

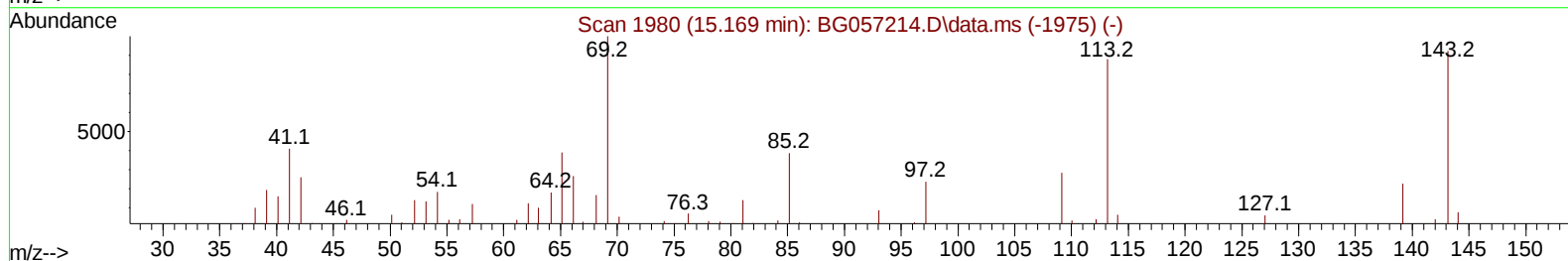
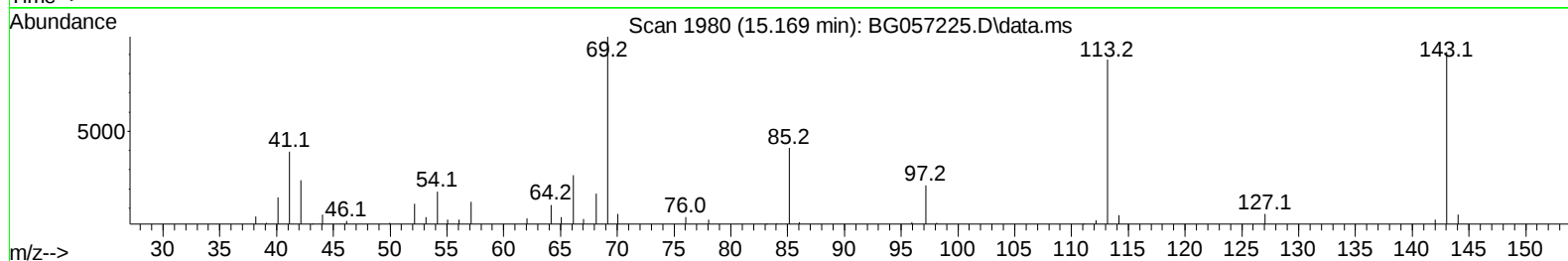
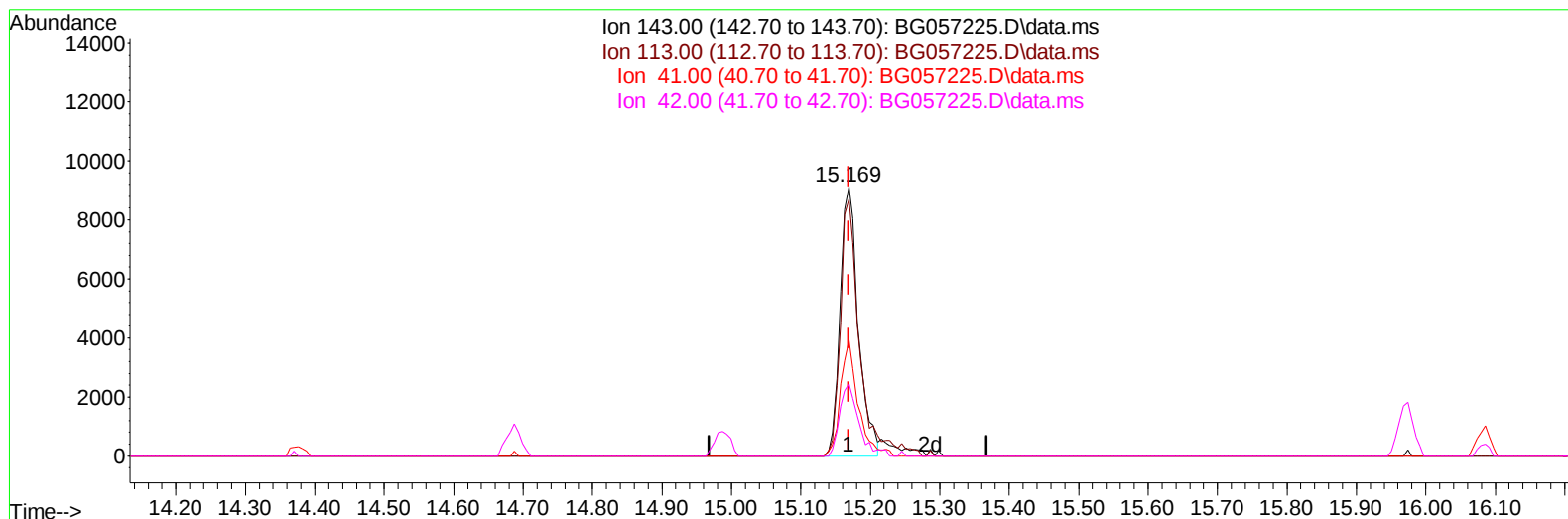
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042823\
Data File : BG057225.D
Acq On : 28 Apr 2023 20:10
Operator : CG/JU
Sample : PB152400BL
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK400

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/01/2023
Supervised By :Jagrut Upadhyay 05/01/2023

Quant Time: Apr 29 01:19:16 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 28 23:37:02 2023
Response via : Initial Calibration



TIC: BG057225.D\data.ms

(54) 4-Nitrophenol-d4 (S)**15.169min (+ 0.001) 30.38 ng/u1****response 16561**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.60	95.42
41.00	44.40	43.02
42.00	28.40	26.83

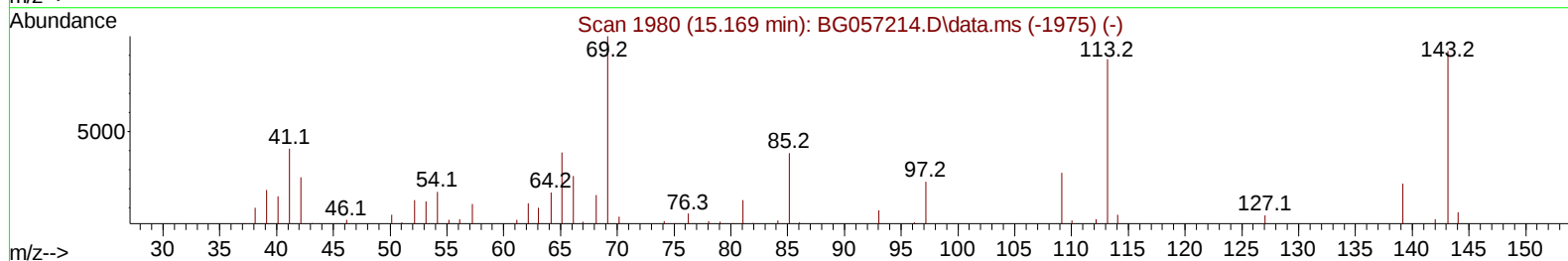
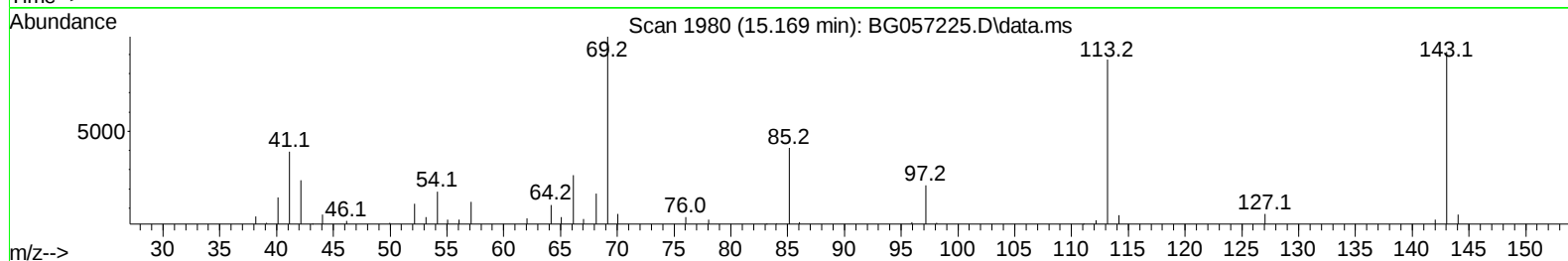
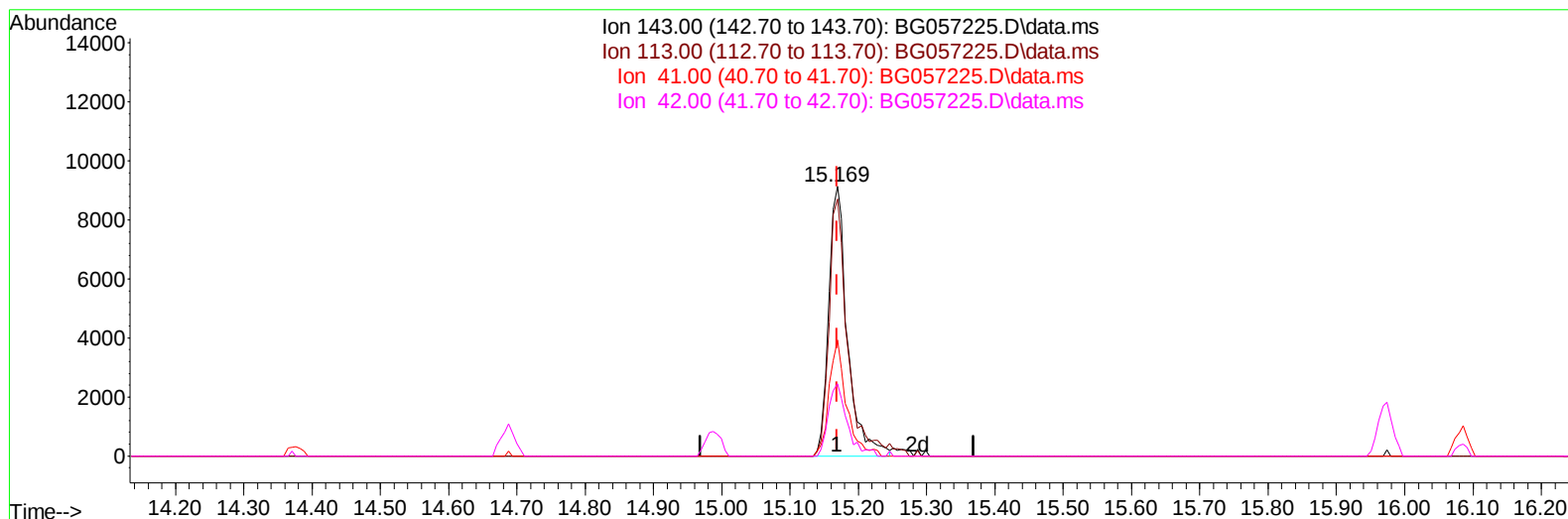
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TIC: BG057225.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.169min (+ 0.001) 31.79 ng/ul m

response 17328

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.60	95.42
41.00	44.40	43.02
42.00	28.40	26.83

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 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/01/2023
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Quant Time: Apr 29 01:19:45 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 28 23:37:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.378	152	13697	20.000	ng/ul	0.00
20) Naphthalene-d8	11.216	136	57621	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.987	164	47256	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.724	188	128922	20.000	ng/ul	0.00
79) Chrysene-d12	22.036	240	119039	20.000	ng/ul	0.00
88) Perylene-d12	25.543	264	121566	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.655	96	1763	5.664	ng/uL	0.00
4) Pyridine-d5	4.096	84	25894	26.388	ng/ul	0.00
7) Phenol-d5	7.509	99	37393	28.867	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.679	67	24185	31.744	ng/ul	0.00
11) 2-Chlorophenol-d4	7.902	132	26914	31.529	ng/ul	0.00
15) 4-Methylphenol-d8	9.071	113	29639	30.105	ng/ul	0.00
21) Nitrobenzene-d5	9.553	128	15034	32.763	ng/ul	0.00
24) 2-Nitrophenol-d4	10.282	143	17595	32.856	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.828	165	31148	30.148	ng/ul	0.00
31) 4-Chloroaniline-d4	11.339	131	39446	28.806	ng/ul	0.00
46) Dimethylphthalate-d6	14.376	166	131747	35.185	ng/ul	0.00
49) Acenaphthylene-d8	14.687	160	136598	32.458	ng/ul	0.00
54) 4-Nitrophenol-d4	15.169	143	17328m	31.792	ng/ul	0.00
60) Fluorene-d10	15.974	176	110376	31.643	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.086	200	22486	29.438	ng/ul	0.00
73) Anthracene-d10	17.824	188	199760	33.946	ng/ul	0.00
81) Pyrene-d10	20.092	212	249363	35.292	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.302	264	219032	35.365	ng/ul	0.00

Target Compounds Qvalue

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

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