

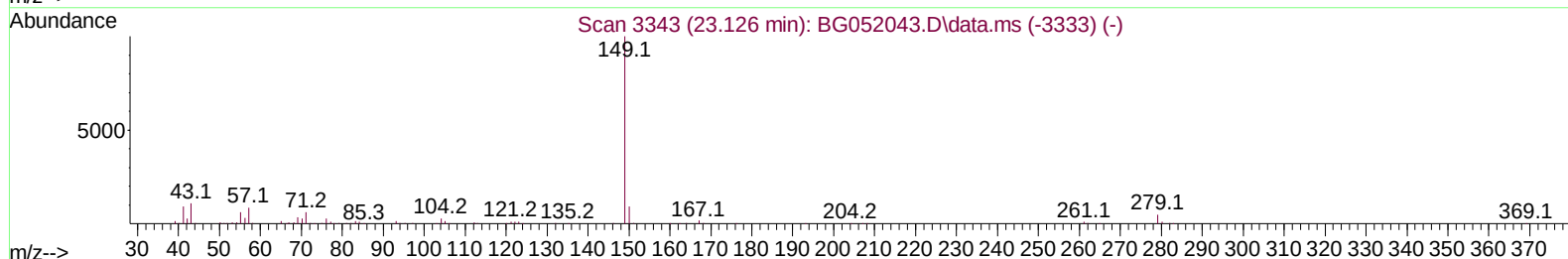
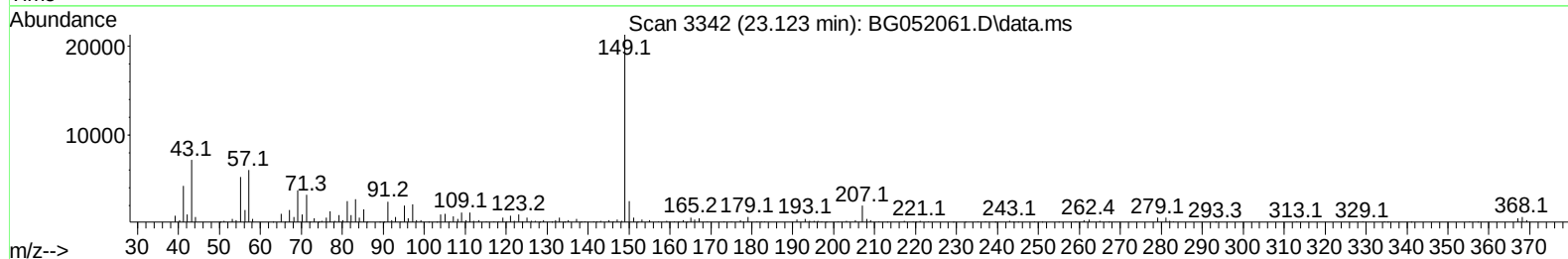
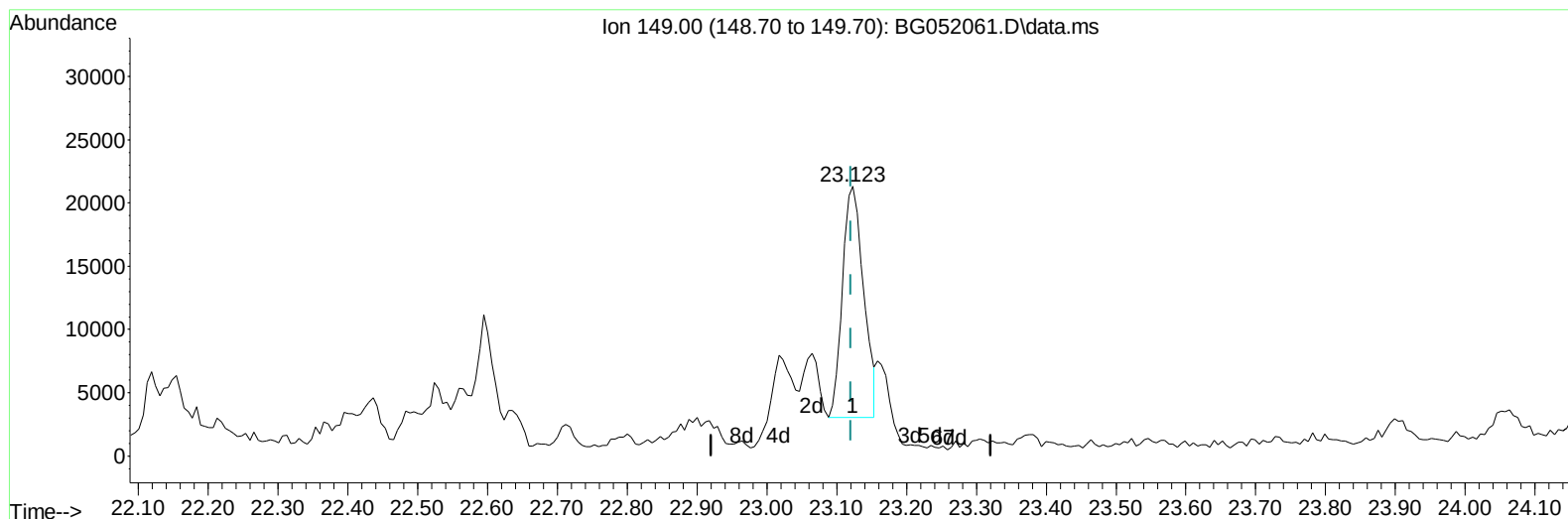
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011722\
 Data File : BG052061.D
 Acq On : 18 Jan 2022 20:35
 Operator : CG/JU
 Sample : N1100-14ME
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLS6ME

Manual IntegrationsAPPROVED

Quant Time: Jan 19 00:32:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/19/2022
 Supervised By :mohammad ahmed 01/21/2022



TIC: BG052061.D\data.ms

(89) Di-n-octyl phthalate

23.123min (+ 0.002) 4.27 ng/ul

response 38258

Ion	Exp%	Act%
149.00	100.00	100.00
0.00	0.00	0.00
0.00	0.00	0.00
0.00	0.00	0.00

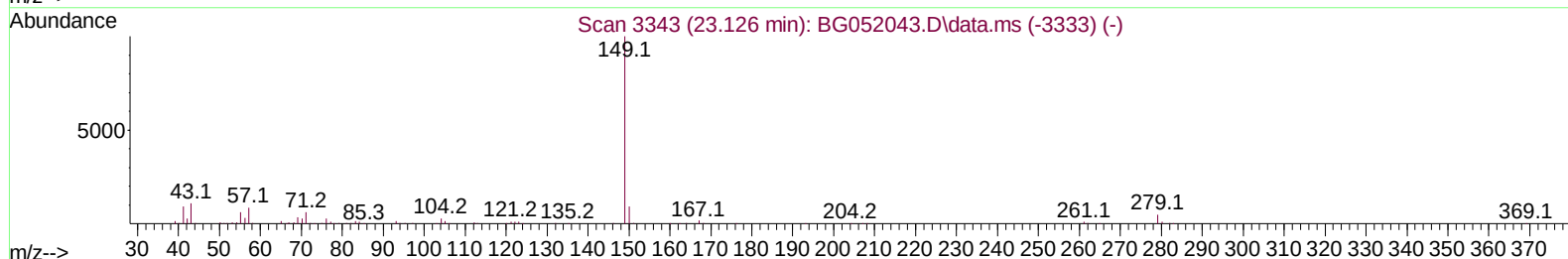
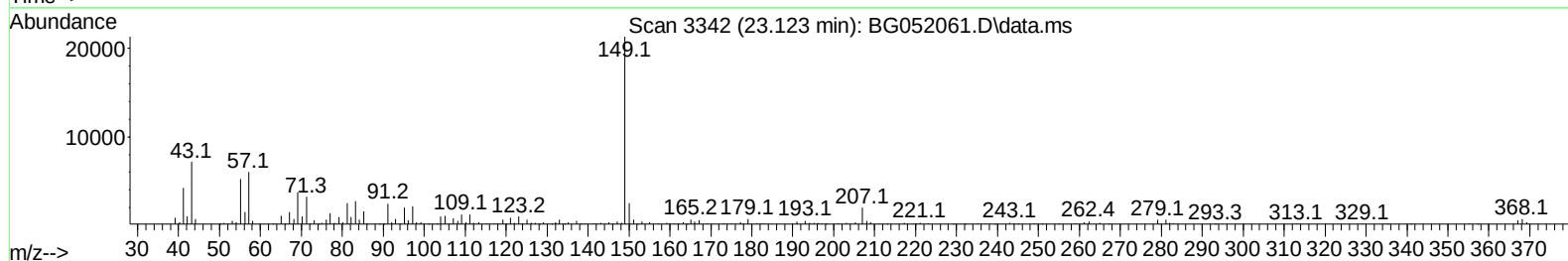
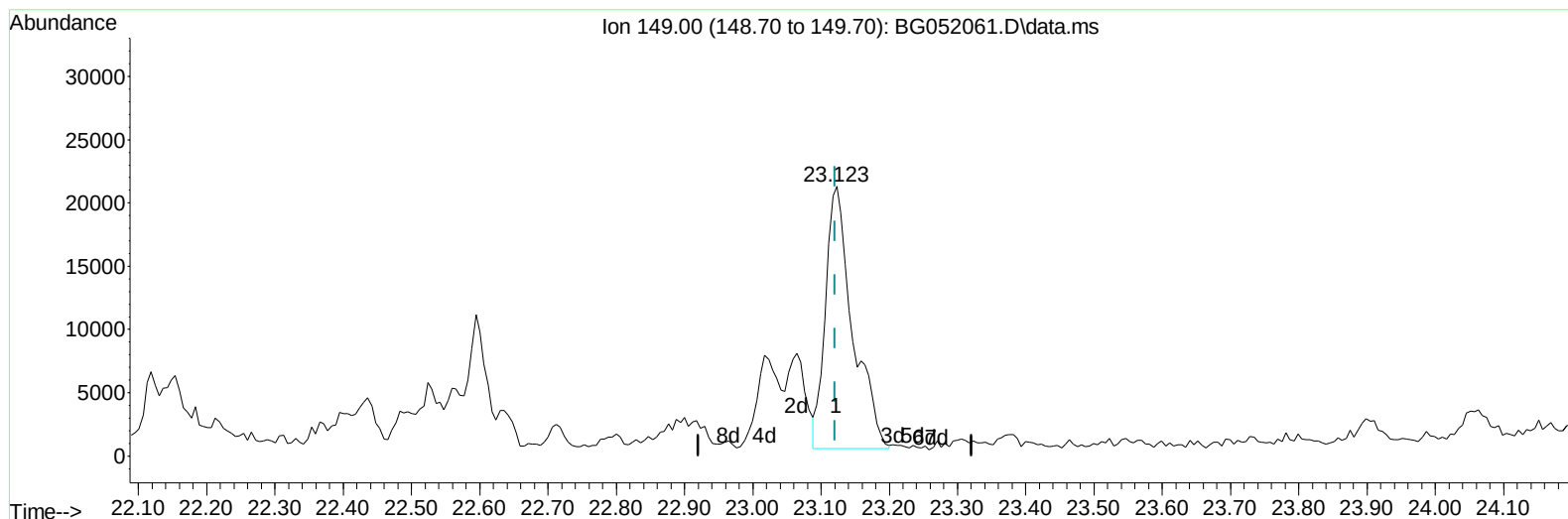
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(89) Di-n-octyl phthalate

23.123min (+ 0.002) 6.39 ng/ul m

response 57278

Ion	Exp%	Act%
149.00	100.00	100.00
0.00	0.00	0.00
0.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.243	152	24717	20.000	ng/ul	0.00
20) Naphthalene-d8	11.081	136	108465	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.887	164	77038	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.642	188	161869	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.960	240	145642	20.000	ng/ul	# 0.00
88) Perylene-d12	25.444	264	162239	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.567	96	1718	2.484	ng/uL	0.00
4) Pyridine-d5	4.002	84	11161	5.191	ng/ul	0.00
7) Phenol-d5	7.380	99	41054	16.971	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.544	67	26528	17.274	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.768	132	30333	18.938	ng/ul	0.00
15) 4-Methylphenol-d8	8.942	113	32751	17.416	ng/ul	0.00
21) Nitrobenzene-d5	9.412	128	16119	18.807	ng/ul	0.00
24) 2-Nitrophenol-d4	10.147	143	19082	20.131	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.699	165	35072	18.929	ng/ul	0.00
31) 4-Chloroaniline-d4	11.204	131	45895	17.787	ng/ul	0.00
46) Dimethylphthalate-d6	14.276	166	115601	18.154	ng/ul	0.00
49) Acenaphthylene-d8	14.582	160	146180	19.004	ng/ul	0.00
54) 4-Nitrophenol-d4	15.064	143	17147	16.106	ng/ul	0.00
60) Fluorene-d10	15.880	176	103846	18.704	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.986	200	13472	13.193	ng/ul	0.00
73) Anthracene-d10	17.736	188	165117	21.503	ng/ul	0.00
81) Pyrene-d10	20.016	212	184988	22.070	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.203	264	176422	20.656	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	11.134	128	7362	1.253	ng/ul#	90
36) 2-Methylnaphthalene	12.726	142	4628	1.105	ng/ul#	80
37) 1-Methylnaphthalene	12.949	142	9462	2.202	ng/ul#	96
43) 1,1'-Biphenyl	13.724	154	7877	1.324	ng/ul	97
52) Acenaphthene	14.952	153	14377	2.855	ng/ul	95
56) Dibenzofuran	15.287	168	15323	2.095	ng/ul	96
61) Fluorene	15.933	166	9972	1.636	ng/ul#	90
72) Phenanthrene	17.683	178	20486	2.415	ng/ul#	95
86) Bis(2-ethylhexyl)phtha...	21.831	149	1543304	317.955	ng/ul#	92
89) Di-n-octyl phthalate	23.123	149	57278m	6.391	ng/ul	

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

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