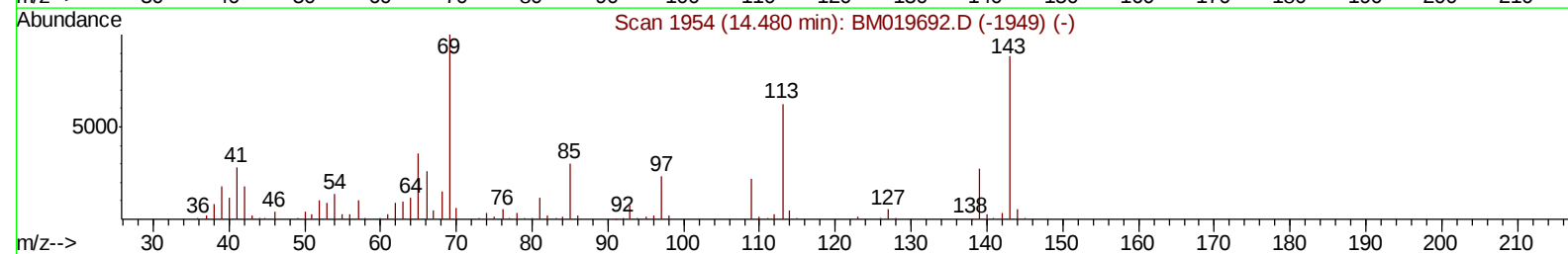
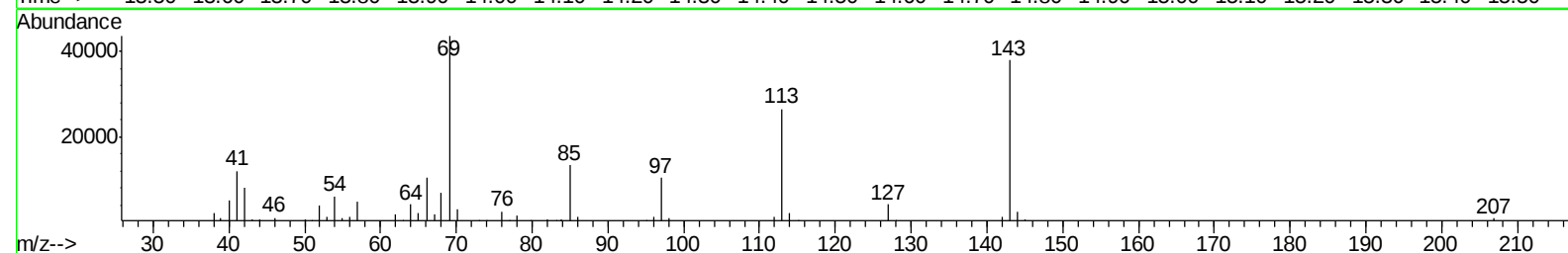
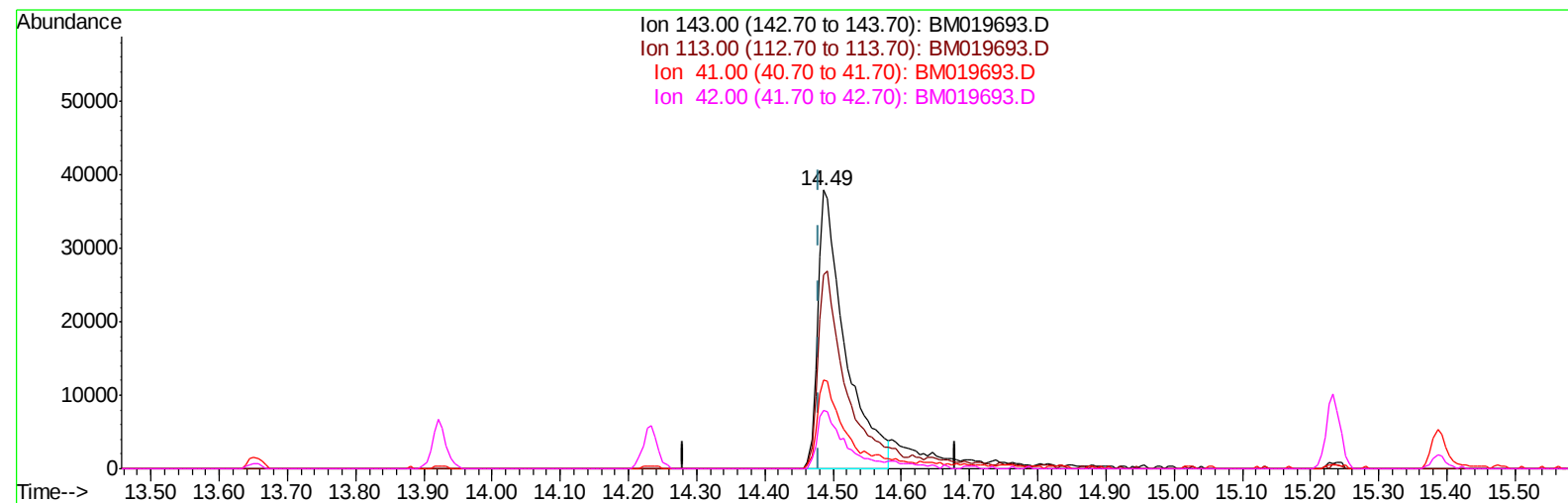


Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM040219\
Data File : BM019693.D
Acq On : 02 Apr 2019 16:11
Operator : JU/SJ
Sample : PB118447BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 03 01:21:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 03 01:13:19 2019
Response via : Initial Calibration



TIC: BM019693.D

(52) 4-Nitrophenol-d4 (S)

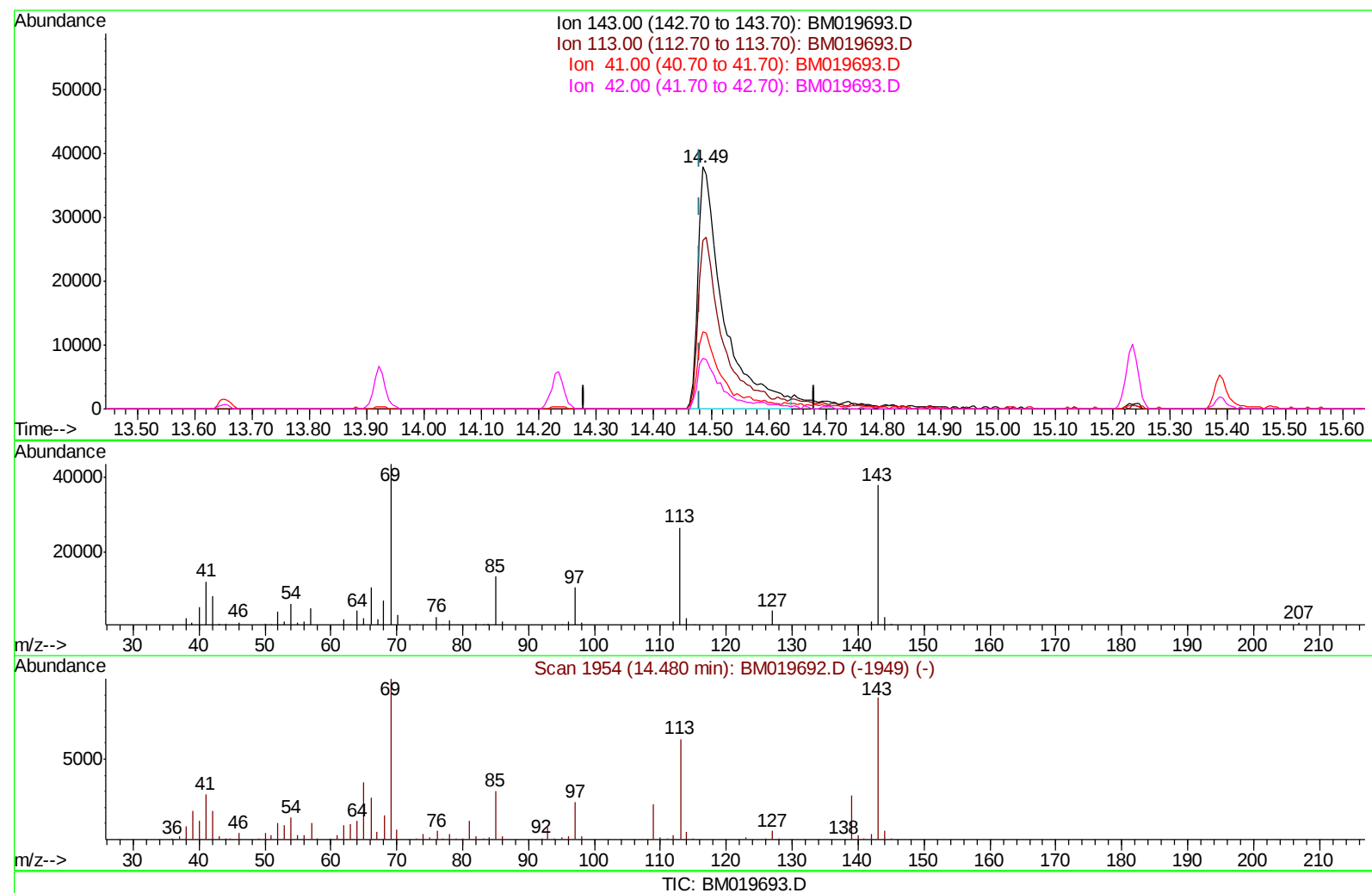
14.486min (+0.006) 26.30ng/ul

response 105437

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	69.59
41.00	32.60	31.65
42.00	22.30	21.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM040219\
Data File : BM019693.D
Acq On : 02 Apr 2019 16:11
Operator : JU/SJ
Sample : PB118447BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 03 01:21:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 03 01:13:19 2019
Response via : Initial Calibration



(52) 4-Nitrophenol-d4 (S)

14.486min (+0.006) 28.66ng/ul m

response 114930

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	69.59
41.00	32.60	31.65
42.00	22.30	21.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM040219\
 Data File : BM019693.D
 Acq On : 02 Apr 2019 16:11
 Operator : JU/SJ
 Sample : PB118447BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Apr 03 02:54:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 03 01:13:19 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	112629	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.35	136	520380	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.23	164	317021	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.99	188	721955	20.00	ng/ul	0.00
78) Chrysene-d12	21.20	240	806053	20.00	ng/ul	0.00
86) Perylene-d12	23.41	264	921658	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	14253	4.95	ng/uL	0.00
5) Phenol-d5	6.76	99	253001	25.22	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.92	67	165884	24.70	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.10	132	205788	27.07	ng/ul	0.00
13) 4-Methylphenol-d8	8.29	113	216859	28.80	ng/ul	0.00
19) Nitrobenzene-d5	8.75	128	101609	27.37	ng/ul	0.00
22) 2-Nitrophenol-d4	9.46	143	121276	29.35	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.99	165	186627	23.94	ng/ul	0.00
29) 4-Chloroaniline-d4	10.52	131	304483	32.52	ng/ul	0.00
44) Dimethylphthalate-d6	13.65	166	830304	32.46	ng/ul	-0.01
47) Acenaphthylene-d8	13.92	160	943398	29.49	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.49	143	114930m	28.66	ng/ul	0.00
58) Fluorene-d10	15.23	176	660917	29.99	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	114320	23.95	ng/ul	0.00
71) Anthracene-d10	17.09	188	1084542	31.30	ng/ul	0.00
79) Pyrene-d10	19.40	212	1205666	32.38	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.27	264	1457032	29.82	ng/ul	0.00

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

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