

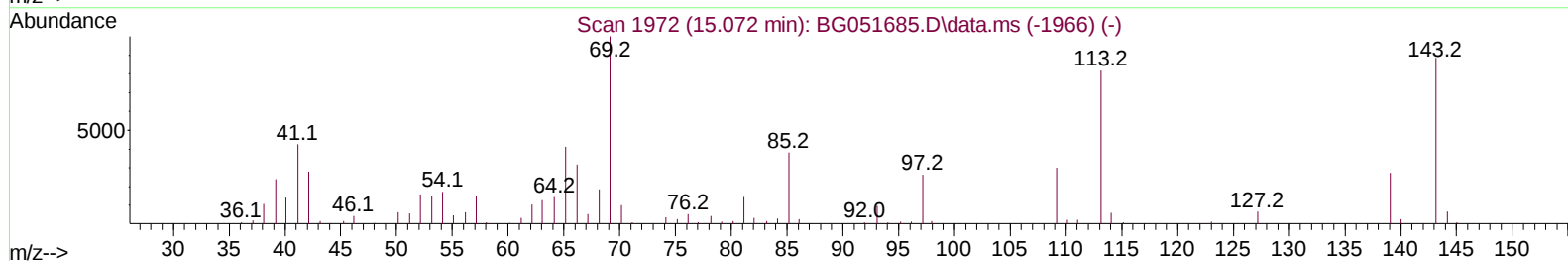
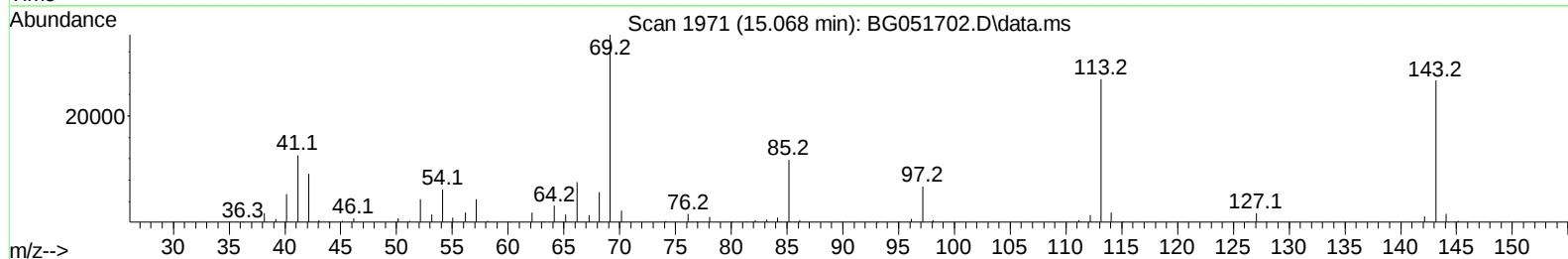
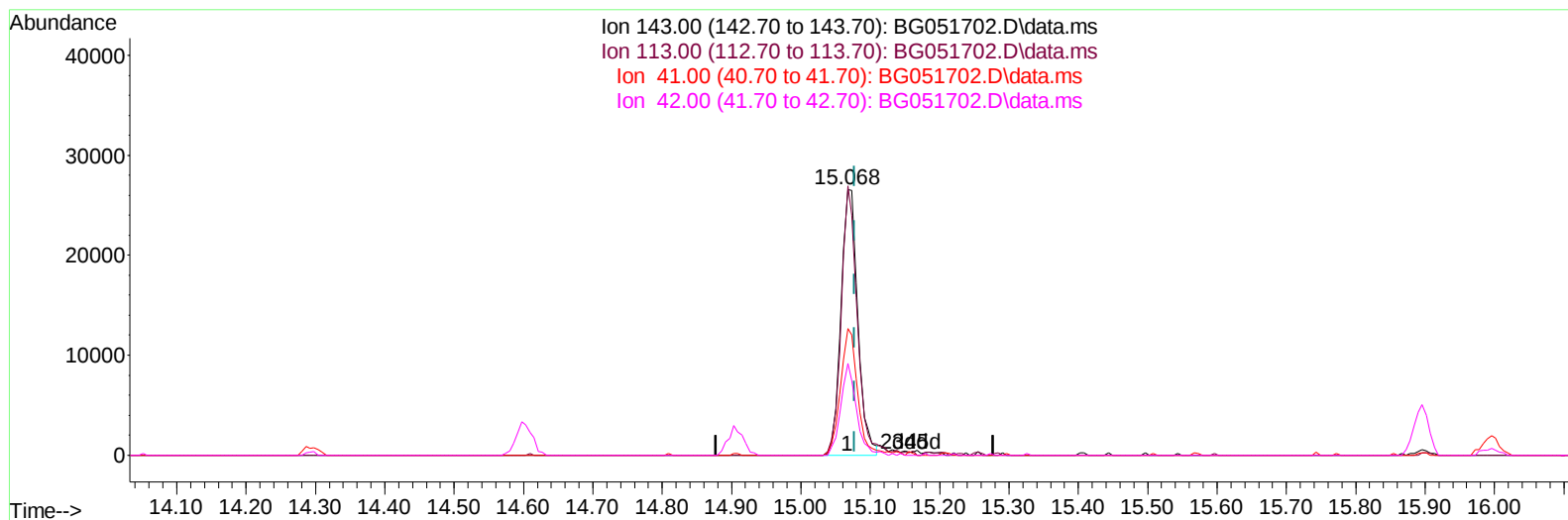
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
 Data File : BG051702.D
 Acq On : 21 Dec 2021 5:29
 Operator : CG/JU
 Sample : PB141554BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK554

Manual IntegrationsAPPROVED

Quant Time: Dec 21 06:02:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 14 14:19:57 2021
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/21/2021
 Supervised By :mohammad ahmed 12/28/2021



TIC: BG051702.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.068min (-0.010) 28.35 ng/u1

response 44577

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	101.26#
41.00	44.40	47.69
42.00	29.70	34.49

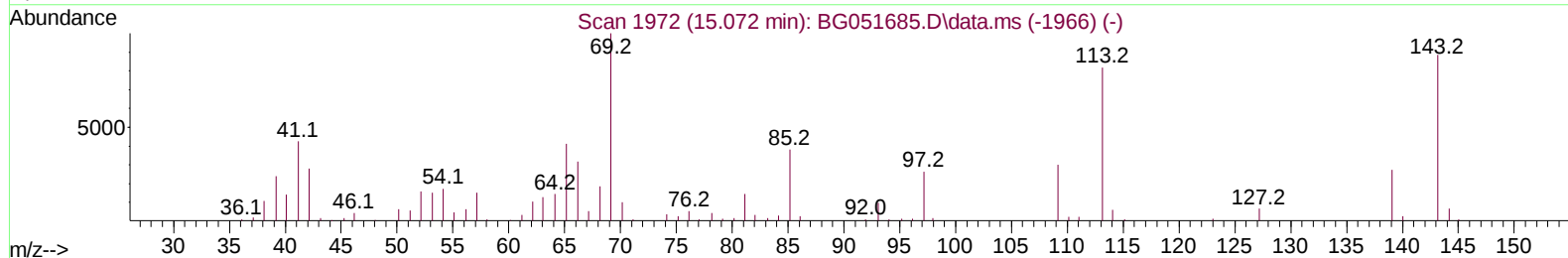
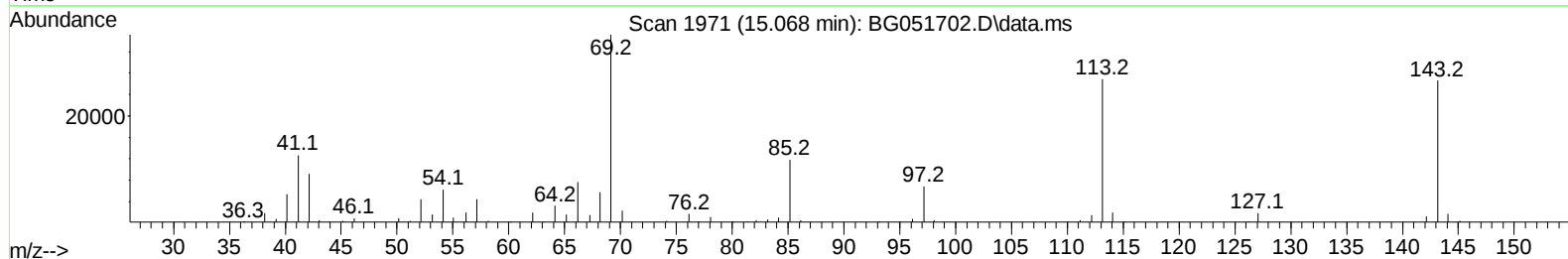
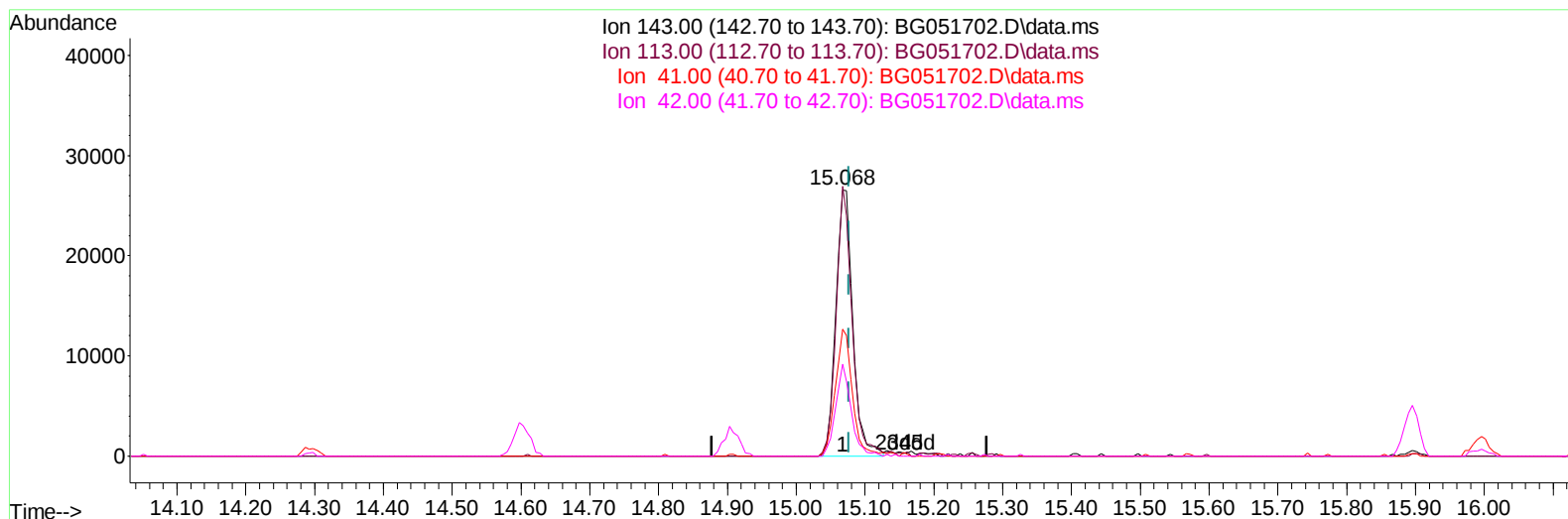
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TIC: BG051702.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.068min (-0.010) 28.77 ng/ul m

response 45243

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.30	101.26#
41.00	44.40	47.69
42.00	29.70	34.49

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.271	152	38596	20.000	ng/ul	-0.02
20) Naphthalene-d8	11.108	136	163910	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.909	164	120802	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.652	188	282879	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.976	240	273518	20.000	ng/ul	# 0.00
88) Perylene-d12	25.471	264	265382	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.595	96	5900	6.349	ng/uL	-0.01
4) Pyridine-d5	4.024	84	94813	33.766	ng/ul	-0.02
7) Phenol-d5	7.401	99	119900	34.793	ng/ul	-0.02
9) Bis-(2-Chloroethyl)eth...	7.578	67	69170	33.764	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.795	132	85150	35.527	ng/ul	-0.02
15) 4-Methylphenol-d8	8.964	113	93610	35.142	ng/ul	-0.02
21) Nitrobenzene-d5	9.440	128	43595	36.277	ng/ul	-0.02
24) 2-Nitrophenol-d4	10.174	143	49387	37.594	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.715	165	96650	33.926	ng/ul	-0.02
31) 4-Chloroaniline-d4	11.232	131	126296	33.672	ng/ul	0.00
46) Dimethylphthalate-d6	14.298	166	318417	32.932	ng/ul	-0.02
49) Acenaphthylene-d8	14.603	160	398200	33.160	ng/ul	0.00
54) 4-Nitrophenol-d4	15.068	143	45243m	28.774	ng/ul	0.00
60) Fluorene-d10	15.896	176	280087	32.332	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.996	200	35198	23.882	ng/ul	-0.02
73) Anthracene-d10	17.752	188	459272	34.035	ng/ul	-0.01
81) Pyrene-d10	20.031	212	543367	33.008	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.230	264	497293	35.614	ng/ul	-0.02

Target Compounds Qvalue

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

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