

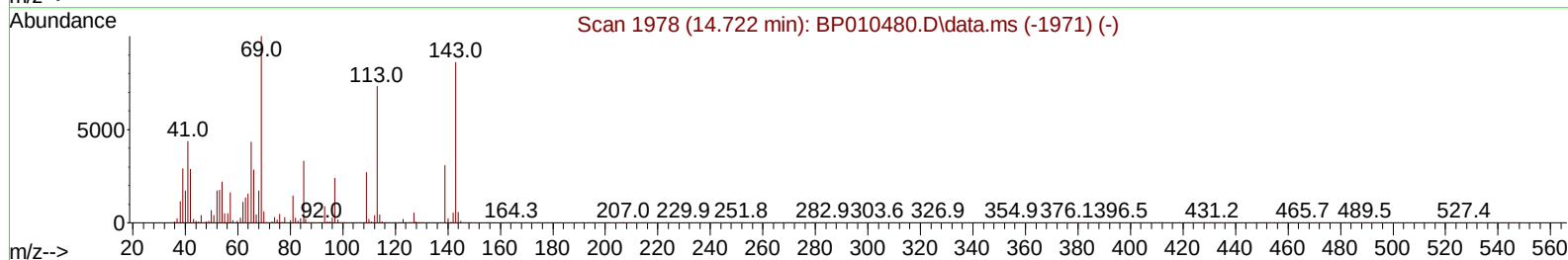
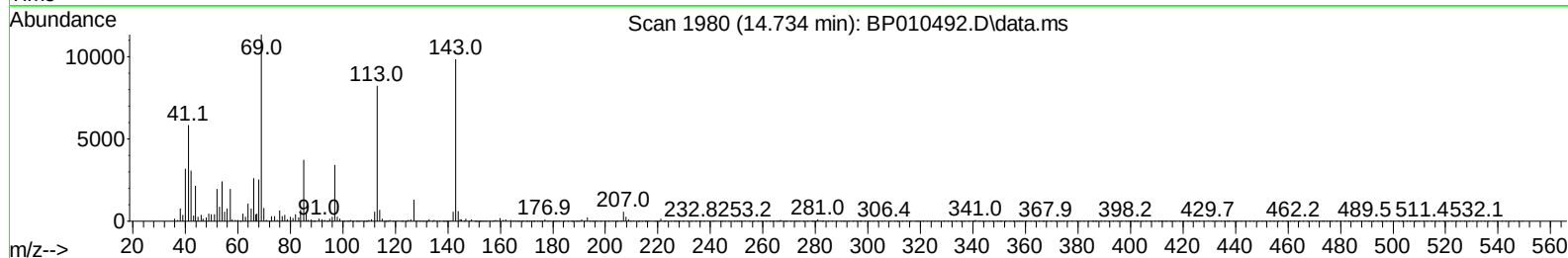
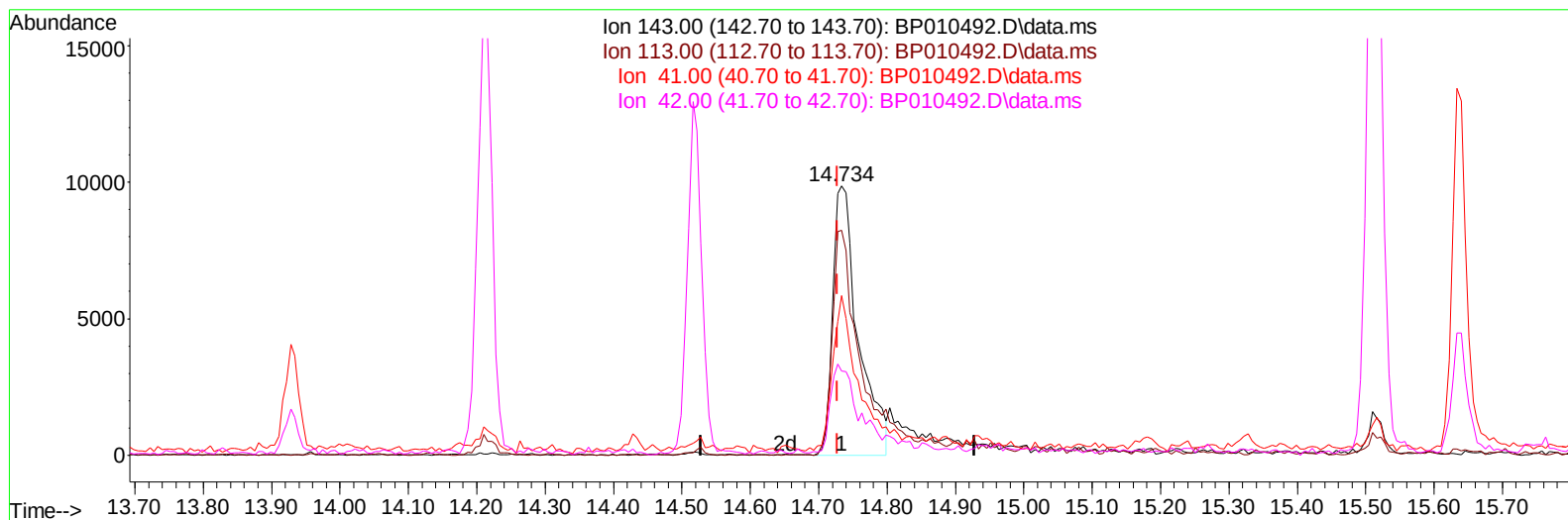
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
Data File : BP010492.D
Acq On : 23 May 2022 23:58
Operator : CG/JU
Sample : N2918-24
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/24/2022
Supervised By :Yogesh Patel 05/26/2022

Quant Time: May 24 00:52:58 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration



TIC: BP010492.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 4.27 ng/ul

response 25592

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.55#
41.00	23.20	59.22#
42.00	18.00	31.34#

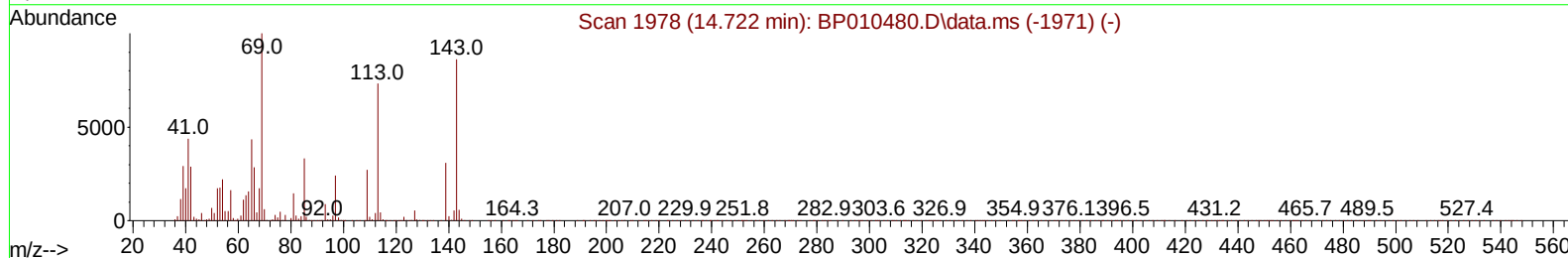
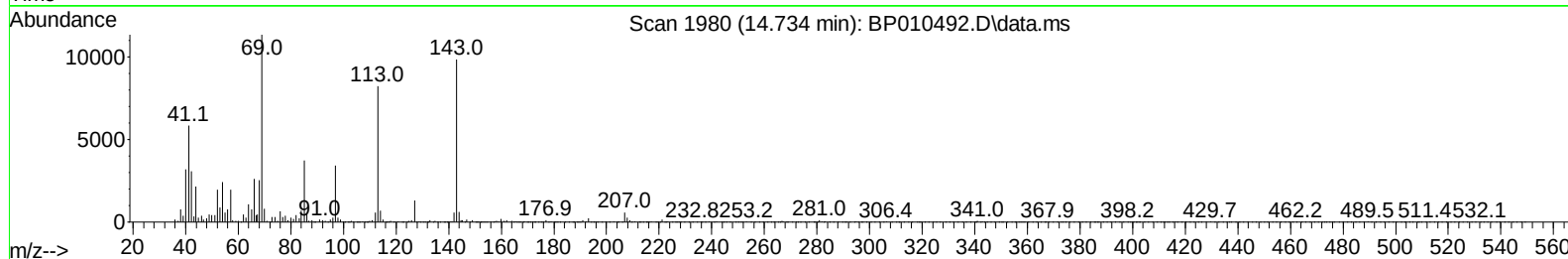
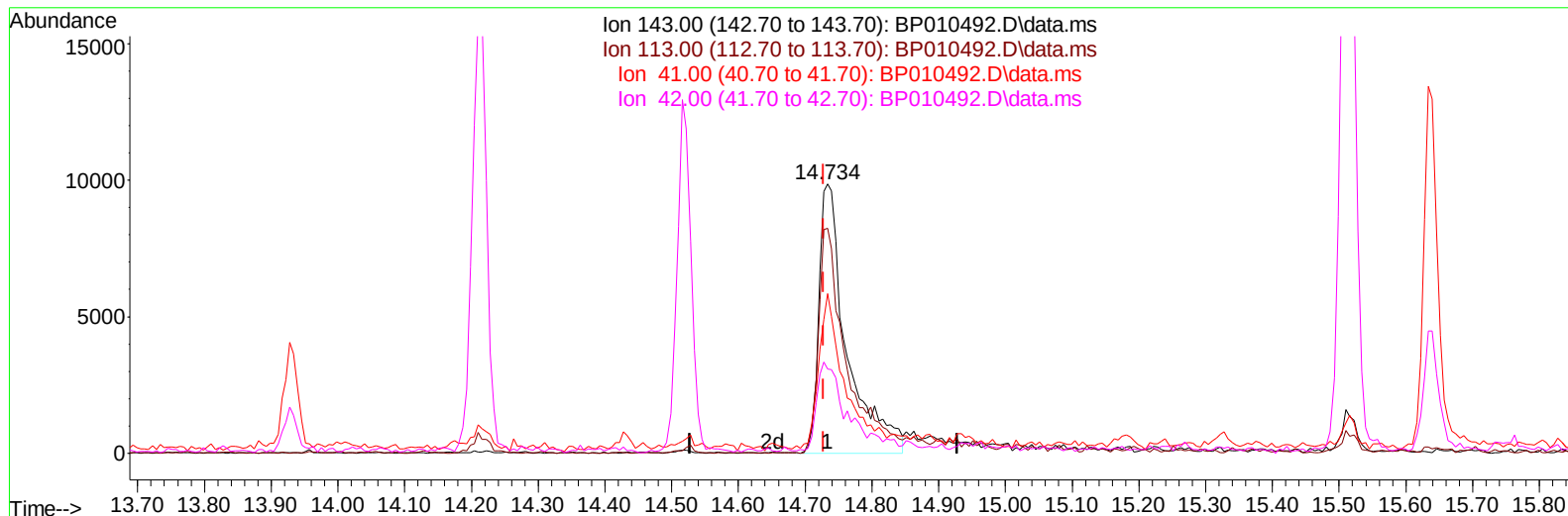
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TIC: BP010492.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 4.78 ng/ul m

response 28636

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.55#
41.00	23.20	59.22#
42.00	18.00	31.34#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	148358	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.687	136	668919	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.522	164	456876	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.269	188	974764	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.351	240	870676	20.000	ng/ul	#-0.01
88) Perylene-d12	23.774	264	691714	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	17058	4.901	ng/uL	0.00
4) Pyridine-d5	3.770	84	61630	6.236	ng/ul	0.00
7) Phenol-d5	7.058	99	86774	6.969	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	304736	36.425	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	257794	27.056	ng/ul	-0.01
15) 4-Methylphenol-d8	8.593	113	184710	17.455	ng/ul	-0.02
21) Nitrobenzene-d5	9.046	128	196969	38.047	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.769	143	202557	36.920	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	331302	31.936	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	462561	29.988	ng/ul	-0.01
46) Dimethylphthalate-d6	13.928	166	1425694	42.354	ng/ul	-0.01
49) Acenaphthylene-d8	14.216	160	1605455	41.349	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	28636m	4.780	ng/ul	0.00
60) Fluorene-d10	15.516	176	1239913	42.340	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	177455	31.128	ng/ul	0.00
73) Anthracene-d10	17.375	188	2000450	45.322	ng/ul	0.00
81) Pyrene-d10	19.604	212	2397120	51.842	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.622	264	1651457	47.546	ng/ul	-0.01

Target Compounds Qvalue

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

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