

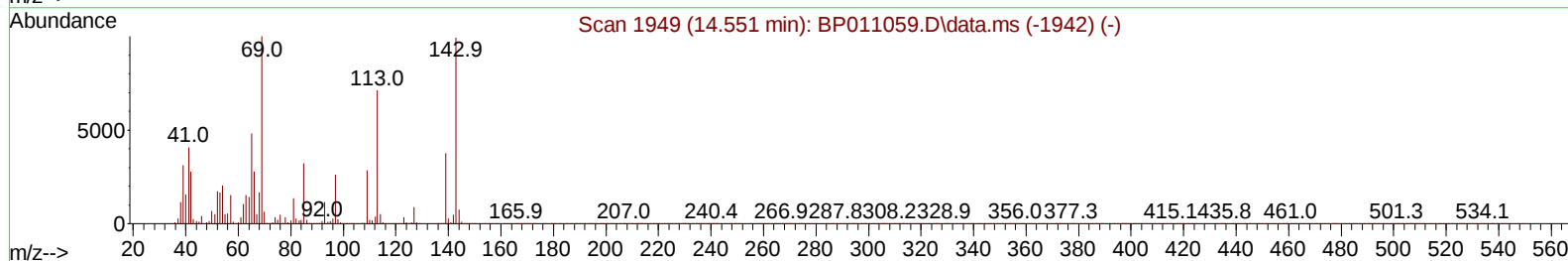
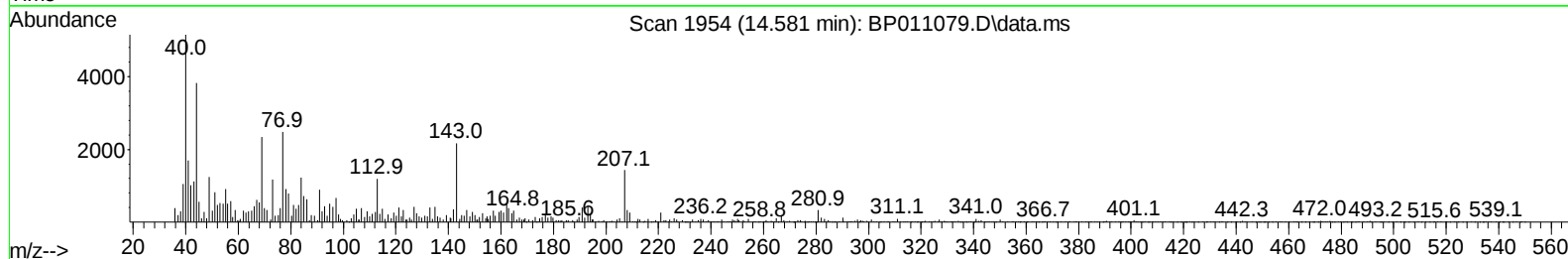
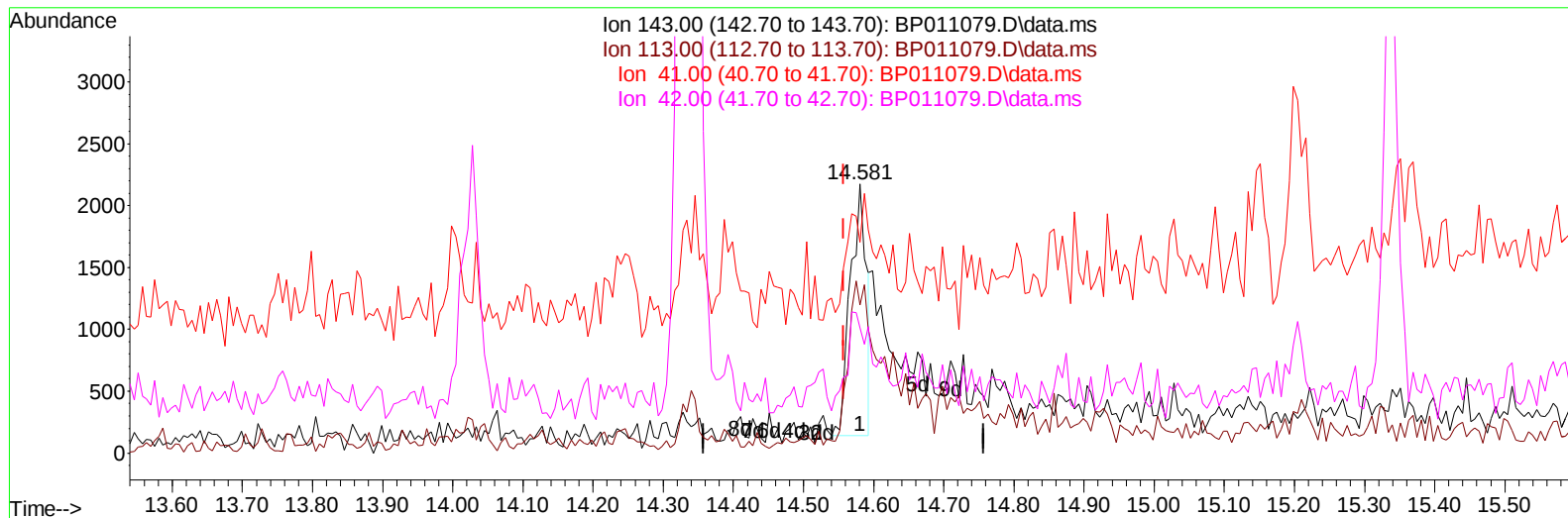
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072122\
Data File : BP011079.D
Acq On : 22 Jul 2022 01:24
Operator : CG/JU
Sample : N3709-14 10X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0B28

Manual IntegrationsAPPROVED

Quant Time: Jul 22 02:30:13 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

Reviewed By :Jagrut Upadhyay 07/22/2022
Supervised By :mohammad ahmed 07/22/2022



TIC: BP011079.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.581min (+ 0.024) 0.19 ng/u1

response 3278

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	54.99#
41.00	59.00	78.30#
42.00	41.10	46.57

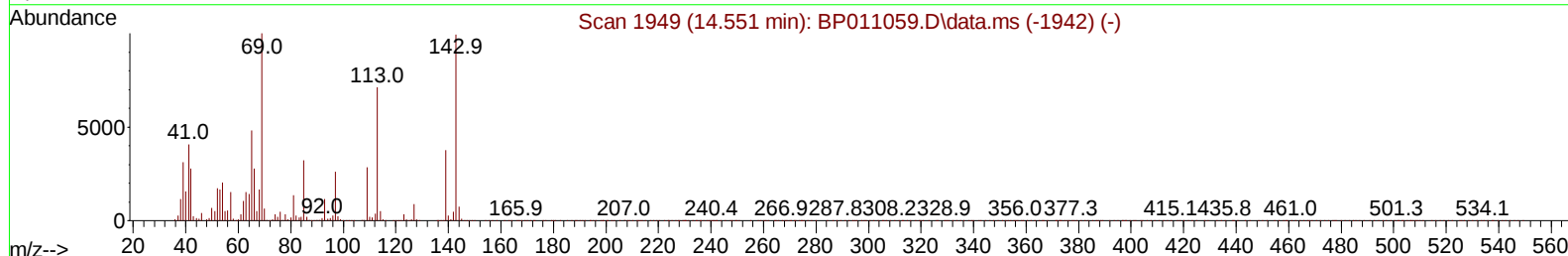
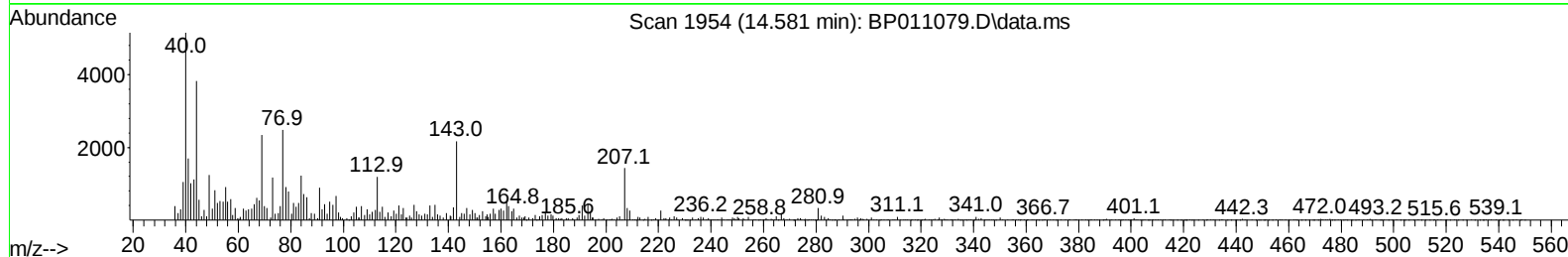
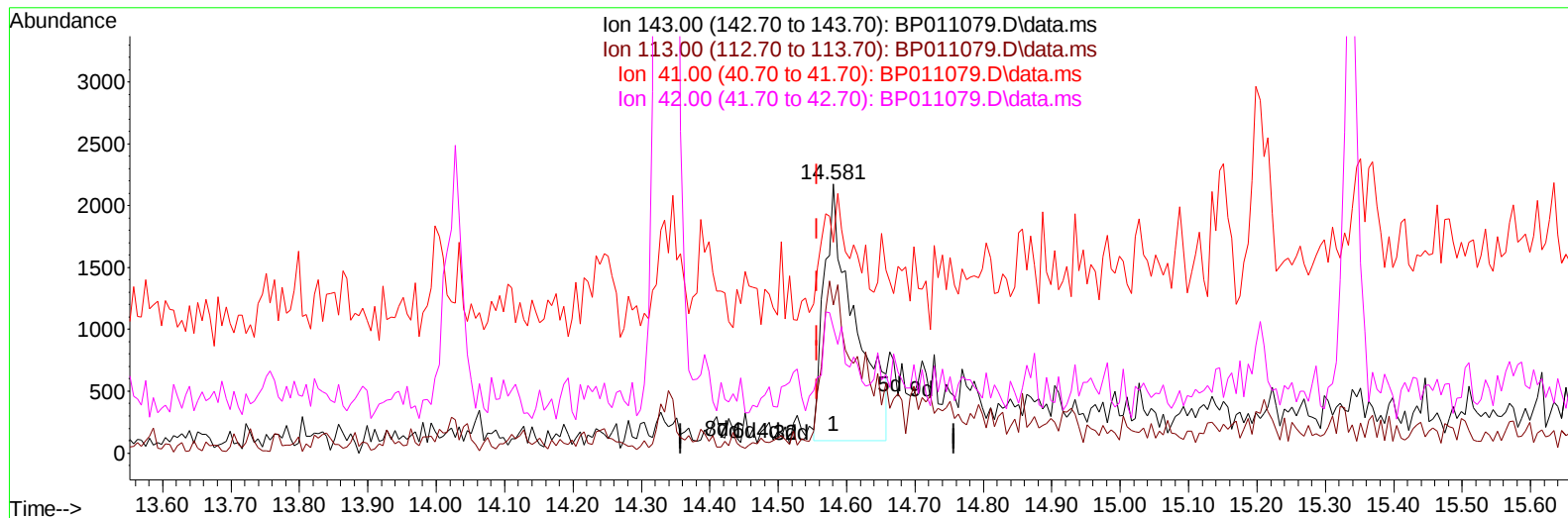
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TIC: BP011079.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.581min (+ 0.024) 0.37 ng/ul m

response 6438

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	54.99#
41.00	59.00	78.30#
42.00	41.10	46.57

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	500773	20.000	ng/ul	0.00
20) Naphthalene-d8	10.469	136	2128200	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.340	164	1247838	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	2510534	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.227	240	1889316	20.000	ng/ul	# 0.00
88) Perylene-d12	23.586	264	1968509	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	3161	0.228	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.858	99	78639	1.913	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	53846	2.037	ng/ul	0.00
11) 2-Chlorophenol-d4	7.211	132	66935	2.015	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	64868	1.971	ng/ul	0.00
21) Nitrobenzene-d5	8.834	128	32126	2.071	ng/ul	0.00
24) 2-Nitrophenol-d4	9.558	143	24621	1.604	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	52839	1.883	ng/ul	0.00
31) 4-Chloroaniline-d4	10.616	131	24401	0.527	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166	177035	2.076	ng/ul	0.00
49) Acenaphthylene-d8	14.028	160	245350	2.438	ng/ul	0.00
54) 4-Nitrophenol-d4	14.581	143	6438m	0.373	ng/ul	0.02
60) Fluorene-d10	15.340	176	180409	2.489	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	2146	0.173	ng/ul	0.00
73) Anthracene-d10	17.204	188	277475	2.600	ng/ul	0.00
81) Pyrene-d10	19.457	212	291943	2.942	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.427	264	237759	2.515	ng/ul	0.00

Target Compounds Qvalue

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

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