

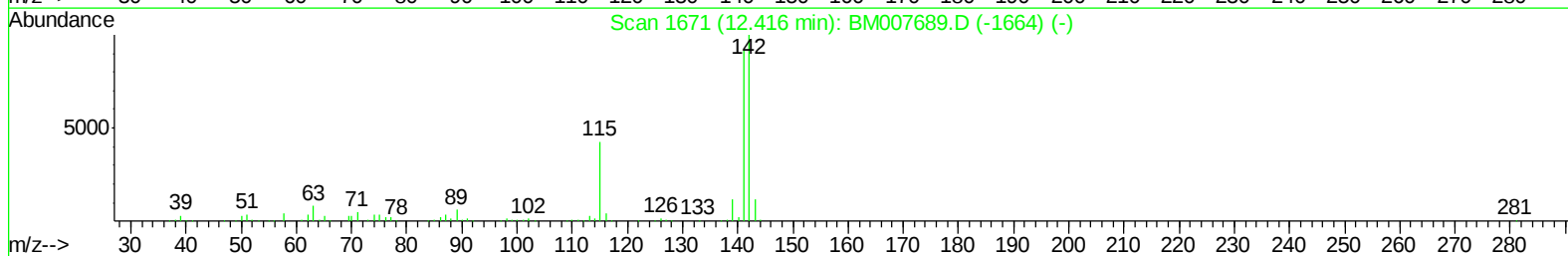
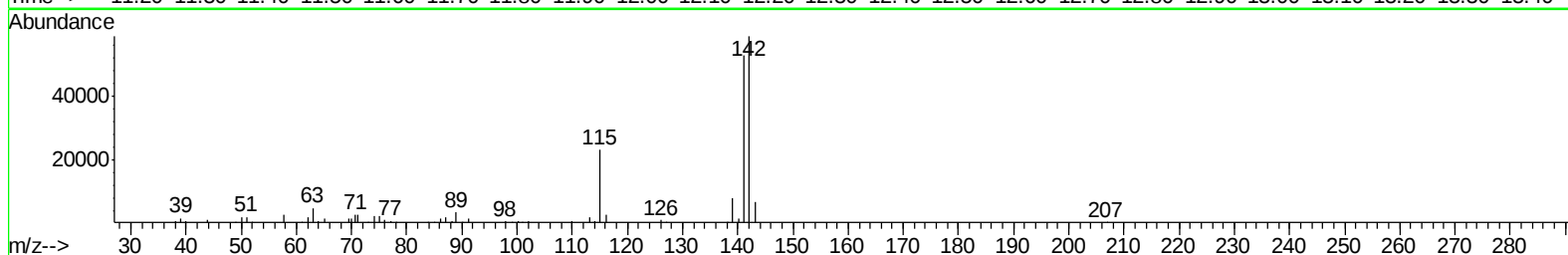
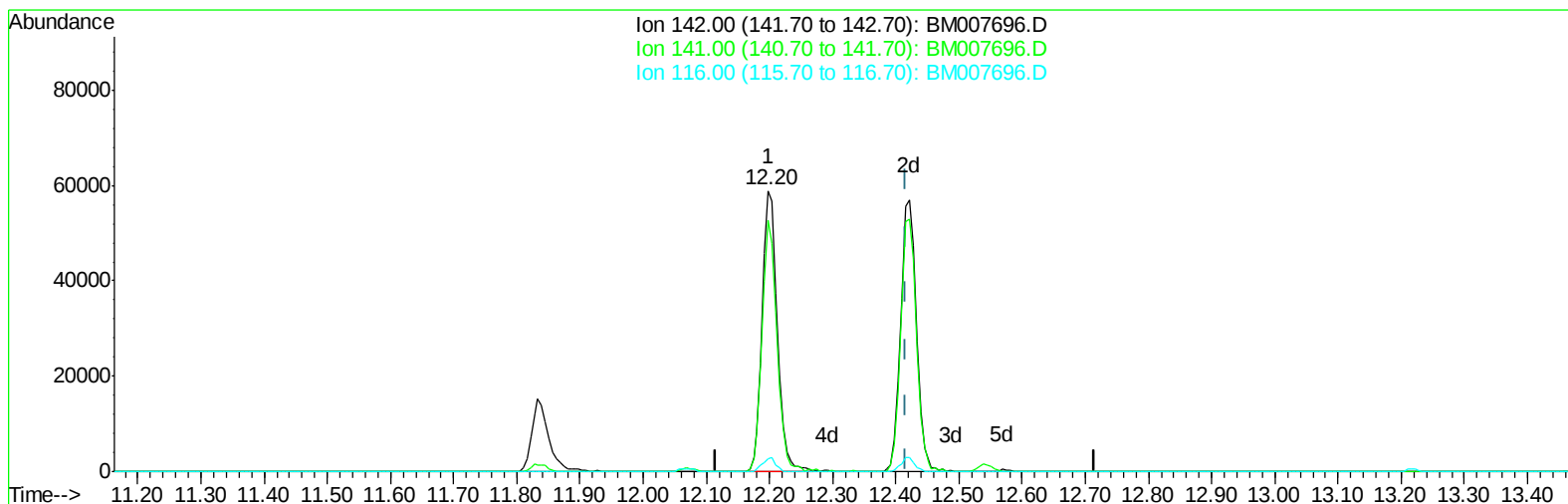
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101916\
Data File : BM007696.D
Acq On : 20 Oct 2016 06:35
Operator : UM/SJ
Sample : MDL-06-W
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
MDL-06-W

Manual Integrations
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Quant Time: Oct 20 07:59:05 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM101916-MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 20 03:26:07 2016
Response via : Initial Calibration



TIC: BM007696.D

(12) 1-Methylnaphthalene

12.198min (-0.218) 3.02ng/ul

response 96373

Ion	Exp%	Act%
142.00	100	100
141.00	93.60	89.49
116.00	3.90	4.54
0.00	0.00	0.00

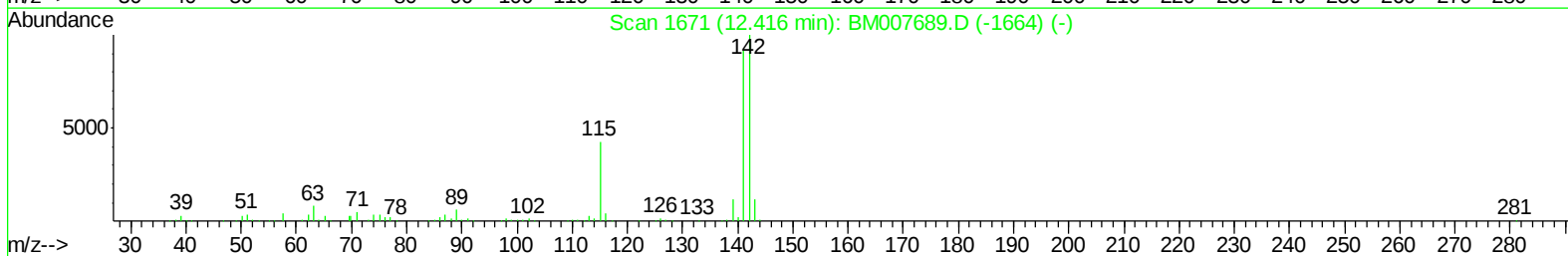
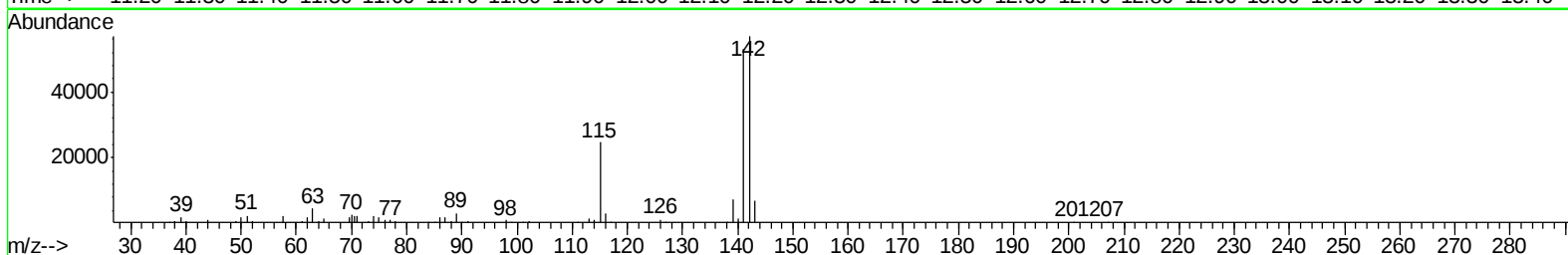
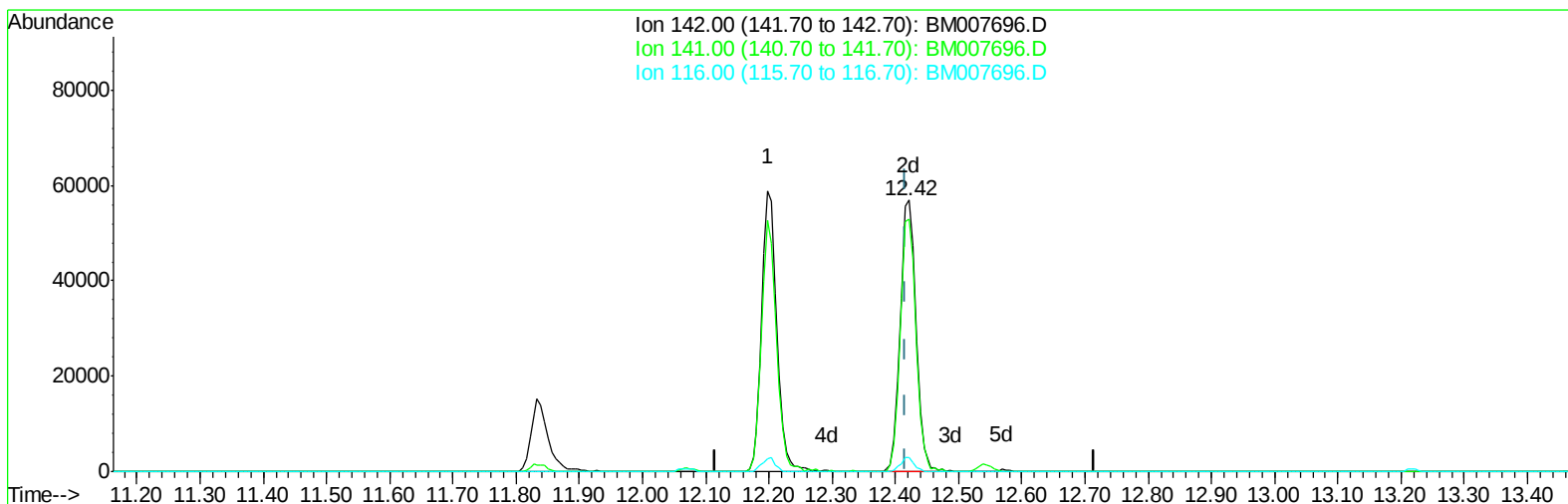
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TIC: BM007696.D

(12) 1-Methylnaphthalene

12.422min (+0.006) 3.02ng/ul m

response 96513

Ion	Exp%	Act%
142.00	100	100
141.00	93.60	93.31
116.00	3.90	5.15#
0.00	0.00	0.00

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 ALS Vial : 25 Sample Multiplier: 1

Instrument :
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Quant Time: Oct 20 08:00:04 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM101916-MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 03:26:07 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	242882	20.00	ng/ul	0.00
7) Naphthalene-d8	10.53	136	928033	20.00	ng/ul	0.00
13) Acenaphthene-d10	14.39	164	584853	20.00	ng/ul	0.00
18) Phenanthrene-d10	17.13	188	1149631	20.00	ng/ul	0.00
23) Chrysene-d12	21.32	240	1449900	20.00	ng/ul	0.00
25) Perylene-d12	23.57	264	1476913	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.29	96	40182	7.06	ng/uL	0.00
3) Phenol-d5	6.93	99	606223	32.21	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.09	67	364515	32.68	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	481799	33.35	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	468759	31.70	ng/ul	0.00
8) Nitrobenzene-d5	8.91	128	225605	34.42	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	246829	34.50	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.16	165	450965	32.03	ng/ul	0.00
11) 4-Chloroaniline-d4	10.67	131	567839	35.54	ng/ul	0.00
14) Dimethylphthalate-d6	13.80	166	1325833	32.50	ng/ul	0.00
15) Acenaphthylene-d8	14.08	160	1601136	33.48	ng/ul	0.00
16) 4-Nitrophenol-d4	14.60	143	185368	27.63	ng/ul	0.00
17) Fluorene-d10	15.38	176	1141344	32.42	ng/ul	0.00
19) 4,6-Dinitro-2-methylphenol	15.51	200	198356	28.47	ng/ul	0.00
20) Anthracene-d10	17.23	188	1700373	32.53	ng/ul	0.00
24) Pyrene-d10	19.53	212	1989897	33.90	ng/ul	0.00
26) Benzo(a)pyrene-d12	23.43	264	2129911	33.11	ng/ul	0.00

Target Compounds						Qvalue
12) 1-Methylnaphthalene	12.42	142	96513m	3.02	ng/ul	
21) 1,2,3,4-Tetrachlorobenzene	13.18	216	54076	3.23	ng/uL	93
22) Pentachlorobenzene	14.70	250	55633	3.18	ng/uL	94

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

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