

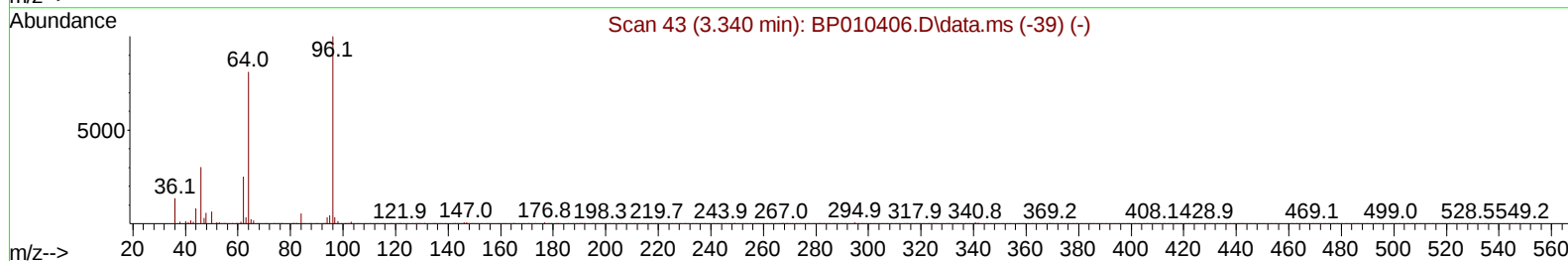
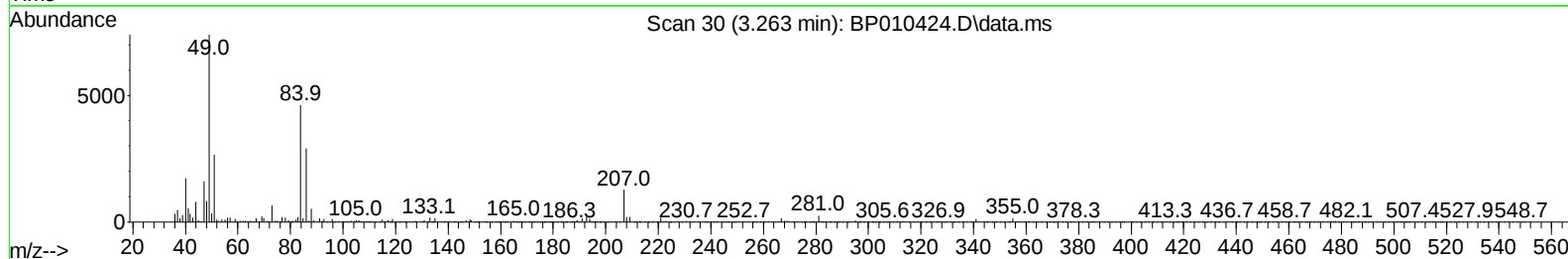
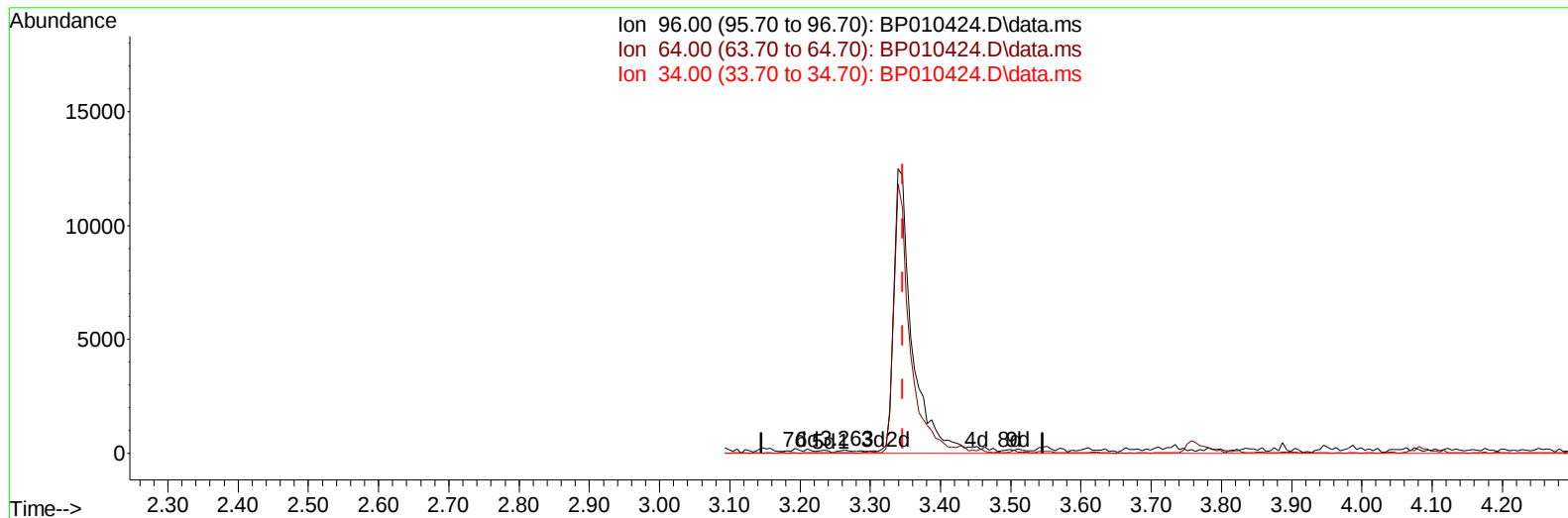
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051922\
 Data File : BP010424.D
 Acq On : 20 May 2022 03:58
 Operator : CG/JU
 Sample : PB144945BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK945

Manual IntegrationsAPPROVED

Quant Time: May 20 05:51:28 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: BP010424.D\data.ms

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

3.263min (-0.083) 0.05 ng/uL

response 163

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	46.40
34.00	0.00	0.00
0.00	0.00	0.00

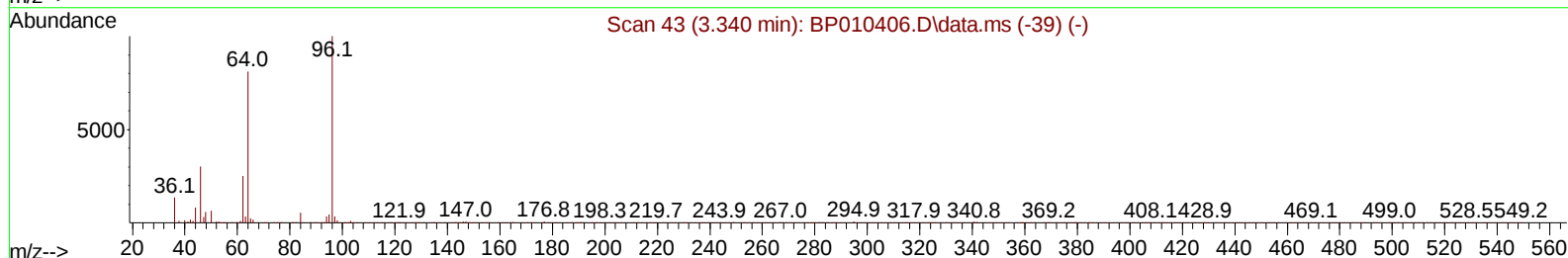
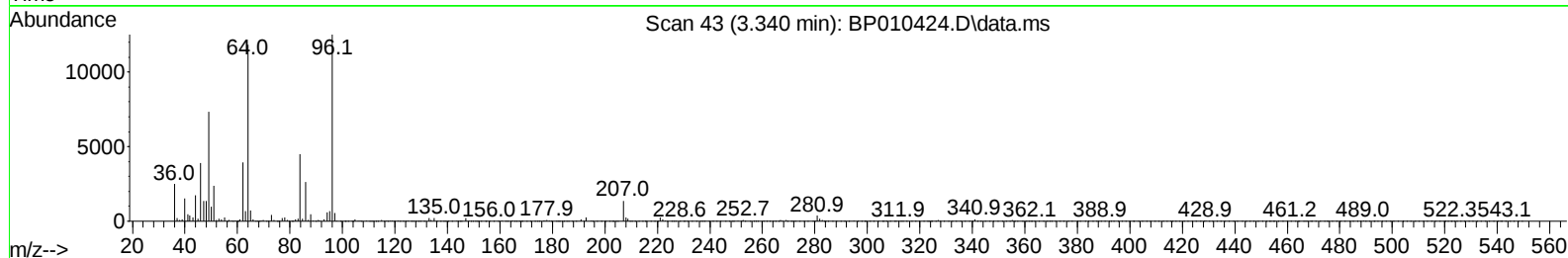
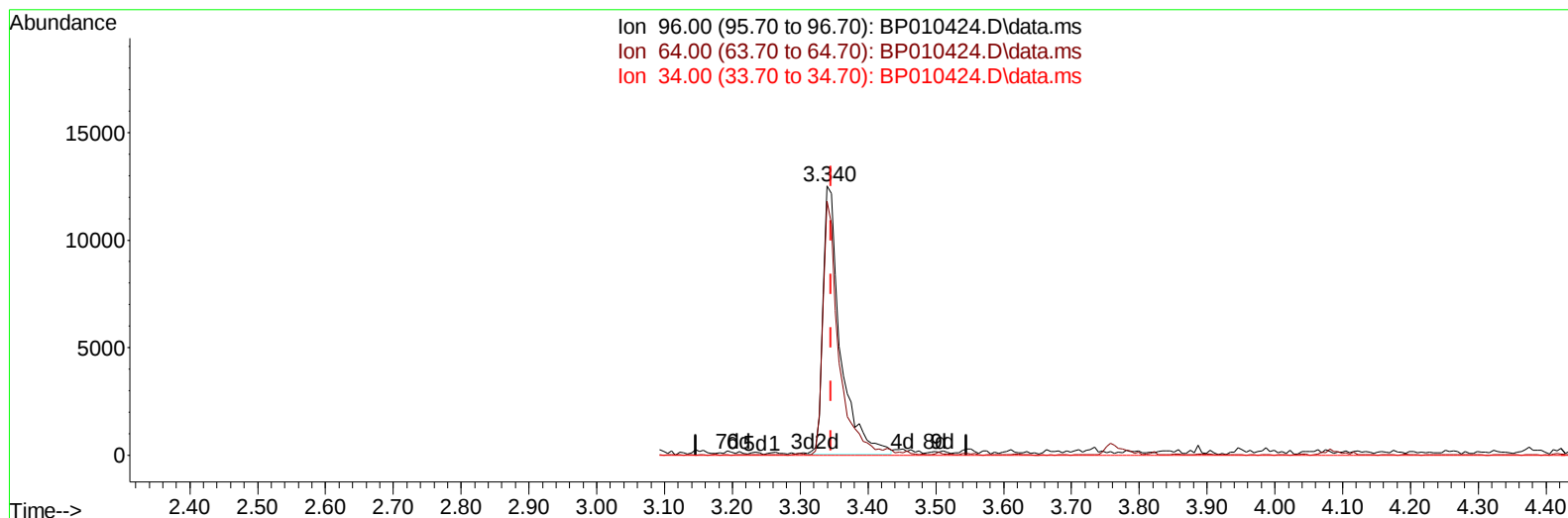
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TIC: BP010424.D\data.ms

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

3.340min (-0.006) 6.30 ng/uL m

response 22227

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	94.47#
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	150357	20.000	ng/ul	0.00
20) Naphthalene-d8	10.692	136	678409	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.522	164	459127	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.274	188	1012758	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	996771	20.000	ng/ul	0.00
88) Perylene-d12	23.786	264	860487	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	22227m	6.301	ng/uL	0.00
4) Pyridine-d5	3.757	84	332924	33.240	ng/ul	0.00
7) Phenol-d5	7.057	99	424101	33.606	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	296005	34.911	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	339927	35.202	ng/ul	-0.01
15) 4-Methylphenol-d8	8.598	113	369740	34.475	ng/ul	-0.01
21) Nitrobenzene-d5	9.051	128	183986	35.042	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	202305	36.358	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	343139	32.614	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	514075	32.862	ng/ul	-0.01
46) Dimethylphthalate-d6	13.933	166	1225915	36.241	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1388405	35.583	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	176033	29.237	ng/ul	0.00
60) Fluorene-d10	15.516	176	1057348	35.929	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	168971	28.528	ng/ul	0.00
73) Anthracene-d10	17.374	188	1698312	37.033	ng/ul	0.00
81) Pyrene-d10	19.610	212	2029725	38.343	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.627	264	1656175	38.330	ng/ul	0.00

Target Compounds Qvalue

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

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