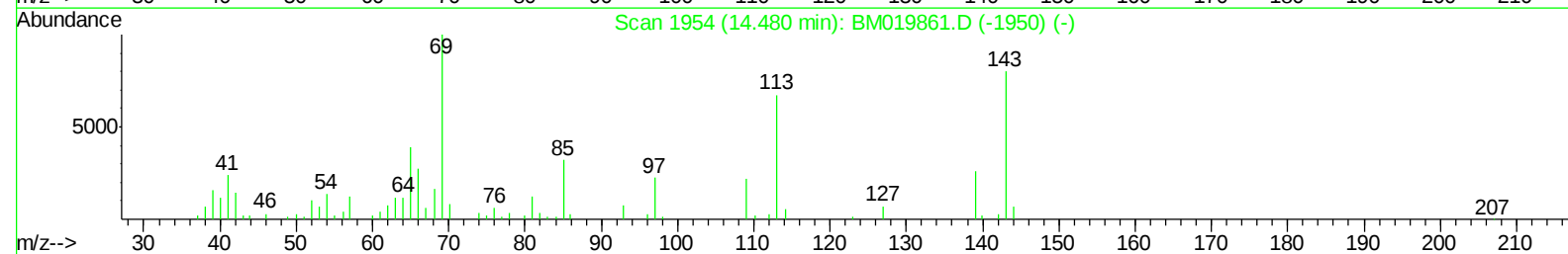
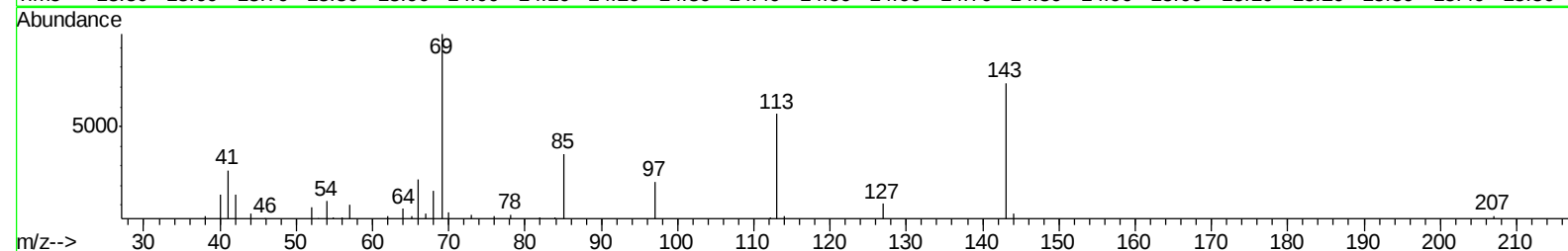
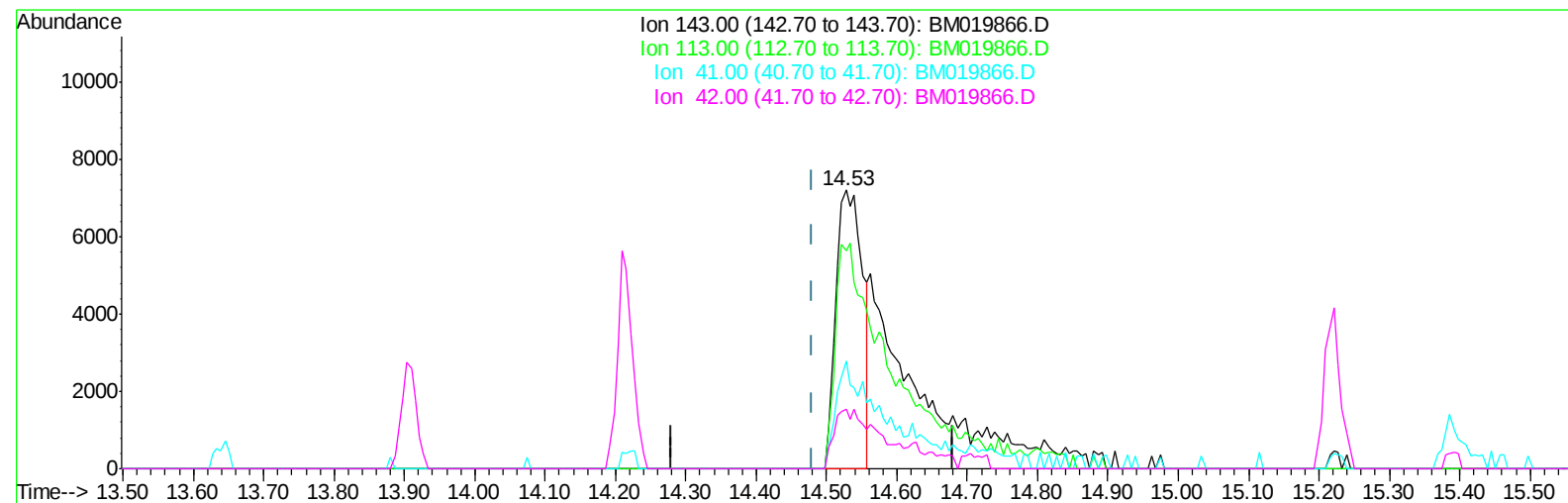


Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
Data File : BM019866.D
Acq On : 10 Apr 2019 16:35
Operator : JU/SJ
Sample : PB118673BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Apr 11 03:23:06 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



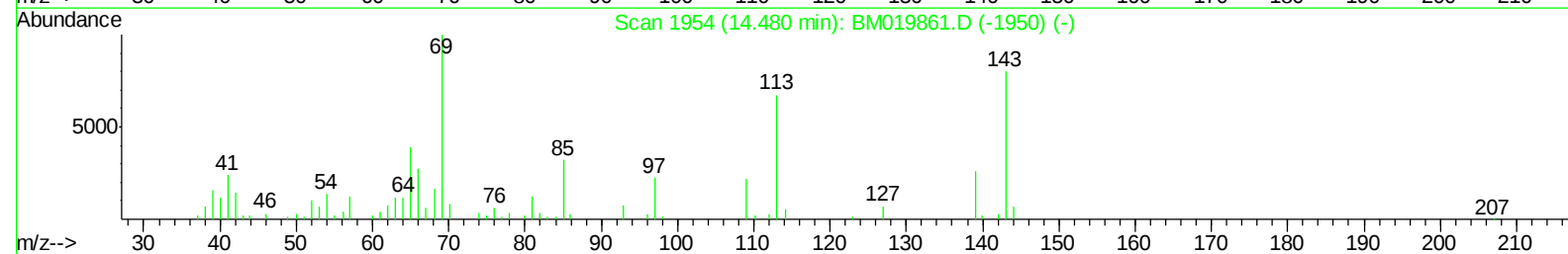
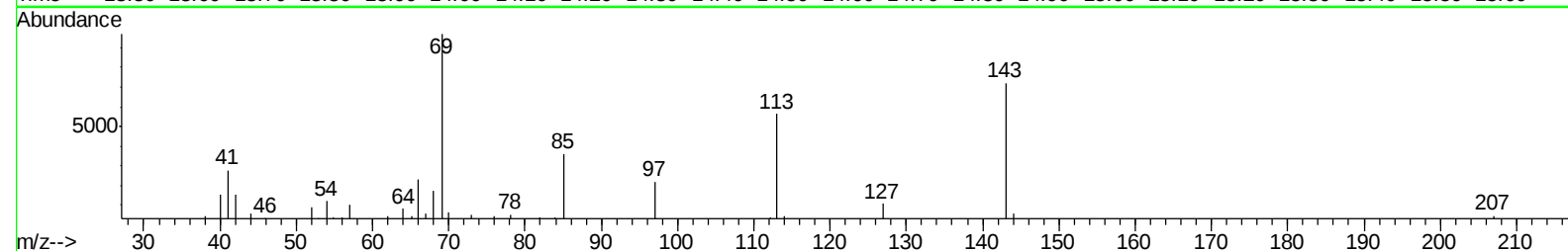
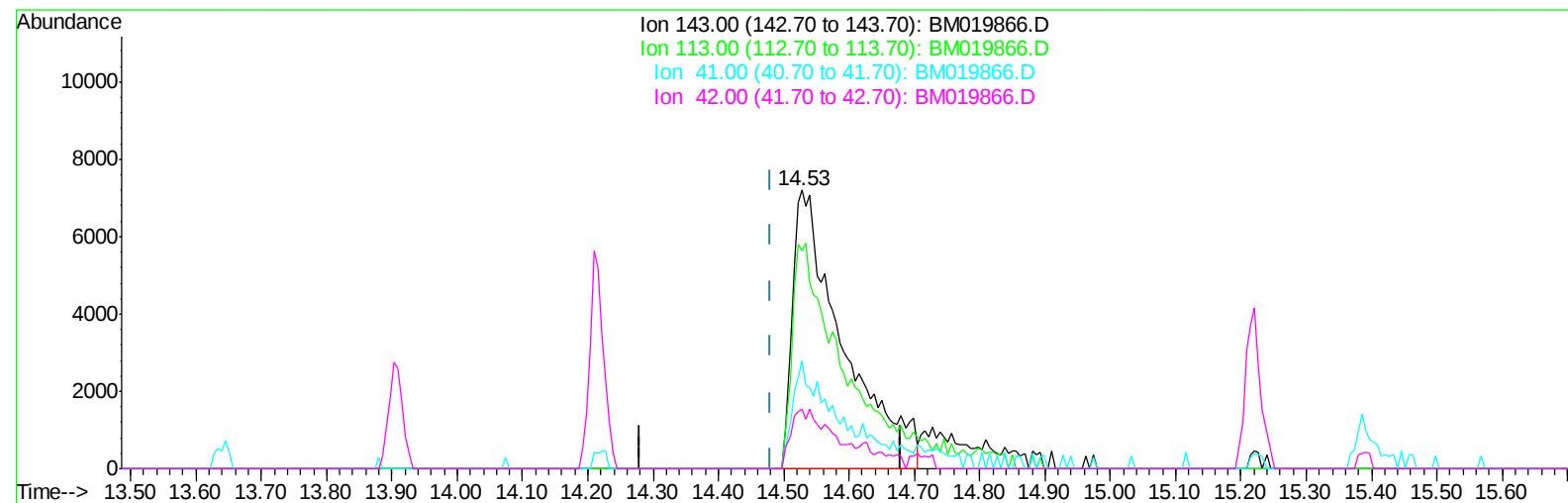
TIC: BM019866.D

(52) 4-Nitrophenol-d4 (S)
14.527min (+0.047) 4.77ng/ul
response 18807

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	78.06
41.00	32.60	38.57
42.00	22.30	21.58

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
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Acq On : 10 Apr 2019 16:35
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Sample : PB118673BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Apr 11 03:23:06 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: BM019866.D

(52) 4-Nitrophenol-d4 (S)

14.527min (+0.047) 9.77ng/ul m

response 38472

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	78.06
41.00	32.60	38.57
42.00	22.30	21.58

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041019\
 Data File : BM019866.D
 Acq On : 10 Apr 2019 16:35
 Operator : JU/SJ
 Sample : PB118673BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Apr 11 03:56:41 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	151772	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	597605	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.22	164	311443	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.97	188	603618	20.00	ng/ul	0.00
78) Chrysene-d12	21.19	240	659091	20.00	ng/ul	0.00
86) Perylene-d12	23.40	264	763455	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	12826	3.31	ng/uL	0.01
5) Phenol-d5	6.75	99	179086	13.25	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.91	67	116336	12.86	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	153724	15.01	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	133610	13.17	ng/ul	0.00
19) Nitrobenzene-d5	8.73	128	64541	15.14	ng/ul	0.00
22) 2-Nitrophenol-d4	9.44	143	74319	15.66	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.98	165	120852	13.50	ng/ul	0.01
29) 4-Chloroaniline-d4	10.50	131	171385	15.94	ng/ul	0.00
44) Dimethylphthalate-d6	13.64	166	364217	14.49	ng/ul	0.00
47) Acenaphthylene-d8	13.91	160	455597	14.50	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	38472m	9.77	ng/ul	0.05
58) Fluorene-d10	15.22	176	304072	14.04	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	29370	7.36	ng/ul	0.01
71) Anthracene-d10	17.08	188	443702	15.32	ng/ul	0.00
79) Pyrene-d10	19.39	212	468765	15.40	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.25	264	588818	14.55	ng/ul	0.00

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

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