

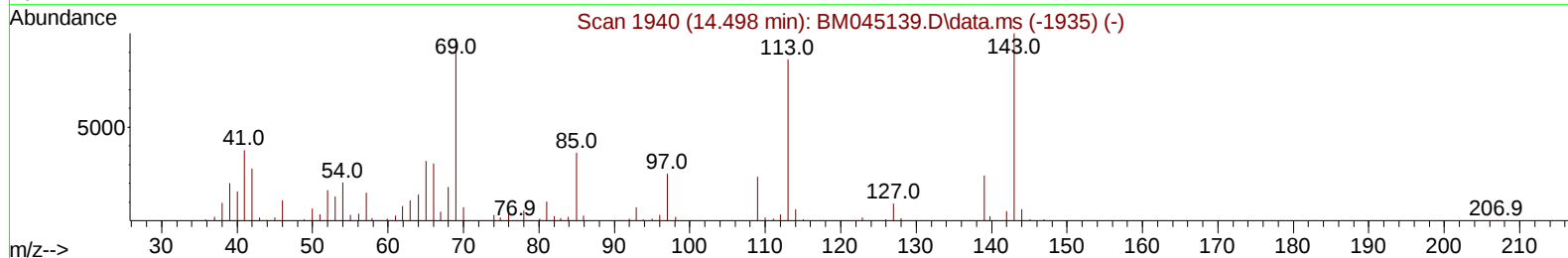
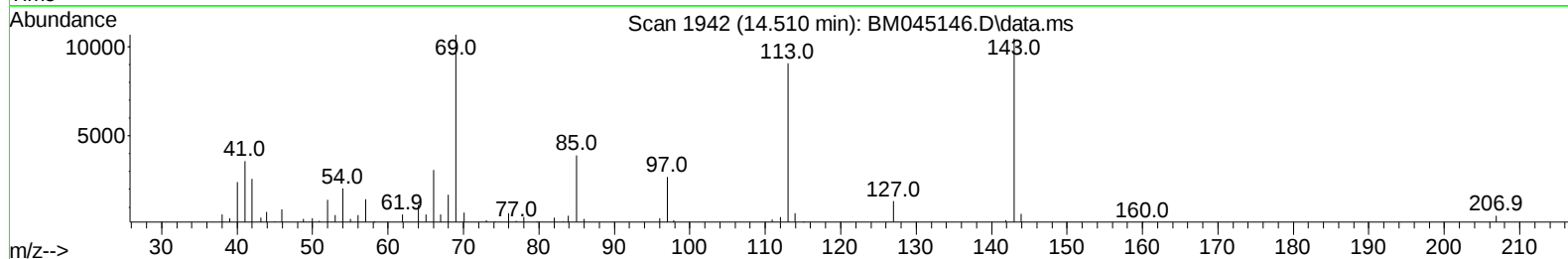
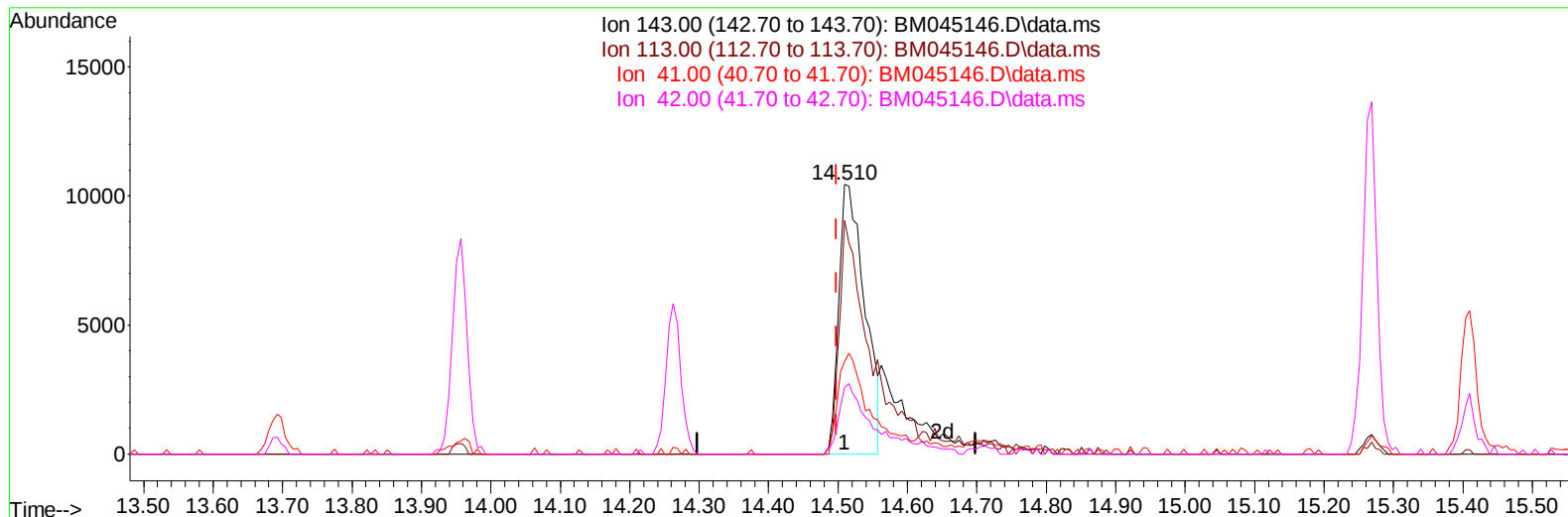
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041124\
Data File : BM045146.D
Acq On : 11 Apr 2024 22:22
Operator : MA/JU
Sample : P1982-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
YCSQ6

Manual IntegrationsAPPROVED

Quant Time: Apr 11 23:01:17 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: BM045146.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.012) 9.92 ng/u1

response 26819

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	86.64
41.00	36.50	34.42
42.00	23.30	24.67

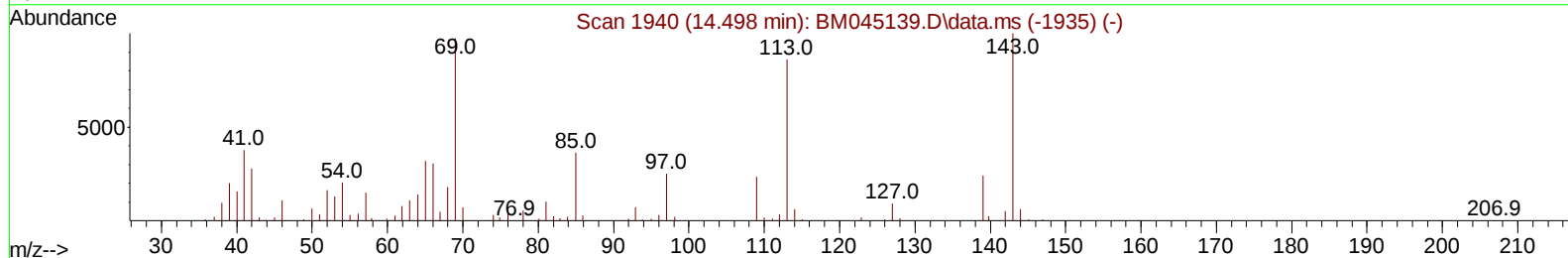
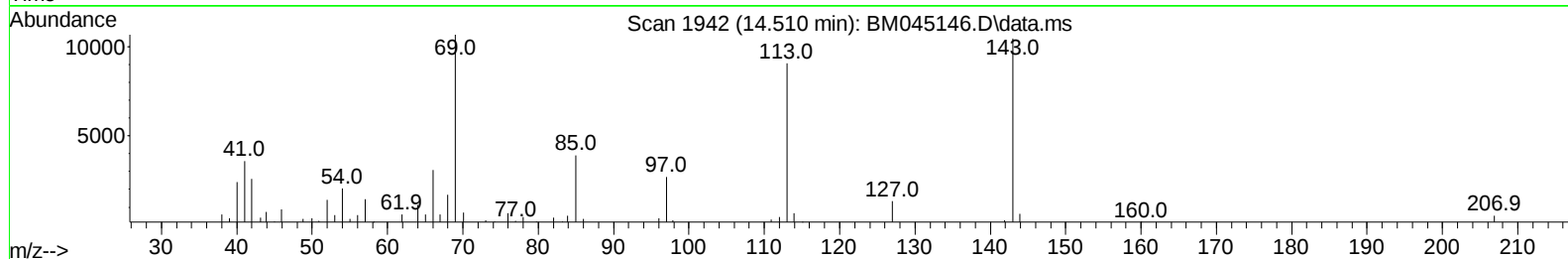
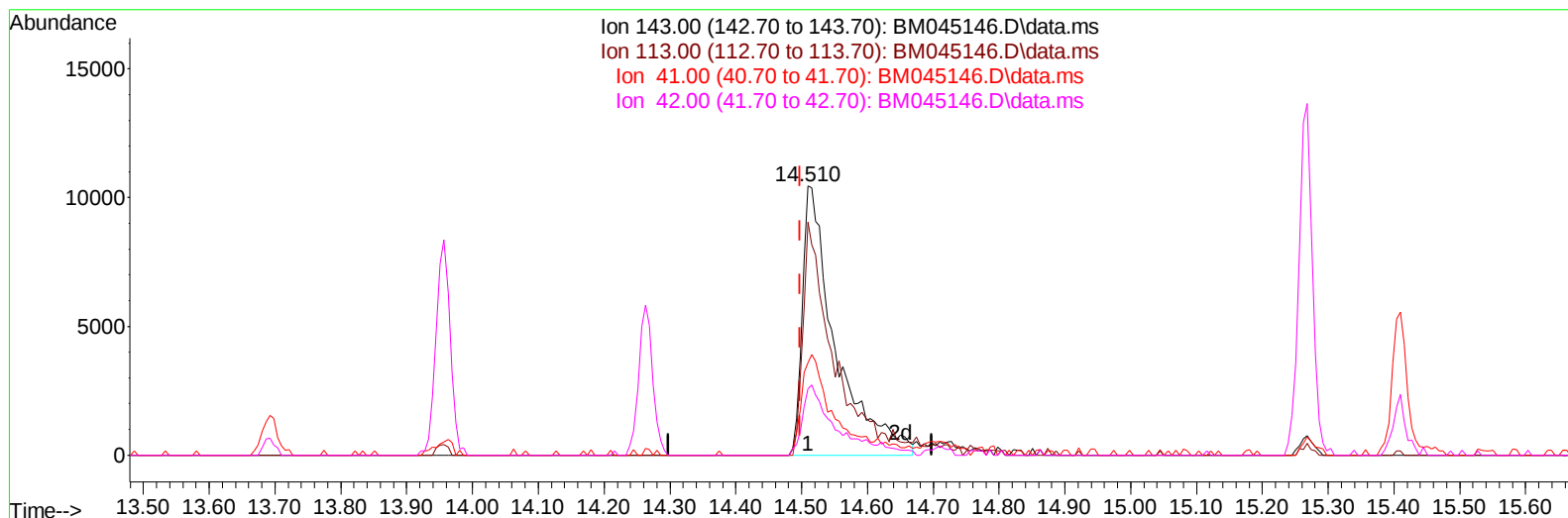
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TIC: BM045146.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.012) 13.49 ng/ul m

response 36469

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	86.64
41.00	36.50	34.42
42.00	23.30	24.67

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/16/2024
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Quant Time: Apr 11 23:02:08 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	61887	20.000	ng/ul	0.00
20) Naphthalene-d8	10.392	136	272532	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.263	164	186072	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.021	188	392526	20.000	ng/ul	0.00
79) Chrysene-d12	21.233	240	336872	20.000	ng/ul	0.00
88) Perylene-d12	23.450	264	361164	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	7468	4.506	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.804	99	73394	13.989	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	131280	41.999	ng/ul	0.00
11) 2-Chlorophenol-d4	7.151	132	163416	39.757	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	125242	30.497	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	91362	43.643	ng/ul	0.00
24) 2-Nitrophenol-d4	9.492	143	102265	46.283	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	172000	42.457	ng/ul	0.00
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	13.692	166	657042	50.715	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	653280	42.683	ng/ul	0.00
54) 4-Nitrophenol-d4	14.510	143	36469m	13.488	ng/ul	0.01
60) Fluorene-d10	15.268	176	550075	47.624	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	100372	43.127	ng/ul	0.00
73) Anthracene-d10	17.127	188	860943	48.757	ng/ul	0.00
81) Pyrene-d10	19.433	212	1023413	61.698	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.309	264	922549	52.362	ng/ul	0.00

Target Compounds Qvalue

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

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