

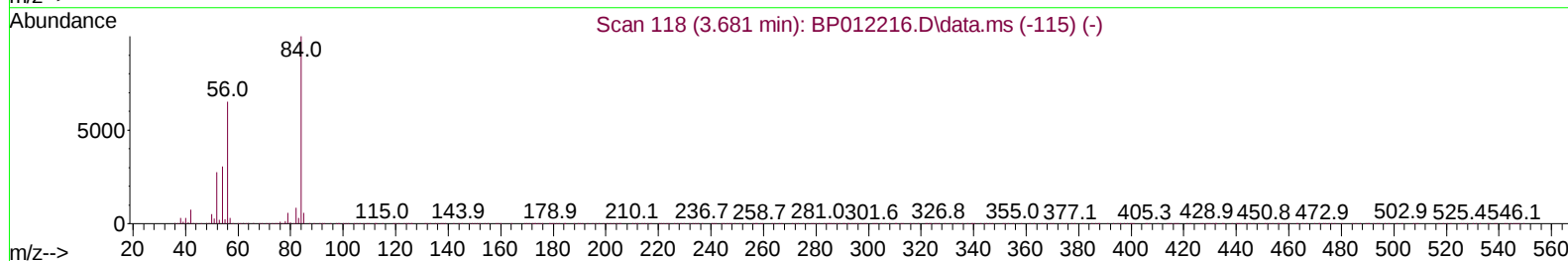
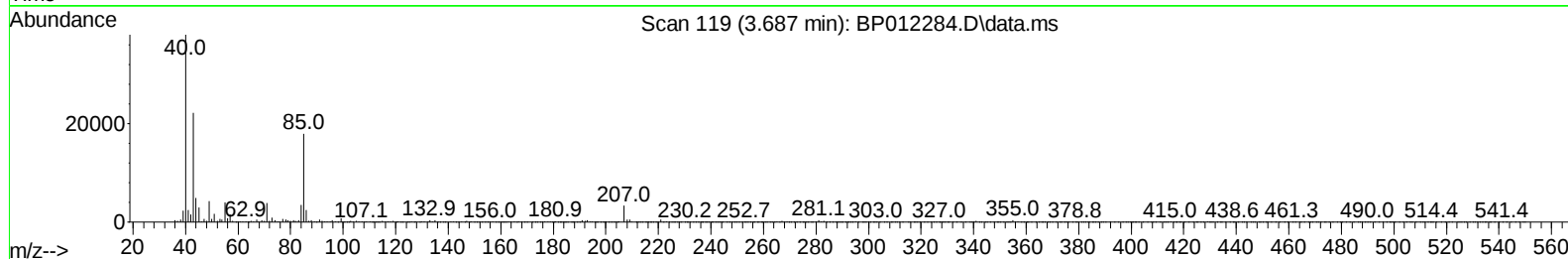
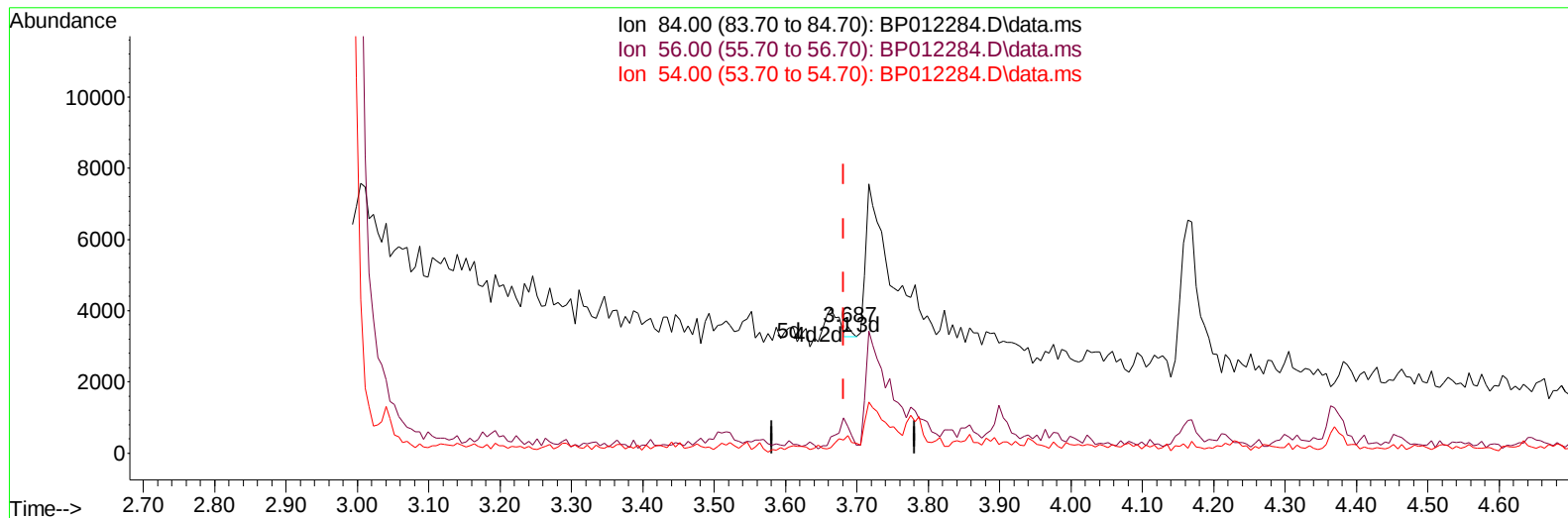
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
Data File : BP012284.D
Acq On : 24 Oct 2022 14:30
Operator : CG/JU
Sample : N5070-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BM2

Manual IntegrationsAPPROVED

Quant Time: Oct 25 05:49:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Oct 22 01:24:49 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/25/2022
Supervised By :mohammad ahmed 10/27/2022



TIC: BP012284.D\data.ms

(4) Pyridine-d5 (S)

3.687min (+ 0.006) 0.00 ng/u1

response 137

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	20.99#
54.00	29.70	13.96#
0.00	0.00	0.00

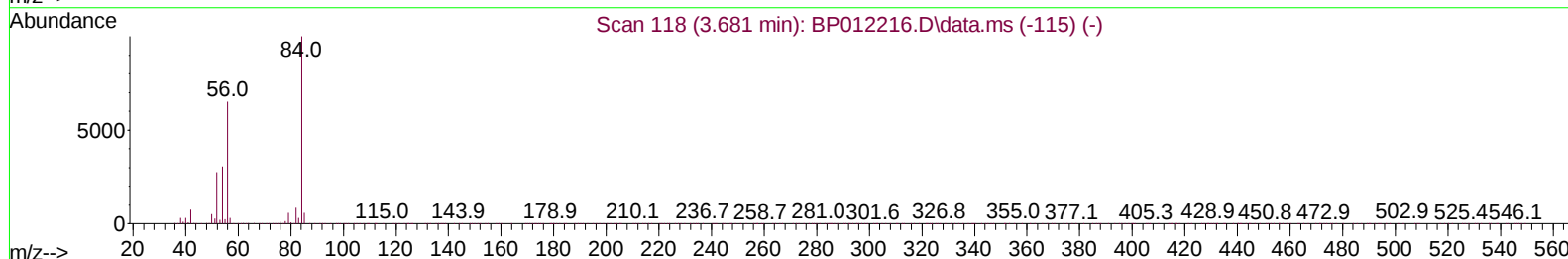
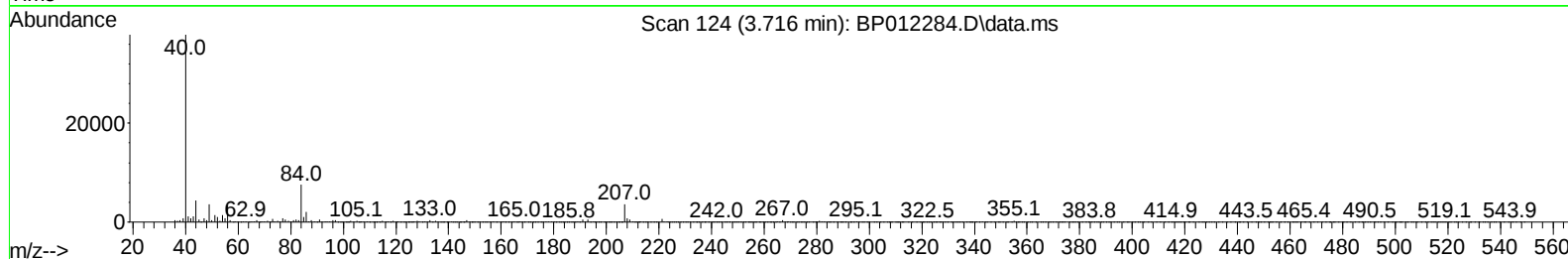
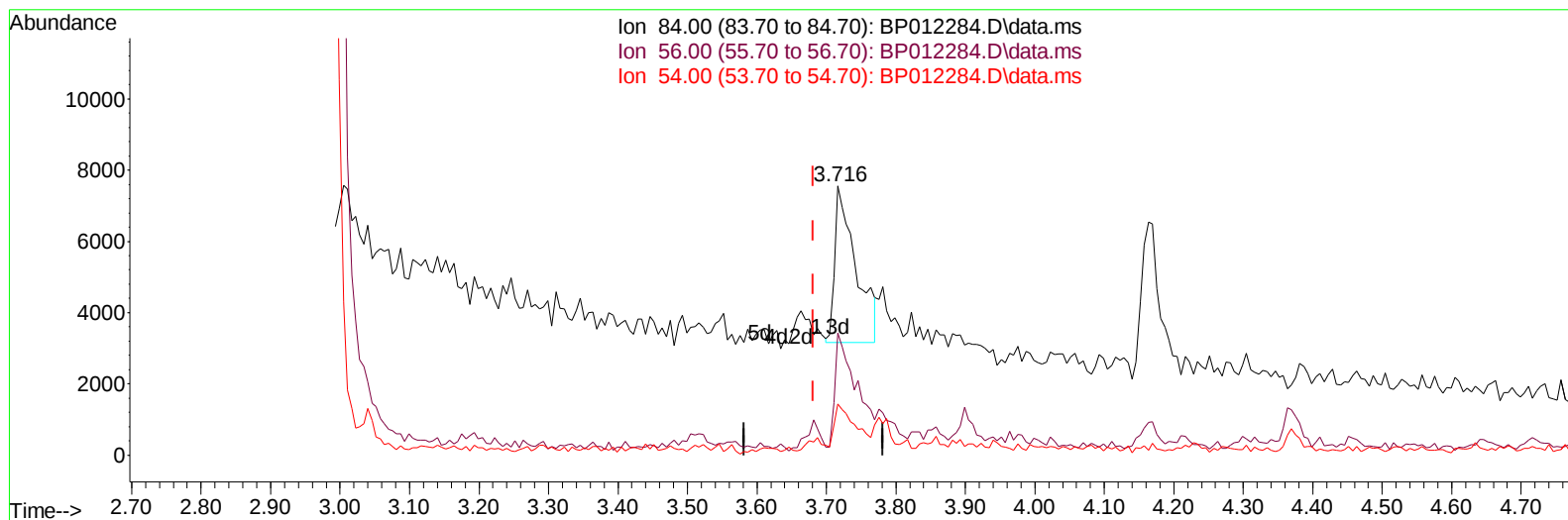
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(4) Pyridine-d5 (S)

3.716min (+ 0.035) 0.33 ng/u1 m

response 9192

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	45.44#
54.00	29.70	18.90#
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	7.816	152	386856	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	1535988	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	852212	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1660305	20.000	ng/ul	0.00
79) Chrysene-d12	21.321	240	1726457	20.000	ng/ul	0.00
88) Perylene-d12	23.715	264	1878201	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	40017	4.186	ng/uL	0.00
4) Pyridine-d5	3.716	84	9192m	0.334	ng/ul	0.04
7) Phenol-d5	6.987	99	656767	19.665	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	470028	23.206	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	565287	22.236	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	184096	7.092	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	292836	25.756	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	300710	24.514	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	467361	20.312	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	359868	11.008	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1522149	25.677	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	1712035	25.301	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	182726	16.453	ng/ul	0.00
60) Fluorene-d10	15.463	176	1311527	25.716	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	114970	12.172	ng/ul	0.00
73) Anthracene-d10	17.328	188	1711339	23.515	ng/ul	0.00
81) Pyrene-d10	19.569	212	2278702	23.182	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.562	264	2306747	24.491	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.016	94	45803	1.311	ng/ul	91
80) Fluoranthene	19.239	202	165292	1.361	ng/ul	99
82) Pyrene	19.592	202	154322	1.242	ng/ul	97
87) Chrysene	21.357	228	145580	1.386	ng/ul	98
90) Benzo(b)fluoranthene	22.986	252	179366	1.433	ng/ul#	92
93) Benzo(a)pyrene	23.610	252	107311	1.003	ng/ul#	95

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

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