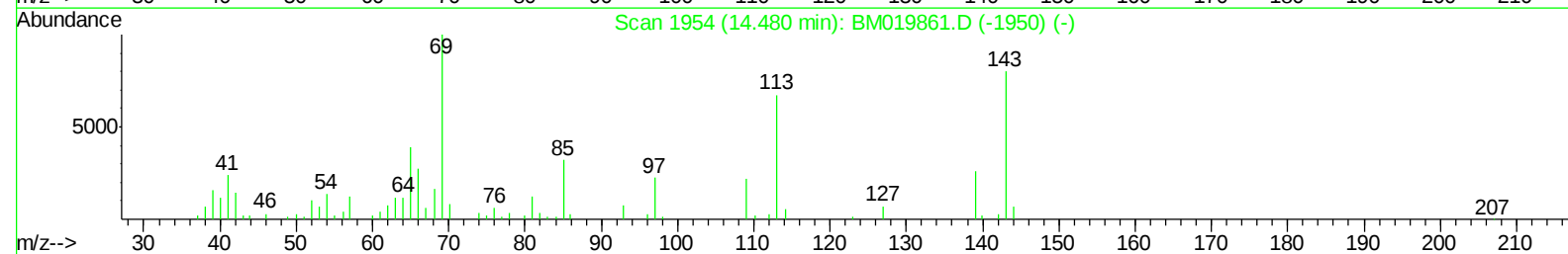
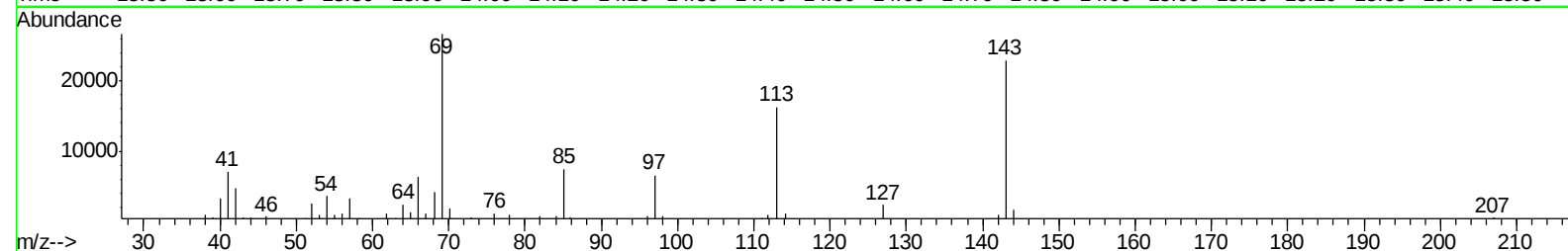
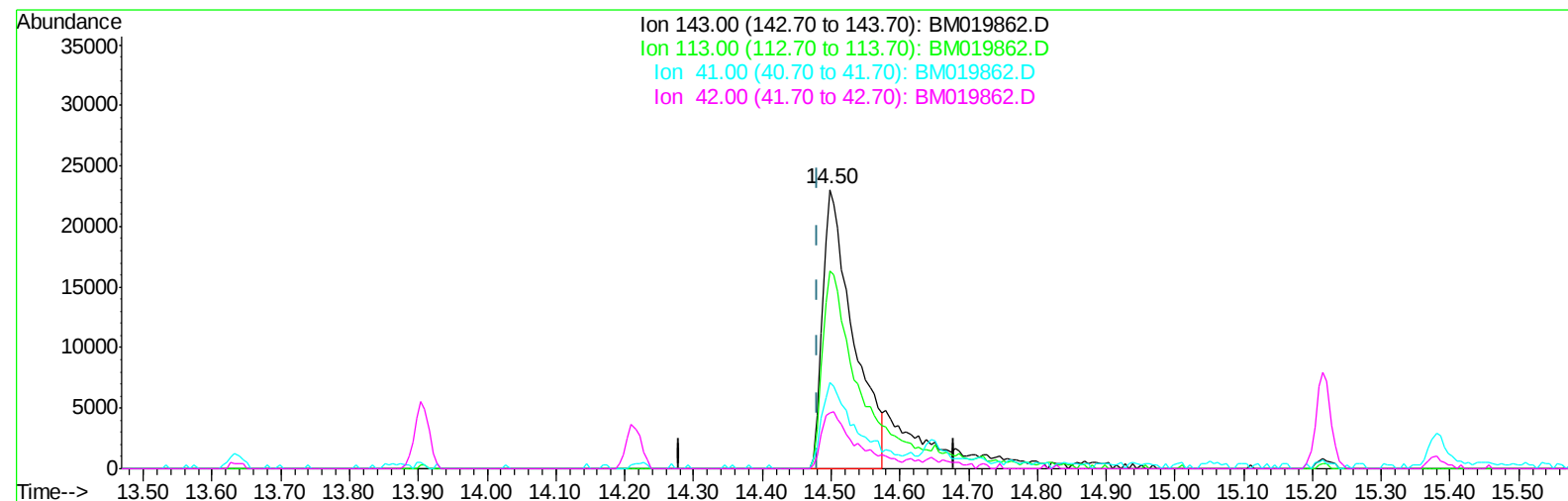


Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
Data File : BM019862.D
Acq On : 10 Apr 2019 14:08
Operator : JU/SJ
Sample : K2218-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 11 03:22:46 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



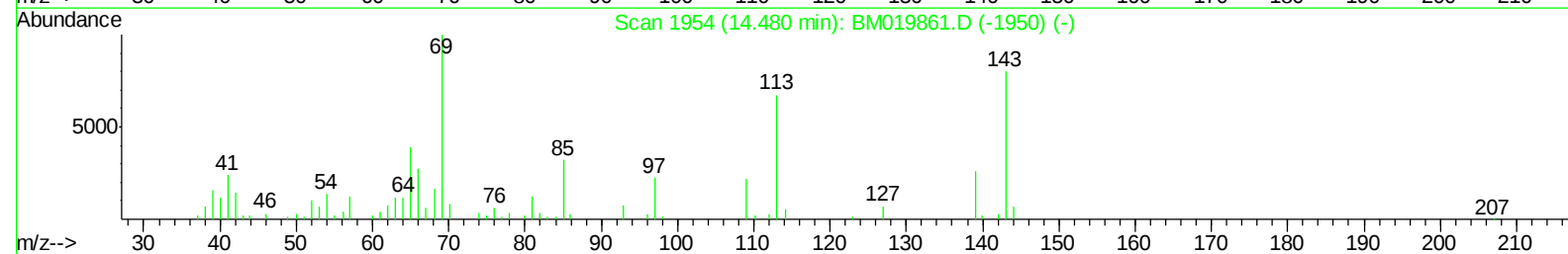
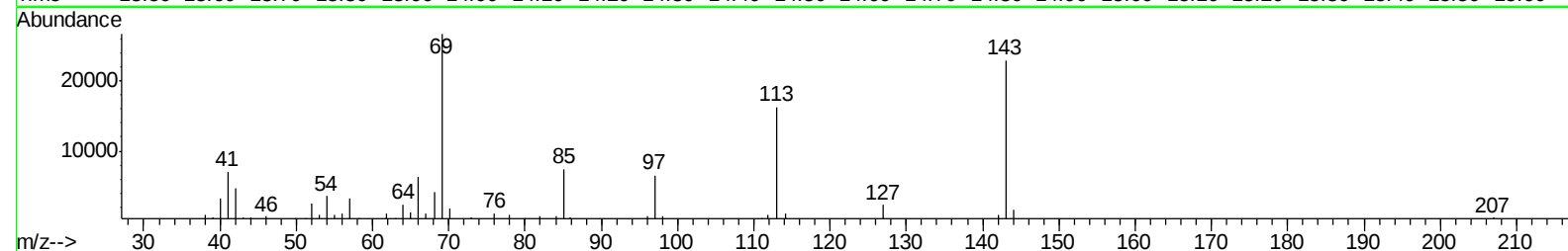
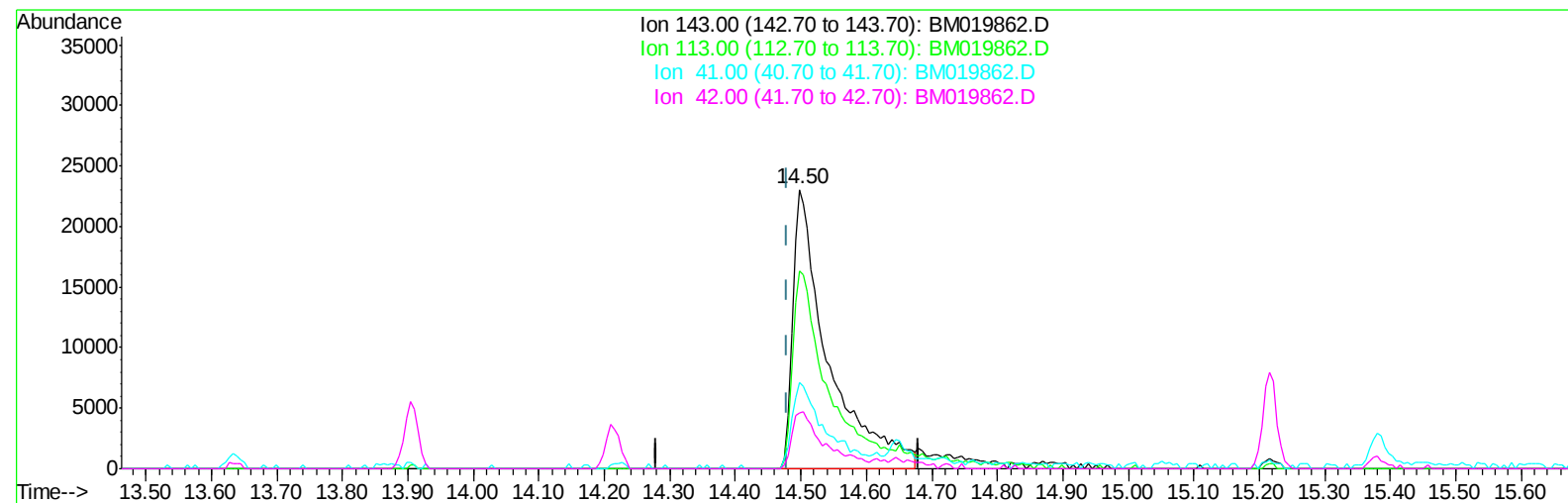
TIC: BM019862.D

(52) 4-Nitrophenol-d4 (S)
14.498min (+0.017) 28.24ng/ul
response 70908

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	70.92
41.00	32.60	30.79
42.00	22.30	20.23

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
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Acq On : 10 Apr 2019 14:08
Operator : JU/SJ
Sample : K2218-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 11 03:22:46 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: BM019862.D

(52) 4-Nitrophenol-d4 (S)

14.498min (+0.017) 34.51ng/ul m

response 86647

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	70.92
41.00	32.60	30.79
42.00	22.30	20.23

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
 Data File : BM019862.D
 Acq On : 10 Apr 2019 14:08
 Operator : JU/SJ
 Sample : K2218-05
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 11 03:49:59 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	89326	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	372934	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.22	164	198491	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.97	188	433000	20.00	ng/ul	0.00
78) Chrysene-d12	21.19	240	429702	20.00	ng/ul	0.00
86) Perylene-d12	23.39	264	457231	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.11	96	23859	10.46	ng/uL	0.00
5) Phenol-d5	6.75	99	281820	35.43	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.90	67	177865	33.40	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	225656	37.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	218346	36.57	ng/ul	0.00
19) Nitrobenzene-d5	8.73	128	104095	39.13	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	122844	41.49	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.97	165	203604	36.45	ng/ul	0.00
29) 4-Chloroaniline-d4	10.50	131	289017	43.08	ng/ul	0.00
44) Dimethylphthalate-d6	13.64	166	656192	40.97	ng/ul	0.00
47) Acenaphthylene-d8	13.90	160	825374	41.21	ng/ul	0.00
52) 4-Nitrophenol-d4	14.50	143	86647m	34.51	ng/ul	0.02
58) Fluorene-d10	15.22	176	560191	40.60	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.38	200	74728	26.10	ng/ul	0.00
71) Anthracene-d10	17.07	188	862643	41.52	ng/ul	0.00
79) Pyrene-d10	19.39	212	918586	46.28	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.26	264	952527	39.30	ng/ul	0.00

Target Compounds				Qvalue		
6) Phenol	6.77	94	27259	3.284	ng/ul#	89
45) Dimethylphthalate	13.69	163	92125	5.733	ng/ul	99

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

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