

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010341.D
 Acq On : 13 May 2022 01:20
 Operator : CG/JU
 Sample : N2801-05
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

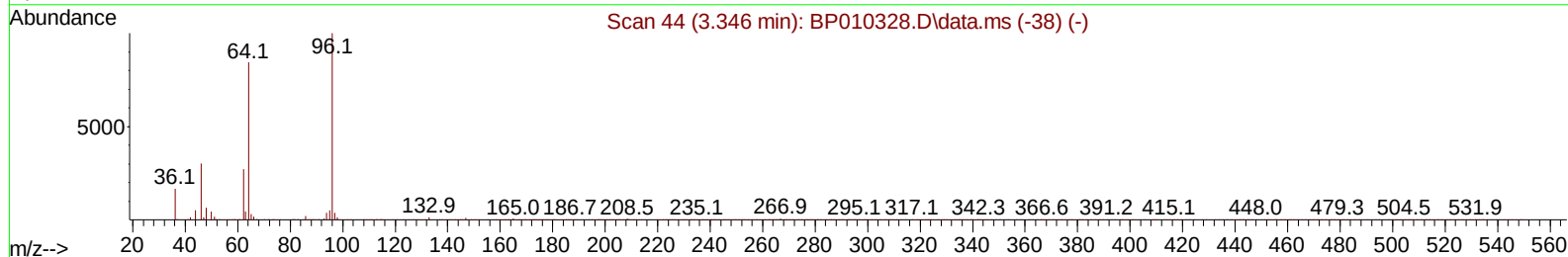
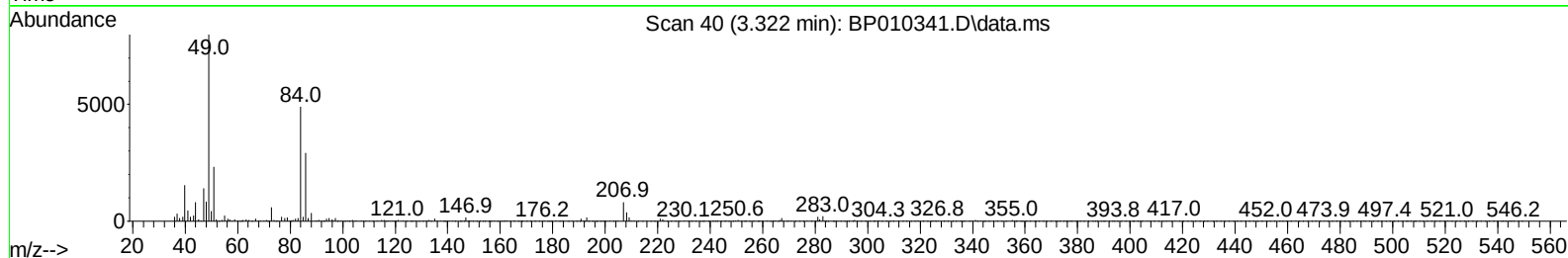
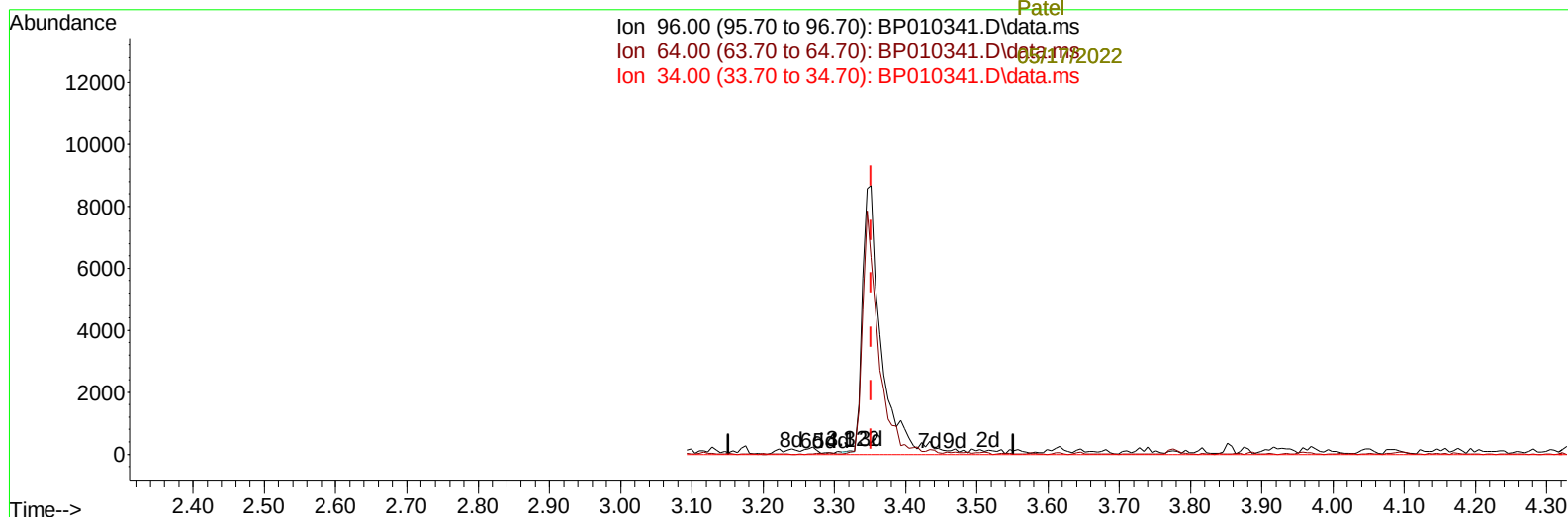
Instrument :
 BNA_P
 ClientSampleId :
 EXLB8

Manual IntegrationsAPPROVED

Quant Time: May 13 01:57:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut
 Upadhyay

05/16/2022
 Supervised By :Yogesh
 Patel



TIC: BP010341.D\data.ms

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

3.322min (-0.029) 0.01 ng/uL

response 22

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	63.46
34.00	0.00	0.00
0.00	0.00	0.00

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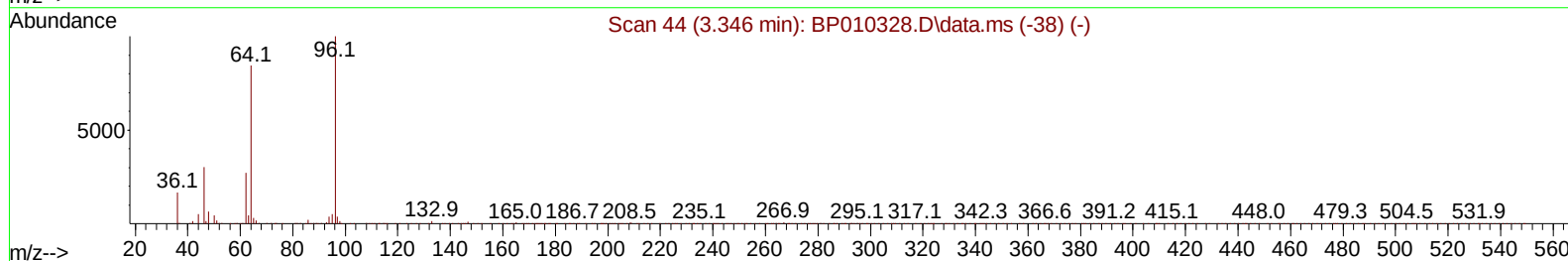
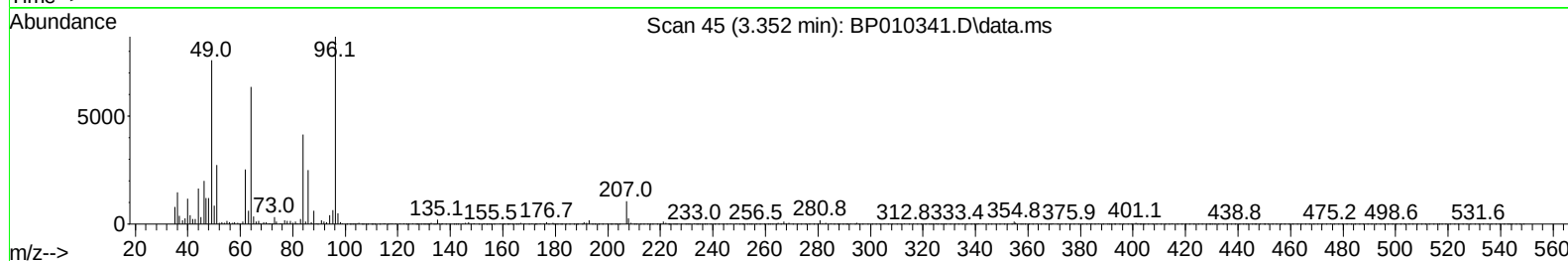
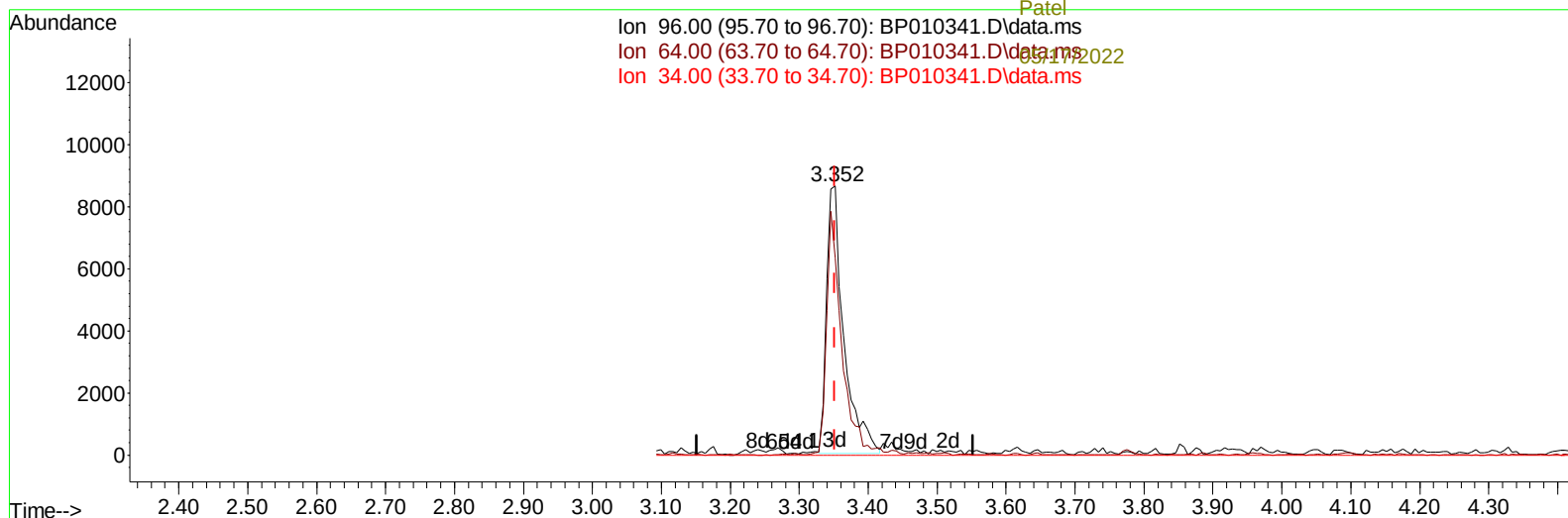
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(3) 1,4-Dioxane-d8 (S)

3.352min (+ 0.000) 4.49 ng/uL m

response 15000

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.27#
34.00	0.00	0.00
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----05/17/2022							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.905	152		142067	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136		625854	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164		436078	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188		964152	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.363	240		981321	20.000	ng/ul	#-0.01
88) Perylene-d12	23.792	264		888391	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.352	96		15000m	4.485	ng/uL	0.00
4) Pyridine-d5	3.769	84		89638	9.632	ng/ul	0.00
7) Phenol-d5	7.069	99		78853	5.976	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67		268513	32.895	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132		237266	24.416	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113		161276	14.867	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128		165882	33.056	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.787	143		175659	30.701	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165		284403	27.310	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131		401067	26.087	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166		1174289	33.801	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160		1326906	32.035	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143		27811m	4.173	ng/ul	0.00
60) Fluorene-d10	15.528	176		1029163	35.095	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200		172141	26.072	ng/ul	-0.01
73) Anthracene-d10	17.381	188		1691265	35.821	ng/ul	-0.01
81) Pyrene-d10	19.616	212		2053897	37.875	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.633	264		1730889	36.534	ng/ul	-0.02

Target Compounds Qvalue

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

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