

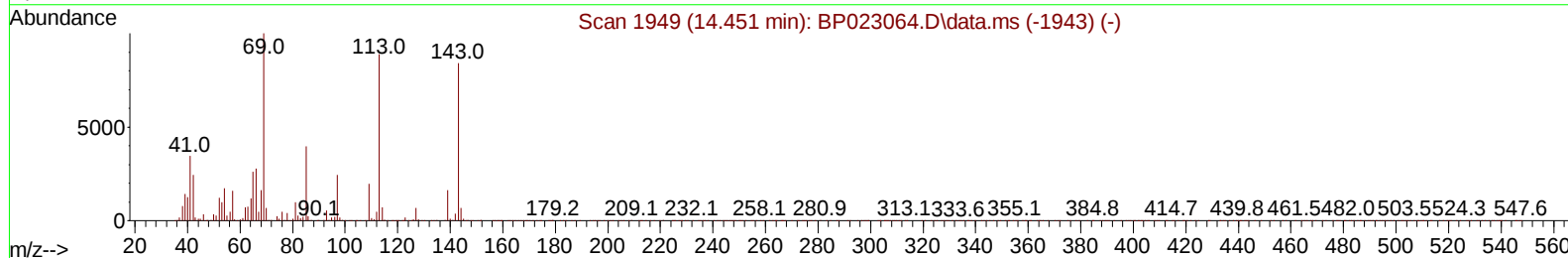
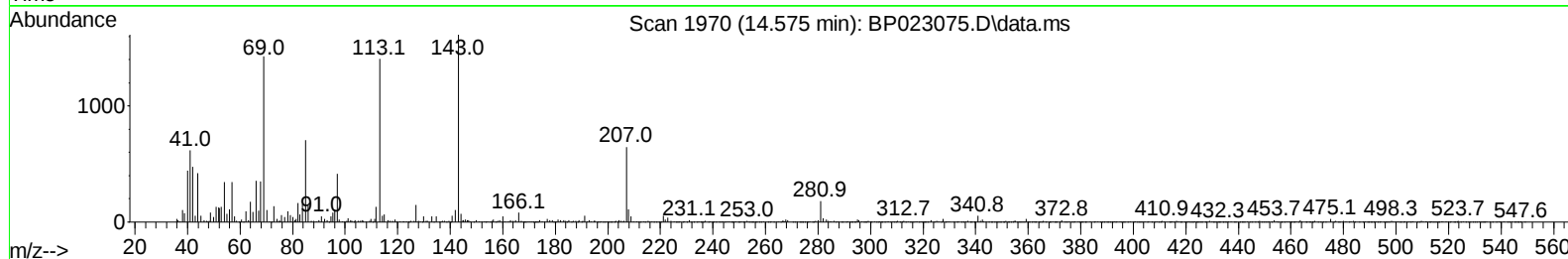
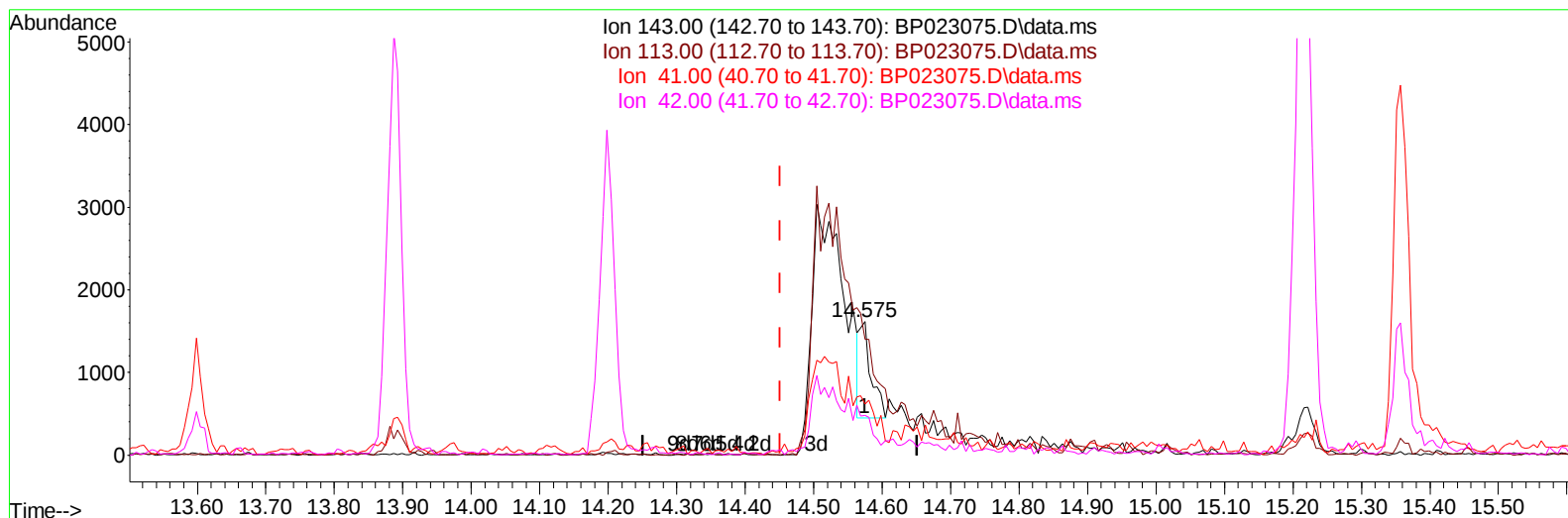
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111224\
Data File : BP023075.D
Acq On : 13 Nov 2024 10:47
Operator : RC/JU
Sample : P4687-14
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0CD7

Manual IntegrationsAPPROVED

Quant Time: Nov 14 00:41:32 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 13 04:49:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2024
Supervised By :mohammad ahmed 11/14/2024



TIC: BP023075.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.575min (+ 0.124) 0.66 ng/u1

response 1362

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	86.99
41.00	39.20	38.17
42.00	25.90	29.62

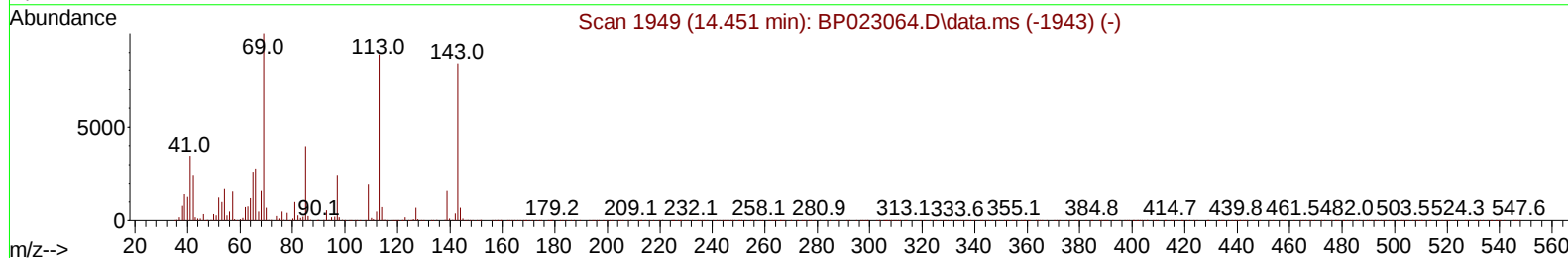
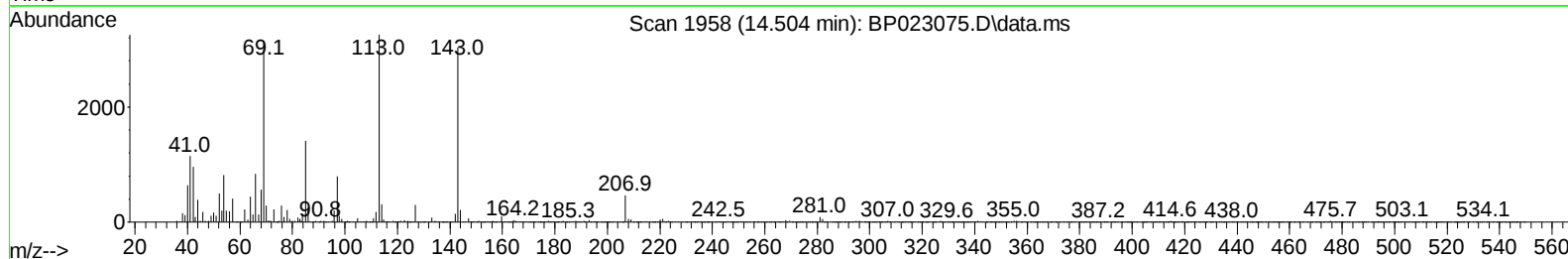
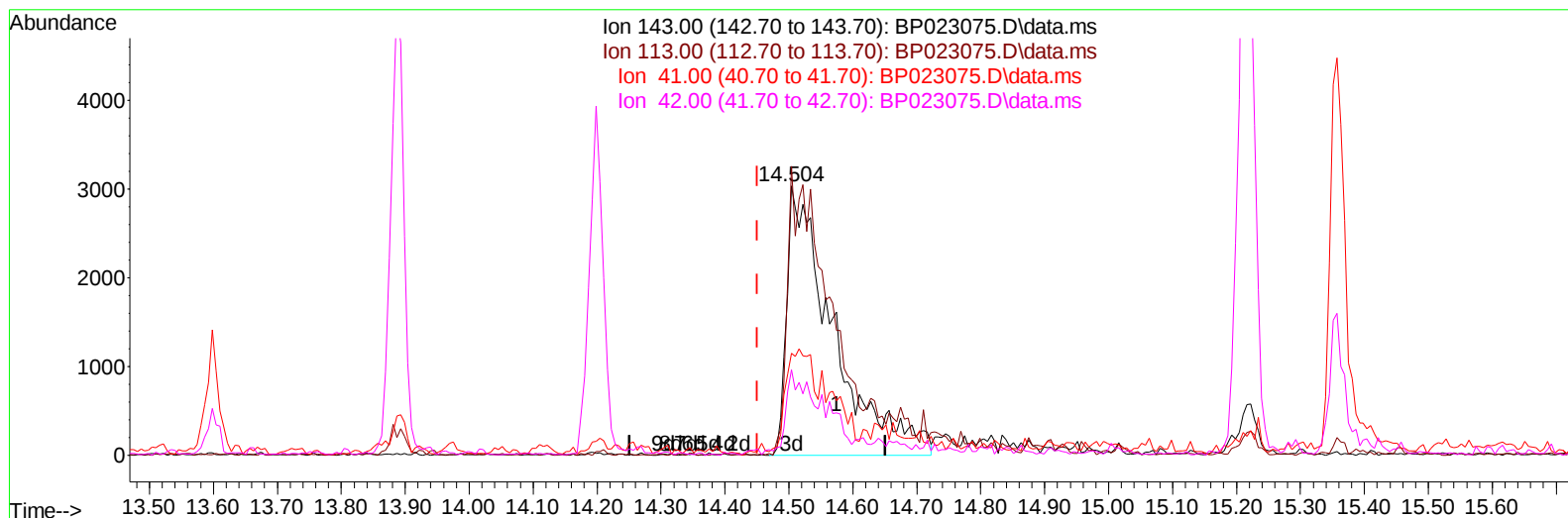
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(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.053) 7.38 ng/ul m

response 15201

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	107.32
41.00	39.20	37.80
42.00	25.90	31.81#

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 Misc :
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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/14/2024
 Supervised By :mohammad ahmed 11/14/2024

Quant Time: Nov 14 01:04:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	61993	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136	238541	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.204	164	190246	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.010	188	486581	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	594290	20.000	ng/ul	-0.01
88) Perylene-d12	24.639	264	720701	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	6707	5.200	ng/uL	0.00
4) Pyridine-d5	3.564	84	32512	8.485	ng/ul	0.02
7) Phenol-d5	6.793	99	40470	8.830	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.928	67	105160	39.052	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	107930	32.154	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	85713	22.726	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	65807	39.627	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	77589	39.190	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	152836	36.863	ng/ul	0.01
31) 4-Chloroaniline-d4	10.481	131	145576	29.431	ng/ul	0.00
46) Dimethylphthalate-d6	13.598	166	579525	42.265	ng/ul	-0.01
49) Acenaphthylene-d8	13.887	160	569026	39.772	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.504	143	15201m	7.377	ng/ul	0.05
60) Fluorene-d10	15.216	176	501919	41.751	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.357	200	108591	37.254	ng/ul	0.00
73) Anthracene-d10	17.104	188	950907	46.155	ng/ul	-0.01
81) Pyrene-d10	19.486	212	1258681	45.982	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.421	264	1578899	45.992	ng/ul	-0.02

Target Compounds Qvalue

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

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