

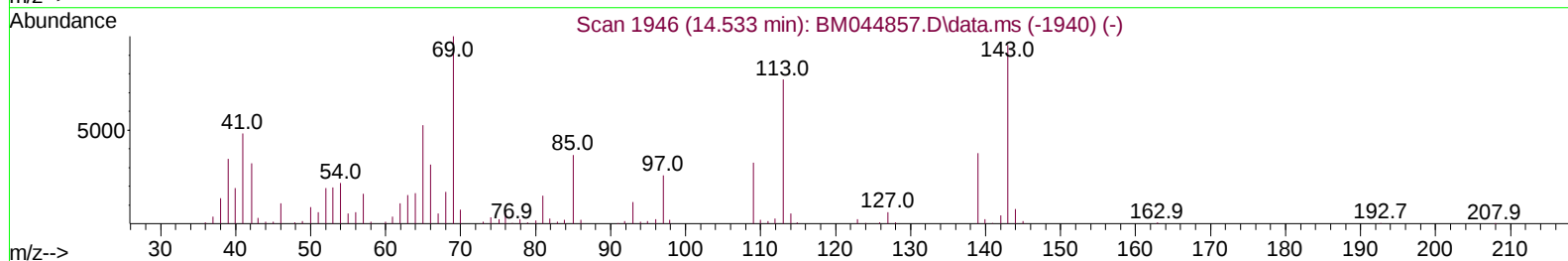
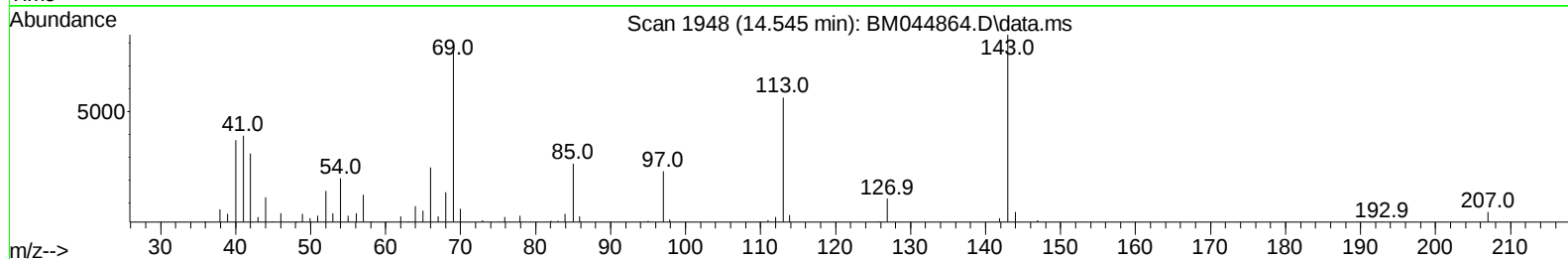
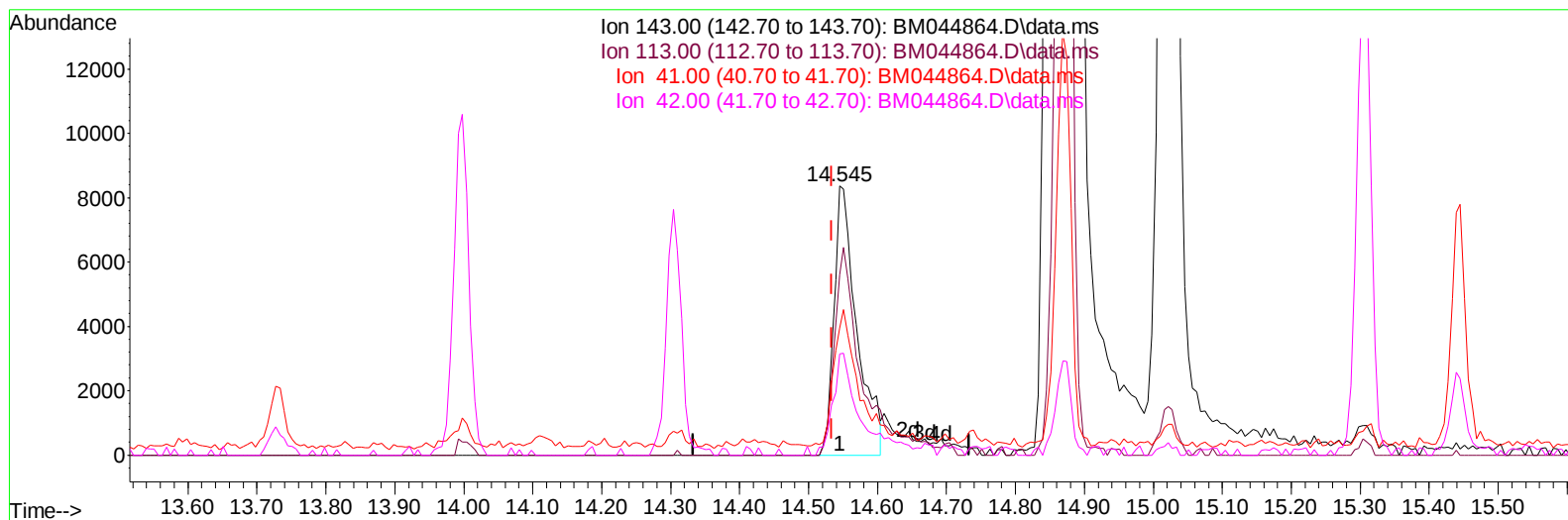
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040124\
Data File : BM044864.D
Acq On : 01 Apr 2024 14:23
Operator : MA/JU
Sample : P1925-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
PMW-2(20240327)

Manual IntegrationsAPPROVED

Quant Time: Apr 01 14:52:08 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Mar 31 13:54:26 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/02/2024
Supervised By :mohammad ahmed 04/02/2024



TIC: BM044864.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.545min (+ 0.012) 5.92 ng/u1

response 19237

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	67.17
41.00	38.40	47.25#
42.00	26.60	37.61#

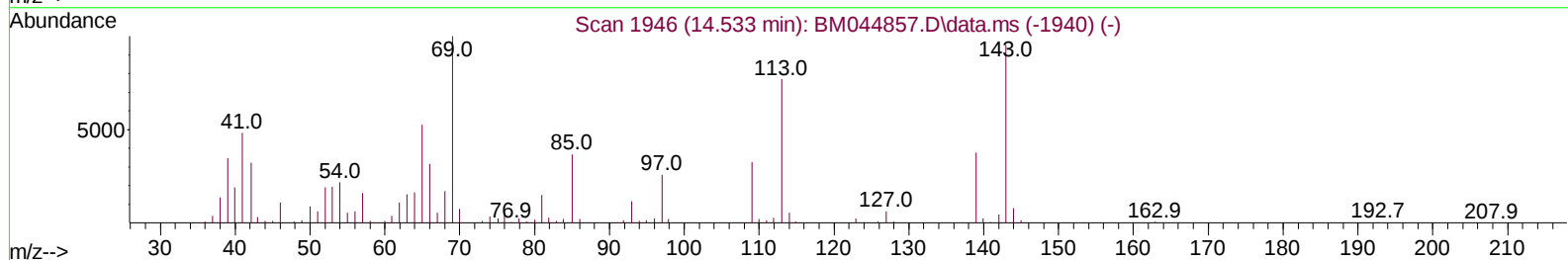
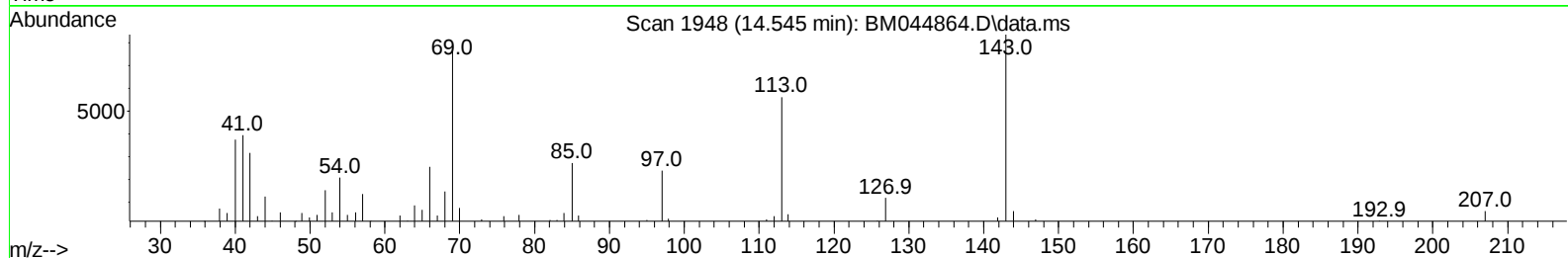
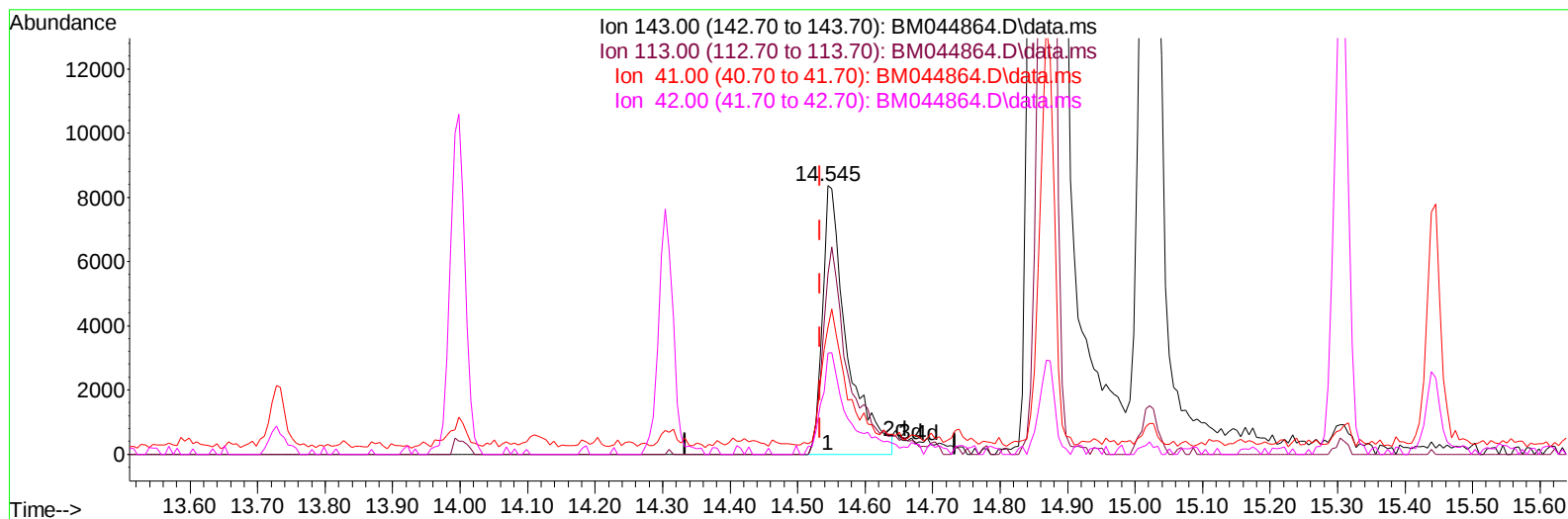
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TIC: BM044864.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.545min (+ 0.012) 6.46 ng/ul m

response 20980

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	67.17
41.00	38.40	47.25#
42.00	26.60	37.61#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/02/2024
 Supervised By :mohammad ahmed 04/02/2024

Quant Time: Apr 01 23:28:34 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Mar 31 13:54:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.657	152	87945	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.434	136	356991	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.304	164	199726	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.063	188	368765	20.000	ng/ul	0.00
79) Chrysene-d12	21.262	240	299120	20.000	ng/ul	0.00
88) Perylene-d12	23.492	264	355019	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	14686	5.664	ng/uL	0.00
4) Pyridine-d5	3.634	84	69762	9.141	ng/ul	0.00
7) Phenol-d5	6.840	99	73252	8.102	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.010	67	194190	34.770	ng/ul	0.00
11) 2-Chlorophenol-d4	7.193	132	184647	27.790	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	128909	18.154	ng/ul	0.00
21) Nitrobenzene-d5	8.822	128	111953	37.242	ng/ul	0.00
24) 2-Nitrophenol-d4	9.534	143	120376	37.146	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	199136	36.106	ng/ul	0.00
31) 4-Chloroaniline-d4	10.587	131	270258	29.504	ng/ul	0.00
46) Dimethylphthalate-d6	13.727	166	641941	40.973	ng/ul	0.00
49) Acenaphthylene-d8	13.998	160	714967	39.902	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	20980m	6.457	ng/ul	0.01
60) Fluorene-d10	15.304	176	533904	40.552	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.445	200	87408	36.515	ng/ul	0.00
73) Anthracene-d10	17.163	188	799803	45.374	ng/ul	0.00
81) Pyrene-d10	19.462	212	861866	46.978	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.356	264	855225	45.831	ng/ul	0.00

Target Compounds Qvalue

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

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