

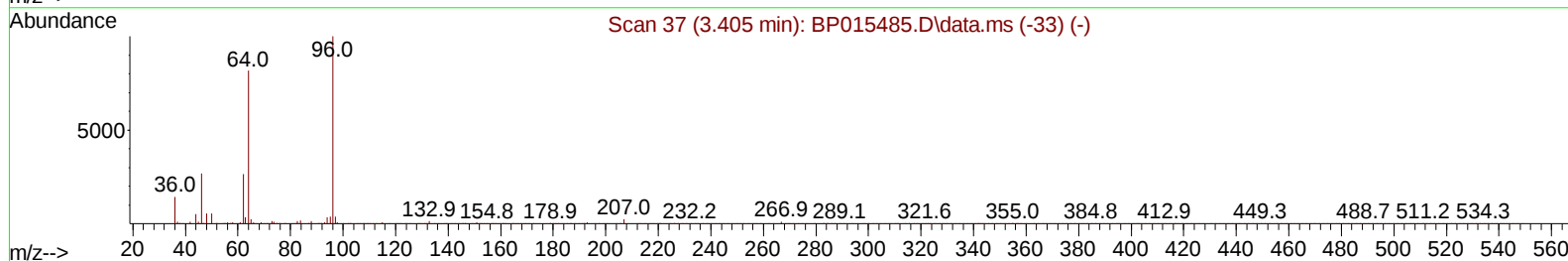
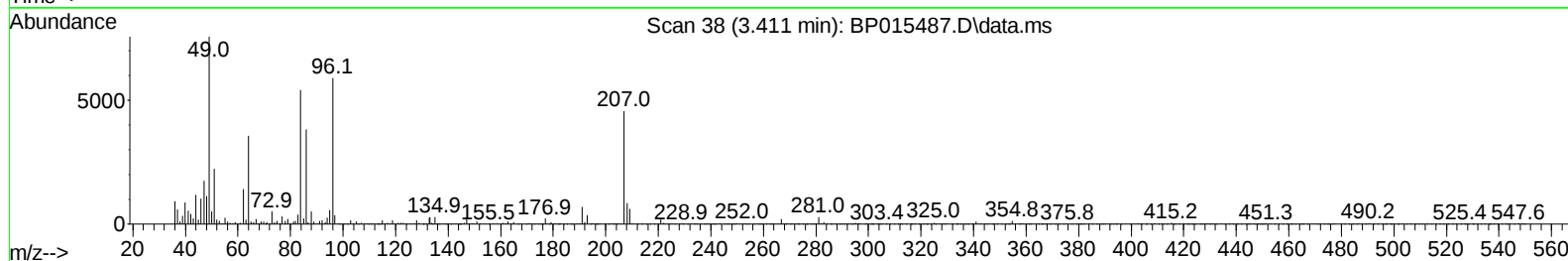
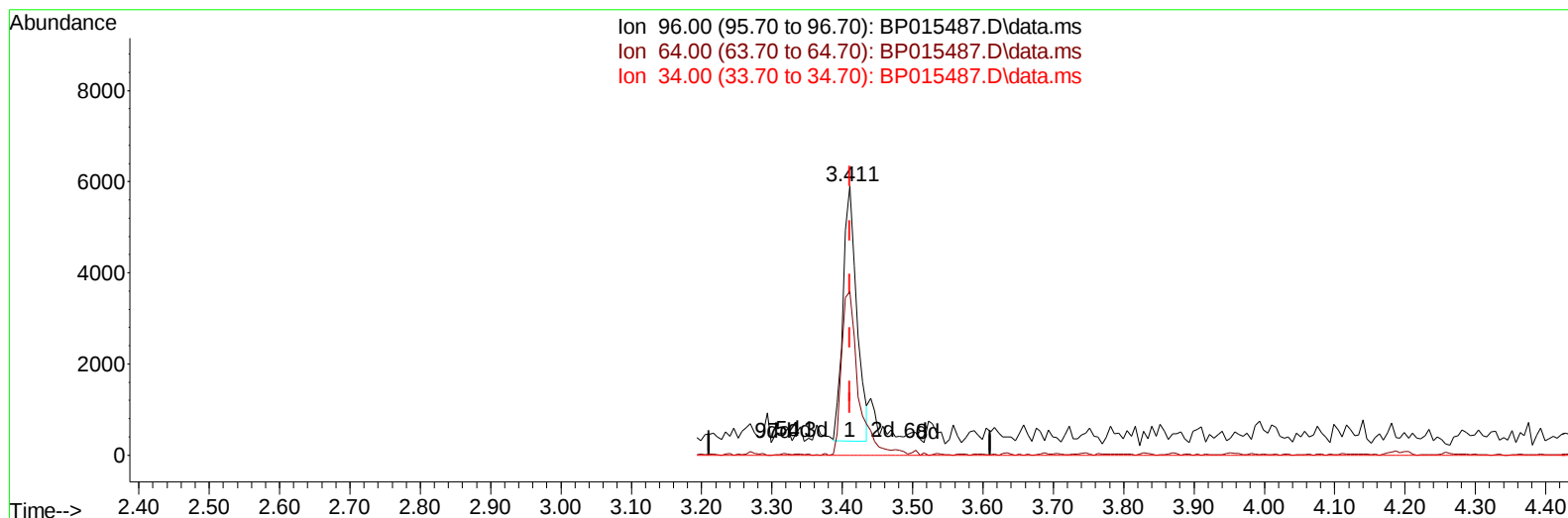
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061323\
Data File : BP015487.D
Acq On : 13 Jun 2023 09:55
Operator : MA/JU
Sample : 03078-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ESQW9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/14/2023
Supervised By :mohammad ahmed 06/16/2023

Quant Time: Jun 14 00:12:29 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration



TIC: BP015487.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.411min (-0.000) 1.75 ng/uL****response 7527**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.60	60.66
34.00	0.00	0.00
0.00	0.00	0.00

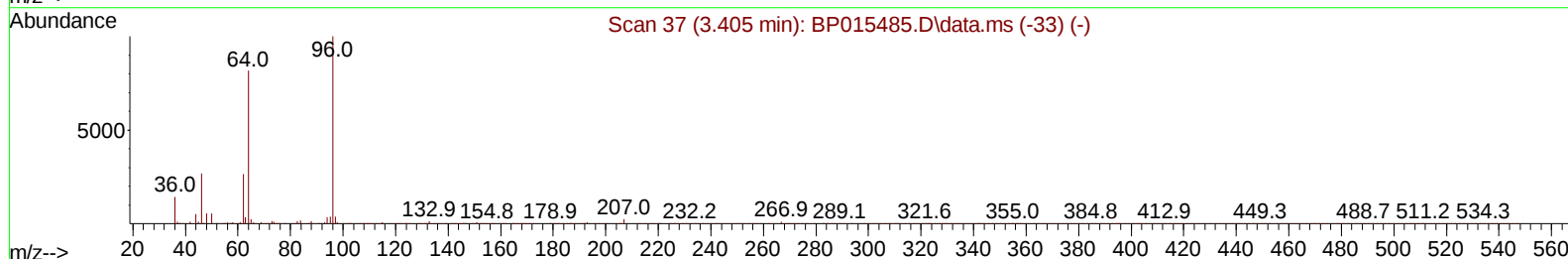
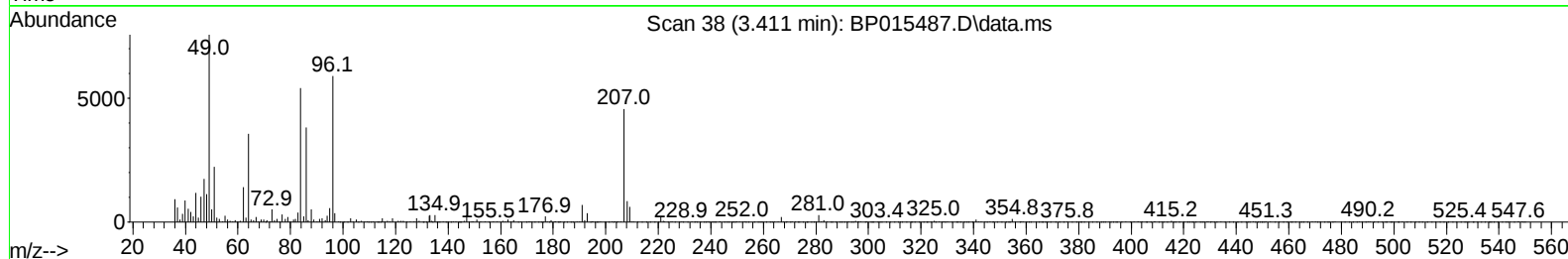
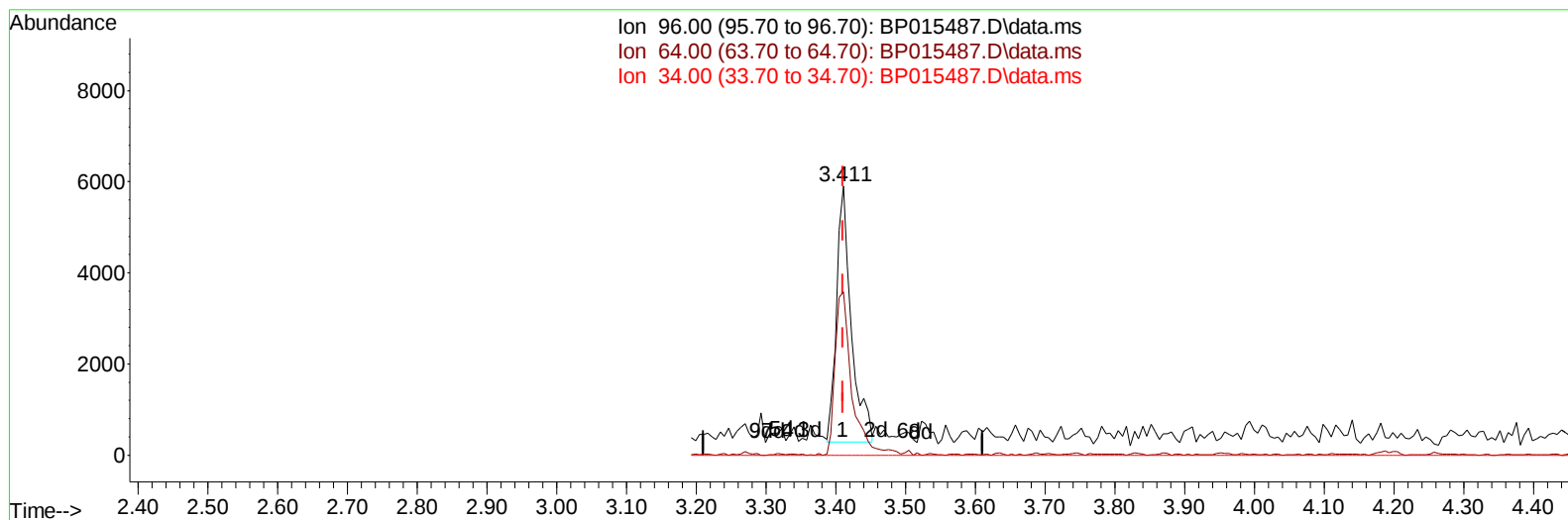
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TIC: BP015487.D\data.ms

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

3.411min (-0.000) 1.91 ng/uL m

response 8230

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.60	60.66
34.00	0.00	0.00
0.00	0.00	0.00

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Quant Time: Jun 14 00:22:45 2023
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.093	152		159455	20.000	ng/ul	0.00
20) Naphthalene-d8	10.922	136		753380	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.739	164		532420	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.492	188		1260302	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240		1207854	20.000	ng/ul	0.00
88) Perylene-d12	24.121	264		1254156	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.411	96		8230m	1.911	ng/uL	0.00
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	7.246	99		124216	8.456	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67		88336	10.069	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132		101637	9.543	ng/ul	0.00
15) 4-Methylphenol-d8	8.810	113		60059	4.982	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128		52199	8.825	ng/ul	0.00
24) 2-Nitrophenol-d4	10.004	143		59063	8.744	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.551	165		102415	8.312	ng/ul	0.01
31) 4-Chloroaniline-d4	11.081	131		39228	2.218	ng/ul	0.02
46) Dimethylphthalate-d6	14.140	166		430455	10.253	ng/ul	0.00
49) Acenaphthylene-d8	14.434	160		454338	9.897	ng/ul	0.00
54) 4-Nitrophenol-d4	14.963	143		30077m	4.000	ng/ul	0.02
60) Fluorene-d10	15.728	176		369573	10.439	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200		44860	5.619	ng/ul	0.00
73) Anthracene-d10	17.592	188		585593	10.212	ng/ul	0.00
81) Pyrene-d10	19.816	212		777493	12.027	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264		644411	10.240	ng/ul	0.00
Target Compounds							
						Qvalue	
6) Benzaldehyde	7.234	77		13050	2.324	ng/ul	93
8) Phenol	7.275	94		26751	1.765	ng/ul	97
16) Acetophenone	8.934	105		121842	6.333	ng/ul	97

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

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