

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023139.D
Acq On : 15 Nov 2024 14:44
Operator : RC/JU
Sample : P4784-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

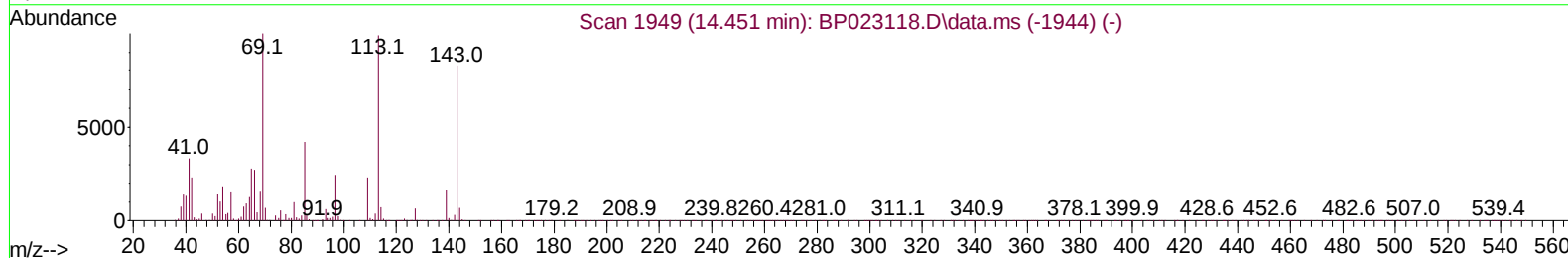
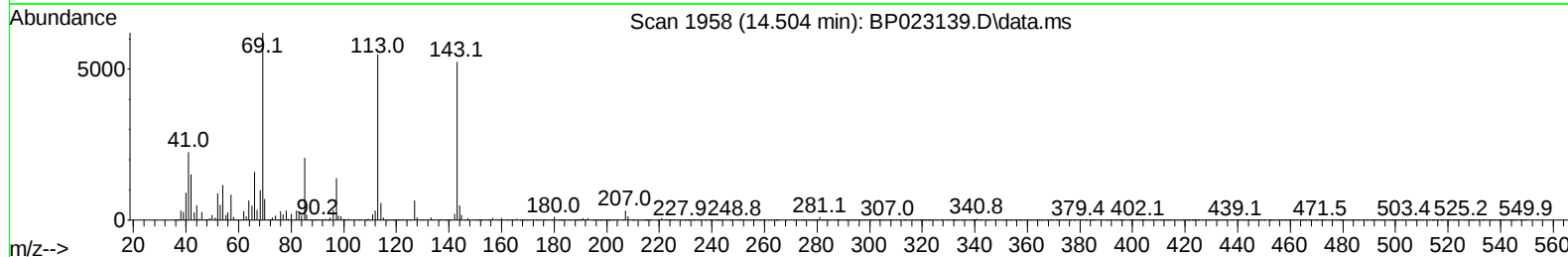
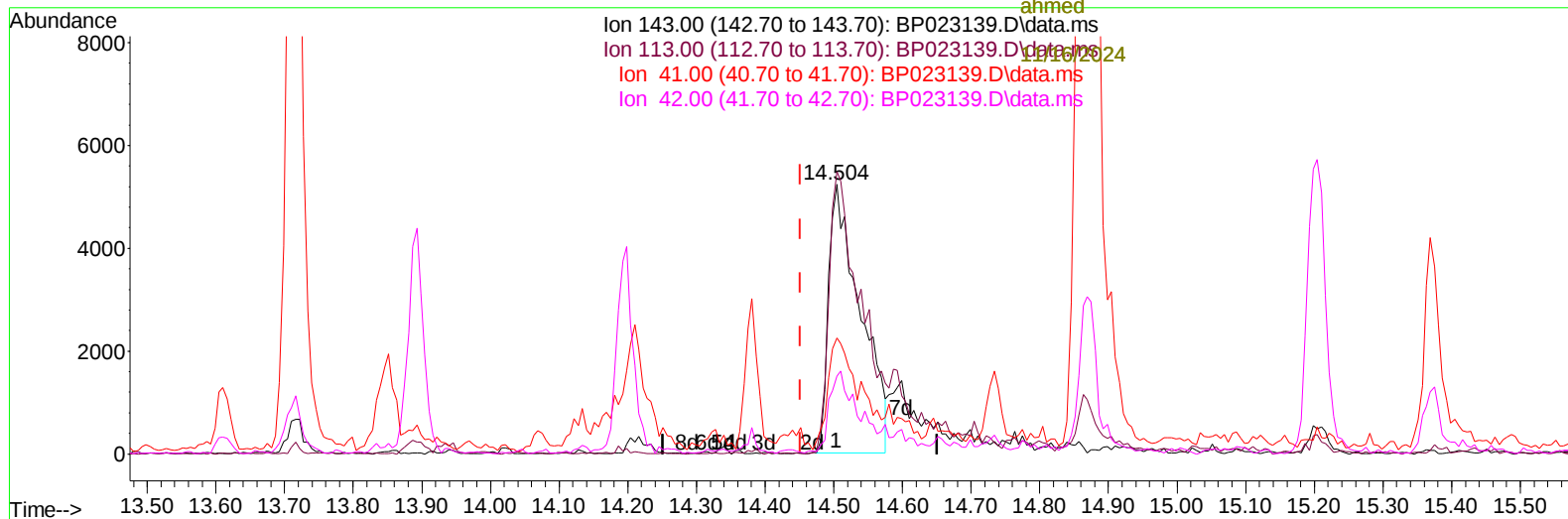
Instrument :
BNA_P
ClientSampleId :
E27N9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2024
Supervised By :mohammad ahmed 11/18/2024

11/16/2024
Supervised By :mohammad
ahmed

Quant Time: Nov 16 00:50:35 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



TIC: BP023139.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.053) 8.04 ng/u1

response 16735

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	104.92
41.00	39.20	43.05
42.00	25.90	28.89

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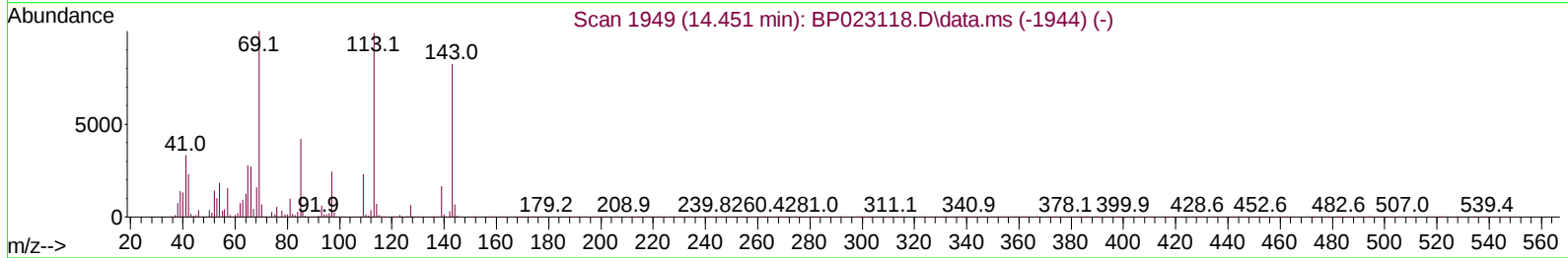
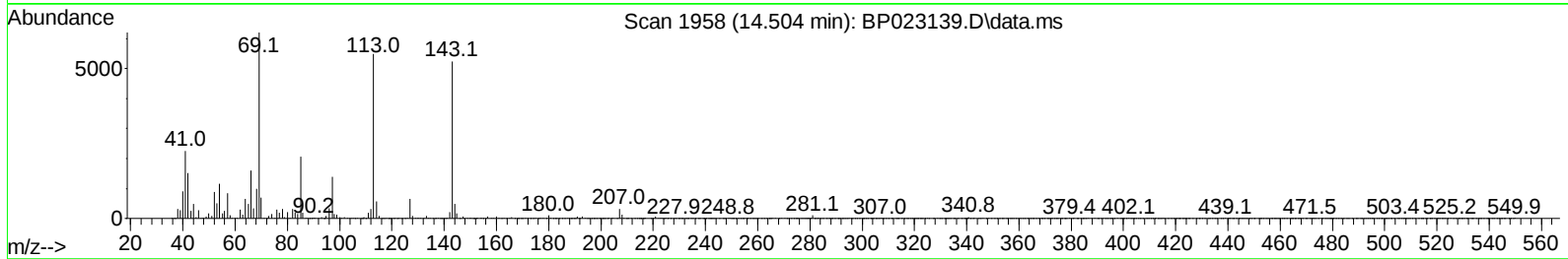
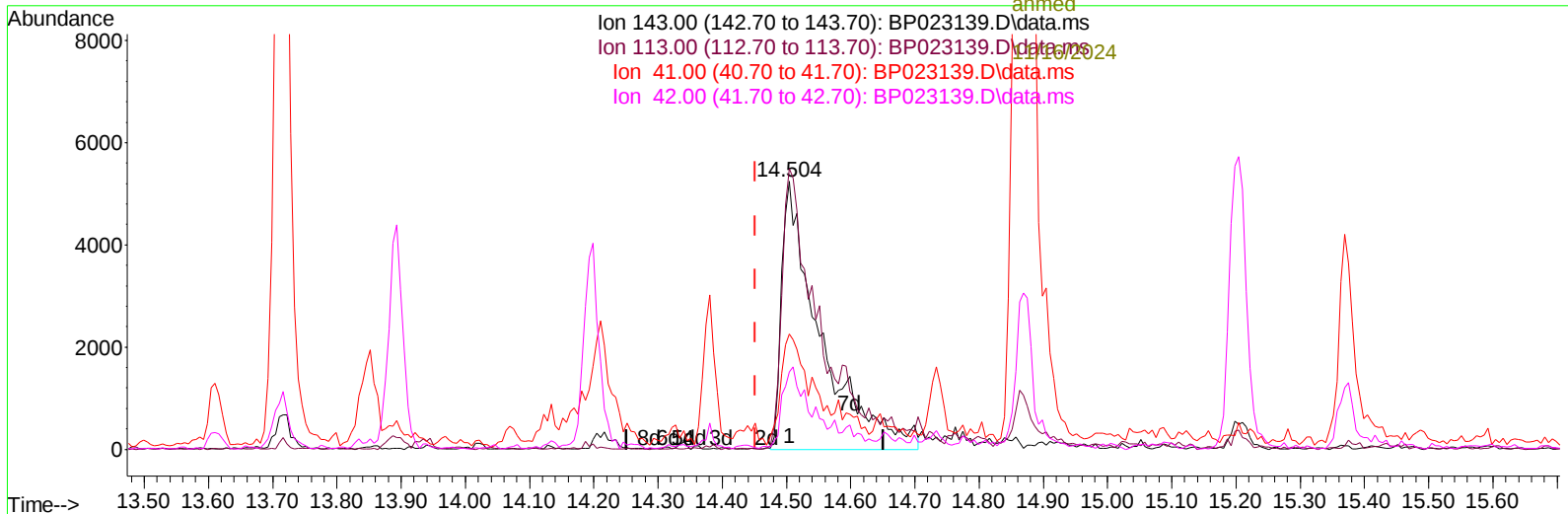
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(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.053) 10.51 ng/ul m

response 21879

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	104.92
41.00	39.20	43.05
42.00	25.90	28.89

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 Response via : Initial Calibration

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/16/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	61591	20.000	ng/ul	0.00
20) Naphthalene-d8	10.328	136	236426	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.198	164	192172	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	488030	20.000	ng/ul	-0.02
79) Chrysene-d12	21.445	240	570870	20.000	ng/ul	0.00
88) Perylene-d12	24.680	264	709075	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	6931	5.409	ng/uL	0.00
4) Pyridine-d5	3.558	84	33000	8.668	ng/ul	0.01
7) Phenol-d5	6.787	99	44398	9.750	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	80268	30.003	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	95955	28.773	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	82470	22.008	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	51431	31.247	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	63238	32.227	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	127888	31.122	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131	116080	23.678	ng/ul	0.00
46) Dimethylphthalate-d6	13.610	166	474692	34.273	ng/ul	0.00
49) Acenaphthylene-d8	13.893	160	467703	32.363	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	21879m	10.511	ng/ul	0.05
60) Fluorene-d10	15.204	176	420264	34.609	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	94944	32.476	ng/ul	0.00
73) Anthracene-d10	17.116	188	829592	40.147	ng/ul	0.00
81) Pyrene-d10	19.492	212	1040713	39.579	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.445	264	1347547	39.897	ng/ul	0.00

Target Compounds Qvalue

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

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