

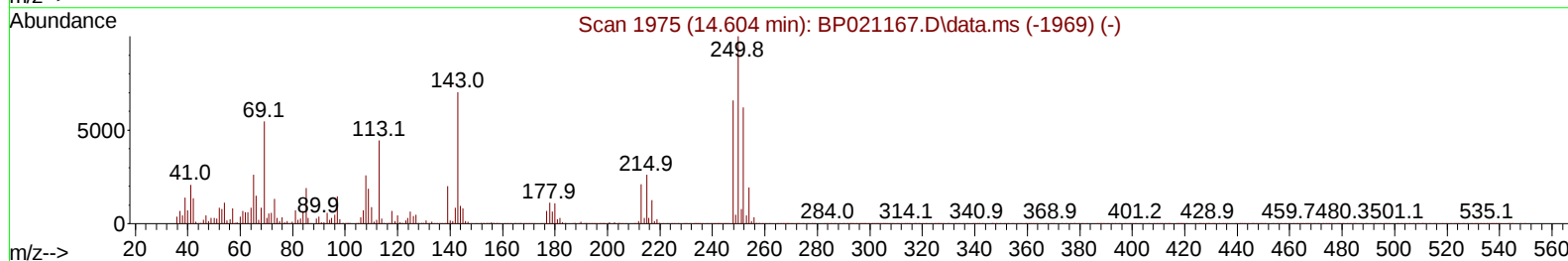
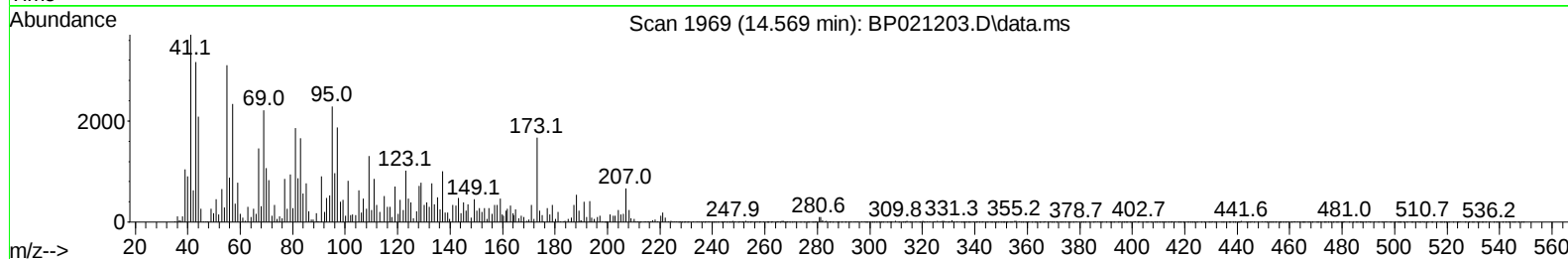
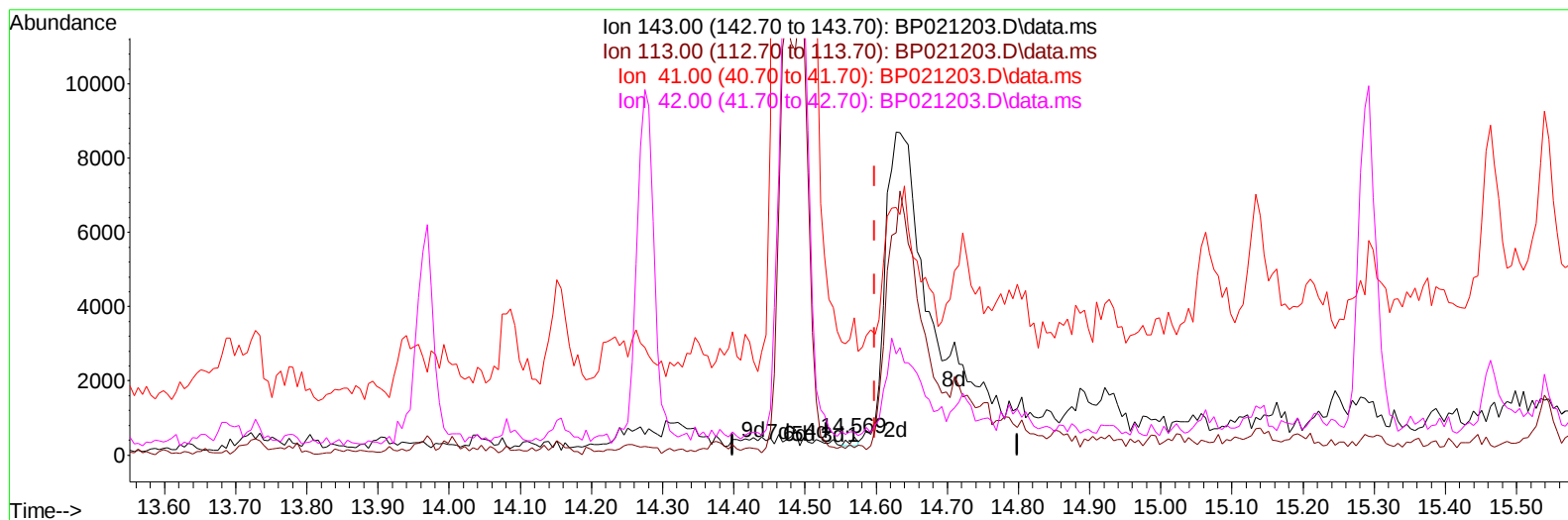
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
Data File : BP021203.D
Acq On : 27 Jul 2024 15:55
Operator : MA/JU
Sample : P3300-11
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G65

Manual IntegrationsAPPROVED

Quant Time: Jul 29 00:44:38 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 25 00:15:24 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/29/2024
Supervised By :mohammad ahmed 07/29/2024



TIC: BP021203.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (-0.030) 0.04 ng/u1

response 202

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	42.83#
41.00	37.90	770.69#
42.00	27.10	130.35#

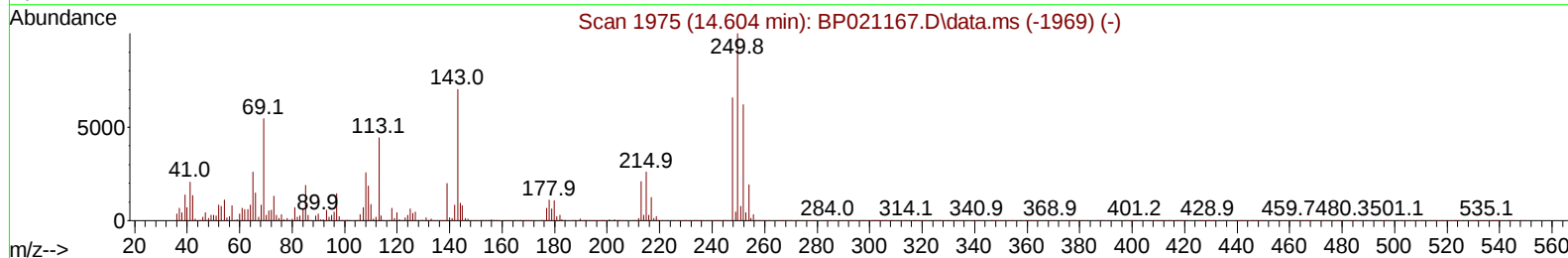
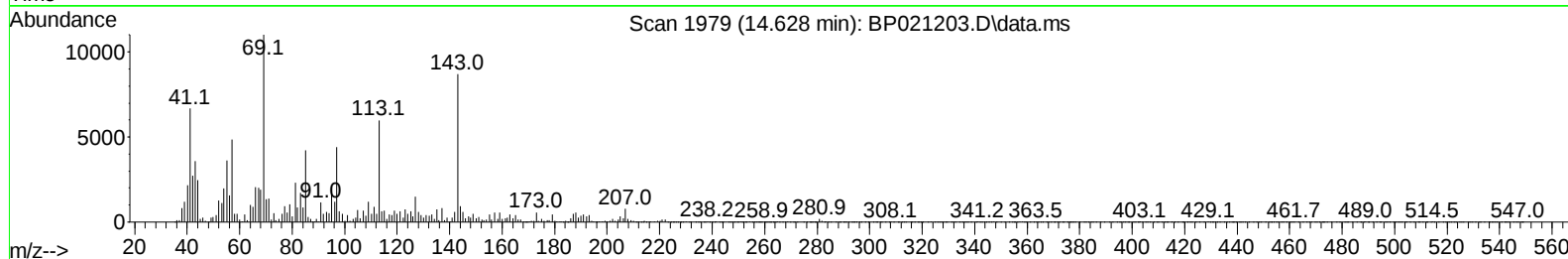
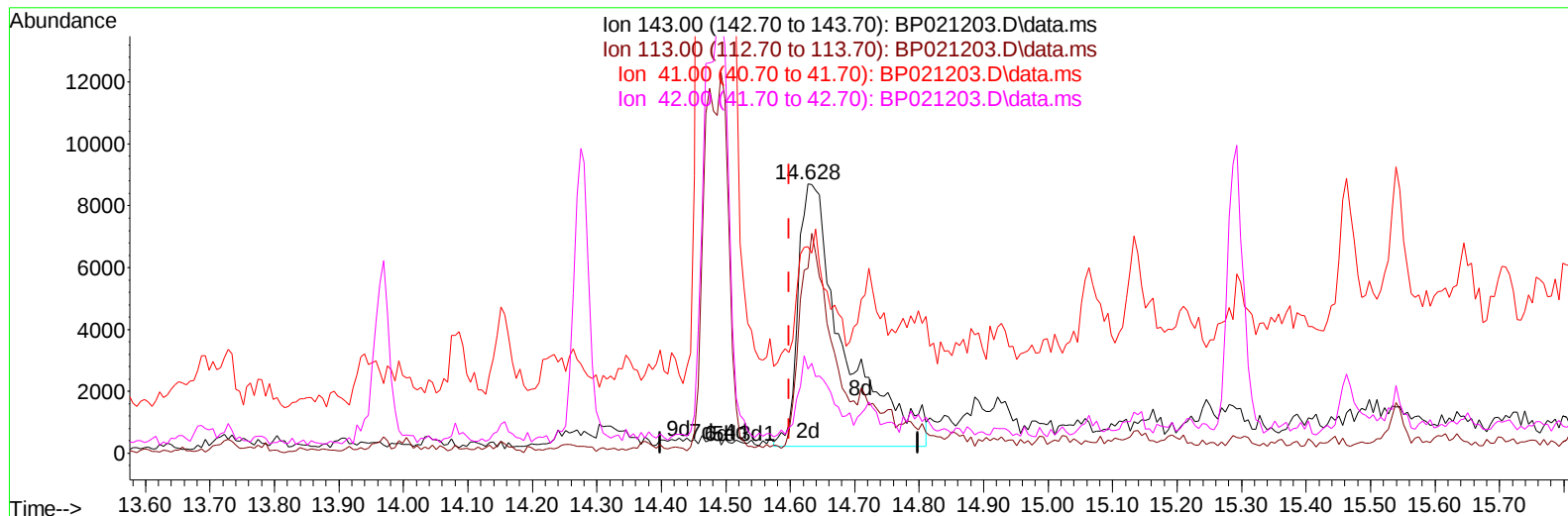
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TIC: BP021203.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.029) 7.73 ng/ul m

response 42226

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	68.76
41.00	37.90	76.71#
42.00	27.10	31.25

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Instrument :
 BNA_P
 ClientSampleId :
 E1G65

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/29/2024
 Supervised By :mohammad ahmed 07/29/2024

Quant Time: Jul 29 01:33:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 25 00:15:24 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	209422	20.000	ng/ul	0.00
20) Naphthalene-d8	10.404	136	807482	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.275	164	528235	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.092	188	1105766	20.000	ng/ul	-0.01
79) Chrysene-d12	21.551	240	1092753	20.000	ng/ul	0.00
88) Perylene-d12	24.857	264	1306417	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	14718	3.198	ng/uL	0.00
4) Pyridine-d5	3.611	84	38288	2.874	ng/ul	0.01
7) Phenol-d5	6.869	99	105167	6.160	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	169630	16.394	ng/ul	0.00
11) 2-Chlorophenol-d4	7.193	132	234600	16.677	ng/ul	-0.01
15) 4-Methylphenol-d8	8.375	113	173989	12.270	ng/ul	0.00
21) Nitrobenzene-d5	8.804	128	101432	16.173	ng/ul	0.00
24) 2-Nitrophenol-d4	9.516	143	117342	16.346	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	222988	17.179	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.575	131	272284	15.834	ng/ul	0.00
46) Dimethylphthalate-d6	13.692	166	702839	18.303	ng/ul	0.00
49) Acenaphthylene-d8	13.969	160	745838	17.298	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.628	143	42226m	7.729	ng/ul	0.03
60) Fluorene-d10	15.292	176	614079	19.493	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.463	200	100307	14.993	ng/ul	0.00
73) Anthracene-d10	17.204	188	974237	19.837	ng/ul	0.00
81) Pyrene-d10	19.586	212	1253387	18.553	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.627	264	1378991	21.402	ng/ul	-0.02
Target Compounds						
78) Di-n-butylphthalate	18.098	149	82335	1.248	ng/ul#	96
86) Bis(2-ethylhexyl)phtha...	21.439	149	53538	1.076	ng/ul#	94

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

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