

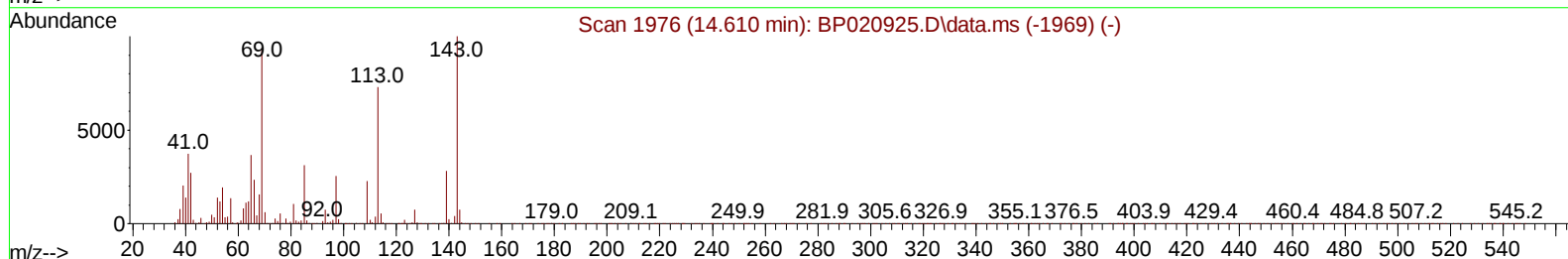
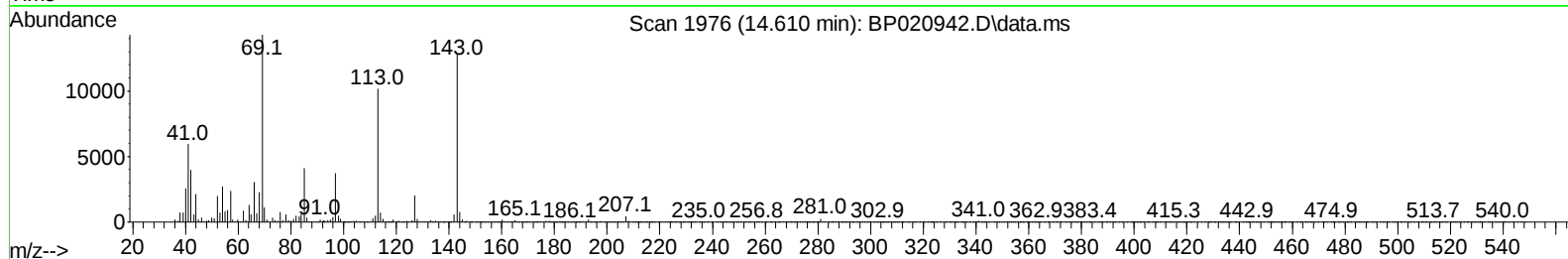
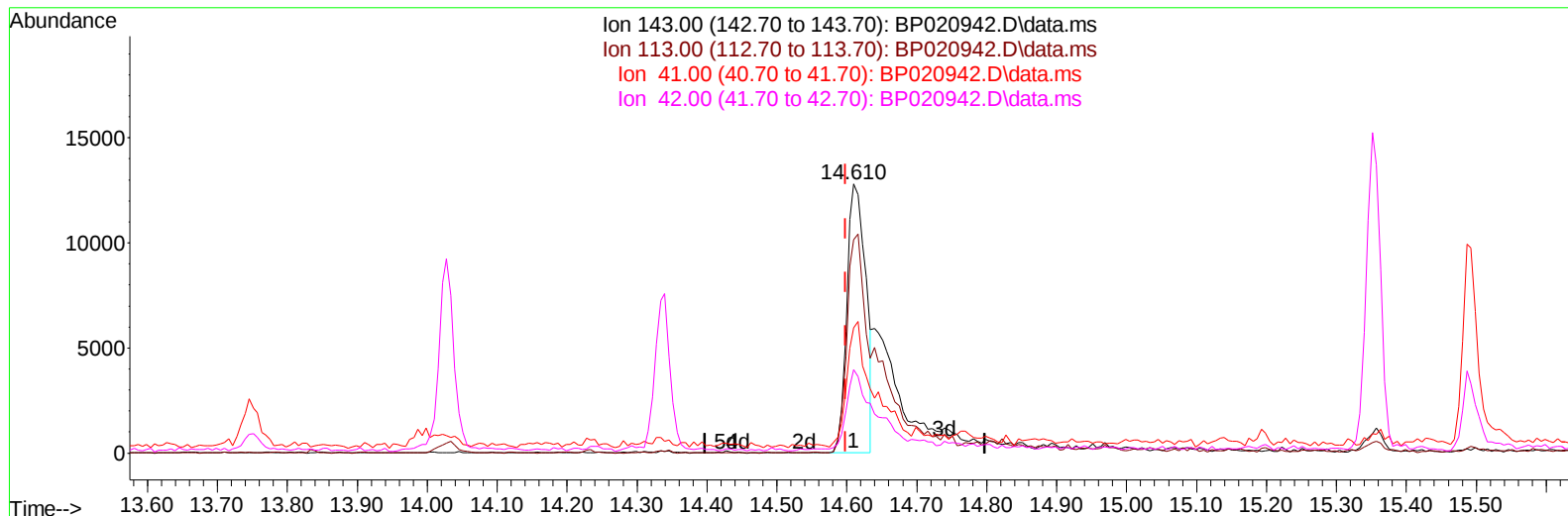
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071124\
Data File : BP020942.D
Acq On : 12 Jul 2024 20:54
Operator : MA/JU
Sample : P3231-03
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E08F2

Manual IntegrationsAPPROVED

Quant Time: Jul 12 23:35:08 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jul 12 23:34:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/13/2024
Supervised By :mohammad ahmed 07/13/2024



TIC: BP020942.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.610min (+ 0.012) 6.11 ng/u1

response 24231

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	79.42
41.00	37.90	46.53#
42.00	27.10	31.09

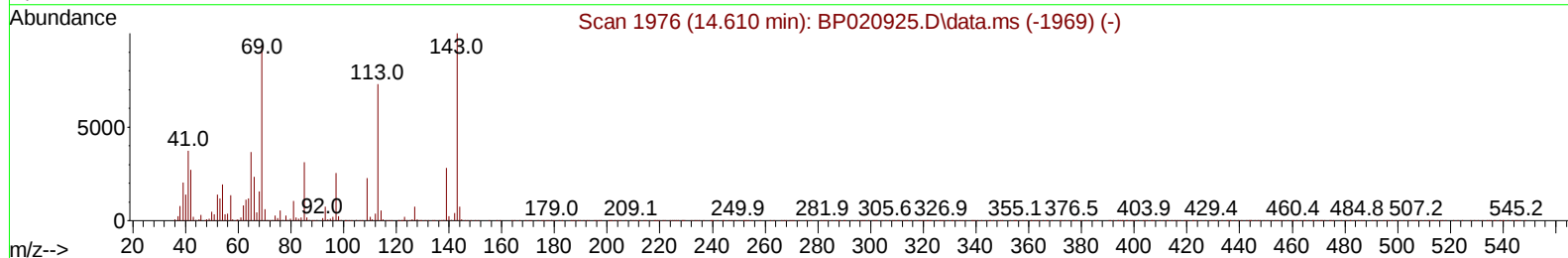
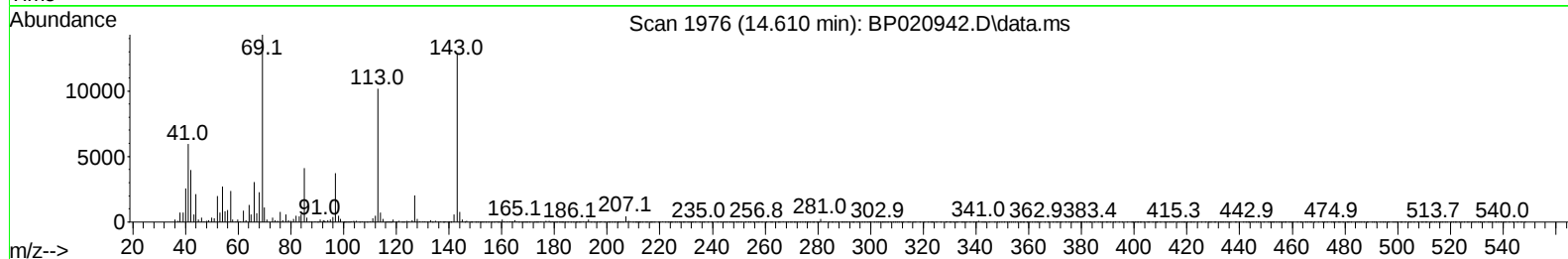
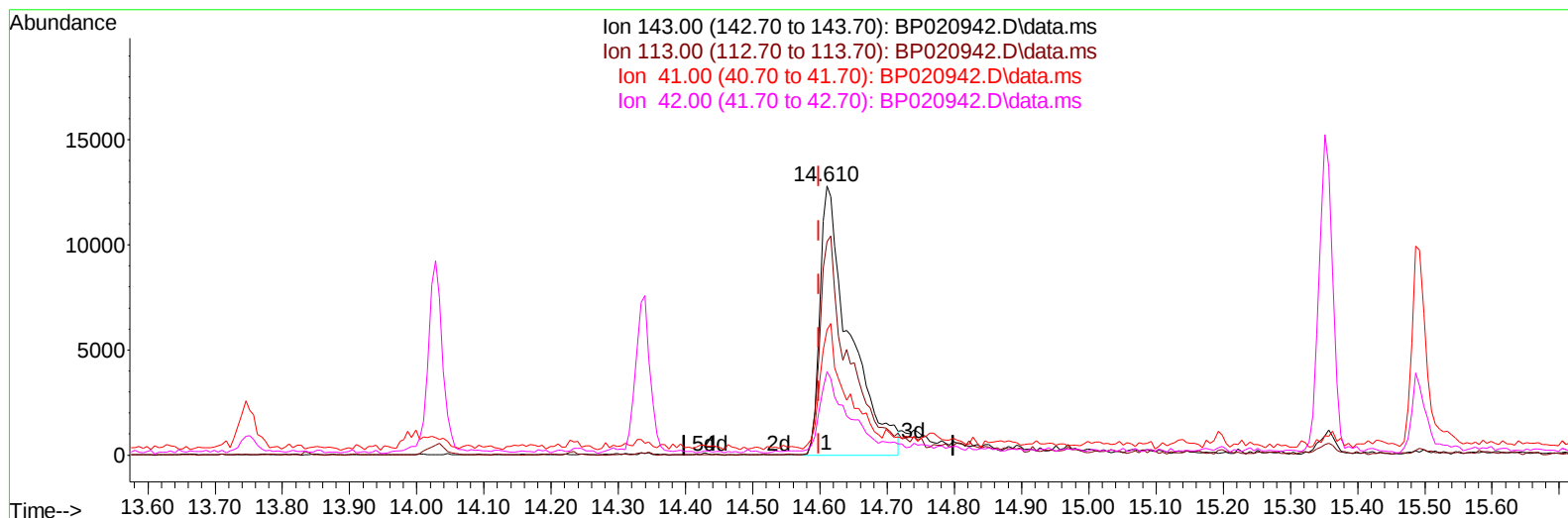
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TIC: BP020942.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.610min (+ 0.012) 9.92 ng/ul m

response 39361

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	79.42
41.00	37.90	46.53#
42.00	27.10	31.09

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/13/2024
 Supervised By :mohammad ahmed 07/13/2024

Quant Time: Jul 12 23:42:34 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 12 23:34:01 2024
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.716	152		144251	20.000	ng/ul	0.00
20) Naphthalene-d8	10.481	136		576397	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.340	164		383462	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.157	188		854124	20.000	ng/ul	0.00
79) Chrysene-d12	21.604	240		868199	20.000	ng/ul	0.00
88) Perylene-d12	24.939	264		984290	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.252	96		15429	4.867	ng/uL	0.00
4) Pyridine-d5	3.670	84		51559	5.618	ng/ul	0.00
7) Phenol-d5	6.911	99		112873	9.598	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67		234324	32.878	ng/ul	0.01
11) 2-Chlorophenol-d4	7.263	132		286660	29.584	ng/ul	0.01
15) 4-Methylphenol-d8	8.428	113		202332	20.715	ng/ul	0.00
21) Nitrobenzene-d5	8.863	128		151510	33.842	ng/ul	0.00
24) 2-Nitrophenol-d4	9.575	143		171906	33.548	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.122	165		291185	31.426	ng/ul	0.01
31) 4-Chloroaniline-d4	10.622	131		326245	26.578	ng/ul	0.00
46) Dimethylphthalate-d6	13.751	166		1072018	38.457	ng/ul	0.00
49) Acenaphthylene-d8	14.028	160		1136550	36.311	ng/ul	0.00
54) 4-Nitrophenol-d4	14.610	143		39361m	9.924	ng/ul	0.01
60) Fluorene-d10	15.351	176		899034	39.312	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200		170234	32.941	ng/ul	0.00
73) Anthracene-d10	17.251	188		1521851	40.116	ng/ul	0.00
81) Pyrene-d10	19.639	212		1853051	34.523	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.710	264		1920985	39.571	ng/ul	0.00

Target Compounds Qvalue

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

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