

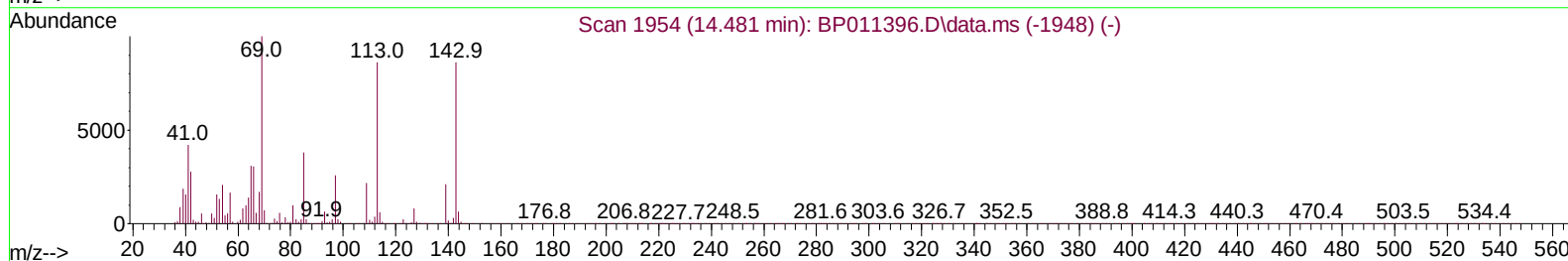
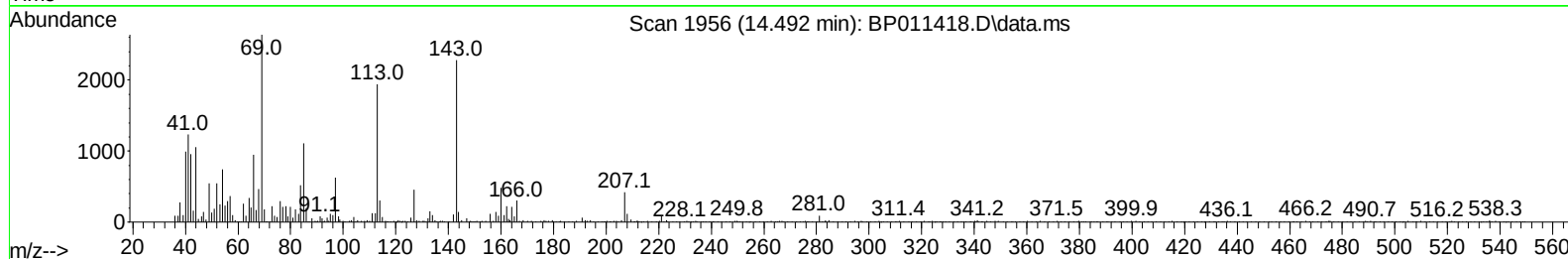
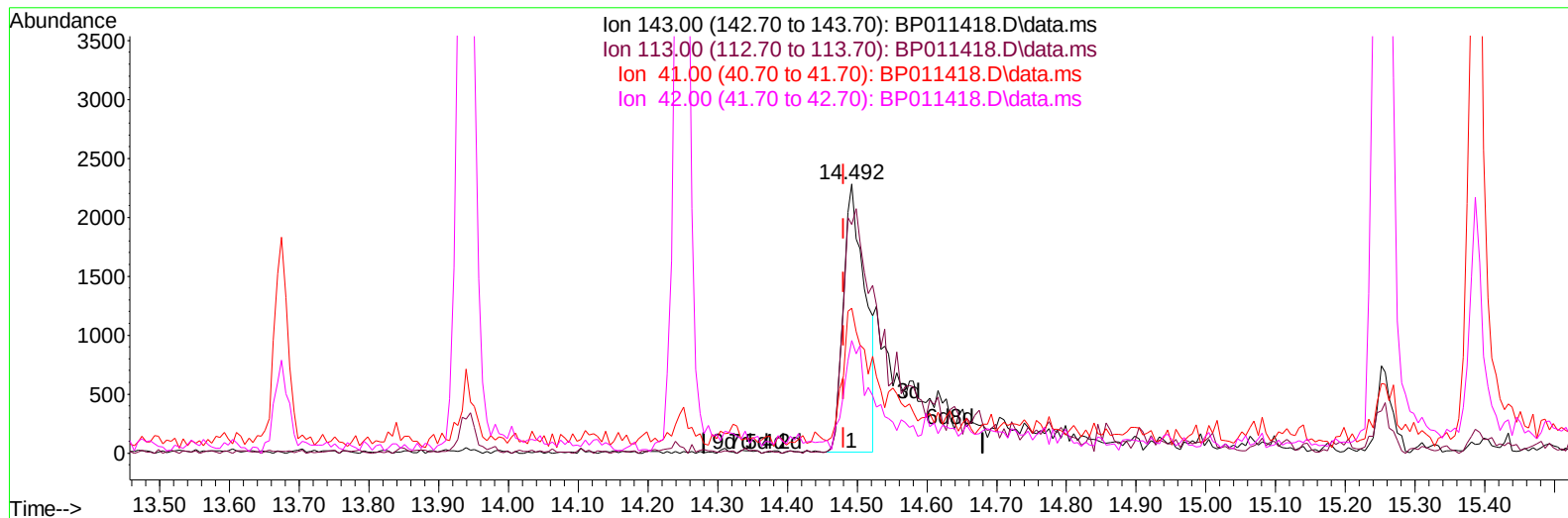
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011418.D
 Acq On : 12 Aug 2022 11:12
 Operator : CG/JU
 Sample : N4129-06
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGBS9

Manual IntegrationsAPPROVED

Quant Time: Aug 12 12:21:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 11 23:05:53 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/12/2022
 Supervised By :Sohil Jodhani 08/13/2022



TIC: BP011418.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.012) 2.21 ng/ul

response 4886

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	84.96
41.00	54.40	54.01
42.00	35.50	41.82

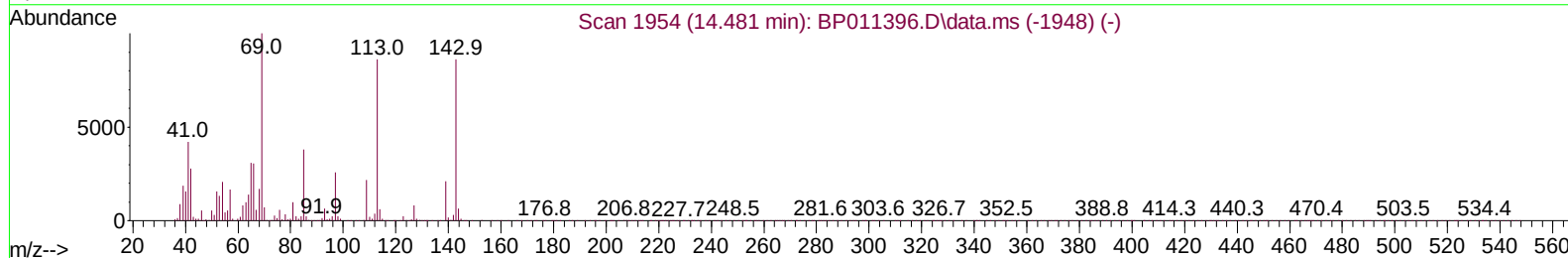
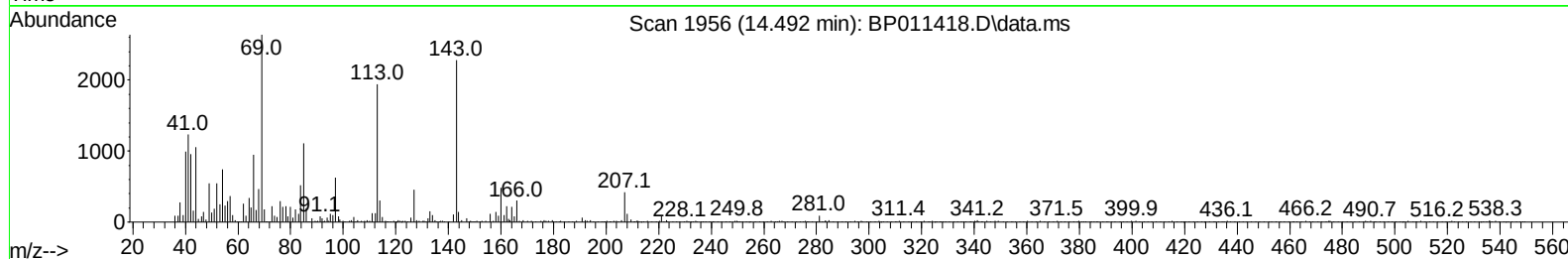
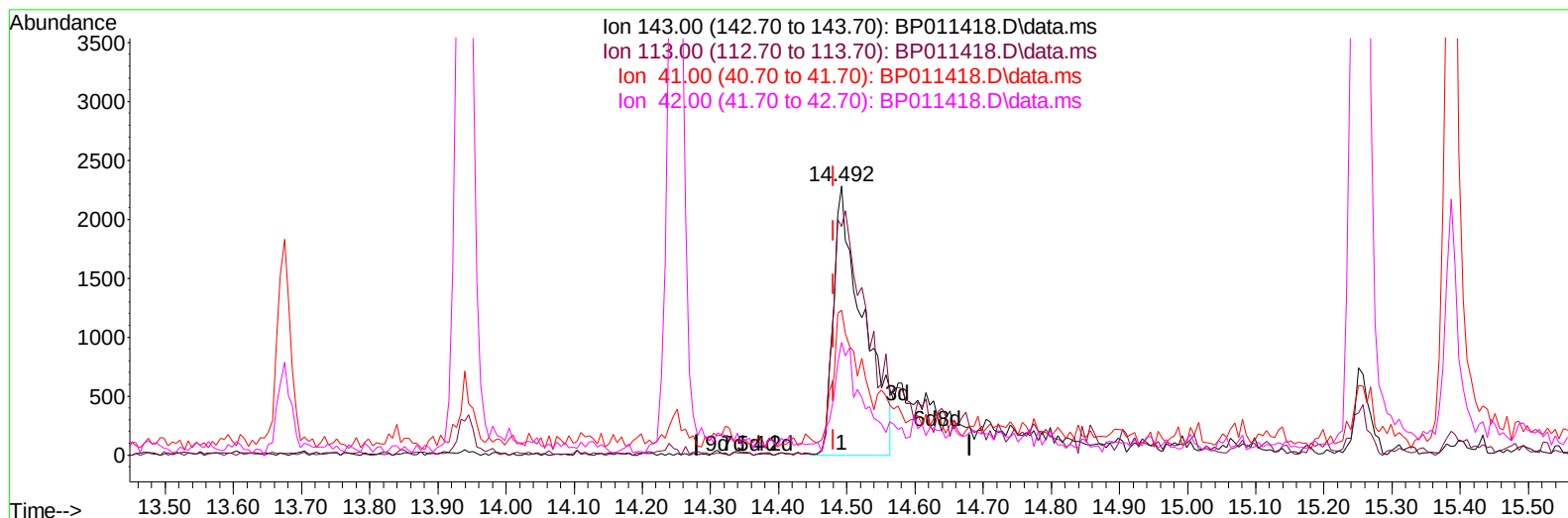
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TIC: BP011418.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (+ 0.012) 3.13 ng/ul m

response 6926

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	84.96
41.00	54.40	54.01
42.00	35.50	41.82

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	70118	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	302052	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	174805	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.022	188	361438	20.000	ng/ul	0.00
79) Chrysene-d12	21.157	240	359209	20.000	ng/ul	# 0.00
88) Perylene-d12	23.462	264	377045	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	12850	5.618	ng/uL	0.00
4) Pyridine-d5	3.493	84	62497	10.832	ng/ul	0.00
7) Phenol-d5	6.758	99	40809	6.577	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	146541	33.306	ng/ul	0.00
11) 2-Chlorophenol-d4	7.110	132	131420	26.992	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	81011	16.627	ng/ul	-0.01
21) Nitrobenzene-d5	8.740	128	83352	40.547	ng/ul	0.00
24) 2-Nitrophenol-d4	9.457	143	76429	40.609	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	122350	32.191	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	189508	28.520	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	509002	41.966	ng/ul	0.00
49) Acenaphthylene-d8	13.939	160	572663	39.223	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	6926m	3.126	ng/ul	0.01
60) Fluorene-d10	15.257	176	422270	40.710	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	58613	34.913	ng/ul	0.00
73) Anthracene-d10	17.122	188	731853	45.049	ng/ul	0.00
81) Pyrene-d10	19.386	212	857890	43.349	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.315	264	849194	44.660	ng/ul	0.00

Target Compounds Qvalue

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

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