

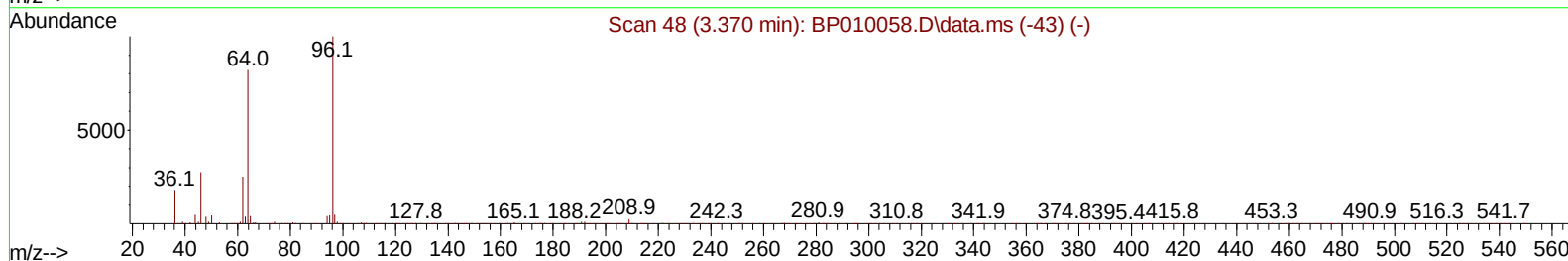
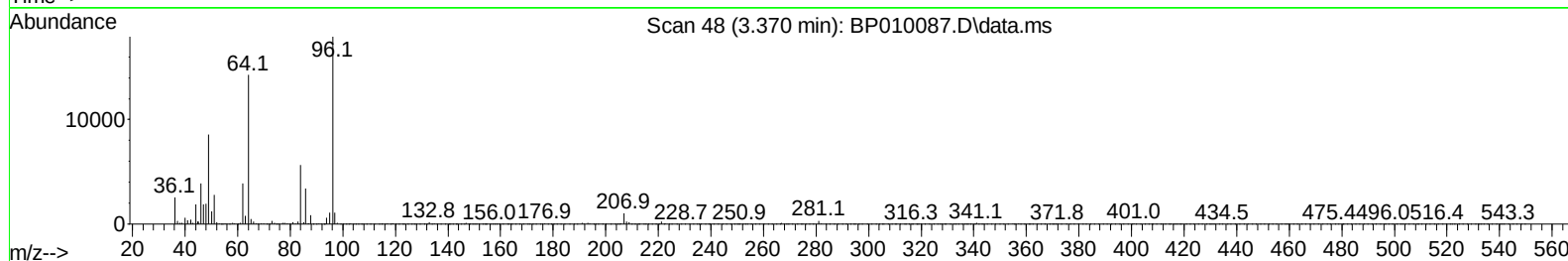
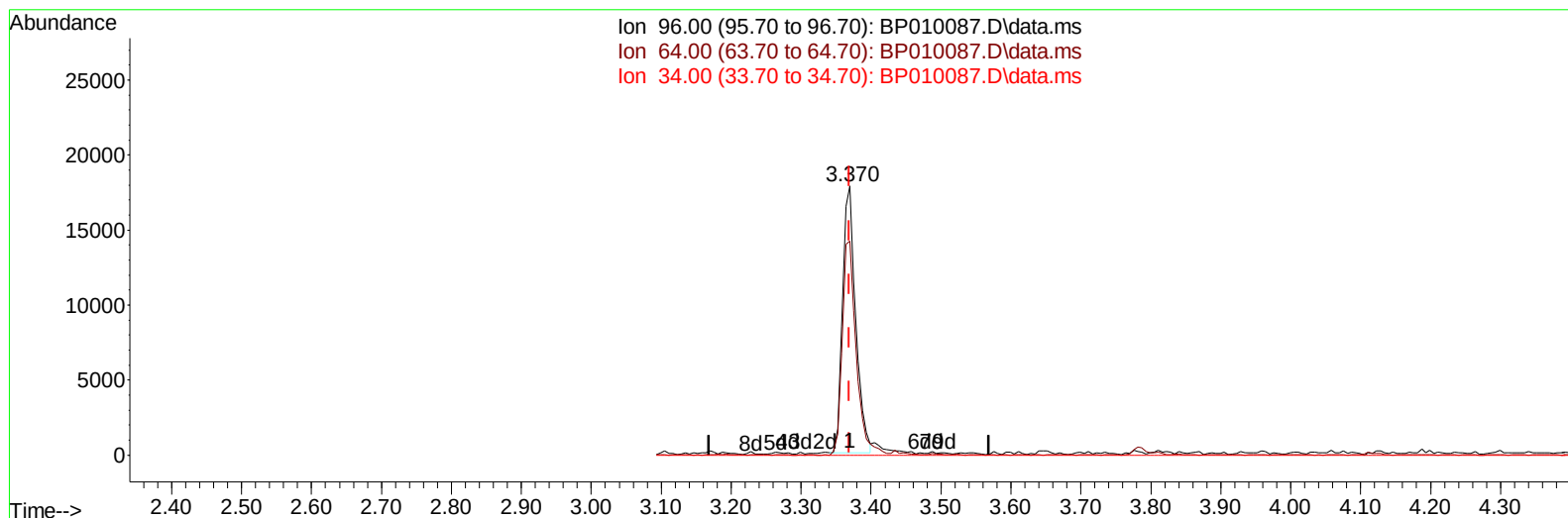
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
 Data File : BP010087.D
 Acq On : 25 Apr 2022 17:40
 Operator : CG/JU
 Sample : PB144305BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK305

Manual IntegrationsAPPROVED

Quant Time: Apr 26 01:03:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022
 Supervised By :Yogesh Patel 04/28/2022



TIC: BP010087.D\data.ms

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

3.370min (+ 0.001) 5.55 ng/uL

response 23470

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	79.65#
34.00	0.00	0.00
0.00	0.00	0.00

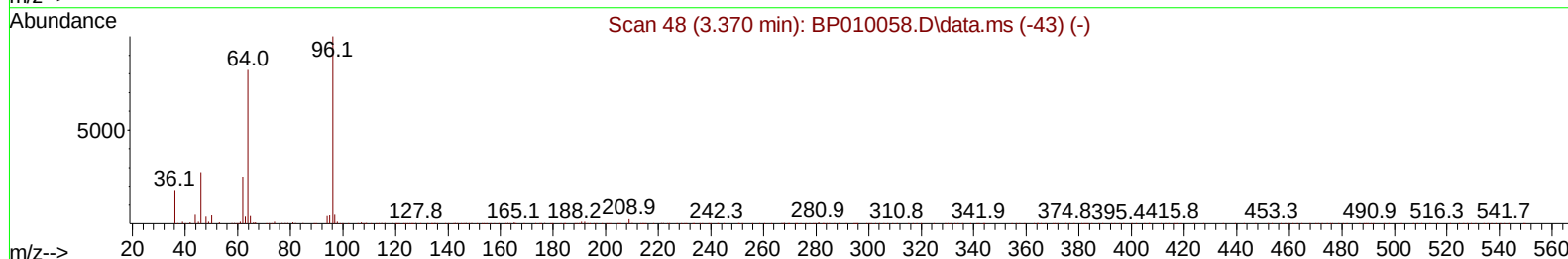
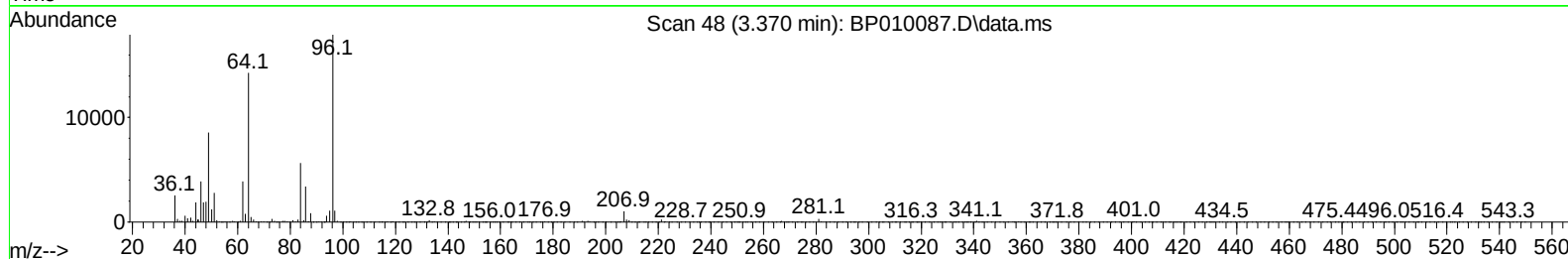
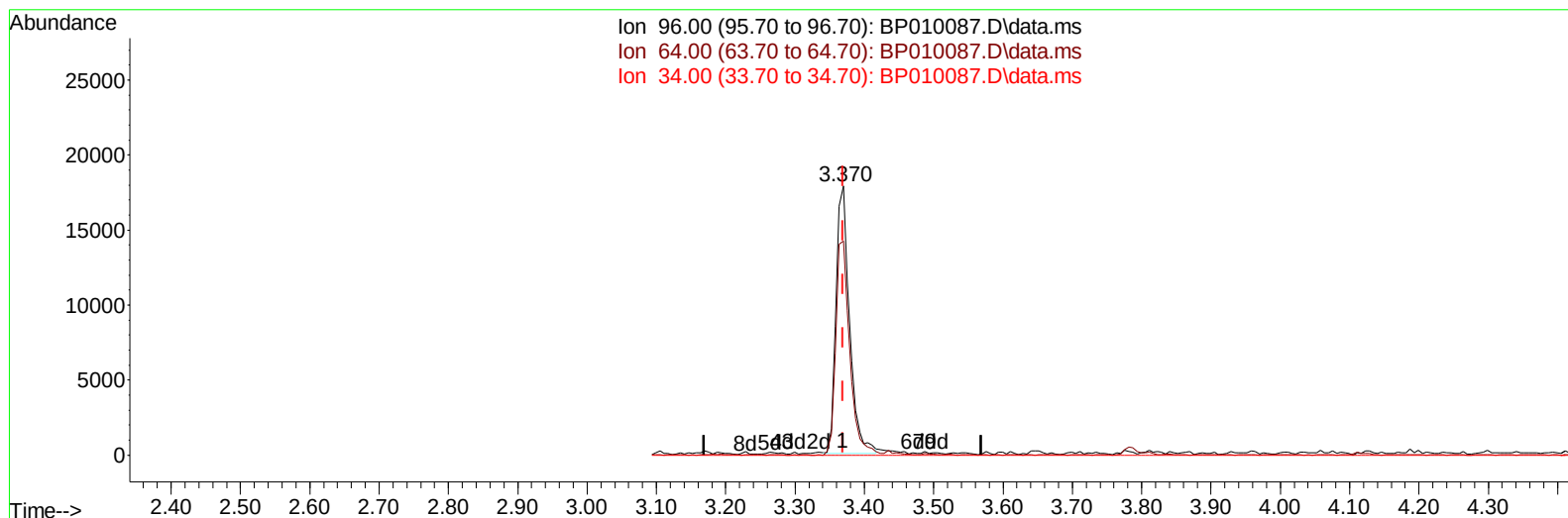
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TIC: BP010087.D\data.ms

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

3.370min (+ 0.001) 5.75 ng/uL m

response 24296

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	79.65#
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	7.934	152	184340	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	790127	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	542588	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.316	188	1165855	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	1029725	20.000	ng/ul	# 0.00
88) Perylene-d12	23.851	264	883713	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	24296m	5.747	ng/uL	0.00
4) Pyridine-d5	3.781	84	303686	25.536	ng/ul	0.00
7) Phenol-d5	7.093	99	412012	24.572	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	292918	29.114	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.464	132	349315	28.130	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	358048	25.773	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	182817	29.091	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	199621	29.086	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.358	165	330559	24.741	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	492520	25.709	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1272371	30.146	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1373120	26.881	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	158678	20.421	ng/ul	0.00
60) Fluorene-d10	15.557	176	1048245	28.872	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	181377	24.604	ng/ul	0.00
73) Anthracene-d10	17.416	188	1660884	29.447	ng/ul	0.00
81) Pyrene-d10	19.645	212	1934297	33.031	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	1389647	29.995	ng/ul	0.00

Target Compounds Qvalue

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

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