

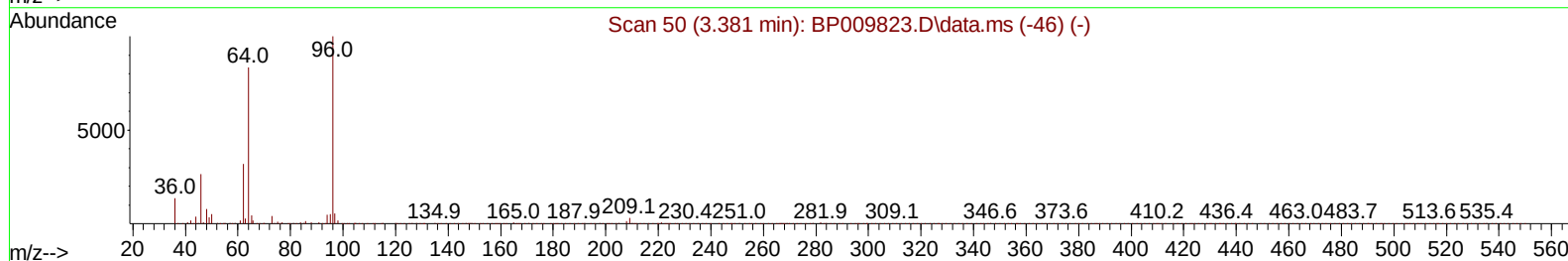
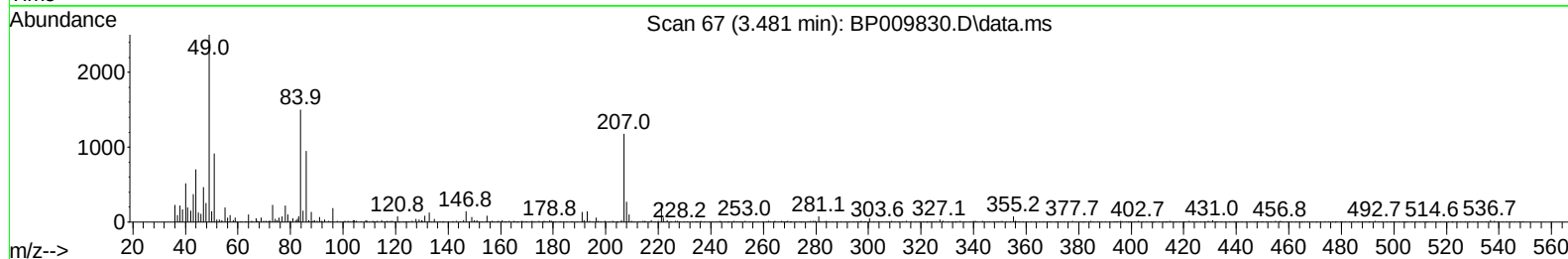
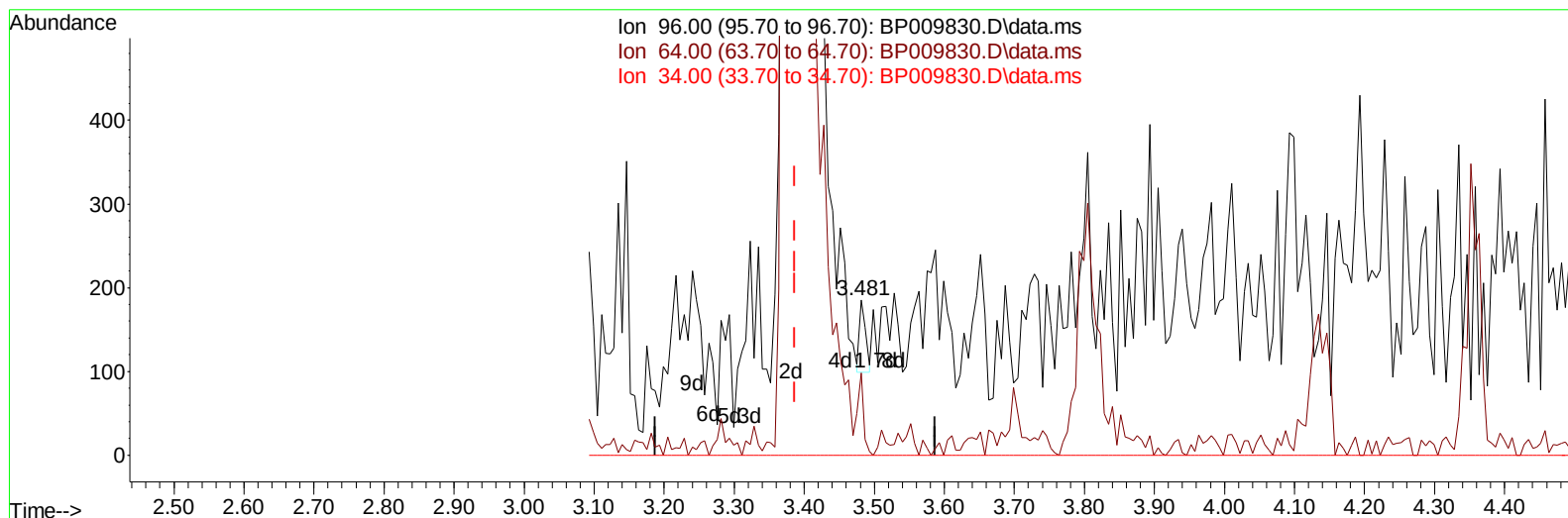
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
 Data File : BP009830.D
 Acq On : 14 Apr 2022 16:23
 Operator : CG/JU
 Sample : N2293-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNN2

Manual IntegrationsAPPROVED

Quant Time: Apr 14 21:46:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 13 02:15:22 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/15/2022
 Supervised By :Yogesh Patel 04/22/2022



TIC: BP009830.D\data.ms

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

3.481min (+ 0.094) 0.02 ng/uL

response 52

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	52.97
34.00	0.00	0.00
0.00	0.00	0.00

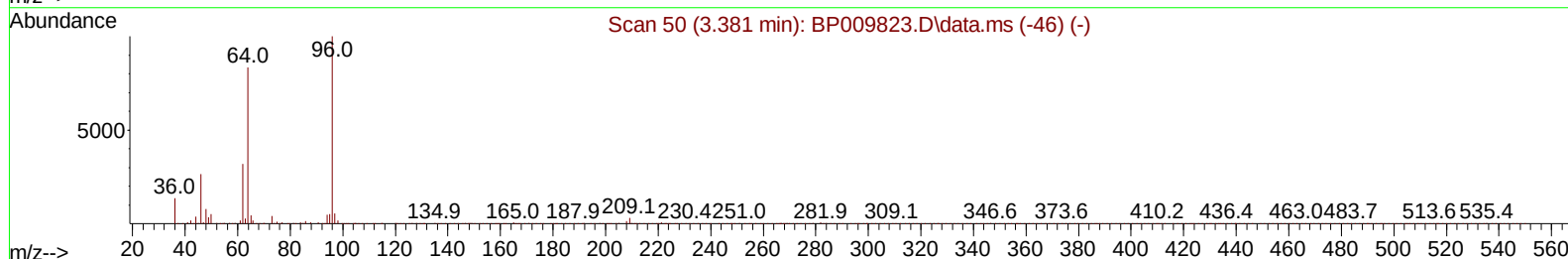
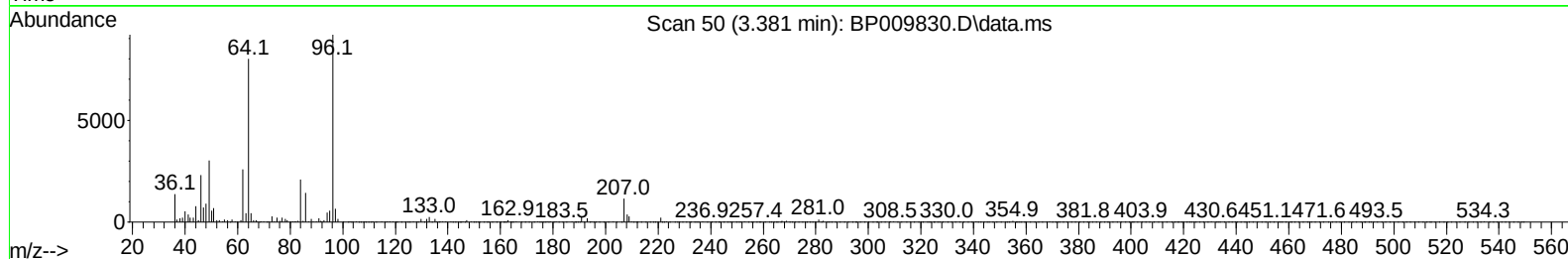
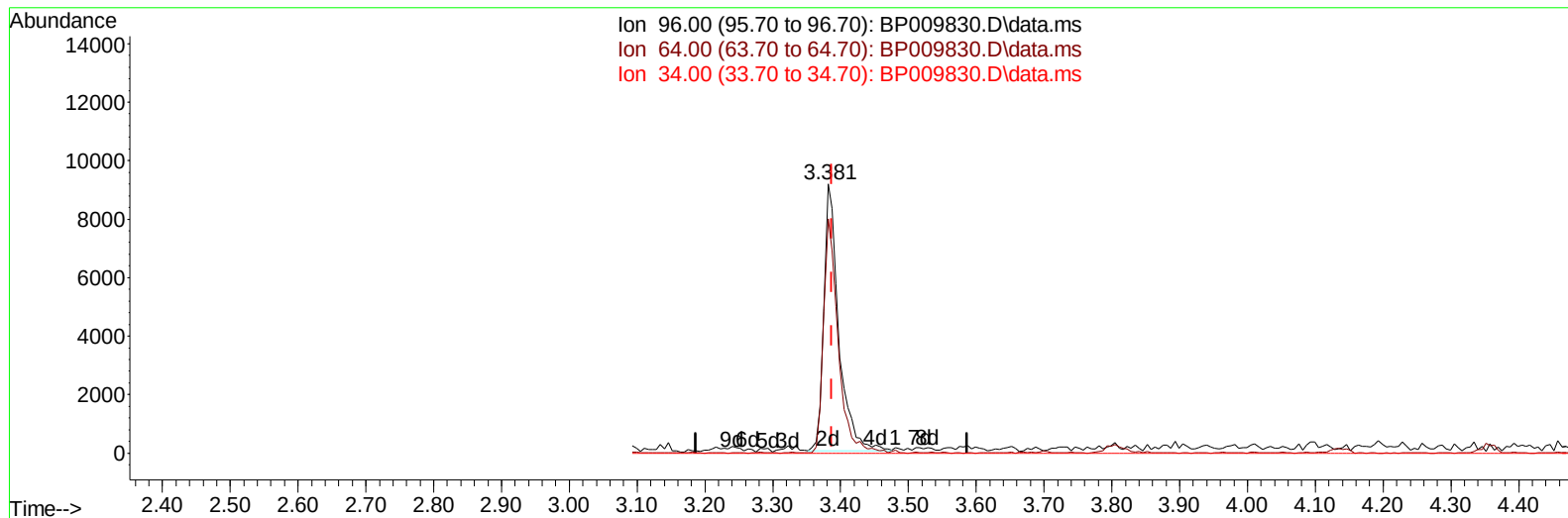
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TIC: BP009830.D\data.ms

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

3.381min (-0.006) 4.42 ng/uL m

response 13980

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	87.23#
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	7.963	152	141053	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	635129	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	464498	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	1046969	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.422	240	1095253	20.000	ng/ul	#-0.01
88) Perylene-d12	23.892	264	1044641	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	13980m	4.423	ng/uL	0.00
4) Pyridine-d5	3.799	84	197306	22.642	ng/ul	0.00
7) Phenol-d5	7.116	99	281839	22.759	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	190091	25.784	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	229057	24.521	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	257952	24.953	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	122647	26.778	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.846	143	133951	26.676	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	249369	23.732	ng/ul	0.00
31) 4-Chloroaniline-d4	10.899	131	340817	22.773	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	935141	25.629	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	1041733	24.099	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	145832	22.039	ng/ul	-0.01
60) Fluorene-d10	15.587	176	793630	25.856	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	122153	19.695	ng/ul	0.00
73) Anthracene-d10	17.445	188	1253231	25.001	ng/ul	0.00
81) Pyrene-d10	19.669	212	1547192	25.865	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	1419889	26.100	ng/ul	0.00

Target Compounds Qvalue

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

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