

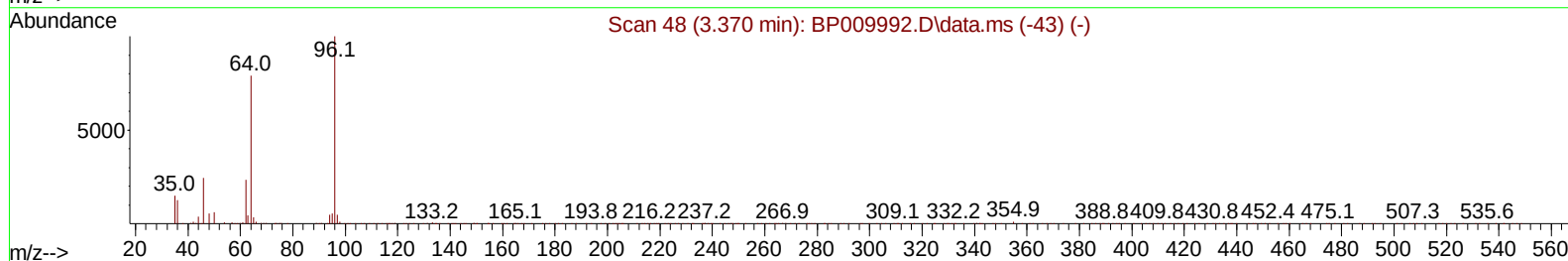
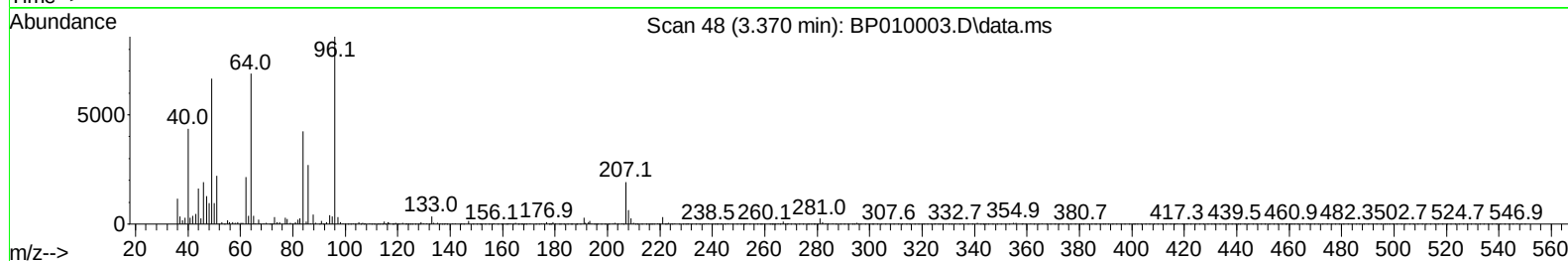
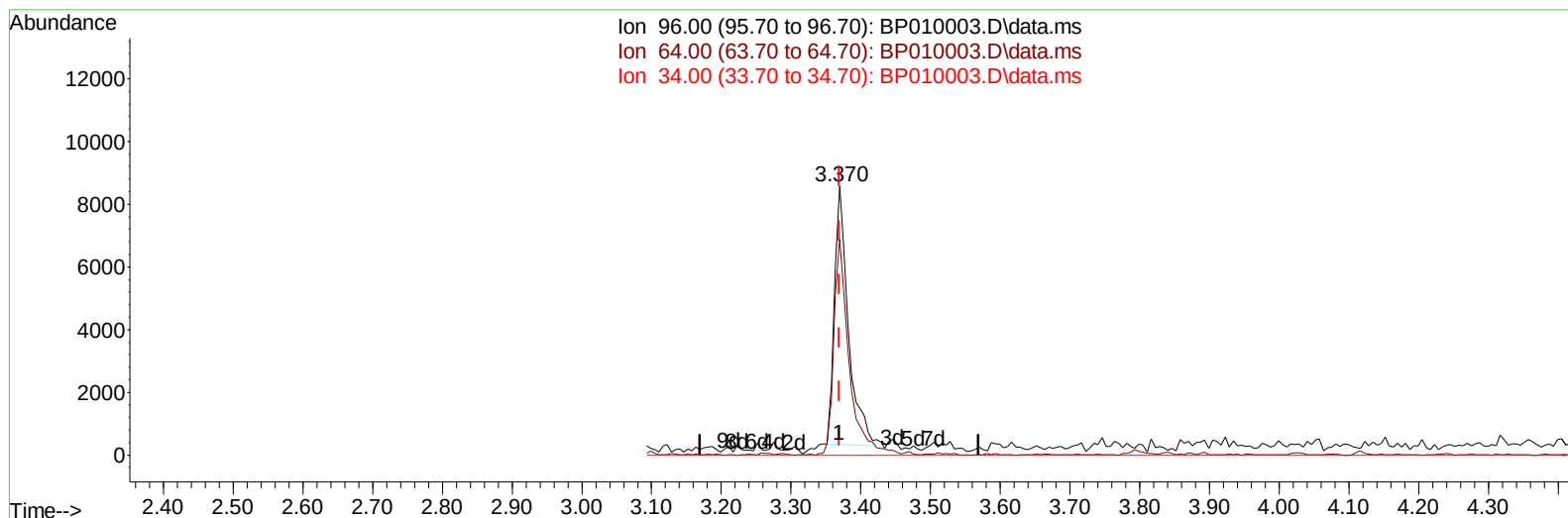
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP010003.D
 Acq On : 21 Apr 2022 06:44
 Operator : CG/JU
 Sample : N2473-03
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J25

Manual IntegrationsAPPROVED

Quant Time: Apr 21 07:25:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/21/2022
 Supervised By :Yogesh Patel 04/25/2022



TIC: BP010003.D\data.ms

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

3.370min (-0.000) 4.25 ng/uL

response 11339

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.24#
34.00	0.00	0.00
0.00	0.00	0.00

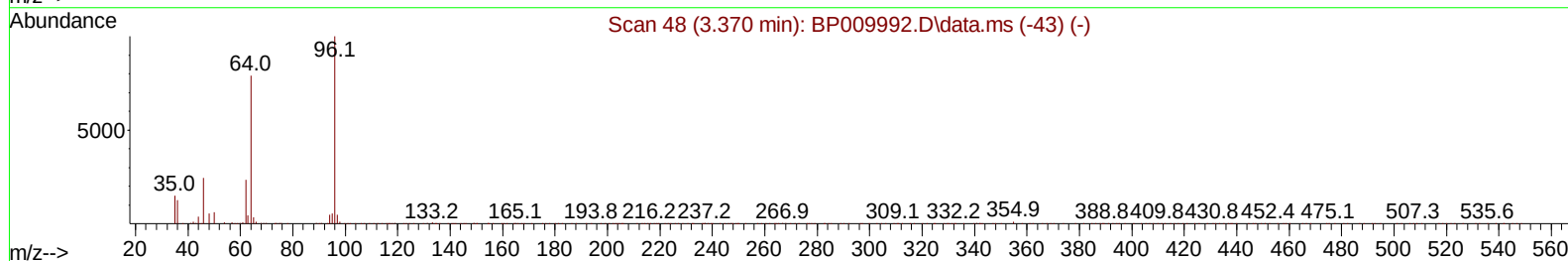
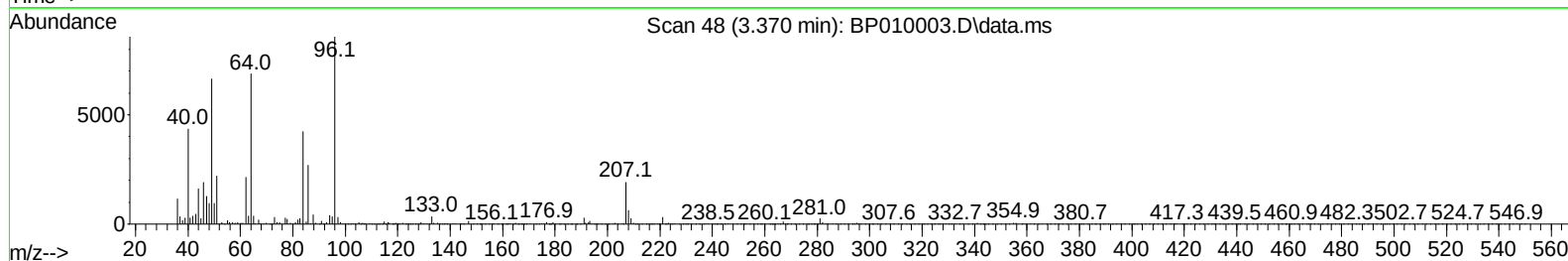
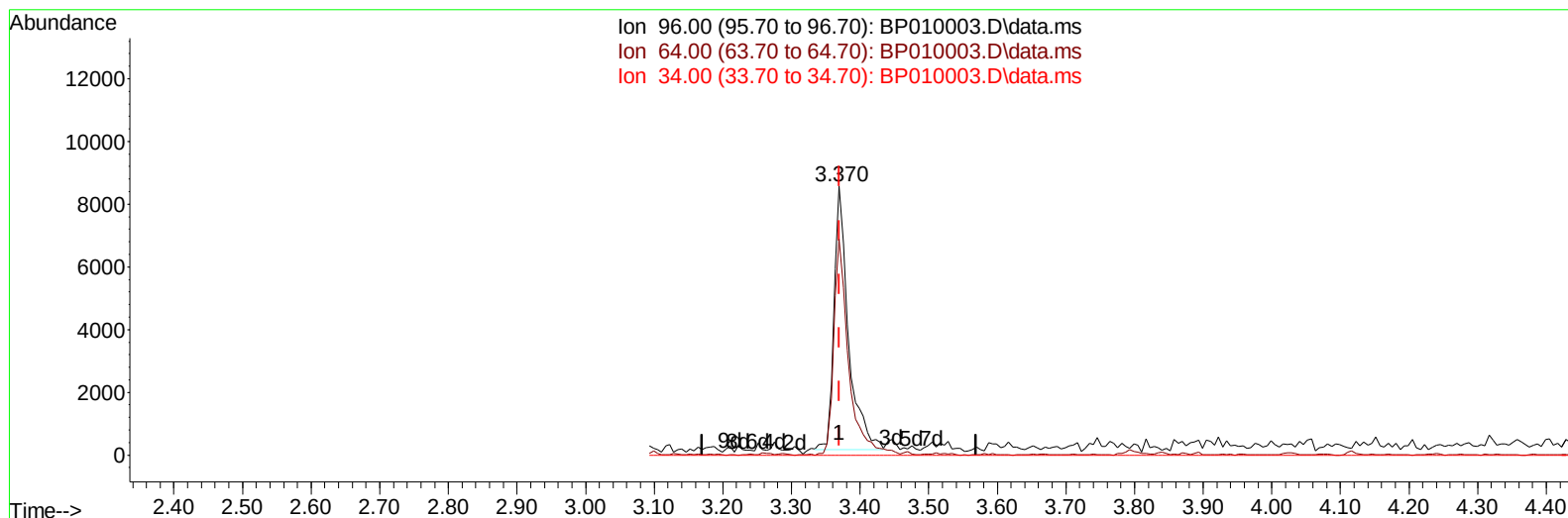
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TIC: BP010003.D\data.ms

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

3.370min (-0.000) 4.61 ng/uL m

response 12293

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.24#
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	7.940	152	118943	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.746	136	557220	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	409007	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	923857	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	935901	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264	848694	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	12293m	4.613	ng/uL	0.00
4) Pyridine-d5	3.793	84	79319	10.794	ng/ul	0.00
7) Phenol-d5	7.099	99	75792	7.258	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	220727	35.505	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	209705	26.623	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	158908	18.229	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	148773	37.024	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	158137	35.896	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	275177	29.850	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	344326	26.224	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1176549	36.621	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1260369	33.113	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	38414	6.593	ng/ul	0.00
60) Fluorene-d10	15.563	176	1026460	37.979	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	185698	33.930	ng/ul	-0.01
73) Anthracene-d10	17.422	188	1809536	40.909	ng/ul	0.00
81) Pyrene-d10	19.651	212	2329862	45.581	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1919062	43.421	ng/ul	-0.01

Target Compounds Qvalue

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

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