

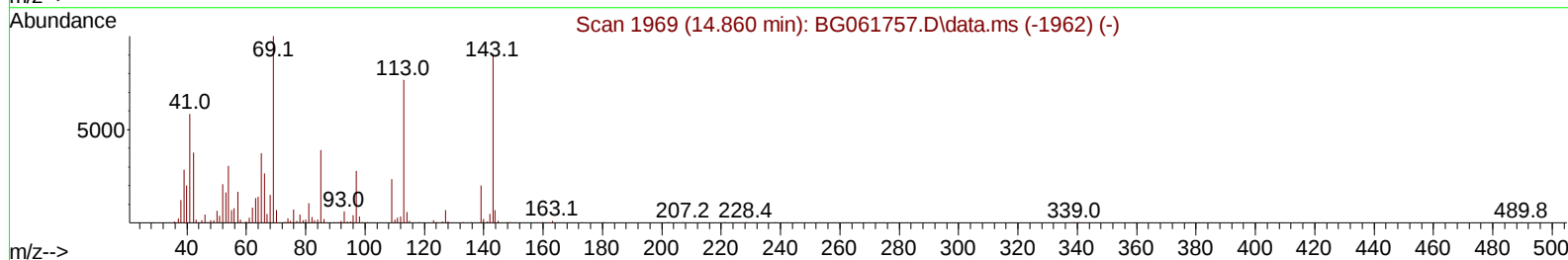
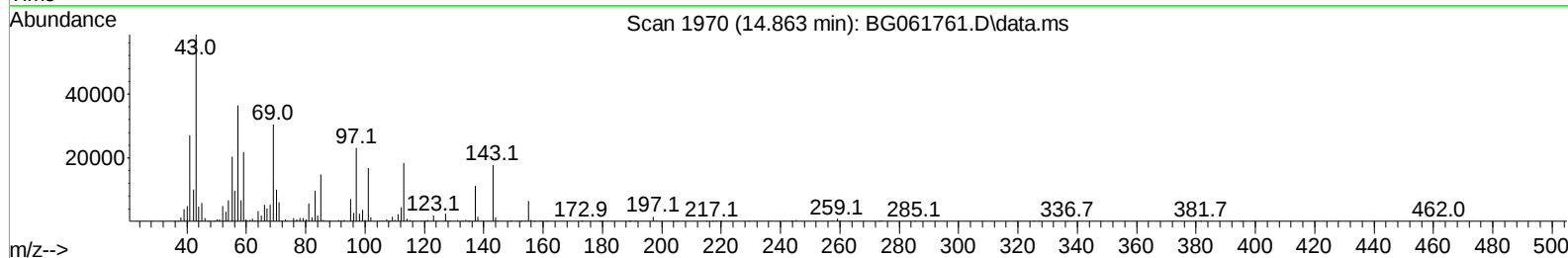
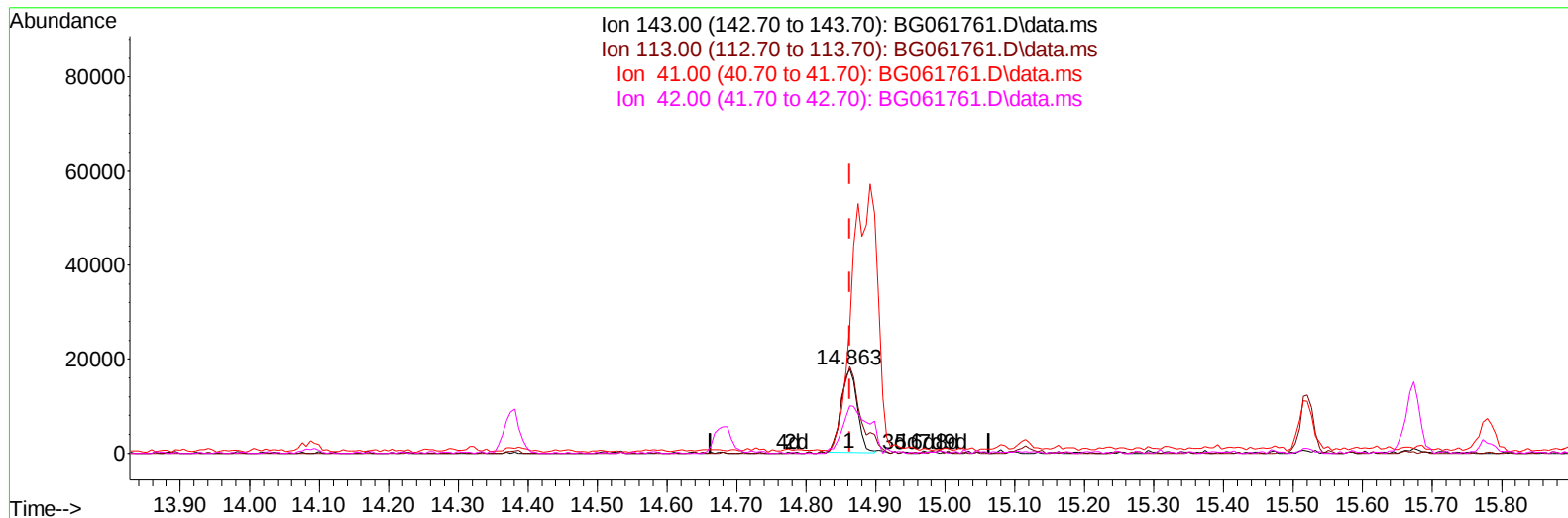
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062724\
Data File : BG061761.D
Acq On : 28 Jun 2024 6:33
Operator : MA/JU
Sample : P2898-08
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E1GB1

Manual IntegrationsAPPROVED

Quant Time: Jun 28 07:08:31 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 06/28/2024
Supervised By :mohammad ahmed 06/29/2024



TIC: BG061761.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (-0.000) 10.27 ng/u1

response 28954

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	92.40	102.91
41.00	58.60	152.24#
42.00	35.10	56.45#

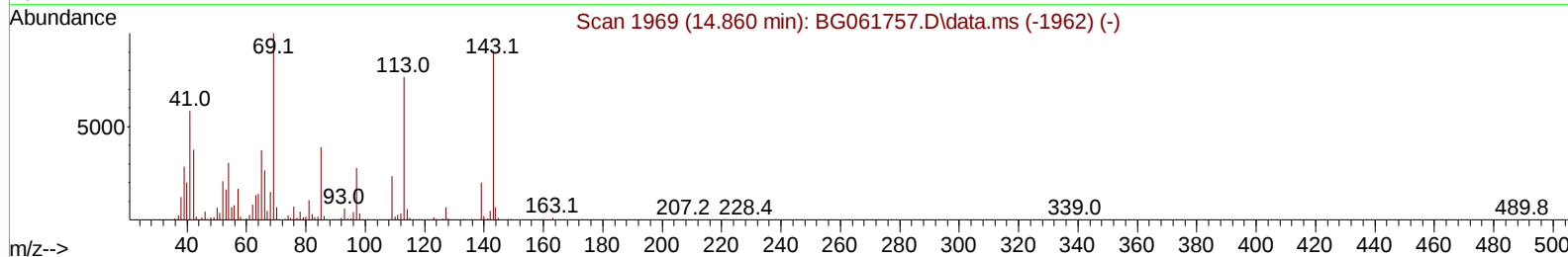
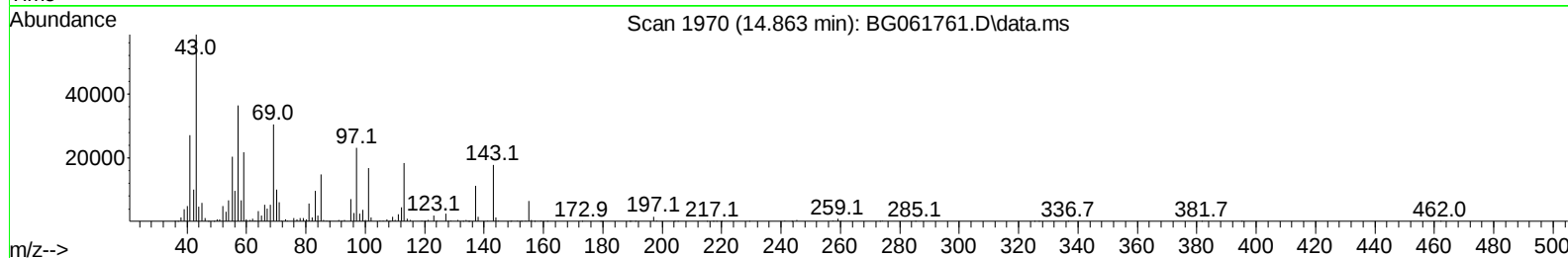
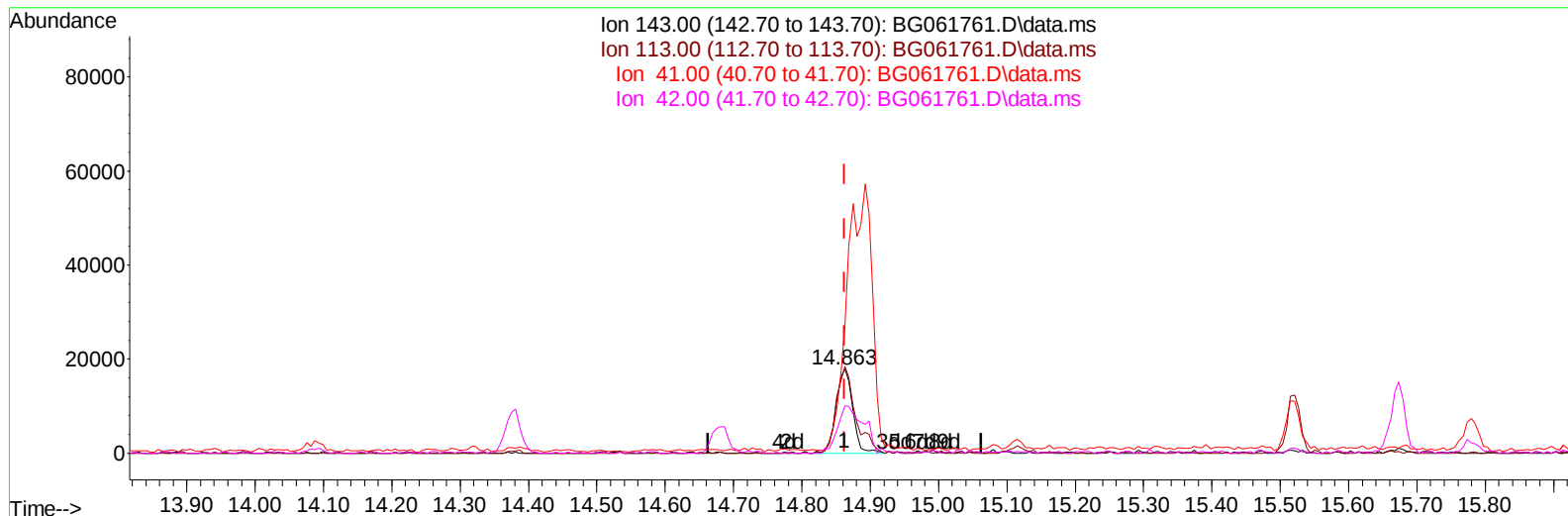
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(54) 4-Nitrophenol-d4 (S)

14.863min (-0.000) 10.72 ng/ul m

response 30205

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	92.40	102.91
41.00	58.60	152.24#
42.00	35.10	56.45#

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/28/2024
 Supervised By :mohammad ahmed 06/29/2024

Quant Time: Jun 28 07:09:26 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 27 12:23:53 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.059	152	70160	20.000	ng/ul	0.00
20) Naphthalene-d8	10.868	136	335393	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.681	164	213276	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.425	188	485637	20.000	ng/ul	0.00
79) Chrysene-d12	21.684	240	375850	20.000	ng/ul	0.00
88) Perylene-d12	24.892	264	434698	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.459	96	10729	5.524	ng/uL	0.00
4) Pyridine-d5	3.876	84	79553	13.548	ng/ul	0.00
7) Phenol-d5	7.201	99	75954	9.972	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.378	67	197260	41.546	ng/ul	0.00
11) 2-Chlorophenol-d4	7.583	132	175782	33.632	ng/ul	0.00
15) 4-Methylphenol-d8	8.752	113	126907	21.553	ng/ul	0.00
21) Nitrobenzene-d5	9.222	128	110580	48.558	ng/ul	0.00
24) 2-Nitrophenol-d4	9.939	143	122346	53.045	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.480	165	212194	38.853	ng/ul	0.00
31) 4-Chloroaniline-d4	10.997	131	290734	35.161	ng/ul	0.00
46) Dimethylphthalate-d6	14.087	166	718009	43.748	ng/ul	0.00
49) Acenaphthylene-d8	14.381	160	812862	45.193	ng/ul	0.00
54) 4-Nitrophenol-d4	14.863	143	30205m	10.716	ng/ul	0.00
60) Fluorene-d10	15.674	176	634886	44.187	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.779	200	77103	37.129	ng/ul	0.00
73) Anthracene-d10	17.524	188	1007560	45.452	ng/ul	0.00
81) Pyrene-d10	19.804	212	1079960	51.441	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.681	264	974899	46.957	ng/ul	0.00

Target Compounds Qvalue

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

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