

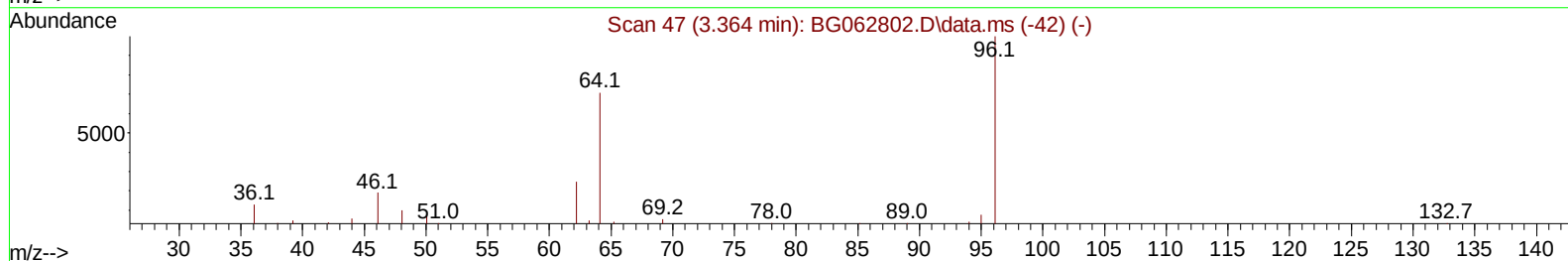
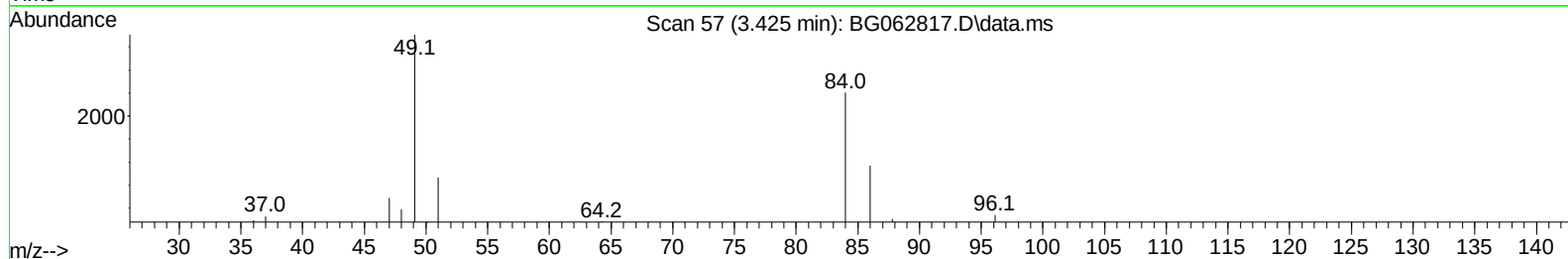
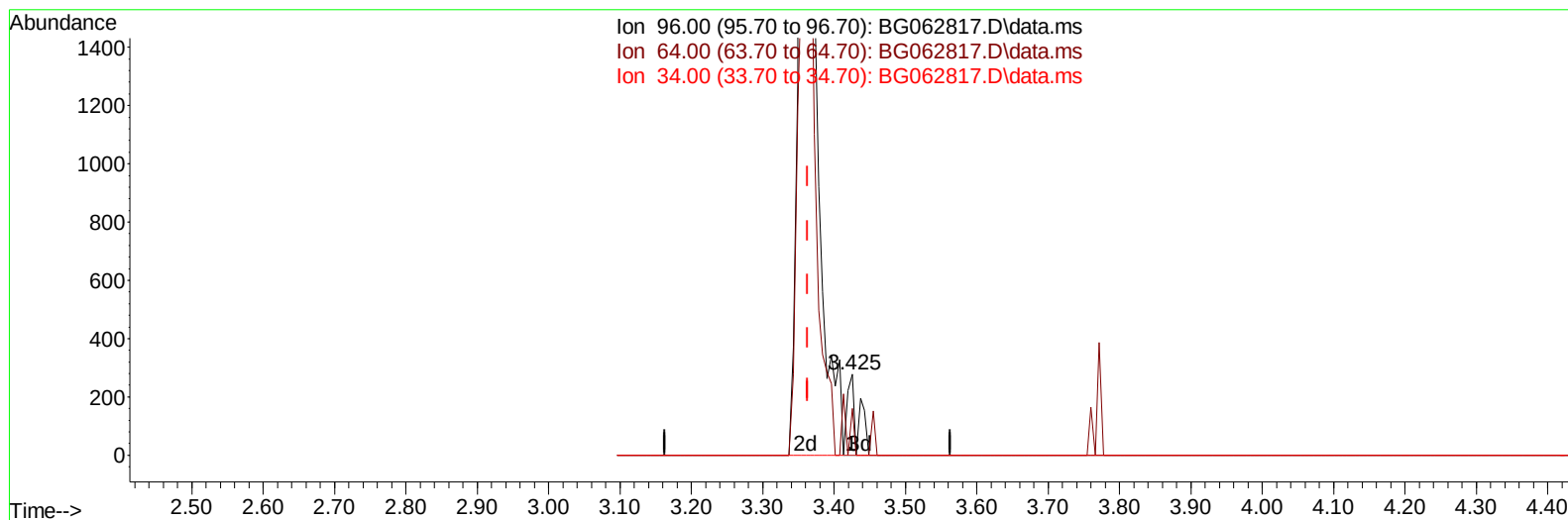
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062817.D
Acq On : 14 Sep 2024 18:41
Operator : JU/RC
Sample : PB163390BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK390

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/16/2024
Supervised By :mohammad ahmed 09/17/2024

Quant Time: Sep 14 23:17:04 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 10 01:28:17 2024
Response via : Initial Calibration



TIC: BG062817.D\data.ms

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

3.425min (+ 0.062) 0.18 ng/uL

response 176

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	62.00	57.91
34.00	0.00	0.00
0.00	0.00	0.00

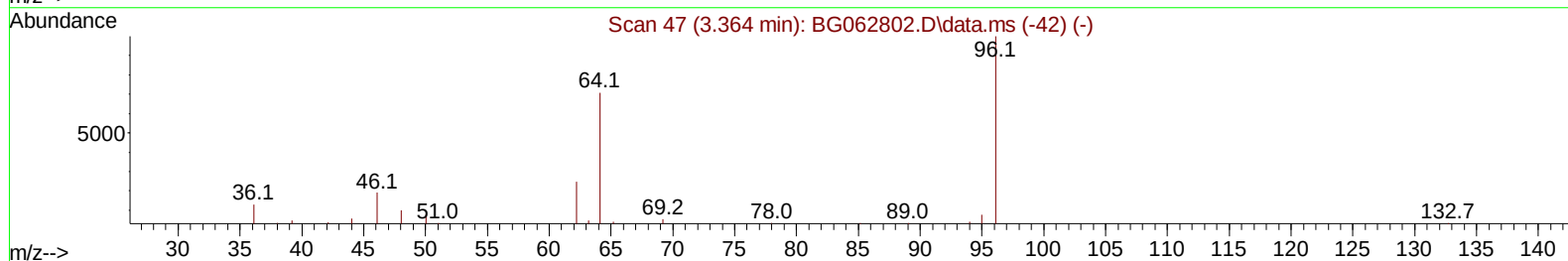
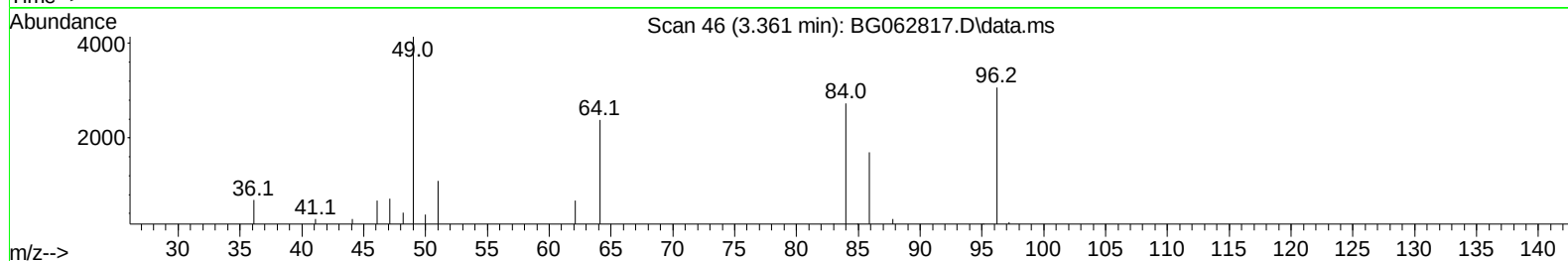
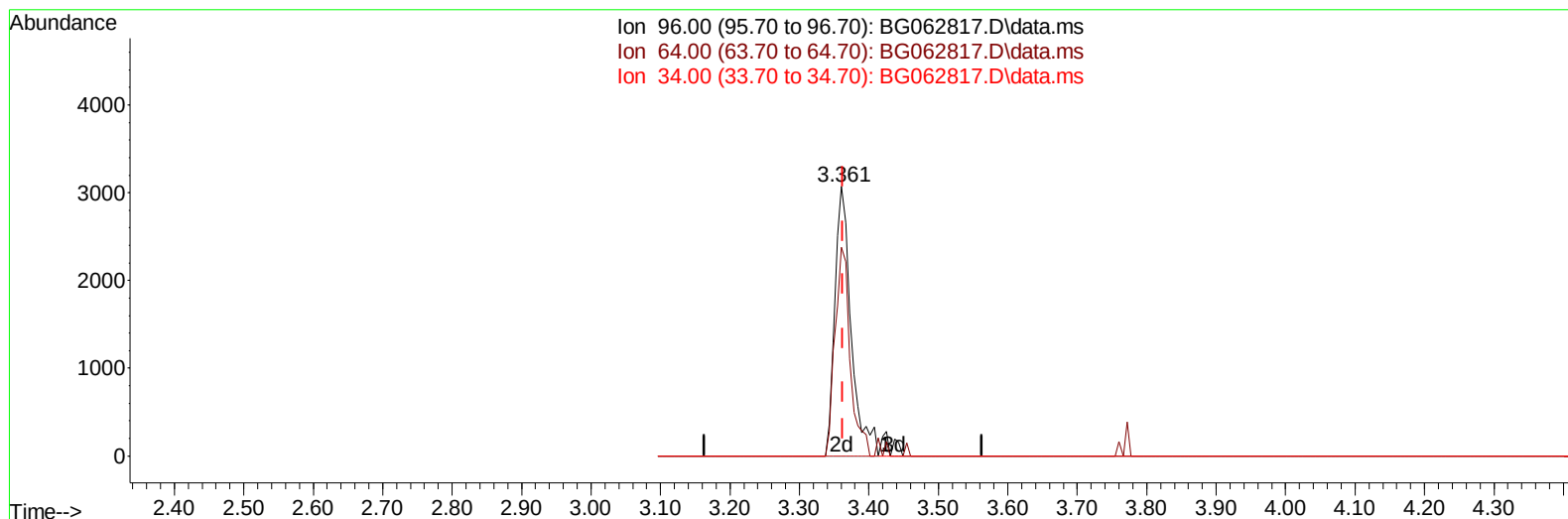
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TIC: BG062817.D\data.ms

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

3.361min (-0.002) 5.20 ng/uL m

response 4999

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	62.00	77.50#
34.00	0.00	0.00
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/16/2024
 Supervised By :mohammad ahmed 09/17/2024

Quant Time: Sep 15 00:54:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 01:28:17 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.902	152	43452	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	177907	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.530	164	151964	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.274	188	406409	20.000	ng/ul	0.00
79) Chrysene-d12	21.522	240	457877	20.000	ng/ul	0.00
88) Perylene-d12	24.583	264	504153	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.361	96	4999m	5.201	ng/uL	0.00
4) Pyridine-d5	3.766	84	74003	24.185	ng/ul	0.00
7) Phenol-d5	7.068	99	92007	24.147	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.233	67	55565	24.324	ng/ul	0.00
11) 2-Chlorophenol-d4	7.438	132	72558	26.143	ng/ul	0.00
15) 4-Methylphenol-d8	8.607	113	78756	24.804	ng/ul	0.00
21) Nitrobenzene-d5	9.054	128	38460	30.247	ng/ul	0.00
24) 2-Nitrophenol-d4	9.777	143	46931	33.749	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.317	165	80590	26.611	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	109469	26.297	ng/ul	0.00
46) Dimethylphthalate-d6	13.936	166	321028	27.433	ng/ul	0.00
49) Acenaphthylene-d8	14.224	160	346766	26.681	ng/ul	0.00
54) 4-Nitrophenol-d4	14.730	143	51703	26.008	ng/ul	0.00
60) Fluorene-d10	15.523	176	290412	27.540	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	56008	23.958	ng/ul	0.00
73) Anthracene-d10	17.374	188	538509	28.076	ng/ul	0.00
81) Pyrene-d10	19.665	212	723156	28.066	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.377	264	735220	27.618	ng/ul	0.00

Target Compounds Qvalue

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

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