

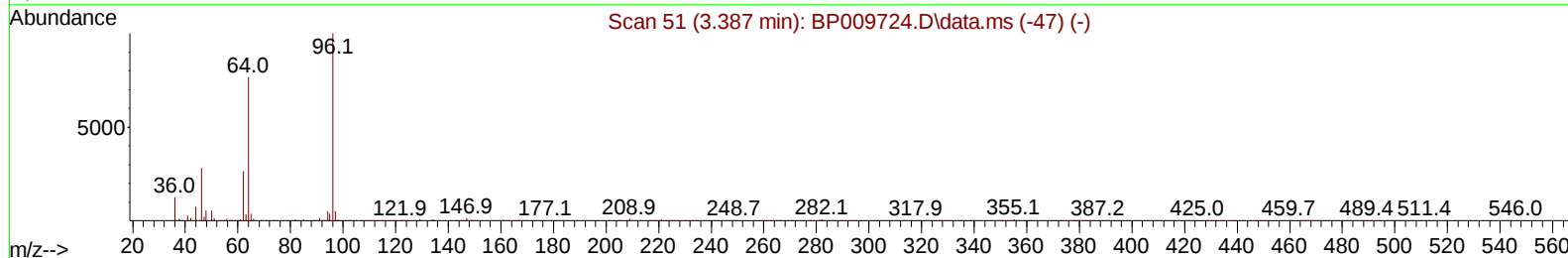
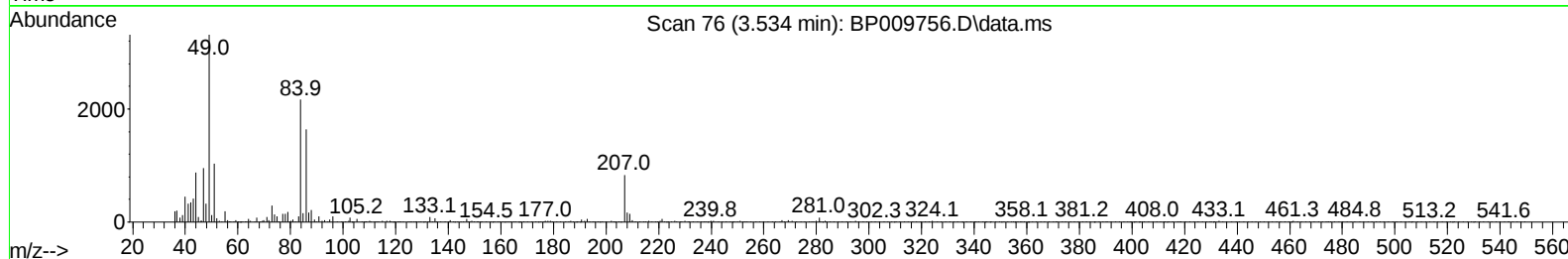
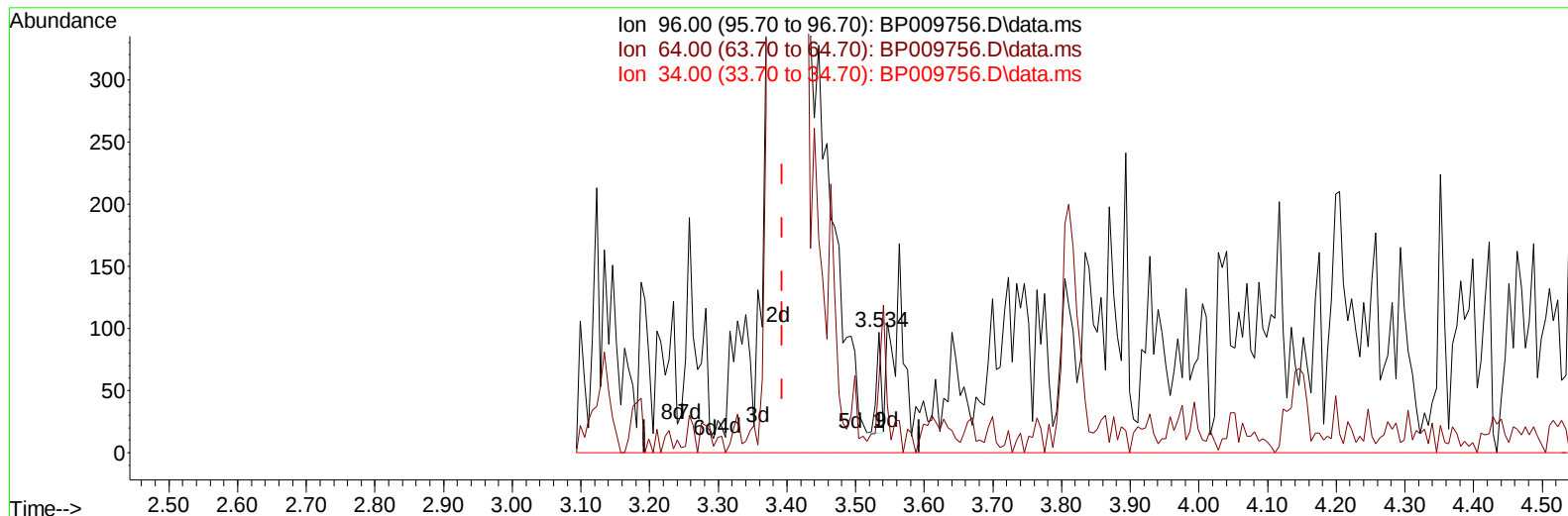
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009756.D
 Acq On : 11 Apr 2022 17:38
 Operator : CG/JU
 Sample : N2266-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YAZ22

Manual IntegrationsAPPROVED

Quant Time: Apr 12 01:49:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009756.D\data.ms

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

3.534min (+ 0.141) 0.02 ng/uL

response 37

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	60.82
34.00	0.00	0.00
0.00	0.00	0.00

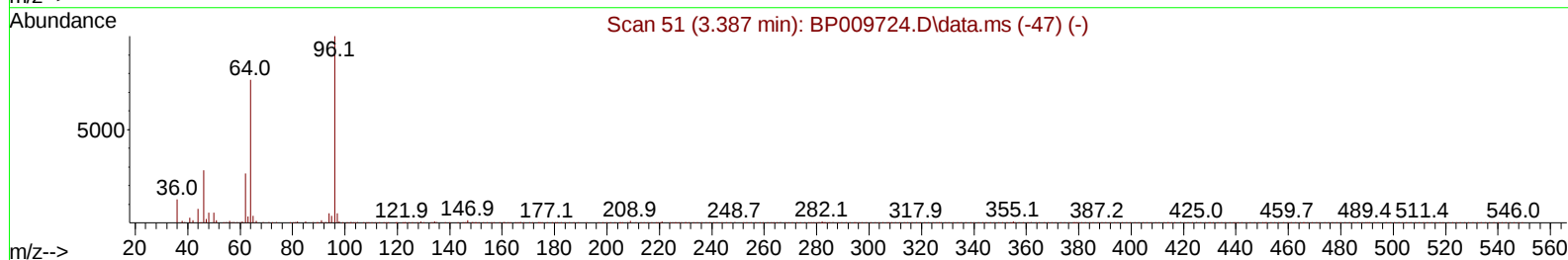
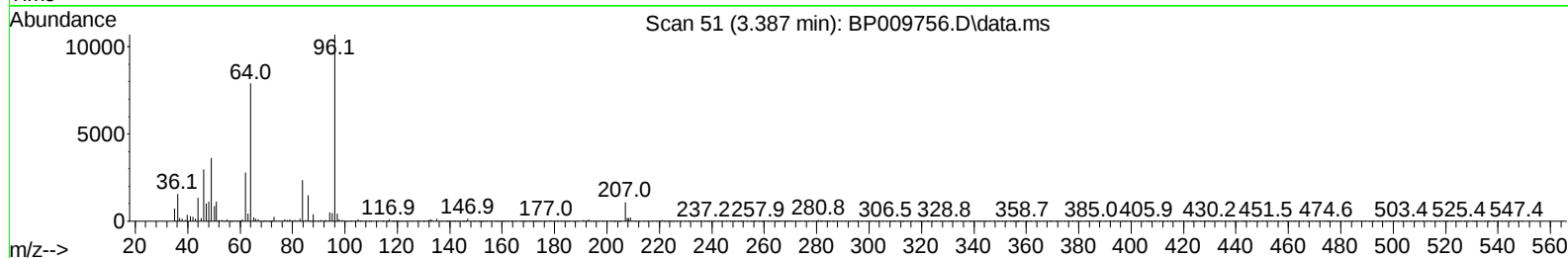
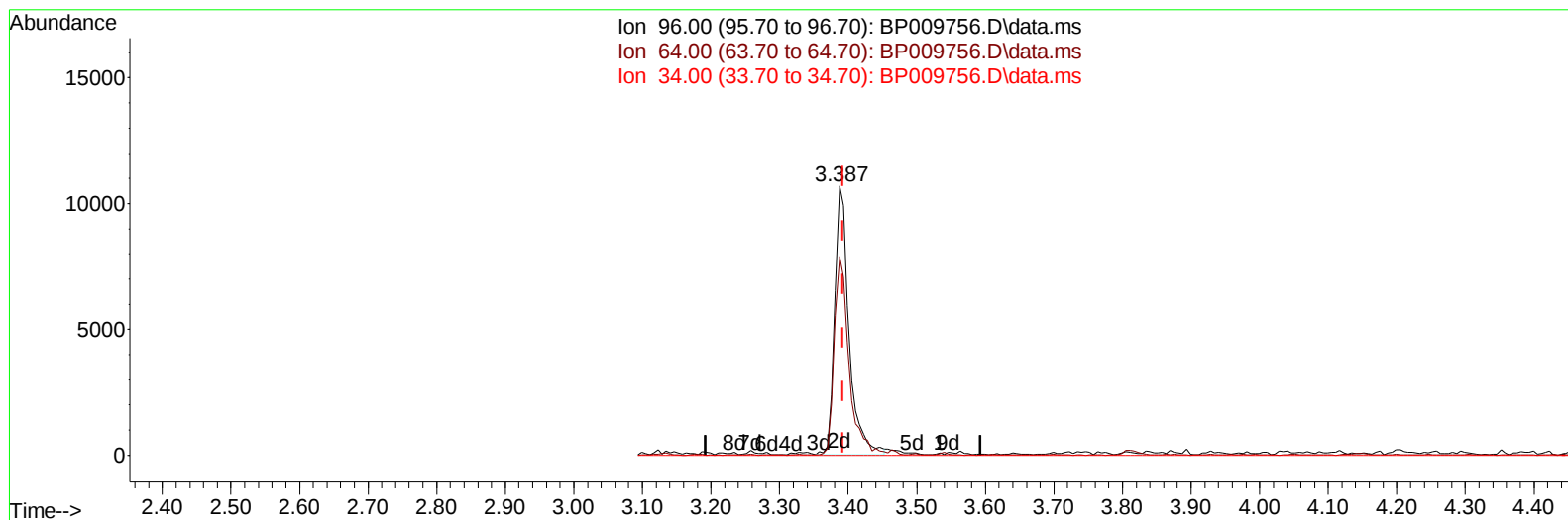
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TIC: BP009756.D\data.ms

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

3.387min (-0.006) 6.75 ng/uL m

response 15624

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.82#
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	7.969	152	103328	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.781	136	465315	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.598	164	336082	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.351	188	755093	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.433	240	860014	20.000	ng/ul	#-0.01
88) Perylene-d12	23.904	264	875801	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	15624m	6.748	ng/uL	0.00
4) Pyridine-d5	3.805	84	85566	13.404	ng/ul	0.00
7) Phenol-d5	7.122	99	81644	9.000	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	224600	41.588	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	218093	31.872	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	161430	21.317	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	147854	44.063	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.852	143	157248	42.744	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	264705	34.385	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.905	131	368106	33.573	ng/ul	-0.01
46) Dimethylphthalate-d6	14.010	166	1159850	43.934	ng/ul	-0.01
49) Acenaphthylene-d8	14.293	160	1236086	39.522	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143	47595	9.941	ng/ul	-0.01
60) Fluorene-d10	15.593	176	990737	44.611	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	183993	41.132	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1701480	47.063	ng/ul	-0.01
81) Pyrene-d10	19.681	212	2128667	45.320	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.745	264	2184337	47.893	ng/ul	-0.02
Target Compounds						
						Qvalue
50) Acenaphthylene	14.322	152	201032	6.513	ng/ul	96
52) Acenaphthene	14.663	153	97454	4.822	ng/ul	97
56) Dibenzofuran	14.998	168	200376	6.737	ng/ul	97
61) Fluorene	15.651	166	122620	5.005	ng/ul	90
74) Anthracene	17.487	178	44895	1.099	ng/ul	95

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

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