

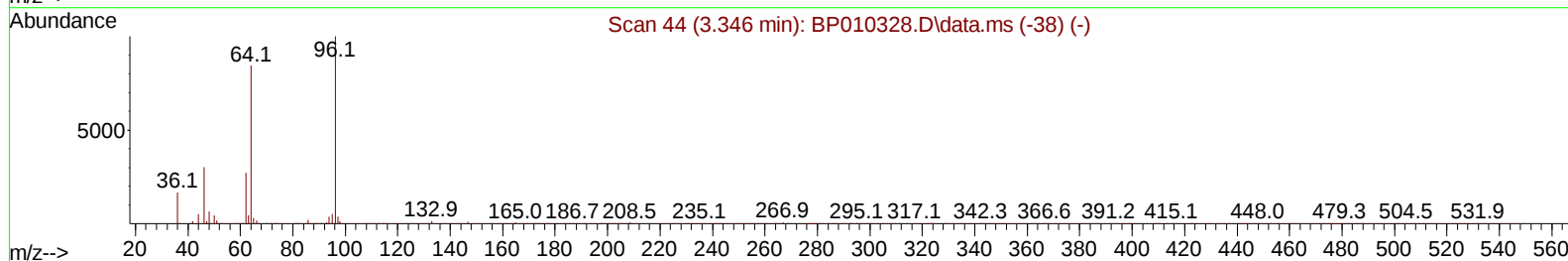
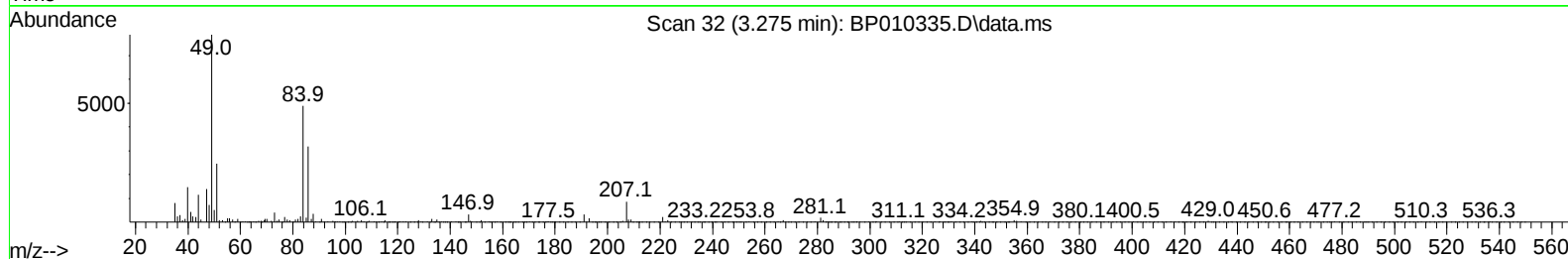
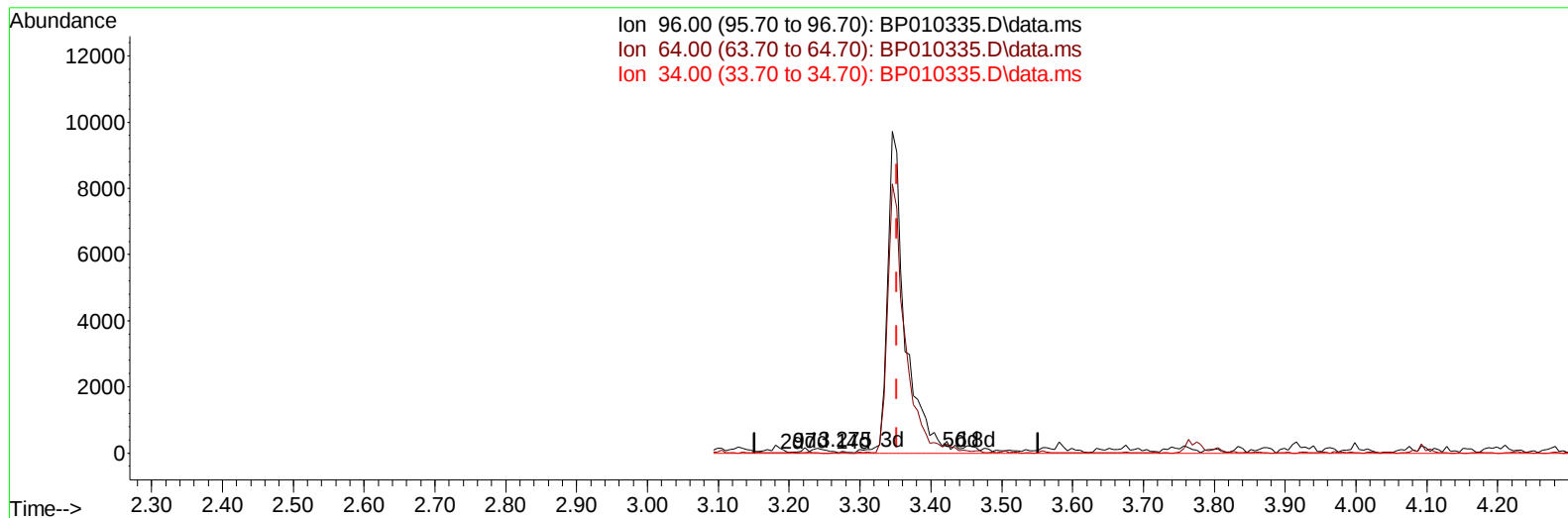
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010335.D
 Acq On : 12 May 2022 21:57
 Operator : CG/JU
 Sample : N2746-03
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBSK6

Manual IntegrationsAPPROVED

Quant Time: May 13 00:03:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/16/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010335.D\data.ms

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

3.275min (-0.077) 0.01 ng/uL

response 30

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	51.16
34.00	0.00	0.00
0.00	0.00	0.00

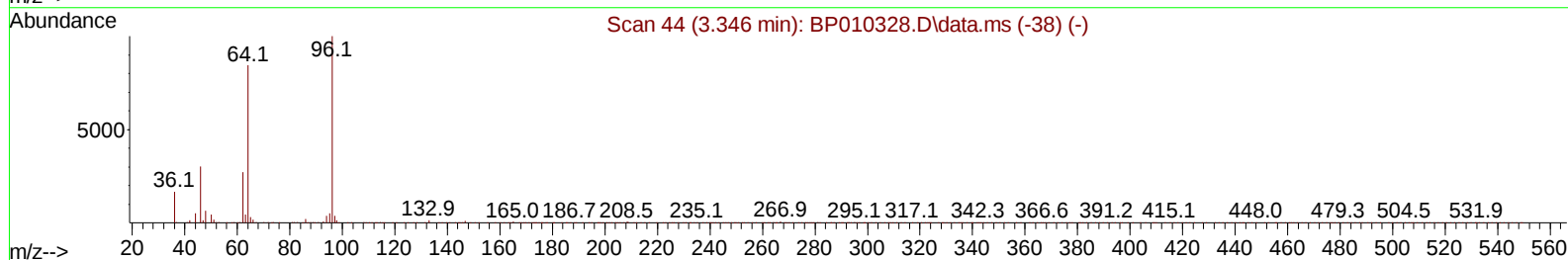
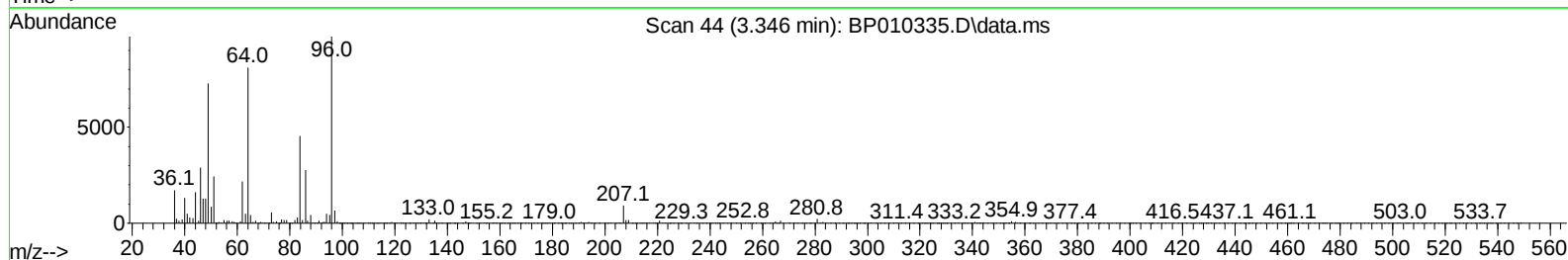
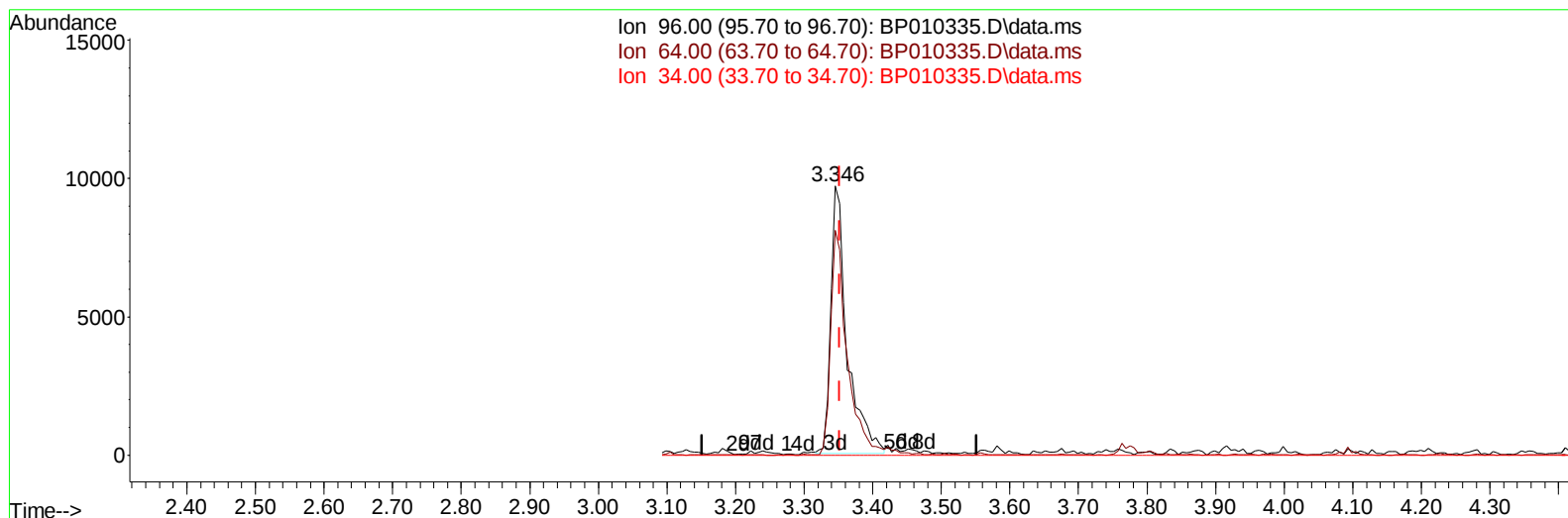
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TIC: BP010335.D\data.ms

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

3.346min (-0.006) 5.39 ng/uL m

response 15977

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.55#
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	7.904	152	125939	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	568100	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.533	164	406749	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	902937	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.368	240	926088	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	864495	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	15977m	5.389	ng/uL	0.00
4) Pyridine-d5	3.763	84	227157	27.535	ng/ul	0.00
7) Phenol-d5	7.069	99	309913	26.495	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	219562	30.343	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	249921	29.011	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	269661	28.041	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	135802	29.813	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	150638	29.005	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	255891	27.070	ng/ul	0.00
31) 4-Chloroaniline-d4	10.839	131	381729	27.354	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	964656	29.769	ng/ul	-0.01
49) Acenaphthylene-d8	14.227	160	1080191	27.959	ng/ul	0.00
54) 4-Nitrophenol-d4	14.727	143	134561	21.649	ng/ul	-0.01
60) Fluorene-d10	15.527	176	830654	30.368	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	123433	19.963	ng/ul	-0.01
73) Anthracene-d10	17.380	188	1310566	29.639	ng/ul	-0.01
81) Pyrene-d10	19.615	212	1617467	31.606	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1431658	31.053	ng/ul	-0.02

Target Compounds Qvalue

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

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