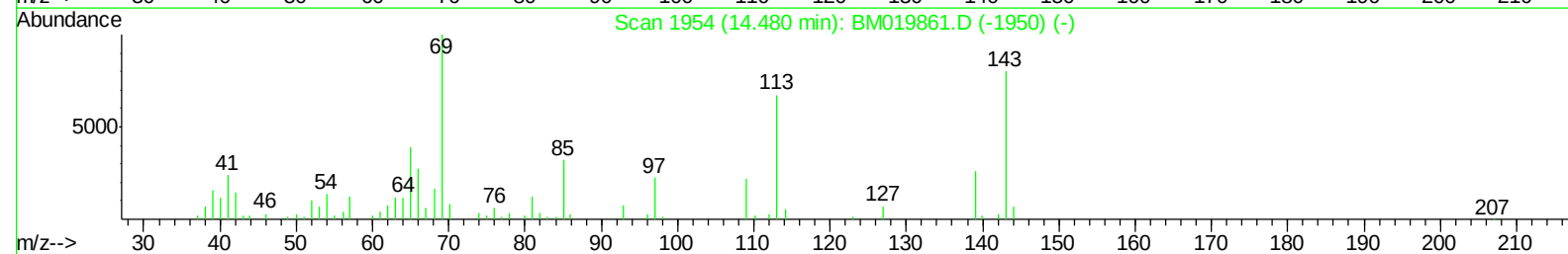
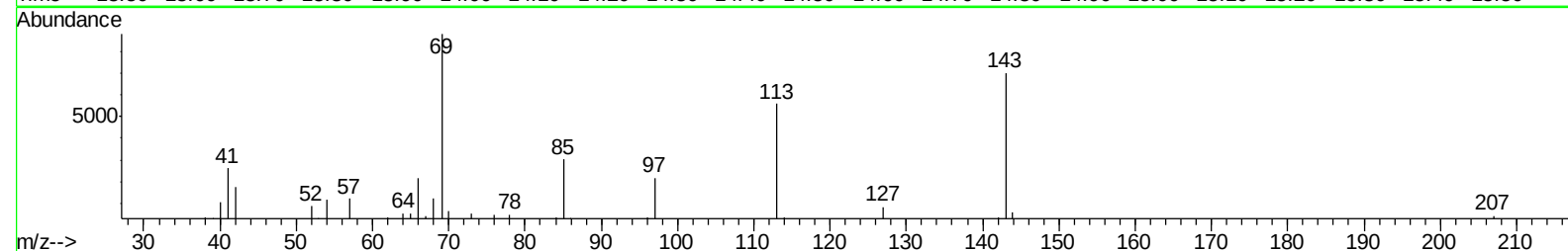
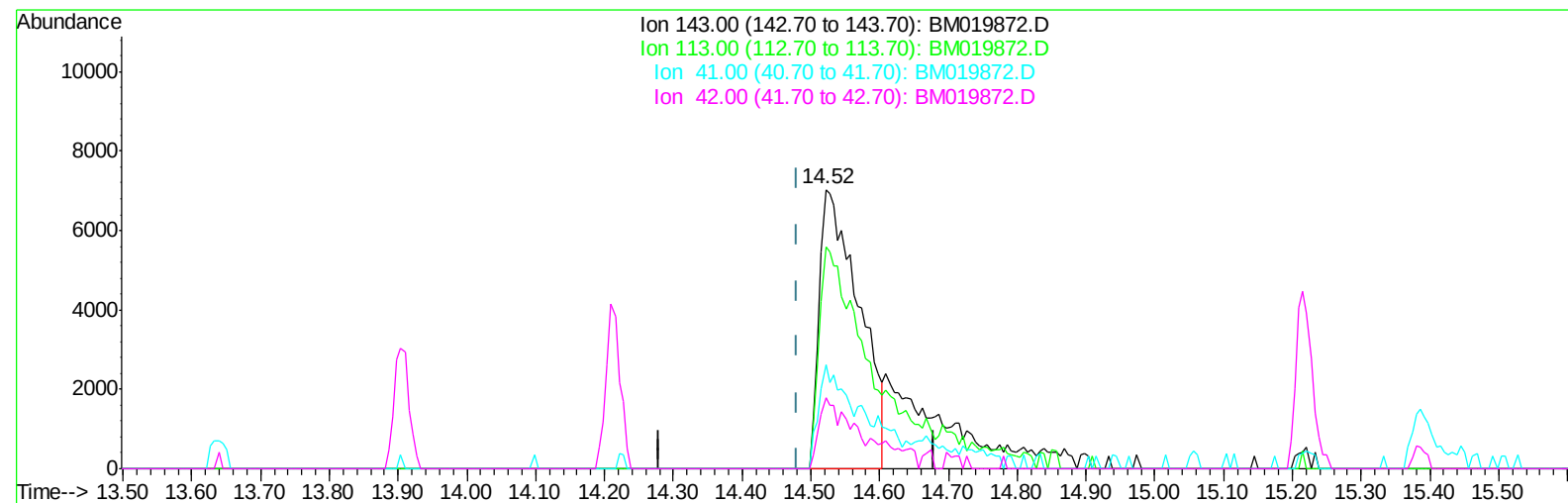


Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
Data File : BM019872.D
Acq On : 10 Apr 2019 20:18
Operator : JU/SJ
Sample : K2303-14
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Apr 11 03:23:40 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



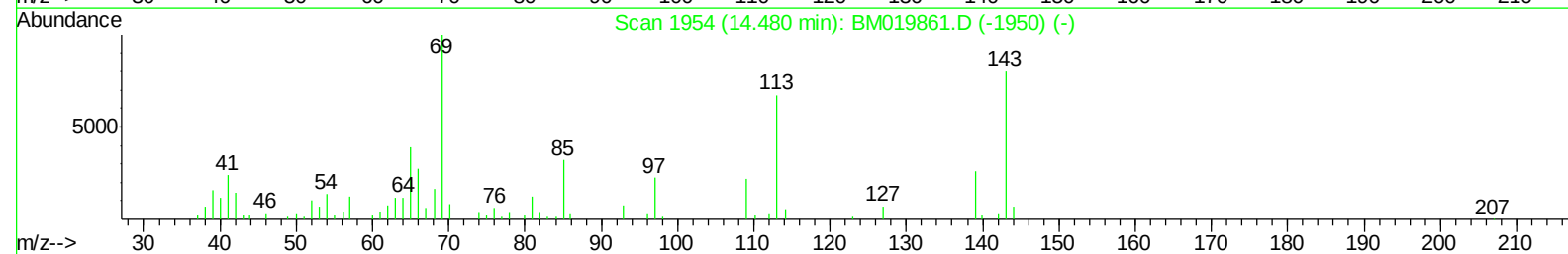
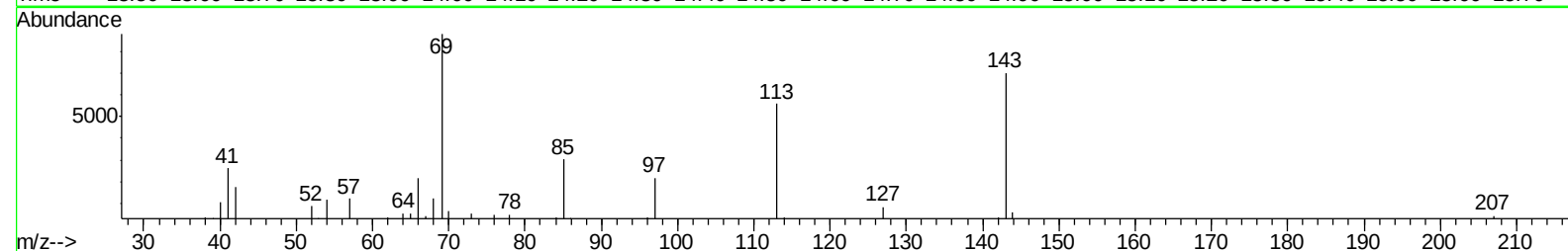
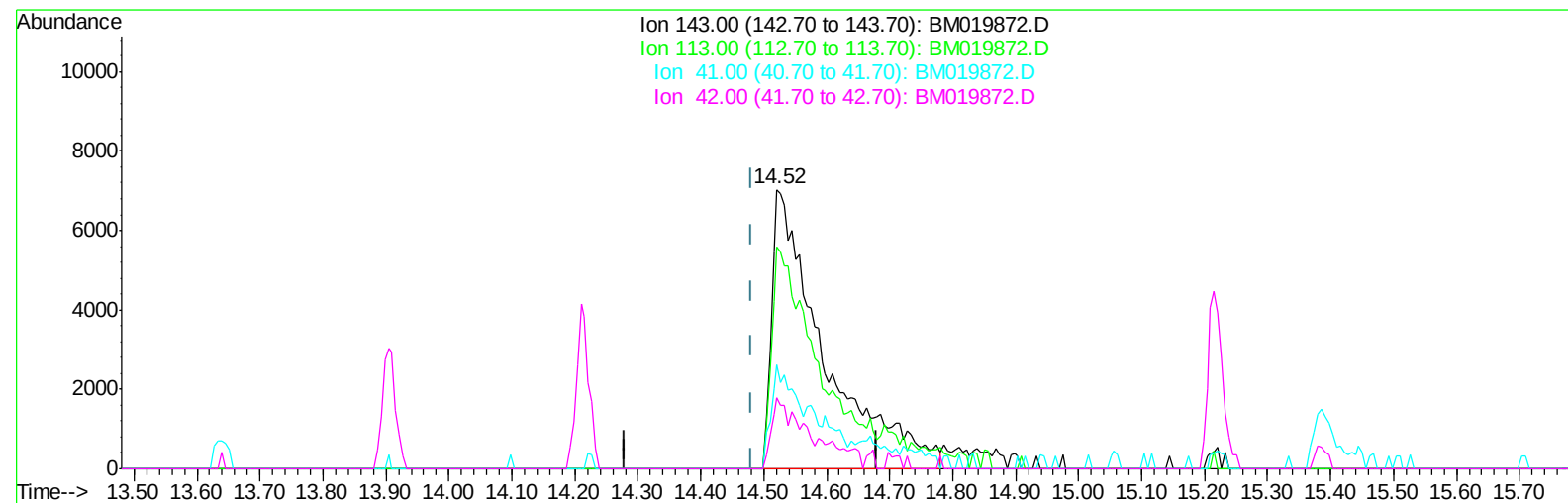
TIC: BM019872.D

(52) 4-Nitrophenol-d4 (S)
14.521min (+0.041) 11.10ng/ul
response 28104

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	79.35
41.00	32.60	37.23
42.00	22.30	25.32

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
Data File : BM019872.D
Acq On : 10 Apr 2019 20:18
Operator : JU/SJ
Sample : K2303-14
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Apr 11 03:23:40 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: BM019872.D

(52) 4-Nitrophenol-d4 (S)

14.521min (+0.041) 16.07ng/ul m

response 40676

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	79.35
41.00	32.60	37.23
42.00	22.30	25.32

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
 Data File : BM019872.D
 Acq On : 10 Apr 2019 20:18
 Operator : JU/SJ
 Sample : K2303-14
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Apr 11 04:07:51 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	105880	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	409712	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.22	164	200178	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.97	188	397253	20.00	ng/ul	0.00
78) Chrysene-d12	21.19	240	433049	20.00	ng/ul	0.00
86) Perylene-d12	23.39	264	521880	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.12	96	21762	8.05	ng/uL	0.00
5) Phenol-d5	6.75	99	188337	19.97	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.91	67	134594	21.32	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	164070	22.96	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	133441	18.85	ng/ul	0.00
19) Nitrobenzene-d5	8.73	128	73082	25.01	ng/ul	0.00
22) 2-Nitrophenol-d4	9.44	143	79385	24.40	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.98	165	114229	18.61	ng/ul	0.01
29) 4-Chloroaniline-d4	10.51	131	166096	22.53	ng/ul	0.01
44) Dimethylphthalate-d6	13.64	166	409601	25.36	ng/ul	0.00
47) Acenaphthylene-d8	13.90	160	474273	23.48	ng/ul	0.00
52) 4-Nitrophenol-d4	14.52	143	40676m	16.07	ng/ul	0.04
58) Fluorene-d10	15.22	176	324693	23.33	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	37019	14.09	ng/ul	0.00
71) Anthracene-d10	17.07	188	450912	23.65	ng/ul	0.00
79) Pyrene-d10	19.39	212	480186	24.01	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.25	264	603238	21.80	ng/ul	0.00

Target Compounds					Qvalue
45) Dimethylphthalate	13.69	163	58555	3.613	ng/ul 99

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

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