

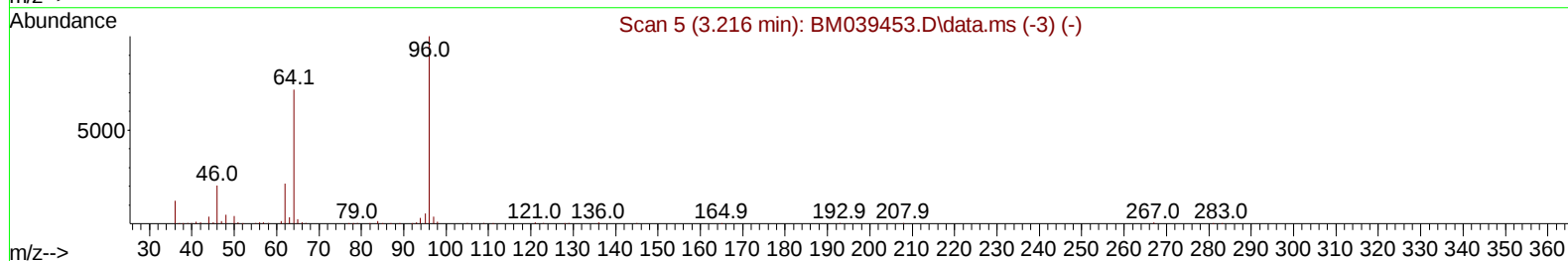
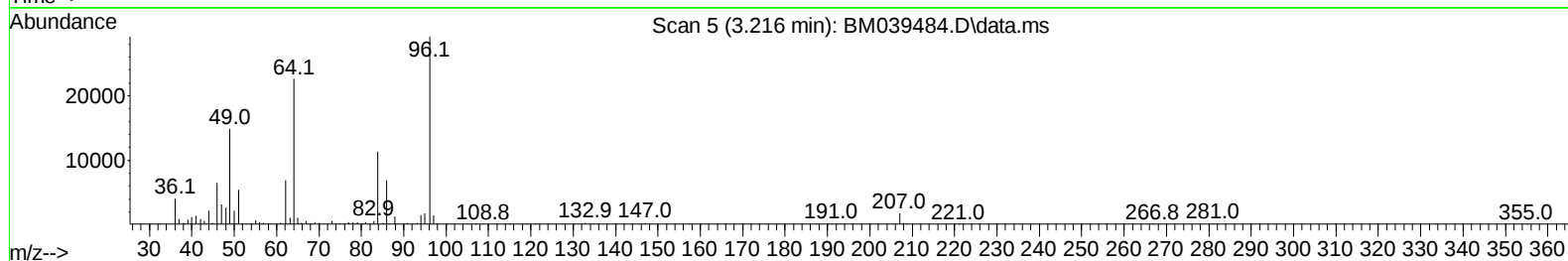
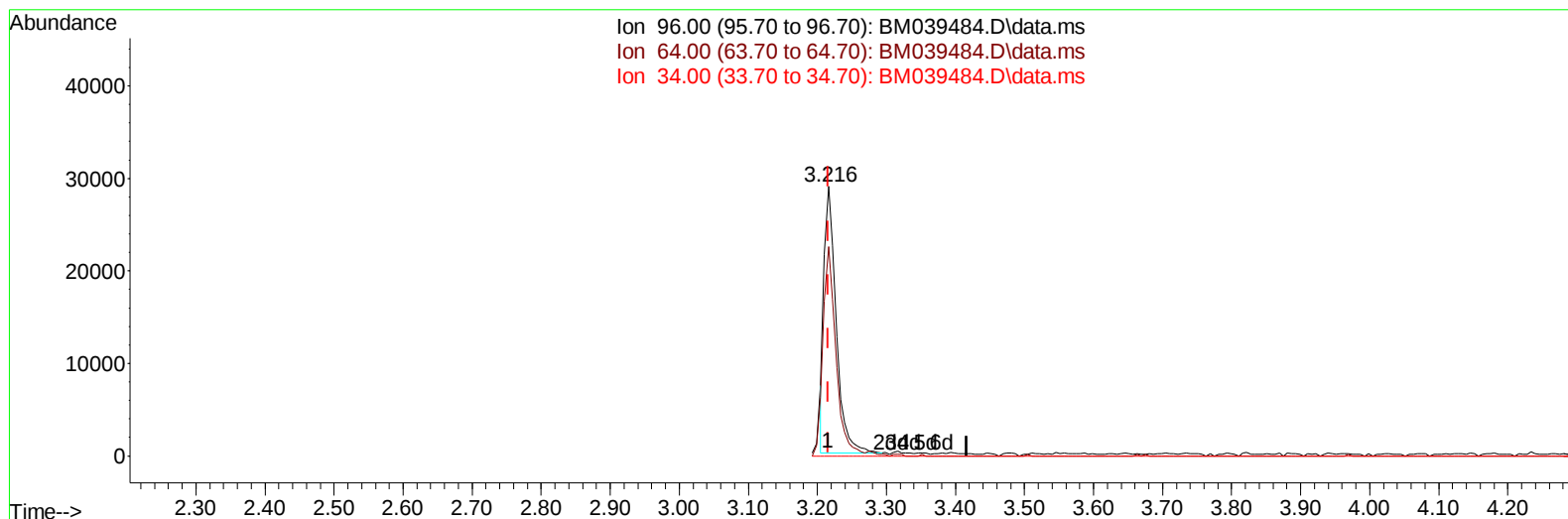
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
 Data File : BM039484.D
 Acq On : 19 Apr 2023 08:17
 Operator : CG/JU
 Sample : PB152174BL
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK174

Manual IntegrationsAPPROVED

Quant Time: Apr 19 08:49:07 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 19 08:18:41 2023
 Response via : Initial Calibration

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



TIC: BM039484.D\data.ms

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

3.216min (-0.000) 5.77 ng/uL

response 35027

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.10	77.70
34.00	0.00	0.00
0.00	0.00	0.00

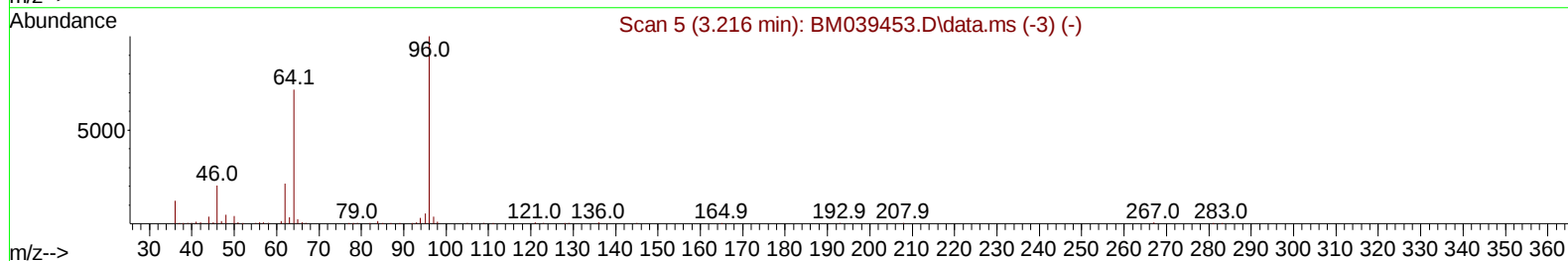
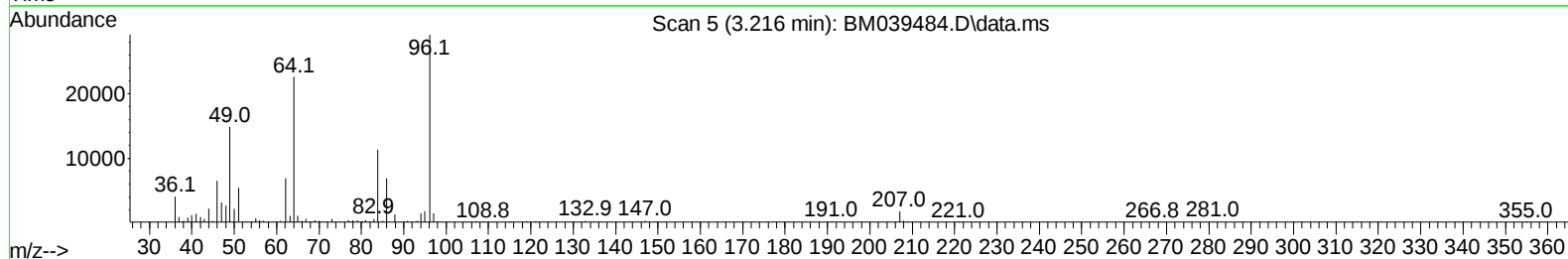
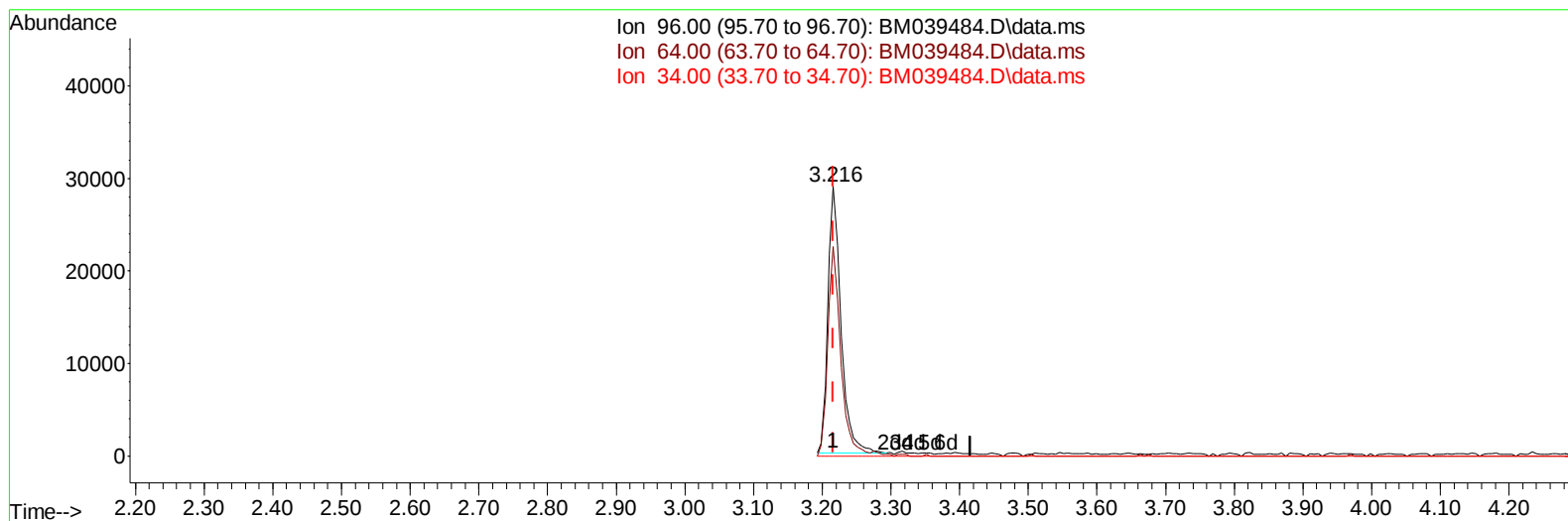
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TIC: BM039484.D\data.ms

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

3.216min (-0.000) 6.30 ng/uL m

response 38252

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.10	77.70
34.00	0.00	0.00
0.00	0.00	0.00

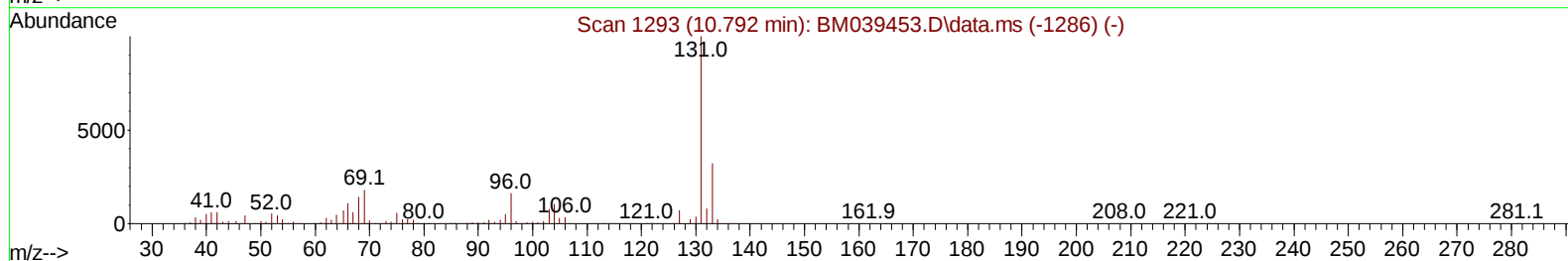
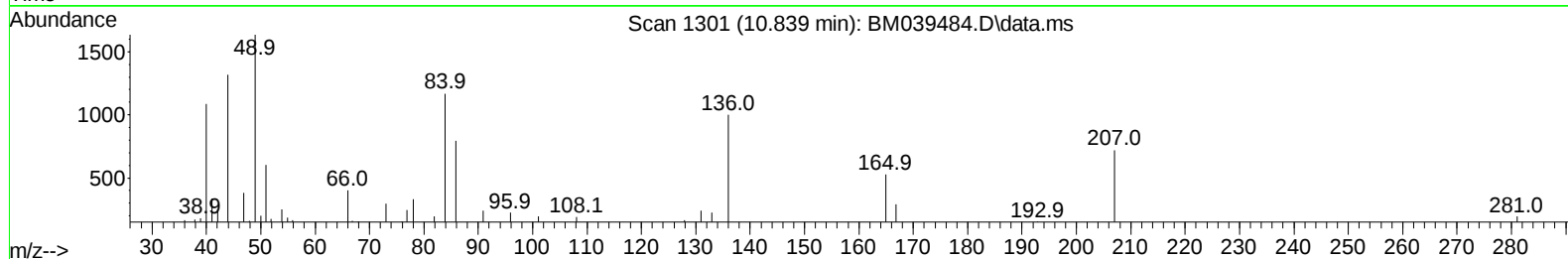
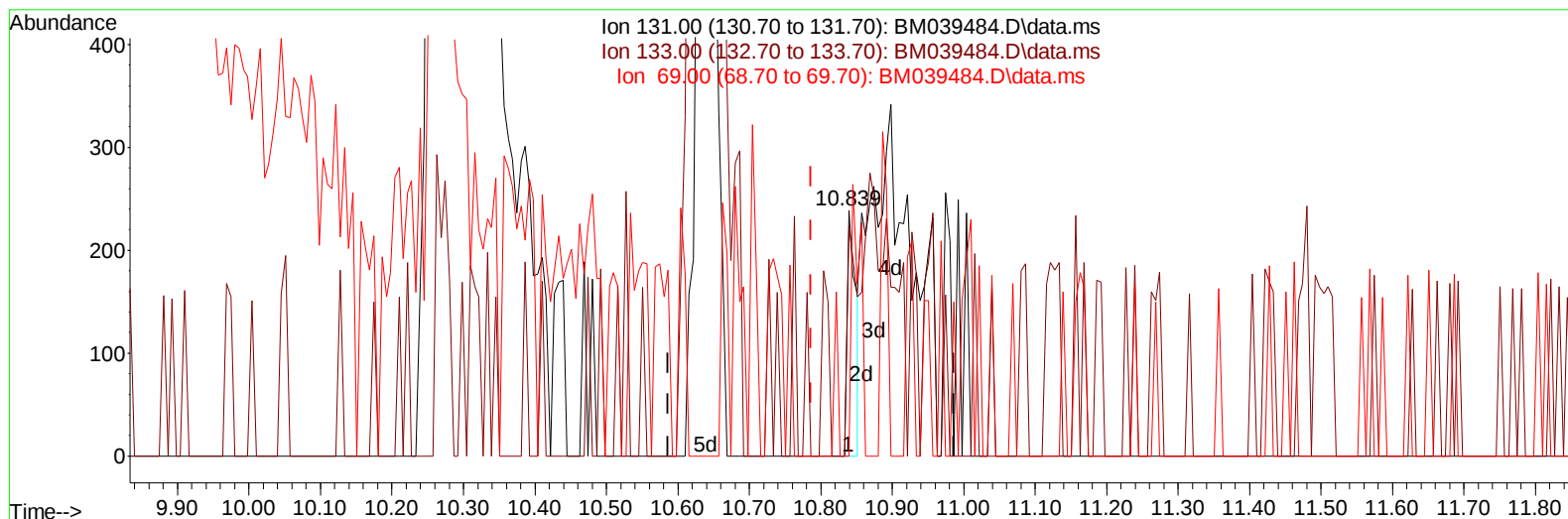
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Manual IntegrationsAPPROVED

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TIC: BM039484.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.839min (+ 0.053) 0.01 ng/ul

response 202

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	93.72#
69.00	18.10	0.00#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/19/2023
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Quant Time: Apr 19 08:51:24 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	233213	20.000	ng/ul	0.00
20) Naphthalene-d8	10.633	136	1059170	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.486	164	700123	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	1497399	20.000	ng/ul	0.00
79) Chrysene-d12	21.427	240	1257440	20.000	ng/ul	0.00
88) Perylene-d12	23.797	264	1163063	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	38252m	6.296	ng/uL	0.00
4) Pyridine-d5	3.657	84	53152	3.113	ng/ul	0.02
7) Phenol-d5	7.016	99	529591	24.811	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.163	67	435097	30.840	ng/ul	0.00
11) 2-Chlorophenol-d4	7.363	132	506779	30.267	ng/ul	0.00
15) 4-Methylphenol-d8	8.557	113	418505	24.340	ng/ul	0.00
21) Nitrobenzene-d5	8.998	128	259932	30.504	ng/ul	0.00
24) 2-Nitrophenol-d4	9.722	143	292770	30.598	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.269	165	446825	27.139	ng/ul	0.00
31) 4-Chloroaniline-d4	10.898	131	1629m	0.066	ng/ul	0.11
46) Dimethylphthalate-d6	13.904	166	1836427	32.524	ng/ul	0.00
49) Acenaphthylene-d8	14.180	160	1668478	26.303	ng/ul	0.00
54) 4-Nitrophenol-d4	14.727	143	188204	22.024	ng/ul	0.00
60) Fluorene-d10	15.480	176	1518337	32.821	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.615	200	277303	27.868	ng/ul	0.00
73) Anthracene-d10	17.339	188	2290773	31.811	ng/ul	0.00
81) Pyrene-d10	19.633	212	2688004	32.170	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.644	264	1845360	29.388	ng/ul	0.00

Target Compounds Qvalue

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

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