

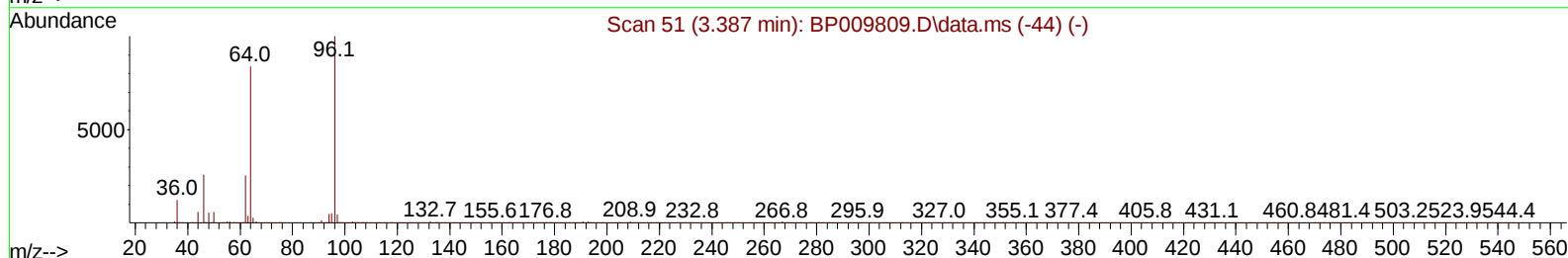
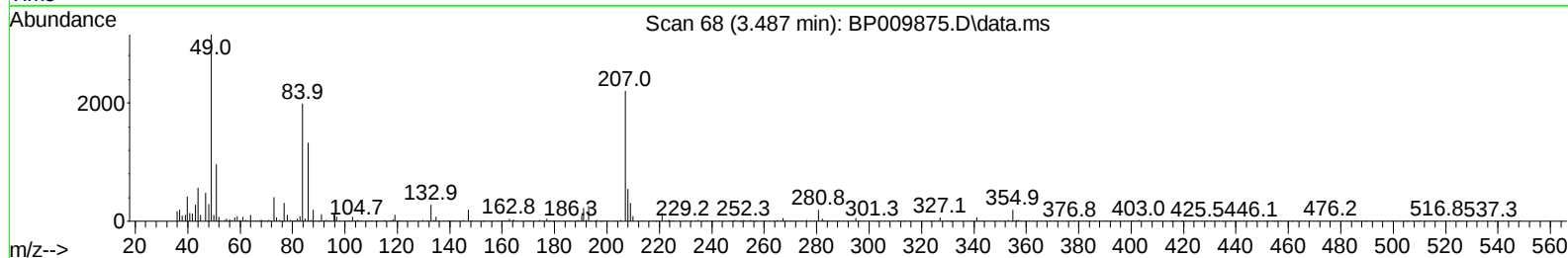
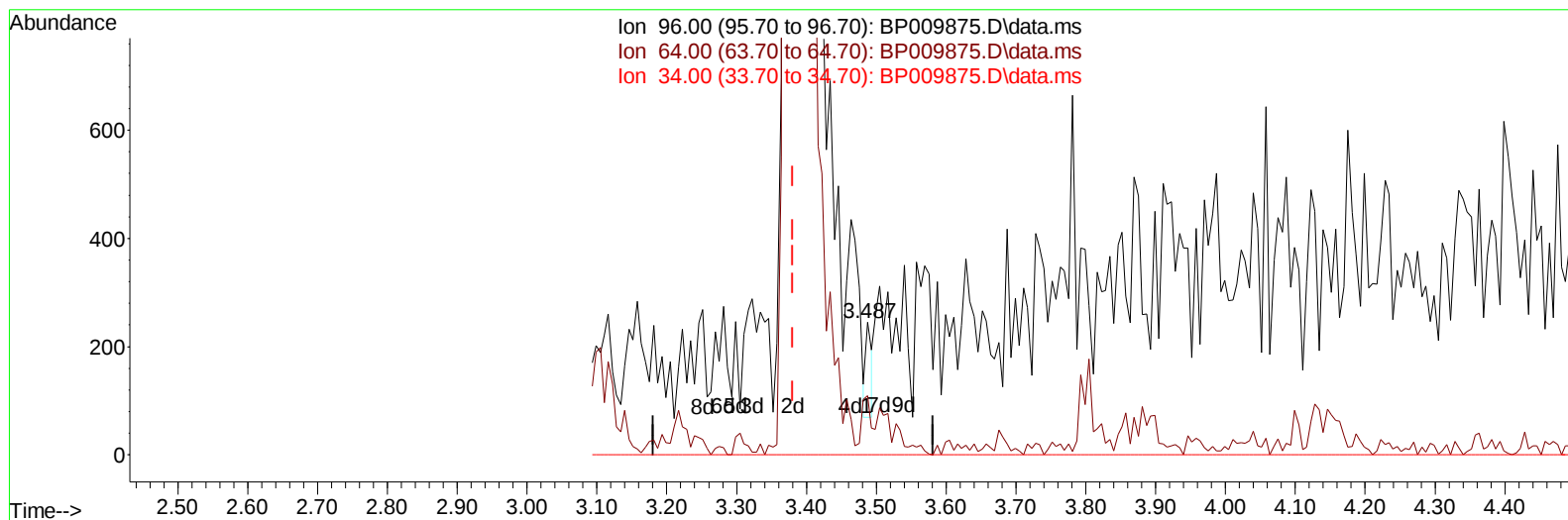
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009875.D
 Acq On : 15 Apr 2022 23:29
 Operator : CG/JU
 Sample : N2421-06
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0HY6

Manual IntegrationsAPPROVED

Quant Time: Apr 18 01:06:19 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: BP009875.D\data.ms

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

3.487min (+ 0.106) 0.04 ng/uL

response 106

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	44.49
34.00	0.00	0.00
0.00	0.00	0.00

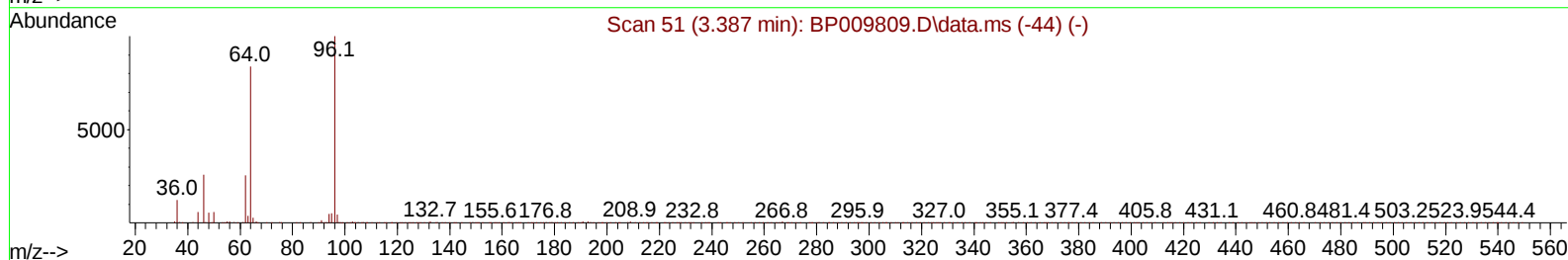
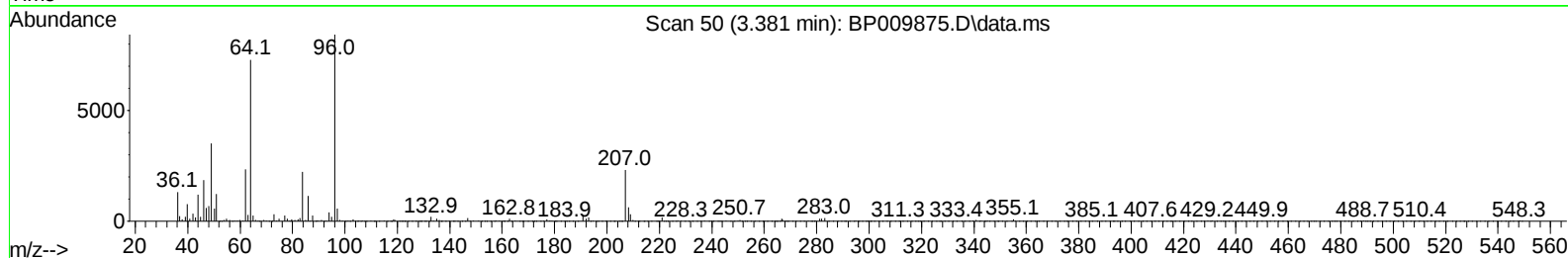
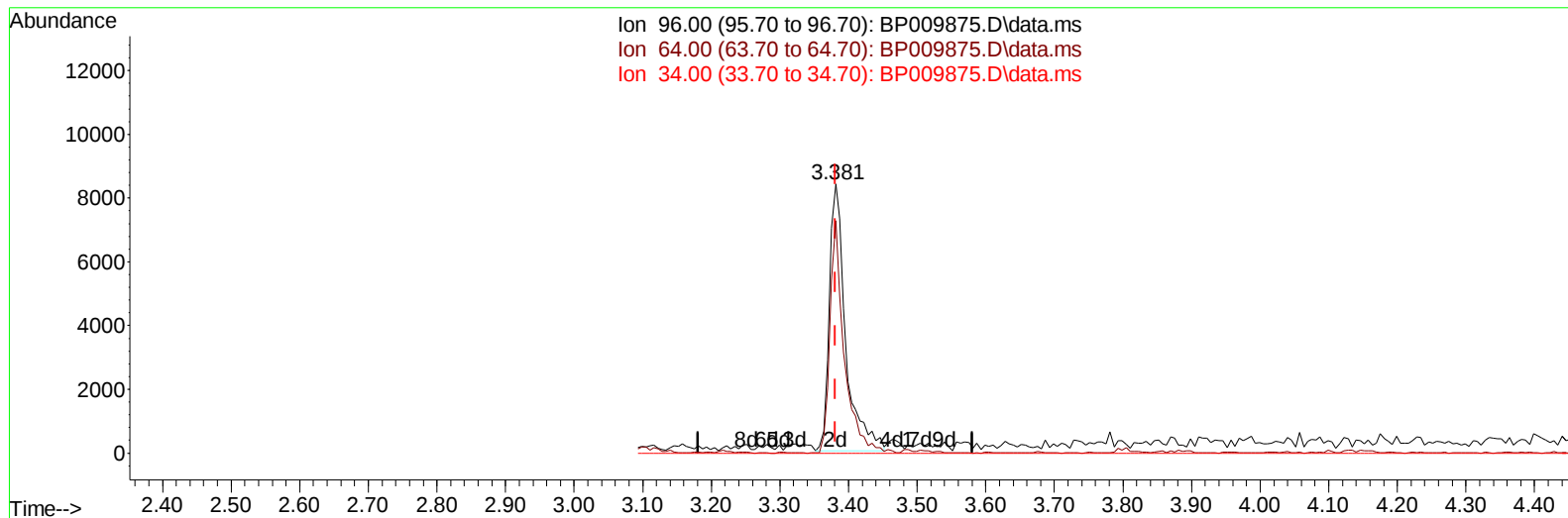
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(3) 1,4-Dioxane-d8 (S)

3.381min (-0.000) 5.24 ng/uL m

response 13967

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.49#
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	7.957	152	119062	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	526685	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.586	164	380144	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	859675	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	959356	20.000	ng/ul	0.00
88) Perylene-d12	23.880	264	956395	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	13967m	5.235	ng/uL	0.00
4) Pyridine-d5	3.799	84	89922	12.225	ng/ul	0.00
7) Phenol-d5	7.110	99	96622	9.244	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	193240	31.053	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	205730	26.092	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	173781	19.915	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	127221	33.496	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	134176	32.223	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	255012	29.266	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	323728	26.085	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	1036803	34.721	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1081164	30.561	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	56854	10.499	ng/ul	0.00
60) Fluorene-d10	15.581	176	891395	35.485	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	168613	33.109	ng/ul	0.00
73) Anthracene-d10	17.439	188	1568881	38.116	ng/ul	0.00
81) Pyrene-d10	19.669	212	2037902	38.895	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	2061380	41.388	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.134	94	17131	1.637	ng/ul#	85

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

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