

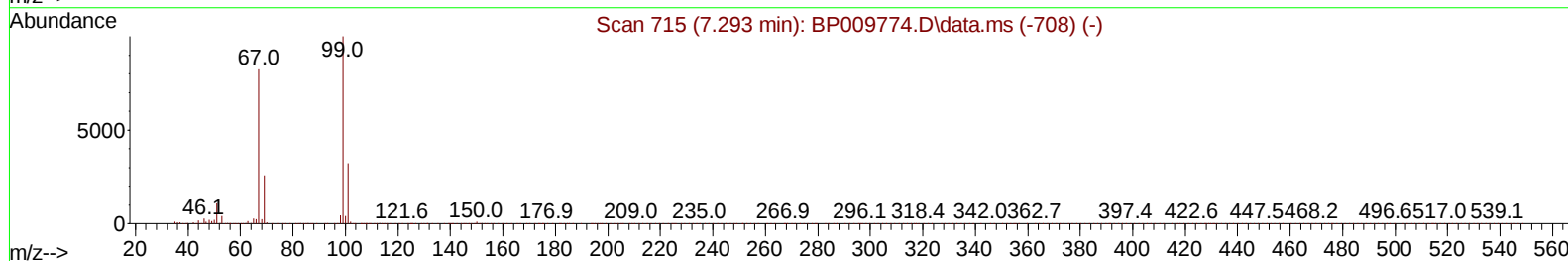
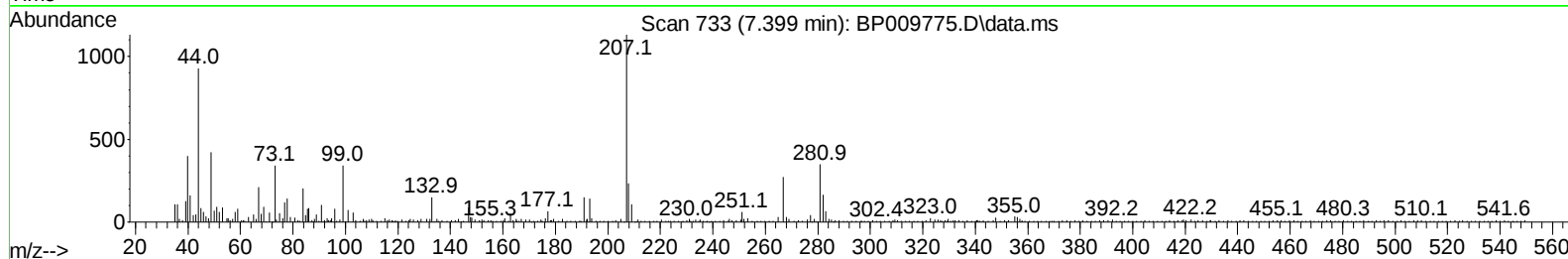
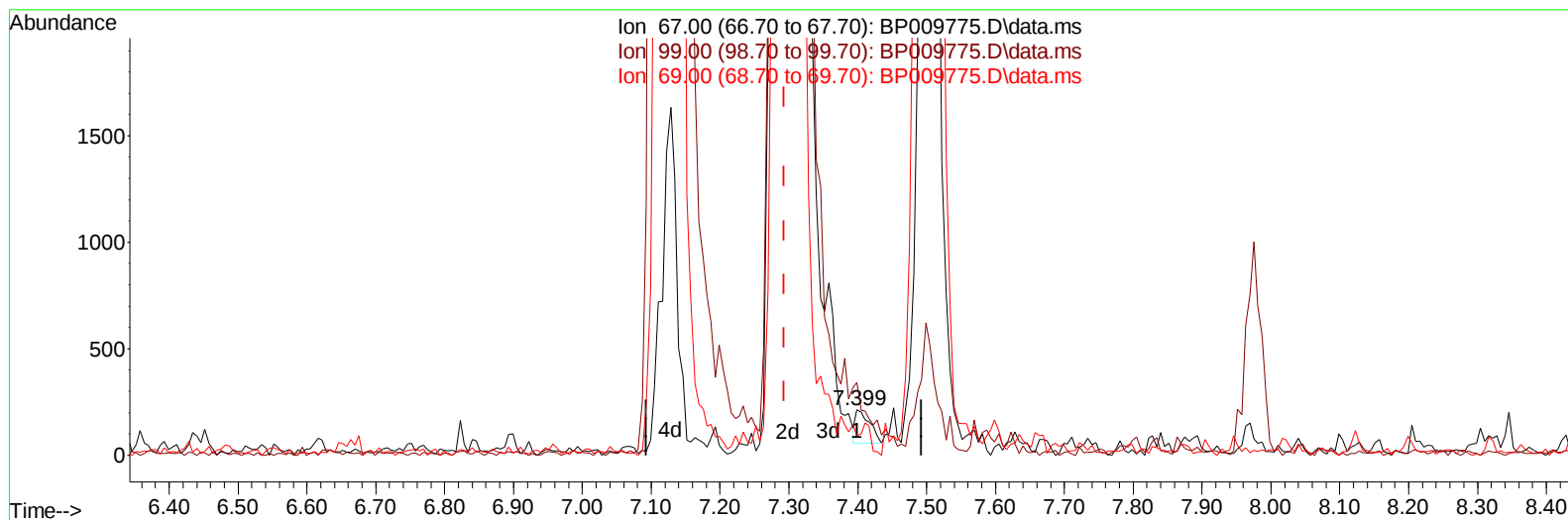
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009775.D
 Acq On : 12 Apr 2022 09:17
 Operator : CG/JU
 Sample : PB144003BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK003

Manual IntegrationsAPPROVED

Quant Time: Apr 13 02:19:06 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 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009775.D\data.ms

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

7.399min (+ 0.106) 0.04 ng/u1

response 220

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	160.85
69.00	40.10	43.40
0.00	0.00	0.00

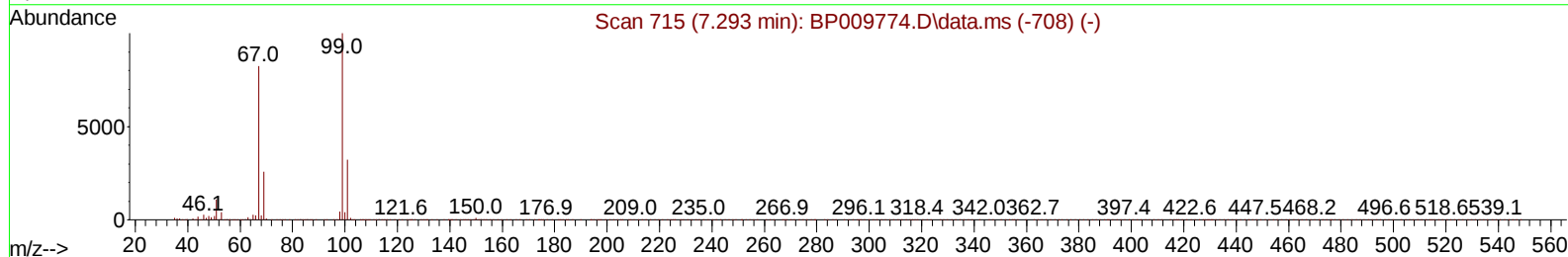
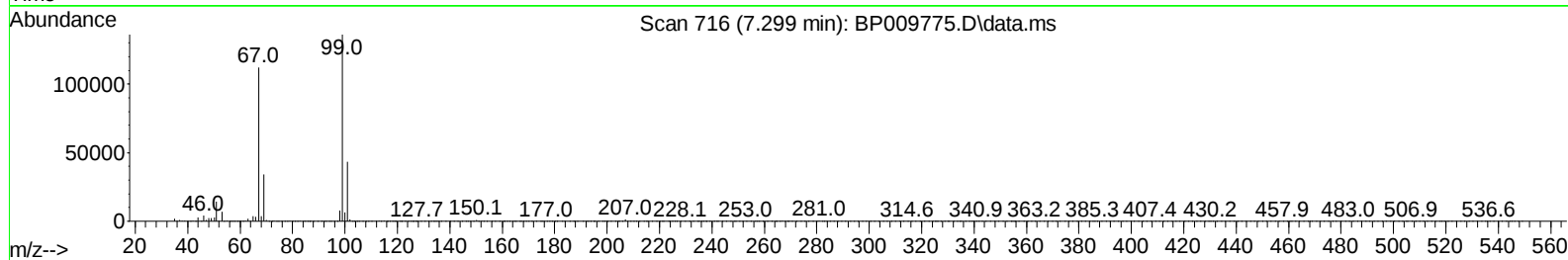
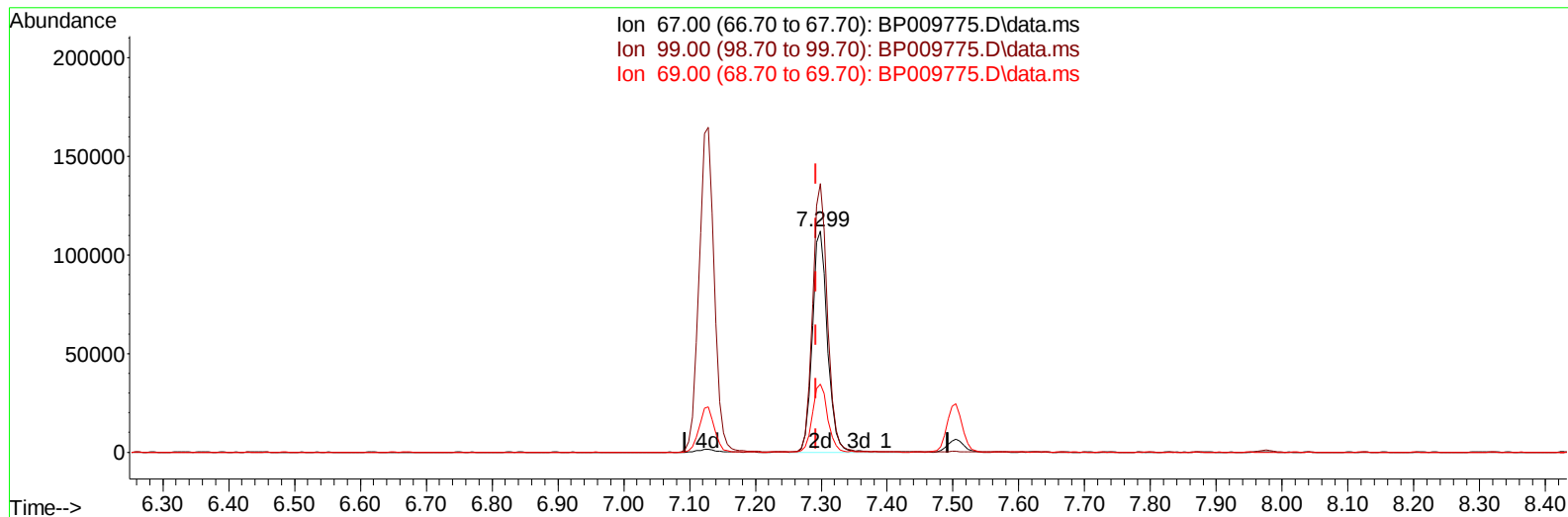
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TIC: BP009775.D\data.ms

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

7.299min (+ 0.006) 33.47 ng/ul m

response 179911

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	121.21#
69.00	40.10	30.62#
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	7.975	152	102835	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	452066	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.610	164	336202	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.363	188	754408	20.000	ng/ul	# 0.01
79) Chrysene-d12	21.439	240	811273	20.000	ng/ul	# 0.00
88) Perylene-d12	23.916	264	762871	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	15611	6.775	ng/uL	0.00
4) Pyridine-d5	3.805	84	212238	33.408	ng/ul	0.00
7) Phenol-d5	7.128	99	264188	29.262	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	179911m	33.473	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	212879	31.259	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	232199	30.809	ng/ul	0.00
21) Nitrobenzene-d5	9.134	128	114043	34.983	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	122891	34.384	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	220005	29.416	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	341254	32.036	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	875123	33.137	ng/ul	0.01
49) Acenaphthylene-d8	14.298	160	946101	30.239	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	143945	30.055	ng/ul	0.00
60) Fluorene-d10	15.604	176	736988	33.173	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.710	200	123631	27.663	ng/ul	0.01
73) Anthracene-d10	17.457	188	1206107	33.391	ng/ul	0.00
81) Pyrene-d10	19.686	212	1481792	33.443	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.757	264	1399167	35.219	ng/ul	0.02

Target Compounds Qvalue

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

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