

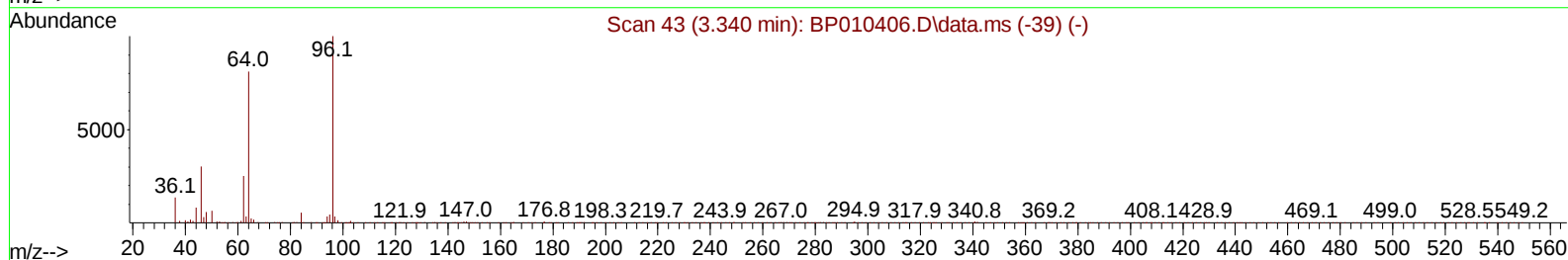
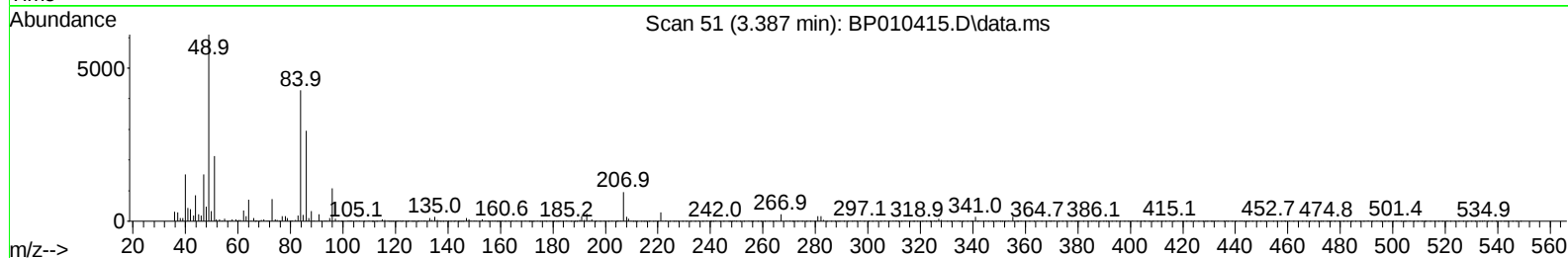
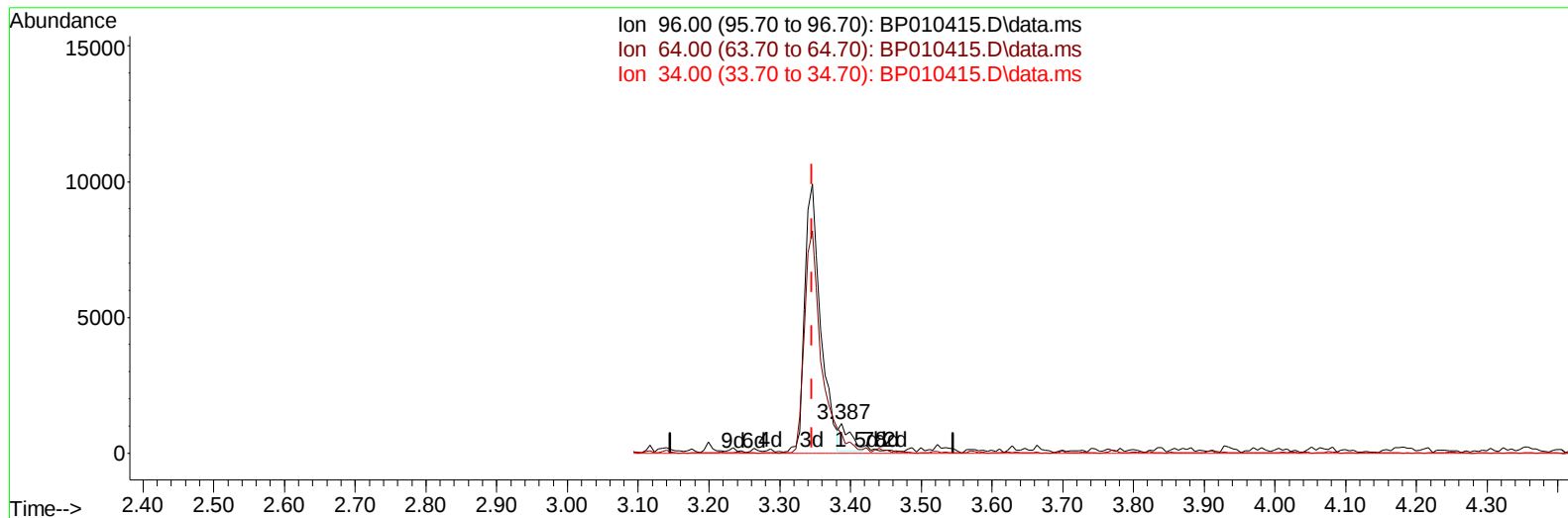
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051922\
 Data File : BP010415.D
 Acq On : 19 May 2022 22:20
 Operator : CG/JU
 Sample : N2918-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTE3

Manual IntegrationsAPPROVED

Quant Time: May 20 00:46:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/20/2022
 Supervised By :Yogesh Patel 05/24/2022



TIC: BP010415.D\data.ms

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

3.387min (+ 0.041) 0.36 ng/uL

response 1122

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	65.31
34.00	0.00	0.00
0.00	0.00	0.00

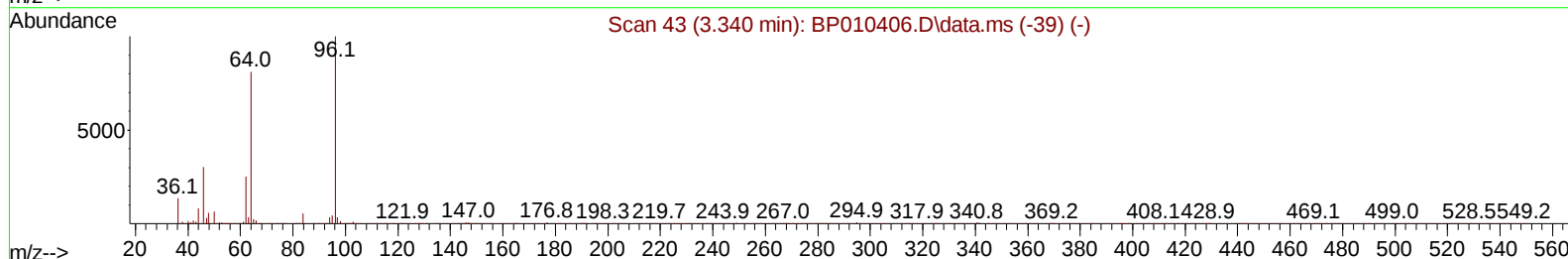
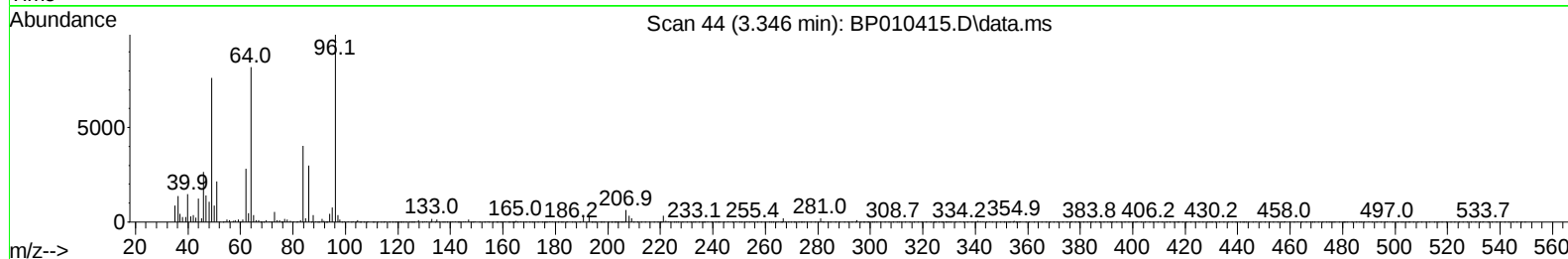
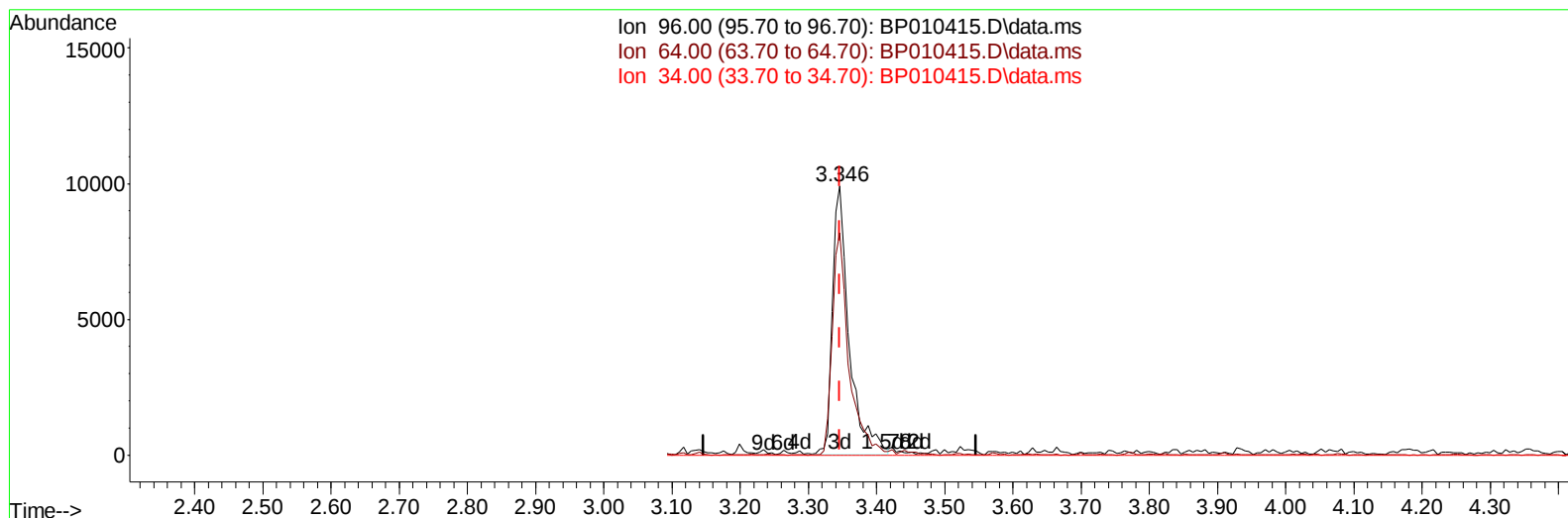
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TIC: BP010415.D\data.ms

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

3.346min (+ 0.000) 5.35 ng/uL m

response 16839

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	82.70#
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	7.893	152	134035	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	596297	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.528	164	404714	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	881932	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	866301	20.000	ng/ul	# 0.00
88) Perylene-d12	23.786	264	749153	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	16839m	5.355	ng/uL	0.00
4) Pyridine-d5	3.770	84	48269	5.406	ng/ul	0.00
7) Phenol-d5	7.058	99	72552	6.449	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	266915	35.314	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	233353	27.108	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	158651	16.594	ng/ul	-0.01
21) Nitrobenzene-d5	9.052	128	165961	35.962	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	173398	35.454	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	282885	30.590	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	345592	25.134	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1136232	38.106	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1265068	36.781	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	27720	5.223	ng/ul	0.00
60) Fluorene-d10	15.516	176	1003960	38.701	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	197866	38.362	ng/ul	0.00
73) Anthracene-d10	17.375	188	1740538	43.584	ng/ul	0.00
81) Pyrene-d10	19.610	212	2103336	45.718	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.627	264	1694089	45.034	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.746	128	1987916	63.827	ng/ul	97
36) 2-Methylnaphthalene	12.352	142	147433	6.687	ng/ul	98
37) 1-Methylnaphthalene	12.569	142	104543	4.696	ng/ul#	97
52) Acenaphthene	14.587	153	137599	5.669	ng/ul	95
56) Dibenzofuran	14.922	168	44564	1.253	ng/ul	96
61) Fluorene	15.575	166	31623	1.087	ng/ul	91
77) Carbazole	17.681	167	184826	4.454	ng/ul#	95

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

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