

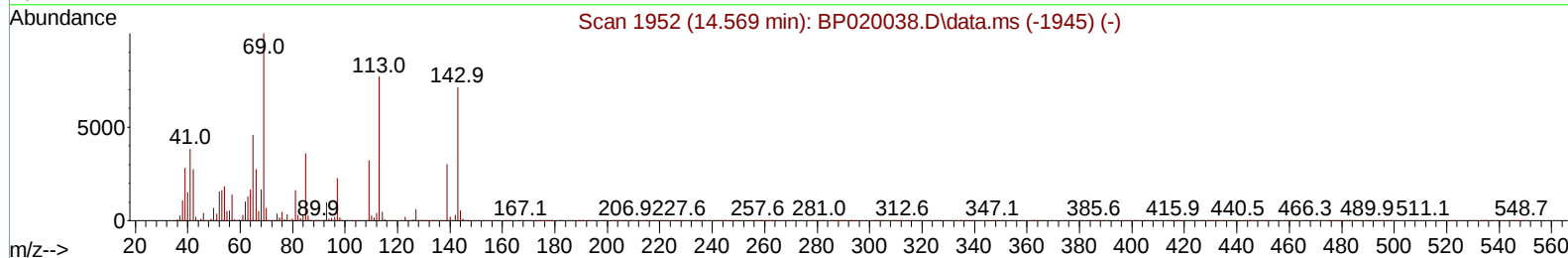
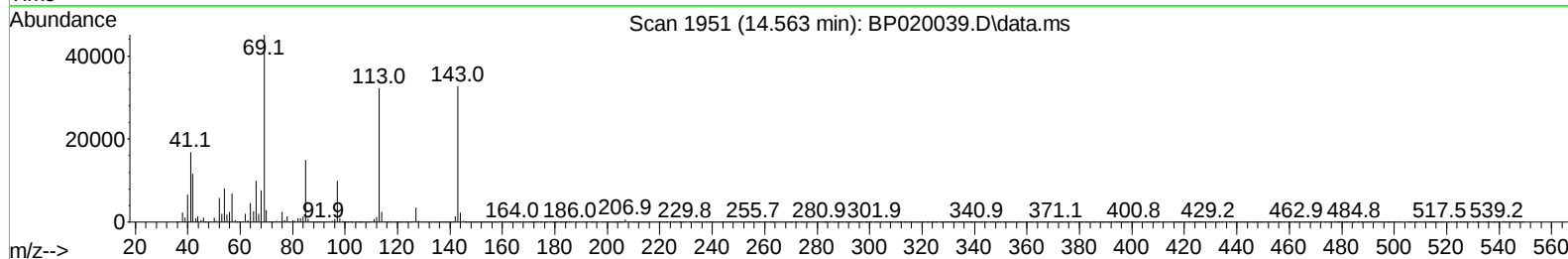
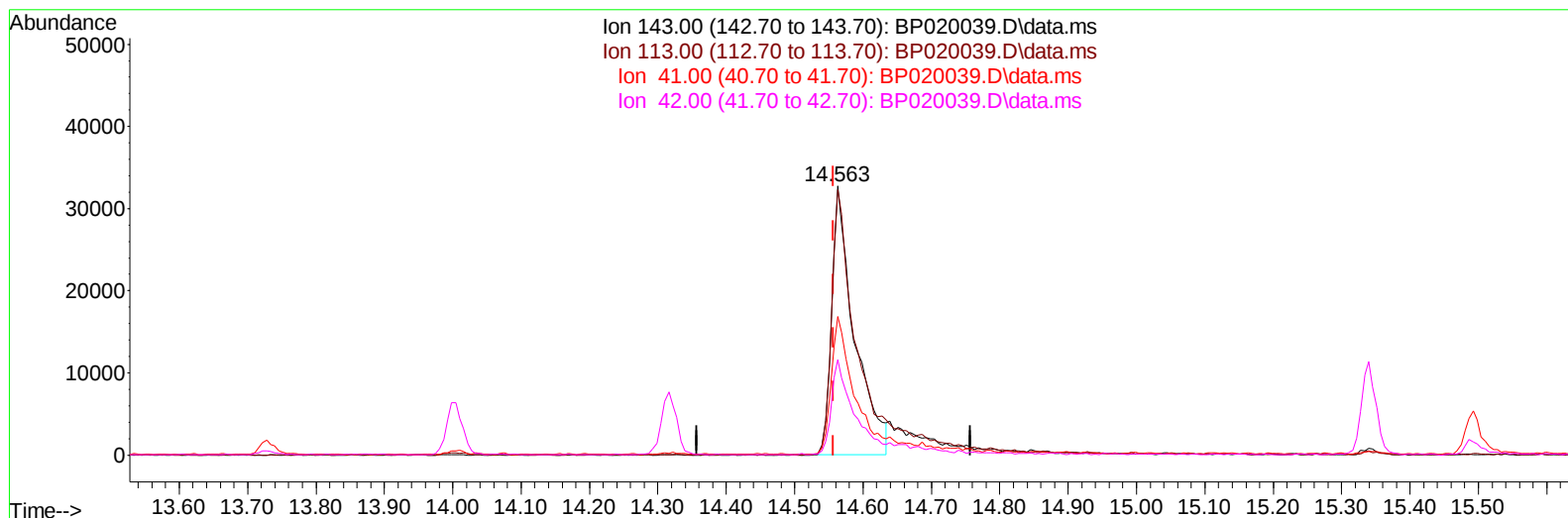
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042024\
Data File : BP020039.D
Acq On : 20 Apr 2024 11:26
Operator : MA/JU
Sample : PB160383BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK383

Manual IntegrationsAPPROVED

Quant Time: Apr 21 23:42:01 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Apr 20 01:32:50 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/24/2024
Supervised By :mohammad ahmed 04/30/2024



TIC: BP020039.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.006) 23.56 ng/u1

response 75335

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.80	98.66
41.00	50.70	51.64
42.00	34.00	35.53

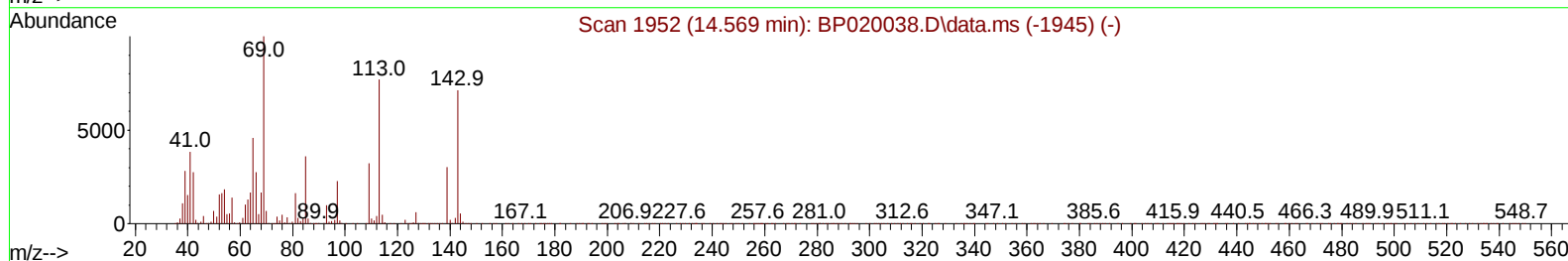
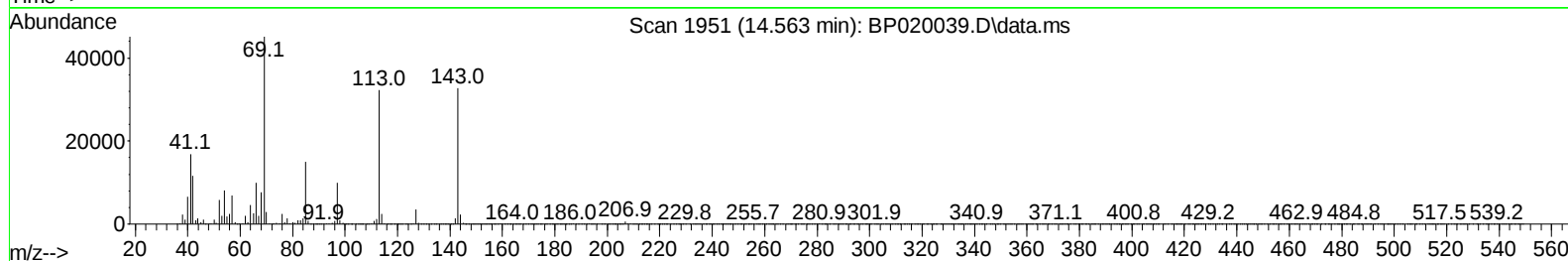
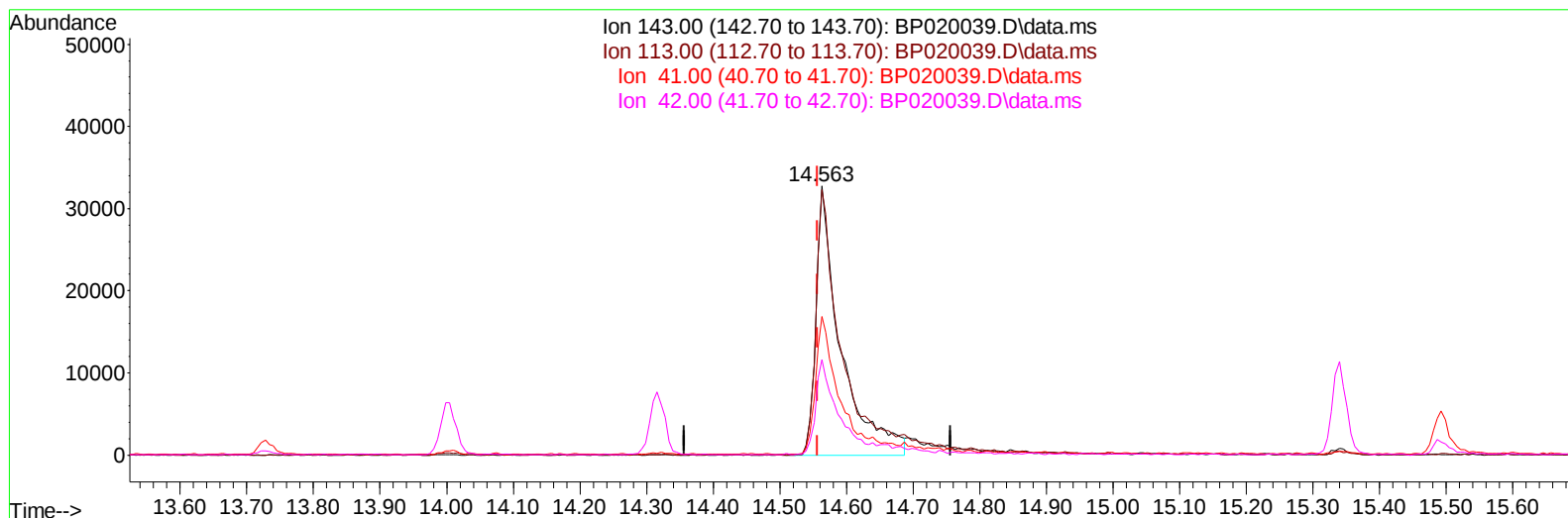
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TIC: BP020039.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.006) 26.40 ng/ul m

response 84423

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.80	98.66
41.00	50.70	51.64
42.00	34.00	35.53

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/24/2024
 Supervised By :mohammad ahmed 04/30/2024

Quant Time: Apr 21 23:51:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 20 01:32:50 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	112103	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136	451451	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	307335	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.163	188	646517	20.000	ng/ul	0.00
79) Chrysene-d12	21.669	240	612005	20.000	ng/ul	0.01
88) Perylene-d12	25.062	264	672203	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	14392	5.260	ng/uL	0.00
4) Pyridine-d5	3.622	84	211850	26.307	ng/ul	-0.01
7) Phenol-d5	6.816	99	256256	26.228	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	181590	26.591	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.169	132	183437	27.498	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	195426	25.606	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	94886	29.219	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	102641	29.241	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.046	165	168244	24.160	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	251640	27.069	ng/ul	0.00
46) Dimethylphthalate-d6	13.728	166	621522	29.591	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	689199	27.530	ng/ul	0.00
54) 4-Nitrophenol-d4	14.563	143	84423m	26.404	ng/ul	0.00
60) Fluorene-d10	15.339	176	532209	28.465	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	90986	24.089	ng/ul	0.00
73) Anthracene-d10	17.263	188	819182	28.282	ng/ul	0.00
81) Pyrene-d10	19.663	212	1003593	30.514	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.821	264	971865	29.597	ng/ul	0.01

Target Compounds Qvalue

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

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