

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072522\
Data File : BP011092.D
Acq On : 25 Jul 2022 15:01
Operator : CG/JU
Sample : N3801-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

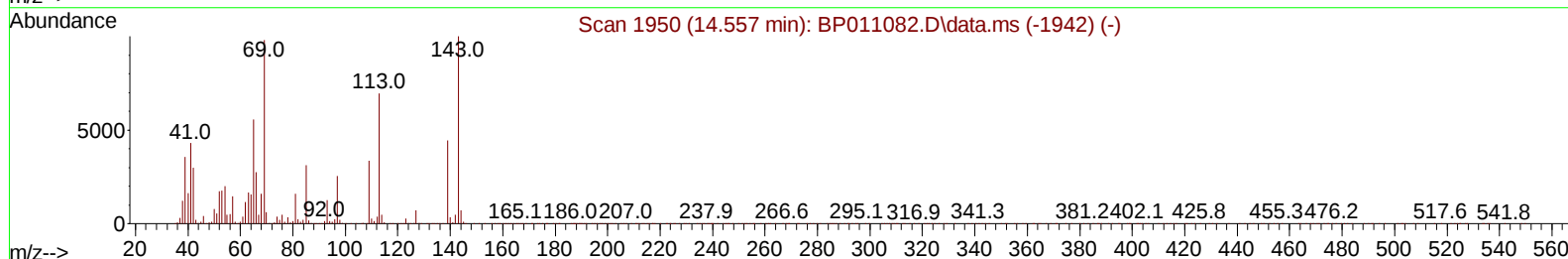
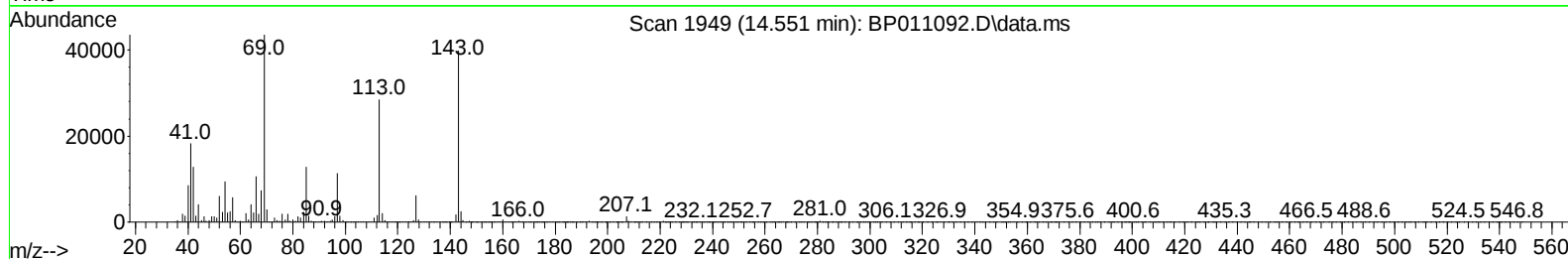
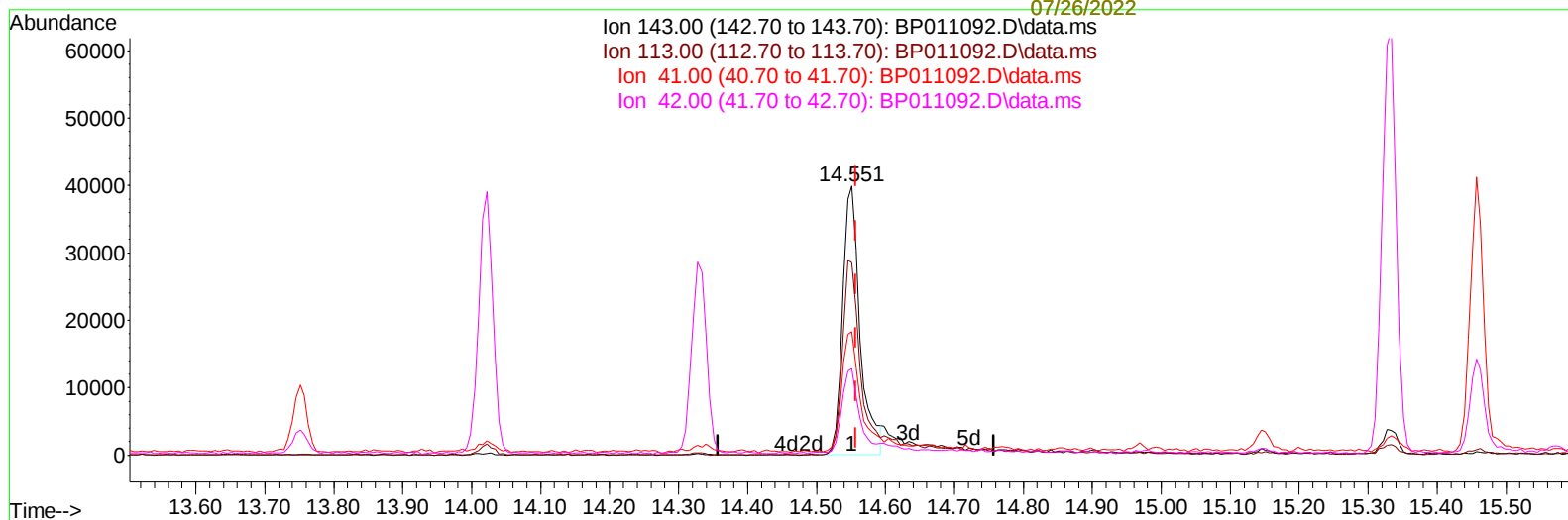
Instrument :
BNA_P
ClientSampleId :
GBHX8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut
Upadhyay
07/26/2022

Supervised By :mohammad
ahmed
07/26/2022

Quant Time: Jul 25 23:21:08 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 19 03:25:54 2022
Response via : Initial Calibration



TIC: BP011092.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.551min (-0.006) 4.15 ng/u1

response 70254

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	71.47#
41.00	59.00	45.87#
42.00	41.10	32.29#

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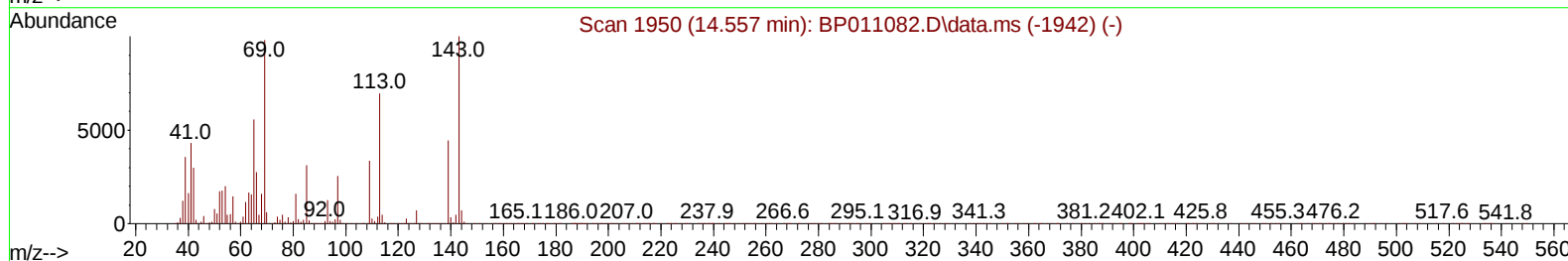
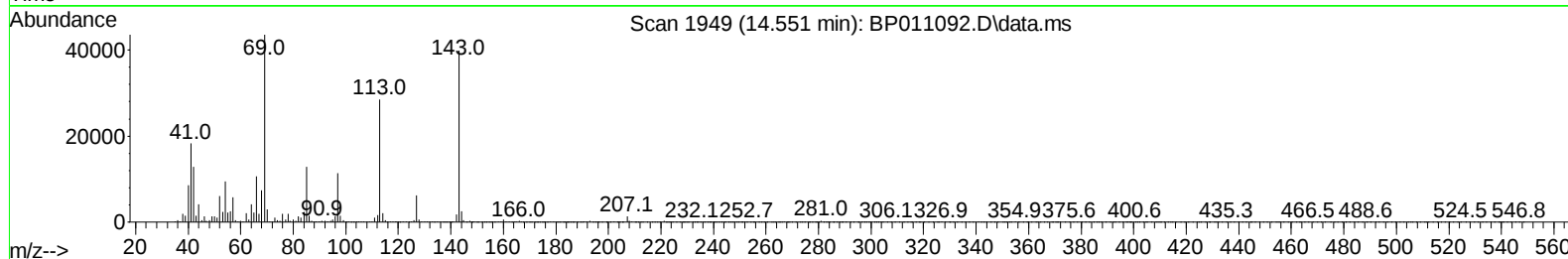
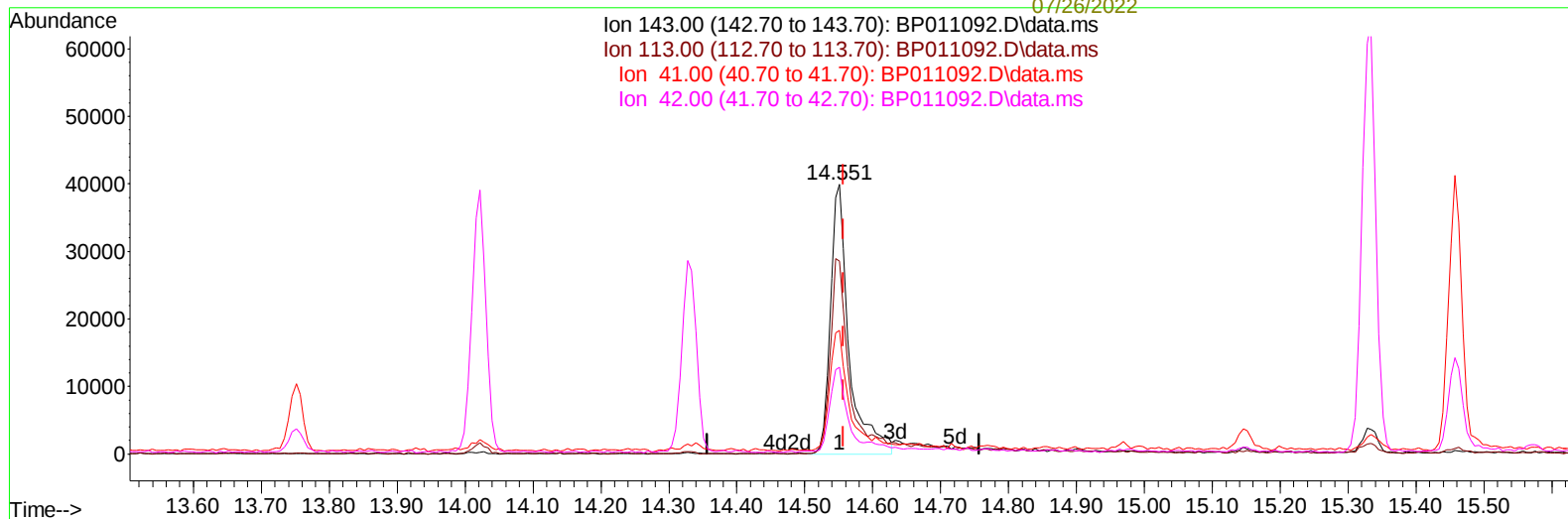
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TIC: BP011092.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.551min (-0.006) 4.51 ng/ul m

response 76273

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	71.47#
41.00	59.00	45.87#
42.00	41.10	32.29#

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.663	152		430066	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.463	136		1991978	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.334	164		1223594	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188		2508127	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.216	240		2104499	20.000	ng/ul	#-0.01
88) Perylene-d12	23.563	264		1791502	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.164	96		52158	4.372	ng/uL	-0.02
4) Pyridine-d5	3.575	84		338730	10.985	ng/ul	-0.02
7) Phenol-d5	6.846	99		251345	7.120	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.010	67		912691	40.205	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.199	132		818736	28.702	ng/ul	-0.01
15) 4-Methylphenol-d8	8.381	113		530657	18.776	ng/ul	-0.02
21) Nitrobenzene-d5	8.828	128		587751	40.488	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.551	143		585471	40.759	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165		885735	33.718	ng/ul	-0.02
31) 4-Chloroaniline-d4	10.610	131		1422394	32.845	ng/ul	0.00
46) Dimethylphthalate-d6	13.751	166		3637365	43.494	ng/ul	-0.01
49) Acenaphthylene-d8	14.022	160		4220658	42.775	ng/ul	0.00
54) 4-Nitrophenol-d4	14.551	143		76273m	4.506	ng/ul	0.00
60) Fluorene-d10	15.334	176		3004859	42.272	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200		497160	40.058	ng/ul	0.00
73) Anthracene-d10	17.198	188		4666017	43.760	ng/ul	0.00
81) Pyrene-d10	19.451	212		5188982	46.947	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.415	264		3794588	44.100	ng/ul	-0.01

Target Compounds Qvalue

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

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