

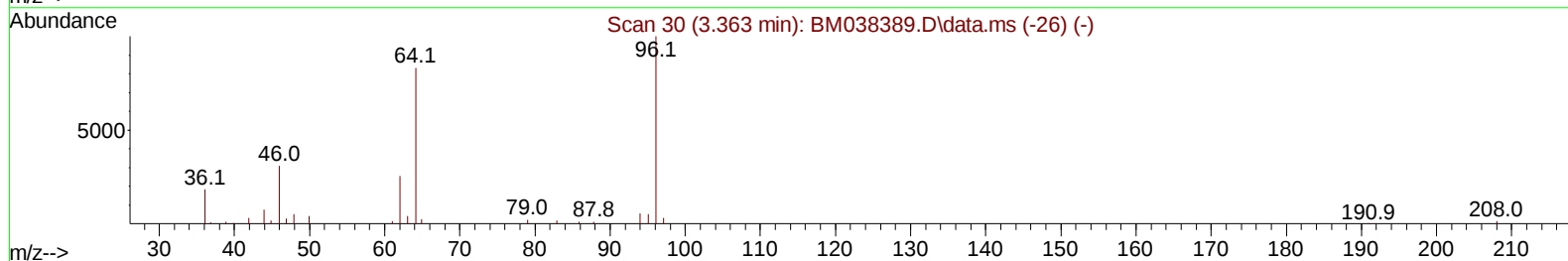
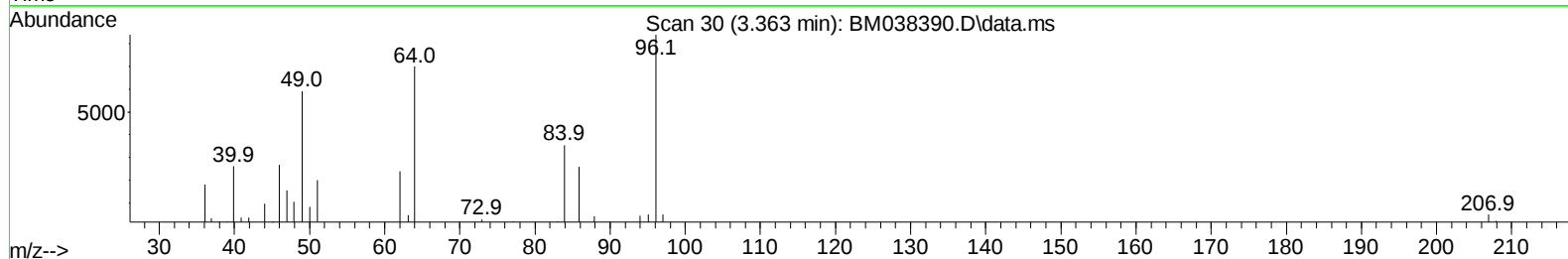
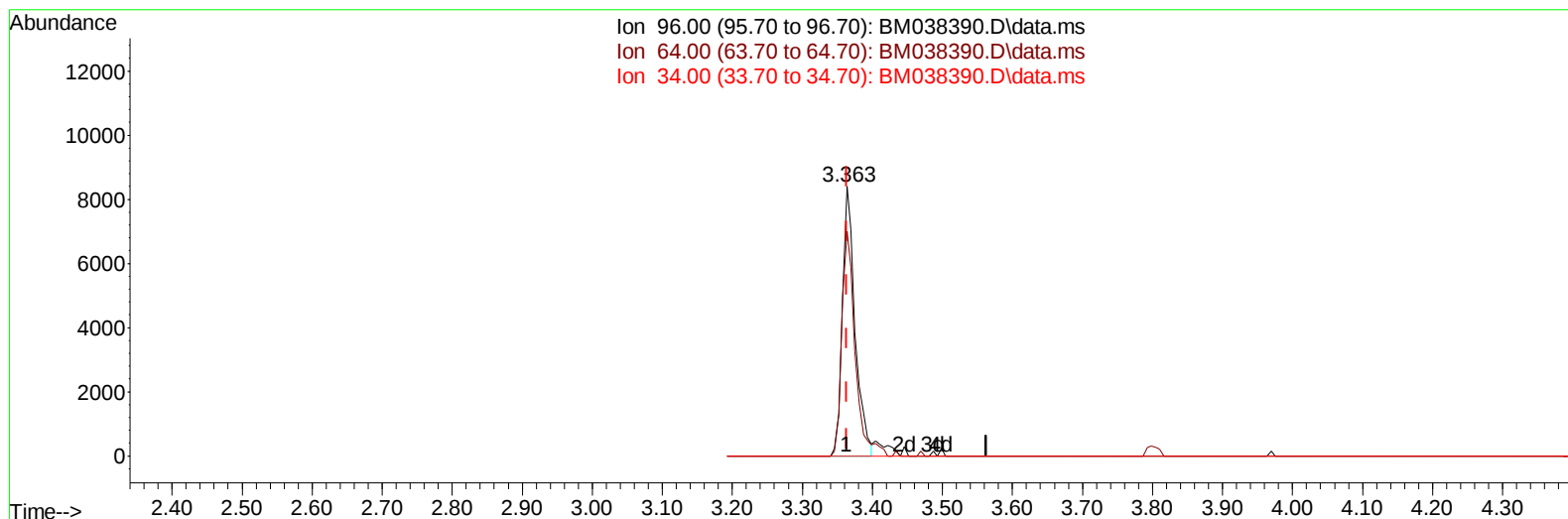
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010623\
Data File : BM038390.D
Acq On : 06 Jan 2023 17:30
Operator : CG/JU
Sample : PB150076BL
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK076

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/09/2023
Supervised By :mohammad ahmed 01/09/2023

Quant Time: Jan 07 00:38:34 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM010423.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jan 05 02:11:50 2023
Response via : Initial Calibration



TIC: BM038390.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.363min (+ 0.000) 4.78 ng/uL

response 10685

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	88.50	83.26
34.00	0.00	0.00
0.00	0.00	0.00

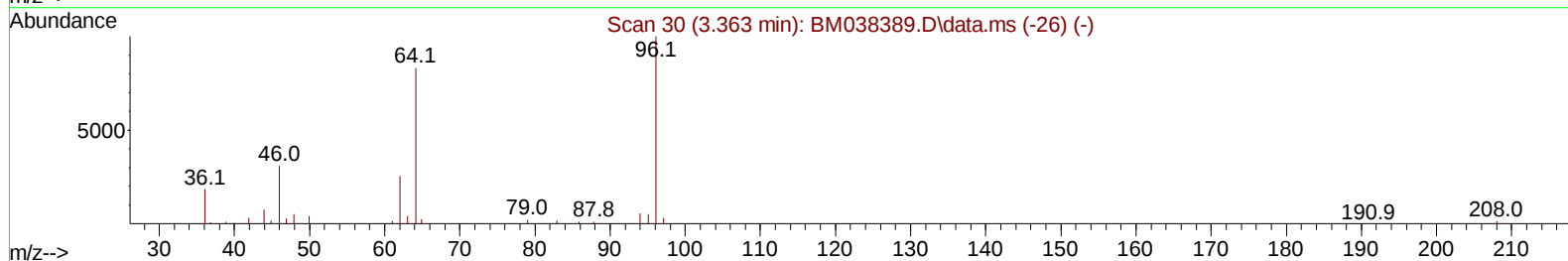
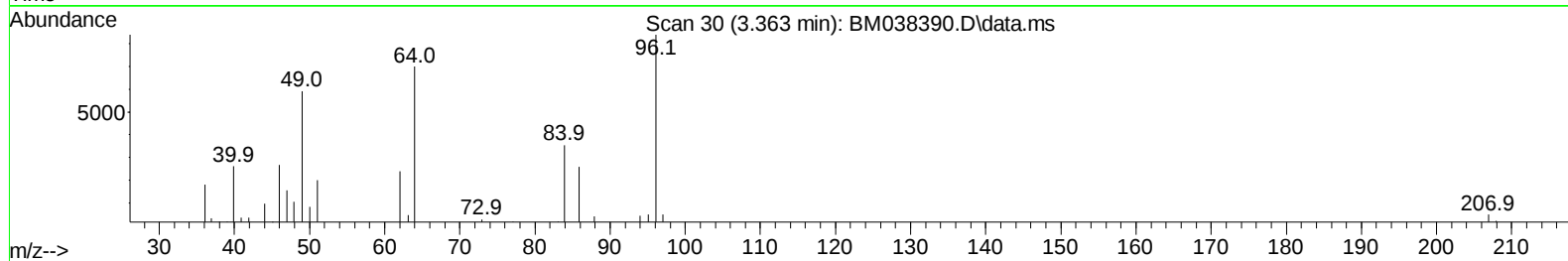
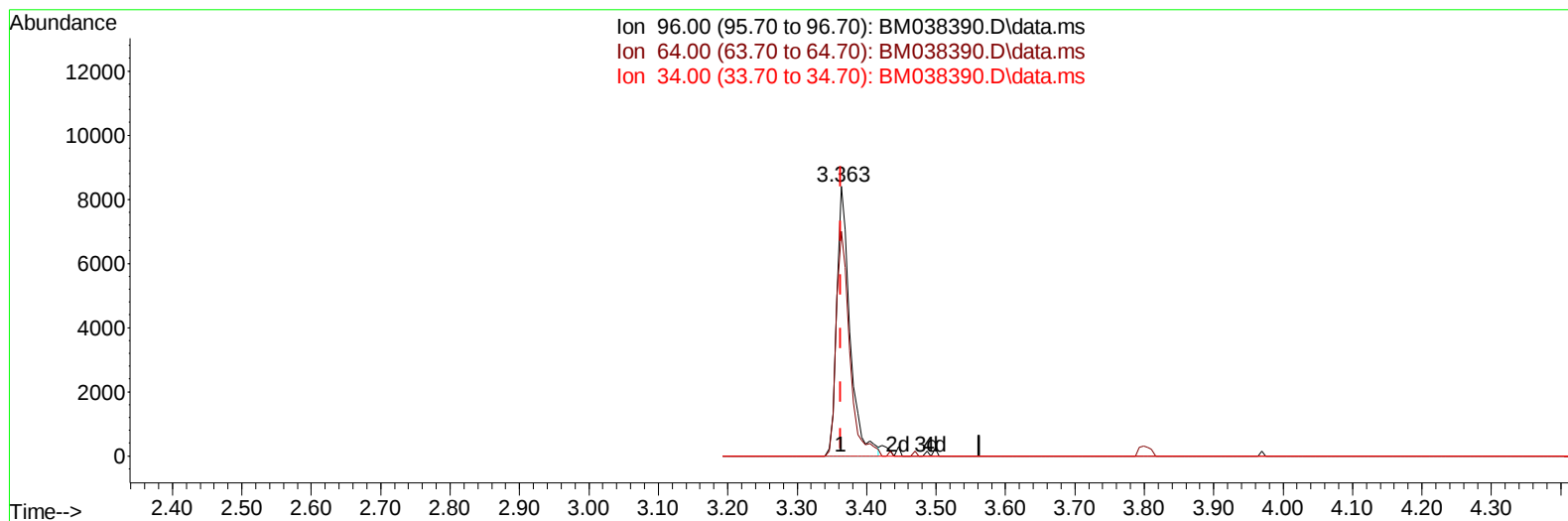
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TIC: BM038390.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.363min (+ 0.000) 4.95 ng/uL m

response 11073

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	88.50	83.26
34.00	0.00	0.00
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.998	152		72510	20.000	ng/ul	0.00
20) Naphthalene-d8	10.810	136		272155	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.633	164		163792	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.368	188		366313	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240		345650	20.000	ng/ul	0.00
88) Perylene-d12	23.968	264		417202	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.363	96		11073m	4.952	ng/uL	0.00
4) Pyridine-d5	3.799	84		160792	24.856	ng/ul	0.00
7) Phenol-d5	7.157	99		175336	26.544	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67		134776	28.363	ng/ul	0.00
11) 2-Chlorophenol-d4	7.528	132		131767	27.841	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113		124353	24.388	ng/ul	0.00
21) Nitrobenzene-d5	9.175	128		56159	26.301	ng/ul	0.00
24) 2-Nitrophenol-d4	9.892	143		64471	26.623	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.434	165		115869	24.778	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131		151467	23.472	ng/ul	0.00
46) Dimethylphthalate-d6	14.039	166		345449	27.849	ng/ul	0.00
49) Acenaphthylene-d8	14.327	160		380899	26.350	ng/ul	0.00
54) 4-Nitrophenol-d4	14.833	143		48120	21.549	ng/ul	0.00
60) Fluorene-d10	15.621	176		295789	26.221	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200		58965	22.210	ng/ul	0.00
73) Anthracene-d10	17.468	188		484897	27.441	ng/ul	0.00
81) Pyrene-d10	19.756	212		589129	30.629	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.815	264		585951	28.278	ng/ul	0.00

Target Compounds Qvalue

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

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