

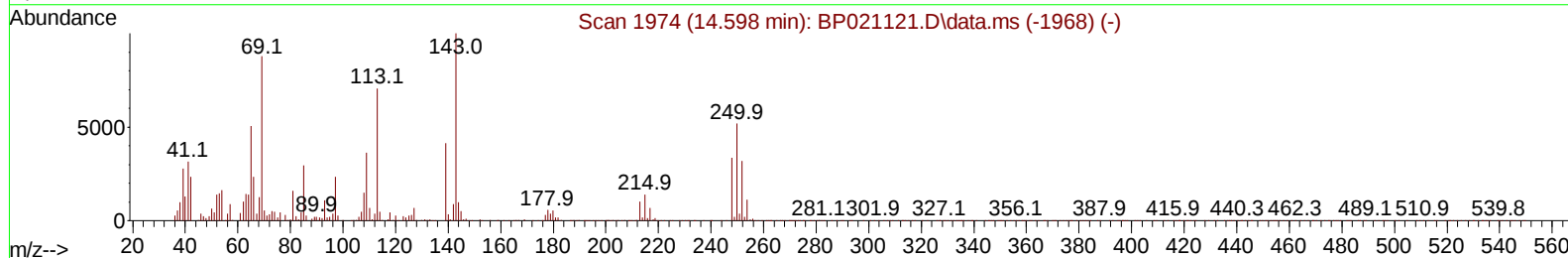
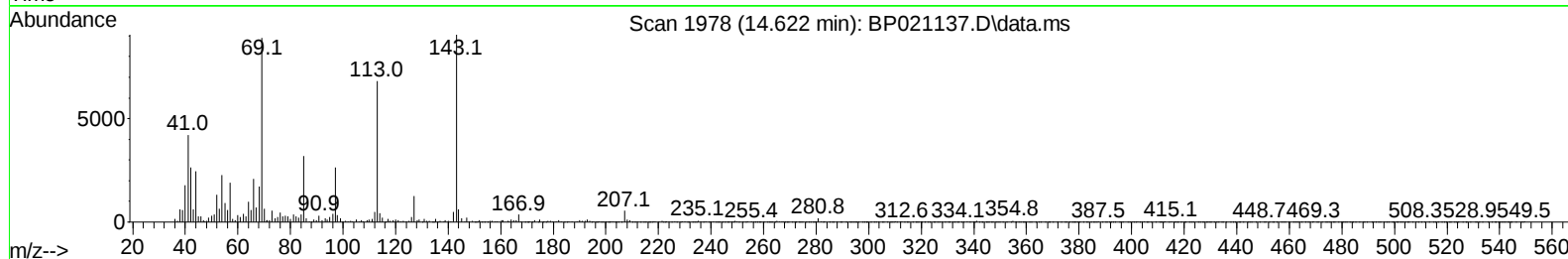
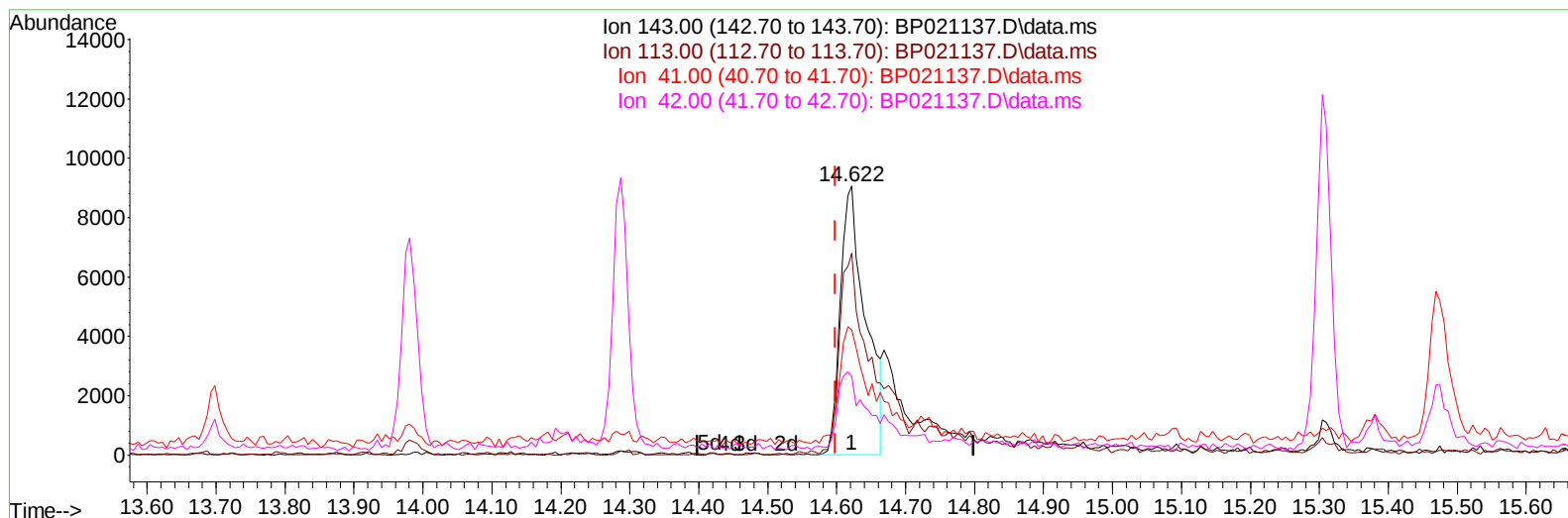
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072424\
Data File : BP021137.D
Acq On : 24 Jul 2024 23:06
Operator : MA/JU
Sample : P3244-12
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZG3

Manual IntegrationsAPPROVED

Quant Time: Jul 25 00:34:48 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 25 00:15:24 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/25/2024
Supervised By :mohammad ahmed 07/25/2024



TIC: BP021137.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.622min (+ 0.023) 4.60 ng/u1

response 22790

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	75.19
41.00	37.90	46.58#
42.00	27.10	29.13

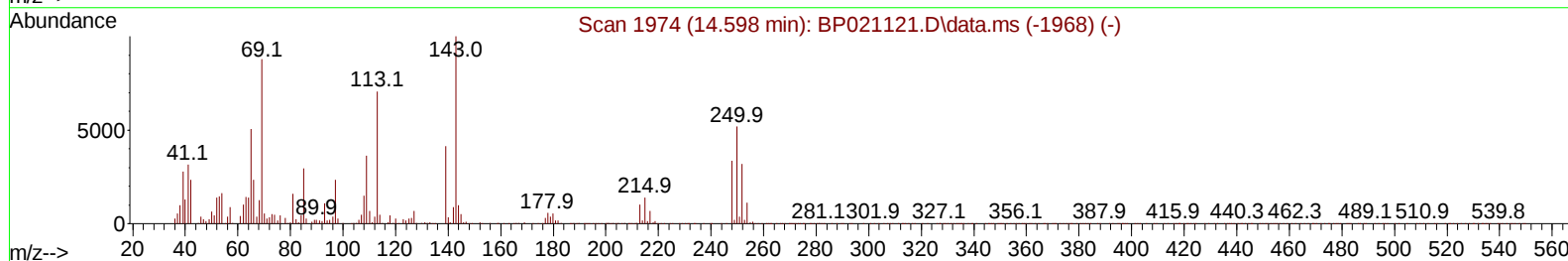
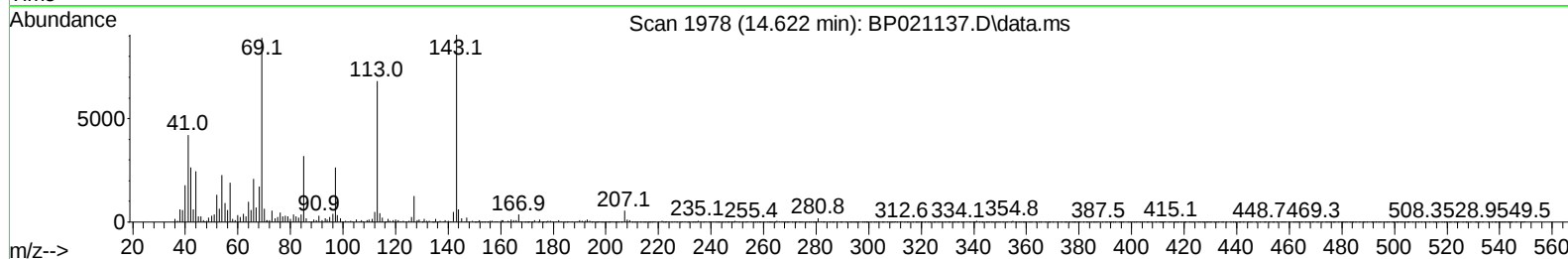
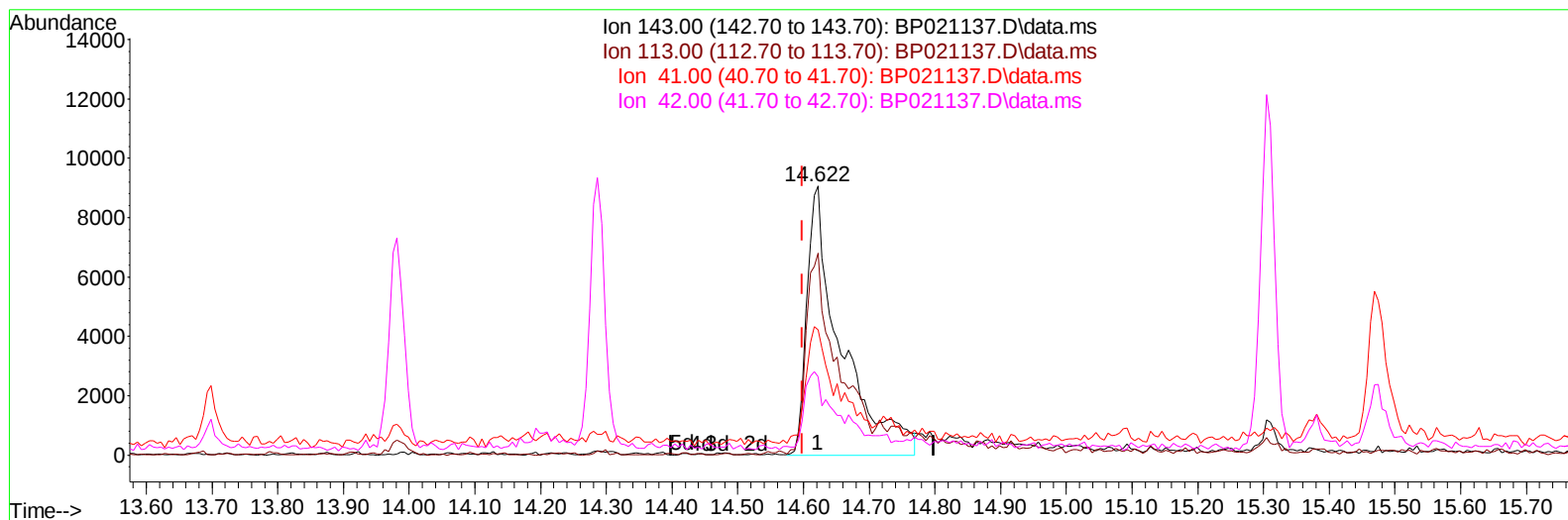
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TIC: BP021137.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.622min (+ 0.023) 6.53 ng/ul m

response 32362

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	75.19
41.00	37.90	46.58#
42.00	27.10	29.13

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/25/2024
 Supervised By :mohammad ahmed 07/25/2024

Quant Time: Jul 25 00:36:51 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 25 00:15:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	215140	20.000	ng/ul	0.00
20) Naphthalene-d8	10.428	136	811160	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.287	164	479066	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.110	188	942204	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	792886	20.000	ng/ul	-0.01
88) Perylene-d12	24.845	264	1076499	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	18620	3.938	ng/uL	0.00
4) Pyridine-d5	3.605	84	111470	8.144	ng/ul	0.00
7) Phenol-d5	6.881	99	120597	6.876	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.999	67	239074	22.491	ng/ul	0.00
11) 2-Chlorophenol-d4	7.211	132	309917	21.446	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113	196681	13.501	ng/ul	0.00
21) Nitrobenzene-d5	8.810	128	154121	24.462	ng/ul	0.00
24) 2-Nitrophenol-d4	9.522	143	181543	25.175	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	308853	23.686	ng/ul	0.00
31) 4-Chloroaniline-d4	10.581	131	254407	14.728	ng/ul	0.01
46) Dimethylphthalate-d6	13.698	166	942109	27.052	ng/ul	0.00
49) Acenaphthylene-d8	13.981	160	1019825	26.080	ng/ul	0.00
54) 4-Nitrophenol-d4	14.622	143	32362m	6.531	ng/ul	0.02
60) Fluorene-d10	15.304	176	768737	26.907	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	124427	21.826	ng/ul	0.00
73) Anthracene-d10	17.210	188	1166037	27.864	ng/ul	0.00
81) Pyrene-d10	19.592	212	1345619	27.451	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.639	264	1608166	30.290	ng/ul	-0.01

Target Compounds Qvalue

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

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