

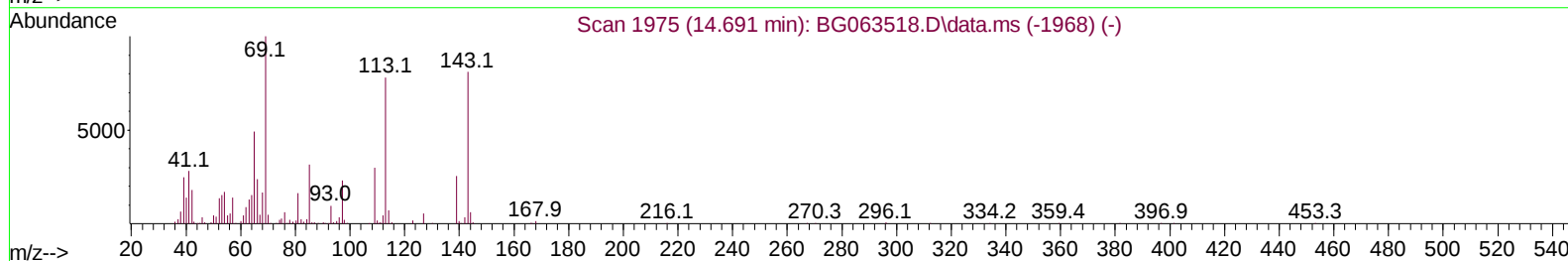
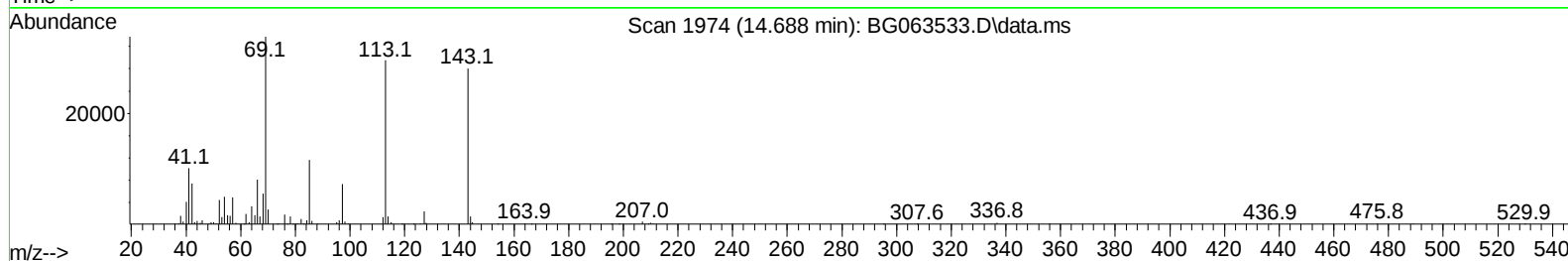
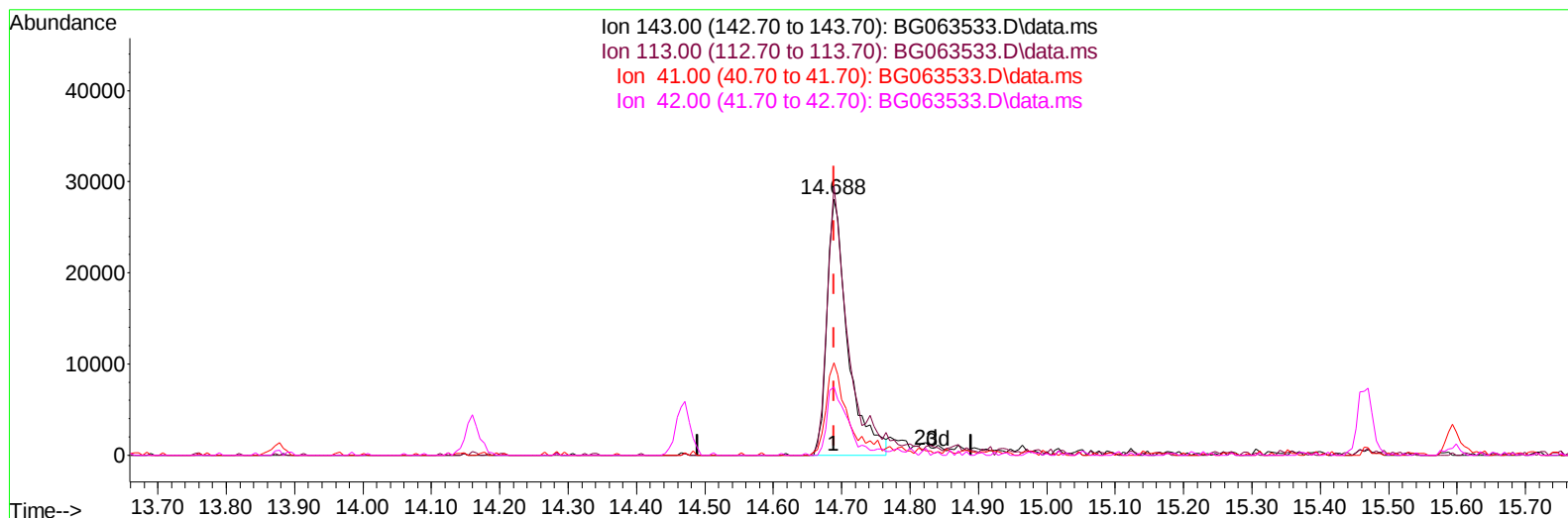
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112224\
Data File : BG063533.D
Acq On : 23 Nov 2024 00:04
Operator : RC/JU
Sample : PB165173BL
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK173

Manual IntegrationsAPPROVED

Quant Time: Nov 23 00:43:57 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 20 14:52:38 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/25/2024
Supervised By :mohammad ahmed 11/27/2024



TIC: BG063533.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.688min (-0.001) 20.09 ng/u1

response 59553

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	105.07
41.00	38.80	36.08
42.00	26.30	26.61

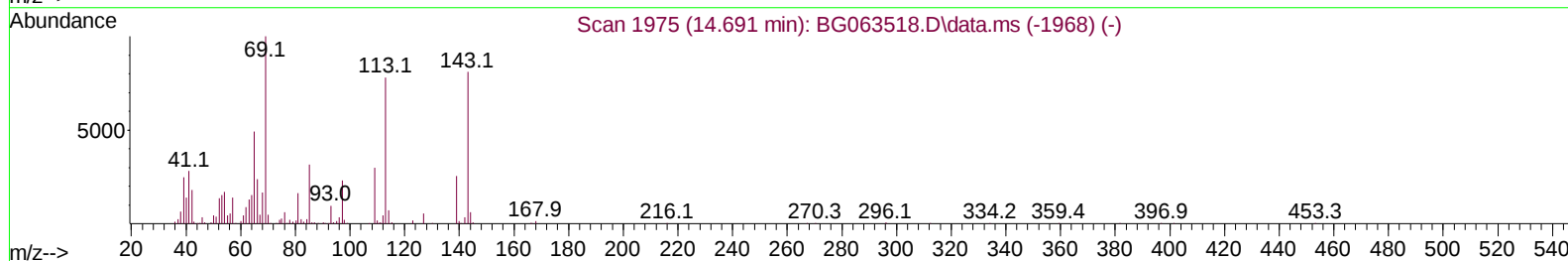
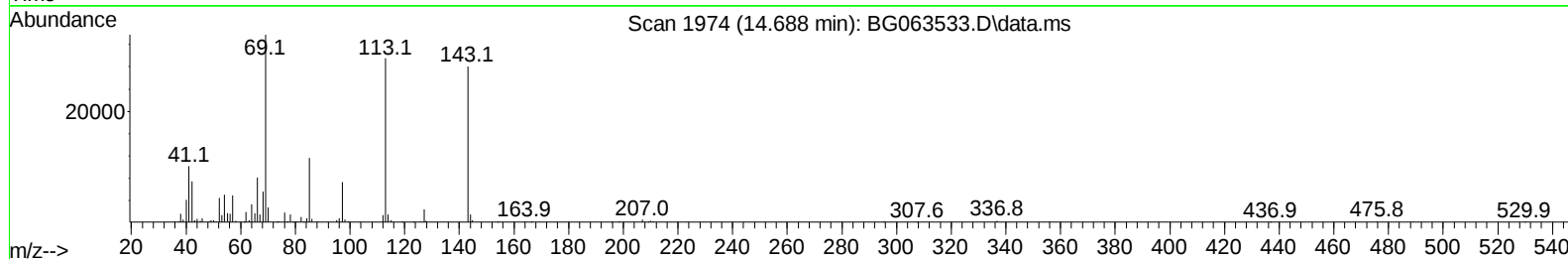
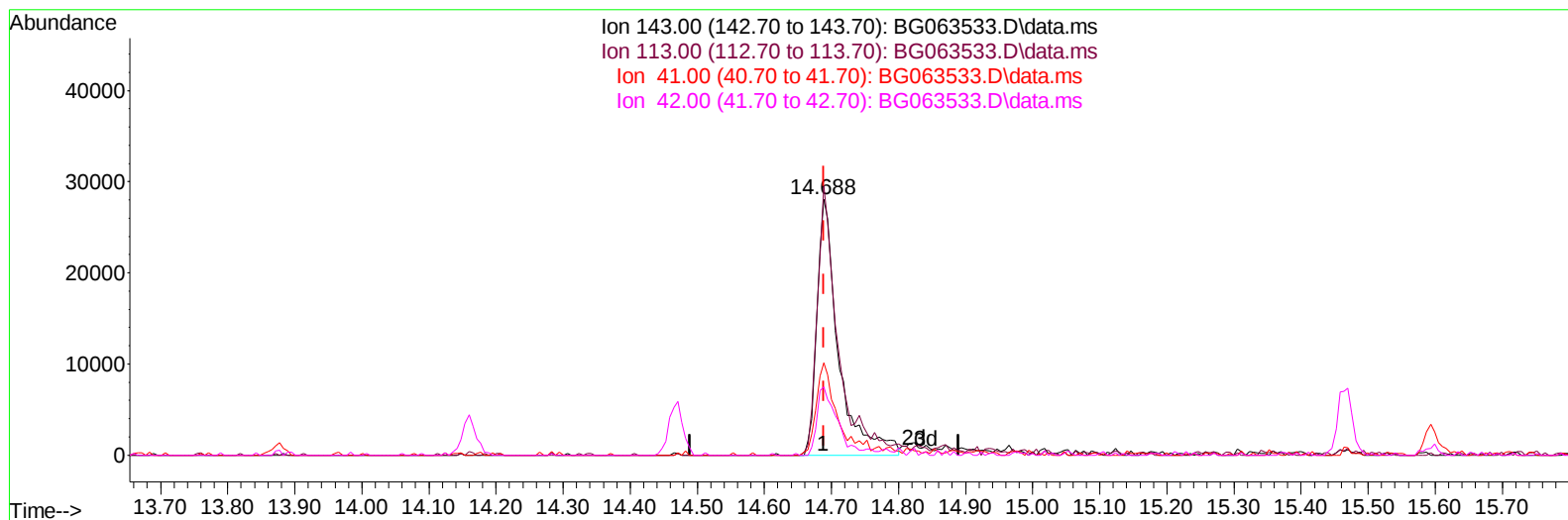
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TIC: BG063533.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.688min (-0.001) 21.18 ng/ul m

response 62779

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	105.07
41.00	38.80	36.08
42.00	26.30	26.61

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/25/2024
 Supervised By :mohammad ahmed 11/27/2024

Quant Time: Nov 23 01:28:13 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 20 14:52:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.826	152	85317	20.000	ng/ul	0.00
20) Naphthalene-d8	10.617	136	357417	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.471	164	293921	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.227	188	739264	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	851696	20.000	ng/ul	0.00
88) Perylene-d12	24.524	264	899499	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.302	96	10937	5.060	ng/uL	0.00
4) Pyridine-d5	3.701	84	147670	21.795	ng/ul	0.00
7) Phenol-d5	7.003	99	176832	25.201	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.162	67	114289	25.395	ng/ul	0.00
11) 2-Chlorophenol-d4	7.367	132	124180	24.972	ng/ul	0.00
15) 4-Methylphenol-d8	8.537	113	148288	25.290	ng/ul	0.00
21) Nitrobenzene-d5	8.983	128	65069	25.351	ng/ul	0.00
24) 2-Nitrophenol-d4	9.706	143	81370	25.662	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	141876	23.442	ng/ul	0.00
31) 4-Chloroaniline-d4	10.758	131	191829	25.060	ng/ul	0.00
46) Dimethylphthalate-d6	13.878	166	538621	26.995	ng/ul	0.00
49) Acenaphthylene-d8	14.160	160	578990	25.446	ng/ul	0.00
54) 4-Nitrophenol-d4	14.688	143	62779m	21.175	ng/ul	0.00
60) Fluorene-d10	15.470	176	491393	27.104	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	86991	20.931	ng/ul	0.00
73) Anthracene-d10	17.321	188	870308	27.098	ng/ul	0.00
81) Pyrene-d10	19.624	212	1180855	27.659	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.318	264	1186607	27.198	ng/ul	0.00

Target Compounds Qvalue

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

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