

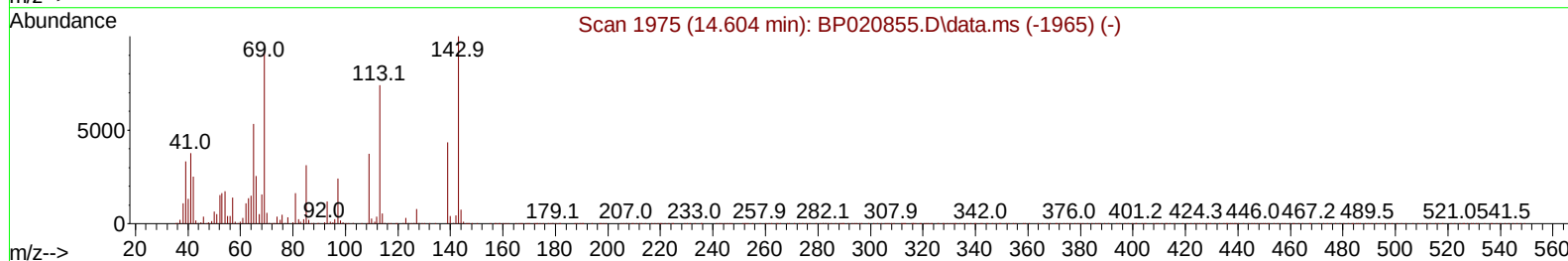
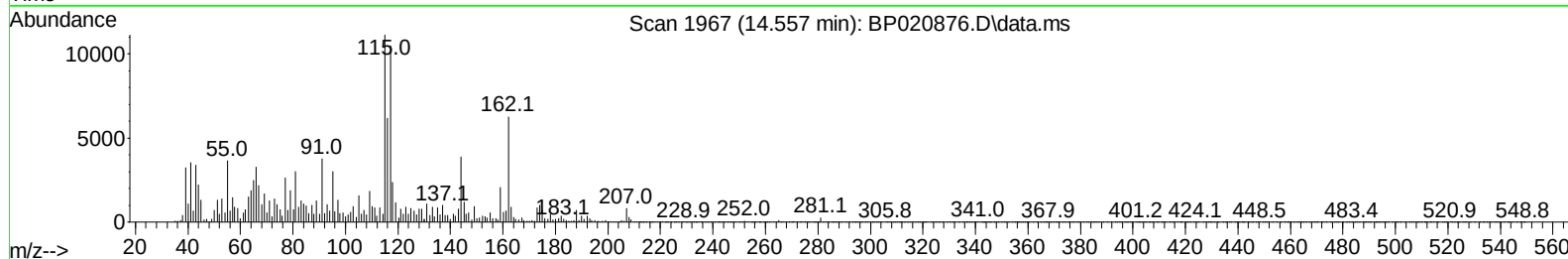
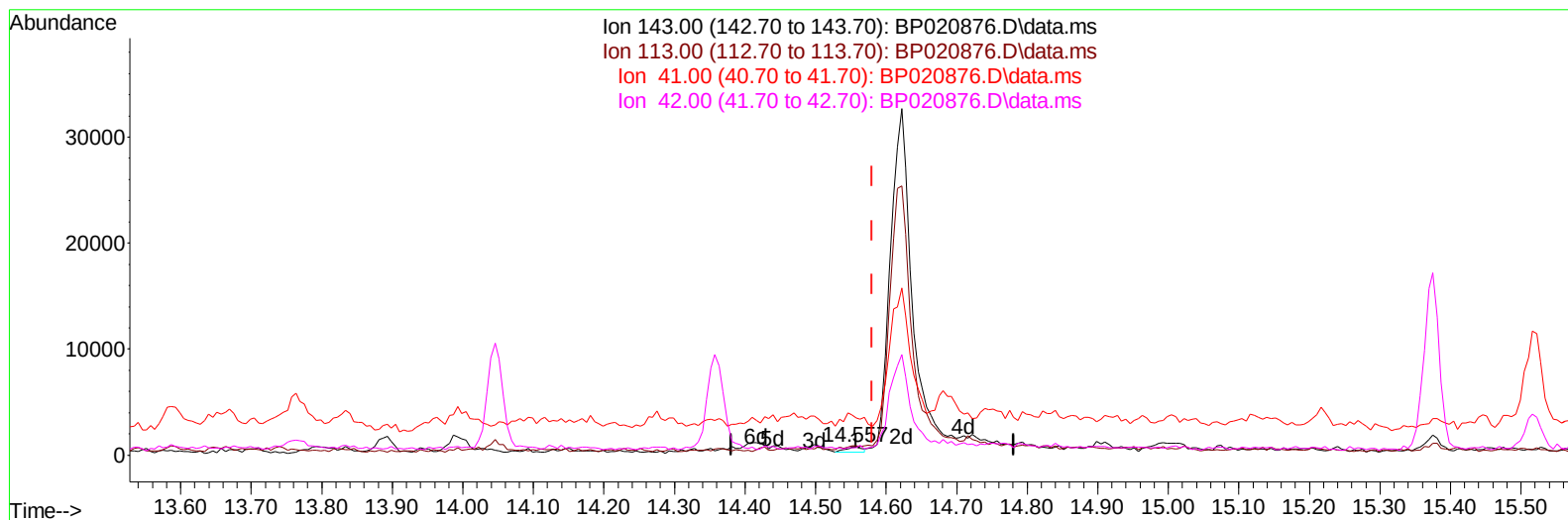
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020876.D
Acq On : 10 Jul 2024 00:14
Operator : MA/JU
Sample : P3109-06
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0HC4

Manual IntegrationsAPPROVED

Quant Time: Jul 10 01:43:39 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jul 03 15:01:56 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/10/2024
Supervised By :mohammad ahmed 07/10/2024



TIC: BP020876.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (-0.023) 0.15 ng/u1

response 848

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.40	109.35#
41.00	40.20	452.24#
42.00	28.00	88.73#

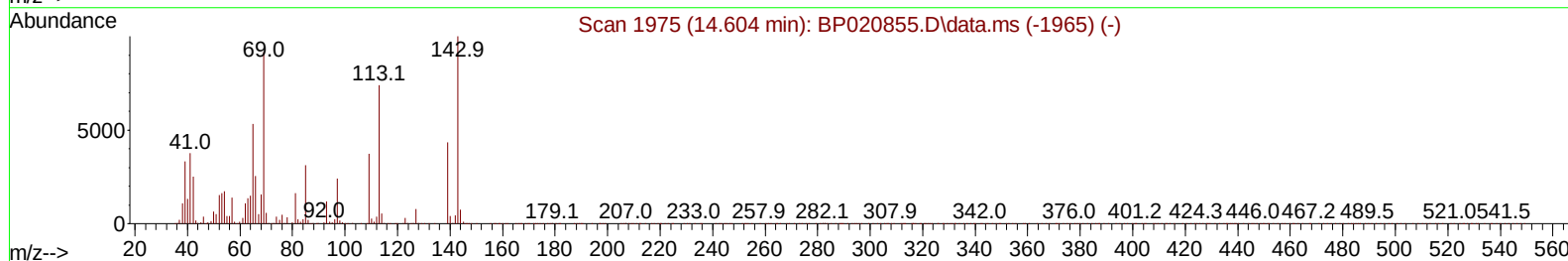
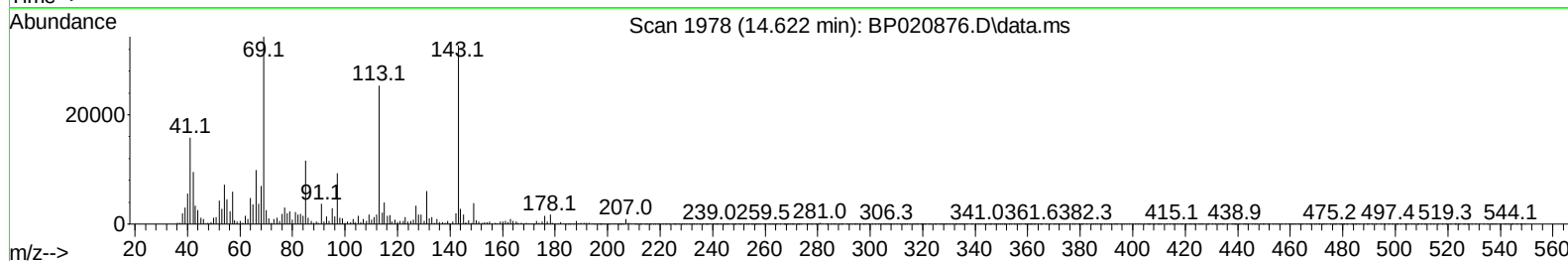
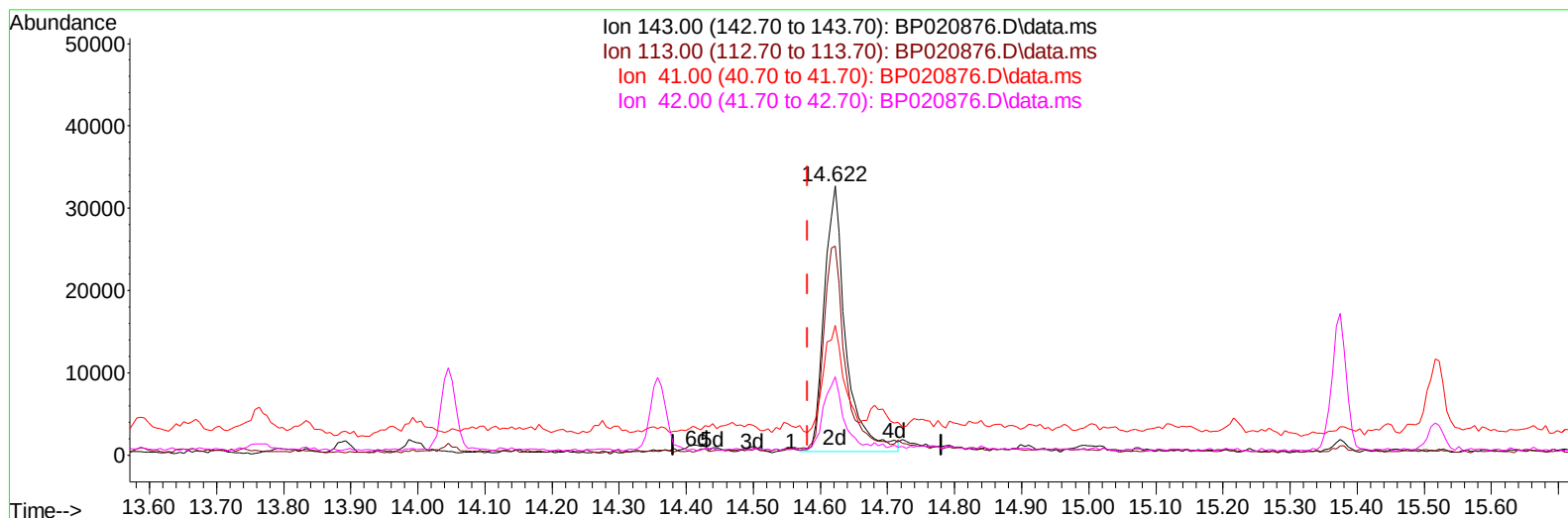
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(54) 4-Nitrophenol-d4 (S)

14.622min (+ 0.041) 12.52 ng/ul m

response 71825

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	79.40	77.65
41.00	40.20	48.32#
42.00	28.00	29.12

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/10/2024
 Supervised By :mohammad ahmed 07/10/2024

Quant Time: Jul 10 02:04:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 03 15:01:56 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	202144	20.000	ng/ul	0.00
20) Naphthalene-d8	10.493	136	817360	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.357	164	539959	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.175	188	1131607	20.000	ng/ul	0.01
79) Chrysene-d12	21.616	240	1092555	20.000	ng/ul	0.00
88) Perylene-d12	24.968	264	1229361	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	20826	4.424	ng/uL	0.00
4) Pyridine-d5	3.681	84	107405	7.829	ng/ul	0.00
7) Phenol-d5	6.922	99	195116	11.768	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	288342	27.433	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.269	132	379879	28.035	ng/ul	0.00
15) 4-Methylphenol-d8	8.440	113	325003	24.136	ng/ul	0.00
21) Nitrobenzene-d5	8.881	128	196701	32.596	ng/ul	0.00
24) 2-Nitrophenol-d4	9.593	143	230342	38.143	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	393249	31.613	ng/ul	0.00
31) 4-Chloroaniline-d4	10.640	131	408115	23.411	ng/ul	0.00
46) Dimethylphthalate-d6	13.763	166	1306633	34.804	ng/ul	0.00
49) Acenaphthylene-d8	14.046	160	1412910	31.540	ng/ul	0.00
54) 4-Nitrophenol-d4	14.622	143	71825m	12.521	ng/ul	0.04
60) Fluorene-d10	15.375	176	1088813	34.465	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.522	200	168333	29.518	ng/ul	0.02
73) Anthracene-d10	17.281	188	1824376	35.922	ng/ul	0.02
81) Pyrene-d10	19.657	212	2237201	34.085	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.727	264	2269449	37.950	ng/ul	-0.01

Target Compounds Qvalue

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

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