

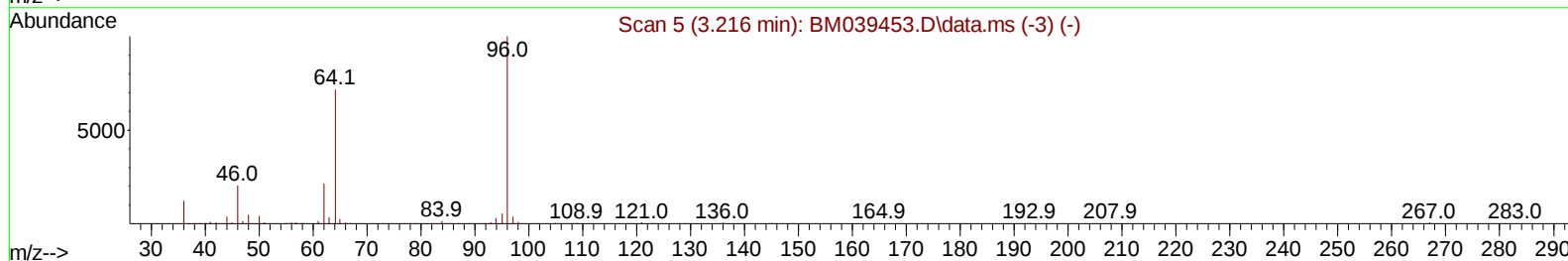
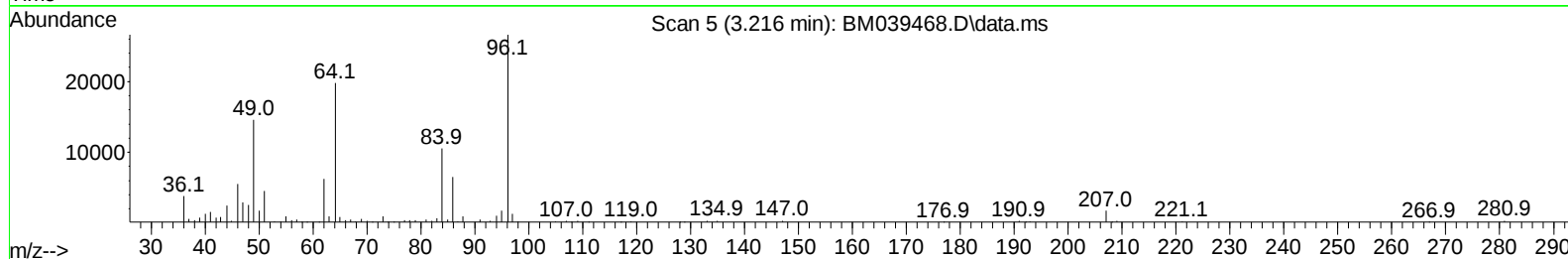
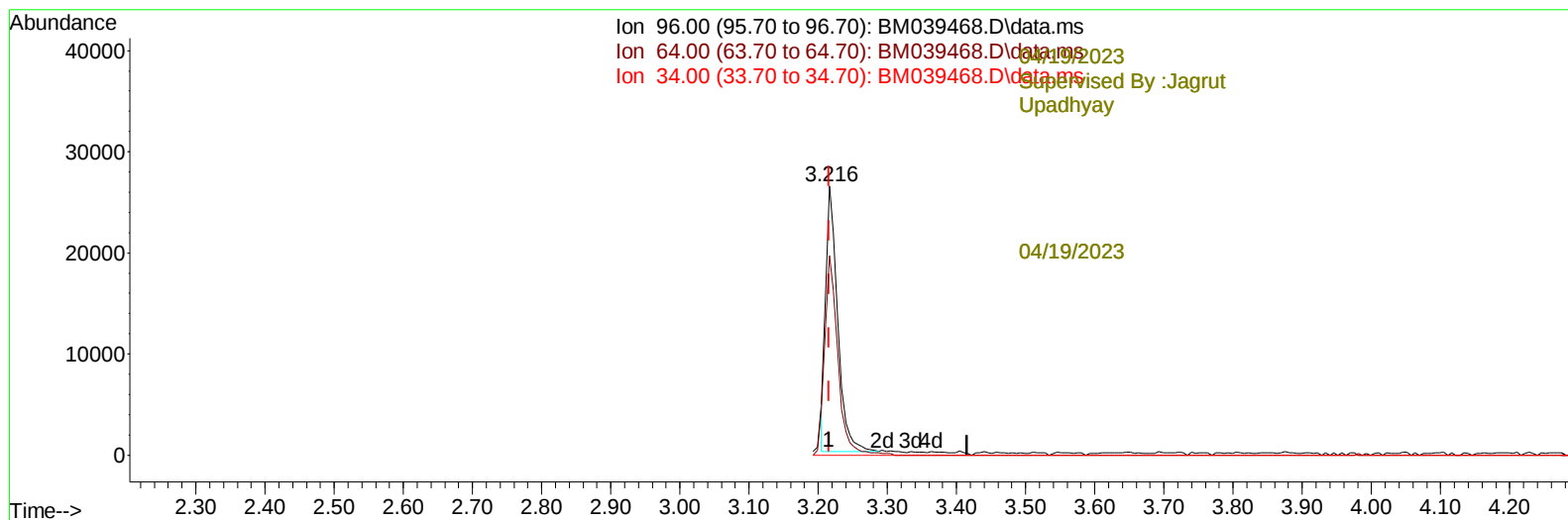
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
 Data File : BM039468.D
 Acq On : 18 Apr 2023 22:04
 Operator : CG/JU
 Sample : PB152152BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK152

Manual IntegrationsAPPROVED

Quant Time: Apr 19 00:15:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 18 23:56:45 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/19/2023
 Supervised By :Jagrut Upadhyay 04/19/2023



TIC: BM039468.D\data.ms

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

3.216min (+ 0.000) 5.99 ng/uL

response 31787

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.10	74.26
34.00	0.00	0.00
0.00	0.00	0.00

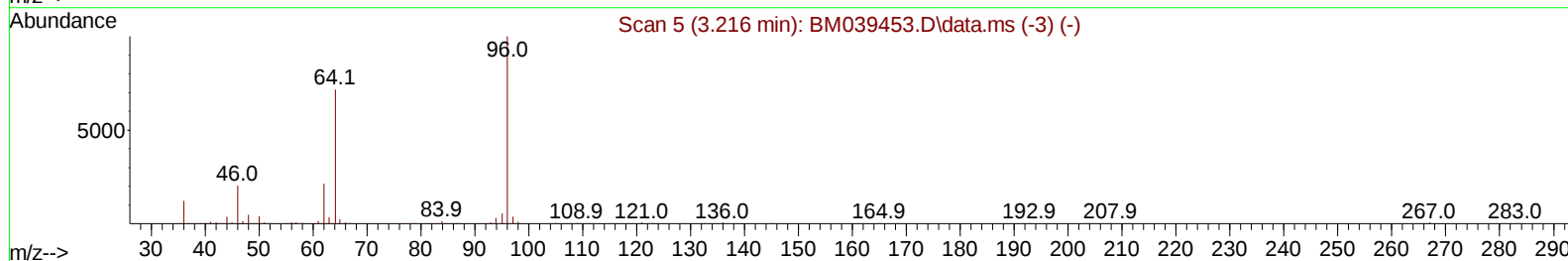
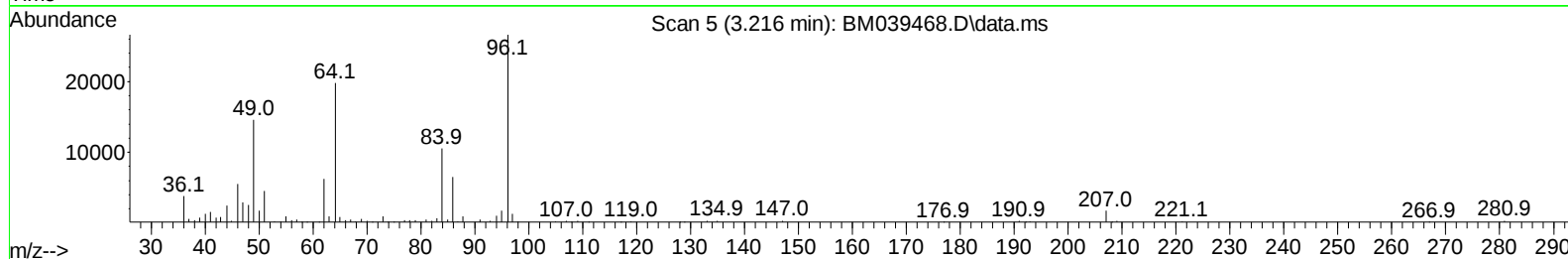
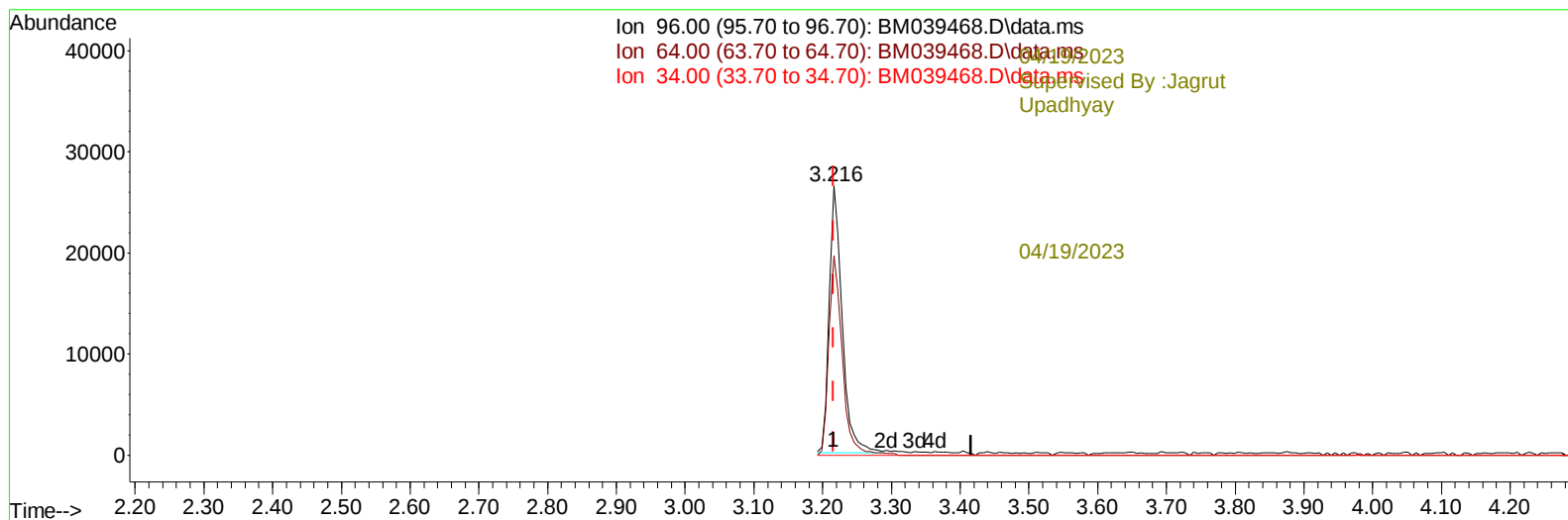
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TIC: BM039468.D\data.ms

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

3.216min (+ 0.000) 6.46 ng/uL m

response 34321

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.10	74.26
34.00	0.00	0.00
0.00	0.00	0.00

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Instrument :
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Quant Time: Apr 19 02:13:23 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----04/19/2023						
Supervised By :Jagrut Upadhyay						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	203822	20.000	ng/ul	0.00
20) Naphthalene-d8	10.639	136	918784	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.486	164	616025	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	1322299	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	1077528	20.000	ng/ul	0.00
88) Perylene-d12	23.803	264	1012540	20.000	ng/ul	0.00
-----04/19/2023						
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	34321m	6.463	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	0.00
7) Phenol-d5	7.016	99	487110	26.112	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.163	67	384685	31.199	ng/ul	0.00
11) 2-Chlorophenol-d4	7.363	132	452167	30.899	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	395685	26.332	ng/ul	0.00
21) Nitrobenzene-d5	9.004	128	228628	30.930	ng/ul	0.00
24) 2-Nitrophenol-d4	9.728	143	259285	31.238	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	390945	27.374	ng/ul	0.00
31) 4-Chloroaniline-d4	10.645	131	1297	0.061	ng/ul	-0.15
46) Dimethylphthalate-d6	13.904	166	1662920	33.472	ng/ul	0.00
49) Acenaphthylene-d8	14.180	160	1668445	29.893	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	164840	21.923	ng/ul	0.00
60) Fluorene-d10	15.486	176	1347547	33.106	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.615	200	248969	28.334	ng/ul	0.00
73) Anthracene-d10	17.339	188	2094679	32.939	ng/ul	0.00
81) Pyrene-d10	19.639	212	2403506	33.568	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.650	264	1687108	30.862	ng/ul	0.00

Target Compounds Qvalue

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

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