

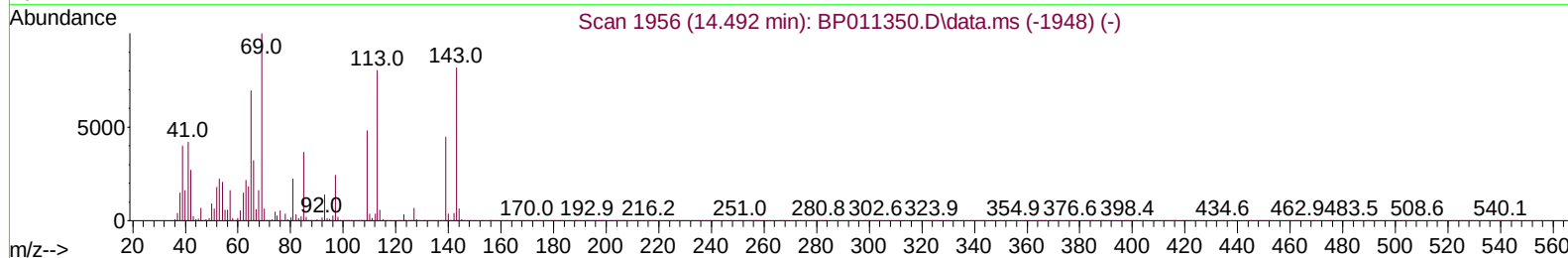
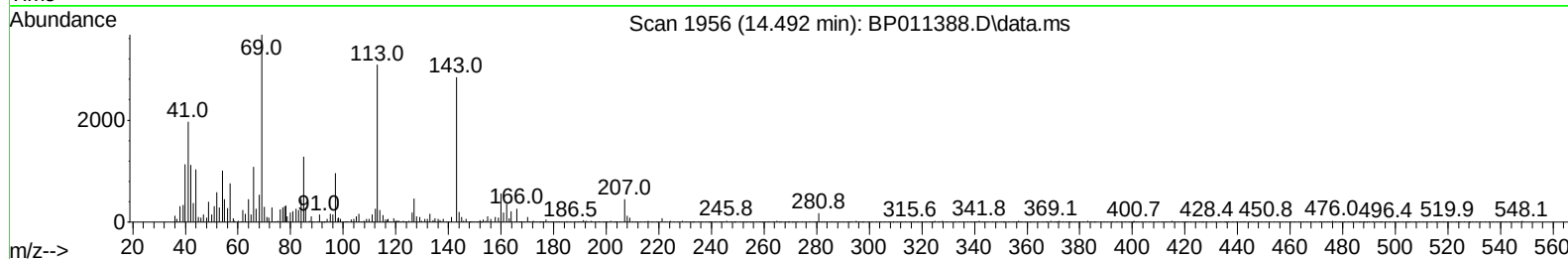
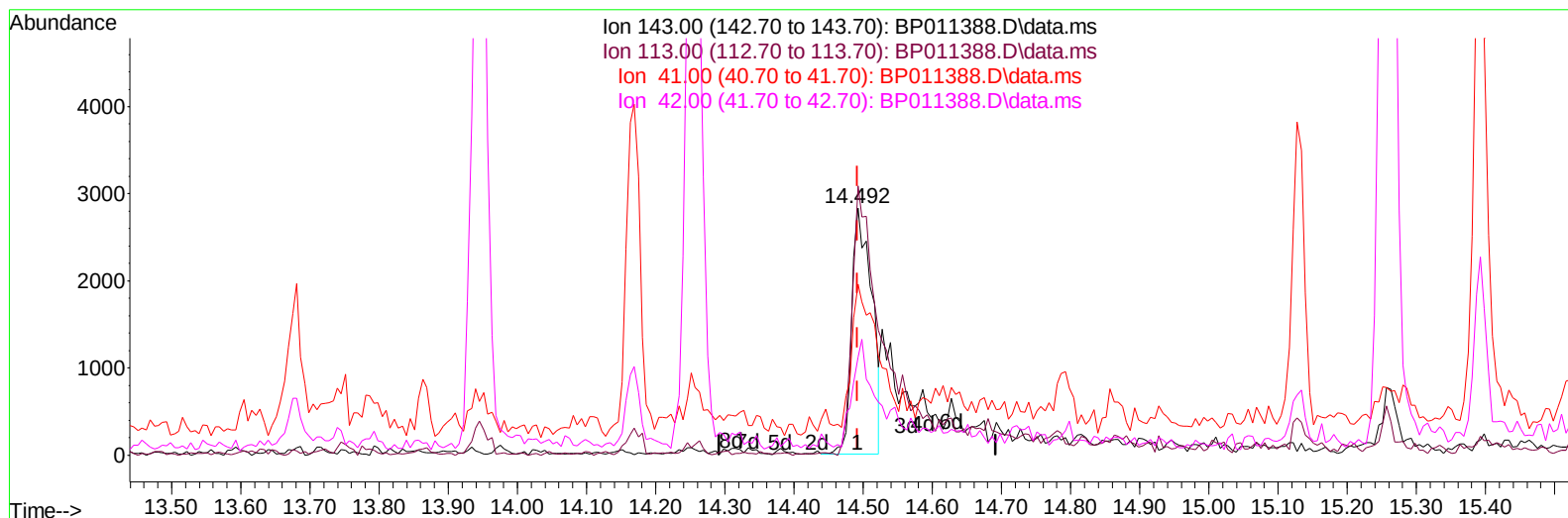
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
 Data File : BP011388.D
 Acq On : 11 Aug 2022 15:36
 Operator : CG/JU
 Sample : N4128-16
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGC10

Manual IntegrationsAPPROVED

Quant Time: Aug 11 16:20:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 08 18:47:19 2022
 Response via : Initial Calibration

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



TIC: BP011388.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (-0.000) 1.98 ng/u1

response 5622

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	108.85#
41.00	54.40	69.26#
42.00	35.50	39.55

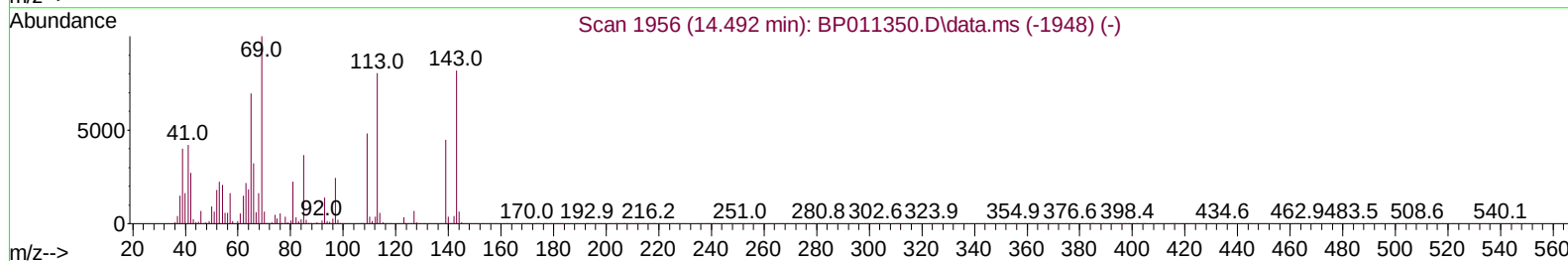
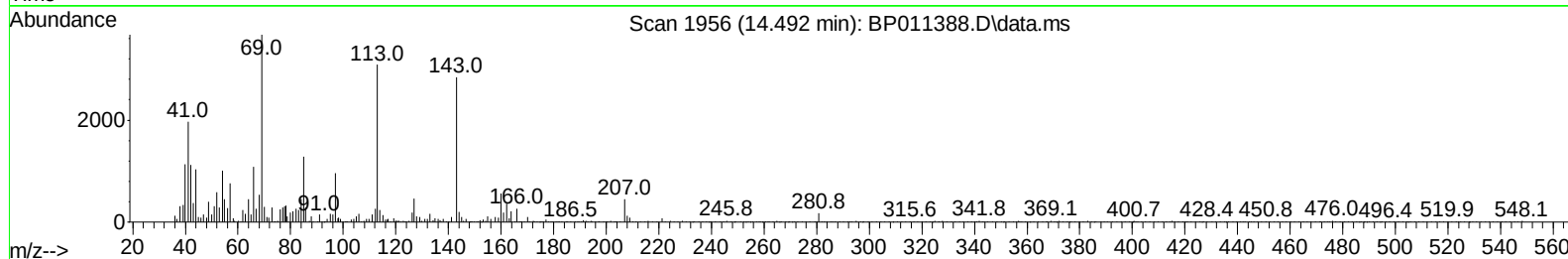
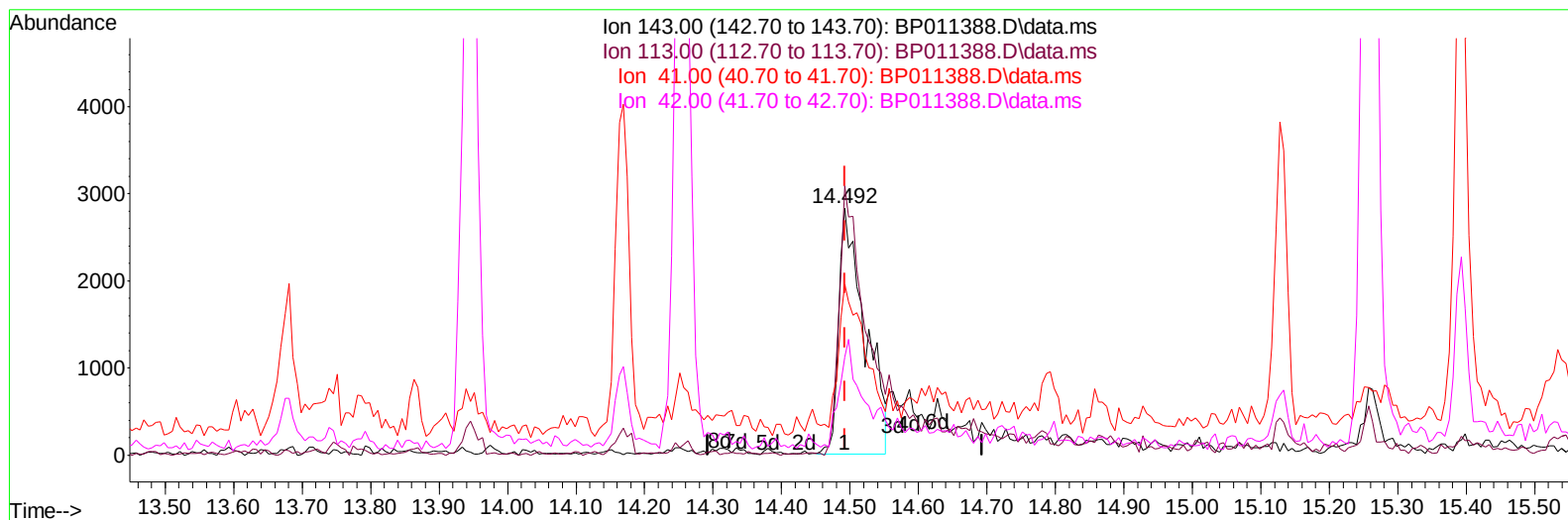
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TIC: BP011388.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (-0.000) 2.61 ng/ul m

response 7413

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	108.85#
41.00	54.40	69.26#
42.00	35.50	39.55

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	93816	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.369	136	389256	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.257	164	224394	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	476473	20.000	ng/ul	-0.01
79) Chrysene-d12	21.163	240	481492	20.000	ng/ul	# 0.00
88) Perylene-d12	23.474	264	501546	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	12495	4.083	ng/uL	0.00
4) Pyridine-d5	3.499	84	44556	5.772	ng/ul	0.00
7) Phenol-d5	6.763	99	44898	5.408	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.928	67	151643	25.760	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.116	132	139135	21.358	ng/ul	0.00
15) 4-Methylphenol-d8	8.304	113	84164	12.911	ng/ul	-0.01
21) Nitrobenzene-d5	8.746	128	83014	31.336	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.463	143	78659	32.431	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.999	165	122127	24.934	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.522	131	159837	18.666	ng/ul	-0.01
46) Dimethylphthalate-d6	13.681	166	503455	32.336	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	570881	30.460	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.492	143	7413m	2.607	ng/ul	0.00
60) Fluorene-d10	15.257	176	413261	31.037	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.392	200	63183	28.549	ng/ul	0.00
73) Anthracene-d10	17.128	188	750802	35.057	ng/ul	-0.01
81) Pyrene-d10	19.392	212	898244	33.861	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	906656	35.846	ng/ul	0.00

Target Compounds Qvalue

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

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