

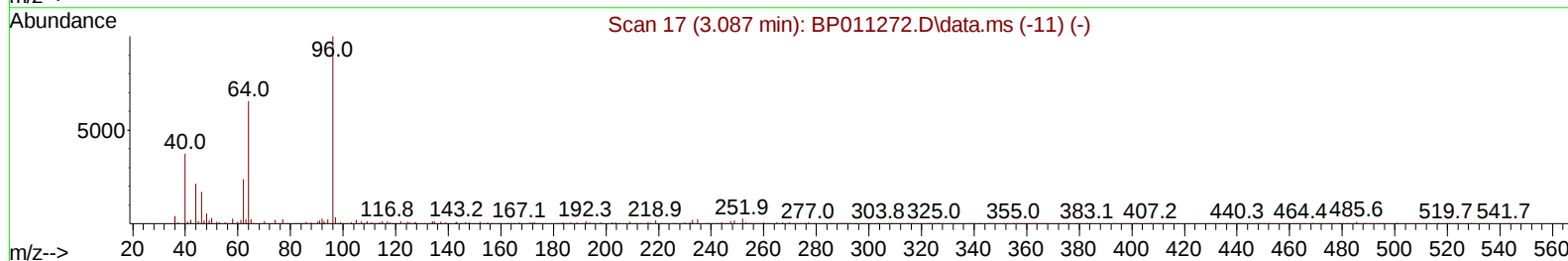
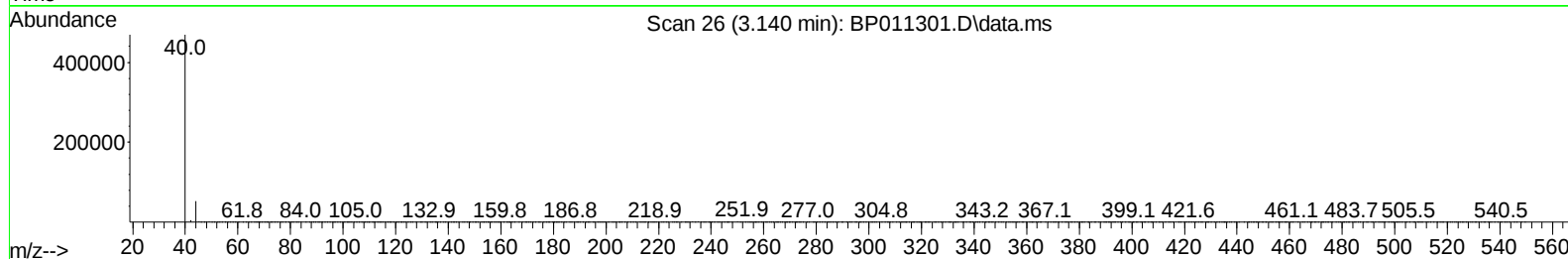
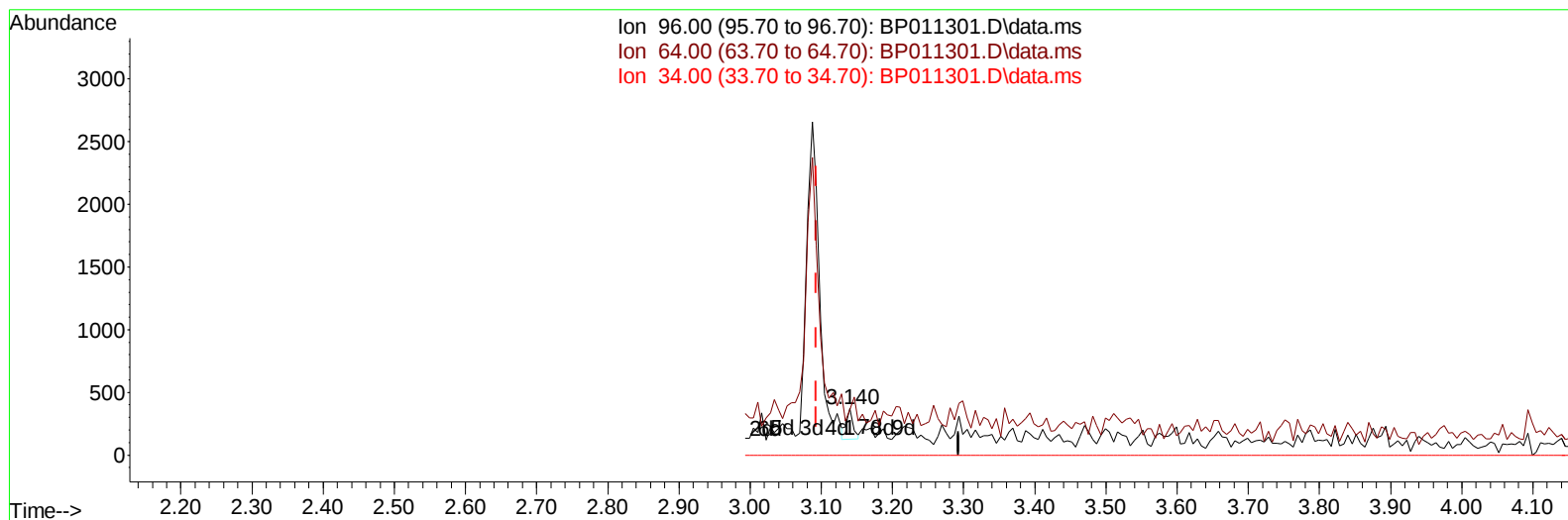
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
 Data File : BP011301.D
 Acq On : 06 Aug 2022 05:31
 Operator : CG/JU
 Sample : PB146785BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK785

Manual IntegrationsAPPROVED

Quant Time: Aug 06 06:48:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 02 23:52:37 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/08/2022
 Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011301.D\data.ms

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

3.140min (+ 0.047) 0.10 ng/uL

response 157

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	66.79
34.00	0.00	0.00
0.00	0.00	0.00

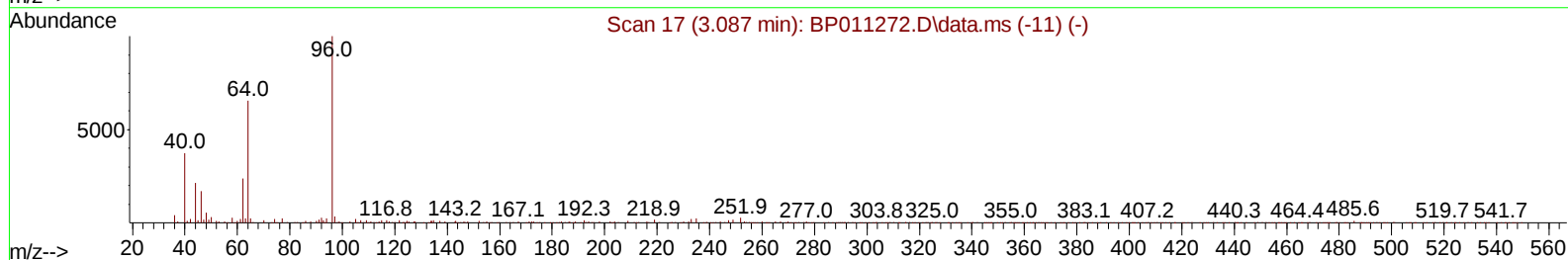
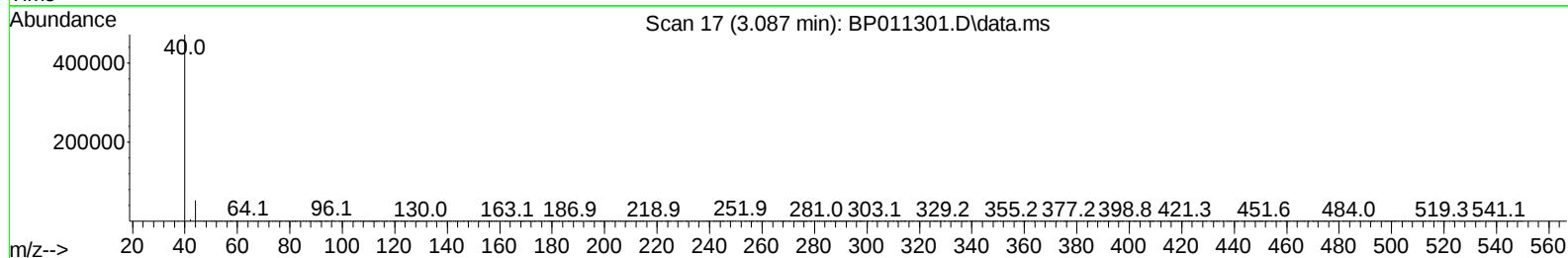
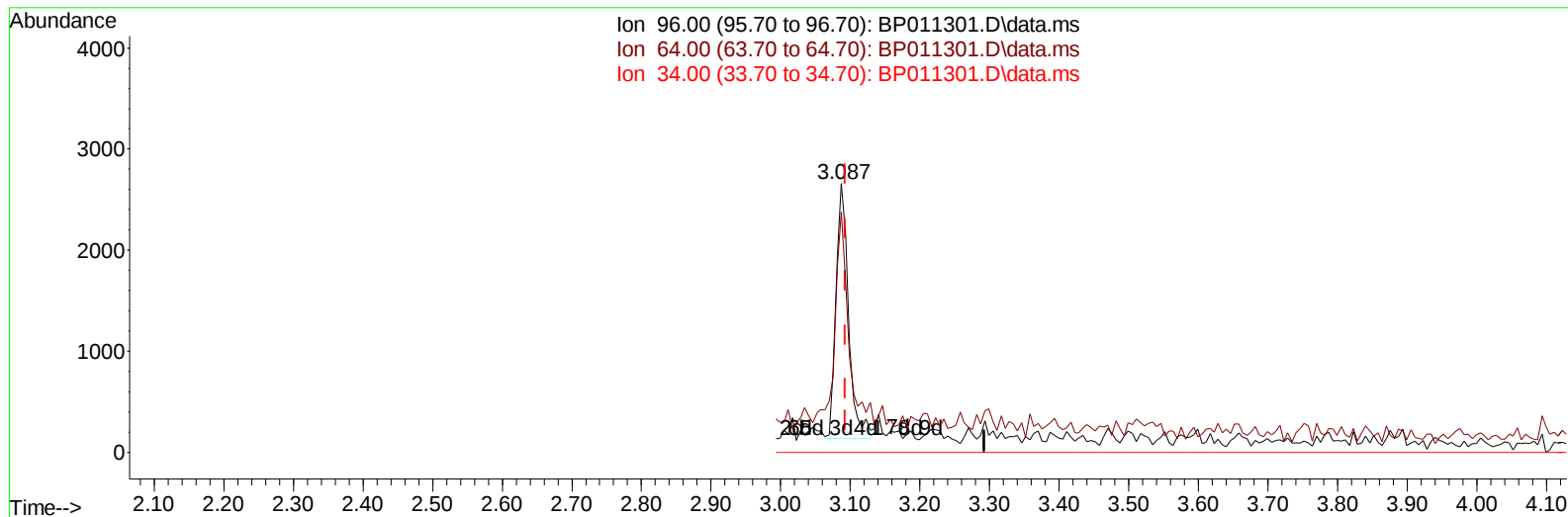
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(3) 1,4-Dioxane-d8 (S)

3.087min (-0.006) 2.01 ng/uL m

response 3148

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	89.42#
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	7.593	152	126695	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.387	136	535491	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.269	164	344466	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.039	188	716431	20.000	ng/ul	0.00
79) Chrysene-d12	21.157	240	847830	20.000	ng/ul	-0.02
88) Perylene-d12	23.468	264	807633	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	3148m	2.009	ng/uL	0.00
4) Pyridine-d5	3.499	84	39158	9.265	ng/ul	0.00
7) Phenol-d5	6.775	99	185998	17.616	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.940	67	188892	24.793	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.128	132	162941	18.933	ng/ul	-0.01
15) 4-Methylphenol-d8	8.316	113	152895	17.599	ng/ul	-0.01
21) Nitrobenzene-d5	8.757	128	137347	35.659	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.475	143	90260	25.000	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.010	165	137000	19.899	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.540	131	161592	15.171	ng/ul	0.00
46) Dimethylphthalate-d6	13.692	166	457859	19.407	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	751141	25.305	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.492	143	74583	18.590	ng/ul	-0.01
60) Fluorene-d10	15.269	176	671546	33.452	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.398	200	79045	23.761	ng/ul	-0.02
73) Anthracene-d10	17.133	188	1240968	39.413	ng/ul	-0.01
81) Pyrene-d10	19.398	212	1309121	30.539	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.321	264	1202163	29.871	ng/ul	-0.02

Target Compounds Qvalue

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

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