

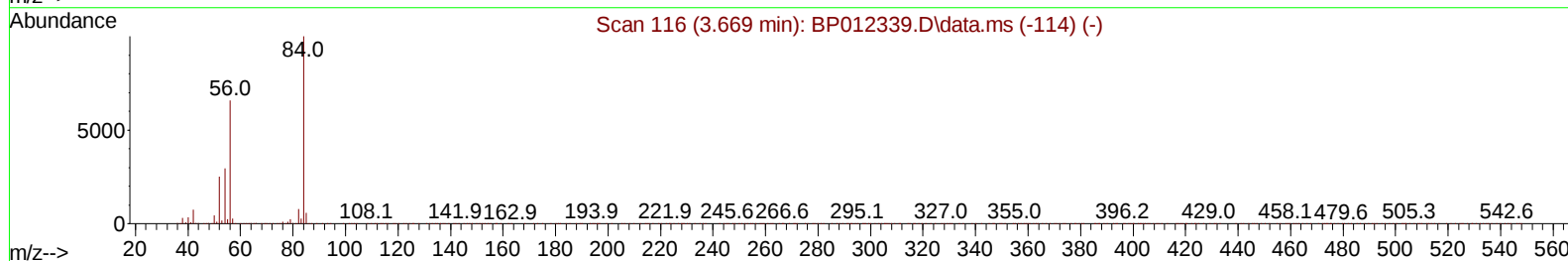
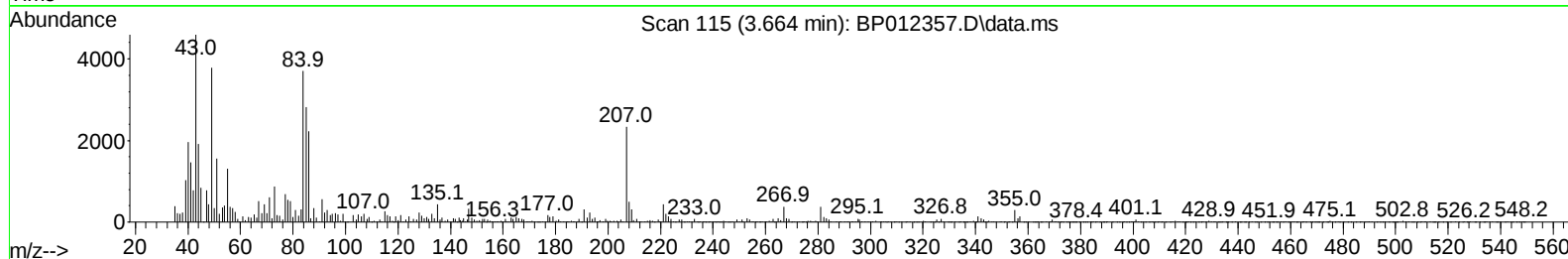
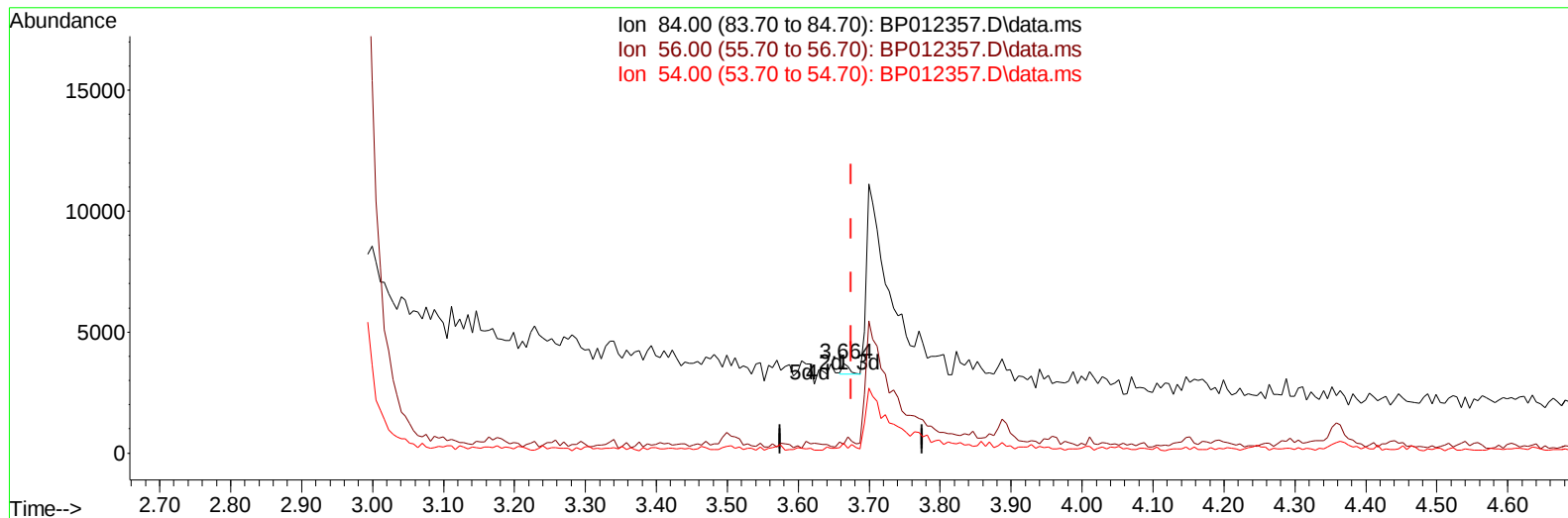
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012357.D
 Acq On : 27 Oct 2022 23:31
 Operator : CG/JU
 Sample : N5144-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BK8

Manual IntegrationsAPPROVED

Quant Time: Oct 28 00:14:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 26 23:40:20 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/28/2022
 Supervised By :mohammad ahmed 10/31/2022



TIC: BP012357.D\data.ms

(4) Pyridine-d5 (S)

3.664min (-0.011) 0.01 ng/u1

response	294	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	9.97#
54.00	29.80	10.75#
0.00	0.00	0.00

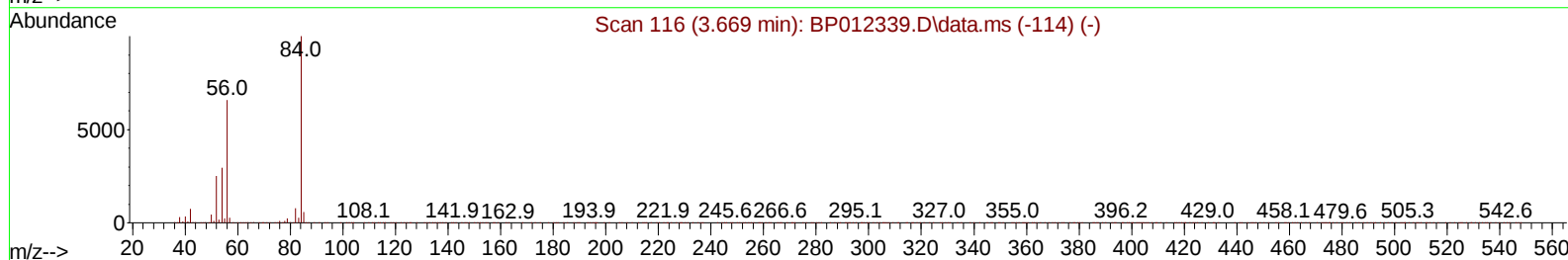
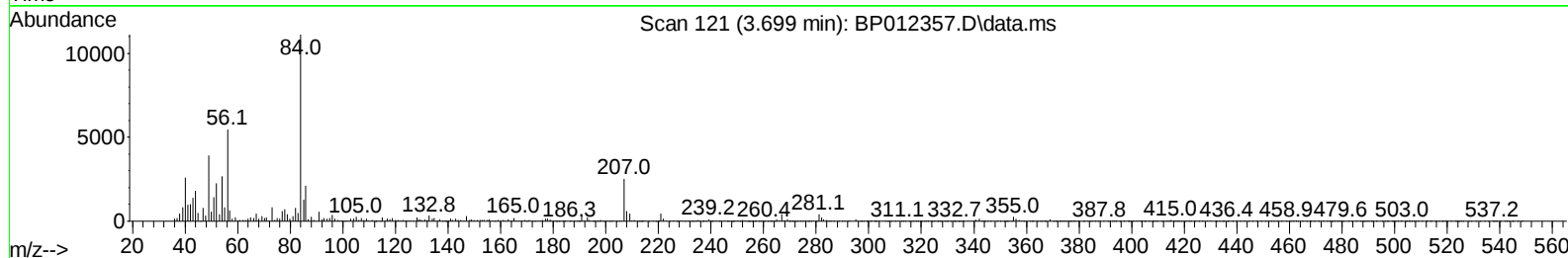
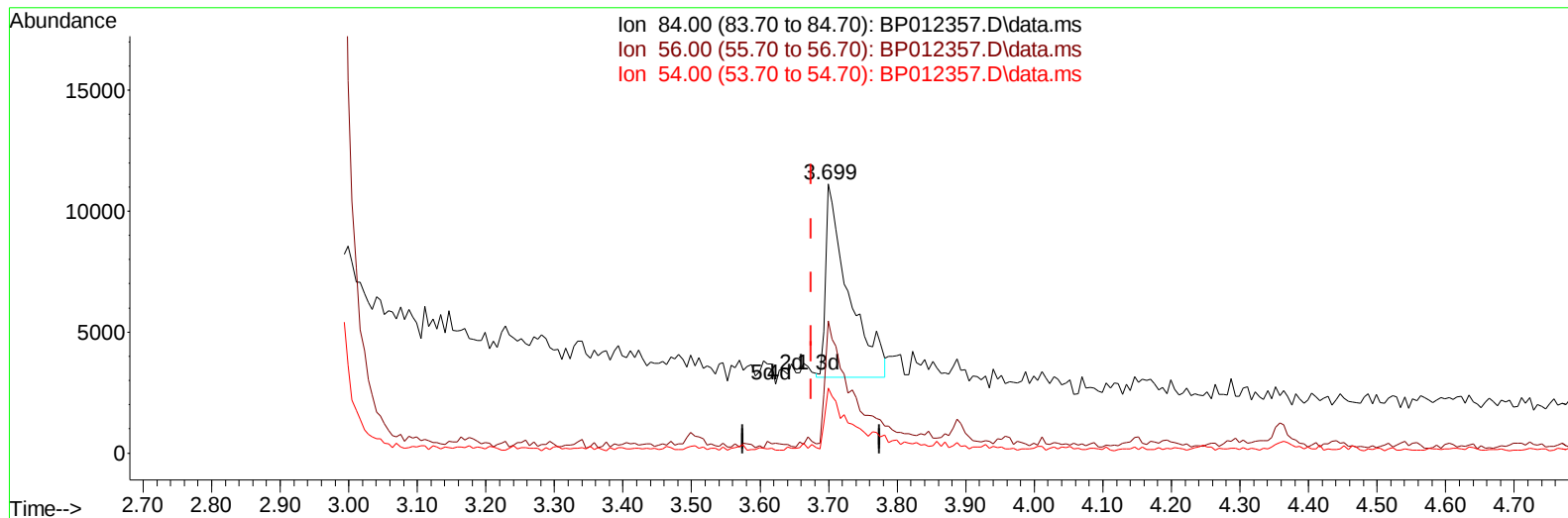
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(4) Pyridine-d5 (S)

3.699min (+ 0.024) 0.81 ng/ul m

response 18408

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	49.29#
54.00	29.80	24.23
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	320051	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.610	136	1473373	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	991049	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	2212979	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	2330577	20.000	ng/ul	0.00
88) Perylene-d12	23.710	264	2149714	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	34631	4.357	ng/uL	0.00
4) Pyridine-d5	3.699	84	18408m	0.806	ng/ul	0.02
7) Phenol-d5	6.975	99	740622	25.800	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	440012	25.677	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.334	132	578910	27.616	ng/ul	-0.01
15) 4-Methylphenol-d8	8.522	113	464121	20.442	ng/ul	-0.01
21) Nitrobenzene-d5	8.975	128	301790	31.586	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.699	143	336506	33.932	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	606856	28.237	ng/ul	0.00
31) 4-Chloroaniline-d4	10.758	131	550361	17.199	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	1975261	29.115	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	2256626	29.107	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	332116	25.866	ng/ul	0.00
60) Fluorene-d10	15.457	176	1759294	30.288	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.587	200	222841	19.341	ng/ul	0.00
73) Anthracene-d10	17.316	188	2674526	28.049	ng/ul	0.00
81) Pyrene-d10	19.557	212	3416459	29.198	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.551	264	3044303	28.983	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.005	94	31820	1.066	ng/ul#	85
80) Fluoranthene	19.234	202	208224	1.441	ng/ul	99
82) Pyrene	19.586	202	185017	1.236	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.216	149	104148	1.221	ng/ul	100
90) Benzo(b)fluoranthene	22.974	252	149961	1.102	ng/ul#	91

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

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