

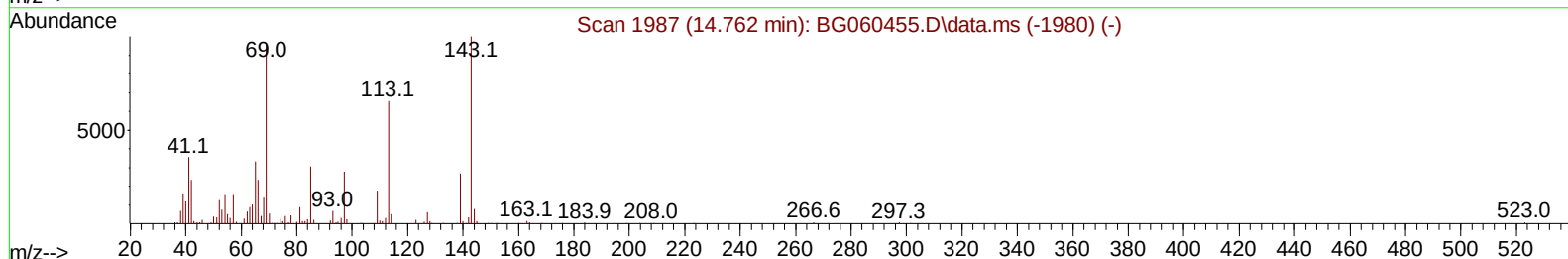
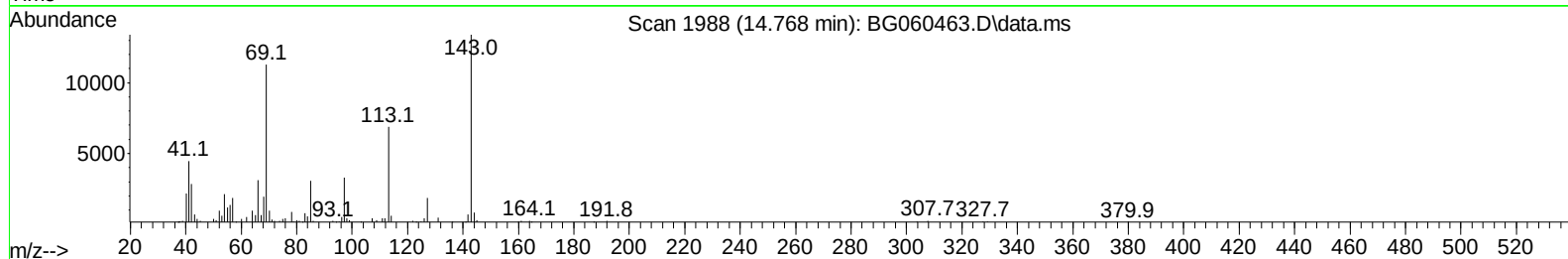
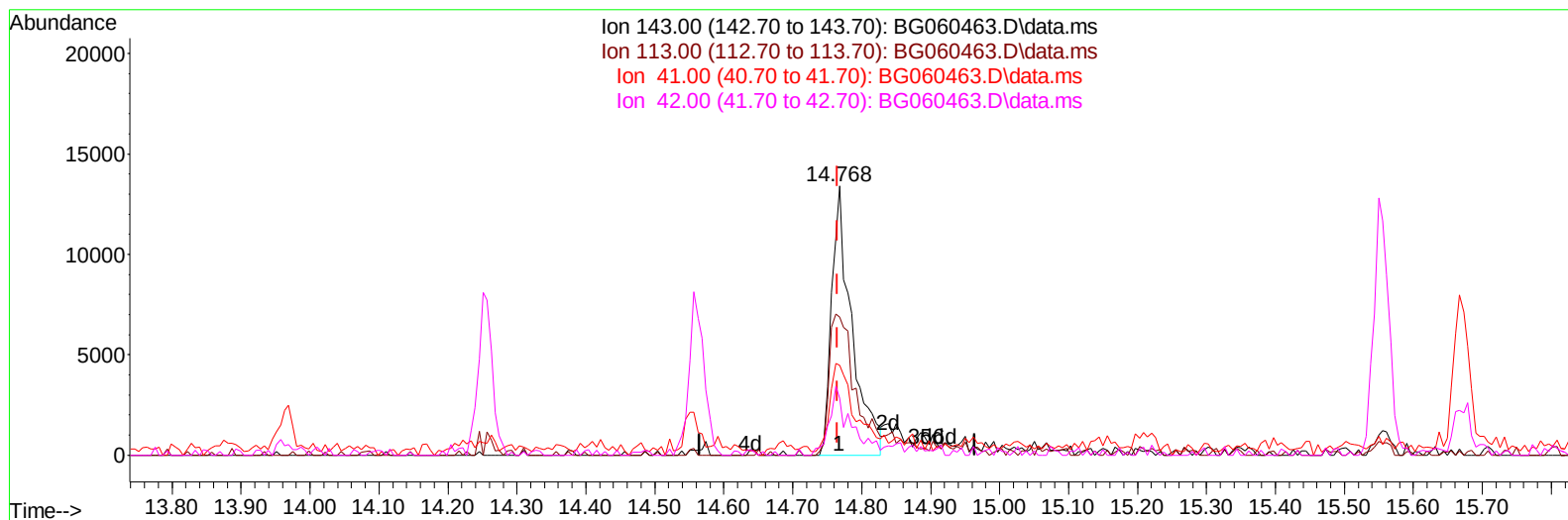
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
Data File : BG060463.D
Acq On : 29 Jan 2024 15:38
Operator : MA/JU
Sample : P1239-11
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0FB5

Manual IntegrationsAPPROVED

Quant Time: Jan 29 16:11:34 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 26 06:06:34 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/30/2024
Supervised By :mohammad ahmed 01/31/2024



TIC: BG060463.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.768min (+ 0.003) 3.65 ng/ul

response 27253

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	51.37
41.00	29.80	33.37
42.00	22.70	21.17

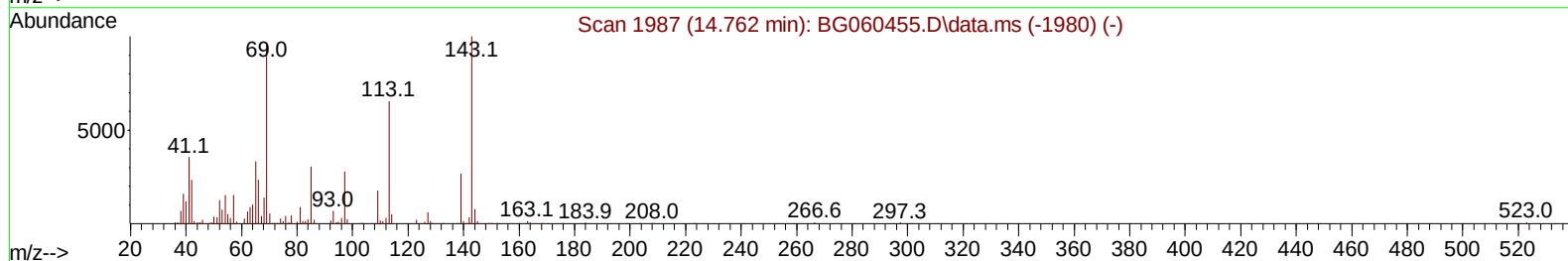
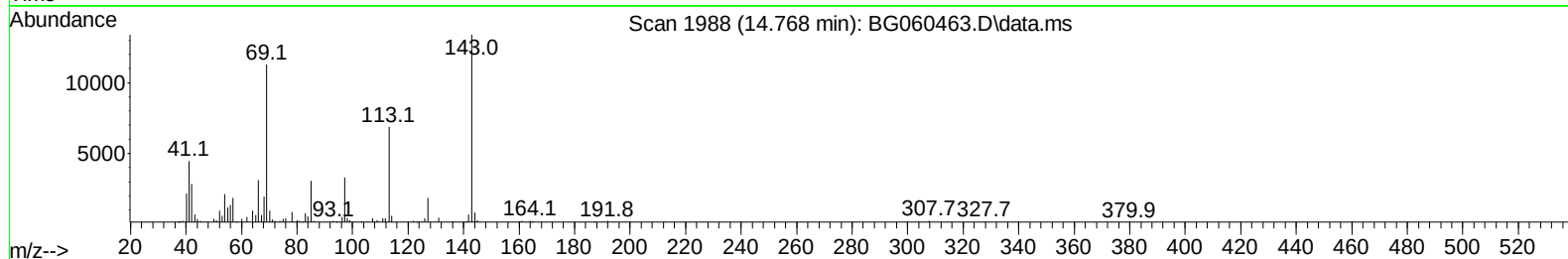
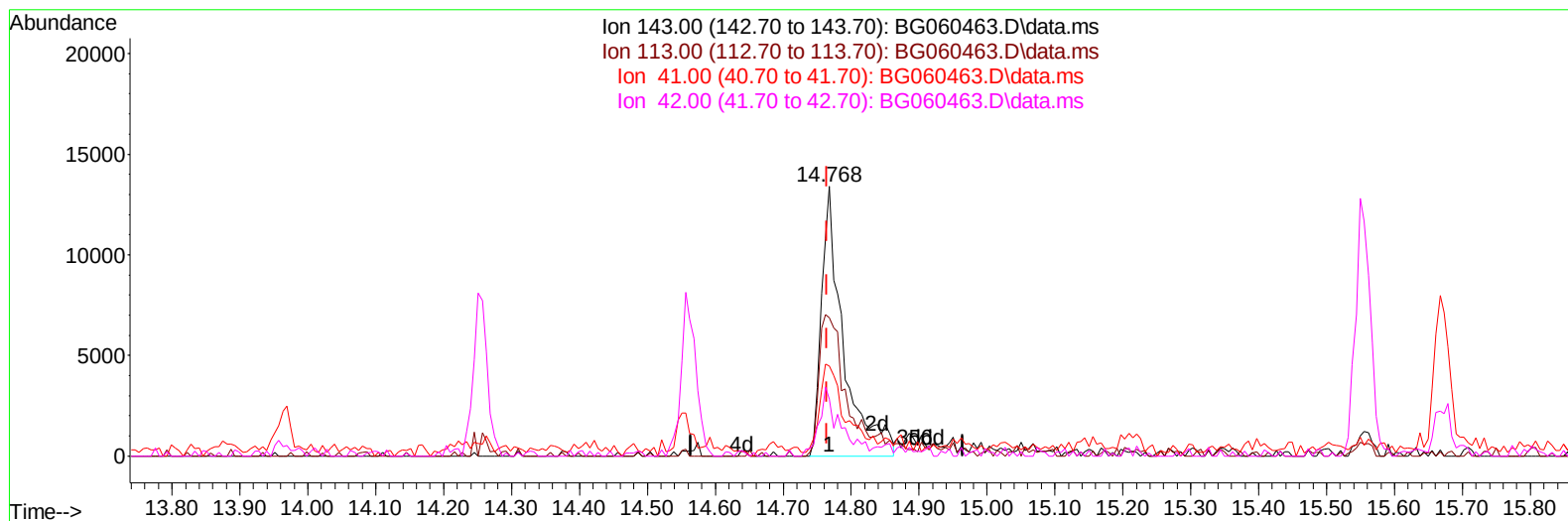
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(54) 4-Nitrophenol-d4 (S)

14.768min (+ 0.003) 4.00 ng/ul m

response 29889

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	51.37
41.00	29.80	33.37
42.00	22.70	21.17

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/30/2024
 Supervised By :mohammad ahmed 01/31/2024

Quant Time: Jan 30 02:44:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 26 06:06:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.923	152	211169	20.000	ng/ul	0.00
20) Naphthalene-d8	10.726	136	842188	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.562	164	636849	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.312	188	1651625	20.000	ng/ul	0.00
79) Chrysene-d12	21.572	240	1577431	20.000	ng/ul	0.00
88) Perylene-d12	24.733	264	1768748	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	14183	3.366	ng/uL	0.00
4) Pyridine-d5	3.775	84	89799	6.736	ng/ul	0.00
7) Phenol-d5	7.089	99	96622	4.818	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.247	67	318865	25.301	ng/ul	0.00
11) 2-Chlorophenol-d4	7.459	132	260469	19.449	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	180201	10.908	ng/ul	0.00
21) Nitrobenzene-d5	9.086	128	167584	26.975	ng/ul	0.00
24) 2-Nitrophenol-d4	9.809	143	191718	26.413	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.350	165	362341	24.004	ng/ul	0.00
31) 4-Chloroaniline-d4	10.861	131	400326	20.216	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	1437241	27.580	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1626594	27.827	ng/ul	0.00
54) 4-Nitrophenol-d4	14.768	143	29889m	4.004	ng/ul	0.00
60) Fluorene-d10	15.555	176	1332489	29.182	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.667	200	228791	22.371	ng/ul	0.00
73) Anthracene-d10	17.412	188	2299005	29.364	ng/ul	0.00
81) Pyrene-d10	19.703	212	2794224	29.555	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.521	264	2730495	29.428	ng/ul	0.00

Target Compounds Qvalue

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

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