Quantitation Report (Qedit)

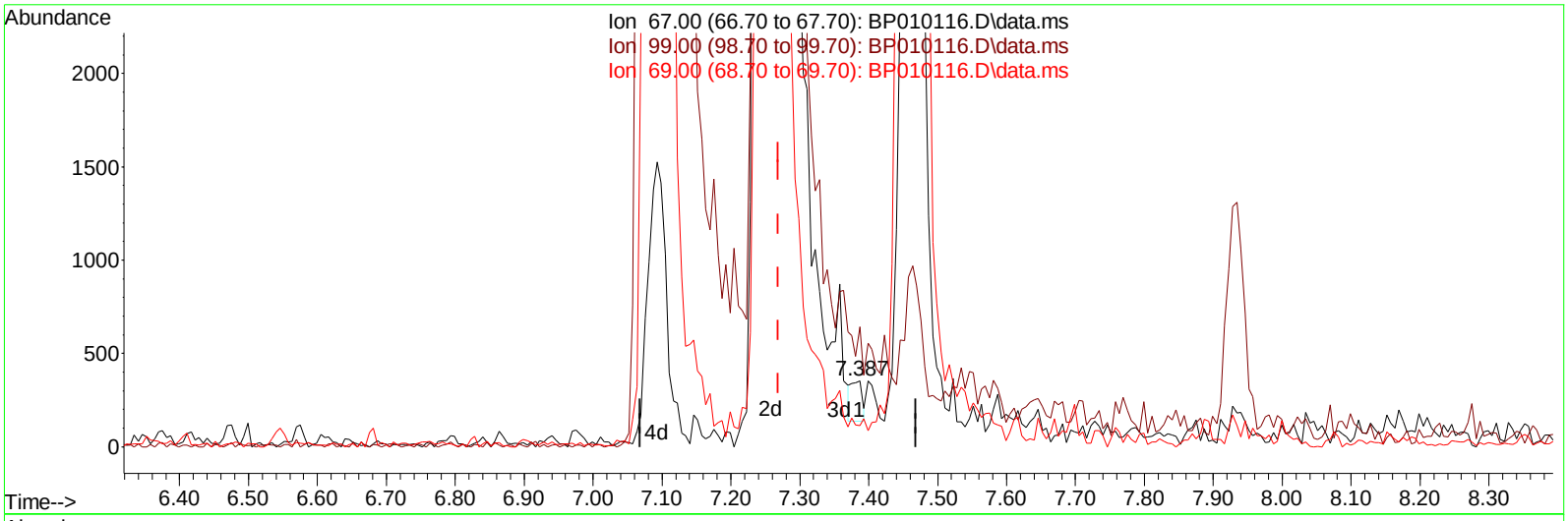
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042722\
Data File : BP010116.D
Acq On : 27 Apr 2022 11:03
Operator : CG/JU
Sample : PB144346BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

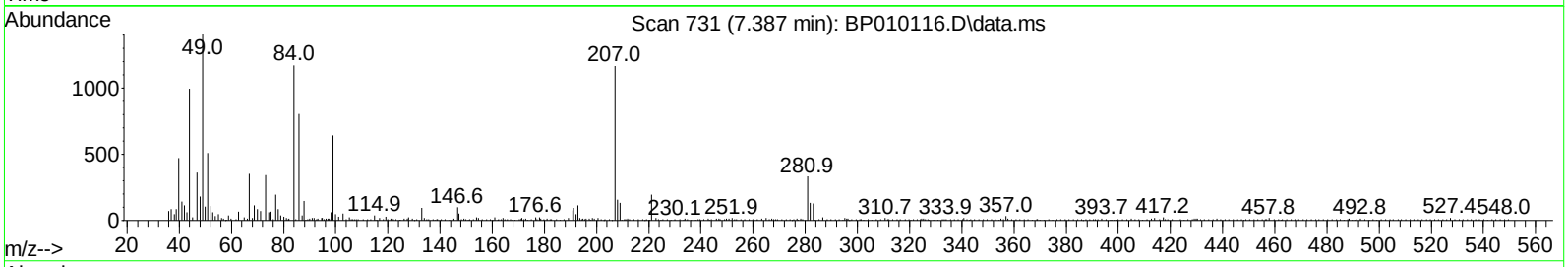
Quant Time: Apr 28 03:22:51 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

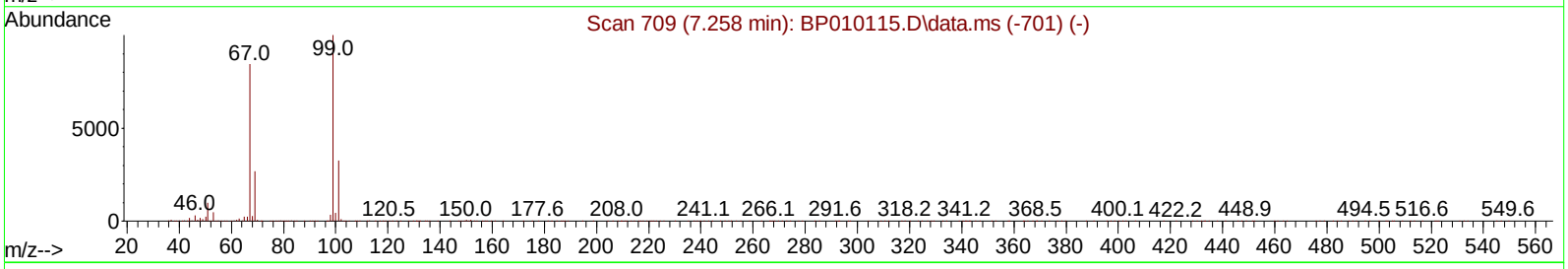
Instrument :
BNA_P
ClientSampleId :
SBLK346

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/28/2022
Supervised By :Yogesh Patel 05/05/2022







TIC: BP010116.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.387min (+ 0.118) 0.03 ng/ul

response 244

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	182.15
69.00	40.10	32.58
0.00	0.00	0.00

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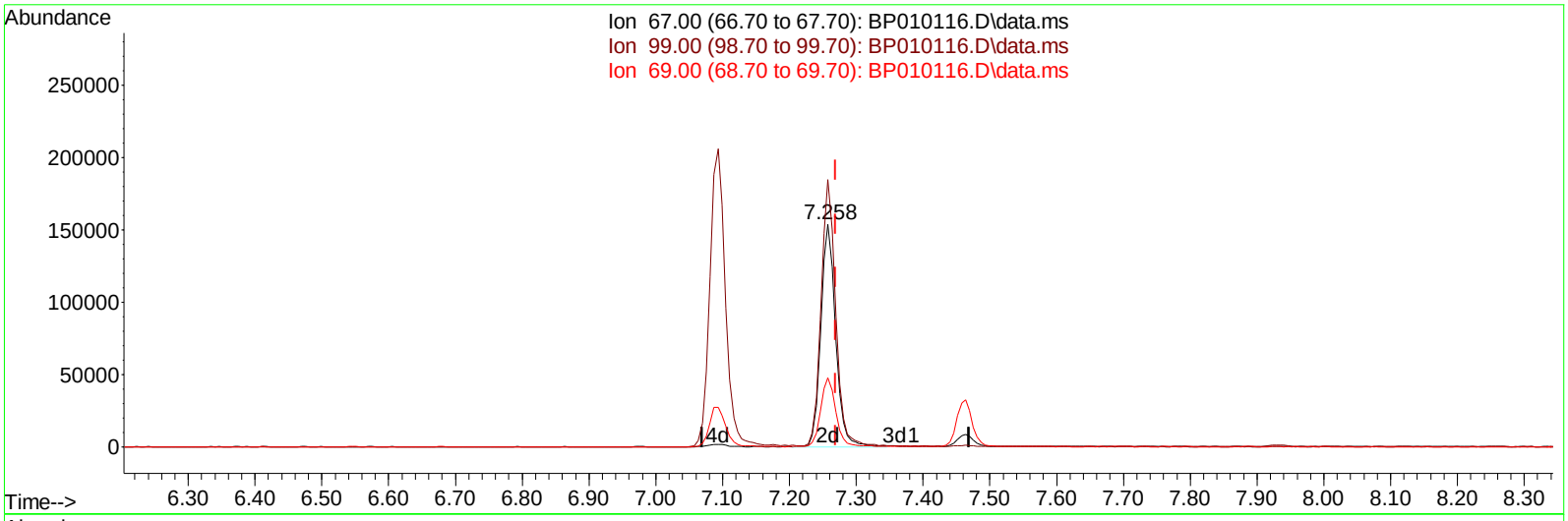

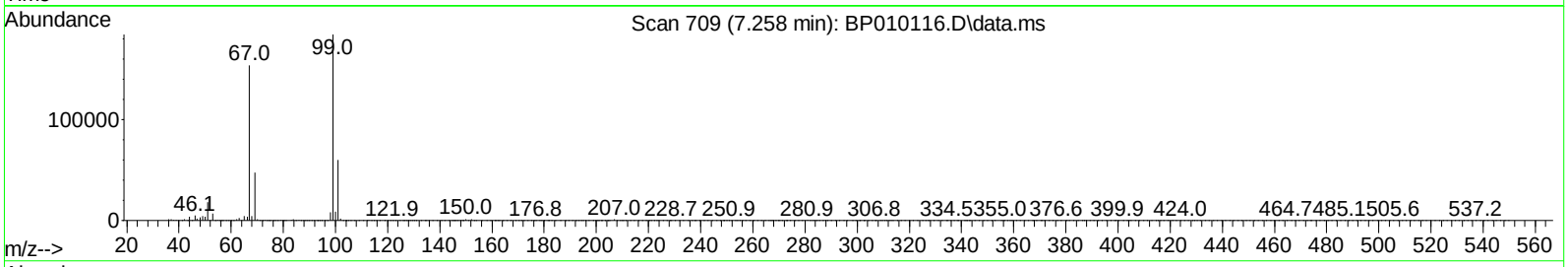

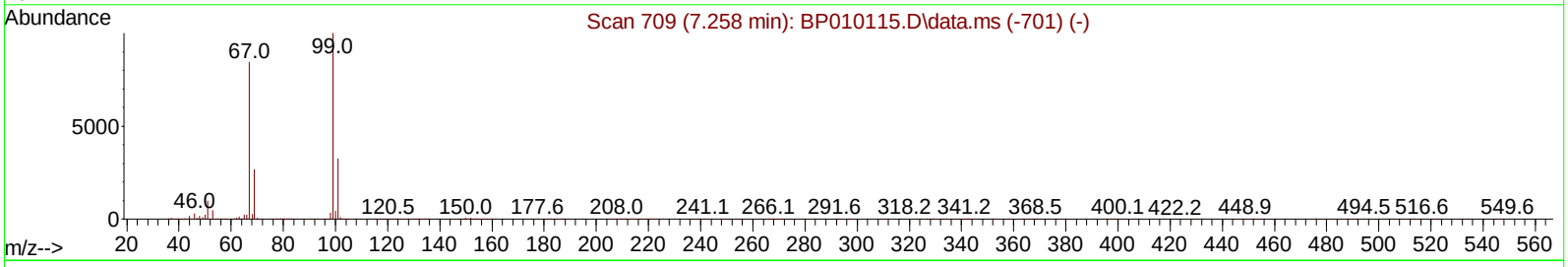

TIC: BP010116.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.258min (-0.012) 28.02 ng/ul m

response 236570

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	119.88#
69.00	40.10	30.89#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	154699	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	672942	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.563	164	469624	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.316	188	1012396	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	879490	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	689229	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	19757	5.569	ng/uL	0.00
4) Pyridine-d5	3.781	84	242753	24.324	ng/ul	0.00
7) Phenol-d5	7.093	99	334121	23.745	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	236570m	28.018	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.464	132	279342	26.805	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	294187	25.234	ng/ul	-0.01
21) Nitrobenzene-d5	9.087	128	147795	27.614	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.816	143	162348	27.774	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.358	165	275258	24.189	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	409067	25.071	ng/ul	-0.01
46) Dimethylphthalate-d6	13.969	166	1089021	29.810	ng/ul	-0.01
49) Acenaphthylene-d8	14.257	160	1176481	26.609	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	136888	20.354	ng/ul	0.00
60) Fluorene-d10	15.557	176	908843	28.921	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	144903	22.636	ng/ul	0.00
73) Anthracene-d10	17.416	188	1483969	30.298	ng/ul	0.00
81) Pyrene-d10	19.651	212	1720644	34.401	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1146922	31.741	ng/ul	0.00

Target Compounds Qvalue

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

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