

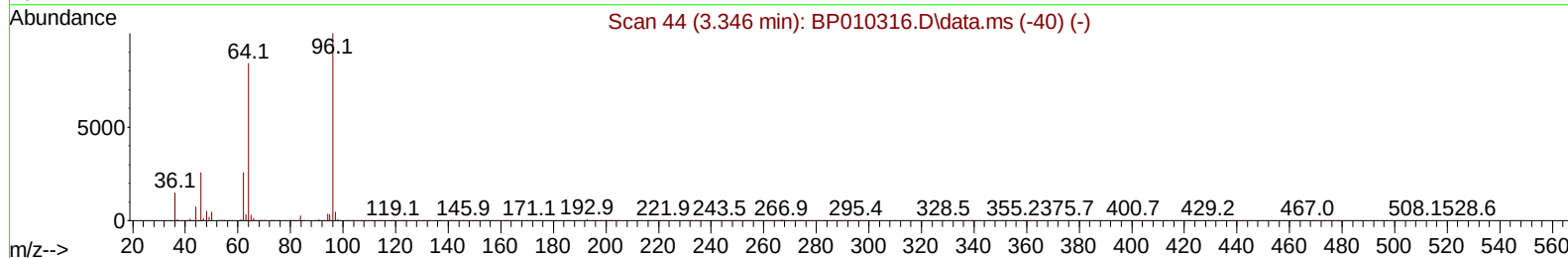
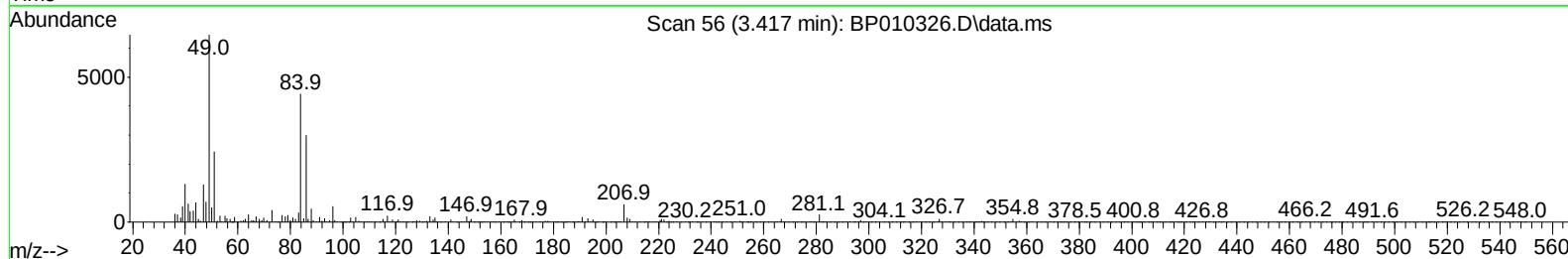
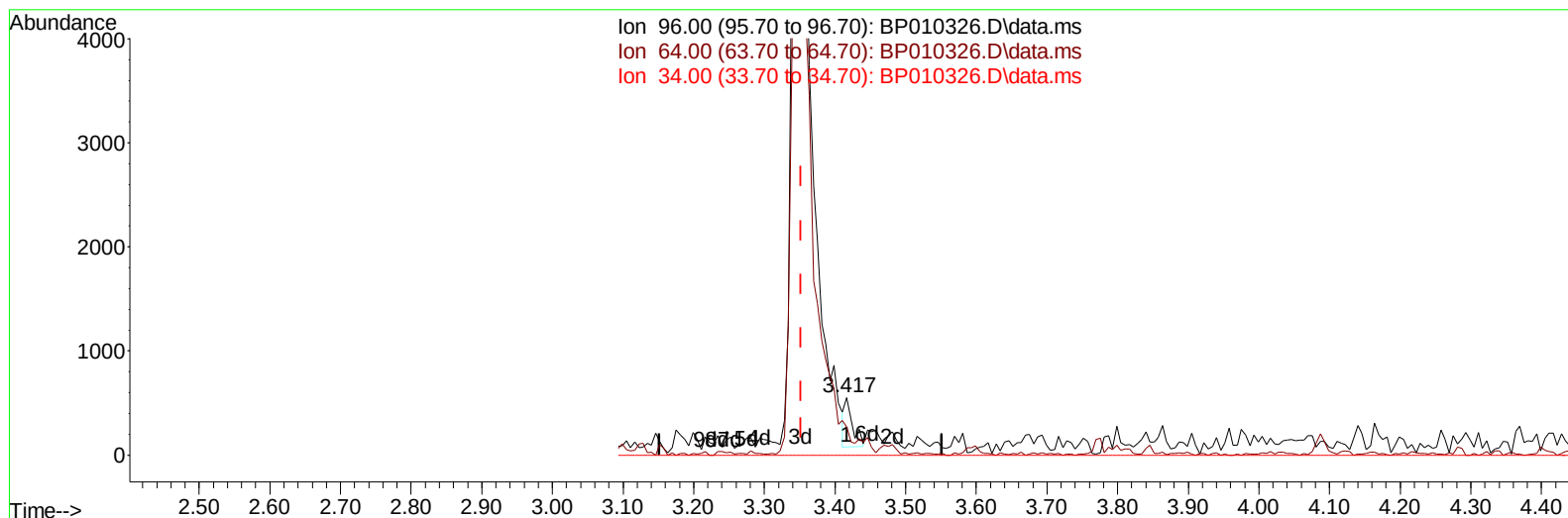
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010326.D
 Acq On : 12 May 2022 16:19
 Operator : CG/JU
 Sample : N2714-06
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW2X7

Manual IntegrationsAPPROVED

Quant Time: May 13 00:02:02 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: BP010326.D\data.ms

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

3.417min (+ 0.065) 0.10 ng/uL

response 352

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	49.46
34.00	0.00	0.00
0.00	0.00	0.00

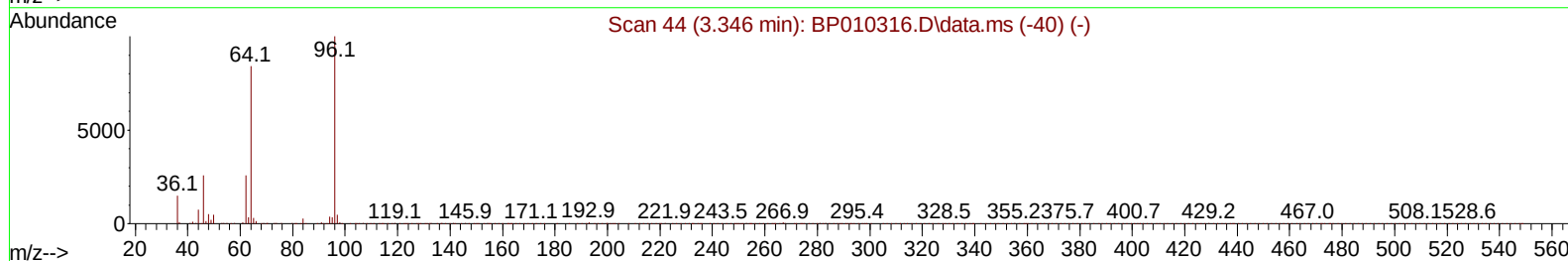
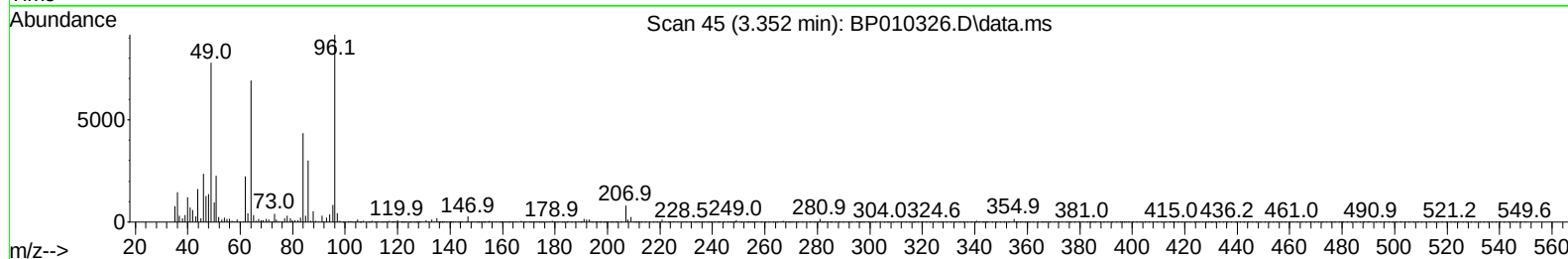
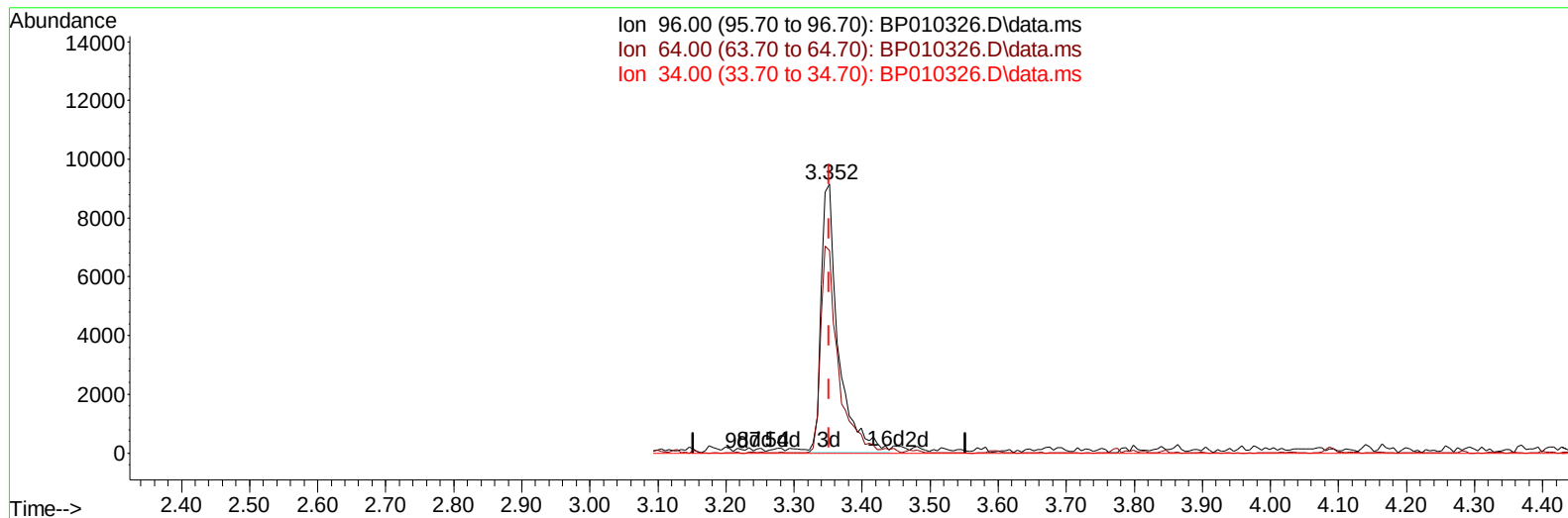
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TIC: BP010326.D\data.ms

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

3.352min (-0.000) 4.34 ng/uL m

response 15934

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	75.44#
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	7.905	152	155827	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	692307	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	497577	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	1088611	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	1078881	20.000	ng/ul	#-0.01
88) Perylene-d12	23.798	264	1034066	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	15934m	4.344	ng/uL	0.00
4) Pyridine-d5	3.775	84	72175	7.071	ng/ul	0.00
7) Phenol-d5	7.069	99	302368	20.892	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	213132	23.805	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	242547	22.755	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	265073	22.277	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	130255	23.465	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	142744	22.554	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	248403	21.563	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	354812	20.863	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	919103	23.186	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1048018	22.175	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	114908	15.112	ng/ul	-0.01
60) Fluorene-d10	15.528	176	793145	23.703	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	112072	15.034	ng/ul	-0.01
73) Anthracene-d10	17.380	188	1248955	23.428	ng/ul	-0.01
81) Pyrene-d10	19.616	212	1518599	25.471	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1322969	23.990	ng/ul	-0.02

Target Compounds Qvalue

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

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