

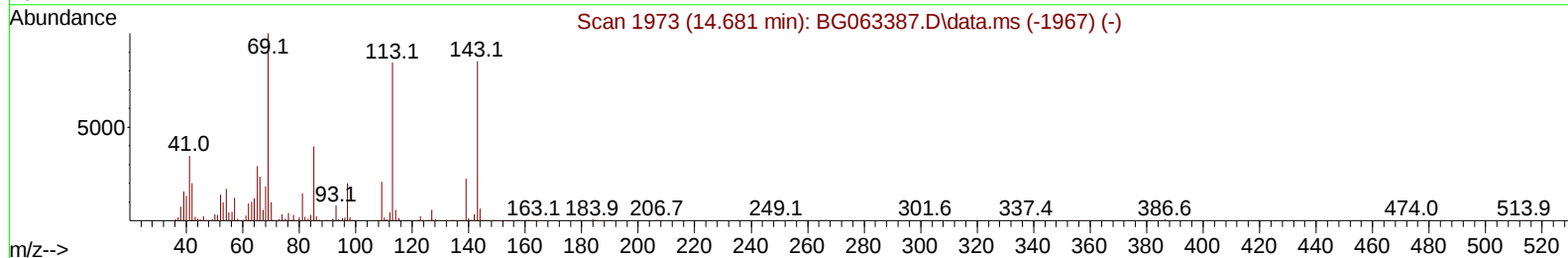
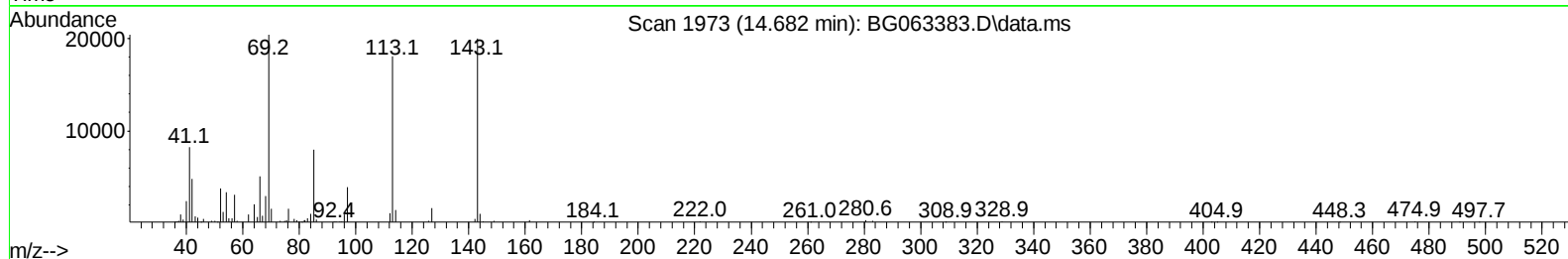
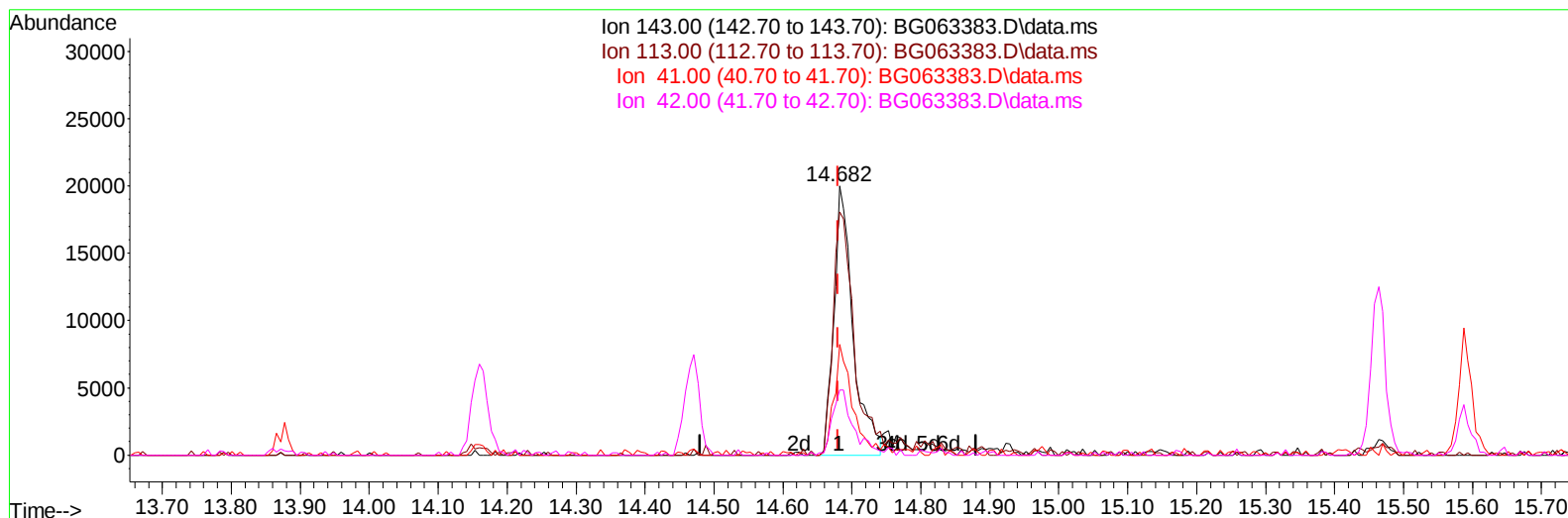
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
Data File : BG063383.D
Acq On : 14 Nov 2024 9:50
Operator : RC/JU
Sample : P4802-09
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GCPD1

Manual IntegrationsAPPROVED

Quant Time: Nov 14 10:22:48 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 06 15:45:52 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/15/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BG063383.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.682min (+ 0.002) 9.47 ng/u1

response 38657

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	90.48#
41.00	42.60	41.16
42.00	30.90	24.12#

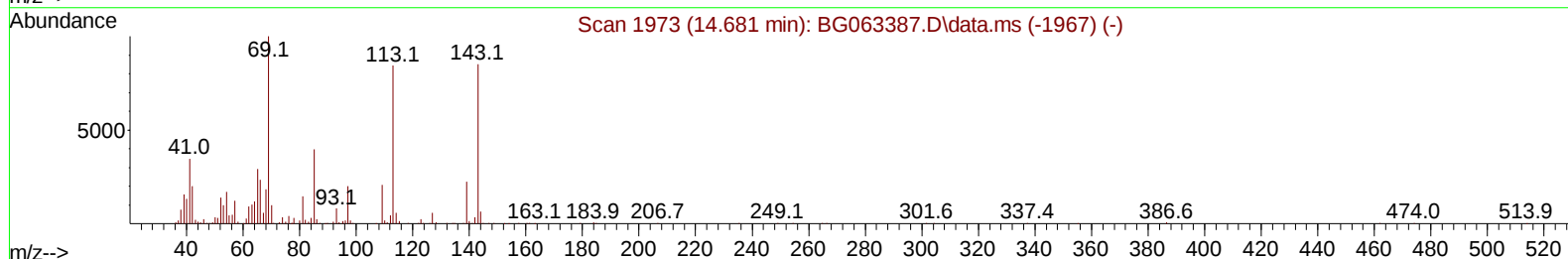
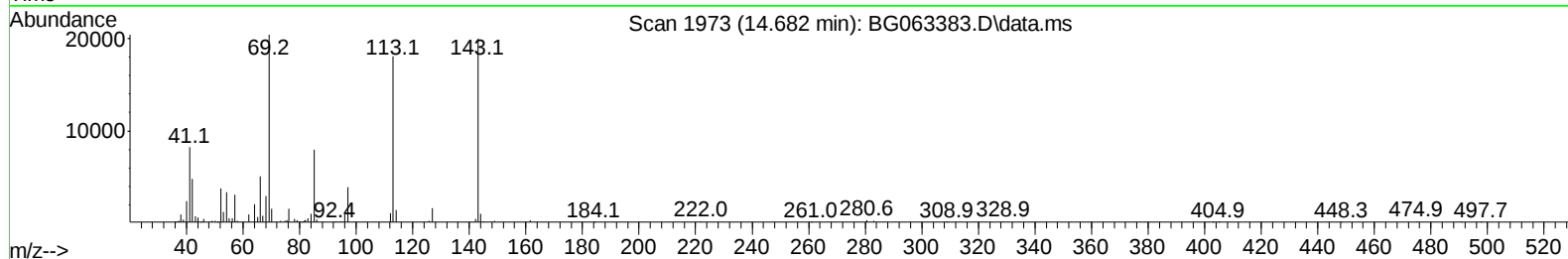
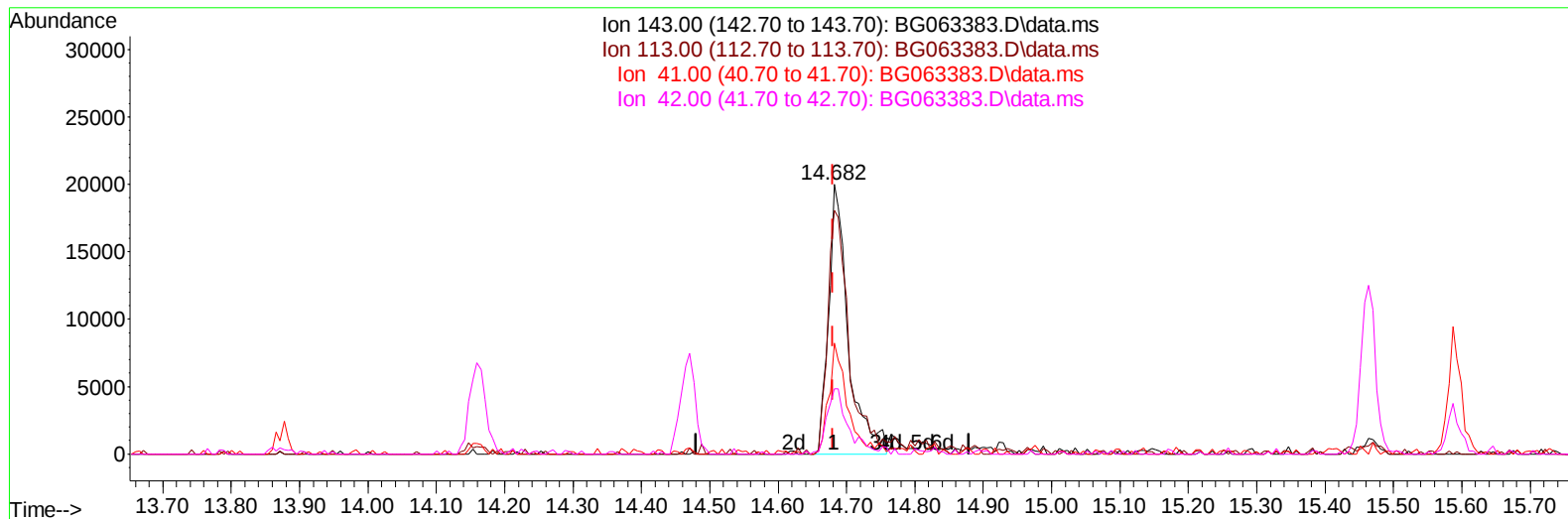
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TIC: BG063383.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.682min (+ 0.002) 9.82 ng/ul m

response 40061

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	90.48#
41.00	42.60	41.16
42.00	30.90	24.12#

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Instrument :
 BNA_G
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 GCPD1

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/15/2024
 Supervised By :mohammad ahmed 11/18/2024

Quant Time: Nov 14 23:39:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 06 15:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.831	152	106476	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	438606	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.471	164	380515	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.220	188	956767	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.474	240	1037997	20.000	ng/ul	0.00
88) Perylene-d12	24.506	264	1083043	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	9684	4.113	ng/uL	0.00
4) Pyridine-d5	3.713	84	41140	5.286	ng/ul	0.00
7) Phenol-d5	7.003	99	85253	8.873	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.162	67	161858	28.721	ng/ul	0.00
11) 2-Chlorophenol-d4	7.367	132	167511	26.861	ng/ul	0.00
15) 4-Methylphenol-d8	8.542	113	150363	18.464	ng/ul	0.00
21) Nitrobenzene-d5	8.983	128	97714	31.689	ng/ul	0.00
24) 2-Nitrophenol-d4	9.706	143	125622	31.521	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	230327	29.450	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	205893	20.212	ng/ul	0.00
46) Dimethylphthalate-d6	13.877	166	925841	31.062	ng/ul	0.00
49) Acenaphthylene-d8	14.159	160	915344	31.376	ng/ul	0.00
54) 4-Nitrophenol-d4	14.682	143	40061m	9.817	ng/ul	0.00
60) Fluorene-d10	15.464	176	822460	32.695	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	192268	31.975	ng/ul	0.00
73) Anthracene-d10	17.320	188	1519283	36.947	ng/ul	0.00
81) Pyrene-d10	19.618	212	1990237	38.243	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.306	264	1974375	37.758	ng/ul	0.00

Target Compounds Qvalue

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

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