

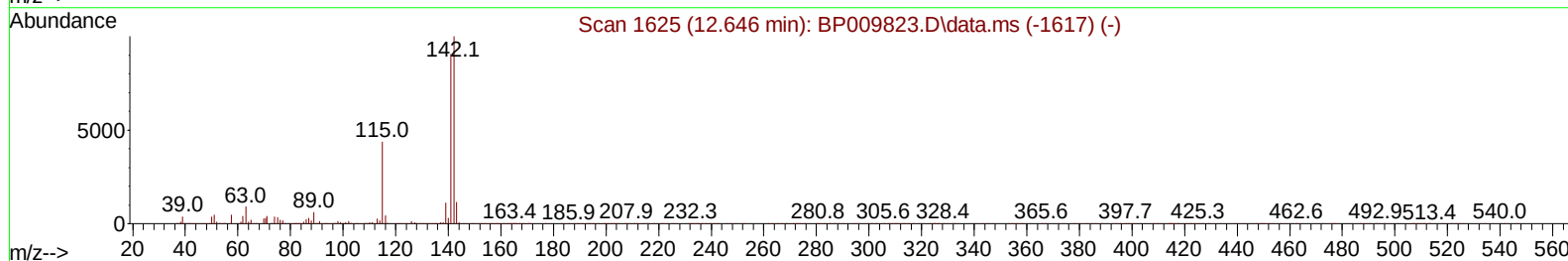
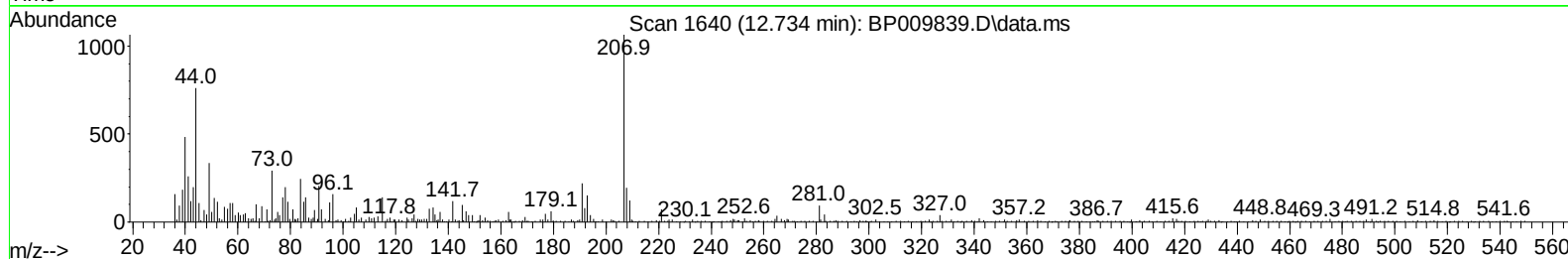
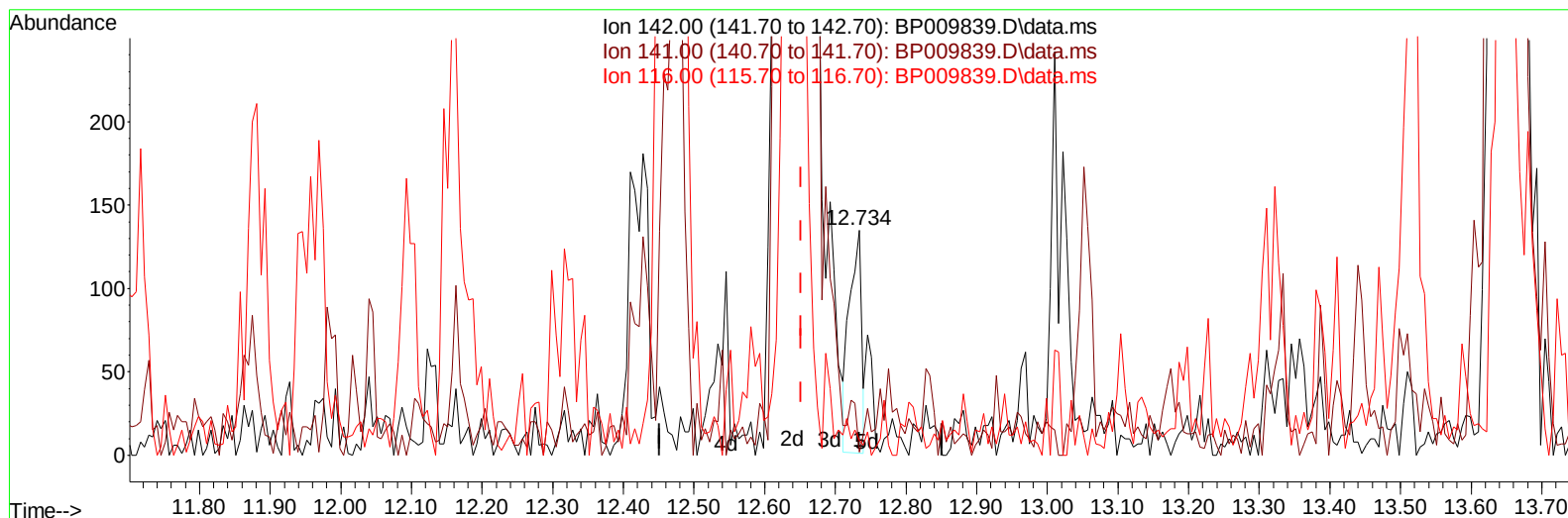
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
 Data File : BP009839.D
 Acq On : 14 Apr 2022 21:28
 Operator : CG/JU
 Sample : N2293-17
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNP5

Manual IntegrationsAPPROVED

Quant Time: Apr 14 22:35:34 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: BP009839.D\data.ms

(37) 1-Methylnaphthalene

12.734min (+ 0.082) 0.01 ng/ul

response 161

Ion	Exp%	Act%
142.00	100.00	100.00
141.00	87.30	90.37
116.00	7.00	6.67
0.00	0.00	0.00

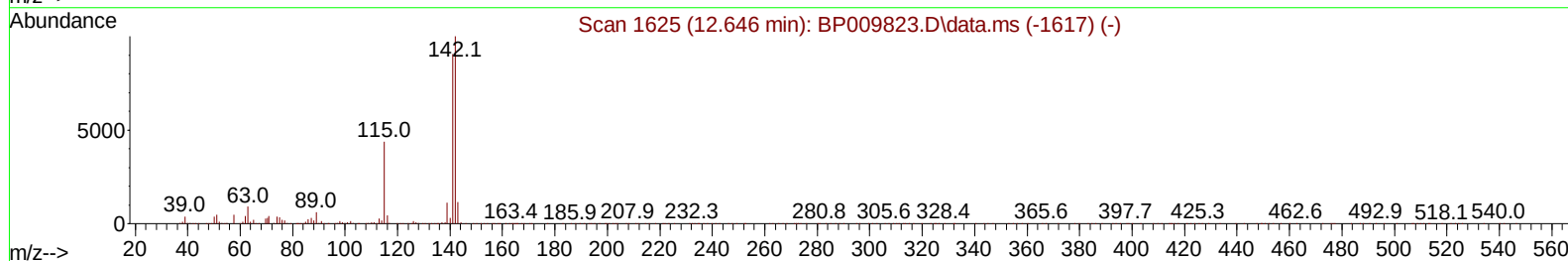
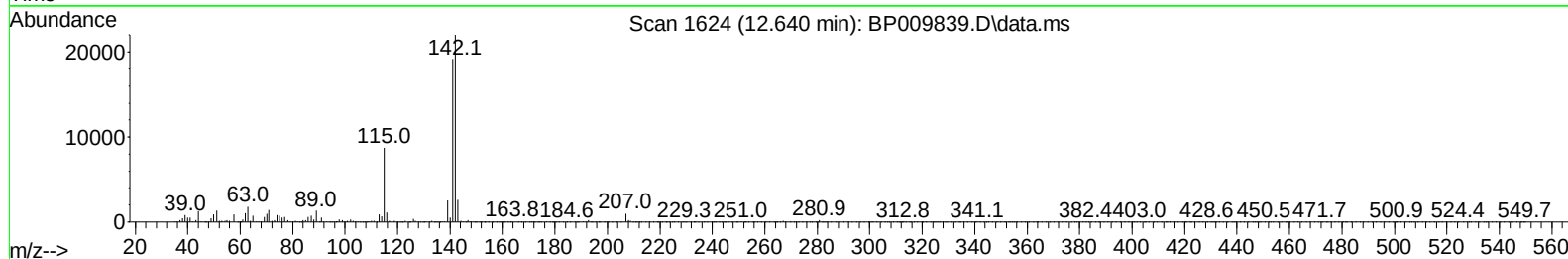
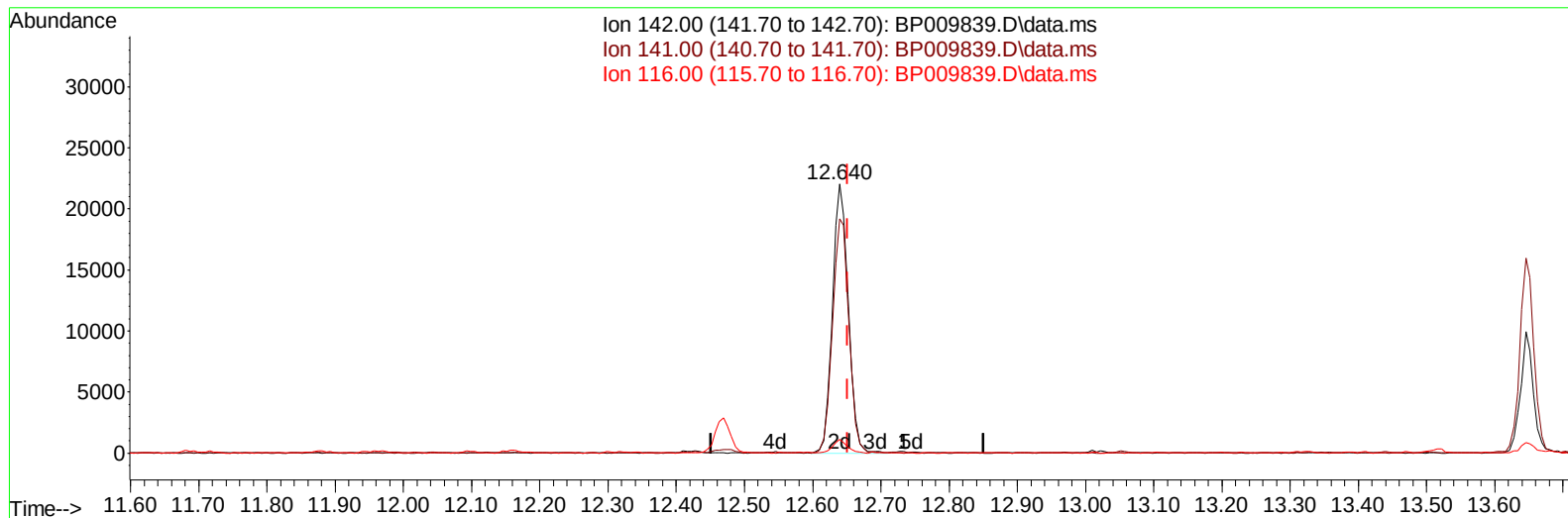
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TIC: BP009839.D\data.ms

(37) 1-Methylnaphthalene

12.640min (-0.012) 2.09 ng/ul m

response 35384

Ion	Exp%	Act%
142.00	100.00	100.00
141.00	87.30	87.18
116.00	7.00	5.19#
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	98286	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.769	136	447286	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	329044	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	765630	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.421	240	875040	20.000	ng/ul	#-0.01
88) Perylene-d12	23.886	264	892584	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	11105	5.043	ng/uL	0.00
4) Pyridine-d5	3.799	84	155177	25.556	ng/ul	0.00
7) Phenol-d5	7.116	99	231698	26.851	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	148249	28.859	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.493	132	182302	28.008	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	205921	28.587	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	96967	30.063	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.846	143	107276	30.336	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	203223	27.463	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.899	131	284279	26.973	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	737478	28.533	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	823718	26.900	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.769	143	116665	24.889	ng/ul	-0.01
60) Fluorene-d10	15.581	176	643485	29.595	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.692	200	96972	21.380	ng/ul	0.00
73) Anthracene-d10	17.439	188	1052388	28.709	ng/ul	-0.01
81) Pyrene-d10	19.669	212	1338830	28.015	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	1371491	29.505	ng/ul	-0.01
Target Compounds						
37) 1-Methylnaphthalene	12.640	142	35384m	2.085	ng/ul	Qvalue

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

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