

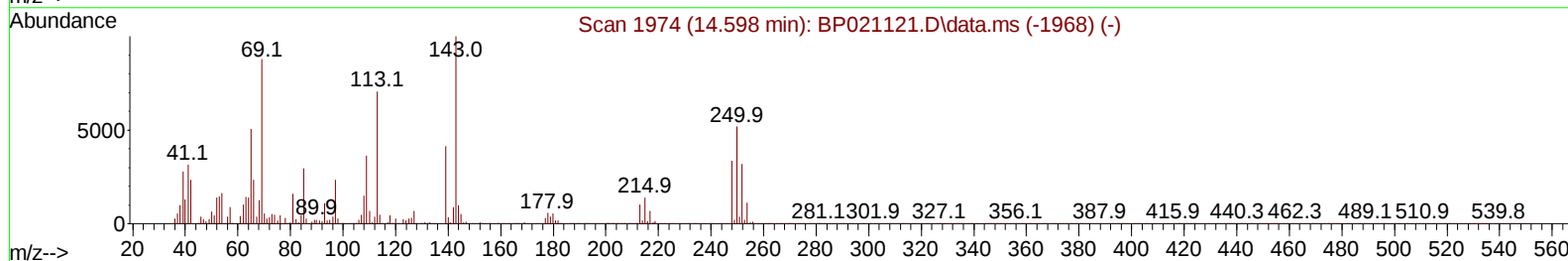
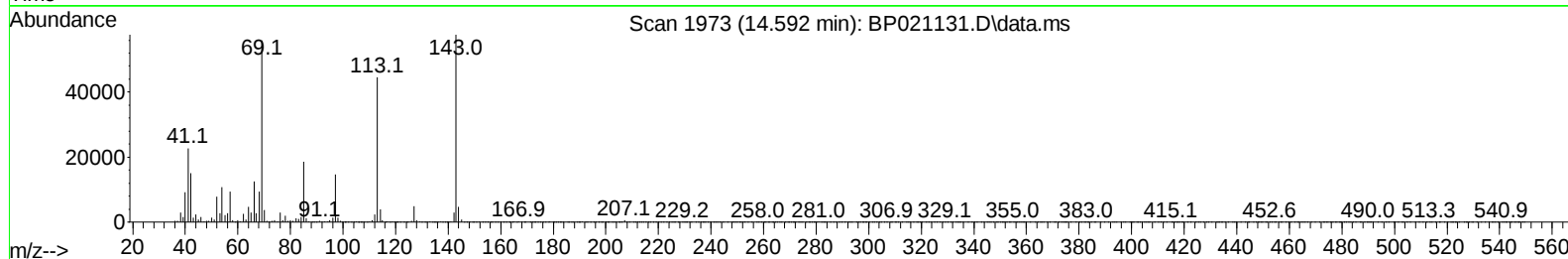
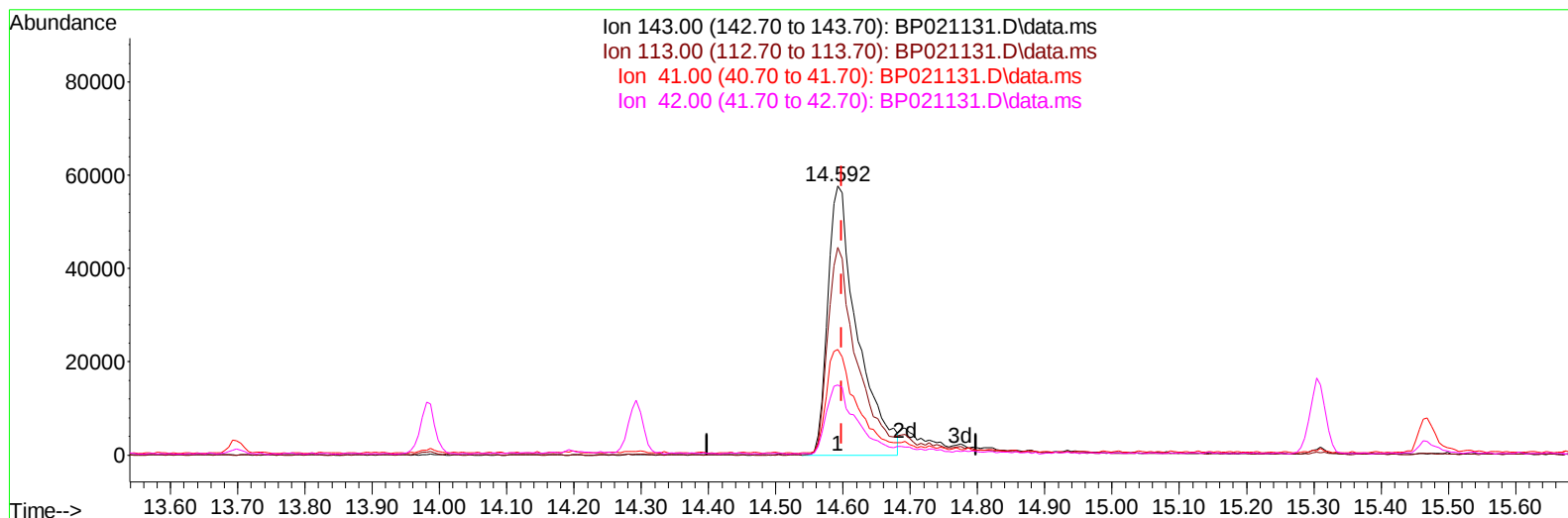
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072424\
Data File : BP021131.D
Acq On : 24 Jul 2024 18:45
Operator : MA/JU
Sample : PB162100BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK100

Manual IntegrationsAPPROVED

Quant Time: Jul 24 23:06:14 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jul 24 23:04:44 2024
Response via : Initial Calibration

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



TIC: BP021131.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.592min (-0.006) 27.13 ng/u1

response 175190

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	77.21
41.00	37.90	39.35
42.00	27.10	26.01

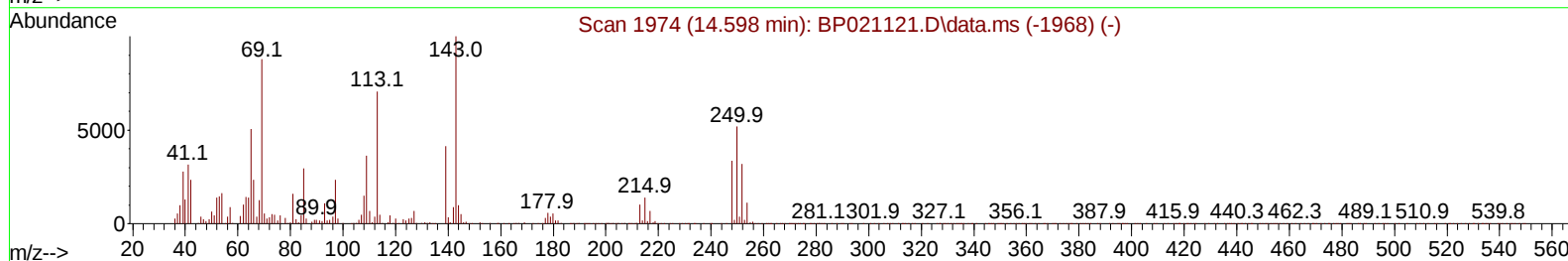
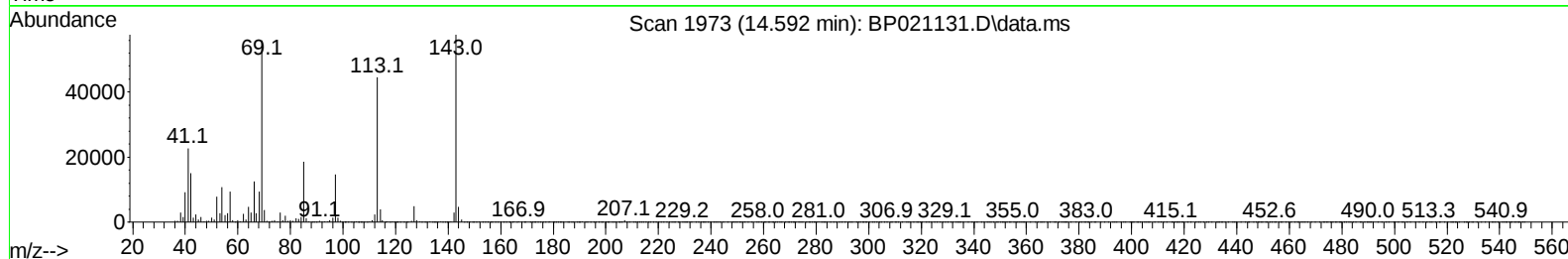
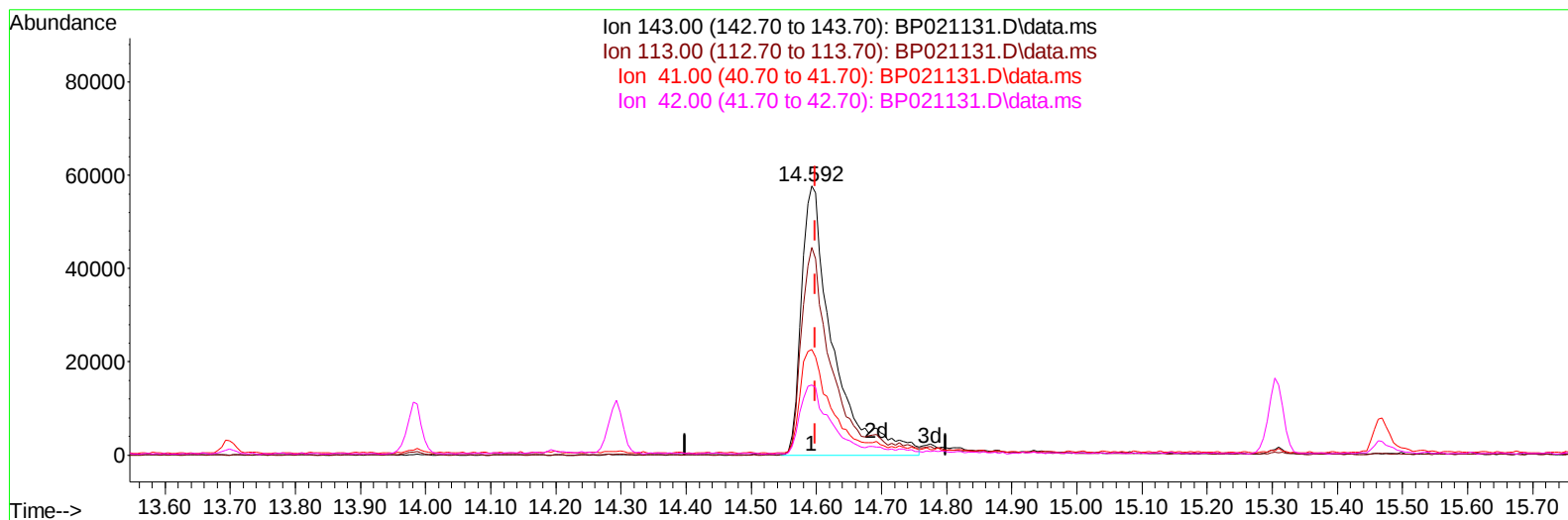
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TIC: BP021131.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.592min (-0.006) 29.64 ng/ul m

response 191431

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	77.21
41.00	37.90	39.35
42.00	27.10	26.01

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 Sample : PB162100BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK100

Manual IntegrationsAPPROVED

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

Quant Time: Jul 25 00:20:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 24 23:04:44 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	246342	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	974502	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.292	164	624410	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	1326082	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	1055651	20.000	ng/ul	0.00
88) Perylene-d12	24.868	264	1085170	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	28904	5.339	ng/uL	0.00
4) Pyridine-d5	3.599	84	346164	22.088	ng/ul	0.00
7) Phenol-d5	6.863	99	537209	26.749	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.999	67	328324	26.976	ng/ul	0.00
11) 2-Chlorophenol-d4	7.199	132	450197	27.207	ng/ul	0.00
15) 4-Methylphenol-d8	8.381	113	434848	26.069	ng/ul	0.00
21) Nitrobenzene-d5	8.810	128	222851	29.442	ng/ul	0.00
24) 2-Nitrophenol-d4	9.516	143	262132	30.258	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	427779	27.308	ng/ul	0.00
31) 4-Chloroaniline-d4	10.575	131	560422	27.005	ng/ul	0.00
46) Dimethylphthalate-d6	13.698	166	1321924	29.123	ng/ul	0.00
49) Acenaphthylene-d8	13.987	160	1498337	29.397	ng/ul	0.00
54) 4-Nitrophenol-d4	14.592	143	191431m	29.641	ng/ul	0.00
60) Fluorene-d10	15.310	176	1131573	30.387	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	176828	22.039	ng/ul	0.00
73) Anthracene-d10	17.204	188	1798640	30.538	ng/ul	0.00
81) Pyrene-d10	19.592	212	2078314	31.844	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.639	264	1601105	29.916	ng/ul	-0.01

Target Compounds Qvalue

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

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