

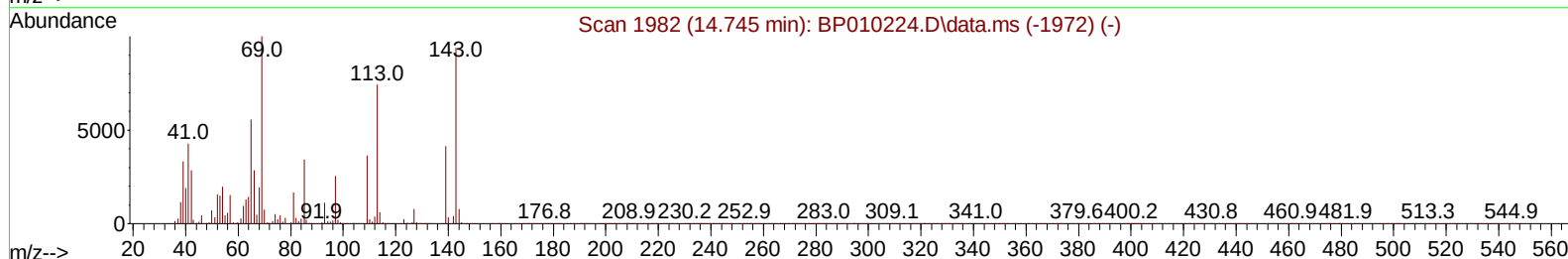
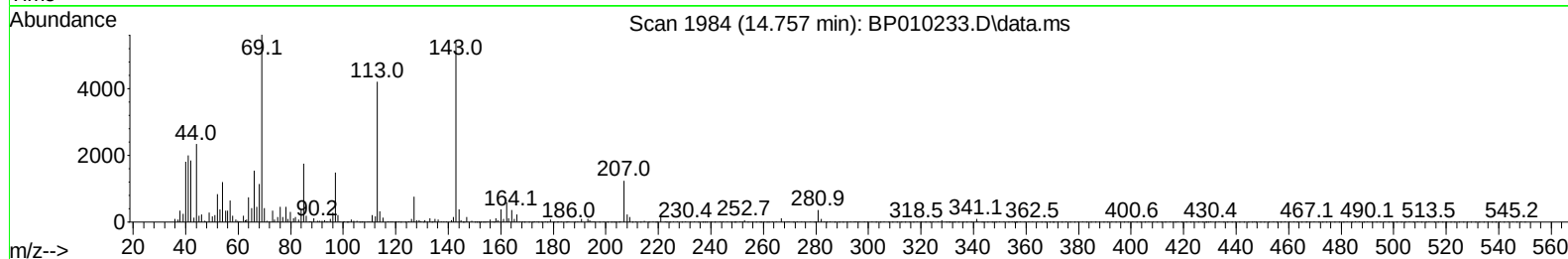
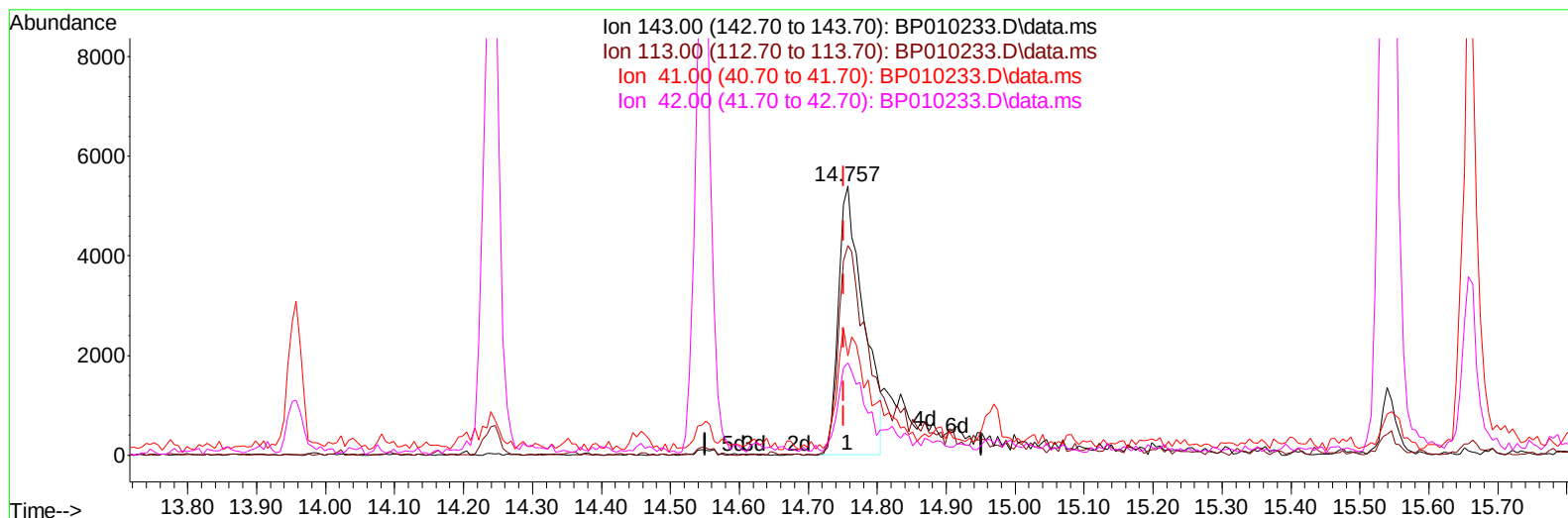
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050422\
Data File : BP010233.D
Acq On : 04 May 2022 20:48
Operator : CG/JU
Sample : N2668-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXL72

Manual IntegrationsAPPROVED

Quant Time: May 05 02:41:35 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 00:58:32 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/05/2022
Supervised By :mohammad ahmed 05/09/2022



TIC: BP010233.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (+ 0.006) 1.91 ng/ul

response 13607

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	78.01#
41.00	23.20	36.93#
42.00	18.00	34.25#

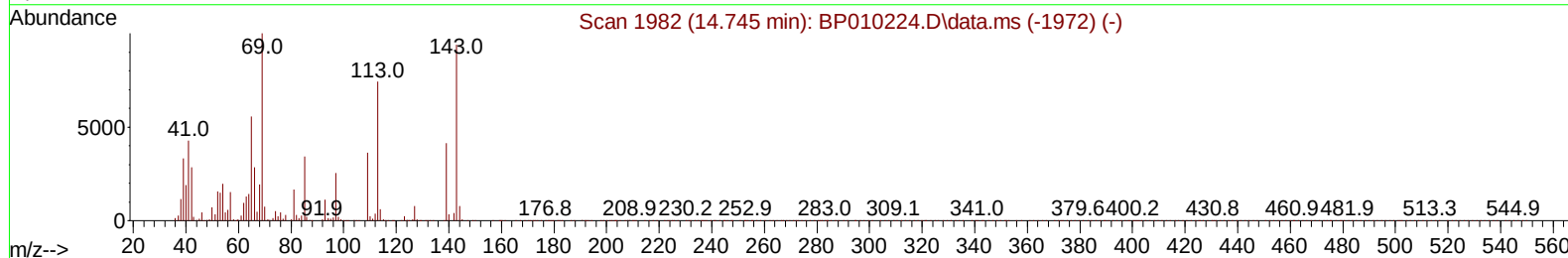
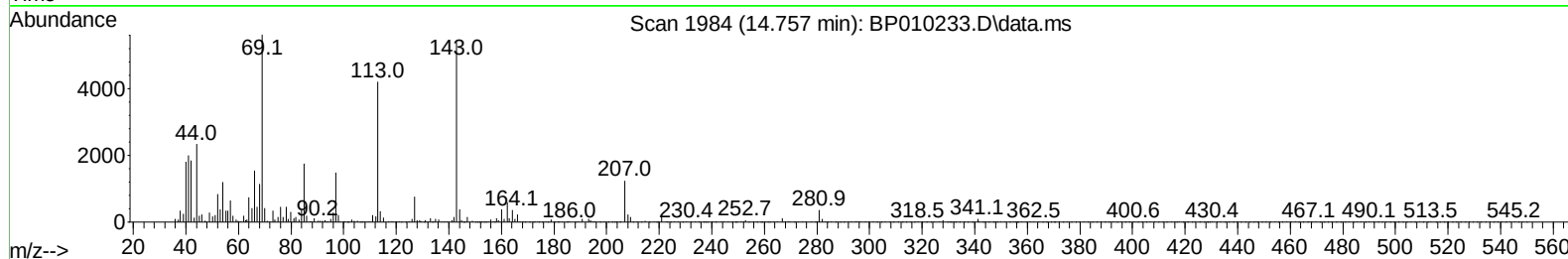
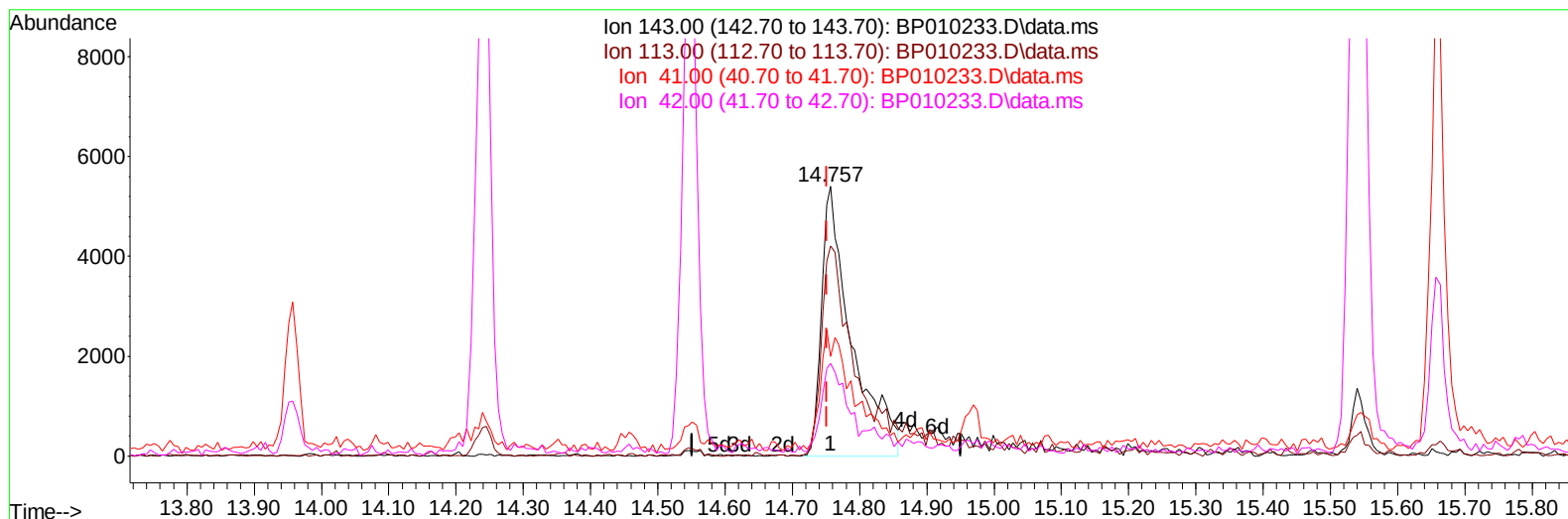
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(54) 4-Nitrophenol-d4 (S)

14.757min (+ 0.006) 2.35 ng/ul m

response 16770

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	78.01#
41.00	23.20	36.93#
42.00	18.00	34.25#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	176011	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	769824	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	498589	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	942207	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.380	240	689738	20.000	ng/ul	#-0.01
88) Perylene-d12	23.827	264	668103	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	16189	4.011	ng/uL	0.00
4) Pyridine-d5	3.781	84	62913	5.541	ng/ul	0.00
7) Phenol-d5	7.081	99	63899	3.991	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	274820	28.607	ng/ul	0.00
11) 2-Chlorophenol-d4	7.446	132	229741	19.376	ng/ul	-0.01
15) 4-Methylphenol-d8	8.622	113	135667	10.228	ng/ul	-0.01
21) Nitrobenzene-d5	9.075	128	168818	27.572	ng/ul	0.00
24) 2-Nitrophenol-d4	9.799	143	170099	25.438	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	269176	20.678	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.857	131	353354	18.931	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1137189	29.321	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1252033	26.673	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.757	143	16770m	2.349	ng/ul	0.00
60) Fluorene-d10	15.540	176	993697	29.785	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.663	200	145496	24.421	ng/ul	0.00
73) Anthracene-d10	17.398	188	1553733	34.086	ng/ul	-0.01
81) Pyrene-d10	19.633	212	1581237	40.311	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	1261784	36.024	ng/ul	-0.01

Target Compounds Qvalue

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

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