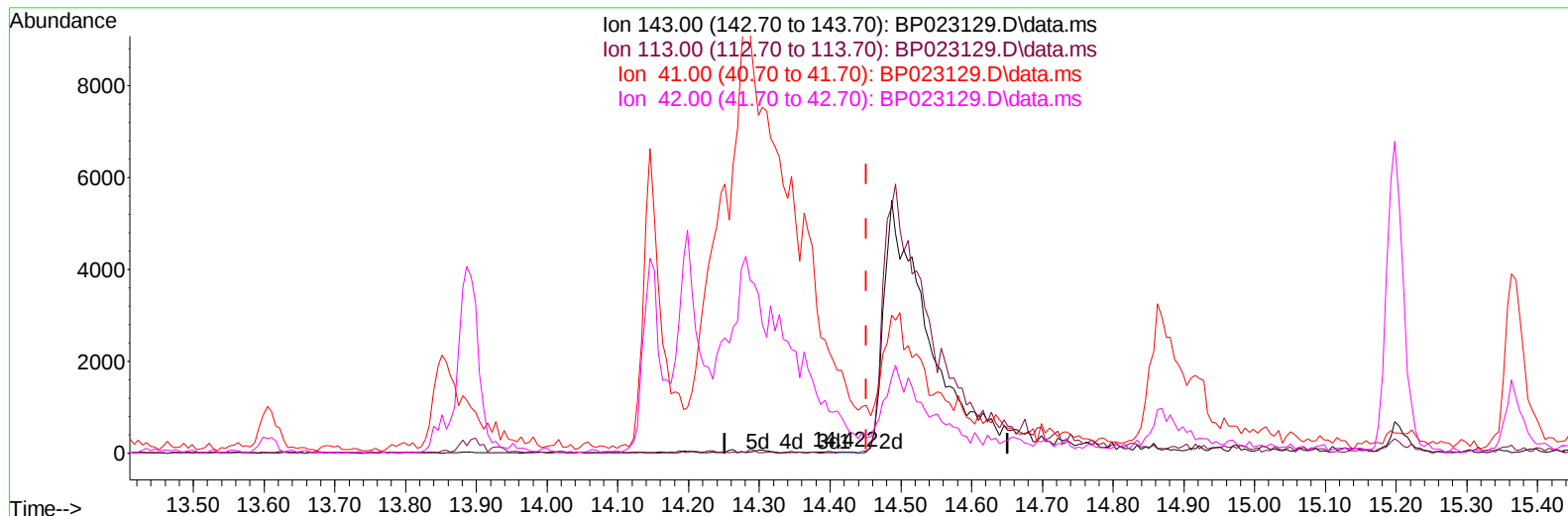
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023129.D
Acq On : 15 Nov 2024 07:58
Operator : RC/JU
Sample : P4784-04
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E27N2

Manual IntegrationsAPPROVED

Quant Time: Nov 15 08:24:05 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: BP023129.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.422min (-0.029) 0.01 ng/u1

response 25

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	72.73#
41.00	39.20	6922.73#
42.00	25.90	2813.64#

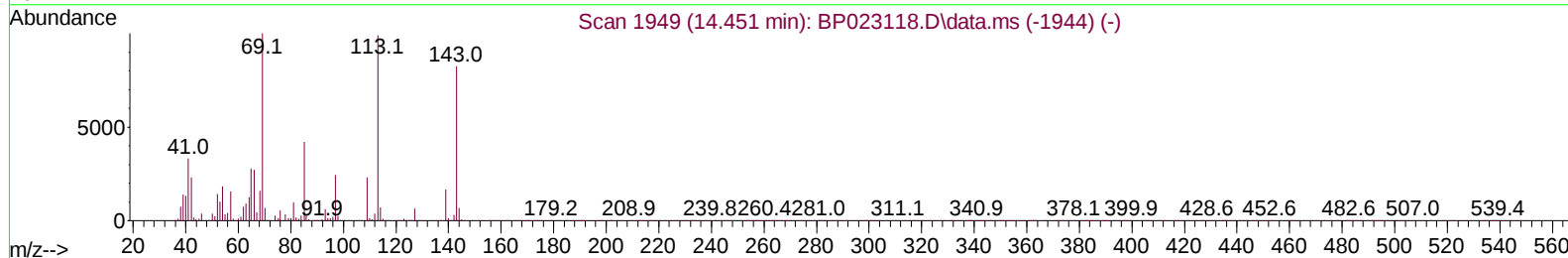
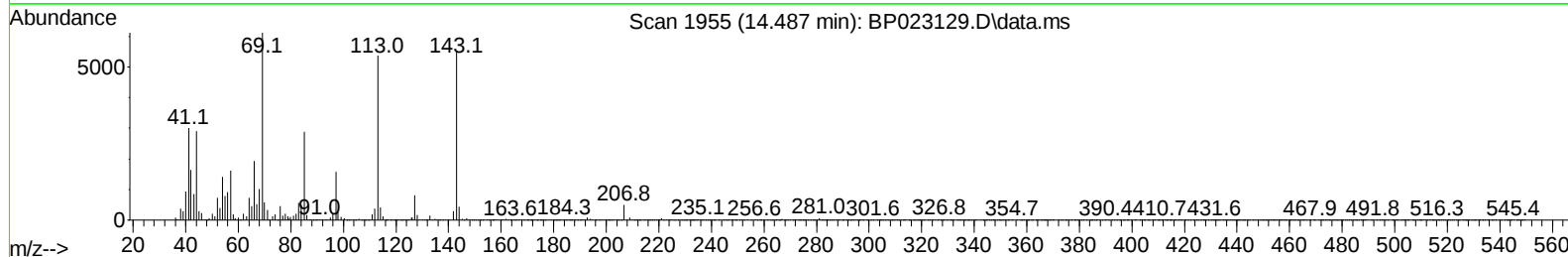
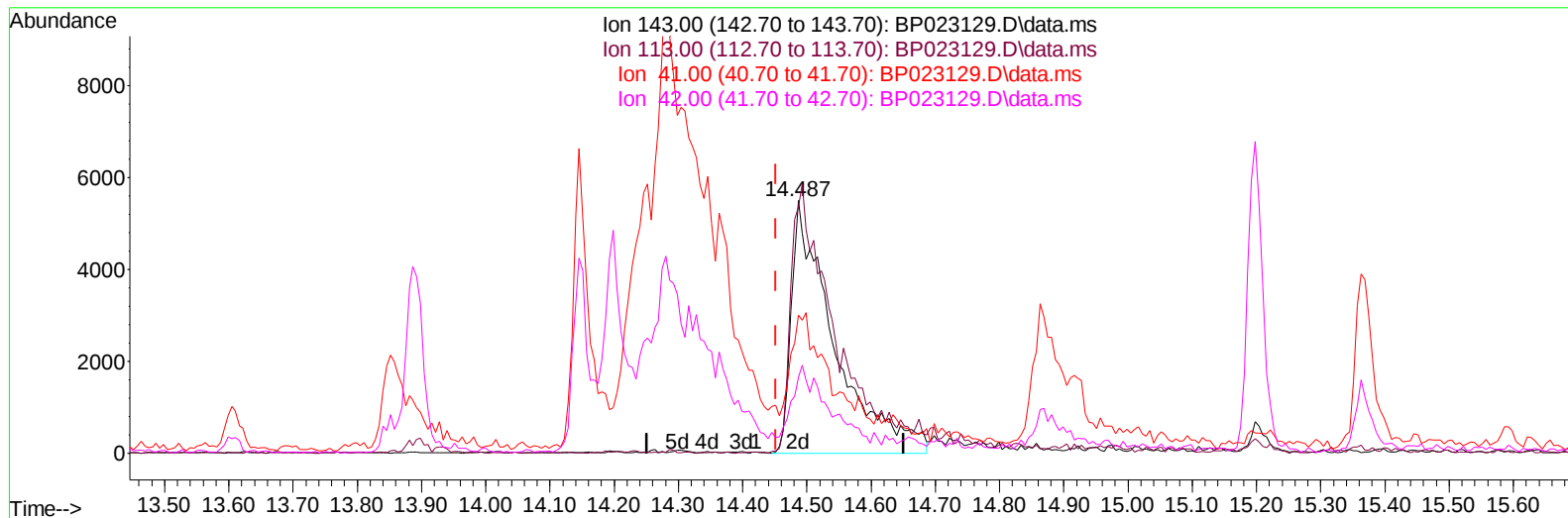
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(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.035) 11.03 ng/ul m

response 25356

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	97.71
41.00	39.20	54.59#
42.00	25.90	29.97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	70033	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136	275584	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.198	164	212294	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	559968	20.000	ng/ul	-0.02
79) Chrysene-d12	21.445	240	707966	20.000	ng/ul	0.00
88) Perylene-d12	24.663	264	849491	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	7664	5.260	ng/uL	0.00
4) Pyridine-d5	3.558	84	36284	8.382	ng/ul	0.01
7) Phenol-d5	6.793	99	56371	10.887	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.922	67	89874	29.544	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	108634	28.648	ng/ul	0.00
15) 4-Methylphenol-d8	8.287	113	98039	23.009	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	57678	30.063	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	69463	30.370	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	149072	31.122	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131	127407	22.296	ng/ul	0.00
46) Dimethylphthalate-d6	13.610	166	528270	34.526	ng/ul	0.00
49) Acenaphthylene-d8	13.893	160	521964	32.694	ng/ul	0.00
54) 4-Nitrophenol-d4	14.487	143	25356m	11.027	ng/ul	0.04
60) Fluorene-d10	15.198	176	467688	34.864	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.363	200	102083	30.432	ng/ul	0.00
73) Anthracene-d10	17.116	188	921423	38.863	ng/ul	0.00
81) Pyrene-d10	19.492	212	1241907	38.084	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.439	264	1552534	38.368	ng/ul	0.00

Target Compounds Qvalue

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

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