

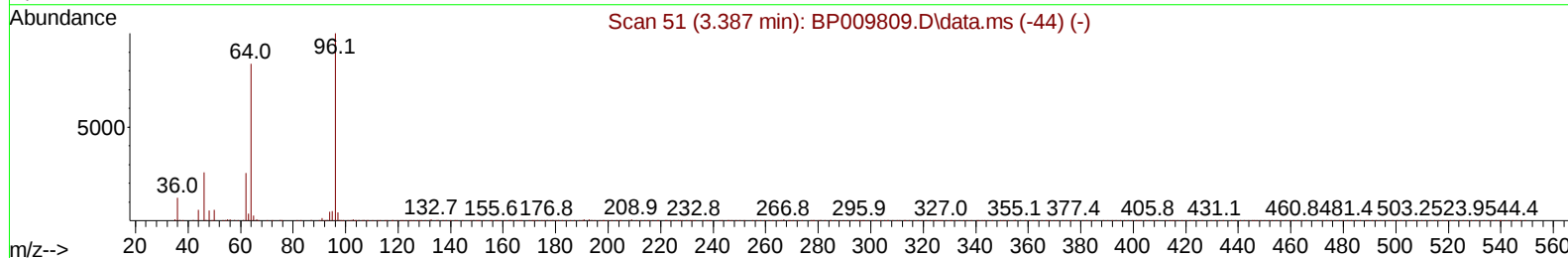
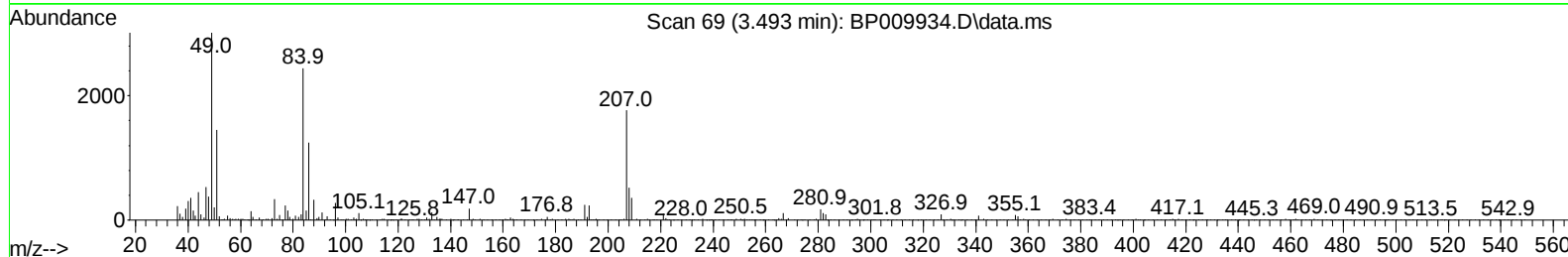
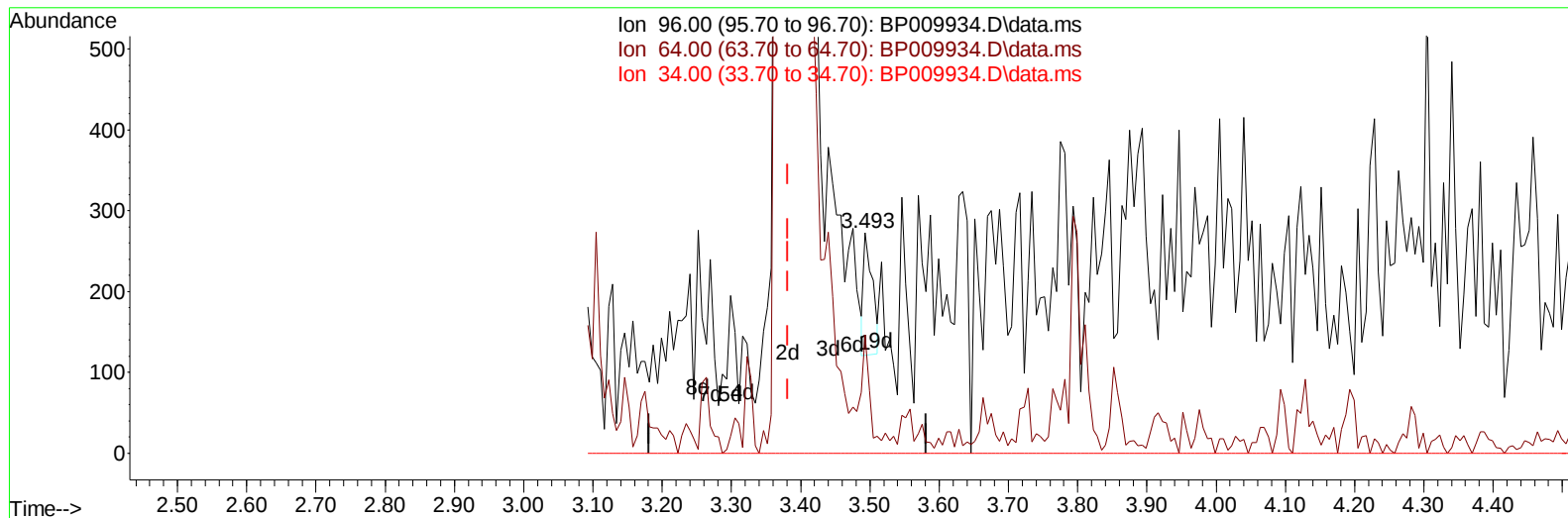
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
Data File : BP009934.D
Acq On : 19 Apr 2022 01:40
Operator : CG/JU
Sample : N2421-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY6

Manual IntegrationsAPPROVED

Quant Time: Apr 19 02:33:58 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 04/19/2022
Supervised By :Yogesh Patel 04/22/2022



TIC: BP009934.D\data.ms

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

3.493min (+ 0.112) 0.04 ng/uL

response 136

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	53.85
34.00	0.00	0.00
0.00	0.00	0.00

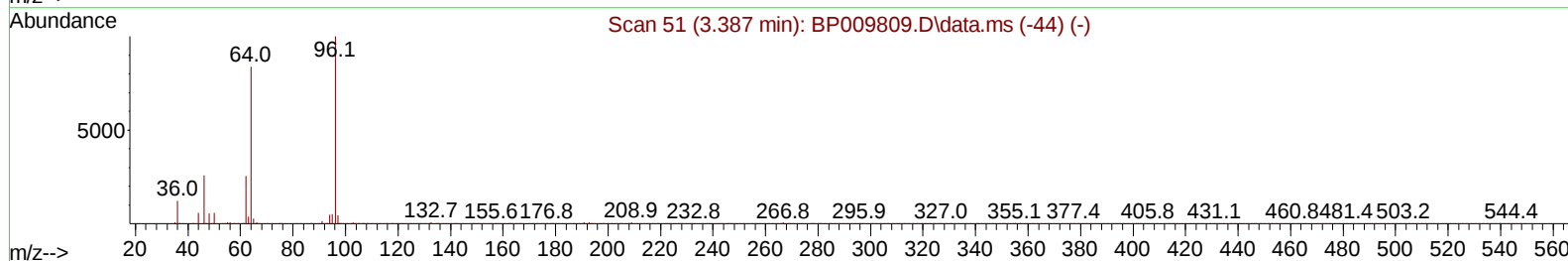
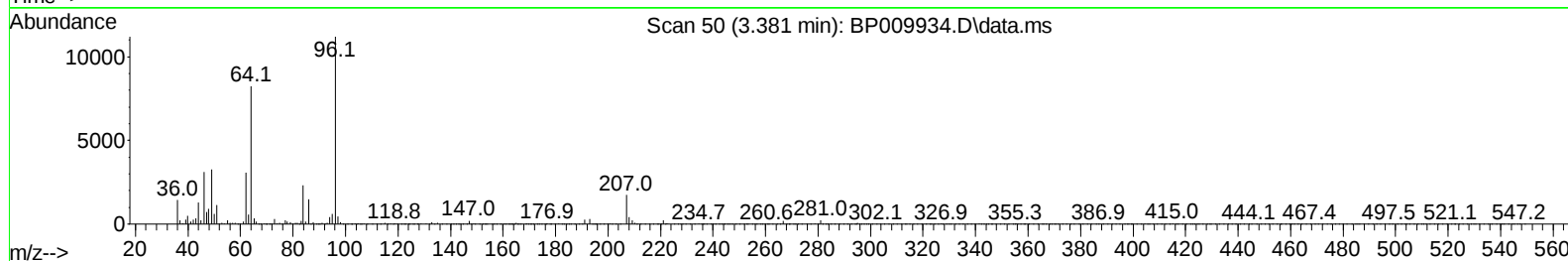
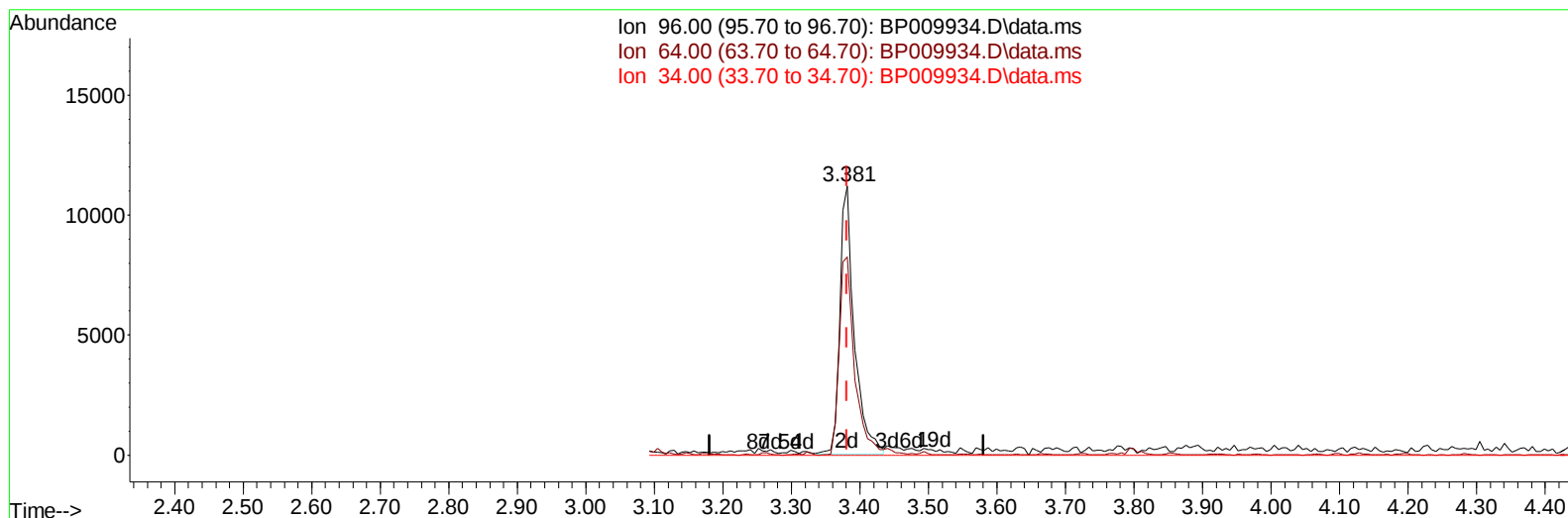
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TIC: BP009934.D\data.ms

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

3.381min (+ 0.000) 5.05 ng/uL m

response 16352

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.58#
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	7.952	152	144609	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.757	136	657299	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.581	164	479356	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.334	188	1060264	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	1060029	20.000	ng/ul	# 0.00
88) Perylene-d12	23.880	264	937113	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	16352m	5.047	ng/uL	0.00
4) Pyridine-d5	3.799	84	95989	10.745	ng/ul	0.00
7) Phenol-d5	7.105	99	104071	8.197	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	206768	27.357	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	217281	22.689	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	189510	17.881	ng/ul	-0.01
21) Nitrobenzene-d5	9.110	128	136526	28.803	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	148680	28.611	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	282791	26.005	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	327413	21.140	ng/ul	0.00
46) Dimethylphthalate-d6	13.987	166	1141028	30.303	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	1187959	26.630	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	57973	8.490	ng/ul	0.00
60) Fluorene-d10	15.575	176	967725	30.551	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	186792	29.739	ng/ul	0.00
73) Anthracene-d10	17.434	188	1678461	33.064	ng/ul	0.00
81) Pyrene-d10	19.663	212	2092332	36.141	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1760938	36.084	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.134	94	17783	1.399	ng/ul#	92

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

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