

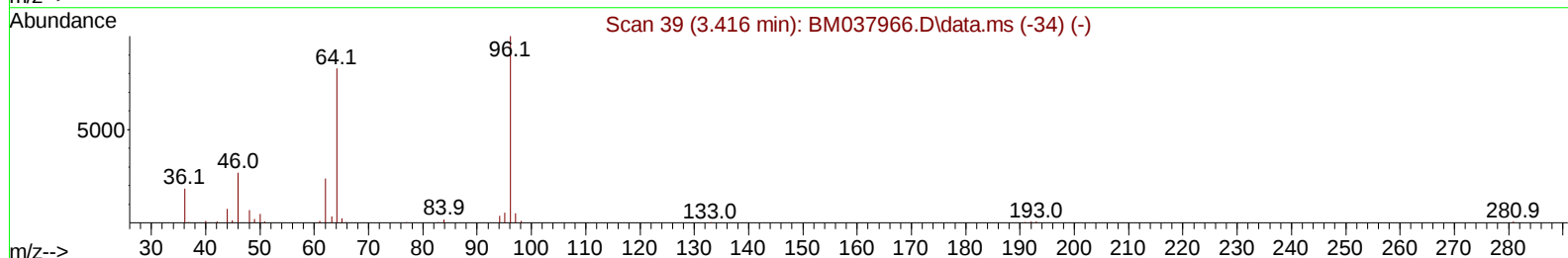
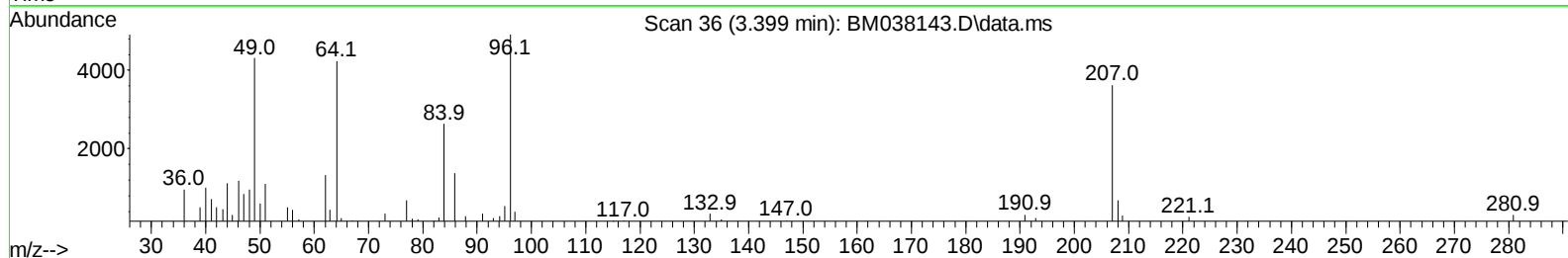
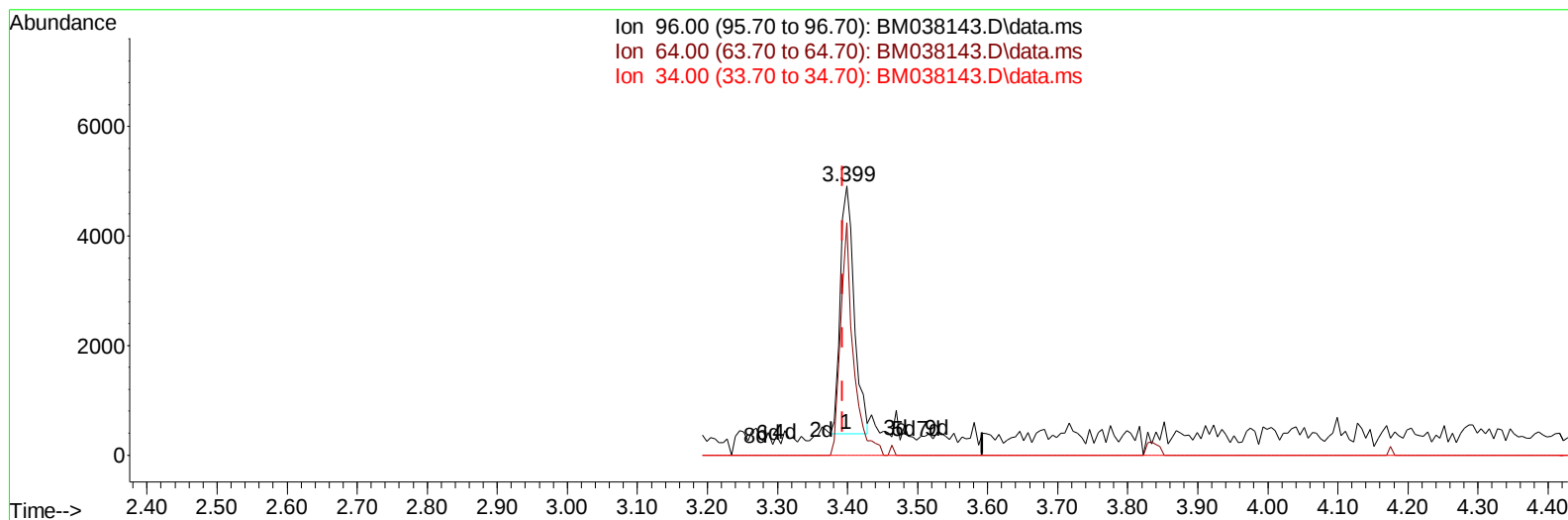
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
 Data File : BM038143.D
 Acq On : 20 Dec 2022 20:27
 Operator : CG/JU
 Sample : N6008-18
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBXY0

Manual IntegrationsAPPROVED

Quant Time: Dec 21 00:16:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 21 00:11:56 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/21/2022
 Supervised By :mohammad ahmed 12/21/2022



TIC: BM038143.D\data.ms

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

3.399min (+ 0.006) 1.86 ng/uL

response 6215

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.40	86.35
34.00	0.00	0.00
0.00	0.00	0.00

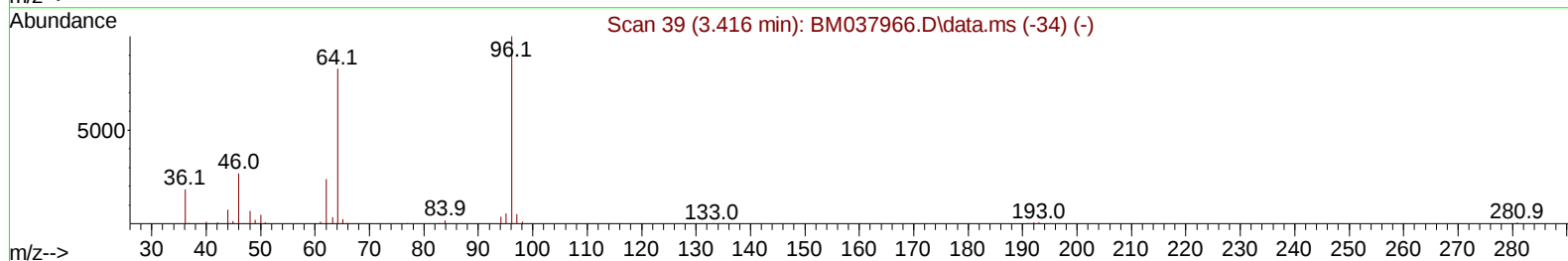
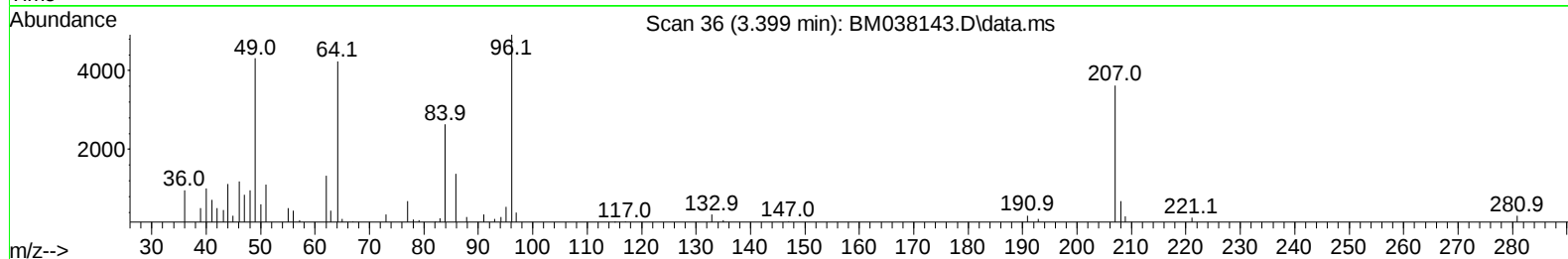
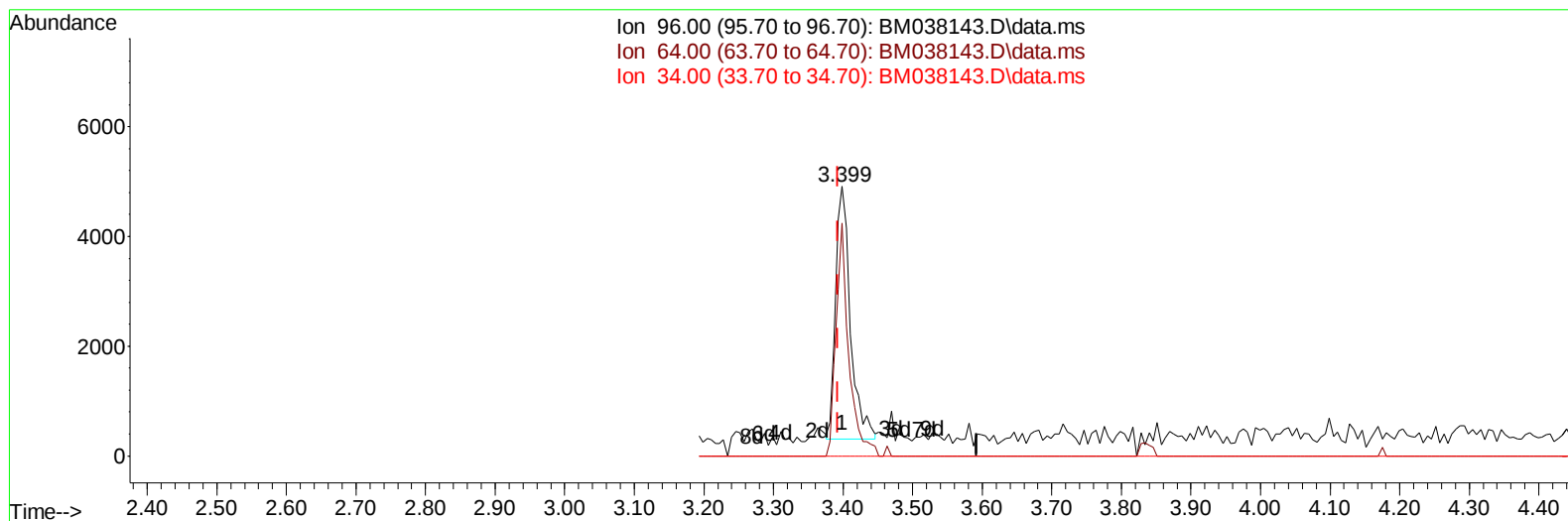
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TIC: BM038143.D\data.ms

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

3.399min (+ 0.006) 2.02 ng/uL m

response 6760

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	76.40	86.35
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.046	152	151777	20.000	ng/ul	0.00
20) Naphthalene-d8	10.863	136	663095	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.675	164	418506	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.410	188	905121	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240	752580	20.000	ng/ul	0.00
88) Perylene-d12	24.027	264	679249	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	6760m	2.022	ng/uL	0.00
4) Pyridine-d5	3.834	84	147583	14.883	ng/ul	0.00
7) Phenol-d5	7.204	99	344784	27.955	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.369	67	160243	21.337	ng/ul	0.00
11) 2-Chlorophenol-d4	7.575	132	263424	26.036	ng/ul	0.00
15) 4-Methylphenol-d8	8.757	113	304533	31.632	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	148753	28.333	ng/ul	0.00
24) 2-Nitrophenol-d4	9.940	143	179862	32.691	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	358080	34.265	ng/ul	0.00
31) 4-Chloroaniline-d4	11.004	131	309481	21.454	ng/ul	0.00
46) Dimethylphthalate-d6	14.080	166	1149198	38.516	ng/ul	0.00
49) Acenaphthylene-d8	14.369	160	1334519	36.596	ng/ul	0.00
54) 4-Nitrophenol-d4	14.869	143	189247	37.217	ng/ul	0.00
60) Fluorene-d10	15.663	176	1017776	38.288	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.780	200	163683	33.825	ng/ul	0.00
73) Anthracene-d10	17.510	188	1579683	37.395	ng/ul	0.00
81) Pyrene-d10	19.792	212	1812226	40.087	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	1304148	38.440	ng/ul	0.00
Target Compounds						
					Qvalue	
80) Fluoranthene	19.457	202	57893	1.085	ng/ul	99
90) Benzo(b)fluoranthene	23.274	252	73969	1.757	ng/ul#	95
93) Benzo(a)pyrene	23.915	252	44642	1.206	ng/ul#	91

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

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