

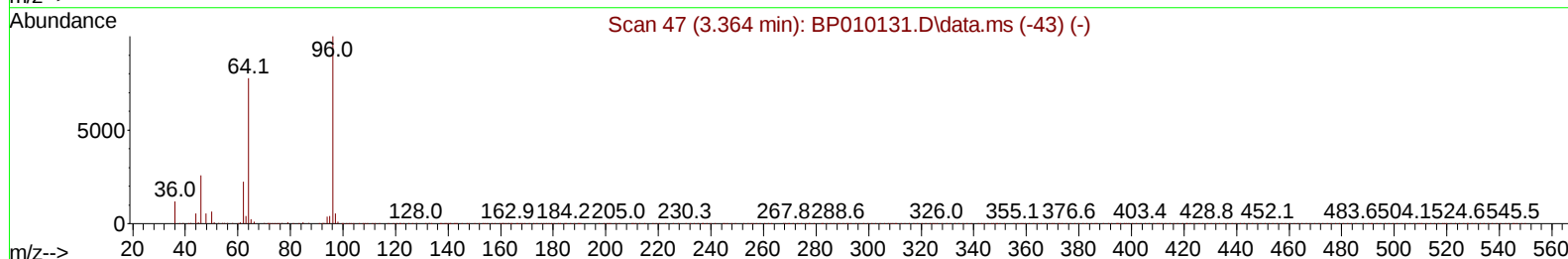
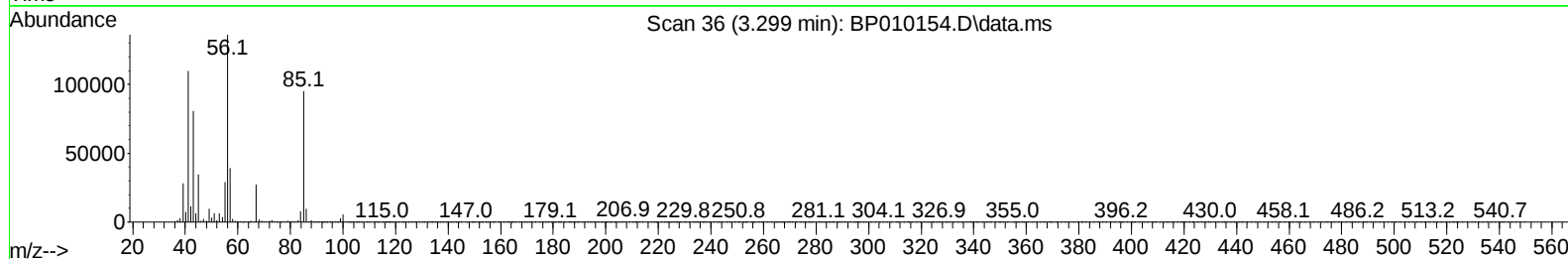
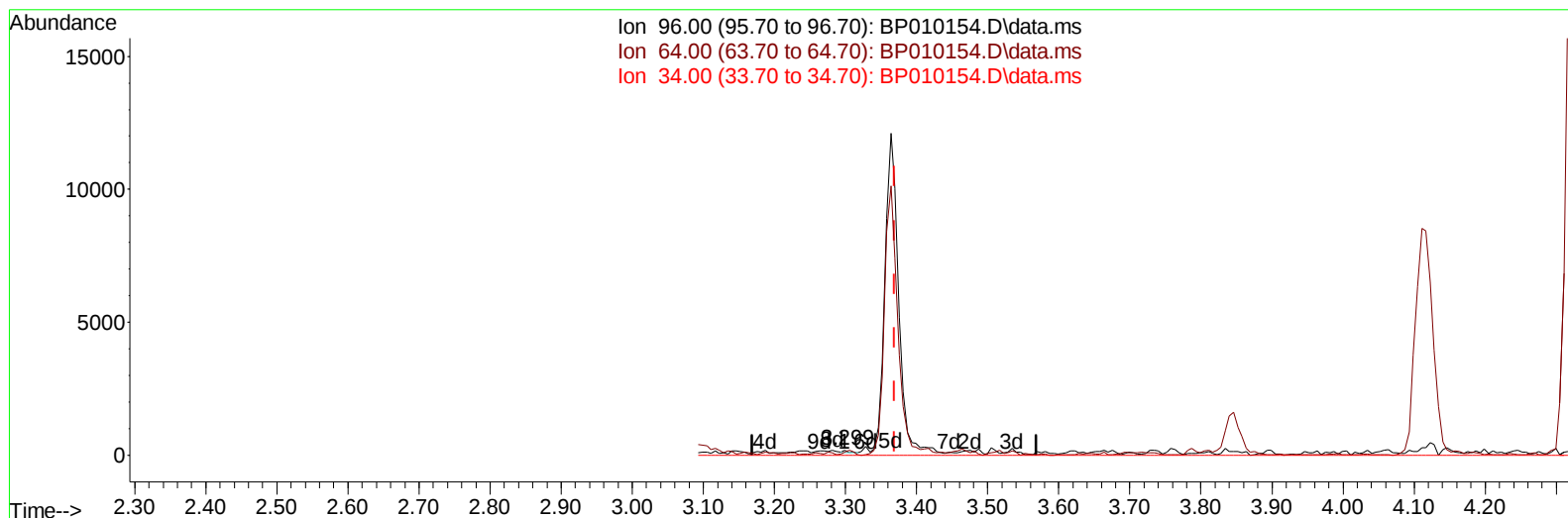
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010154.D
 Acq On : 29 Apr 2022 16:43
 Operator : CG/JU
 Sample : N2587-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXET4

Manual IntegrationsAPPROVED

Quant Time: Apr 30 00:31:12 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 05/02/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010154.D\data.ms

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

3.299min (-0.070) 0.01 ng/uL

response 49

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	64.84
34.00	0.00	0.00
0.00	0.00	0.00

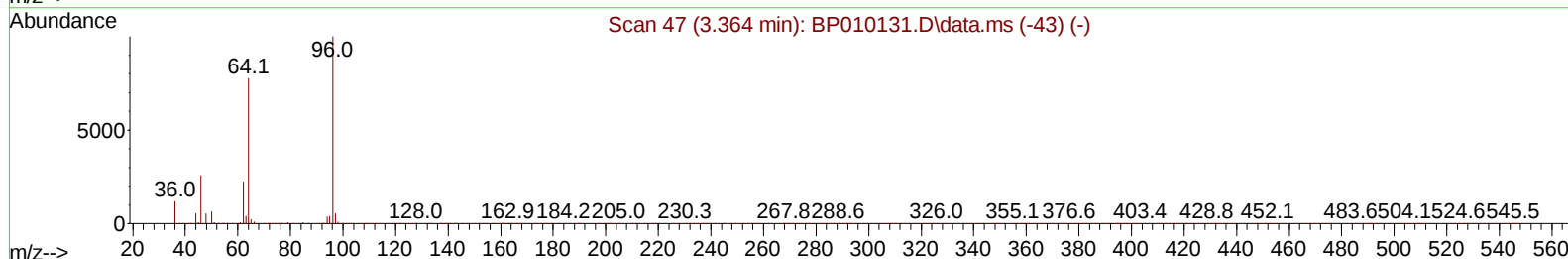
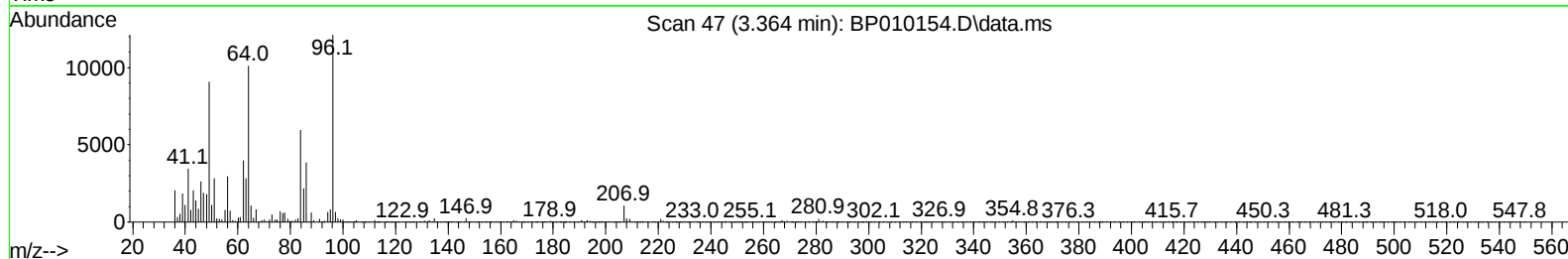
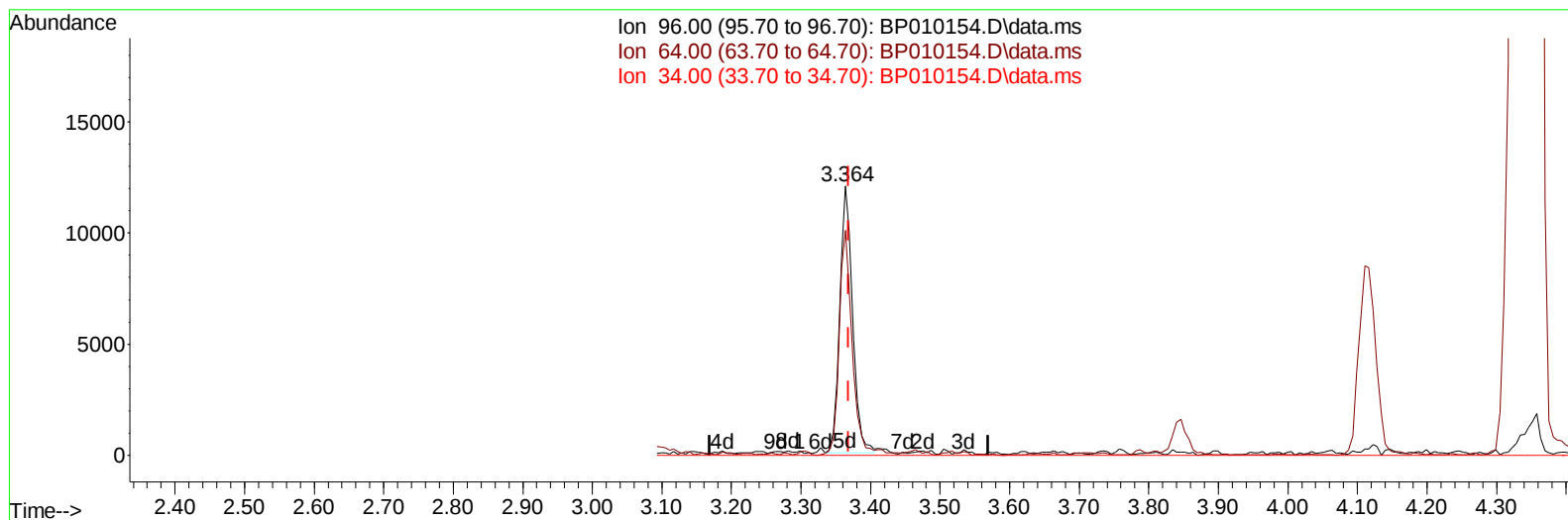
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(3) 1,4-Dioxane-d8 (S)

3.364min (-0.005) 4.72 ng/uL m

response 15614

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	83.51#
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	7.928	152	144405	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.734	136	565447	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.563	164	353185	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.310	188	747427	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.398	240	721319	20.000	ng/ul	#-0.01
88) Perylene-d12	23.845	264	715576	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	15614m	4.715	ng/uL	0.00
4) Pyridine-d5	3.787	84	93113	9.995	ng/ul	0.00
7) Phenol-d5	7.099	99	103169	7.855	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	265110	33.637	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	264290	27.169	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	204266	18.770	ng/ul	-0.01
21) Nitrobenzene-d5	9.087	128	166420	37.005	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.828	143	176419	35.919	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	290105	30.340	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	374059	27.284	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	1045592	38.058	ng/ul	-0.01
49) Acenaphthylene-d8	14.257	160	1224528	36.827	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	36289	7.175	ng/ul	0.00
60) Fluorene-d10	15.551	176	910253	38.516	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.669	200	163055	34.501	ng/ul	-0.01
73) Anthracene-d10	17.410	188	1406689	38.902	ng/ul	-0.01
81) Pyrene-d10	19.645	212	1657172	40.397	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1528986	40.757	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.128	94	13263	1.005	ng/ul#	92

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

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