

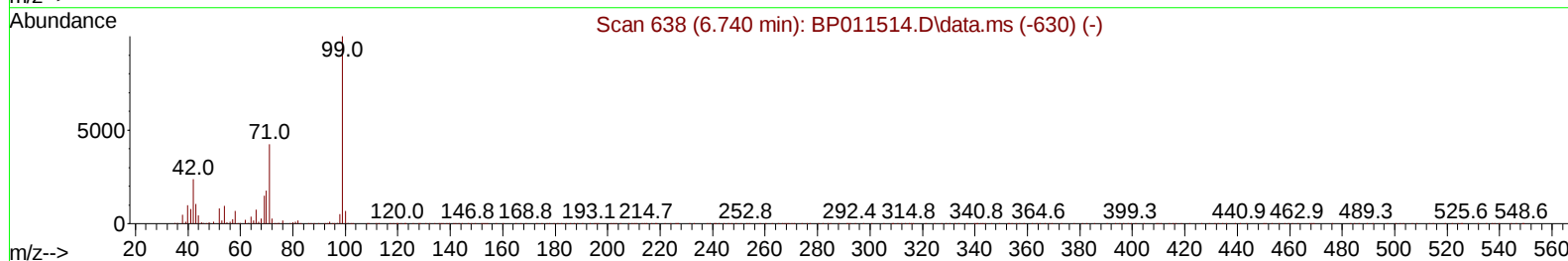
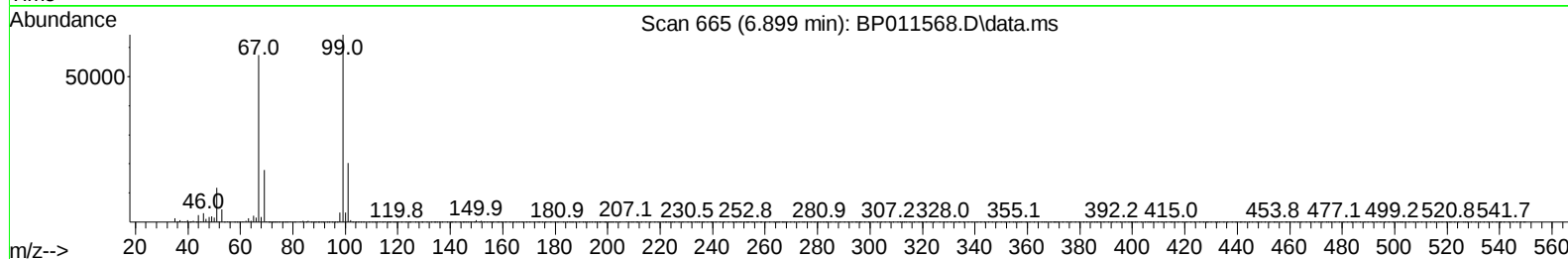
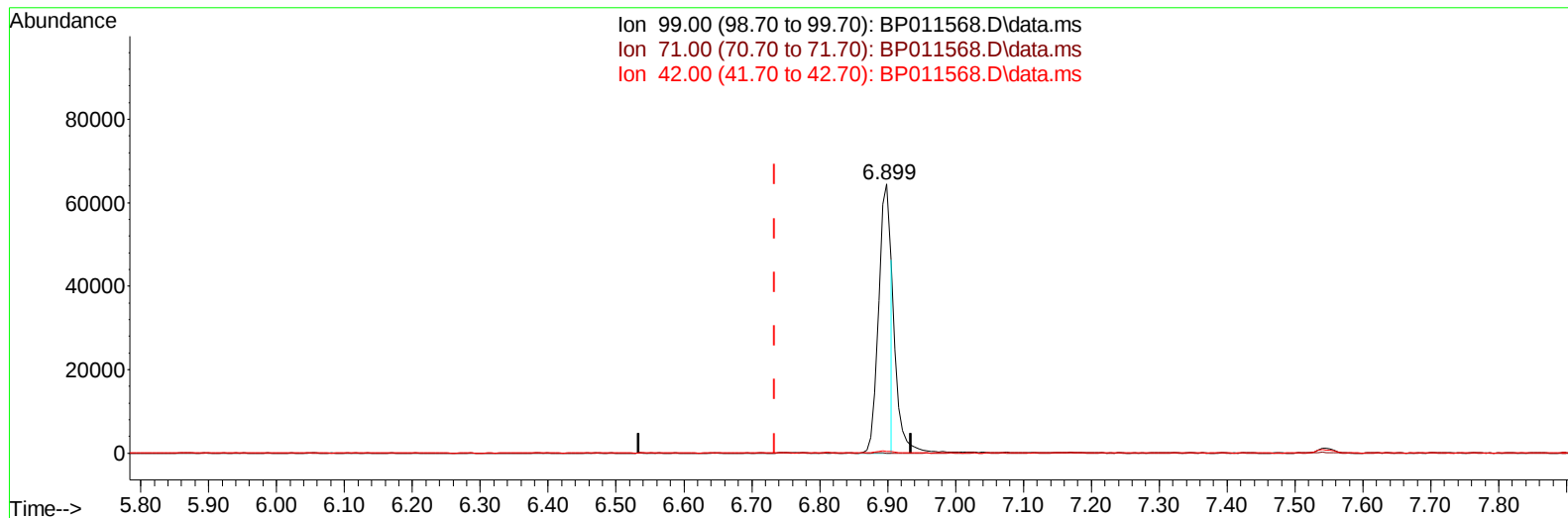
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082422\
Data File : BP011568.D
Acq On : 25 Aug 2022 18:27
Operator : CG/JU
Sample : N4349-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGN64

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 08/26/2022
Supervised By :mohammad ahmed 08/27/2022

Quant Time: Aug 26 09:37:04 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 25 02:14:54 2022
Response via : Initial Calibration



TIC: BP011568.D\data.ms

(7) Phenol-d5 (S)

6.899min (+ 0.165) 16.09 ng/u1

response 79722

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.20	0.02#
42.00	25.30	0.68#
0.00	0.00	0.00

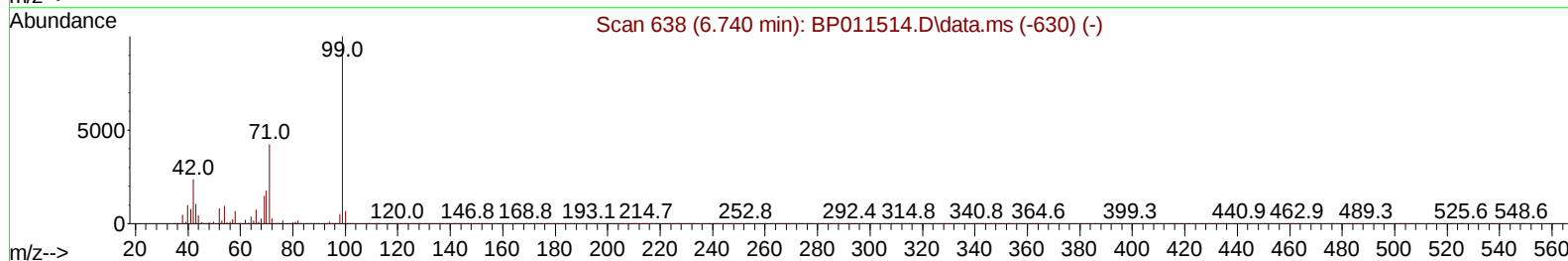
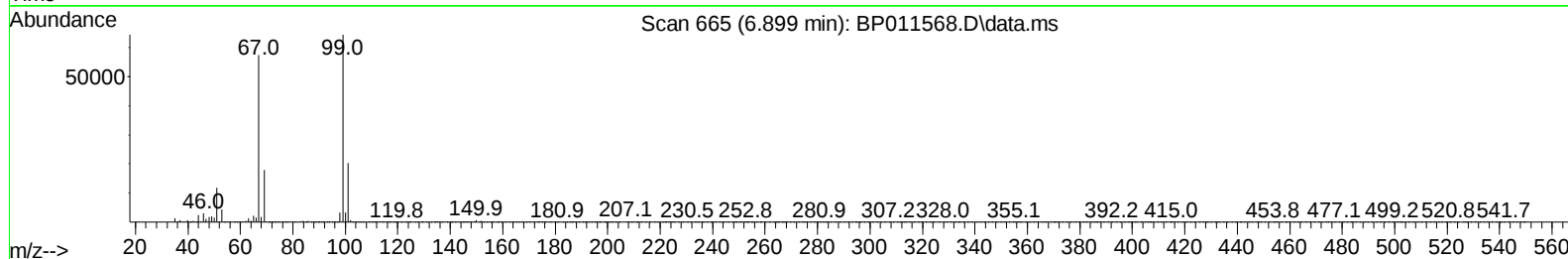
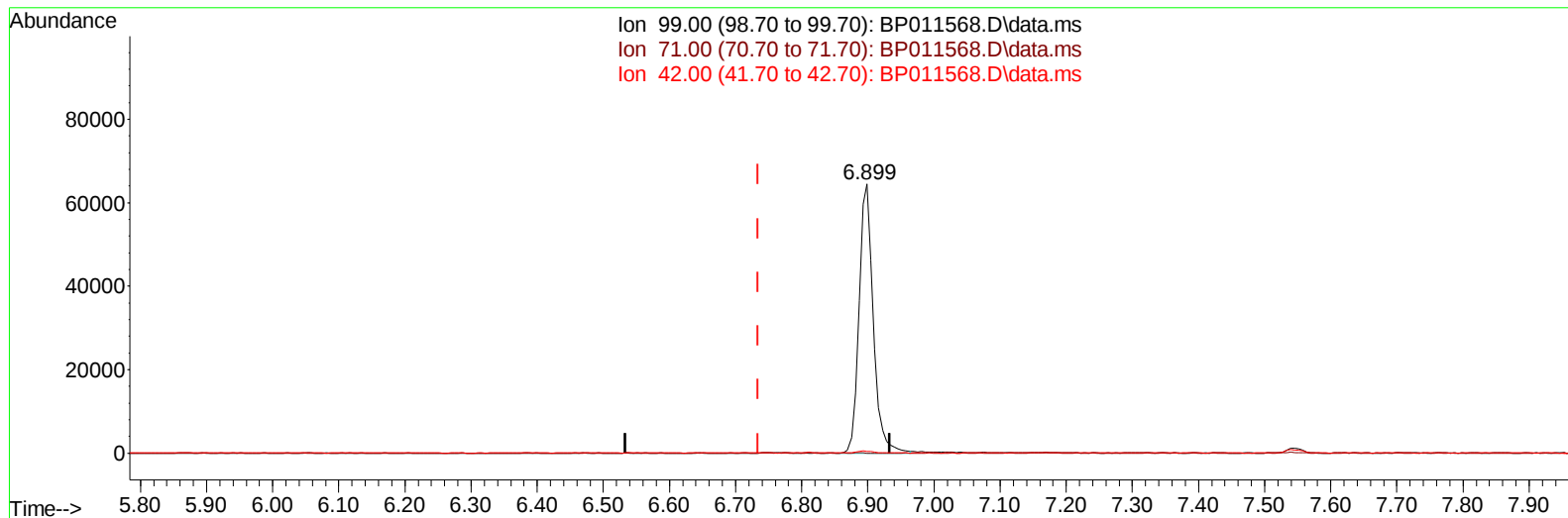
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(7) Phenol-d5 (S)

6.899min (+ 0.165) 19.64 ng/ul m

response 97332

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	43.20	0.01#
42.00	25.30	0.68#
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	7.546	152	53926	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136	226903	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.216	164	139318	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.992	188	283925	20.000	ng/ul	0.00
79) Chrysene-d12	21.121	240	259566	20.000	ng/ul	0.00
88) Perylene-d12	23.404	264	278203	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.064	96	7942	4.570	ng/uL	0.00
4) Pyridine-d5	3.475	84	65008	14.611	ng/ul	0.00
7) Phenol-d5	6.899	99	97332m	19.641	ng/ul	0.16
9) Bis-(2-Chloroethyl)eth...	6.899	67	89487	27.795	ng/ul	0.00
11) 2-Chlorophenol-d4	7.052	132	35	0.010	ng/ul	-0.03
15) 4-Methylphenol-d8	8.275	113	234	0.059	ng/ul	0.00
21) Nitrobenzene-d5	8.710	128	51543	31.553	ng/ul	0.00
24) 2-Nitrophenol-d4	9.446	143	36	0.022	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	9.951	165	7	0.002	ng/ul	-0.02
31) 4-Chloroaniline-d4	10.487	131	125341	24.536	ng/ul	0.00
46) Dimethylphthalate-d6	13.645	166	325349	33.018	ng/ul	0.00
49) Acenaphthylene-d8	13.910	160	345878	31.582	ng/ul	0.00
54) 4-Nitrophenol-d4	14.445	143	7	0.004	ng/ul	-0.01
60) Fluorene-d10	15.222	176	274842	33.154	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.375	200	11	0.007	ng/ul	0.00
73) Anthracene-d10	17.092	188	498452	39.639	ng/ul	0.00
81) Pyrene-d10	19.351	212	549081	41.810	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.257	264	522426	38.604	ng/ul	0.00

Target Compounds Qvalue

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

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