

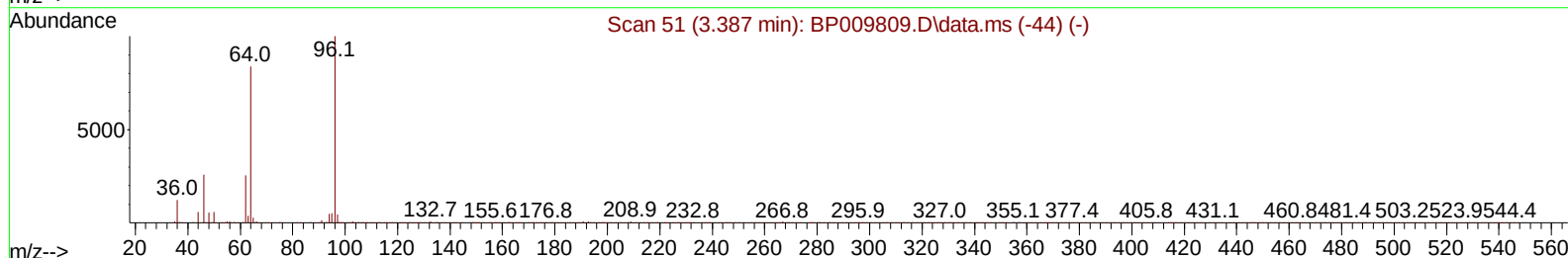
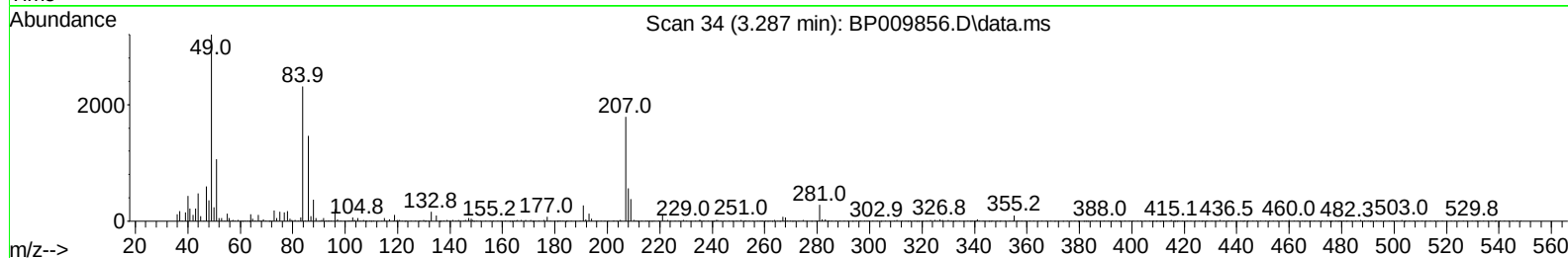
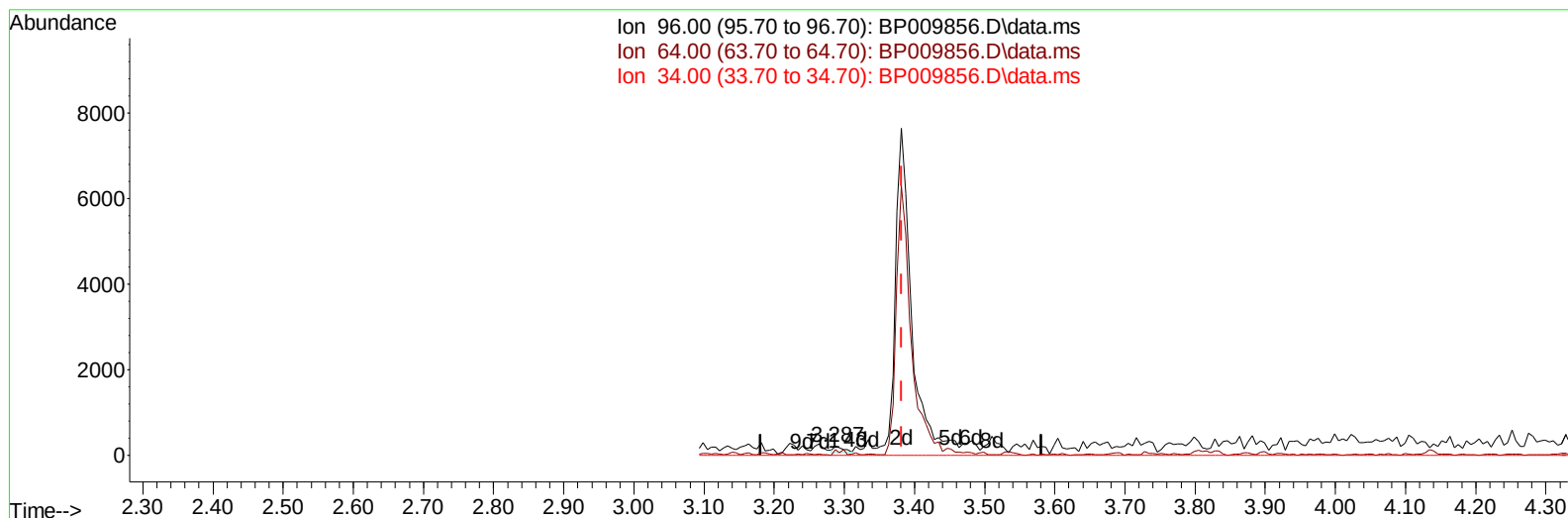
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
Data File : BP009856.D
Acq On : 15 Apr 2022 12:15
Operator : CG/JU
Sample : N2422-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J08

Manual IntegrationsAPPROVED

Quant Time: Apr 18 01:00:24 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: BP009856.D\data.ms

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

3.287min (-0.094) 0.06 ng/uL

response 136

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	59.90
34.00	0.00	0.00
0.00	0.00	0.00

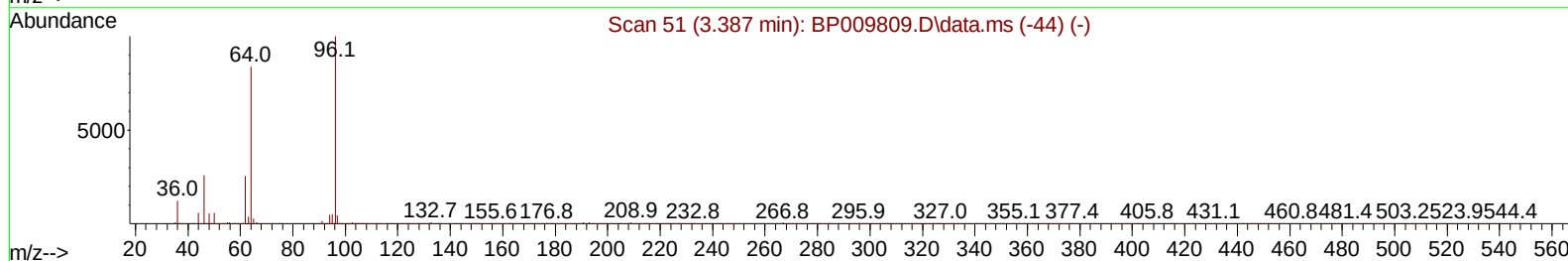
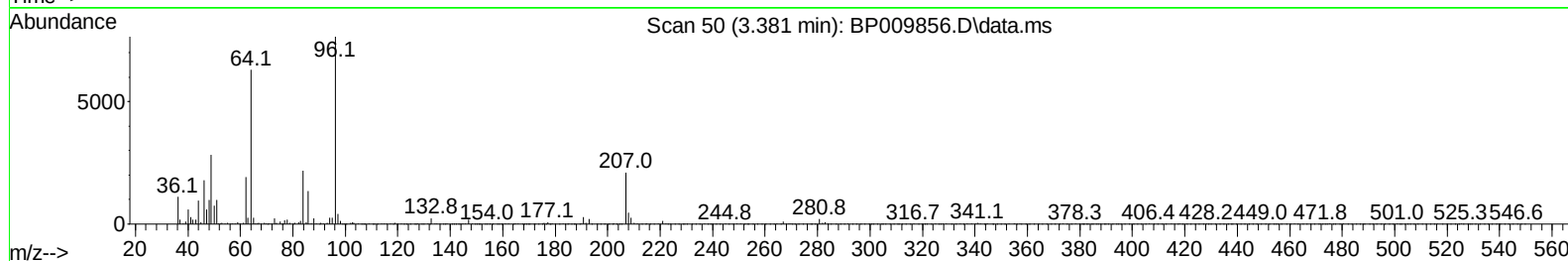
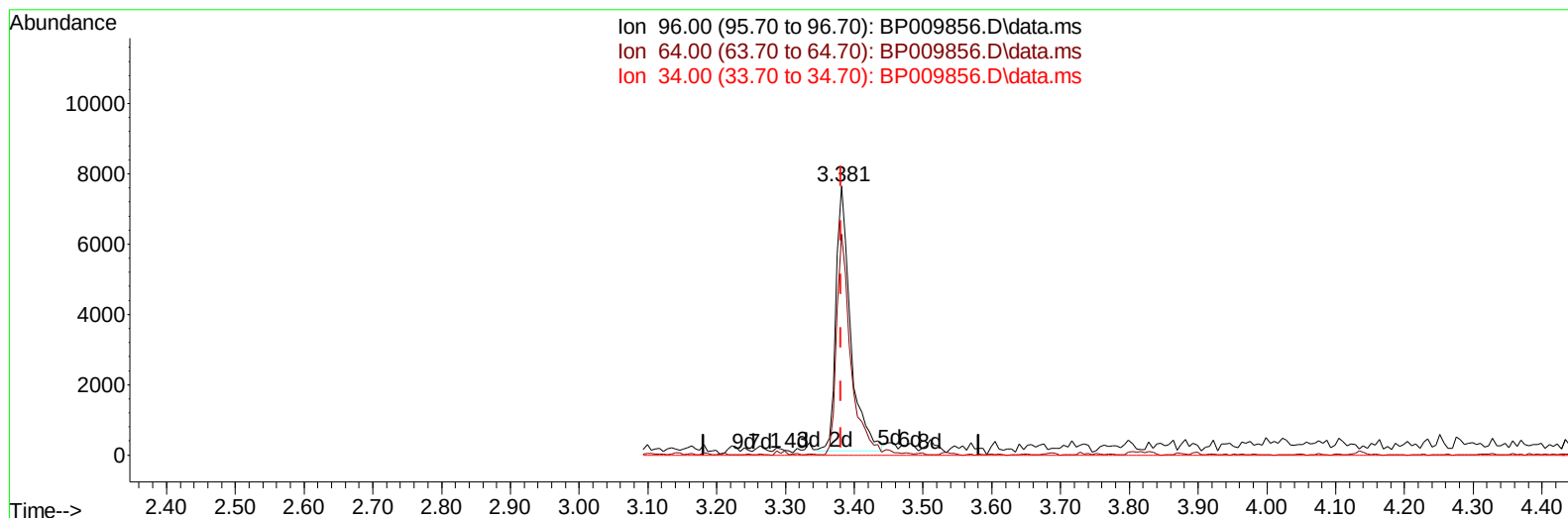
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(3) 1,4-Dioxane-d8 (S)

3.381min (-0.000) 4.82 ng/uL m

response 11170

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	82.25#
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	103471	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	462201	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	336673	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	772993	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.421	240	832020	20.000	ng/ul	# 0.00
88) Perylene-d12	23.886	264	810763	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	11170m	4.818	ng/uL	0.00
4) Pyridine-d5	3.805	84	61727	9.657	ng/ul	0.00
7) Phenol-d5	7.110	99	60106	6.617	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	174182	32.208	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	163274	23.828	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	118266	15.596	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	111521	33.459	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	115090	31.495	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	197917	25.883	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	287732	26.419	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	866688	32.772	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	939223	29.977	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	28570	5.957	ng/ul	0.00
60) Fluorene-d10	15.581	176	771828	34.693	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	127399	27.821	ng/ul	0.00
73) Anthracene-d10	17.439	188	1368672	36.981	ng/ul	0.00
81) Pyrene-d10	19.669	212	1691359	37.221	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	1629014	38.582	ng/ul	0.00

Target Compounds Qvalue

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

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