

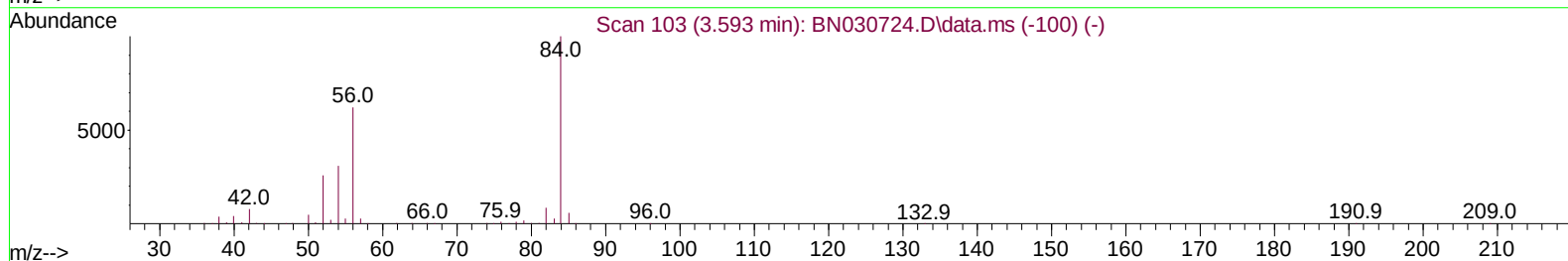
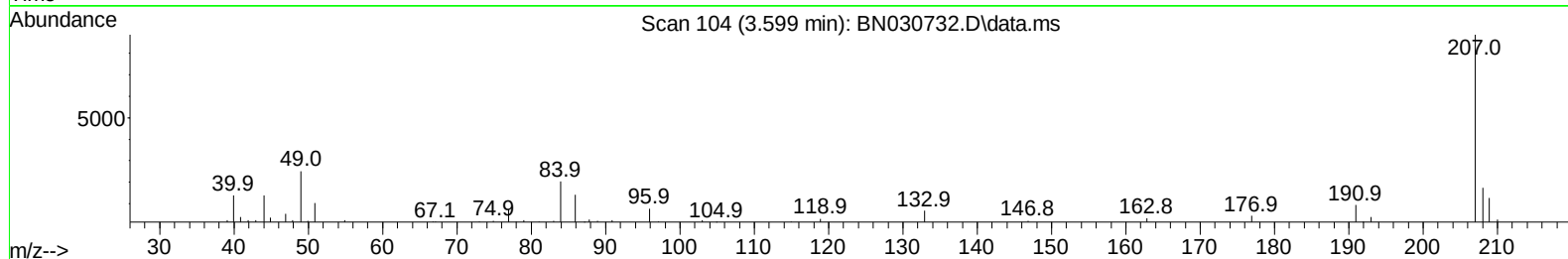
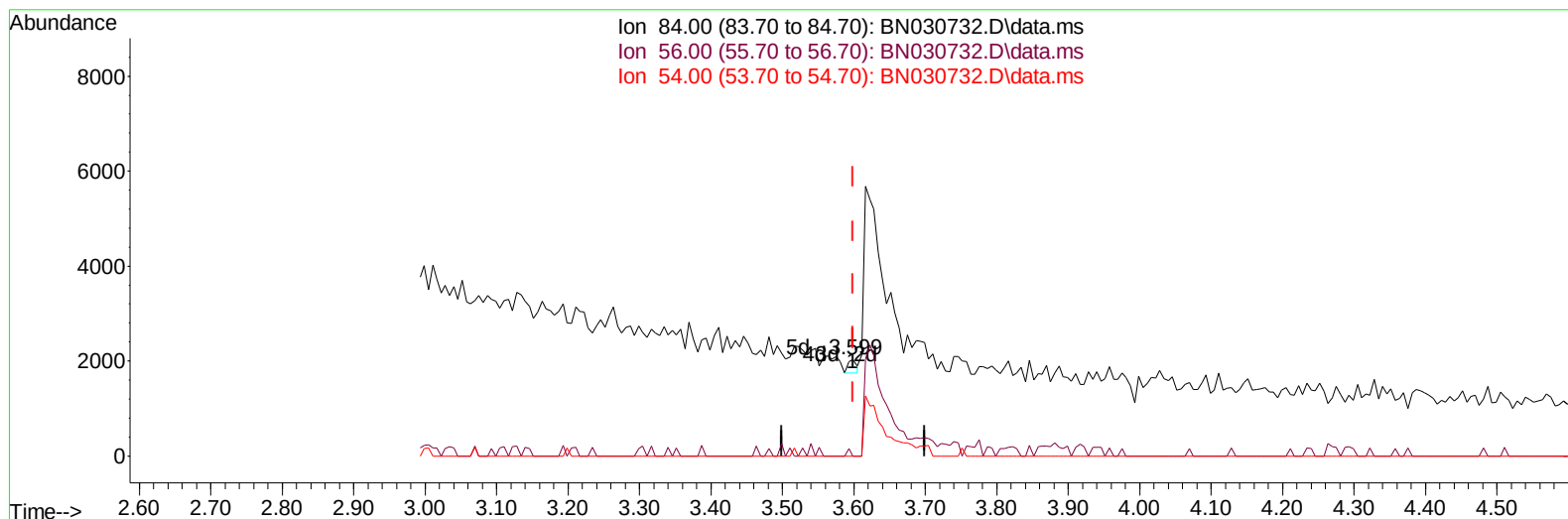
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040624\
Data File : BN030732.D
Acq On : 07 Apr 2024 07:30
Operator : MA/JU
Sample : P1987-10
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
MCORG8

Manual IntegrationsAPPROVED

Quant Time: Apr 07 13:43:55 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Mar 31 12:24:07 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/08/2024
Supervised By :mohammad ahmed 04/08/2024



TIC: BN030732.D\data.ms

(4) Pyridine-d5 (S)

3.599min (-0.000) 0.03 ng/u1

response 229

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.50	0.00#
54.00	31.00	0.00#
0.00	0.00	0.00

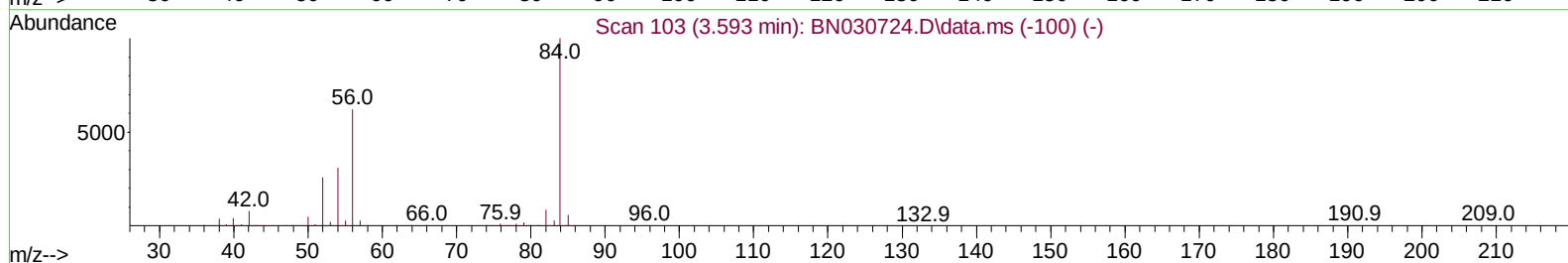
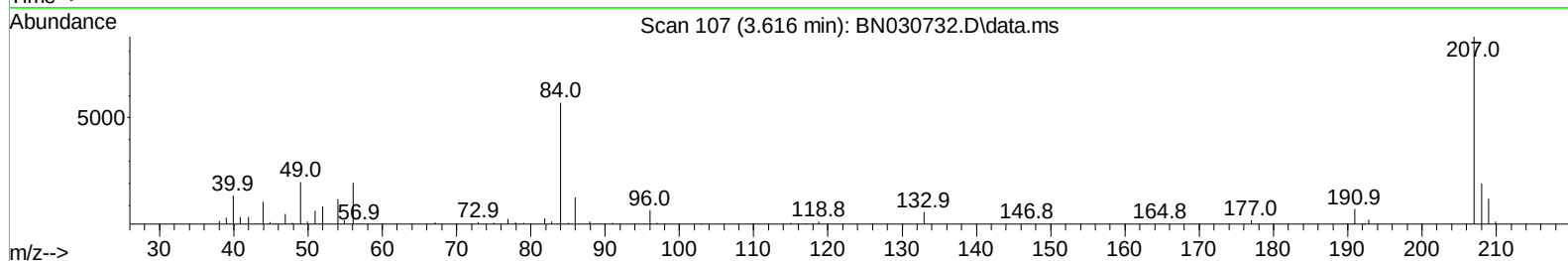
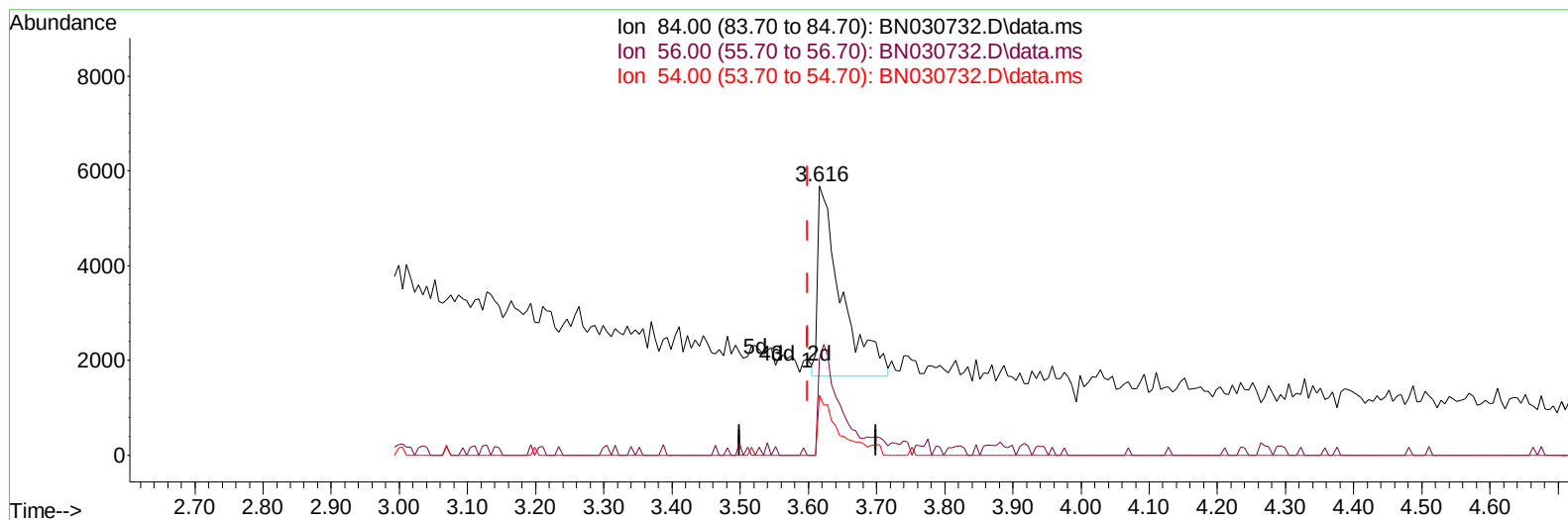
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(4) Pyridine-d5 (S)

3.616min (+ 0.017) 1.06 ng/ul m

response 9599

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.50	35.92#
54.00	31.00	22.27#
0.00	0.00	0.00

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Quant Time: Apr 07 13:44:33 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.687	152	141937	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.469	136	698033	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.328	164	510870	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.074	188	1142490	20.000	ng/ul	-0.01
79) Chrysene-d12	21.315	240	1124977	20.000	ng/ul	-0.01
88) Perylene-d12	24.251	264	1217823	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	15515	5.125	ng/uL	-0.01
4) Pyridine-d5	3.616	84	9599m	1.058	ng/ul	0.02
7) Phenol-d5	6.857	99	77908	6.785	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.028	67	274155	38.658	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.216	132	267545	30.716	ng/ul	-0.01
15) 4-Methylphenol-d8	8.387	113	167221	17.377	ng/ul	-0.01
21) Nitrobenzene-d5	8.840	128	191964	39.644	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.557	143	189005	42.668	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.092	165	337630	34.815	ng/ul	0.00
31) 4-Chloroaniline-d4	10.610	131	69168	4.332	ng/ul	-0.01
46) Dimethylphthalate-d6	13.745	166	1447093	42.426	ng/ul	-0.01
49) Acenaphthylene-d8	14.022	160	1533552	39.104	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.533	143	36161	5.543	ng/ul	0.00
60) Fluorene-d10	15.328	176	1256589	42.100	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.445	200	165572	32.833	ng/ul	0.00
73) Anthracene-d10	17.174	188	2210933	44.914	ng/ul	-0.01
81) Pyrene-d10	19.474	212	2642180	43.684	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.056	264	2700078	48.117	ng/ul	-0.02

Target Compounds Qvalue

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

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