

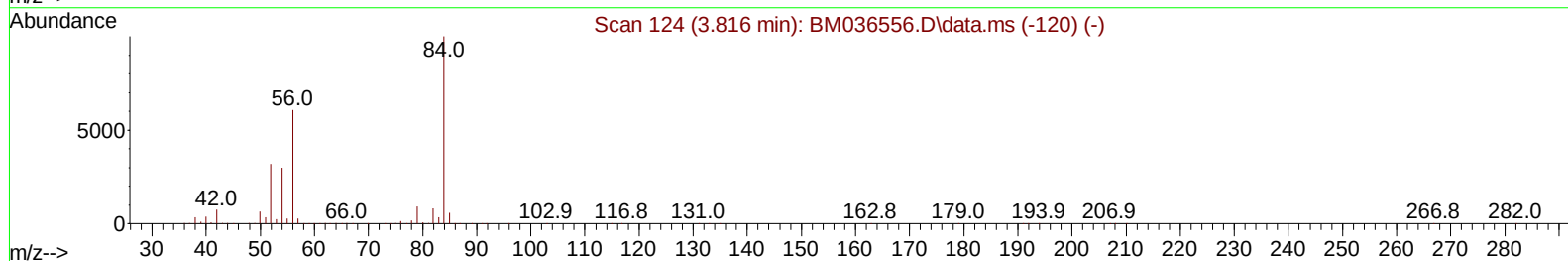
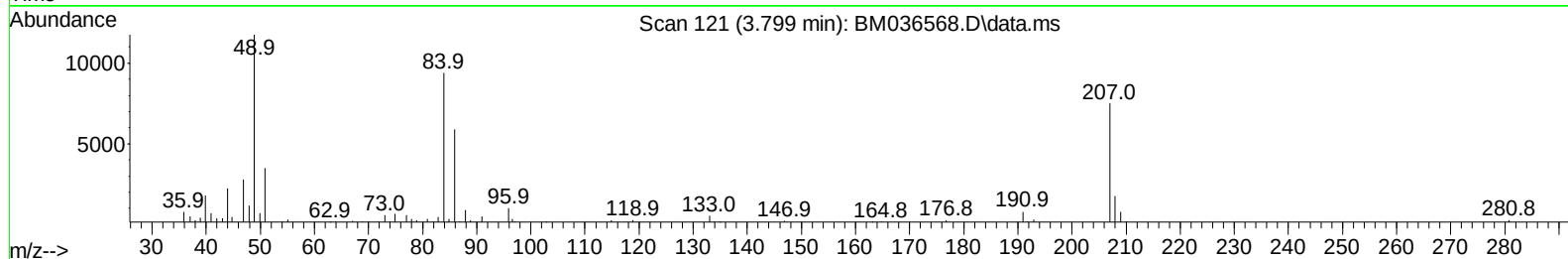
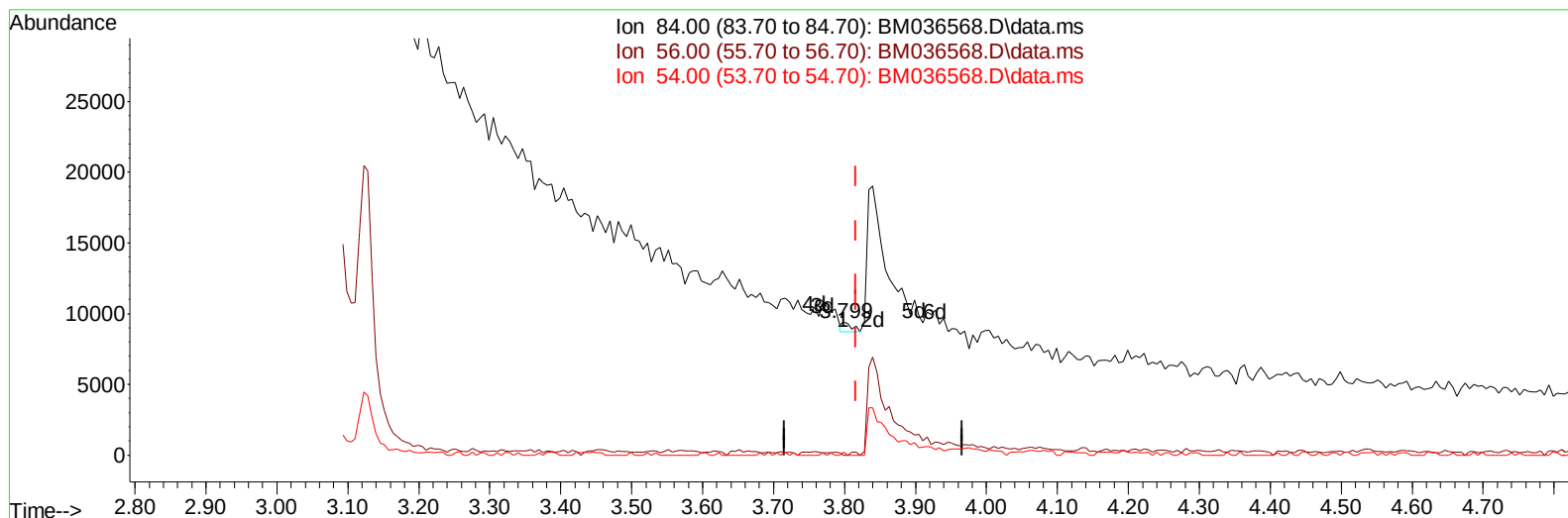
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
 Data File : BM036568.D
 Acq On : 16 Sep 2022 18:23
 Operator : CG/JU
 Sample : N4601-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A5751

Manual IntegrationsAPPROVED

Quant Time: Sep 16 23:01:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/19/2022
 Supervised By :mohammad ahmed 09/19/2022



TIC: BM036568.D\data.ms

(4) Pyridine-d5 (S)

3.799min (-0.017) 0.02 ng/ul

response 627

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.90	0.00#
54.00	30.90	2.00#
0.00	0.00	0.00

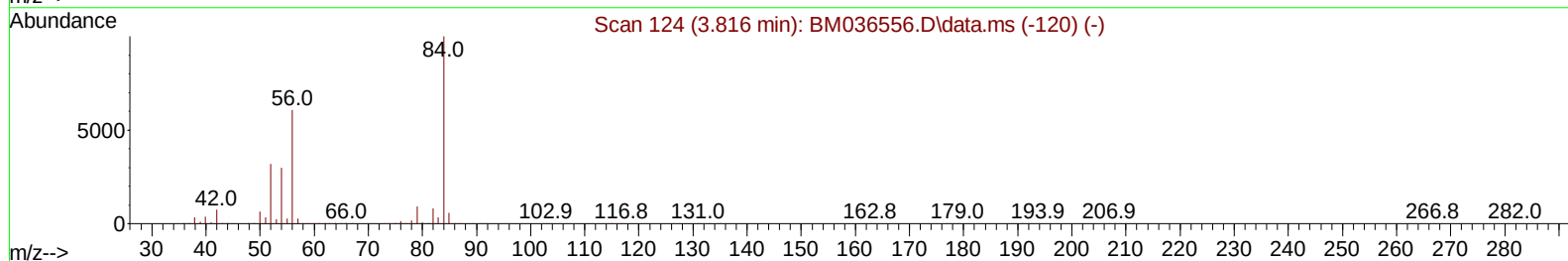
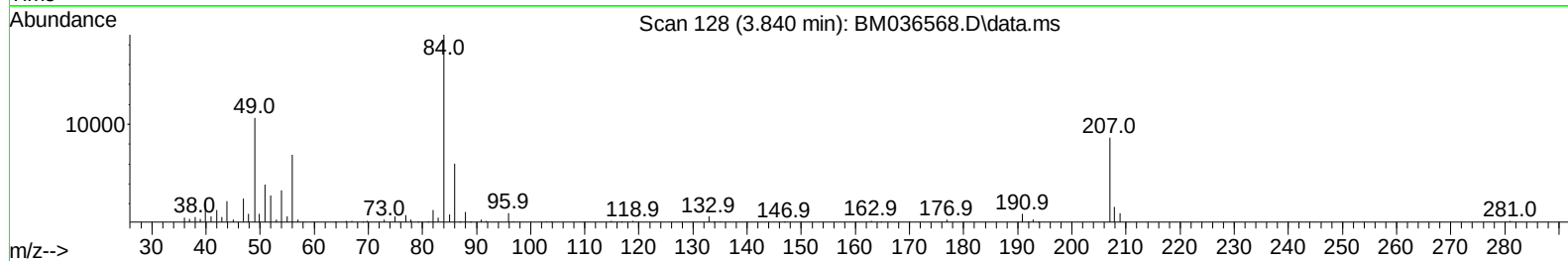
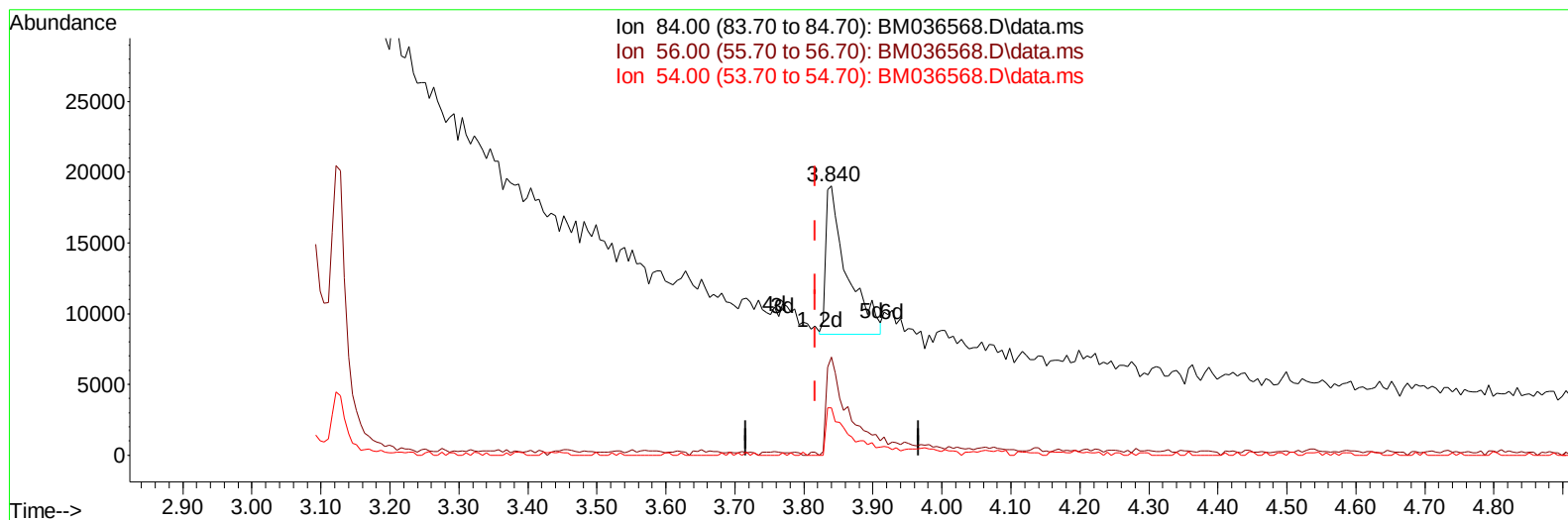
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(4) Pyridine-d5 (S)

3.840min (+ 0.024) 0.74 ng/ul m

response 22199

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.90	36.44#
54.00	30.90	17.70#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	384170	20.000	ng/ul	0.00
20) Naphthalene-d8	10.839	136	1702088	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	1137402	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.386	188	2394143	20.000	ng/ul	0.00
79) Chrysene-d12	21.550	240	2280419	20.000	ng/ul	0.00
88) Perylene-d12	23.991	264	2144978	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	33159	3.236	ng/uL	0.00
4) Pyridine-d5	3.840	84	22199m	0.743	ng/ul	0.02
7) Phenol-d5	7.175	99	523917	15.261	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	362565	17.390	ng/ul	0.00
11) 2-Chlorophenol-d4	7.551	132	431269	17.448	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	403145	15.892	ng/ul	0.00
21) Nitrobenzene-d5	9.192	128	208737	18.119	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	183854	17.817	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	387924	16.221	ng/ul	0.00
31) 4-Chloroaniline-d4	10.969	131	394724	9.927	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	1440693	18.781	ng/ul	0.00
49) Acenaphthylene-d8	14.345	160	1685803	18.527	ng/ul	0.00
54) 4-Nitrophenol-d4	14.821	143	141692	9.523	ng/ul	0.00
60) Fluorene-d10	15.639	176	1320461	19.208	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	98764	10.427	ng/ul	0.00
73) Anthracene-d10	17.486	188	2002468	18.293	ng/ul	0.00
81) Pyrene-d10	19.762	212	2292687	20.436	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	1928290	18.401	ng/ul	0.00

Target Compounds Qvalue

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

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