

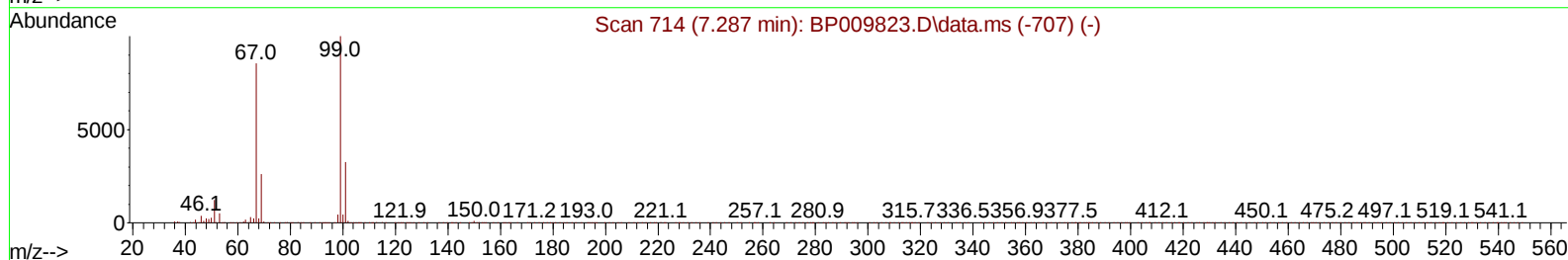
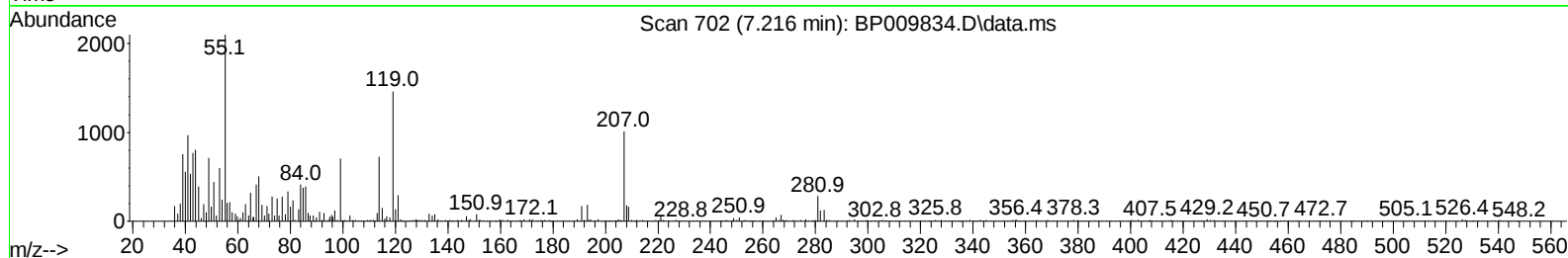
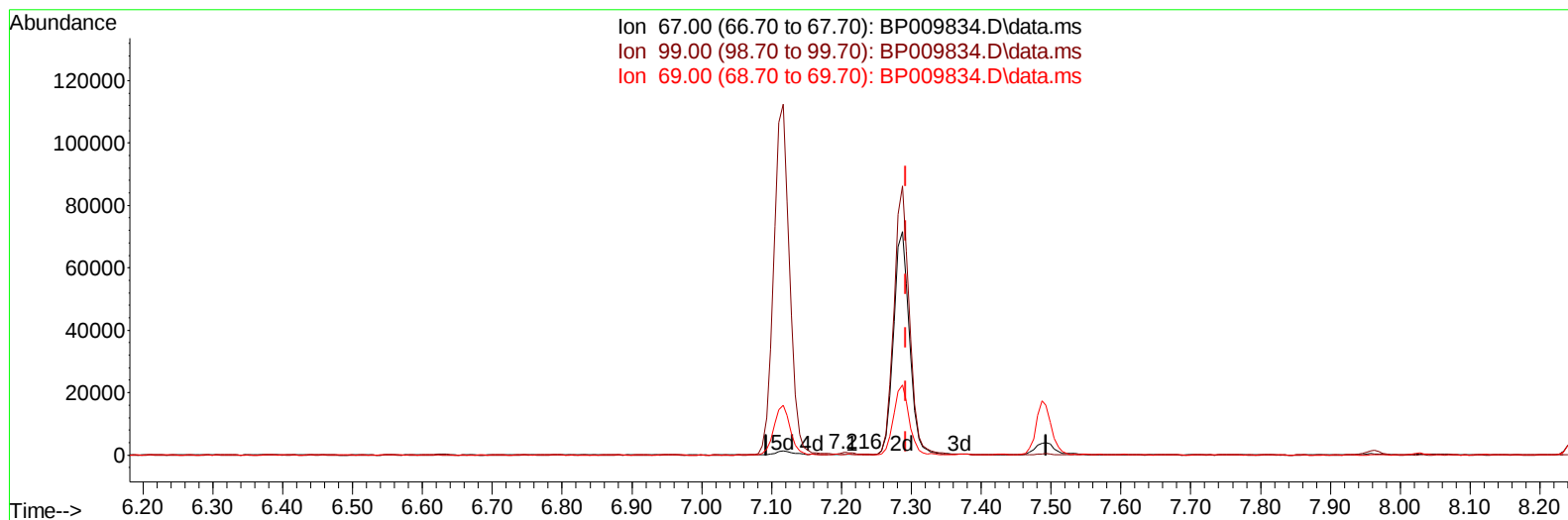
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
 Data File : BP009834.D
 Acq On : 14 Apr 2022 18:39
 Operator : CG/JU
 Sample : N2293-09
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNN9

Manual IntegrationsAPPROVED

Quant Time: Apr 14 21:47:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 13 02:15:22 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/15/2022
 Supervised By :Yogesh Patel 04/22/2022



TIC: BP009834.D\data.ms

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

7.216min (-0.076) 0.06 ng/u1

response 492

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	171.01
69.00	40.10	45.41
0.00	0.00	0.00

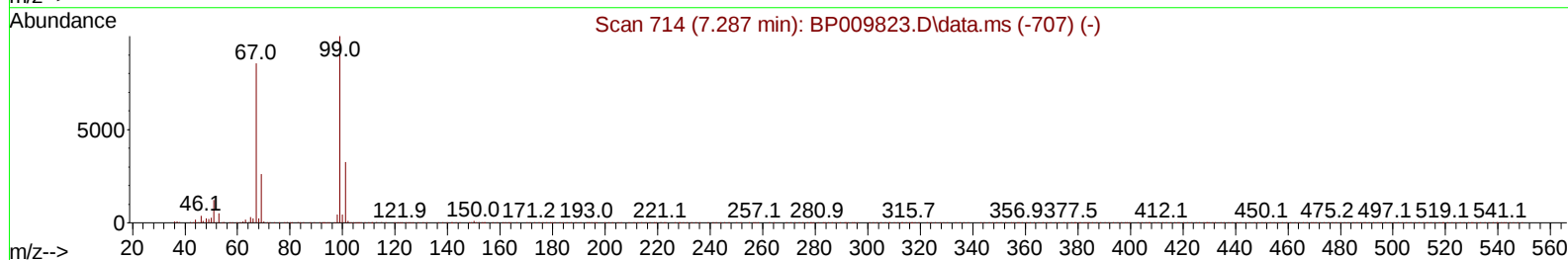
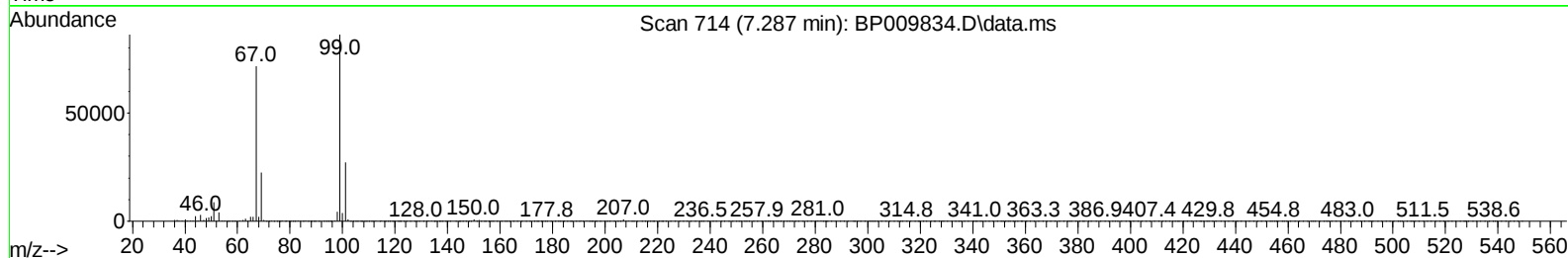
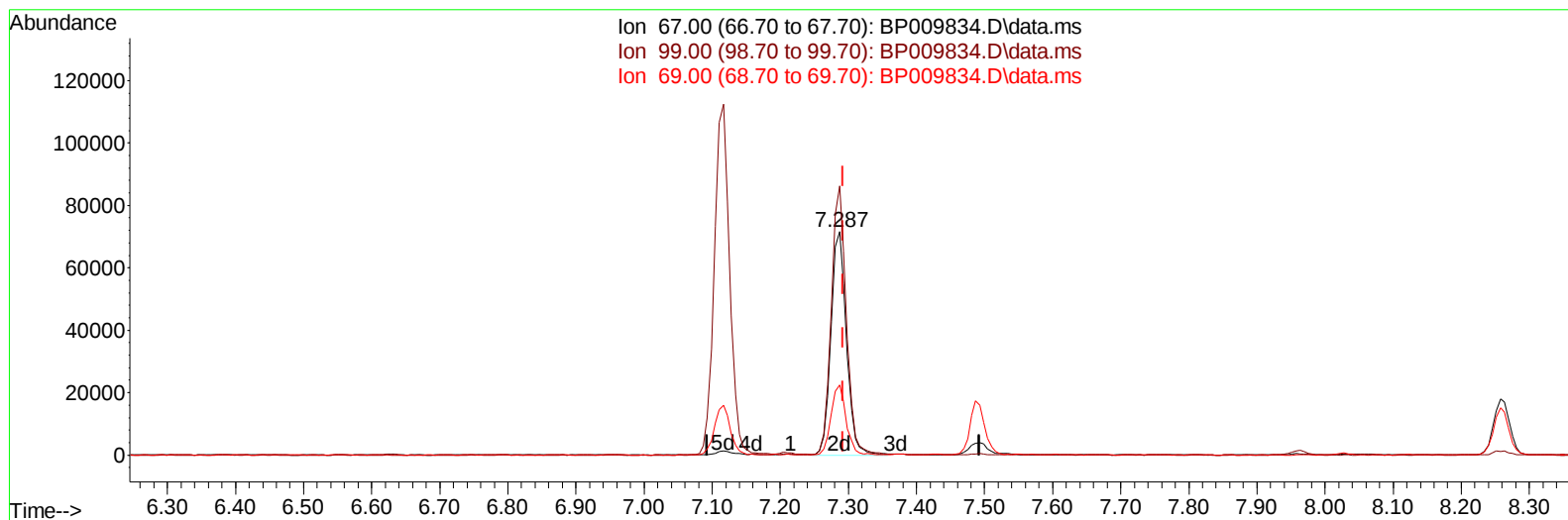
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TIC: BP009834.D\data.ms

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

7.287min (-0.006) 14.97 ng/ul m

response 113450

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	120.22#
69.00	40.10	31.28#
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	7.963	152	144991	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	652252	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.593	164	473384	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	1074465	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.422	240	1125351	20.000	ng/ul	#-0.01
88) Perylene-d12	23.886	264	1027182	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	8370	2.576	ng/uL	0.00
4) Pyridine-d5	3.805	84	62997	7.033	ng/ul	0.00
7) Phenol-d5	7.116	99	176410	13.859	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	113450m	14.971	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	136802	14.247	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	159006	14.963	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	73364	15.597	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.846	143	76559	14.846	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	152682	14.149	ng/ul	0.00
31) 4-Chloroaniline-d4	10.899	131	202594	13.182	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	638452	17.170	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	618160	14.032	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	106310	15.764	ng/ul	-0.01
60) Fluorene-d10	15.581	176	518915	16.589	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.692	200	87786	13.792	ng/ul	0.00
73) Anthracene-d10	17.439	188	932112	18.119	ng/ul	-0.01
81) Pyrene-d10	19.669	212	1188514	19.338	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1027251	19.204	ng/ul	-0.02
Target Compounds						
					Qvalue	
8) Phenol	7.140	94	16553	1.299	ng/ul#	86
18) 4-Methylphenol	8.722	108	11715	1.089	ng/ul	90

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

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