

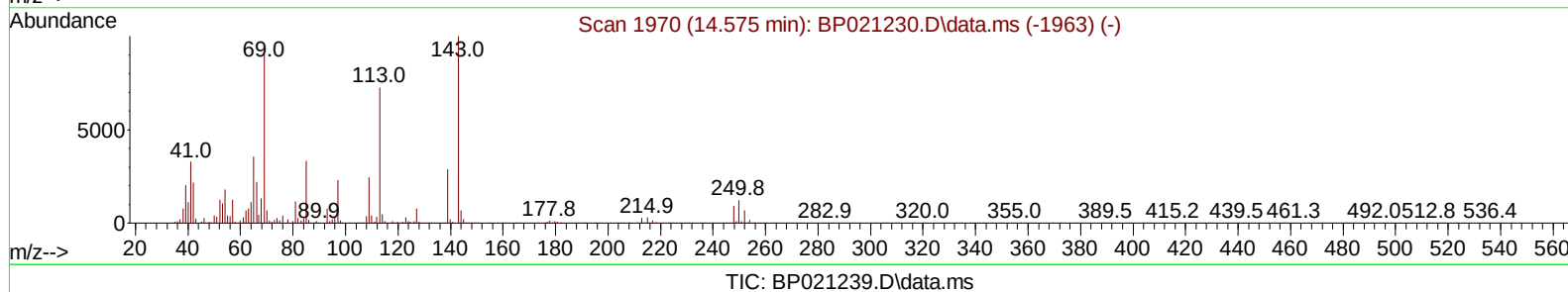
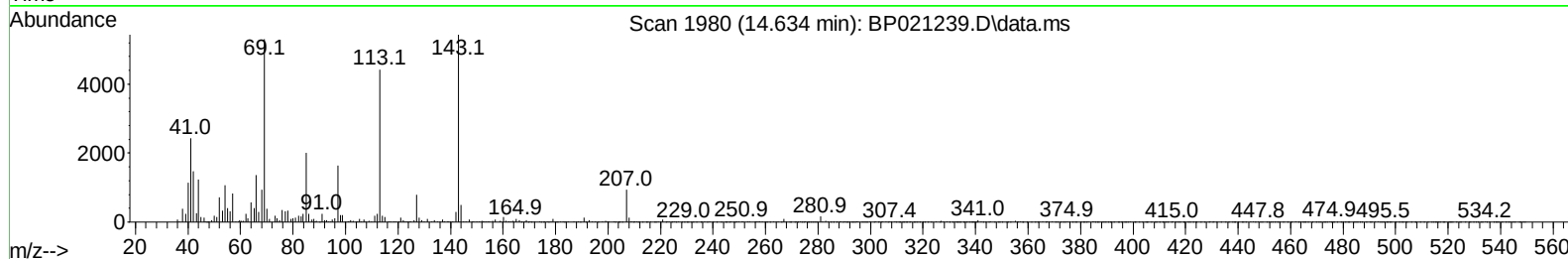
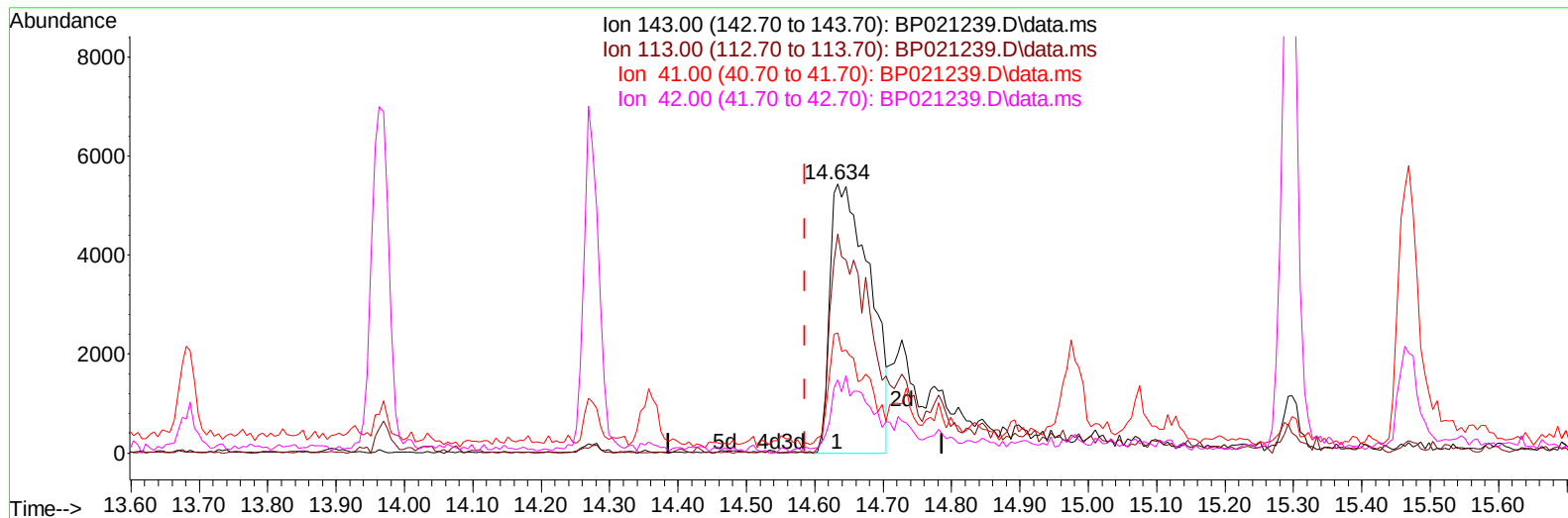
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021239.D
Acq On : 30 Jul 2024 17:47
Operator : MA/JU
Sample : P3359-13
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1GF1

Manual IntegrationsAPPROVED

Quant Time: Jul 30 22:54:44 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/31/2024
Supervised By :mohammad ahmed 09/04/2024



(54) 4-Nitrophenol-d4 (S)

14.634min (+ 0.047) 4.95 ng/ul

response 21878

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	81.47
41.00	37.90	44.73
42.00	27.10	27.14

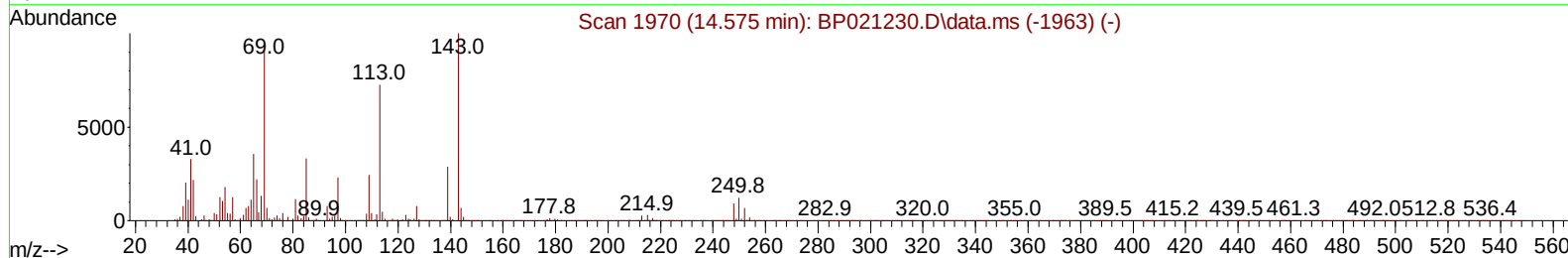
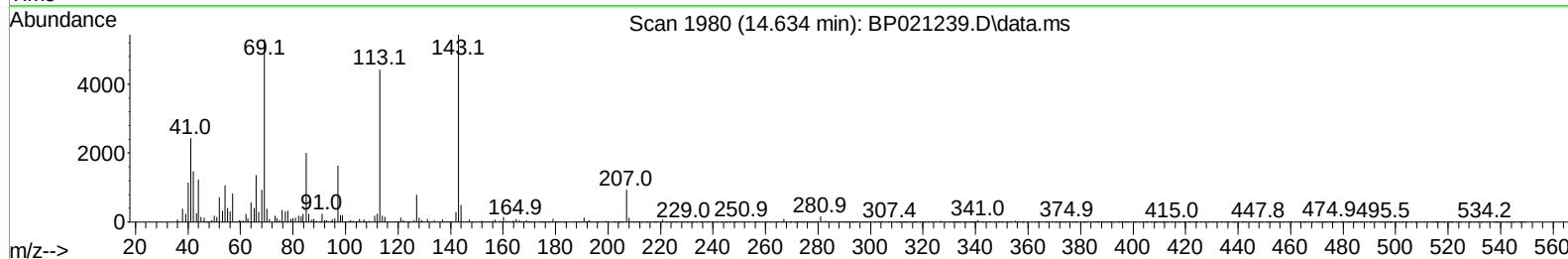
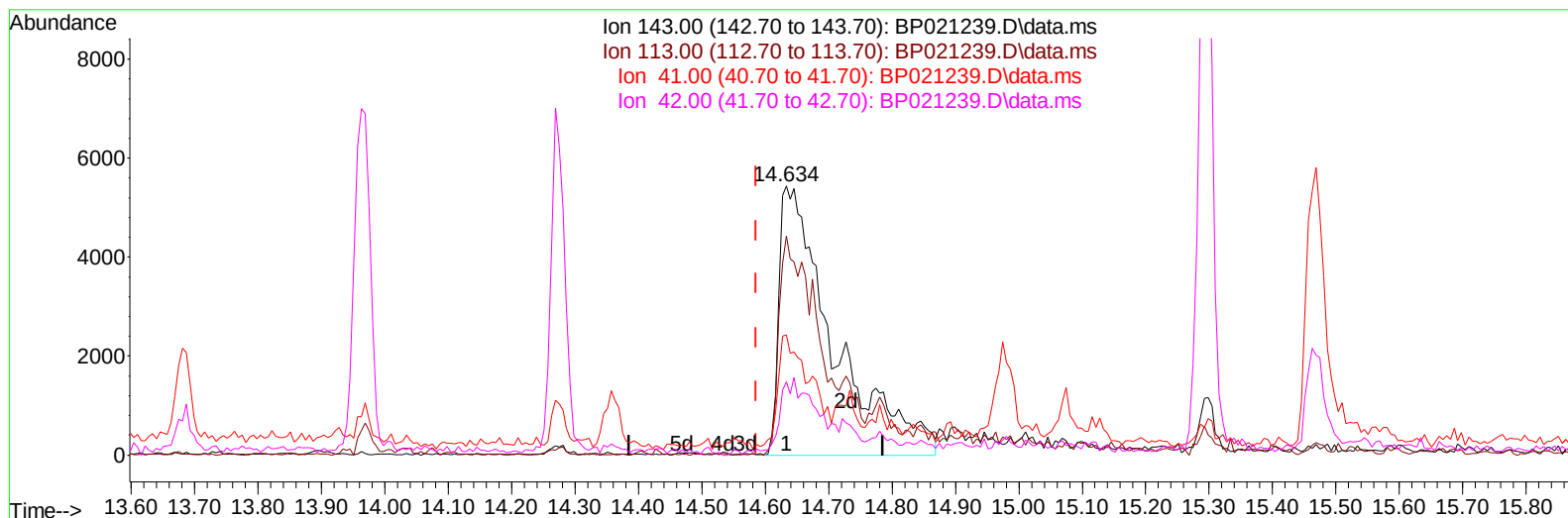
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TIC: BP021239.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.634min (+ 0.047) 7.32 ng/ul m

response 32372

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	81.47
41.00	37.90	44.73
42.00	27.10	27.14

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Quant Time: Jul 30 23:22:25 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 29 12:46:18 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	160084	20.000	ng/ul	0.00
20) Naphthalene-d8	10.398	136	630708	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.275	164	427561	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	957837	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	1014878	20.000	ng/ul	0.01
88) Perylene-d12	24.845	264	1030558	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	17598	5.002	ng/uL	0.00
4) Pyridine-d5	3.605	84	51410	5.048	ng/ul	0.01
7) Phenol-d5	6.863	99	95242	7.298	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	6.981	67	229643	29.034	ng/ul	0.00
11) 2-Chlorophenol-d4	7.187	132	302716	28.152	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	194914	17.982	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	163067	33.287	ng/ul	0.00
24) 2-Nitrophenol-d4	9.504	143	188937	33.697	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.057	165	318435	31.408	ng/ul	0.00
31) 4-Chloroaniline-d4	10.563	131	301556	22.452	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	1101755	35.447	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	1178064	33.755	ng/ul	0.00
54) 4-Nitrophenol-d4	14.634	143	32372m	7.320	ng/ul	0.05
60) Fluorene-d10	15.298	176	932241	36.560	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.469	200	162596	28.056	ng/ul	0.00
73) Anthracene-d10	17.204	188	1598747	37.580	ng/ul	0.01
81) Pyrene-d10	19.586	212	2043886	32.575	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.621	264	2033144	40.001	ng/ul	0.02

Target Compounds Qvalue

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

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