

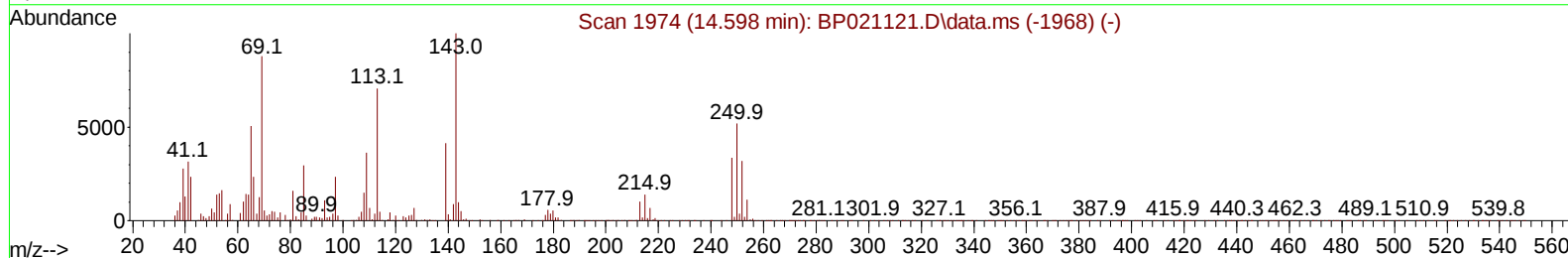
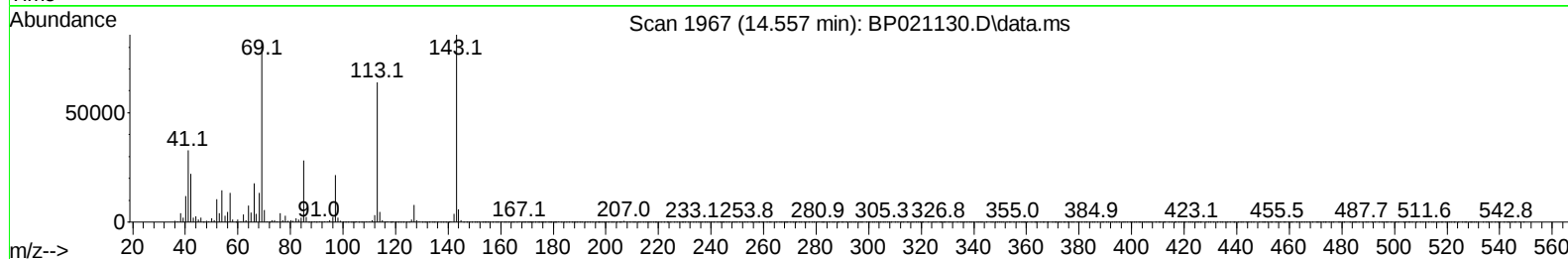
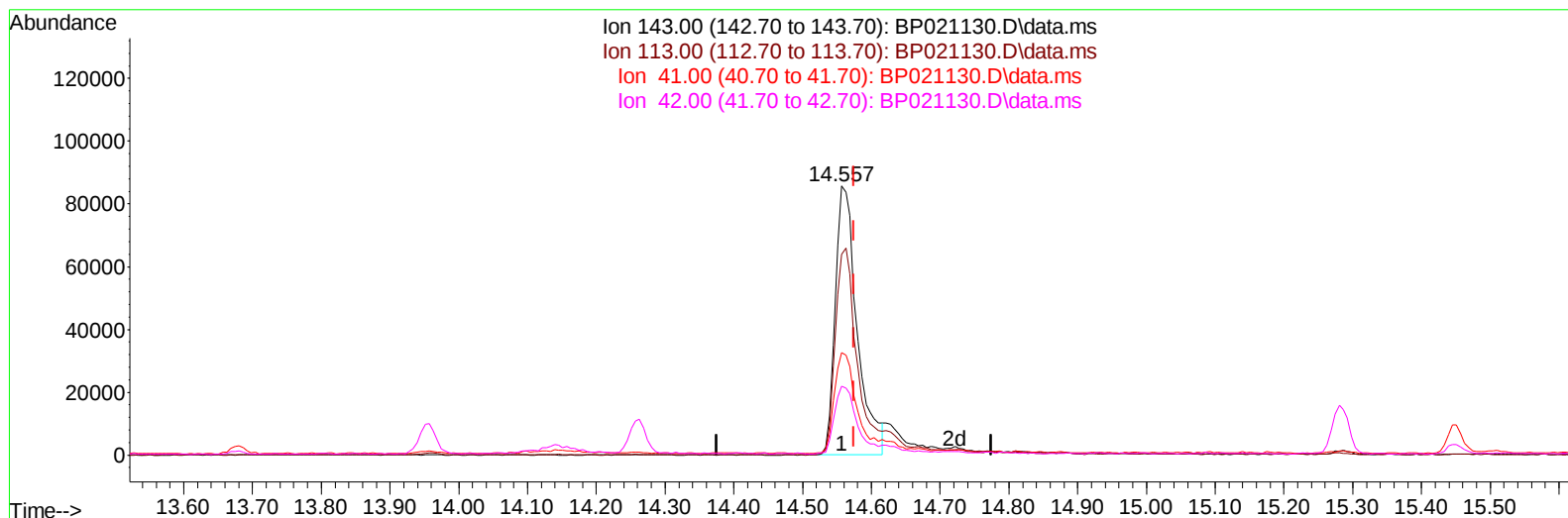
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072424\
Data File : BP021130.D
Acq On : 24 Jul 2024 18:02
Operator : MA/JU
Sample : PB162104BL
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK104

Manual IntegrationsAPPROVED

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

Quant Time: Jul 25 00:15:23 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



TIC: BP021130.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (-0.018) 28.89 ng/u1

response 188741

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	74.72
41.00	37.90	38.20
42.00	27.10	25.63

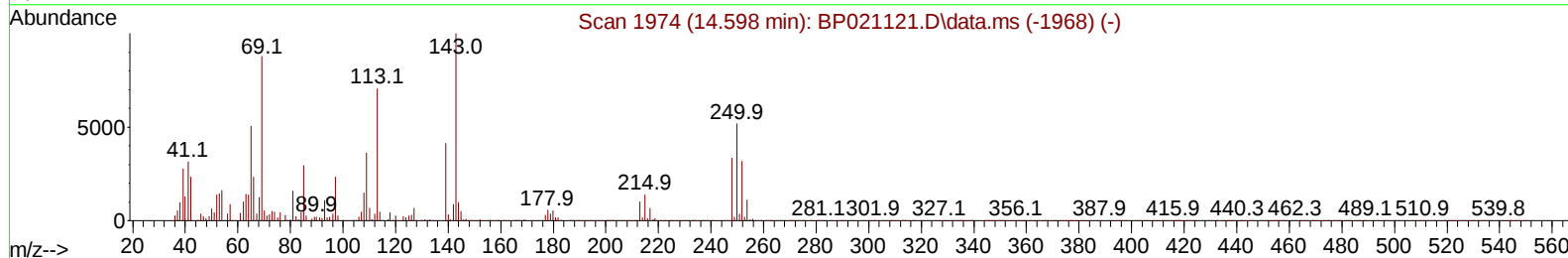
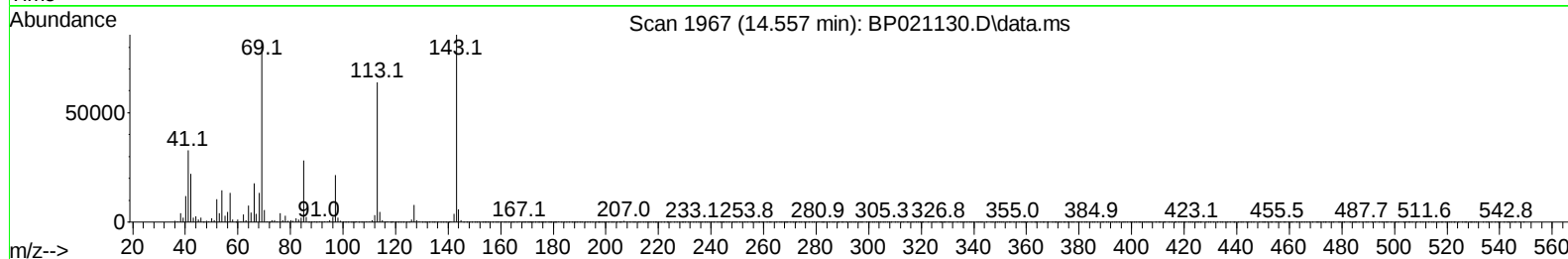
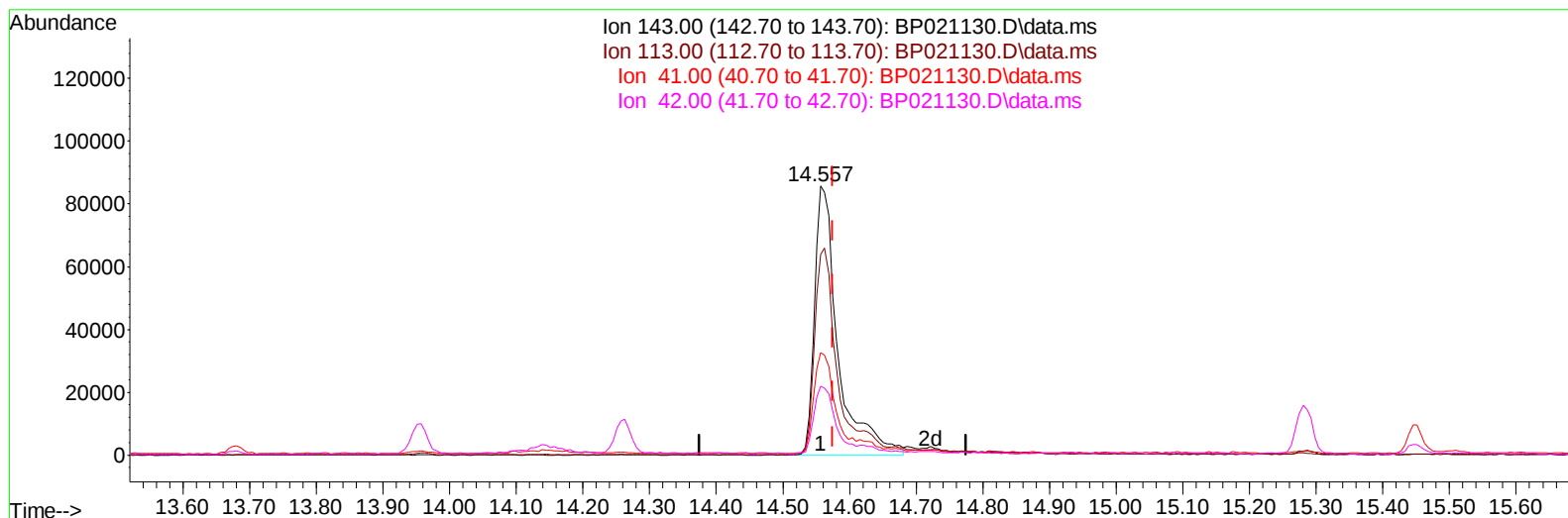
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TIC: BP021130.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (-0.018) 32.19 ng/ul m

response 210310

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	74.72
41.00	37.90	38.20
42.00	27.10	25.63

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072424\
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 Sample : PB162104BL
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 ALS Vial : 11 Sample Multiplier: 1

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 00:17:41 2024
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 QLast Update : Thu Jul 25 00:15:24 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.604	152	246490	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	952718	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.263	164	631727	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	1375105	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	1358605	20.000	ng/ul	0.00
88) Perylene-d12	24.856	264	1398702	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.157	96	30404	5.613	ng/uL	0.00
4) Pyridine-d5	3.575	84	351599	22.421	ng/ul	0.00
7) Phenol-d5	6.816	99	543643	27.053	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.951	67	327542	26.895	ng/ul	0.00
11) 2-Chlorophenol-d4	7.151	132	446173	26.948	ng/ul	-0.01
15) 4-Methylphenol-d8	8.334	113	433969	26.001	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	220254	29.765	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	258813	30.558	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.022	165	437088	28.540	ng/ul	-0.02
31) 4-Chloroaniline-d4	10.534	131	556604	27.434	ng/ul	0.00
46) Dimethylphthalate-d6	13.680	166	1317149	28.682	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	1504371	29.174	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	210310m	32.187	ng/ul	-0.02
60) Fluorene-d10	15.280	176	1133629	30.090	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.451	200	189246	22.746	ng/ul	0.00
73) Anthracene-d10	17.192	188	1820988	29.815	ng/ul	0.00
81) Pyrene-d10	19.586	212	2268612	27.009	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.639	264	2075757	30.091	ng/ul	0.00

Target Compounds Qvalue

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

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