

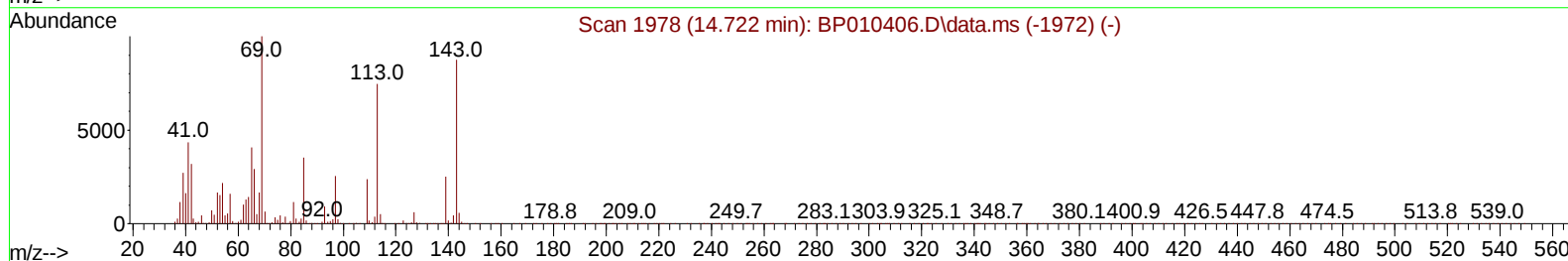
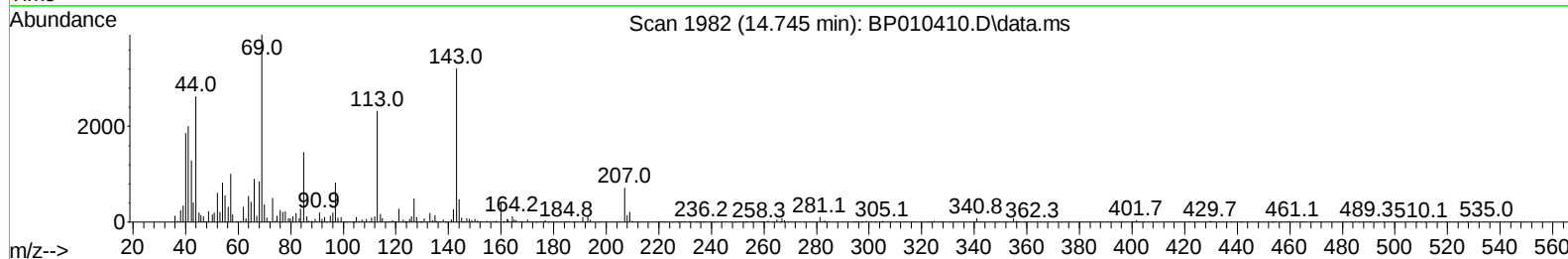
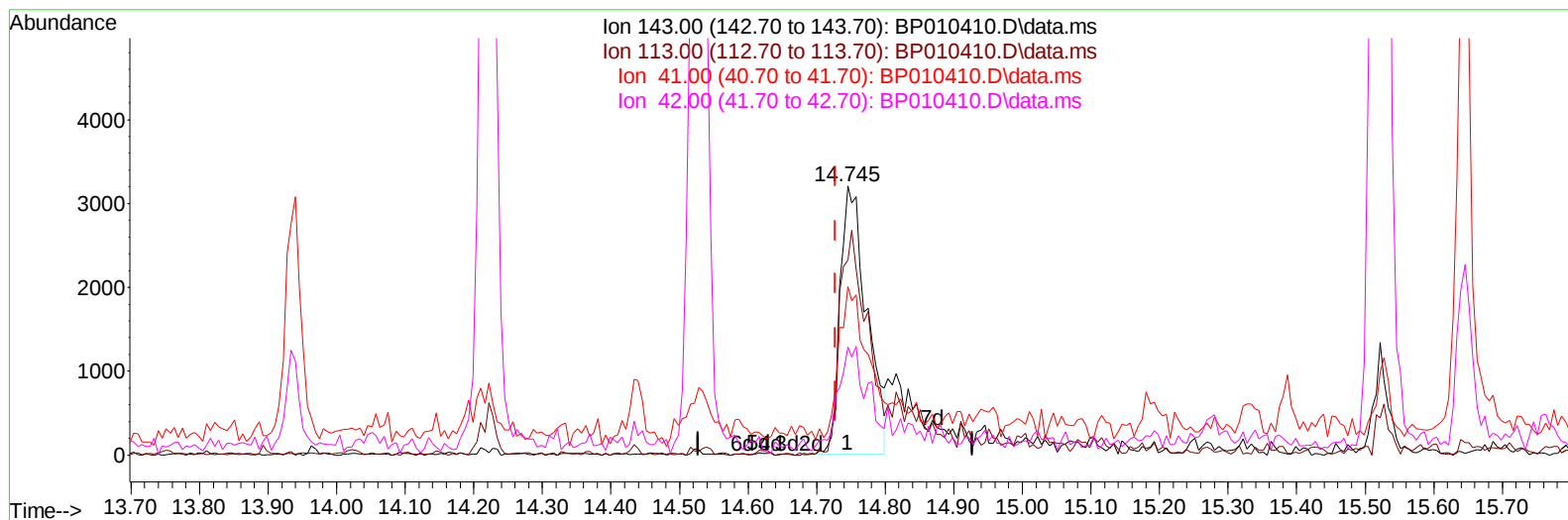
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051922\
Data File : BP010410.D
Acq On : 19 May 2022 19:30
Operator : CG/JU
Sample : N2918-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTE4

Manual IntegrationsAPPROVED

Quant Time: May 20 00:46:02 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

Reviewed By :Jagrut Upadhyay 05/20/2022
Supervised By :Yogesh Patel 05/24/2022



TIC: BP010410.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (+ 0.018) 1.50 ng/u1

response 8667

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.38#
41.00	23.20	62.59#
42.00	18.00	40.15#

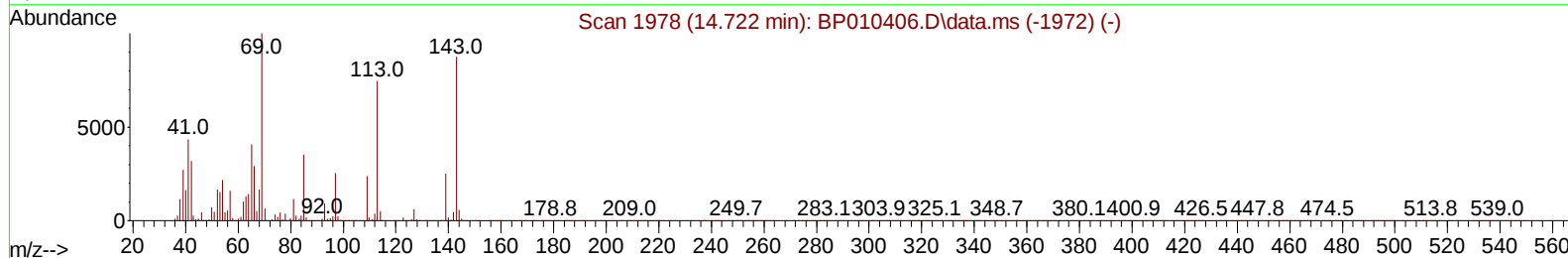
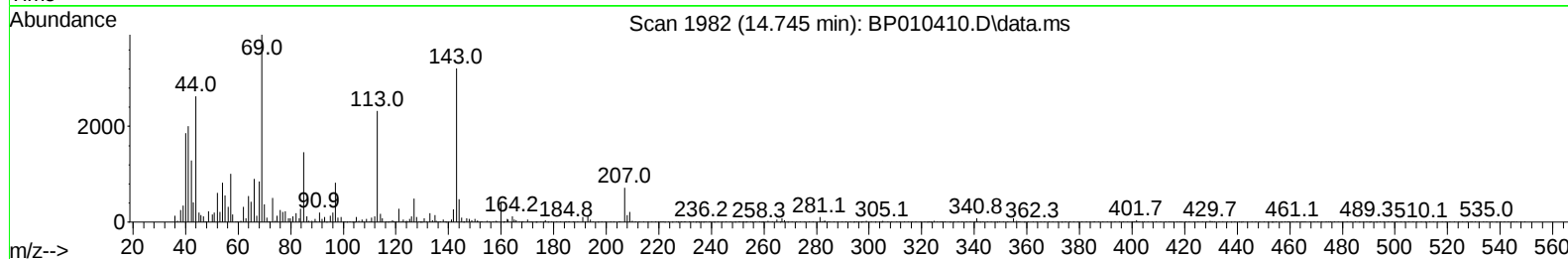
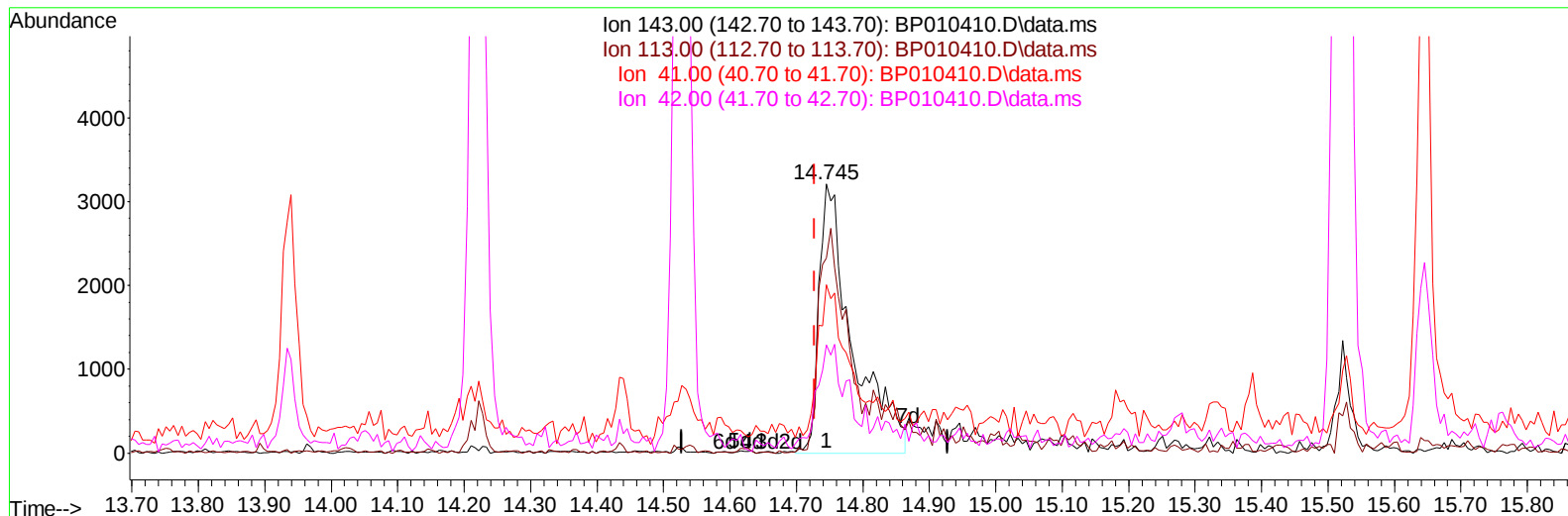
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TIC: BP010410.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (+ 0.018) 1.95 ng/ul m

response 11221

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.38#
41.00	23.20	62.59#
42.00	18.00	40.15#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	150635	20.000	ng/ul	0.00
20) Naphthalene-d8	10.699	136	640578	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	439868	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	958107	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.369	240	918113	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	808842	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	15933	4.508	ng/uL	0.00
4) Pyridine-d5	3.764	84	77449	7.718	ng/ul	0.00
7) Phenol-d5	7.063	99	42844	3.389	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	231456	27.248	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	133406	13.790	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113	100710	9.373	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	145975	29.444	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	91661	17.446	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	162673	16.375	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	343272	23.239	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	1043007	32.184	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1113991	29.800	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	11221m	1.945	ng/ul	0.02
60) Fluorene-d10	15.522	176	909716	32.266	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	83627	14.924	ng/ul	0.00
73) Anthracene-d10	17.381	188	1626552	37.491	ng/ul	0.00
81) Pyrene-d10	19.616	212	1952142	40.037	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1539315	37.900	ng/ul	0.00

Target Compounds Qvalue

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

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