

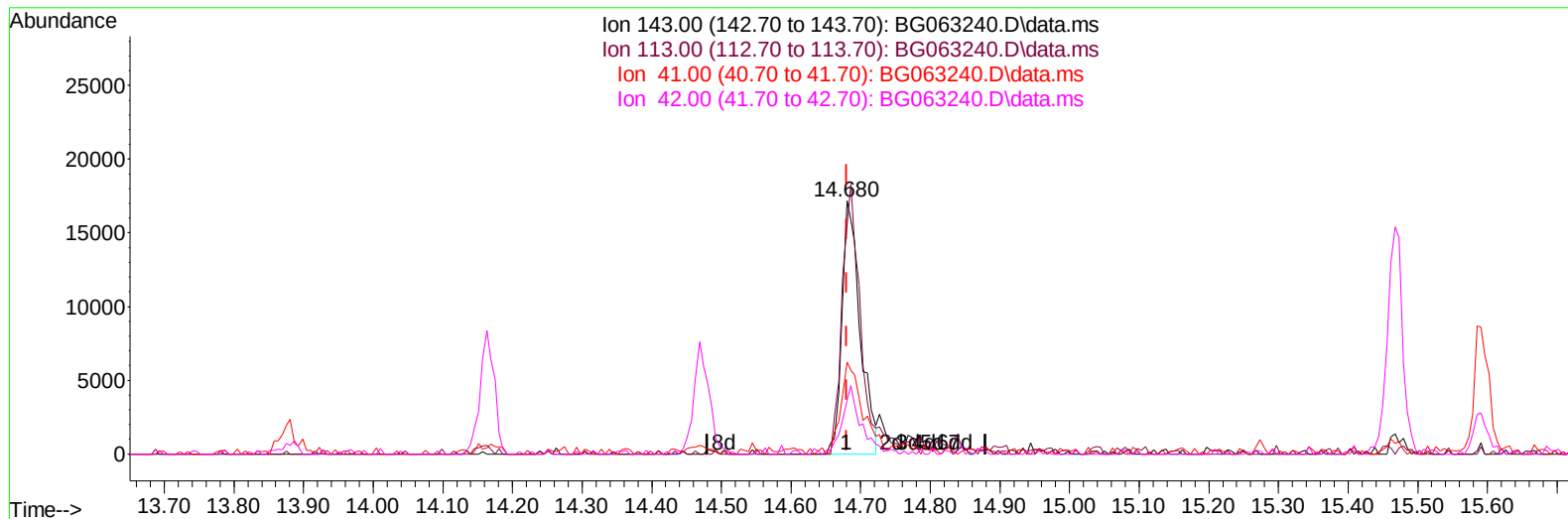
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110824\
Data File : BG063240.D
Acq On : 8 Nov 2024 22:25
Operator : RC/JU
Sample : P4627-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0CD1

Manual IntegrationsAPPROVED

Quant Time: Nov 08 22:23:45 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 11/10/2024
Supervised By :mohammad ahmed 11/11/2024



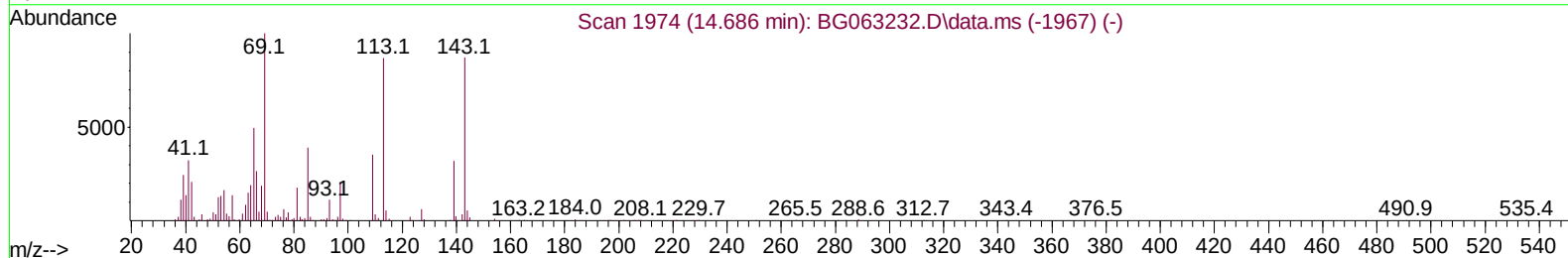
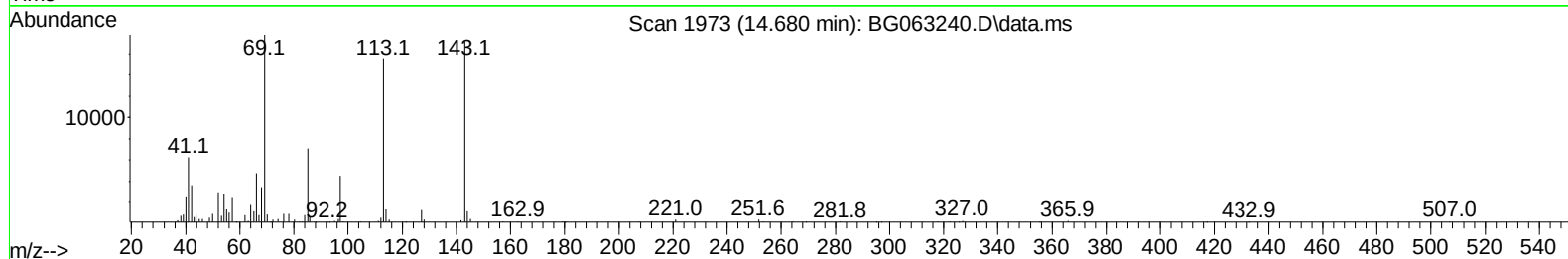
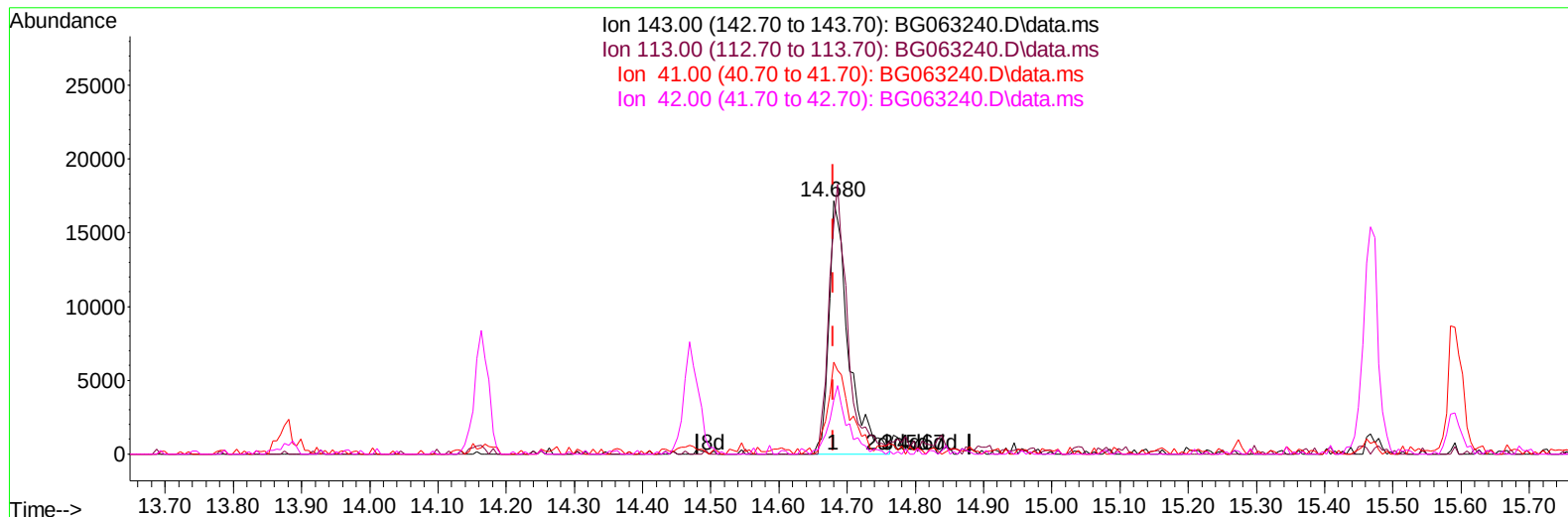
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TIC: BG063240.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.680min (-0.000) 7.29 ng/ul m

response 34202

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	123.00	90.76#
41.00	42.60	36.39
42.00	30.90	21.26#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.835	152	107703	20.000	ng/ul	0.00
20) Naphthalene-d8	10.626	136	500681	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.474	164	437369	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.224	188	1072898	20.000	ng/ul	0.00
79) Chrysene-d12	21.478	240	1089540	20.000	ng/ul	0.00
88) Perylene-d12	24.516	264	1137972	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	9952	4.178	ng/uL	0.00
4) Pyridine-d5	3.711	84	63129	8.019	ng/ul	0.00
7) Phenol-d5	7.007	99	98457	10.131	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.165	67	173522	30.440	ng/ul	0.00
11) 2-Chlorophenol-d4	7.371	132	183236	29.048	ng/ul	0.00
15) 4-Methylphenol-d8	8.546	113	180348	21.894	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	113934	32.368	ng/ul	0.00
24) 2-Nitrophenol-d4	9.709	143	144196	31.696	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.250	165	278544	31.199	ng/ul	0.00
31) 4-Chloroaniline-d4	10.761	131	291540	25.071	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1091365	31.856	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	1134066	33.820	ng/ul	0.00
54) 4-Nitrophenol-d4	14.680	143	34202m	7.292	ng/ul	0.00
60) Fluorene-d10	15.467	176	1001497	34.637	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.591	200	223959	33.214	ng/ul	0.00
73) Anthracene-d10	17.324	188	1803334	39.108	ng/ul	0.00
81) Pyrene-d10	19.621	212	2221660	40.670	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.304	264	2205834	40.148	ng/ul	0.00

Target Compounds Qvalue

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

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