

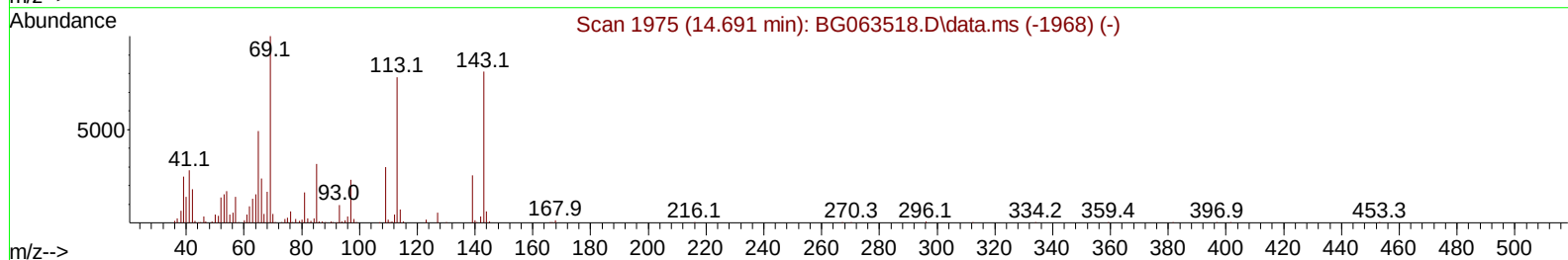
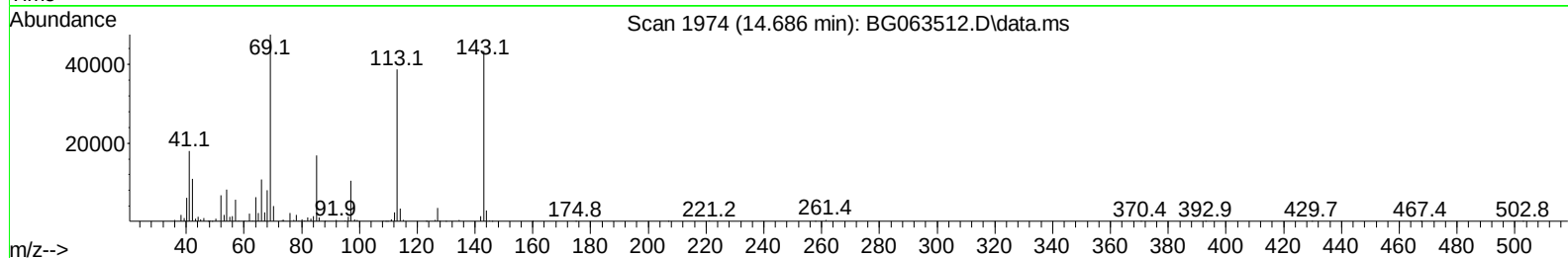
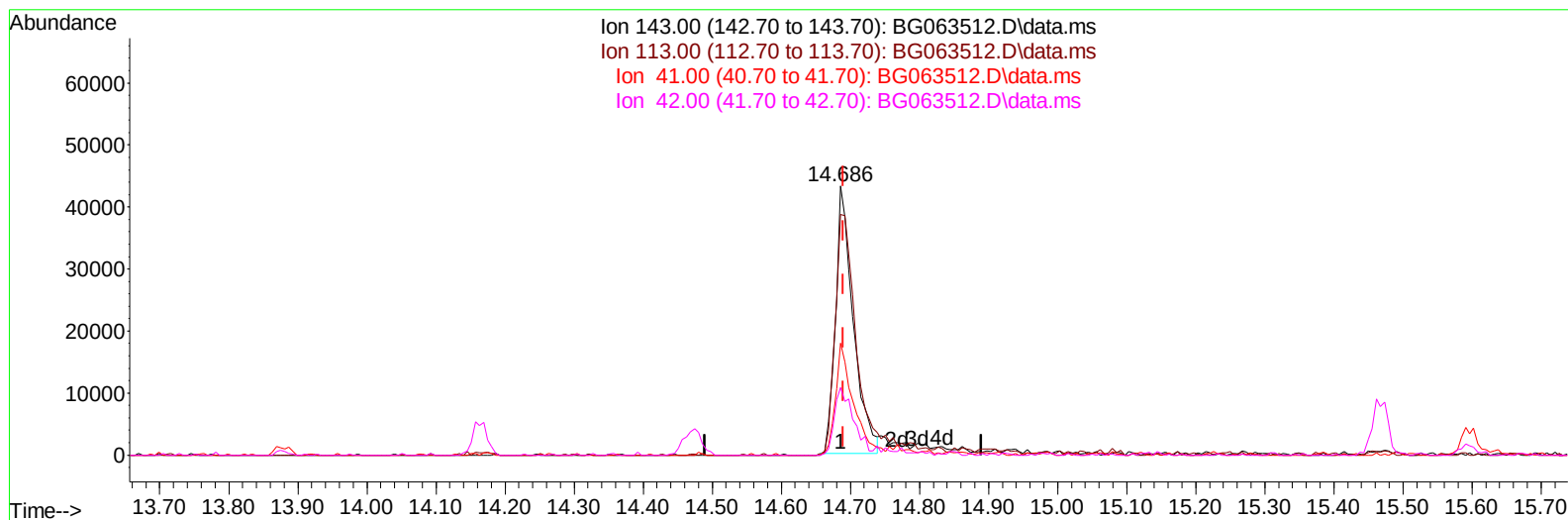
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112224\
Data File : BG063512.D
Acq On : 22 Nov 2024 9:44
Operator : RC/JU
Sample : PB165168BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK168

Manual IntegrationsAPPROVED

Quant Time: Nov 22 12:21:30 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: BG063512.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.686min (-0.004) 27.29 ng/ul

response 77023

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	89.39
41.00	38.80	41.51
42.00	26.30	25.20

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Quant Time: Nov 23 00:16:38 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.829	152	88272	20.000	ng/ul	0.00
20) Naphthalene-d8	10.620	136	354158	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.468	164	279798	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.224	188	717581	20.000	ng/ul	0.00
79) Chrysene-d12	21.484	240	844772	20.000	ng/ul	0.00
88) Perylene-d12	24.533	264	901603	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	14628	6.541	ng/uL	0.00
4) Pyridine-d5	3.699	84	217594	31.040	ng/ul	0.00
7) Phenol-d5	7.007	99	234639	32.320	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.159	67	159588	34.273	ng/ul	0.00
11) 2-Chlorophenol-d4	7.365	132	174386	33.894	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	192149	31.673	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	84333	33.159	ng/ul	0.00
24) 2-Nitrophenol-d4	9.703	143	104270	33.187	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.250	165	189982	31.680	ng/ul	0.00
31) 4-Chloroaniline-d4	10.761	131	241487	31.837	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	684133	36.018	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	751821	34.710	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	84133m	29.810	ng/ul	0.00
60) Fluorene-d10	15.467	176	627432	36.354	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.597	200	114557	28.396	ng/ul	0.00
73) Anthracene-d10	17.324	188	1129947	36.245	ng/ul	0.00
81) Pyrene-d10	19.627	212	1542257	36.421	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.322	264	1559911	35.671	ng/ul	0.00

Target Compounds Qvalue

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

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