

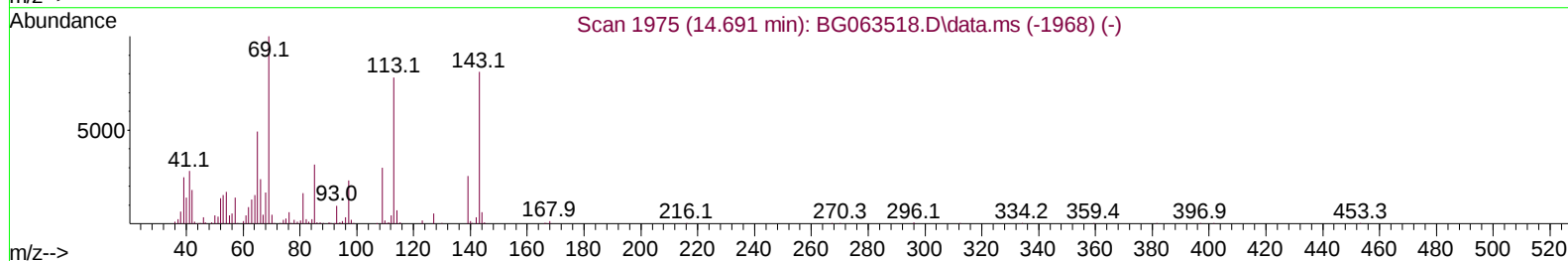
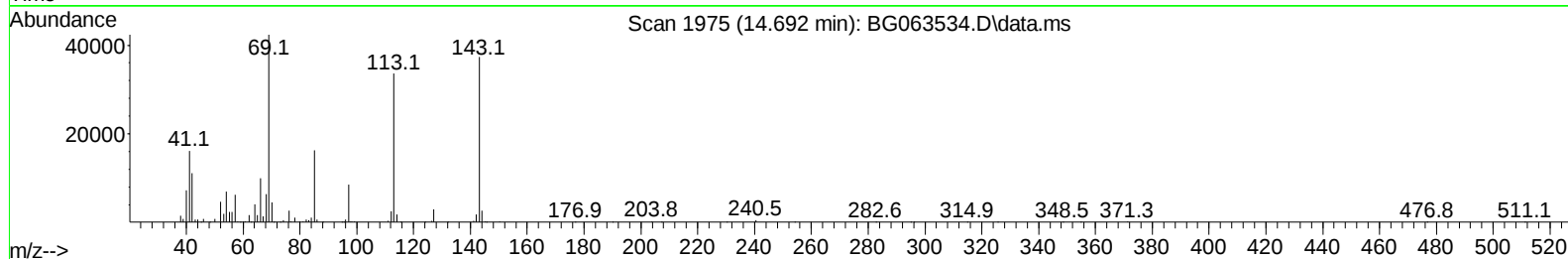
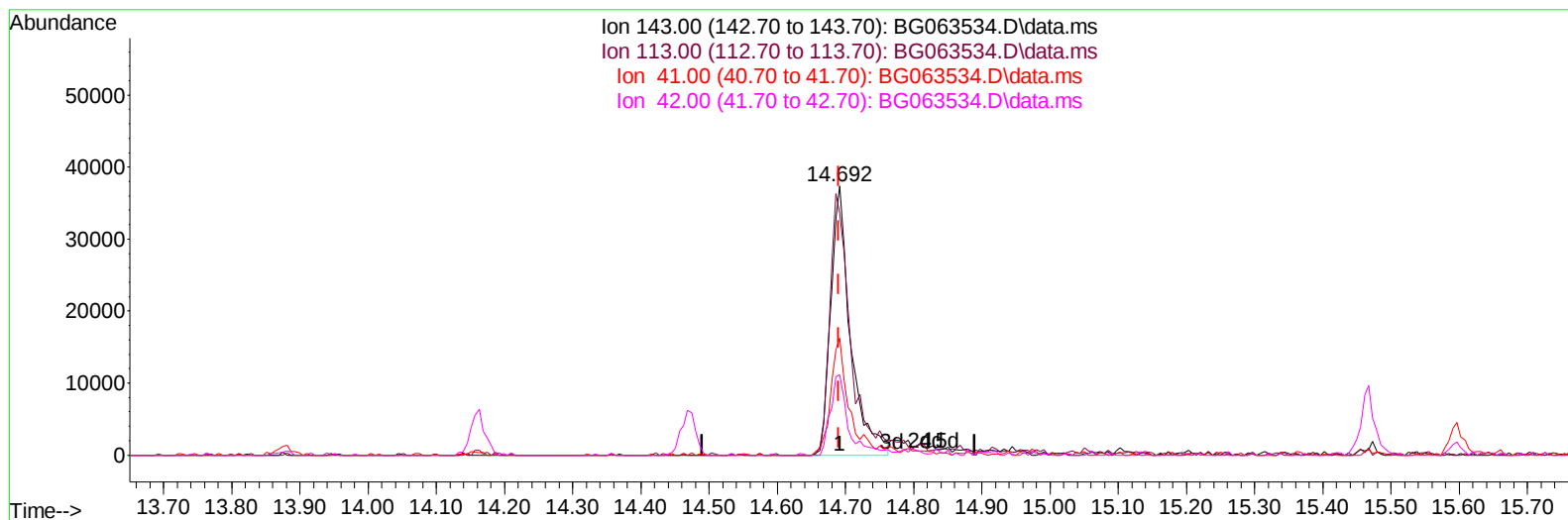
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112224\
 Data File : BG063534.D
 Acq On : 23 Nov 2024 00:42
 Operator : RC/JU
 Sample : PB165140BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK140

Manual IntegrationsAPPROVED

Quant Time: Nov 23 01:31:41 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: BG063534.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.692min (+ 0.002) 22.15 ng/u1

response 75349

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	89.91
41.00	38.80	43.46
42.00	26.30	29.96

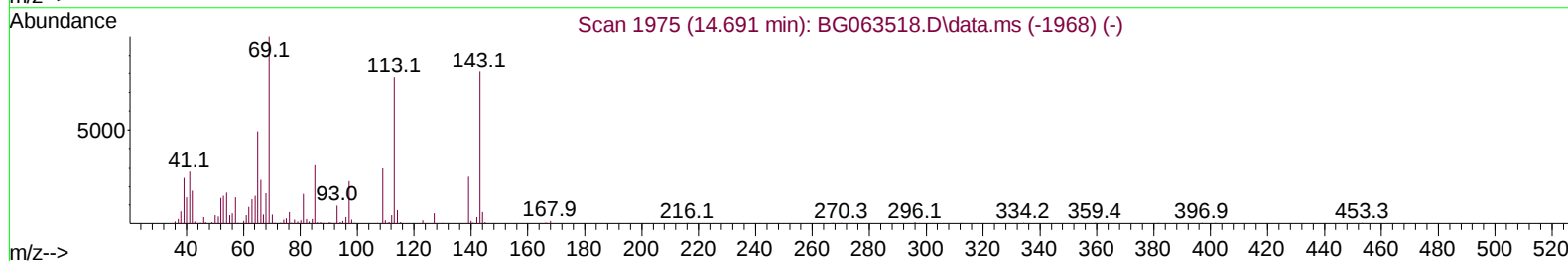
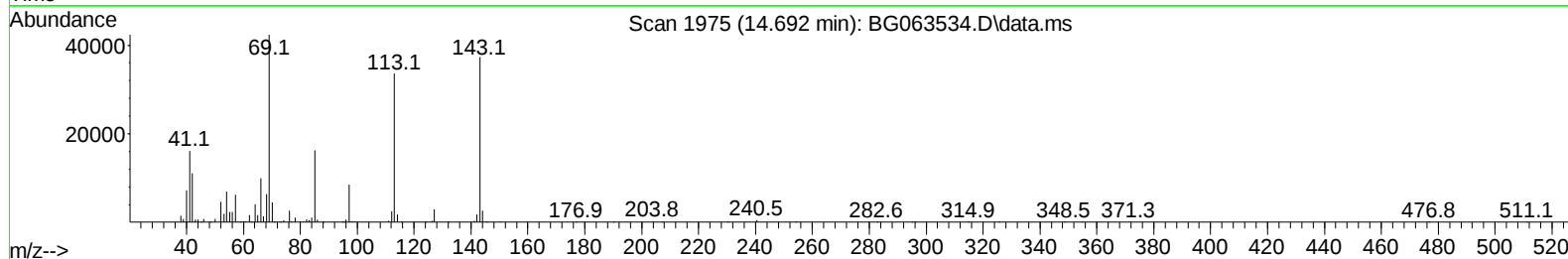
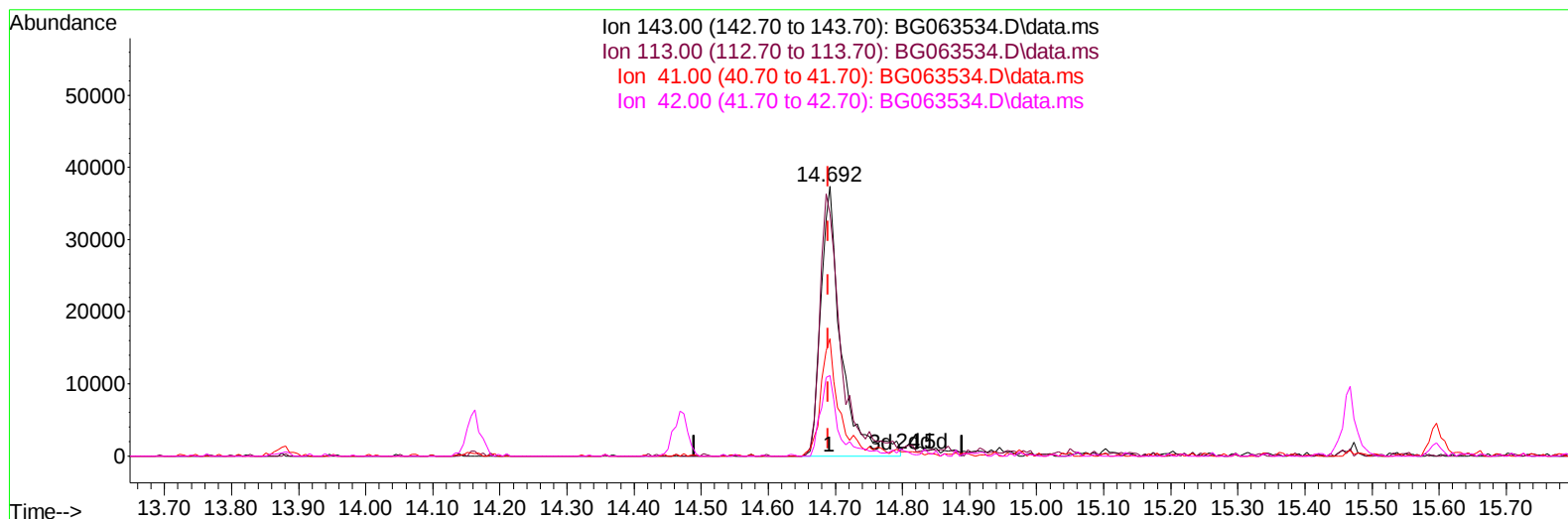
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TIC: BG063534.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.692min (+ 0.002) 23.29 ng/ul m

response 79213

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	89.91
41.00	38.80	43.46
42.00	26.30	29.96

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.829	152	102559	20.000	ng/ul	0.00
20) Naphthalene-d8	10.620	136	435949	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.468	164	337252	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.224	188	873223	20.000	ng/ul	0.00
79) Chrysene-d12	21.484	240	990151	20.000	ng/ul	0.00
88) Perylene-d12	24.527	264	1058576	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	13834	5.324	ng/uL	0.00
4) Pyridine-d5	3.699	84	179392	22.025	ng/ul	0.00
7) Phenol-d5	7.006	99	219039	25.968	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.159	67	145474	26.890	ng/ul	0.00
11) 2-Chlorophenol-d4	7.365	132	158729	26.554	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	180632	25.627	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	80237	25.629	ng/ul	0.00
24) 2-Nitrophenol-d4	9.703	143	97716	25.266	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.244	165	179959	24.378	ng/ul	0.00
31) 4-Chloroaniline-d4	10.755	131	233869	25.048	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	637477	27.844	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	702225	26.897	ng/ul	0.00
54) 4-Nitrophenol-d4	14.692	143	79213m	23.286	ng/ul	0.00
60) Fluorene-d10	15.467	176	582666	28.009	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.596	200	107996	21.998	ng/ul	0.00
73) Anthracene-d10	17.324	188	1028831	27.119	ng/ul	0.00
81) Pyrene-d10	19.627	212	1384079	27.886	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.321	264	1418114	27.619	ng/ul	0.00

Target Compounds Qvalue

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

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