

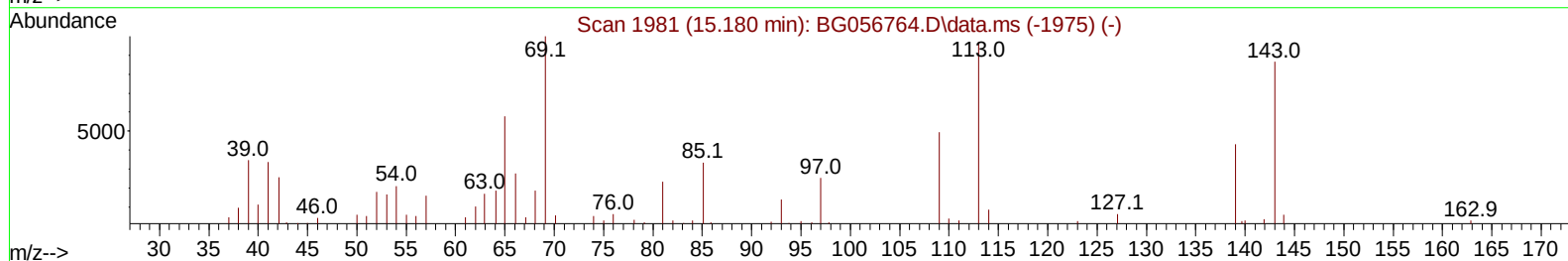
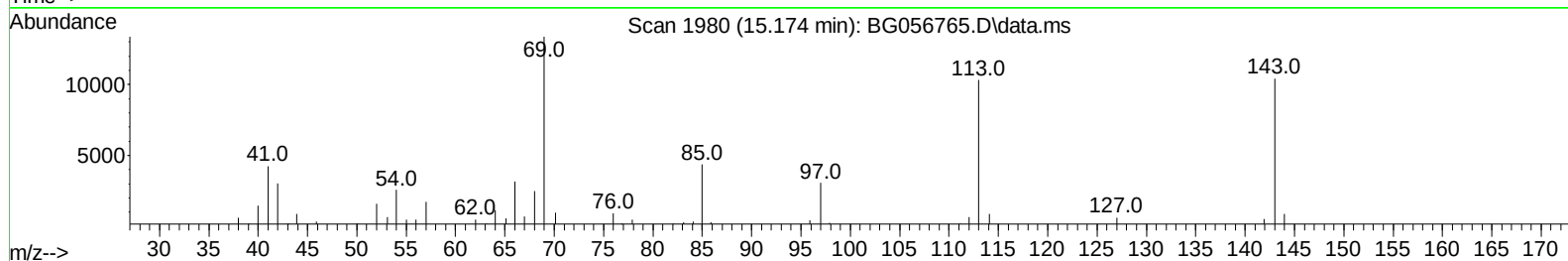
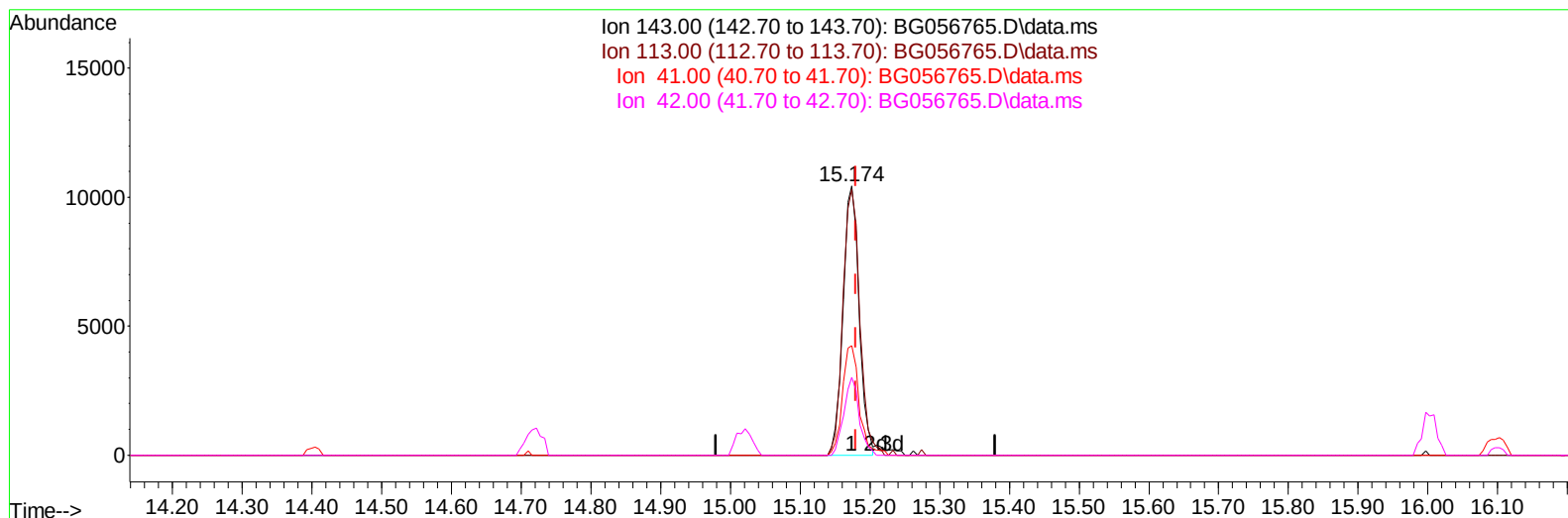
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030123\
 Data File : BG056765.D
 Acq On : 1 Mar 2023 15:46
 Operator : CG/JU
 Sample : PB151097BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK097

Manual IntegrationsAPPROVED

Quant Time: Mar 02 00:14:38 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 02 00:12:28 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/02/2023
 Supervised By :Jagrut Upadhyay 03/02/2023



TIC: BG056765.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.174min (-0.006) 22.44 ng/ul

response 16589

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.30	98.97
41.00	46.10	40.80
42.00	27.10	28.94

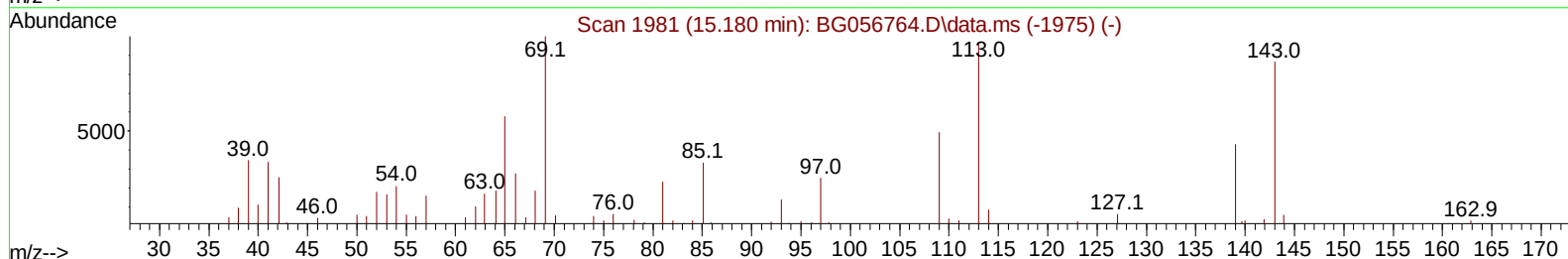
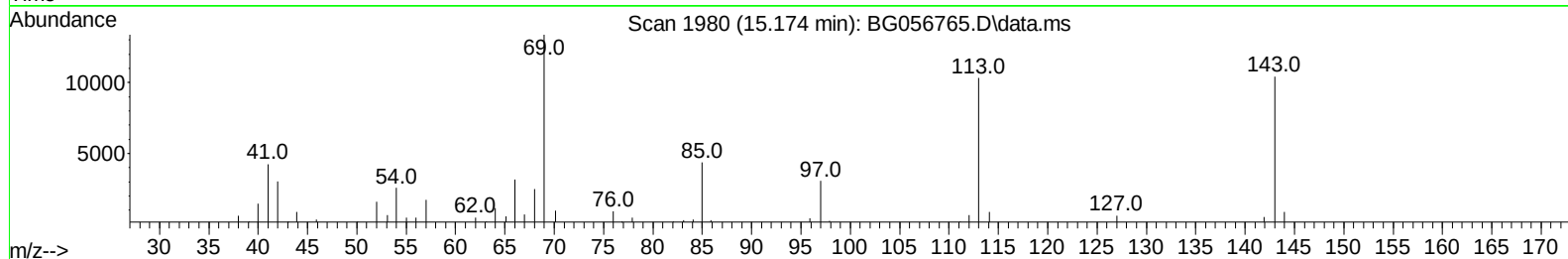
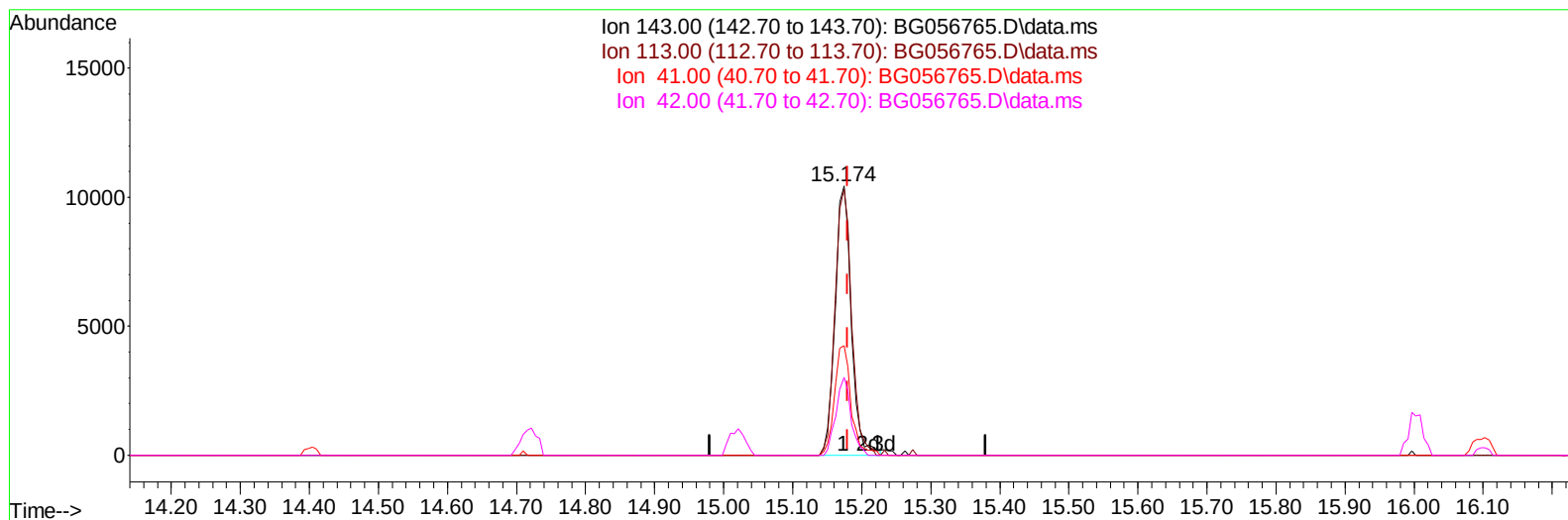
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TIC: BG056765.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.174min (-0.006) 22.95 ng/ul m

response 16971

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.30	98.97
41.00	46.10	40.80
42.00	27.10	28.94

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.417	152	15028	20.000	ng/ul	0.00
20) Naphthalene-d8	11.255	136	67459	20.000	ng/ul	0.00
38) Acenaphthene-d10	15.021	164	53150	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.753	188	136597	20.000	ng/ul	0.00
79) Chrysene-d12	22.060	240	130507	20.000	ng/ul	0.00
88) Perylene-d12	25.585	264	138315	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.699	96	1851	6.832	ng/uL	0.00
4) Pyridine-d5	4.134	84	31509	28.815	ng/ul	0.00
7) Phenol-d5	7.542	99	35738	23.764	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.724	67	23650	25.782	ng/ul	0.00
11) 2-Chlorophenol-d4	7.941	132	25377	25.411	ng/ul	0.00
15) 4-Methylphenol-d8	9.105	113	28540	24.837	ng/ul	0.00
21) Nitrobenzene-d5	9.592	128	14060	24.870	ng/ul	0.00
24) 2-Nitrophenol-d4	10.315	143	15825	23.516	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.862	165	29329	23.134	ng/ul	0.00
31) 4-Chloroaniline-d4	11.379	131	40040	24.215	ng/ul	0.00
46) Dimethylphthalate-d6	14.404	166	112735	26.289	ng/ul	0.00
49) Acenaphthylene-d8	14.722	160	125911	25.834	ng/ul	0.00
54) 4-Nitrophenol-d4	15.174	143	16971m	22.954	ng/ul	0.00
60) Fluorene-d10	16.003	176	101636	25.622	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.102	200	19184	20.992	ng/ul	0.00
73) Anthracene-d10	17.853	188	170309	25.700	ng/ul	0.00
81) Pyrene-d10	20.110	212	207030	23.443	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.339	264	198111	26.536	ng/ul	-0.01

Target Compounds Qvalue

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

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