

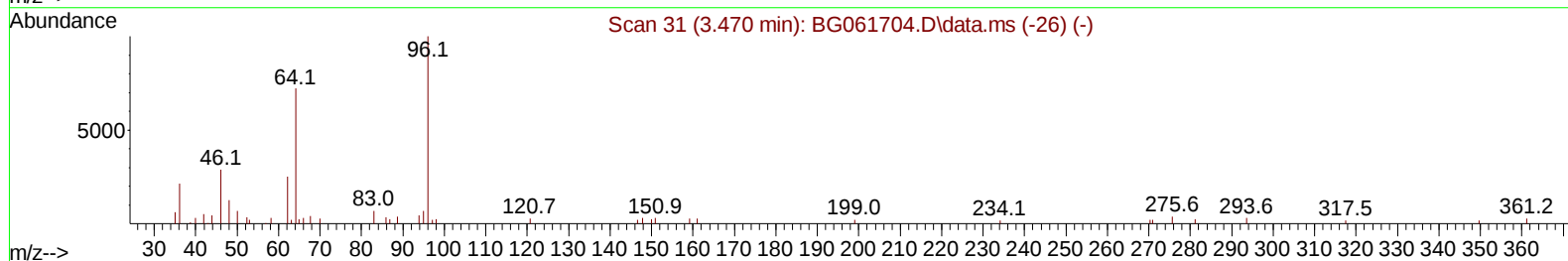
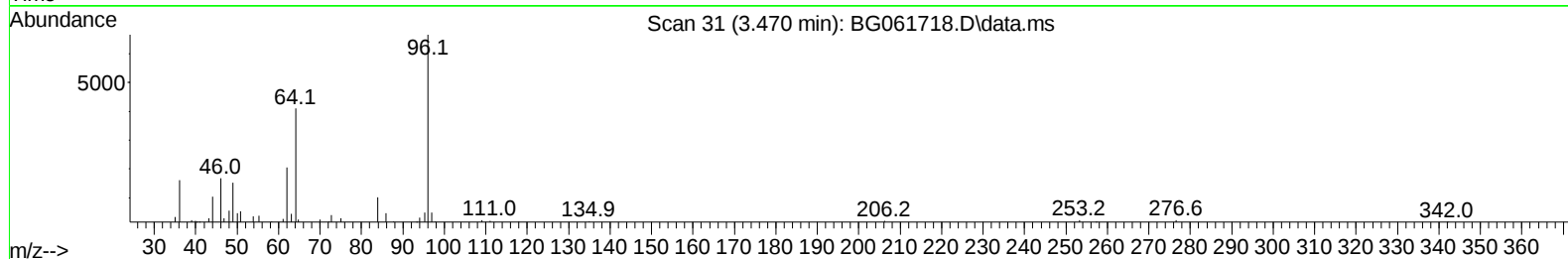
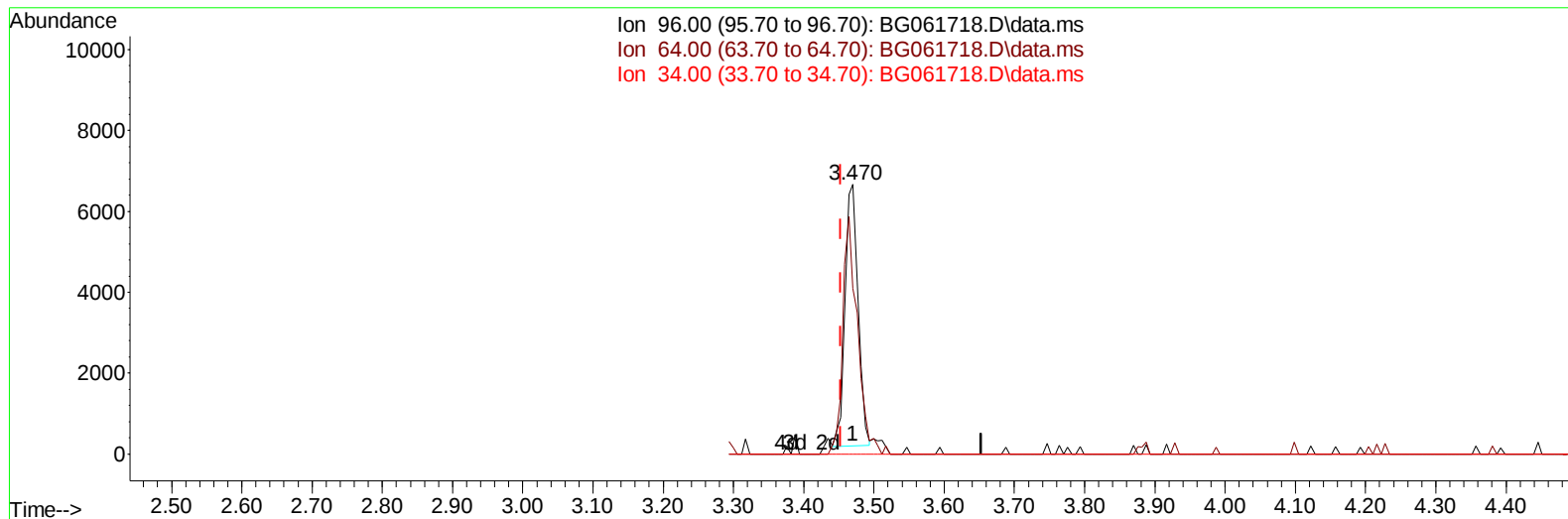
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061718.D
Acq On : 26 Jun 2024 1:26
Operator : MA/JU
Sample : P2873-19
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0SK8

Manual IntegrationsAPPROVED

Quant Time: Jun 28 07:02:38 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 27 12:23:53 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/26/2024
Supervised By :mohammad ahmed 06/27/2024



TIC: BG061718.D\data.ms

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

3.470min (+ 0.017) 5.03 ng/uL

response 8328

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	61.45
34.00	0.00	0.00
0.00	0.00	0.00

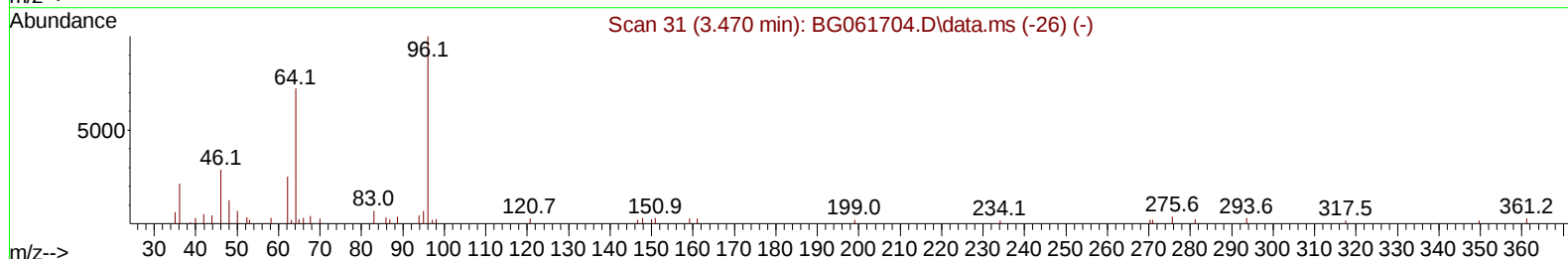
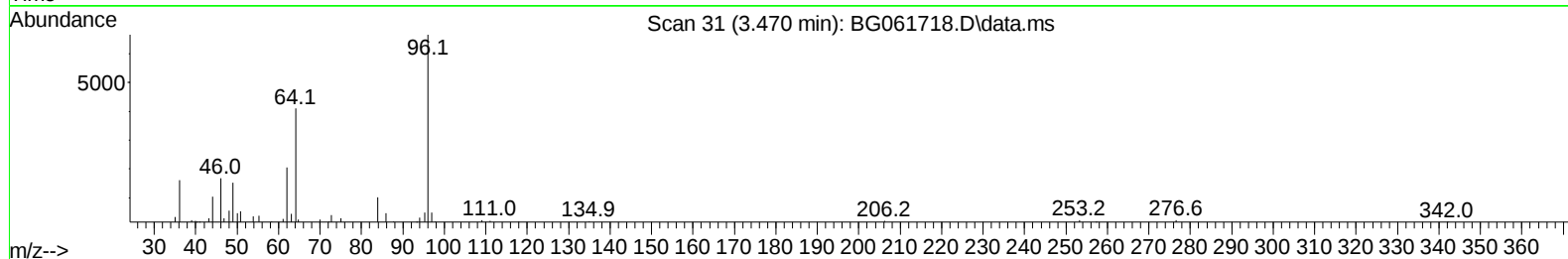
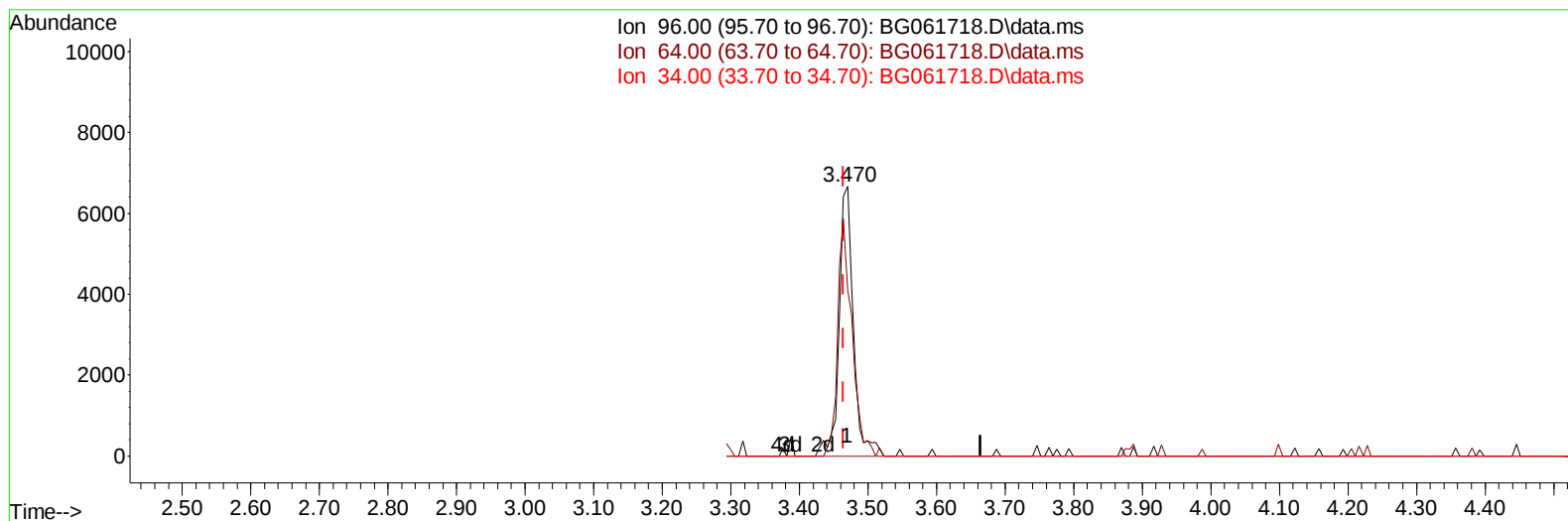
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Manual IntegrationsAPPROVED

Quant Time: Jun 26 02:43:07 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 21 05:03:55 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/26/2024
Supervised By :mohammad ahmed 06/27/2024



TIC: BG061718.D\data.ms

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

3.470min (+ 0.006) 5.84 ng/uL m

response 9664

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	61.45
34.00	0.00	0.00
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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 QLast Update : Fri Jun 21 05:03:55 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.071	152	59789	20.000	ng/ul	0.00
20) Naphthalene-d8	10.885	136	296884	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.698	164	218916	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.442	188	523679	20.000	ng/ul	0.00
79) Chrysene-d12	21.708	240	440111	20.000	ng/ul	0.00
88) Perylene-d12	24.927	264	527882	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.470	96	9664m	5.839	ng/uL	0.00
4) Pyridine-d5	3.887	84	65533	13.096	ng/ul	0.00
7) Phenol-d5	7.213	99	215671	33.228	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.395	67	130389	32.225	ng/ul	0.00
11) 2-Chlorophenol-d4	7.601	132	140204	31.478	ng/ul	0.00
15) 4-Methylphenol-d8	8.764	113	179130	35.700	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	81062	40.213	ng/ul	0.00
24) 2-Nitrophenol-d4	9.957	143	98727	48.357	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.491	165	183662	37.991	ng/ul	0.00
31) 4-Chloroaniline-d4	11.014	131	241411	32.983	ng/ul	0.00
46) Dimethylphthalate-d6	14.105	166	651752	38.688	ng/ul	0.00
49) Acenaphthylene-d8	14.393	160	744278	40.314	ng/ul	0.00
54) 4-Nitrophenol-d4	14.874	143	109079	37.701	ng/ul	0.00
60) Fluorene-d10	15.691	176	607114	41.165	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.797	200	88683	39.603	ng/ul	0.00
73) Anthracene-d10	17.542	188	967502	40.475	ng/ul	0.00
81) Pyrene-d10	19.821	212	1035385	42.117	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.716	264	1055215	41.853	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.932	128	431019	25.947	ng/ul	98
36) 2-Methylnaphthalene	12.536	142	31044	2.648	ng/ul	97
37) 1-Methylnaphthalene	12.753	142	18645	1.491	ng/ul#	90

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

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