

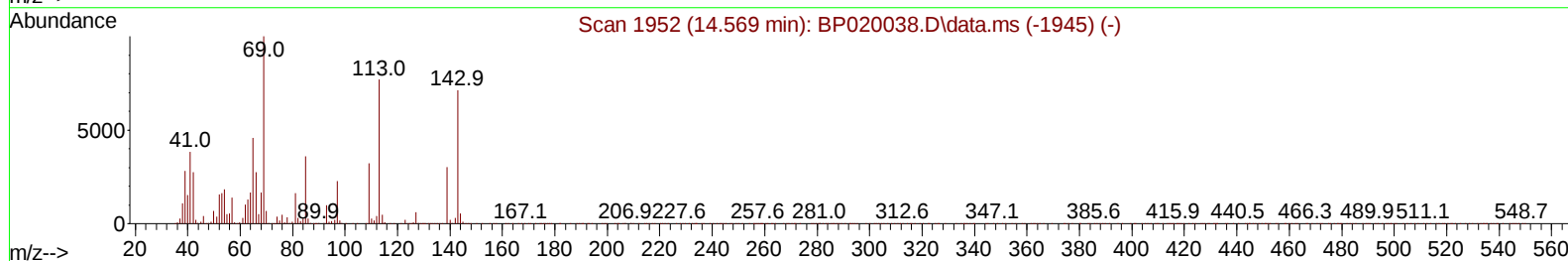
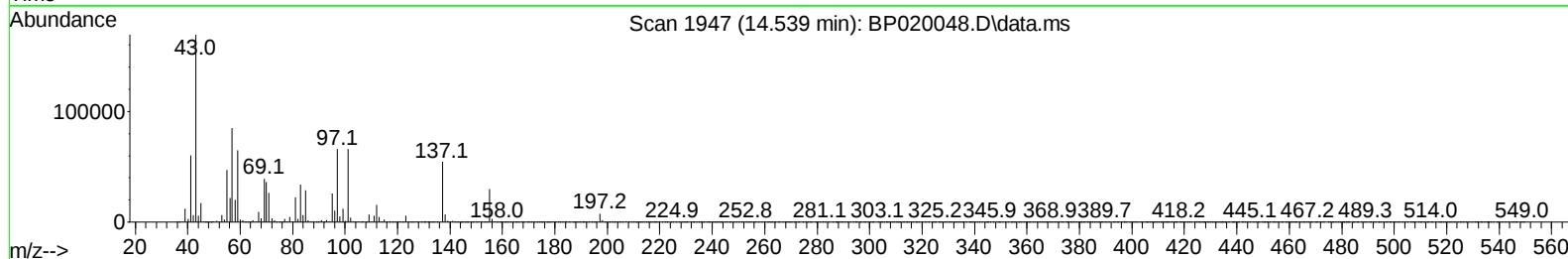
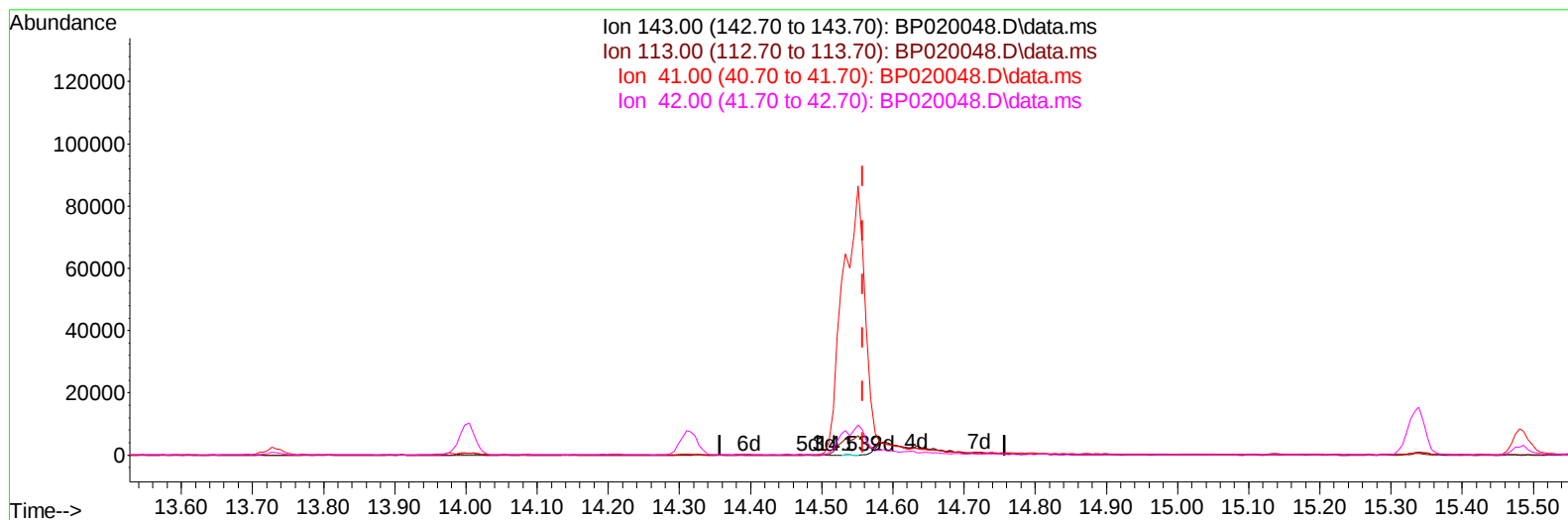
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042024\
Data File : BP020048.D
Acq On : 20 Apr 2024 18:01
Operator : MA/JU
Sample : P2161-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0RC6

Manual IntegrationsAPPROVED

Quant Time: Apr 21 23:43:47 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: BP020048.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.539min (-0.018) 0.01 ng/u1

response 39

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.80	6609.86#
41.00	50.70	84630.99#
42.00	34.00	8823.94#

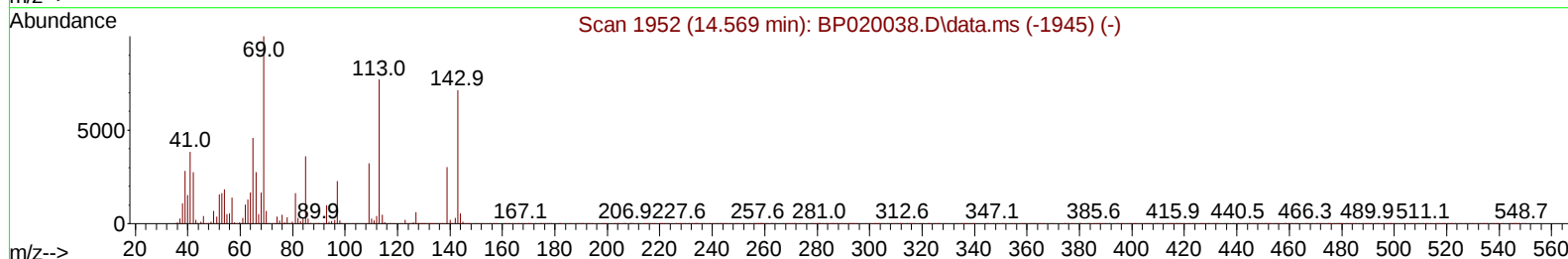
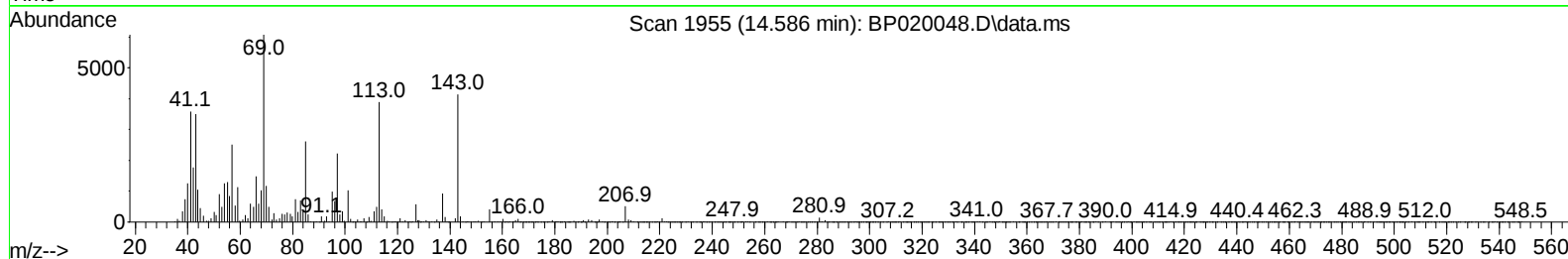
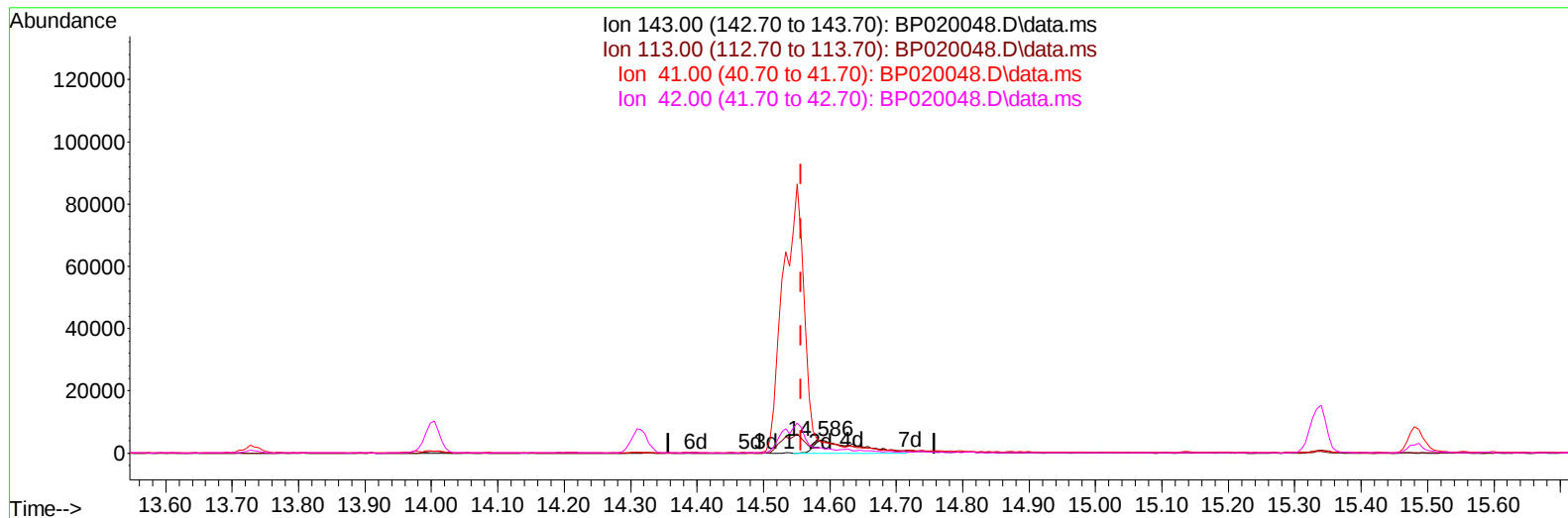
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TIC: BP020048.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.586min (+ 0.029) 5.77 ng/ul m

response 18673

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.80	93.99
41.00	50.70	86.69#
42.00	34.00	42.90#

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 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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

Quant Time: Apr 22 00:12:04 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	7.640	152	100193	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136	434496	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.316	164	311029	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.151	188	663876	20.000	ng/ul	0.00
79) Chrysene-d12	21.663	240	473205	20.000	ng/ul	0.00
88) Perylene-d12	25.056	264	533208	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	8745	3.576	ng/uL	0.00
4) Pyridine-d5	3.658	84	8425	1.171	ng/ul	0.02
7) Phenol-d5	6.822	99	56704	6.494	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	234066	38.349	ng/ul	0.00
11) 2-Chlorophenol-d4	7.175	132	171429	28.753	ng/ul	0.00
15) 4-Methylphenol-d8	8.346	113	114058	16.721	ng/ul	0.00
21) Nitrobenzene-d5	8.799	128	123567	39.536	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	132801	39.309	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.046	165	221933	33.114	ng/ul	0.00
31) 4-Chloroaniline-d4	10.575	131	106776	11.934	ng/ul	0.00
46) Dimethylphthalate-d6	13.728	166	899858	42.334	ng/ul	0.00
49) Acenaphthylene-d8	13.998	160	985470	38.897	ng/ul	0.00
54) 4-Nitrophenol-d4	14.586	143	18673m	5.771	ng/ul	0.03
60) Fluorene-d10	15.339	176	771190	40.757	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	139793	36.044	ng/ul	0.00
73) Anthracene-d10	17.257	188	1248991	41.994	ng/ul	0.00
81) Pyrene-d10	19.663	212	1307266	51.405	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.815	264	1112362	42.706	ng/ul	0.00

Target Compounds Qvalue

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

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