

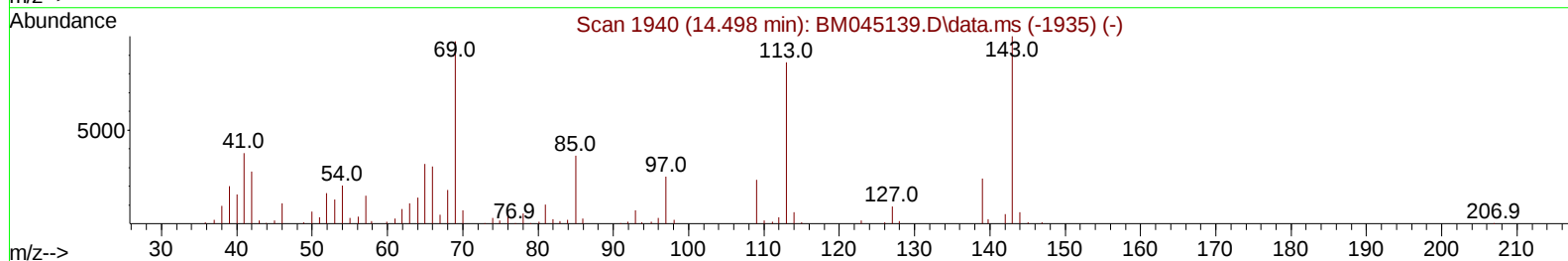
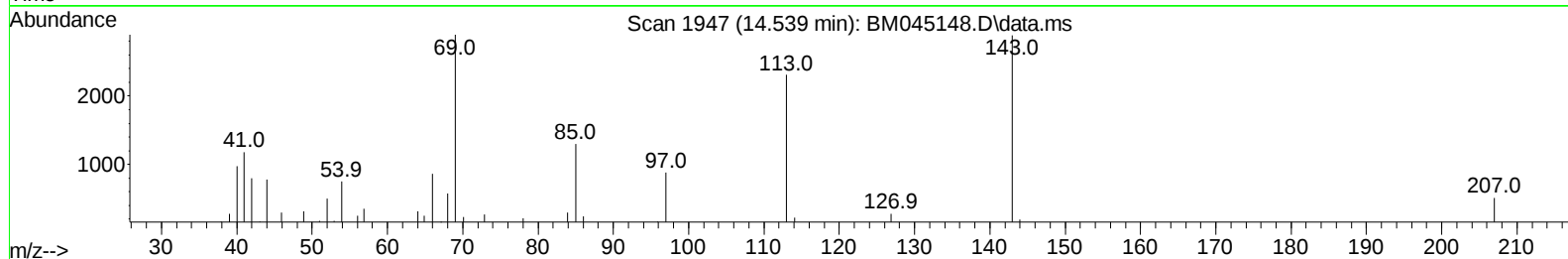
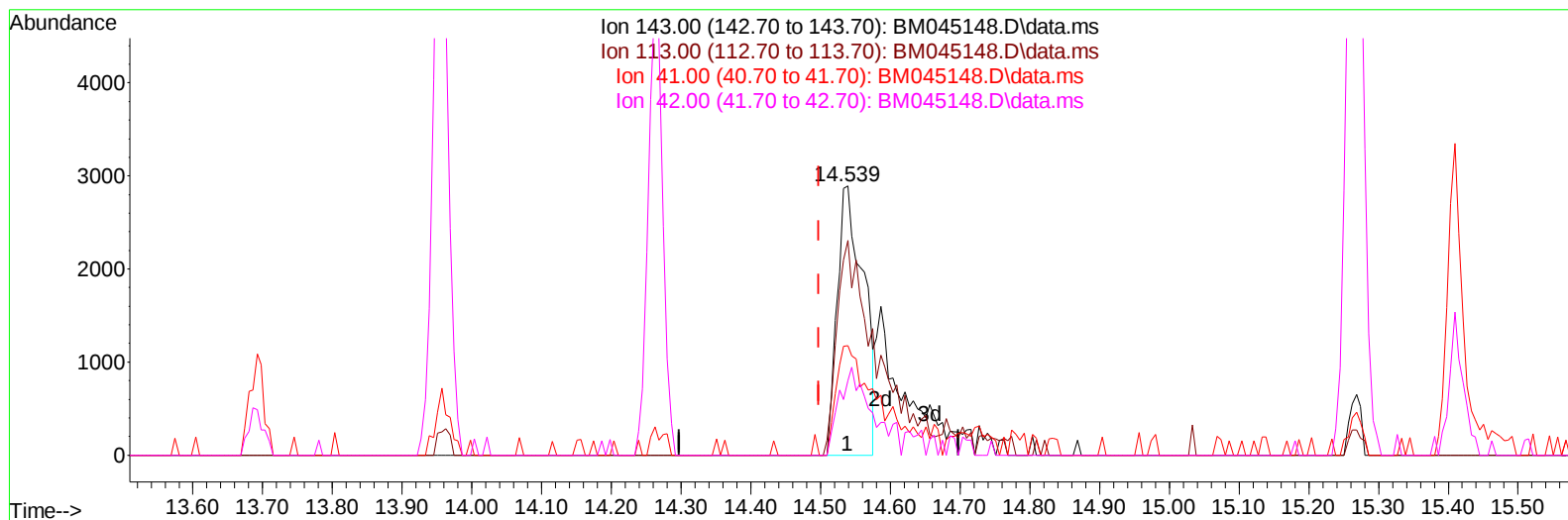
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041124\
Data File : BM045148.D
Acq On : 11 Apr 2024 23:35
Operator : MA/JU
Sample : P2013-13
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EOMB9

Manual IntegrationsAPPROVED

Quant Time: Apr 12 00:08:05 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 11 18:54:49 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/16/2024
Supervised By :mohammad ahmed 04/16/2024



TIC: BM045148.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.539min (+ 0.041) 3.42 ng/u1

response 7468

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	79.93
41.00	36.50	40.80
42.00	23.30	27.61

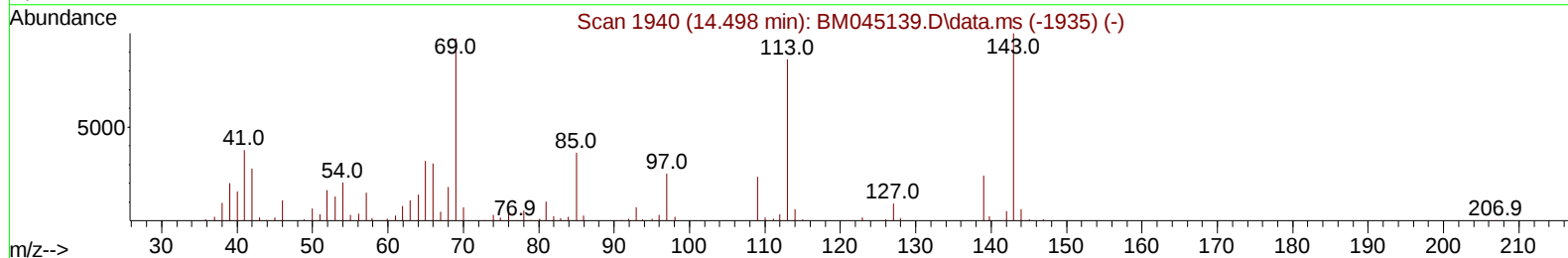
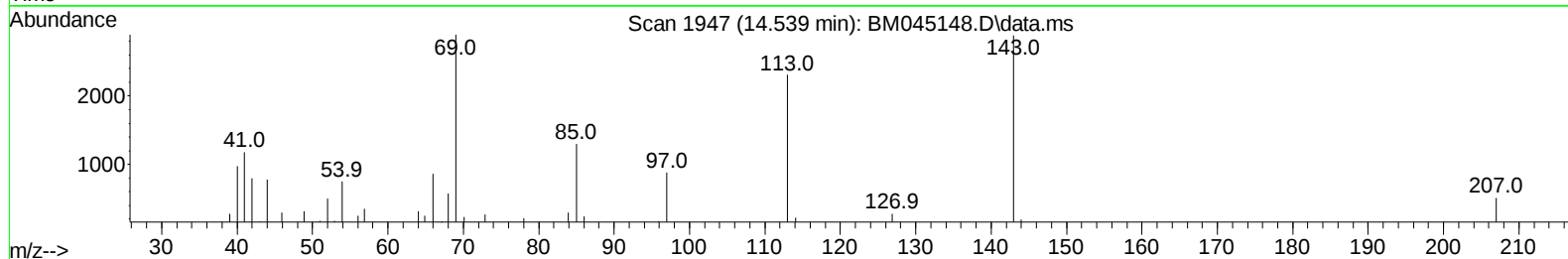
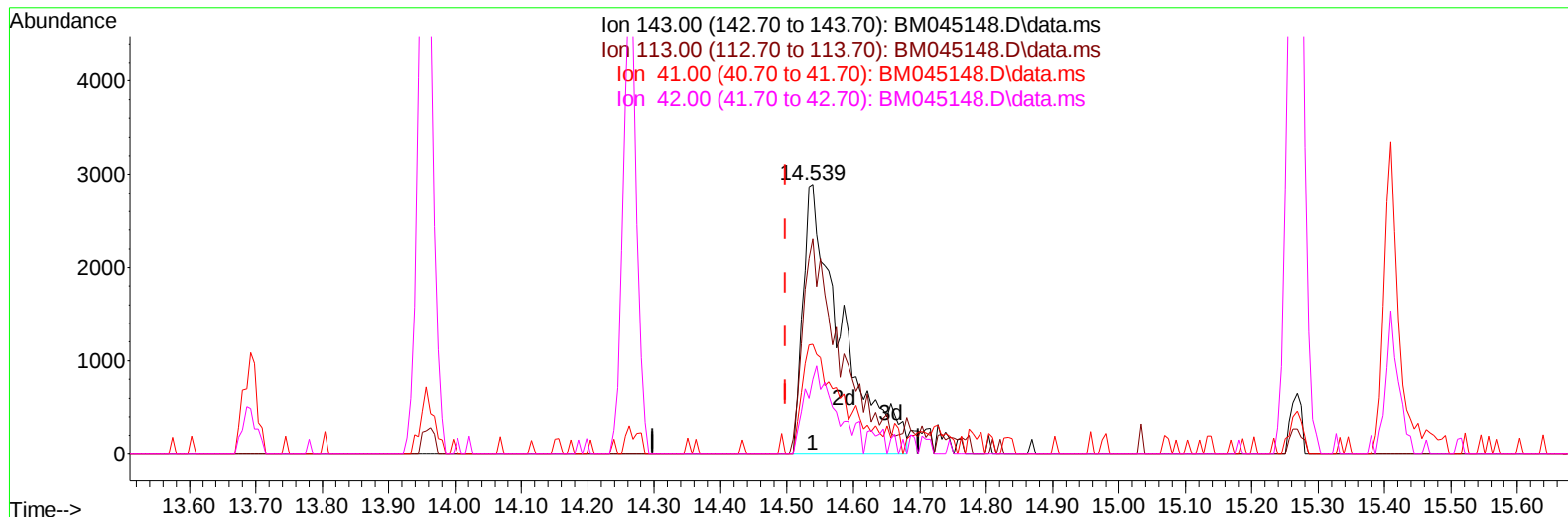
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TIC: BM045148.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.539min (+ 0.041) 5.36 ng/ul m

response 11721

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	79.93
41.00	36.50	40.80
42.00	23.30	27.61

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 Sample : P2013-13
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Instrument :
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 ClientSampleId :
 E0MB9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/16/2024
 Supervised By :mohammad ahmed 04/16/2024

Quant Time: Apr 12 00:08:57 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 11 18:54:49 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	59533	20.000	ng/ul	0.00
20) Naphthalene-d8	10.386	136	251125	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.262	164	150369	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.027	188	294600	20.000	ng/ul	0.00
79) Chrysene-d12	21.233	240	278379	20.000	ng/ul	0.00
88) Perylene-d12	23.450	264	366390	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	7807	4.897	ng/uL	0.00
4) Pyridine-d5	3.610	84	48068	11.087	ng/ul	0.00
7) Phenol-d5	6.804	99	40279	7.981	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	121405	40.376	ng/ul	0.00
11) 2-Chlorophenol-d4	7.151	132	123637	31.269	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	78831	19.955	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	80879	41.929	ng/ul	0.00
24) 2-Nitrophenol-d4	9.492	143	82428	40.485	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	129191	34.609	ng/ul	0.00
31) 4-Chloroaniline-d4	10.551	131	182434	30.696	ng/ul	0.00
46) Dimethylphthalate-d6	13.692	166	452567	43.226	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	540431	43.693	ng/ul	0.00
54) 4-Nitrophenol-d4	14.539	143	11721m	5.364	ng/ul	0.04
60) Fluorene-d10	15.268	176	397149	42.549	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.409	200	55463	31.752	ng/ul	0.00
73) Anthracene-d10	17.127	188	611520	46.143	ng/ul	0.00
81) Pyrene-d10	19.427	212	700201	51.082	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.309	264	819393	45.844	ng/ul	0.00

Target Compounds Qvalue

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

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