

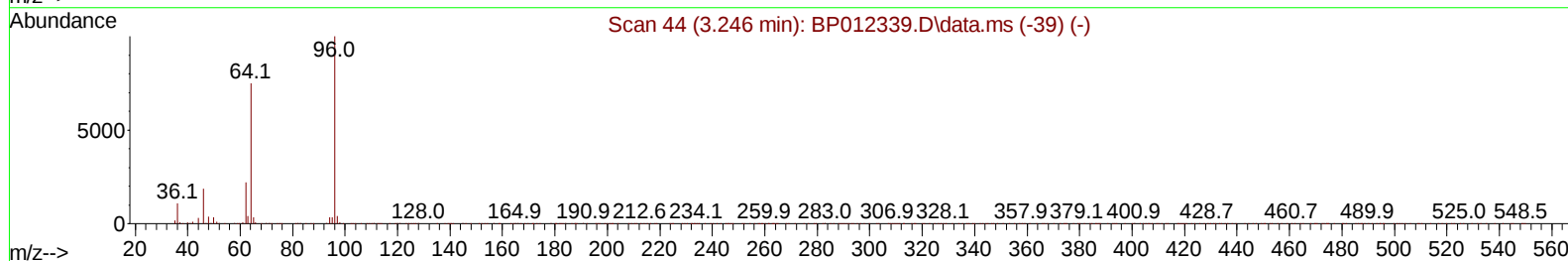
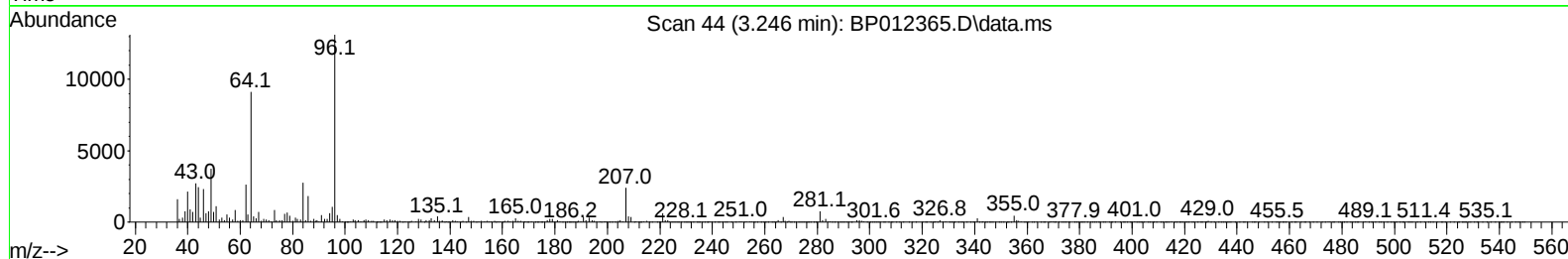
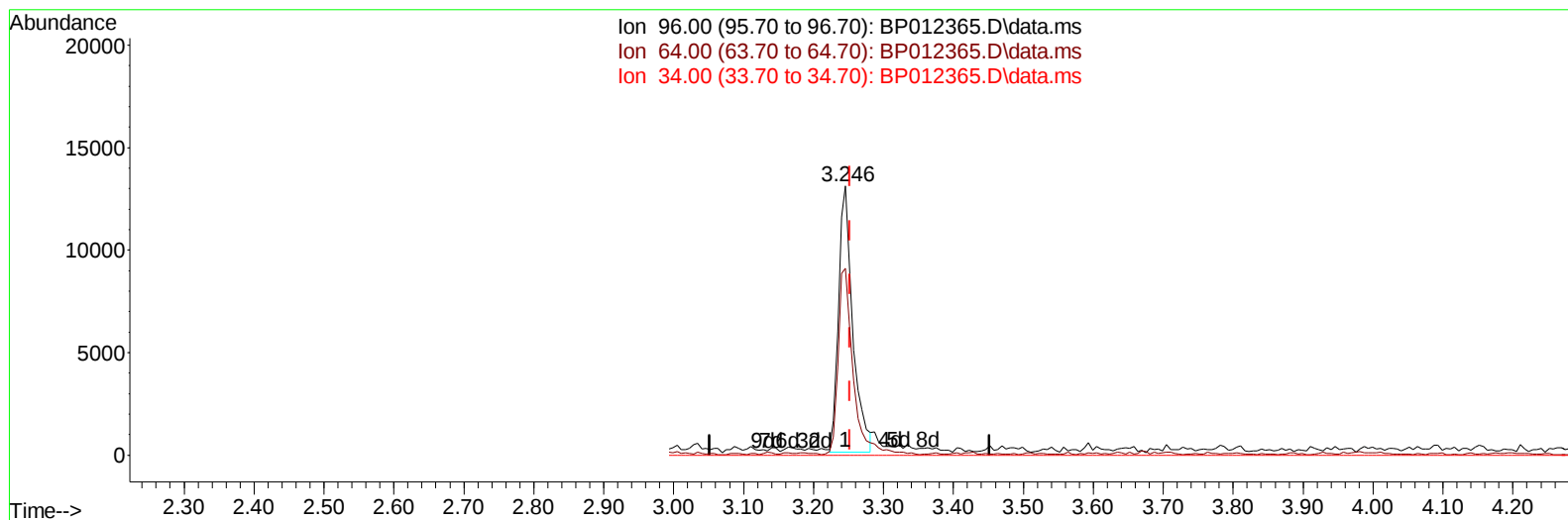
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012365.D
Acq On : 28 Oct 2022 04:03
Operator : CG/JU
Sample : N5136-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8Z7

Manual IntegrationsAPPROVED

Quant Time: Oct 28 05:03:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 26 23:40:20 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/28/2022
Supervised By :mohammad ahmed 10/31/2022



TIC: BP012365.D\data.ms

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

3.246min (-0.006) 3.10 ng/uL

response 18599

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	71.90	69.36
34.00	0.00	0.00
0.00	0.00	0.00

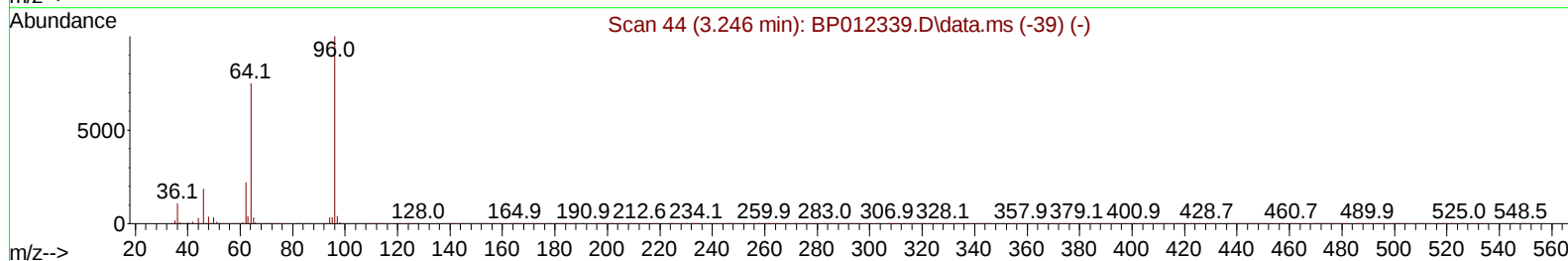
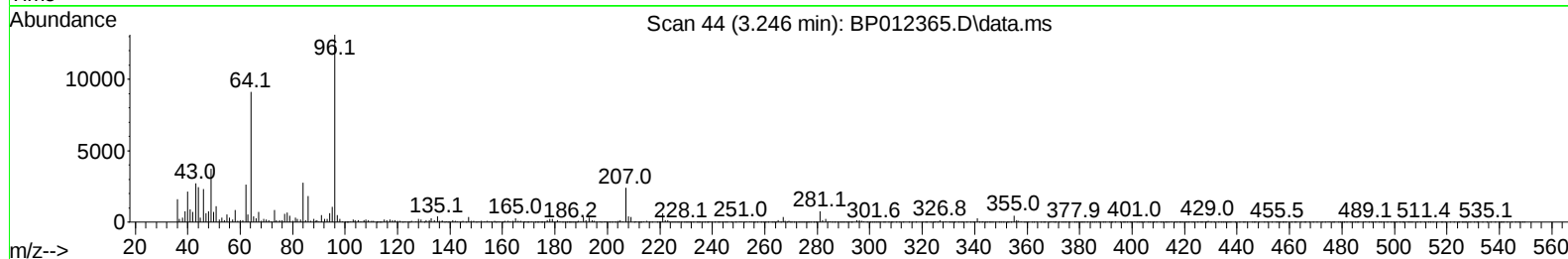
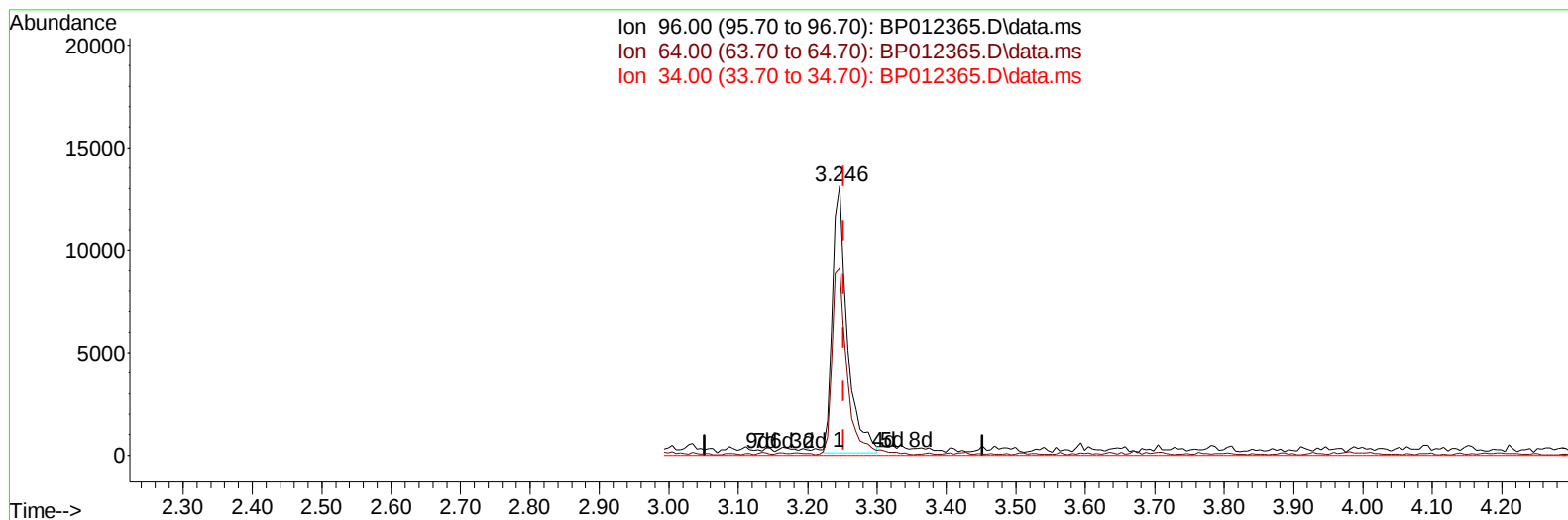
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TIC: BP012365.D\data.ms

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

3.246min (-0.006) 3.24 ng/uL m

response 19417

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	71.90	69.36
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	7.805	152	241203	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.604	136	1102897	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.457	164	785503	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	1819804	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	2079458	20.000	ng/ul	0.00
88) Perylene-d12	23.704	264	2028266	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	19417m	3.241	ng/uL	0.00
4) Pyridine-d5	3.669	84	106489	6.188	ng/ul	0.00
7) Phenol-d5	6.981	99	59611	2.755	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	264983	20.518	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.334	132	160887	10.184	ng/ul	-0.01
15) 4-Methylphenol-d8	8.522	113	133435	7.798	ng/ul	-0.01
21) Nitrobenzene-d5	8.975	128	175318	24.513	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.693	143	101321	13.649	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.234	165	204122	12.688	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	332079	13.863	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	1464061	27.227	ng/ul	-0.01
49) Acenaphthylene-d8	14.145	160	1541161	25.080	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.692	143	18720	1.839	ng/ul	0.01
60) Fluorene-d10	15.451	176	1316543	28.597	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.586	200	102914	10.862	ng/ul	0.00
73) Anthracene-d10	17.316	188	2512837	32.047	ng/ul	0.00
81) Pyrene-d10	19.557	212	3158084	30.249	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.551	264	3256121	32.856	ng/ul	0.00

Target Compounds Qvalue

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

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