

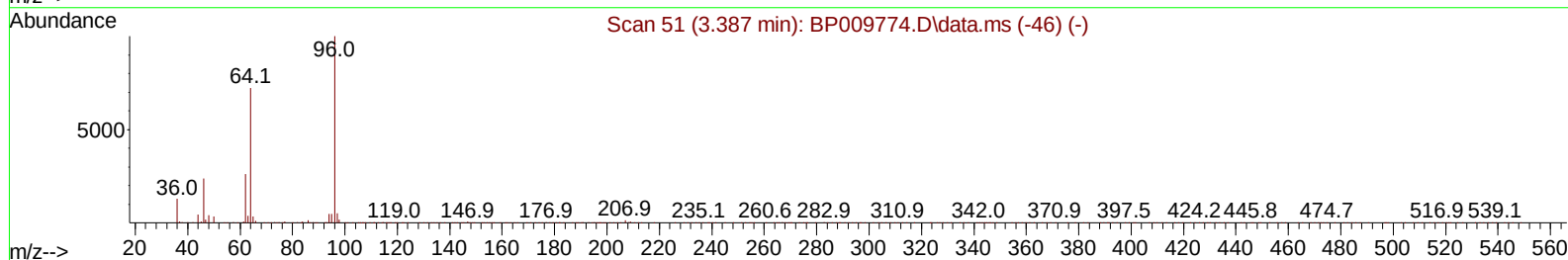
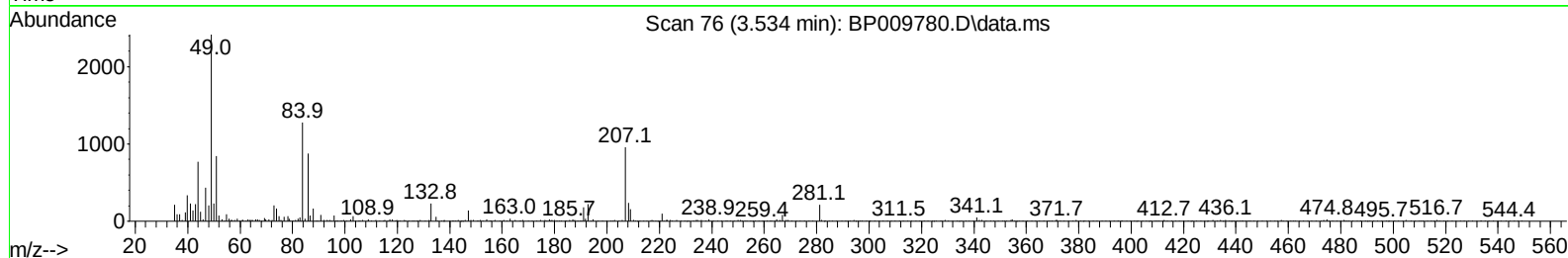
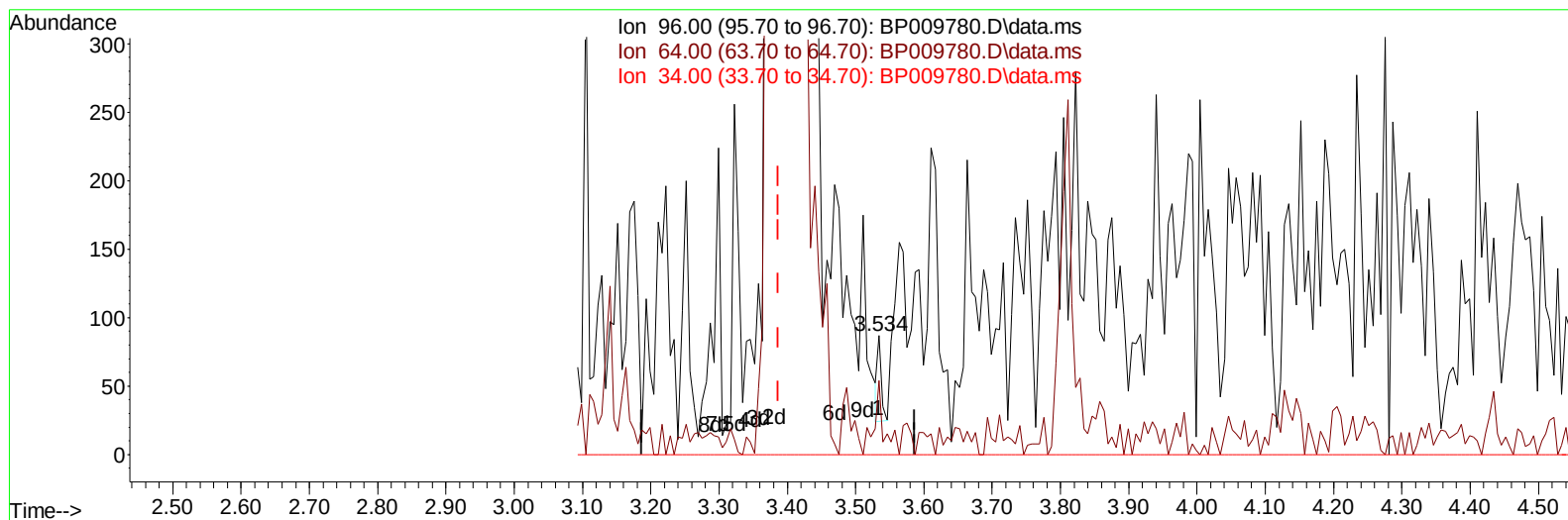
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
Data File : BP009780.D
Acq On : 12 Apr 2022 12:36
Operator : CG/JU
Sample : N2338-08
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J98

Manual IntegrationsAPPROVED

Quant Time: Apr 13 02:20:15 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 04/13/2022
Supervised By :Yogesh Patel 04/13/2022



TIC: BP009780.D\data.ms

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

3.534min (+ 0.147) 0.01 ng/uL

response 26

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	62.07
34.00	0.00	0.00
0.00	0.00	0.00

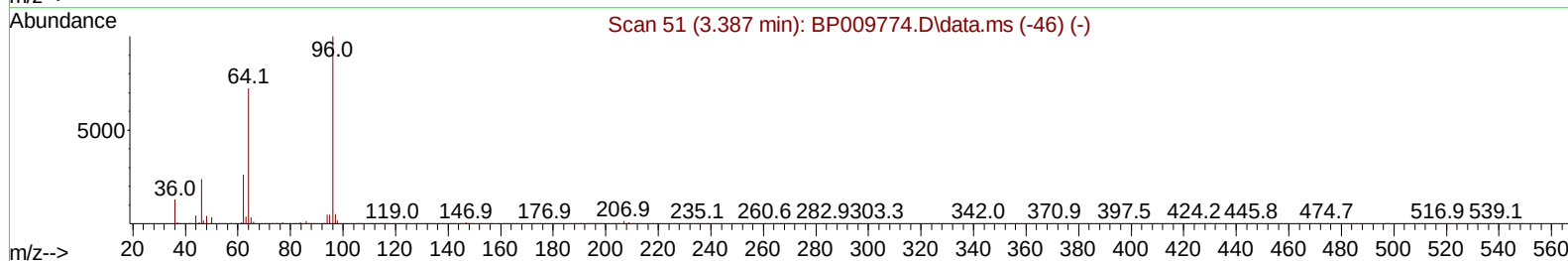
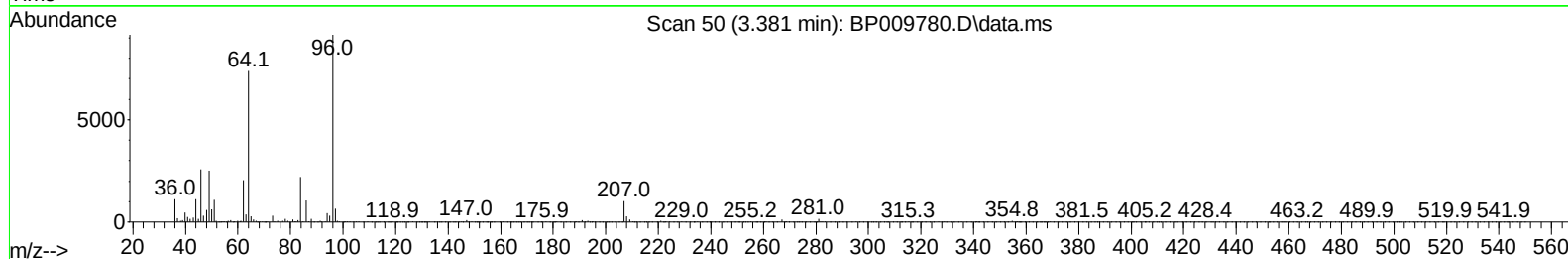
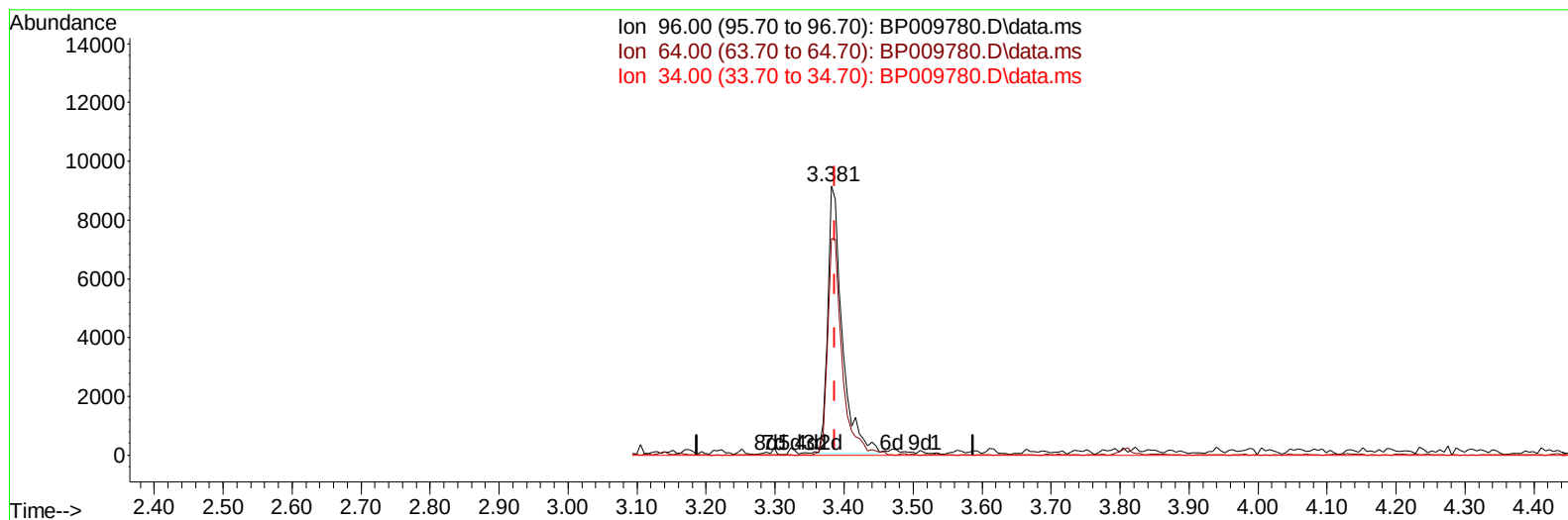
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TIC: BP009780.D\data.ms

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

3.381min (-0.006) 6.08 ng/uL m

response 13413

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.60#
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	7.969	152	98420	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.775	136	442608	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164	317951	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	714199	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.427	240	776099	20.000	ng/ul	# 0.00
88) Perylene-d12	23.898	264	821446	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	13413m	6.082	ng/uL	0.00
4) Pyridine-d5	3.805	84	108319	17.815	ng/ul	0.00
7) Phenol-d5	7.116	99	120127	13.903	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	194752	37.859	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	193396	29.672	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	163989	22.735	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	125517	39.325	ng/ul	0.00
24) 2-Nitrophenol-d4	9.852	143	128763	36.797	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	234189	31.982	ng/ul	0.00
31) 4-Chloroaniline-d4	10.904	131	334556	32.078	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	1039480	41.620	ng/ul	0.00
49) Acenaphthylene-d8	14.292	160	1076936	36.397	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	70476	15.560	ng/ul	0.00
60) Fluorene-d10	15.592	176	873039	41.553	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	159763	37.761	ng/ul	0.00
73) Anthracene-d10	17.451	188	1488898	43.541	ng/ul	0.00
81) Pyrene-d10	19.675	212	1938411	45.731	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.739	264	1997538	46.695	ng/ul	0.00

Target Compounds Qvalue

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

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