

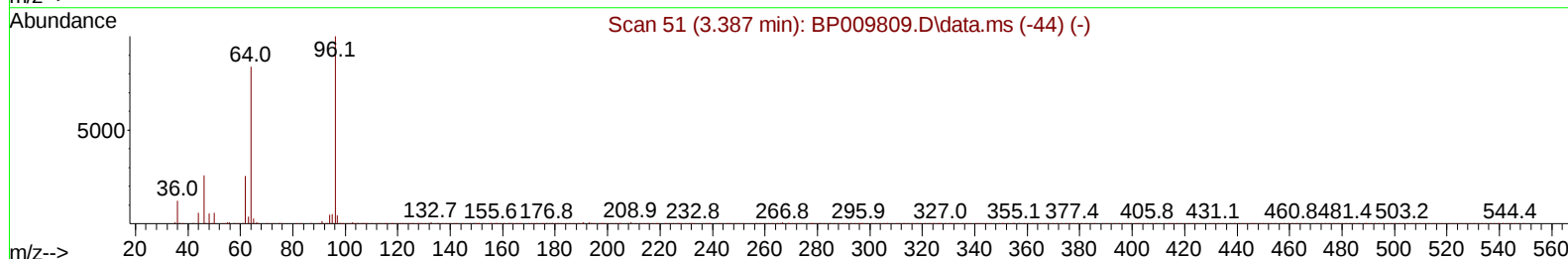
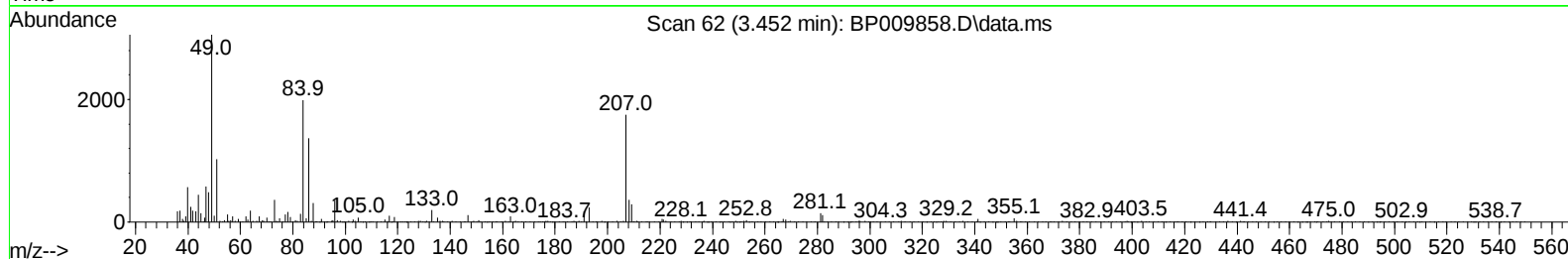
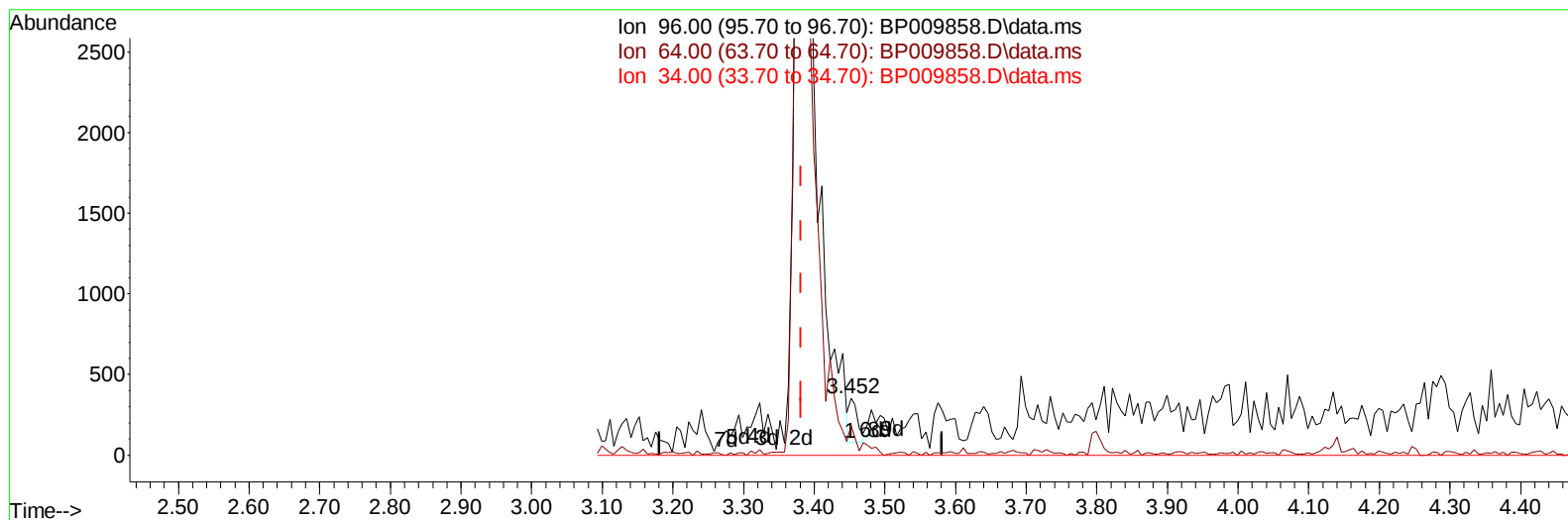
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009858.D
 Acq On : 15 Apr 2022 13:22
 Operator : CG/JU
 Sample : N2422-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J33

Manual IntegrationsAPPROVED

Quant Time: Apr 18 01:00:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/18/2022
 Supervised By :Yogesh Patel 04/18/2022



TIC: BP009858.D\data.ms

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

3.452min (+ 0.071) 0.09 ng/uL

response 228

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	57.51
34.00	0.00	0.00
0.00	0.00	0.00

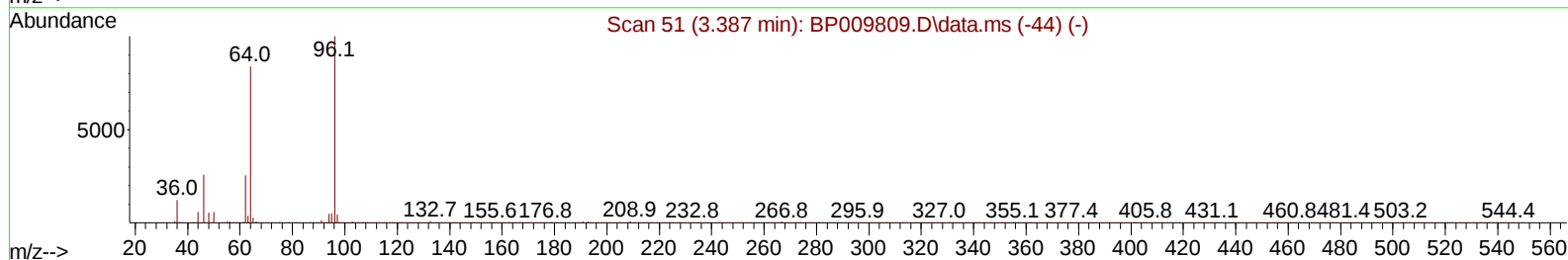
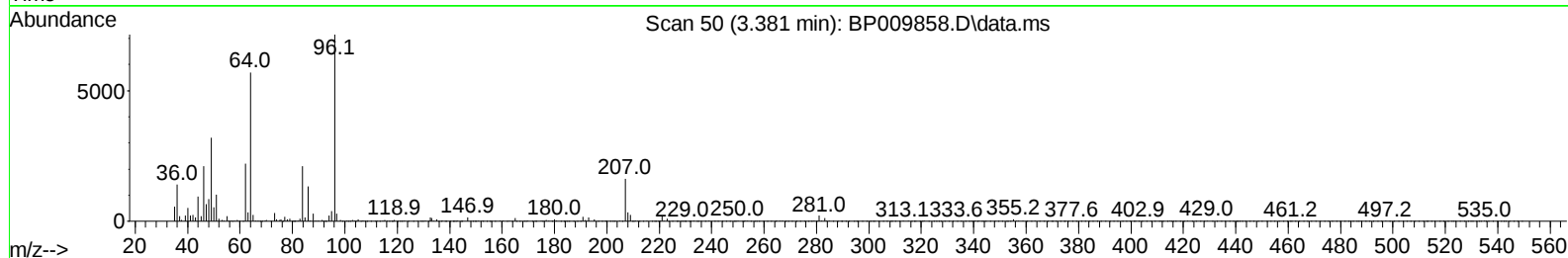
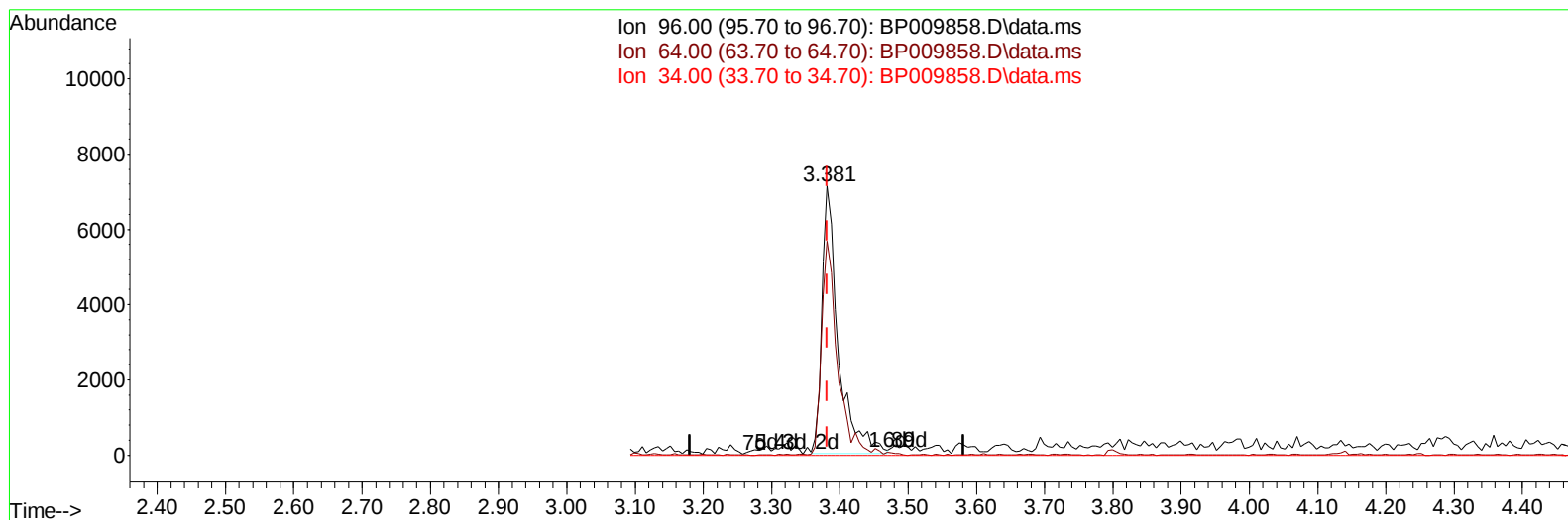
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TIC: BP009858.D\data.ms

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

3.381min (+ 0.000) 4.81 ng/uL m

response 11832

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	79.66#
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	7.958	152	109852	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	501454	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	366782	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	834914	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.422	240	892884	20.000	ng/ul	# 0.00
88) Perylene-d12	23.886	264	867898	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	11832m	4.807	ng/uL	0.00
4) Pyridine-d5	3.805	84	68938	10.158	ng/ul	0.00
7) Phenol-d5	7.111	99	59393	6.158	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	190482	33.176	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	171254	23.541	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	122188	15.177	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	122514	33.880	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	126994	32.032	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	219114	26.412	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	318519	26.957	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	976992	33.910	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1049069	30.735	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	30176	5.775	ng/ul	0.00
60) Fluorene-d10	15.581	176	854680	35.263	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	155050	31.348	ng/ul	0.00
73) Anthracene-d10	17.439	188	1530451	38.285	ng/ul	0.00
81) Pyrene-d10	19.669	212	1898627	38.934	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	1837473	40.655	ng/ul	0.00

Target Compounds Qvalue

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

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