

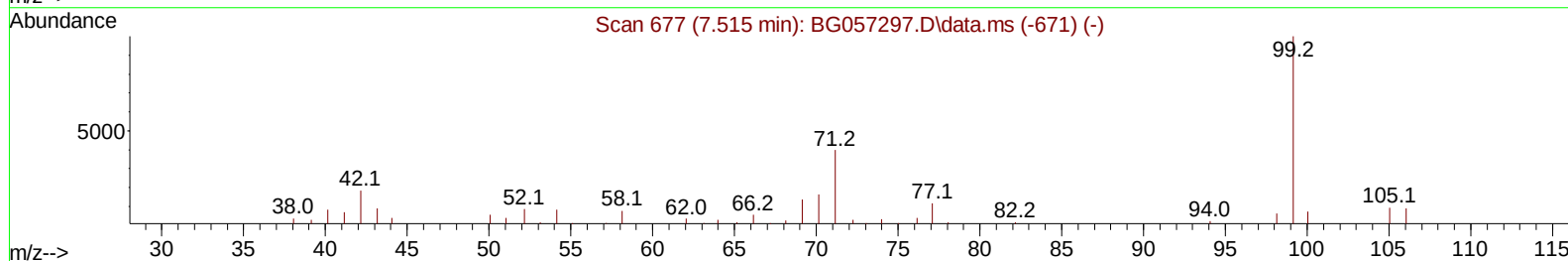
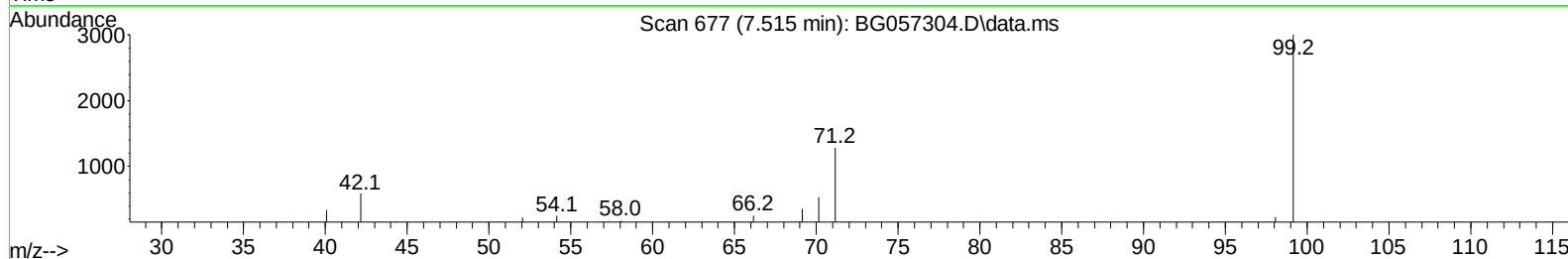
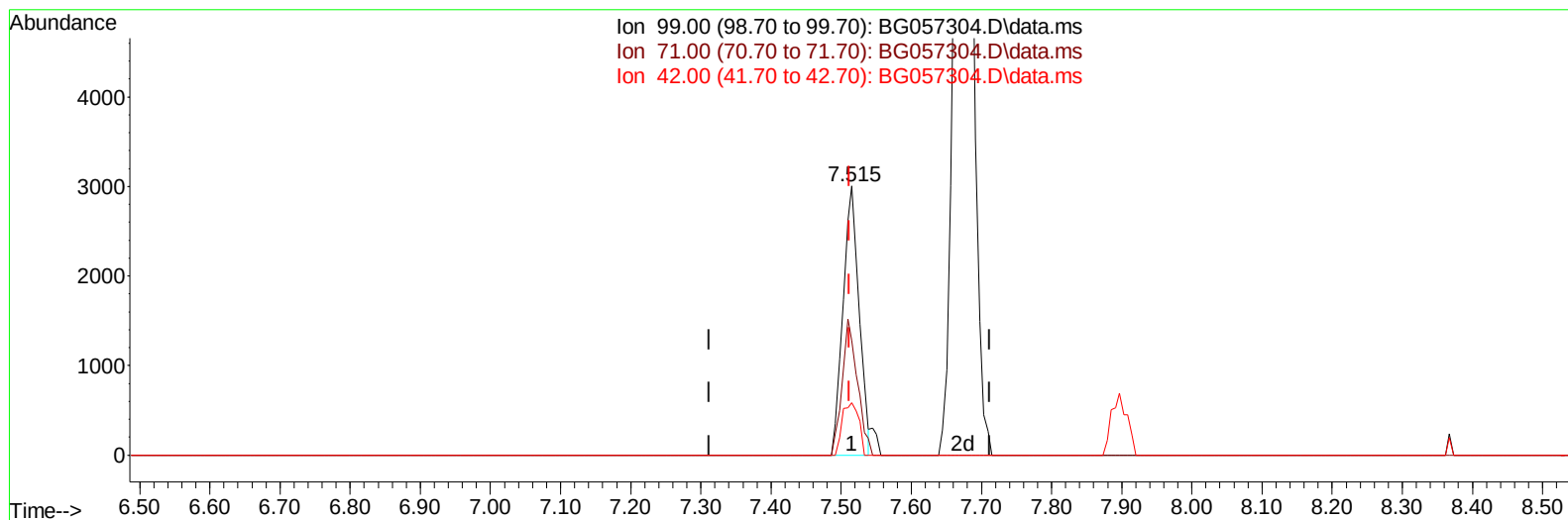
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG050423\
Data File : BG057304.D
Acq On : 4 May 2023 16:47
Operator : CG/JU
Sample : 02613-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AA1

Manual IntegrationsAPPROVED

Quant Time: May 05 01:44:19 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon May 01 12:45:38 2023
Response via : Initial Calibration

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



TIC: BG057304.D\data.ms

(7) Phenol-d5 (S)

7.515min (+ 0.004) 3.89 ng/u1

response 4769

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	42.50	42.64
42.00	18.90	19.54
0.00	0.00	0.00

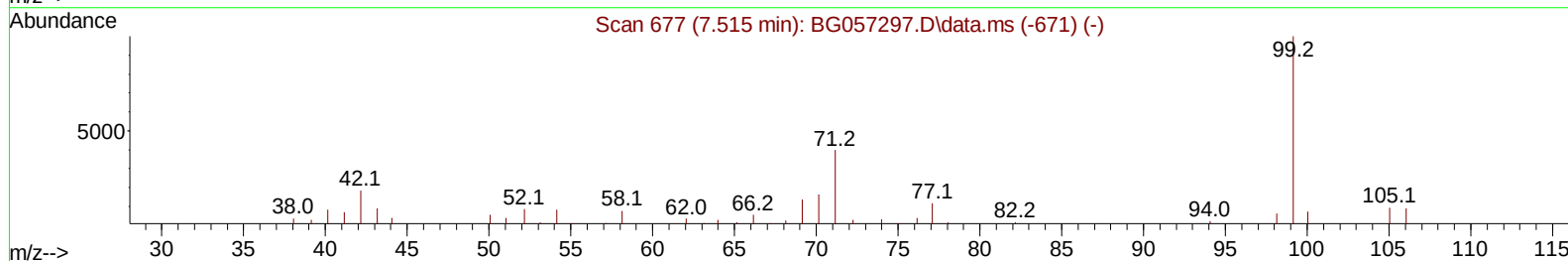
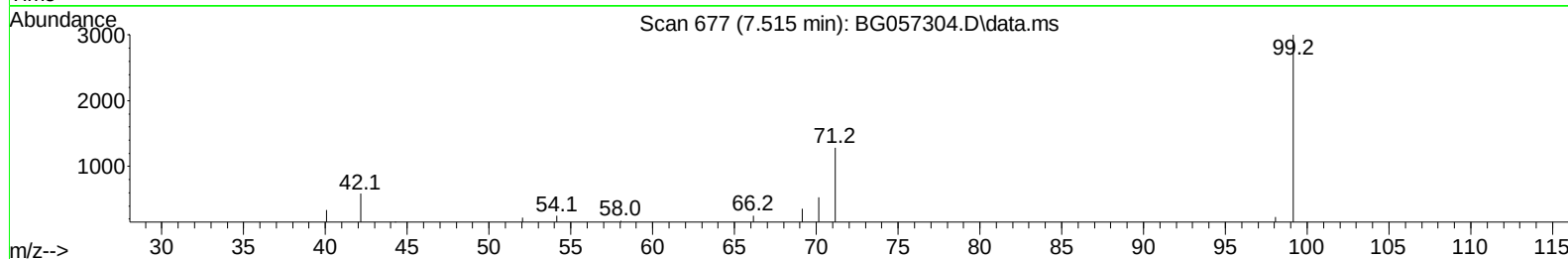
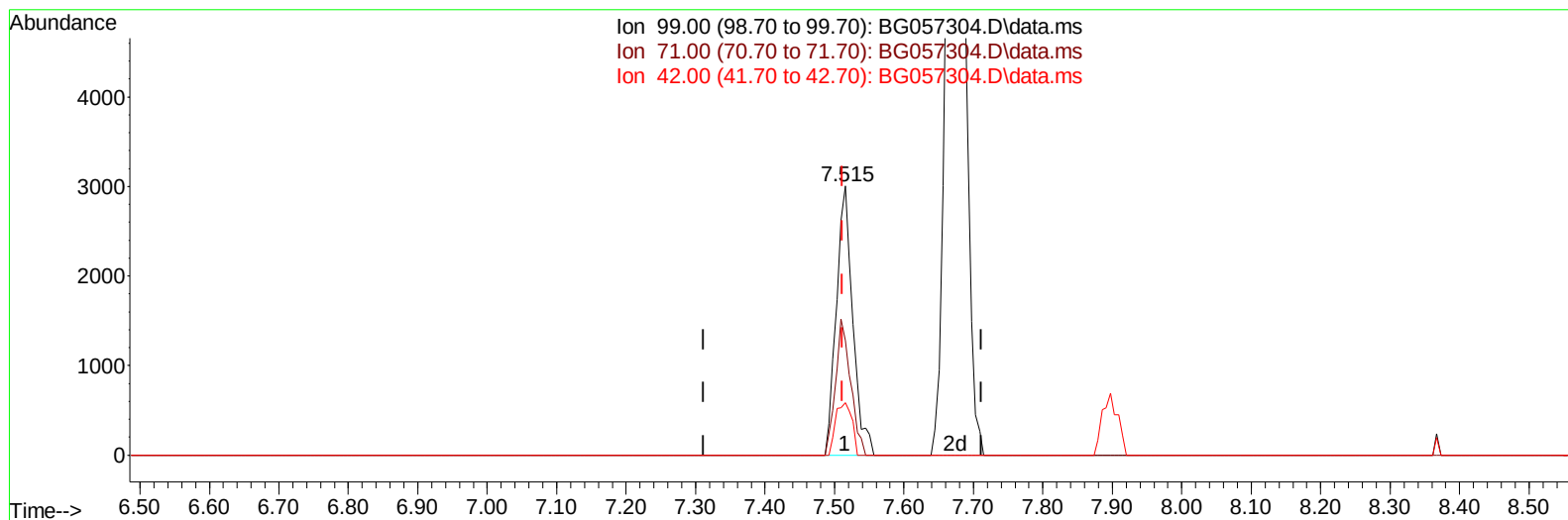
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(7) Phenol-d5 (S)

7.515min (+ 0.004) 4.04 ng/ul m

response 4957

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	42.50	42.64
42.00	18.90	19.54
0.00	0.00	0.00

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Quant Time: May 05 02:27:29 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.373	152	12964	20.000	ng/ul	0.00
20) Naphthalene-d8	11.216	136	56410	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.988	164	49447	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.731	188	133893	20.000	ng/ul	0.00
79) Chrysene-d12	22.043	240	117163	20.000	ng/ul	# 0.00
88) Perylene-d12	25.567	264	120440	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.644	96	1062	3.605	ng/uL	-0.01
4) Pyridine-d5	4.102	84	5436	5.853	ng/ul	0.00
7) Phenol-d5	7.515	99	4957m	4.043	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.674	67	17380	24.102	ng/ul	0.00
11) 2-Chlorophenol-d4	7.897	132	15122	18.717	ng/ul	0.00
15) 4-Methylphenol-d8	9.072	113	10538	11.309	ng/ul	0.00
21) Nitrobenzene-d5	9.554	128	10875	24.208	ng/ul	0.00
24) 2-Nitrophenol-d4	10.282	143	12728	24.278	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.828	165	22730	22.472	ng/ul	0.00
31) 4-Chloroaniline-d4	11.339	131	28634	21.359	ng/ul	0.00
46) Dimethylphthalate-d6	14.377	166	115035	29.360	ng/ul	0.00
49) Acenaphthylene-d8	14.688	160	119383	27.110	ng/ul	0.00
54) 4-Nitrophenol-d4	15.193	143	2478	4.345	ng/ul	0.02
60) Fluorene-d10	15.974	176	103956	28.482	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.098	200	17952	22.630	ng/ul	0.00
73) Anthracene-d10	17.831	188	197517	32.319	ng/ul	0.00
81) Pyrene-d10	20.098	212	225102	32.368	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.326	264	193025	31.457	ng/ul	0.00

Target Compounds Qvalue

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

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