

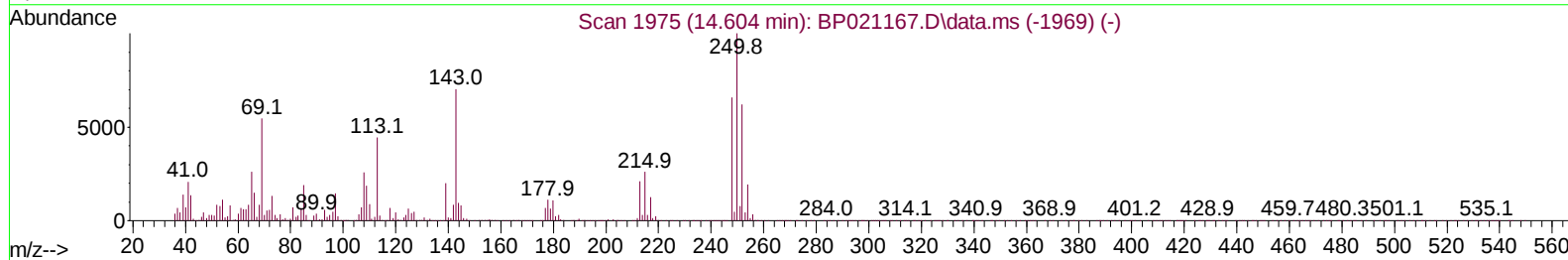
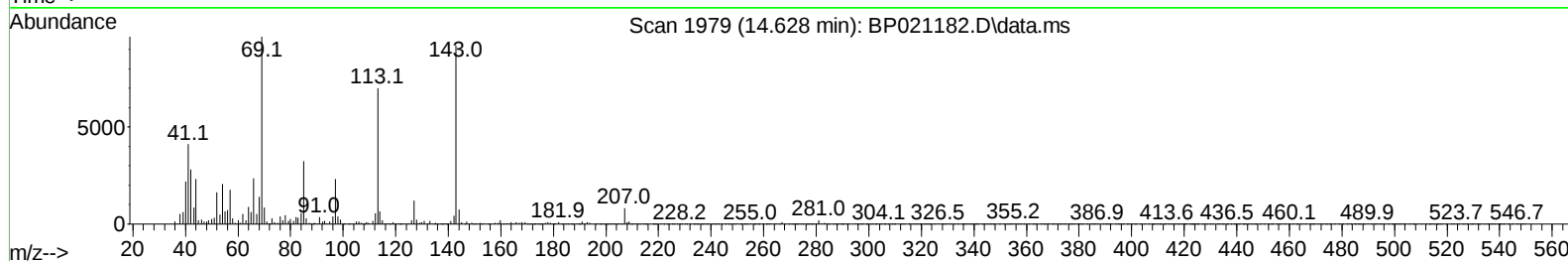
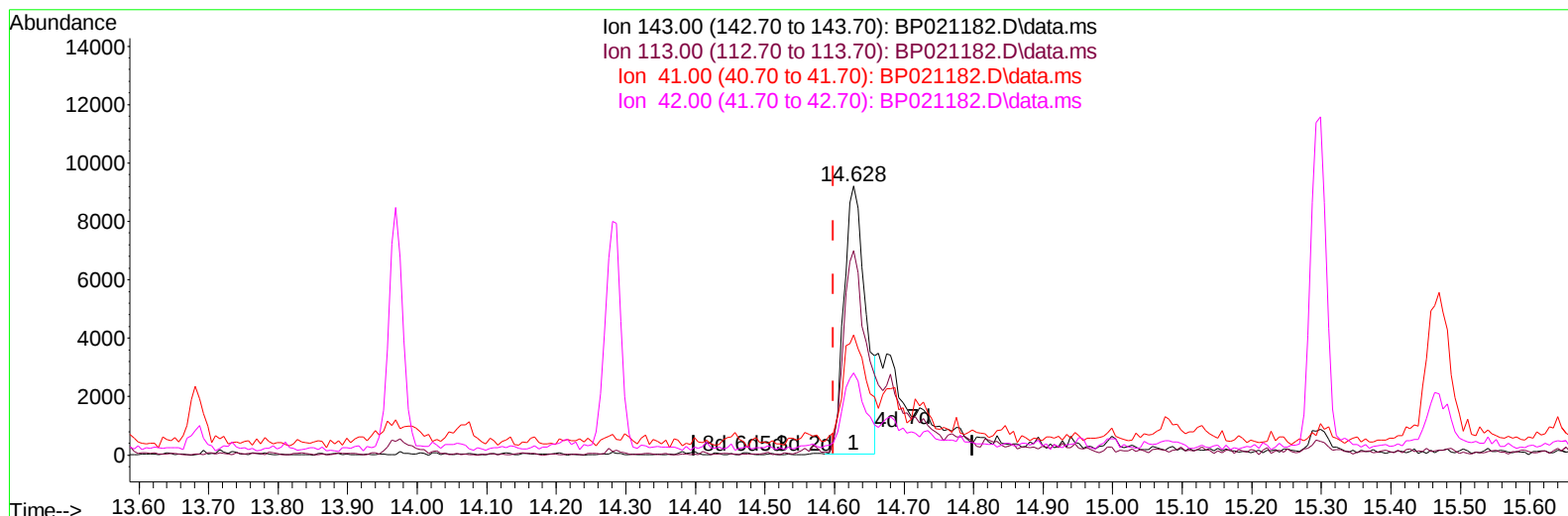
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
Data File : BP021182.D
Acq On : 26 Jul 2024 23:56
Operator : MA/JU
Sample : P3300-20
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G96

Manual IntegrationsAPPROVED

Quant Time: Jul 27 00:24:49 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 07/27/2024
Supervised By :mohammad ahmed 07/29/2024



TIC: BP021182.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.029) 4.51 ng/u1

response 20029

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	76.08
41.00	37.90	44.74
42.00	27.10	30.43

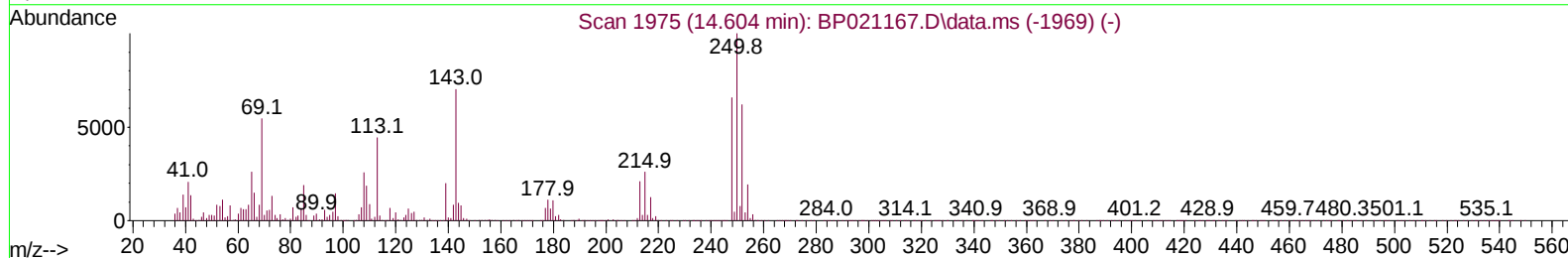
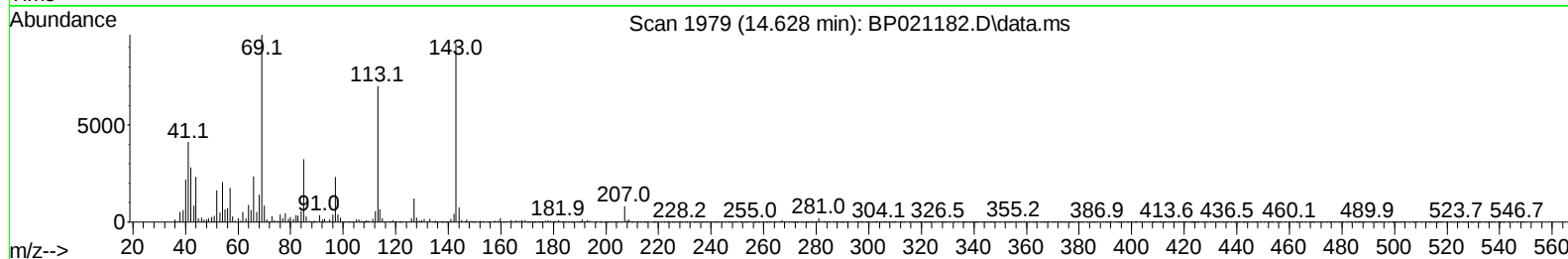
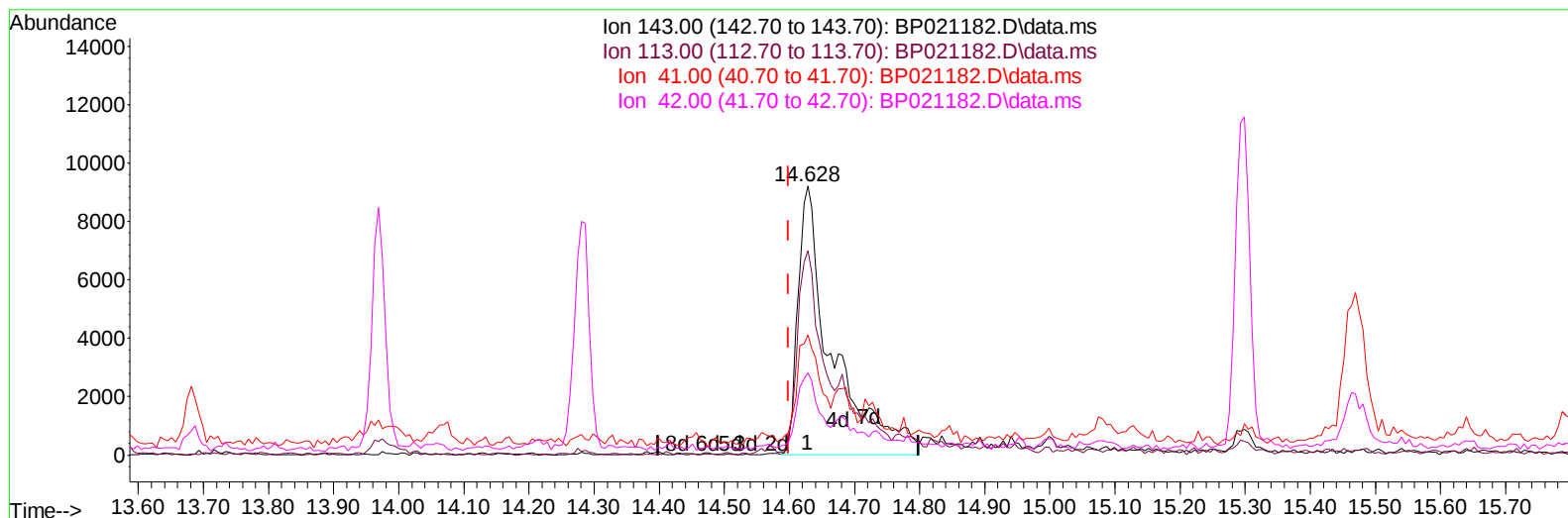
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TIC: BP021182.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.029) 7.50 ng/ul m

response 33304

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	76.08
41.00	37.90	44.74
42.00	27.10	30.43

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
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 Operator : MA/JU
 Sample : P3300-20
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E1G96

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/27/2024
 Supervised By :mohammad ahmed 07/29/2024

Quant Time: Jul 27 00:27:05 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	171525	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	674566	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.287	164	429450	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.092	188	931505	20.000	ng/ul	-0.01
79) Chrysene-d12	21.545	240	931192	20.000	ng/ul	-0.01
88) Perylene-d12	24.863	264	1096816	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.170	96	17203	4.564	ng/uL	0.00
4) Pyridine-d5	3.605	84	93011	8.523	ng/ul	0.00
7) Phenol-d5	6.875	99	105274	7.528	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	226329	26.707	ng/ul	0.00
11) 2-Chlorophenol-d4	7.199	132	284473	24.690	ng/ul	0.00
15) 4-Methylphenol-d8	8.375	113	193859	16.691	ng/ul	0.00
21) Nitrobenzene-d5	8.799	128	144000	27.484	ng/ul	0.00
24) 2-Nitrophenol-d4	9.516	143	166888	27.829	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	283328	26.128	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	336244	23.407	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	909708	29.140	ng/ul	-0.02
49) Acenaphthylene-d8	13.969	160	997060	28.443	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.628	143	33304m	7.498	ng/ul	0.03
60) Fluorene-d10	15.298	176	761898	29.748	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	135242	23.996	ng/ul	0.00
73) Anthracene-d10	17.198	188	1259057	30.432	ng/ul	0.00
81) Pyrene-d10	19.586	212	1608676	27.943	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.633	264	1795821	33.198	ng/ul	-0.02

Target Compounds Qvalue

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

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