

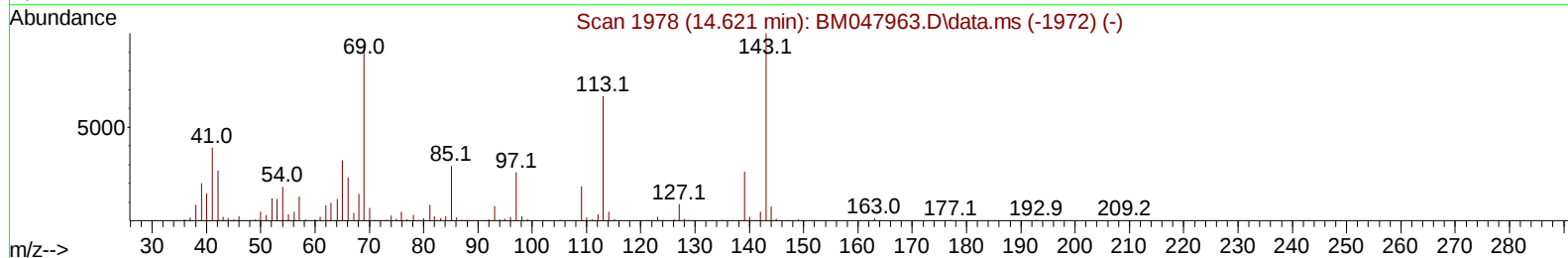
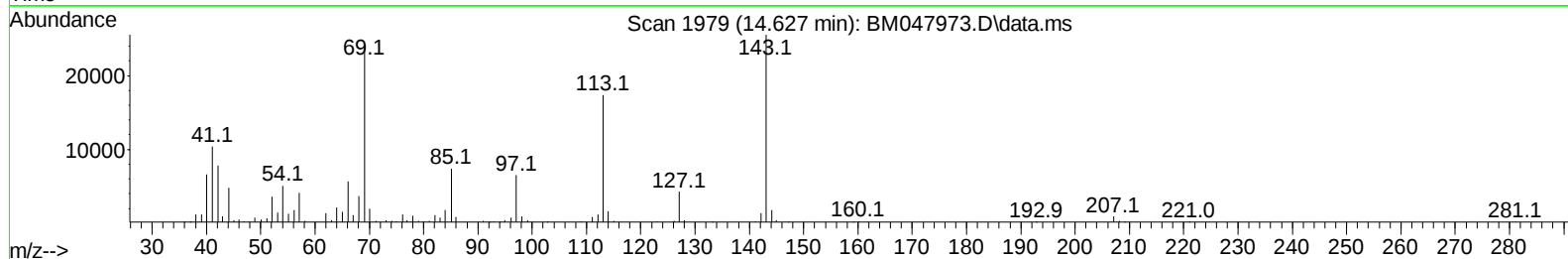
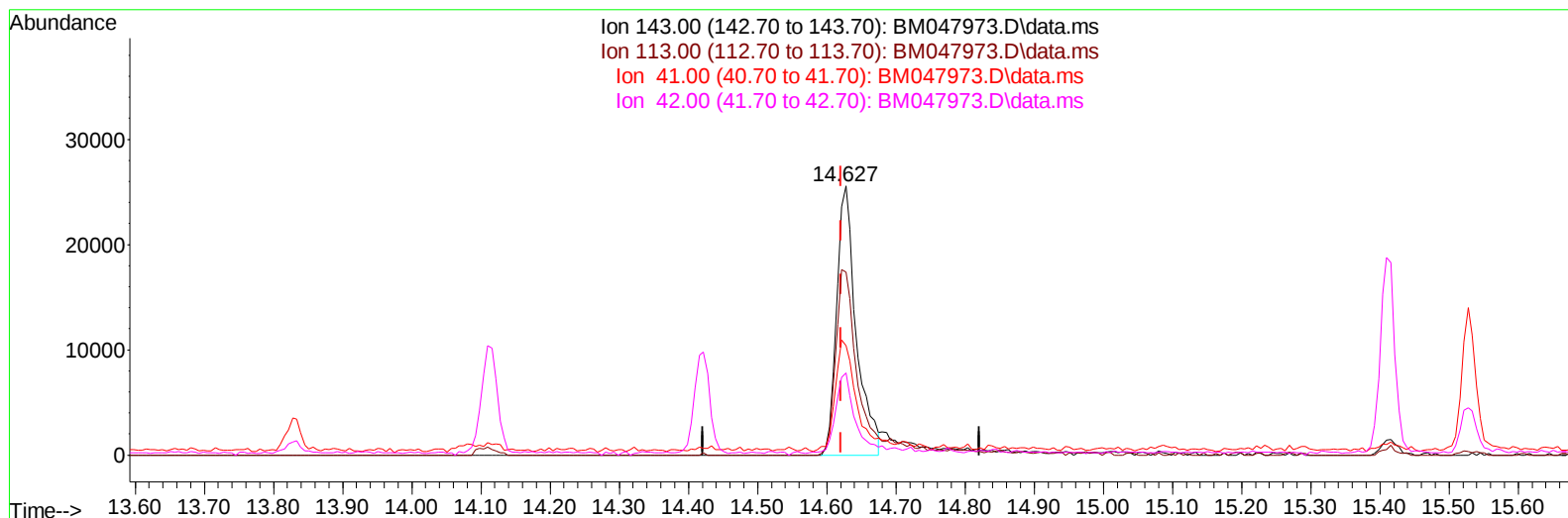
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100924\
Data File : BM047973.D
Acq On : 10 Oct 2024 10:30
Operator : RC/JU
Sample : P4215-02
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EY073

Manual IntegrationsAPPROVED

Quant Time: Oct 10 11:14:09 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 10 05:10:14 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/11/2024
Supervised By :mohammad ahmed 10/11/2024



TIC: BM047973.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.627min (+ 0.006) 6.14 ng/u1

response 50341

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.20	68.03
41.00	60.00	40.71#
42.00	40.50	30.57#

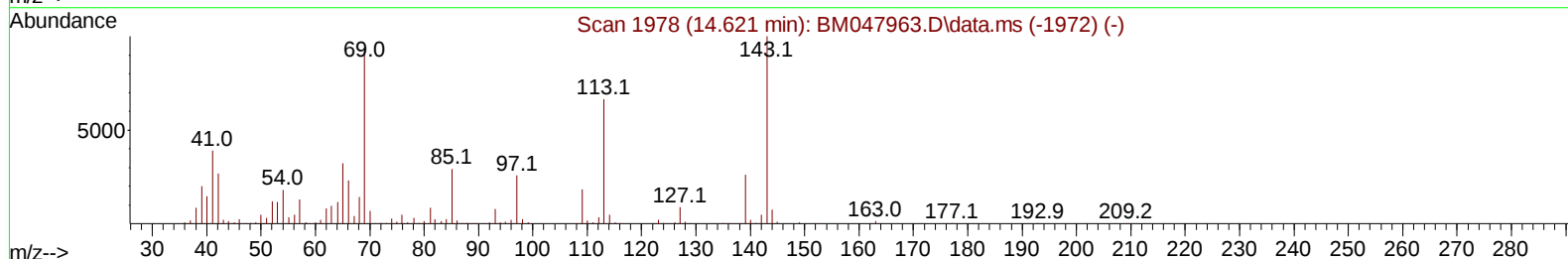
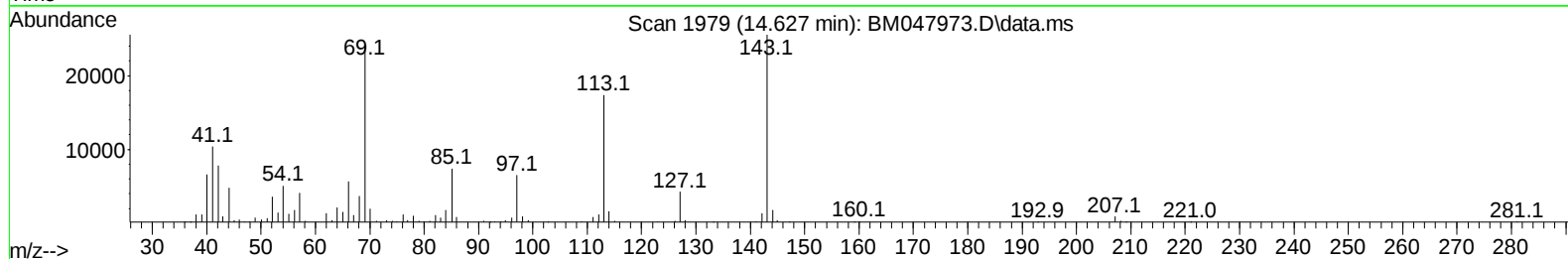
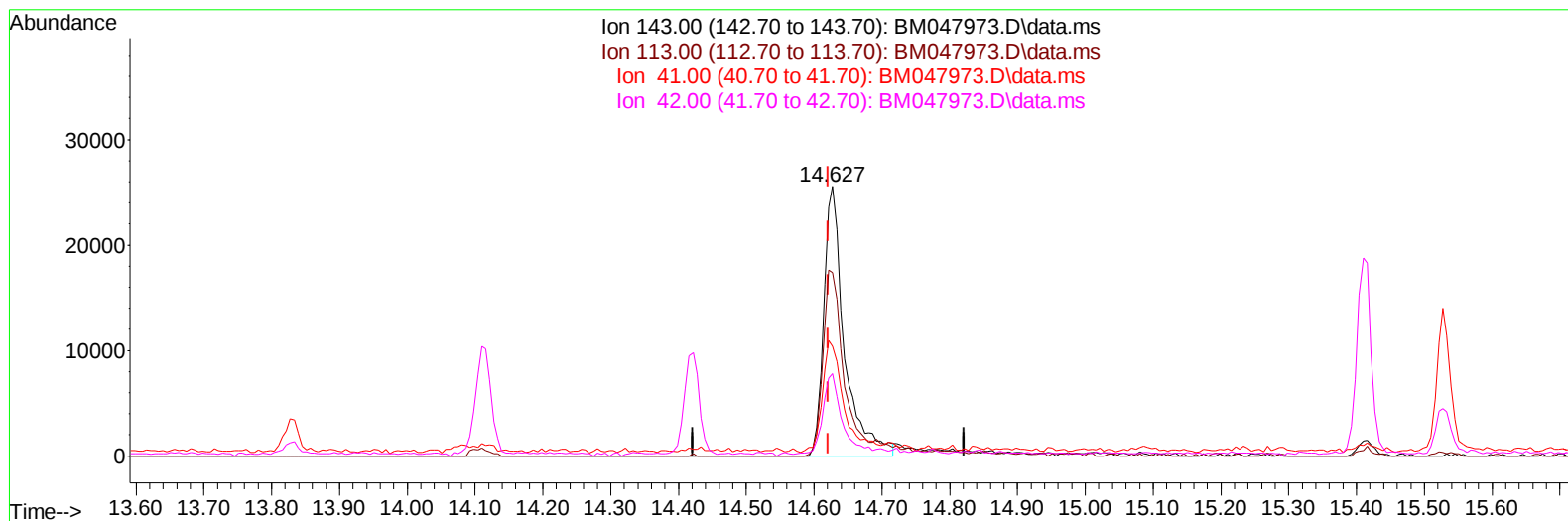
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TIC: BM047973.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.627min (+ 0.006) 6.60 ng/ul m

response 54147

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.20	68.03
41.00	60.00	40.71#
42.00	40.50	30.57#

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/11/2024
 Supervised By :mohammad ahmed 10/11/2024

Quant Time: Oct 10 23:32:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 10 05:10:14 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	240403	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136	948130	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.422	164	568422	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.157	188	1173433	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	1190210	20.000	ng/ul	0.00
88) Perylene-d12	24.374	264	1401297	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	29126	3.915	ng/uL	0.00
4) Pyridine-d5	3.658	84	157190	7.217	ng/ul	0.00
7) Phenol-d5	6.957	99	176731	7.774	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	354056	21.949	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	450940	27.265	ng/ul	0.00
15) 4-Methylphenol-d8	8.493	113	308440	18.249	ng/ul	0.00
21) Nitrobenzene-d5	8.940	128	226719	30.033	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	243923	33.718	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	429349	29.710	ng/ul	0.00
31) 4-Chloroaniline-d4	10.710	131	526250	23.554	ng/ul	0.00
46) Dimethylphthalate-d6	13.827	166	1337557	32.456	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	1474471	30.176	ng/ul	0.00
54) 4-Nitrophenol-d4	14.627	143	54147m	6.599	ng/ul	0.00
60) Fluorene-d10	15.410	176	1173530	32.876	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.527	200	204030	29.713	ng/ul	0.00
73) Anthracene-d10	17.257	188	2045299	35.448	ng/ul	0.00
81) Pyrene-d10	19.545	212	2457929	36.387	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.168	264	2708485	37.253	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.275	88	82414	10.143	ng/uL	91

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

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