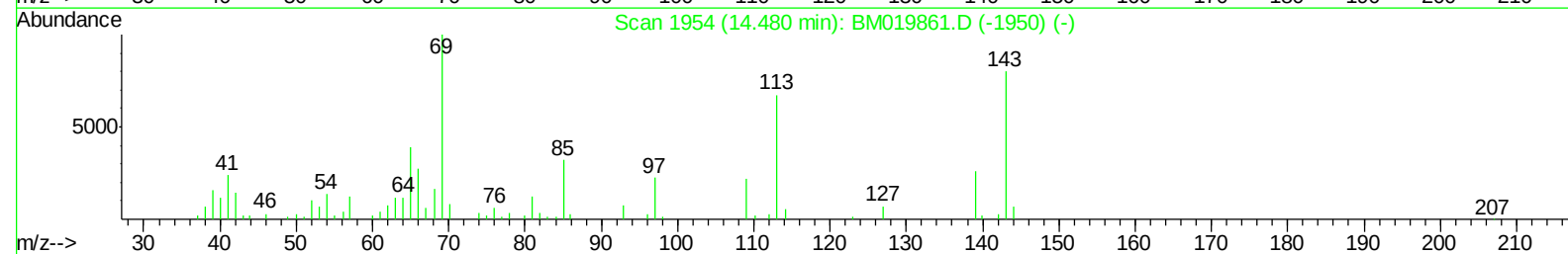
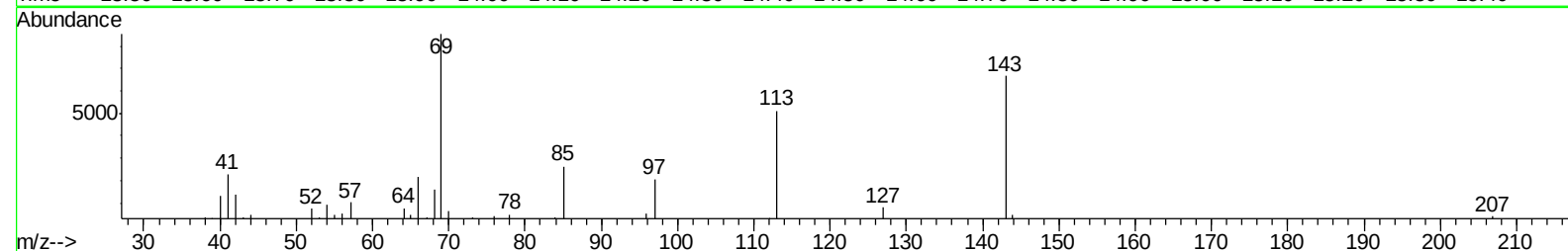
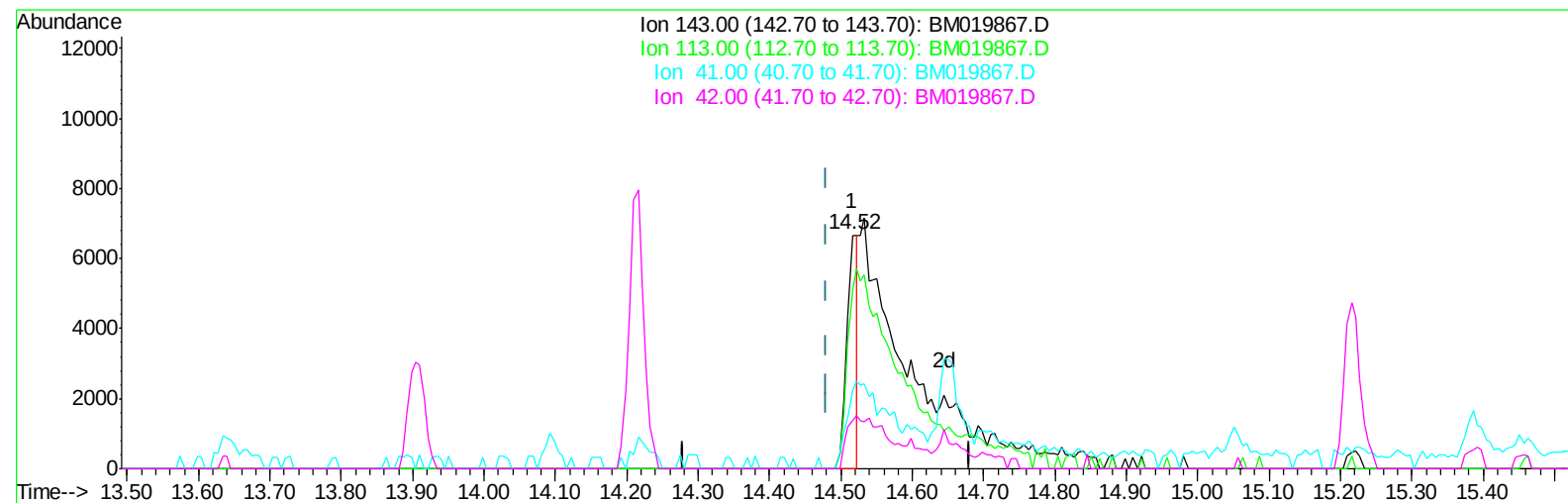


Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
Data File : BM019867.D
Acq On : 10 Apr 2019 17:13
Operator : JU/SJ
Sample : K2303-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Apr 11 03:23:10 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 11 03:09:11 2019
Response via : Initial Calibration



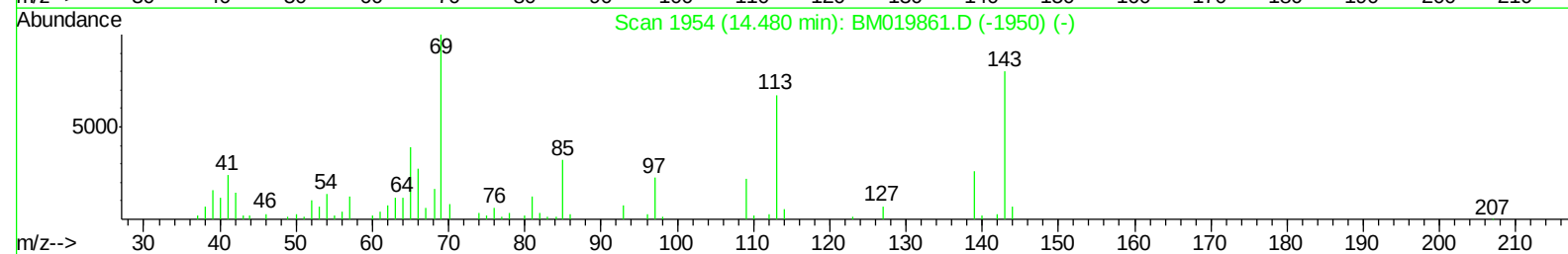
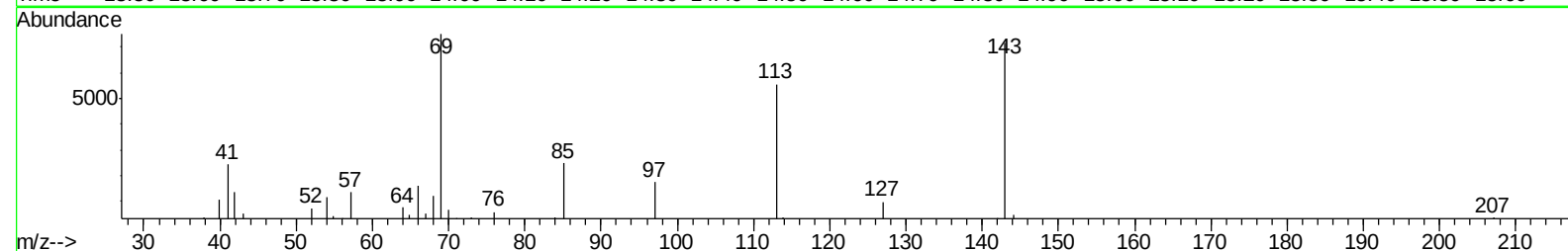
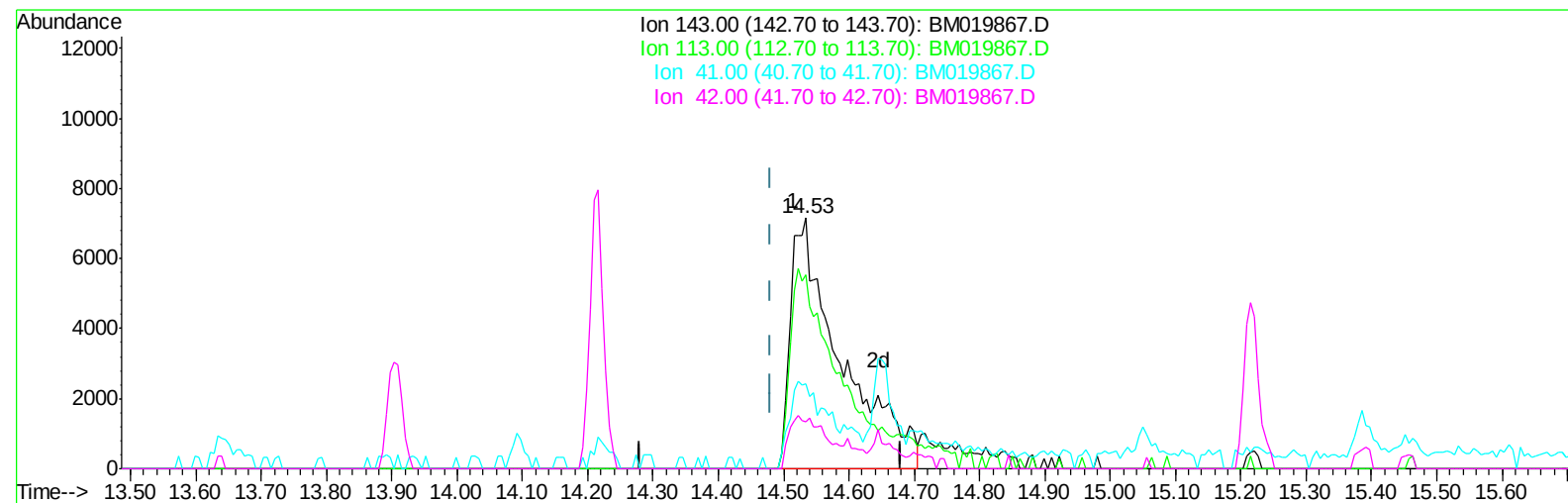
TIC: BM019867.D

(52) 4-Nitrophenol-d4 (S)
14.516min (+0.035) 1.32ng/ul
response 7084

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	76.55
41.00	32.60	33.83
42.00	22.30	20.42

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
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Operator : JU/SJ
Sample : K2303-16
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Apr 11 03:23:10 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 11 03:09:11 2019
Response via : Initial Calibration



TIC: BM019867.D

(52) 4-Nitrophenol-d4 (S)

14.533min (+0.053) 7.10ng/ul m

response 38117

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	77.57
41.00	32.60	34.11
42.00	22.30	18.76

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
 Data File : BM019867.D
 Acq On : 10 Apr 2019 17:13
 Operator : JU/SJ
 Sample : K2303-16
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Apr 11 03:58:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 11 03:09:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	190753	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	798433	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.22	164	424359	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.97	188	849134	20.00	ng/ul	0.00
78) Chrysene-d12	21.19	240	806481	20.00	ng/ul	0.00
86) Perylene-d12	23.39	264	933910	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.12	96	20899	4.29	ng/uL	0.00
5) Phenol-d5	6.75	99	167054	9.83	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.90	67	117277	10.31	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	143175	11.12	ng/ul	0.00
13) 4-Methylphenol-d8	8.28	113	105735	8.29	ng/ul	0.00
19) Nitrobenzene-d5	8.73	128	65231	11.45	ng/ul	0.00
22) 2-Nitrophenol-d4	9.44	143	71845	11.33	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.98	165	113192	9.46	ng/ul	0.01
29) 4-Chloroaniline-d4	10.51	131	146956	10.23	ng/ul	0.01
44) Dimethylphthalate-d6	13.64	166	398111	11.63	ng/ul	0.00
47) Acenaphthylene-d8	13.90	160	460433	10.75	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	38117m	7.10	ng/ul	0.05
58) Fluorene-d10	15.22	176	323497	10.97	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	34201	6.09	ng/ul	0.00
71) Anthracene-d10	17.07	188	431413	10.59	ng/ul	0.00
79) Pyrene-d10	19.39	212	437421	11.74	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.25	264	499223	10.08	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
45) Dimethylphthalate	13.69	163	75610	2.201	ng/ul	99

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

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