

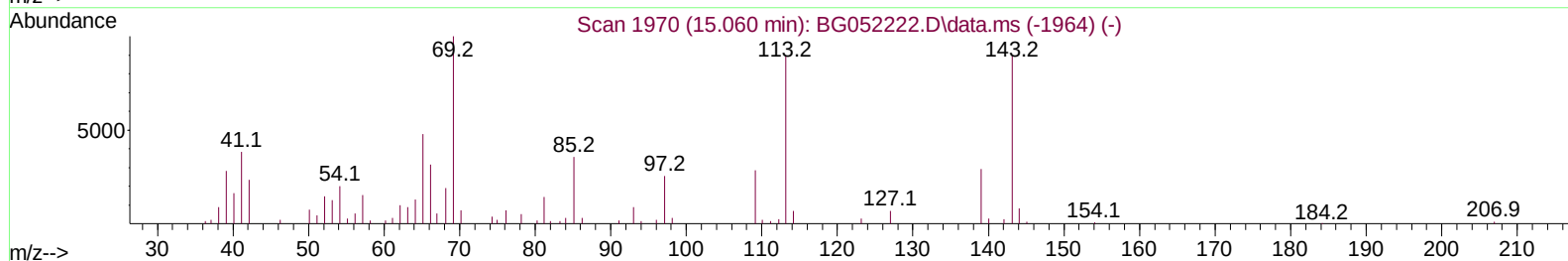
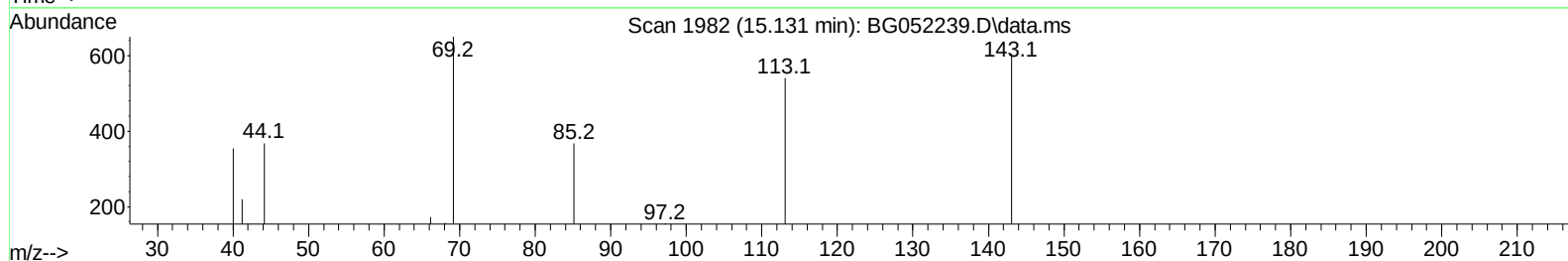
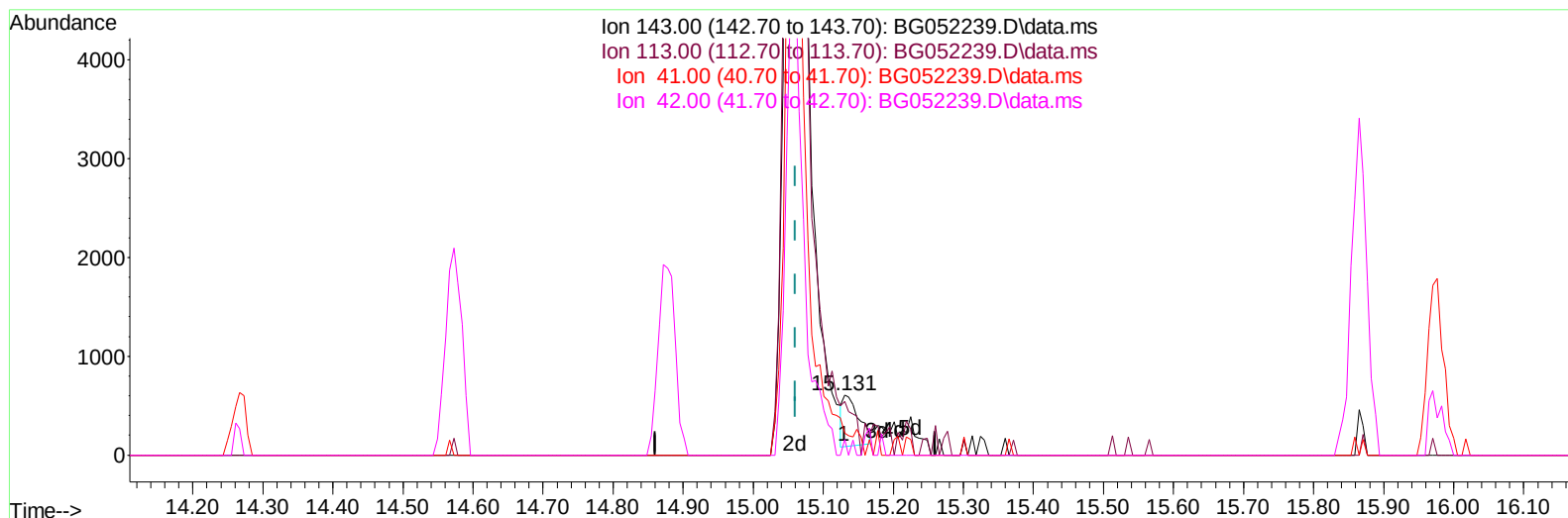
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020122\
Data File : BG052239.D
Acq On : 1 Feb 2022 21:35
Operator : CG/JU
Sample : PB142387BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK387

Manual IntegrationsAPPROVED

Quant Time: Feb 02 03:10:21 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG013122.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jan 31 18:07:10 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/03/2022
Supervised By :mohammad ahmed 02/04/2022



TIC: BG052239.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.131min (+ 0.070) 0.68 ng/ul

response 810

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	89.26
41.00	44.40	36.36
42.00	29.70	25.79

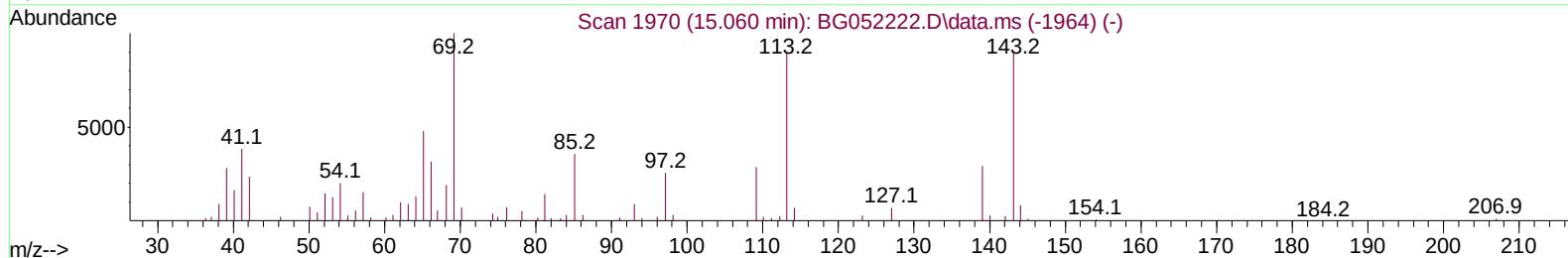
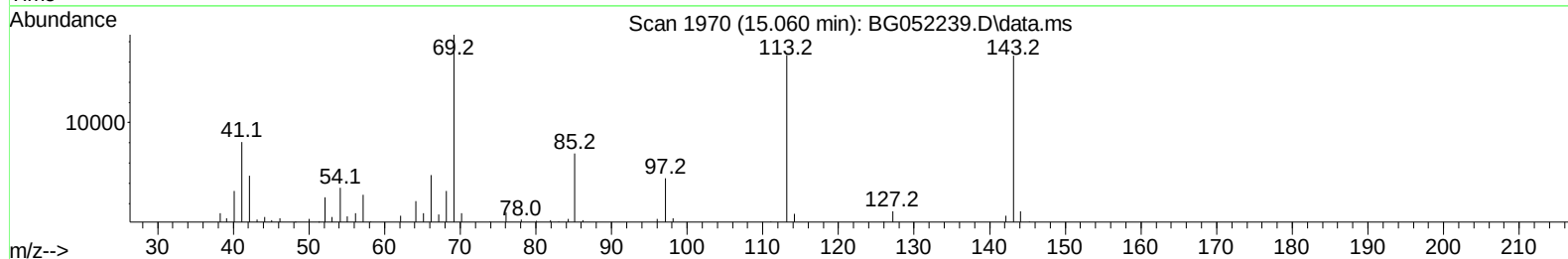
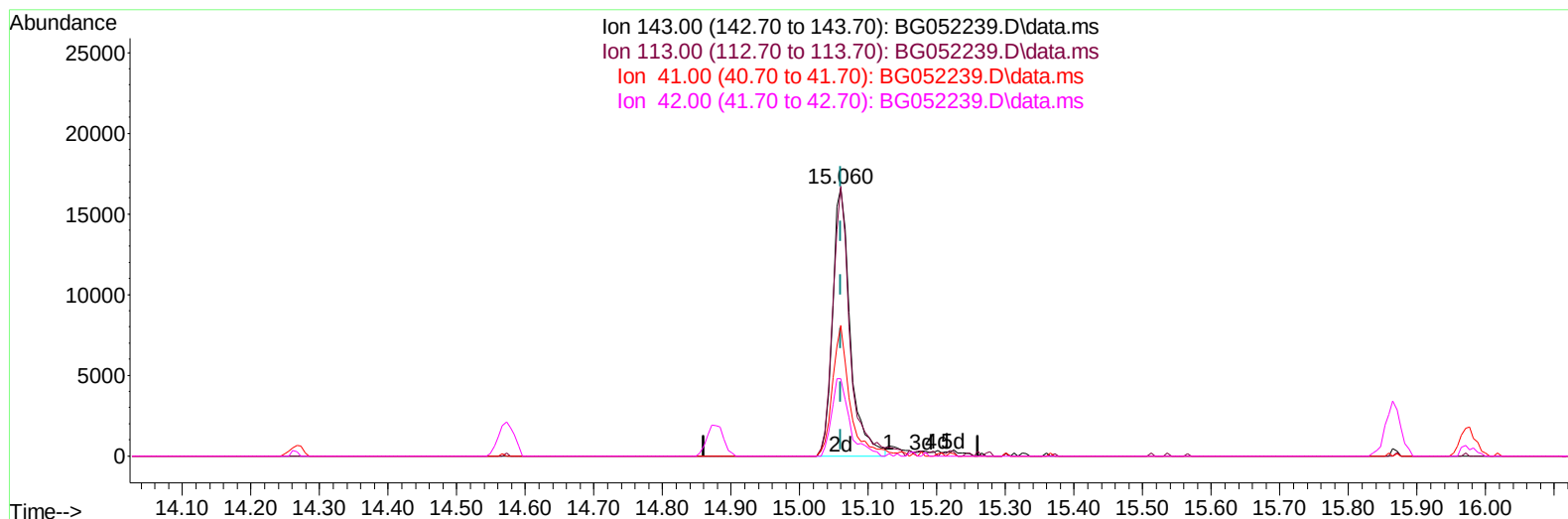
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TIC: BG052239.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.060min (-0.000) 24.77 ng/ul m

response 29715

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	100.53#
41.00	44.40	48.72
42.00	29.70	28.74

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.234	152	33243	20.000	ng/ul	0.00
20) Naphthalene-d8	11.071	136	141375	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.878	164	100788	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.627	188	245964	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.939	240	244317	20.000	ng/ul	# 0.00
88) Perylene-d12	25.411	264	241238	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.558	96	4206	5.104	ng/uL	0.00
4) Pyridine-d5	3.987	84	66811	26.068	ng/ul	0.00
7) Phenol-d5	7.376	99	81735	27.031	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.535	67	52166	28.397	ng/ul	0.00
11) 2-Chlorophenol-d4	7.758	132	60937	28.920	ng/ul	0.00
15) 4-Methylphenol-d8	8.933	113	63734	27.775	ng/ul	0.00
21) Nitrobenzene-d5	9.403	128	32589	28.311	ng/ul	0.00
24) 2-Nitrophenol-d4	10.131	143	35975	28.600	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.684	165	65850	26.853	ng/ul	0.00
31) 4-Chloroaniline-d4	11.195	131	84203	27.070	ng/ul	0.00
46) Dimethylphthalate-d6	14.267	166	235800	29.764	ng/ul	0.00
49) Acenaphthylene-d8	14.572	160	283996	28.760	ng/ul	0.00
54) 4-Nitrophenol-d4	15.060	143	29715m	24.766	ng/ul	0.00
60) Fluorene-d10	15.865	176	202661	29.199	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.976	200	34249	22.554	ng/ul	0.00
73) Anthracene-d10	17.721	188	343308	29.606	ng/ul	0.00
81) Pyrene-d10	20.000	212	414152	28.284	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.176	264	385579	30.086	ng/ul	0.00

Target Compounds Qvalue

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

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