

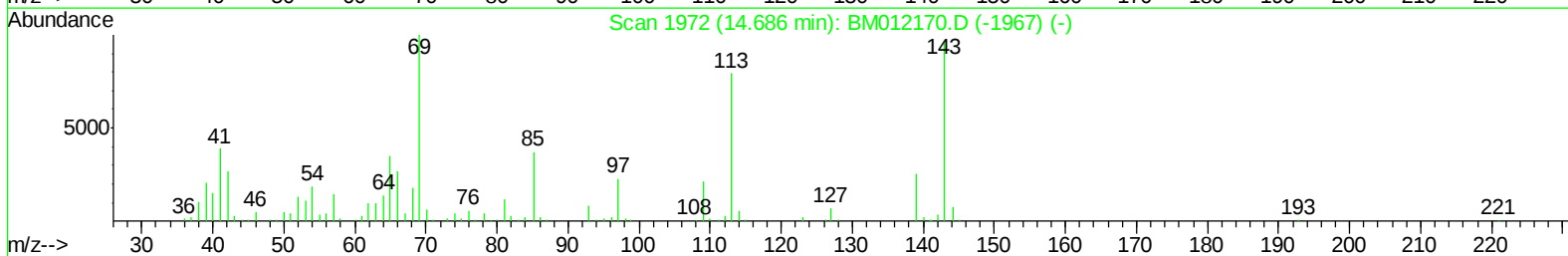
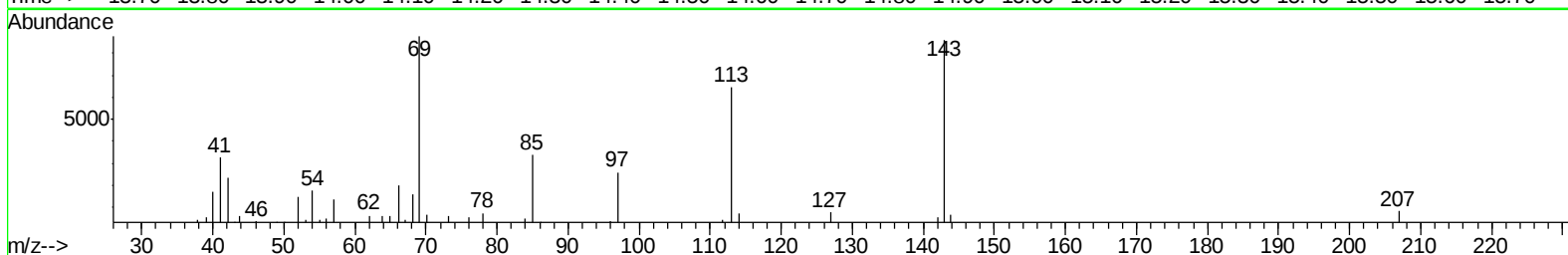
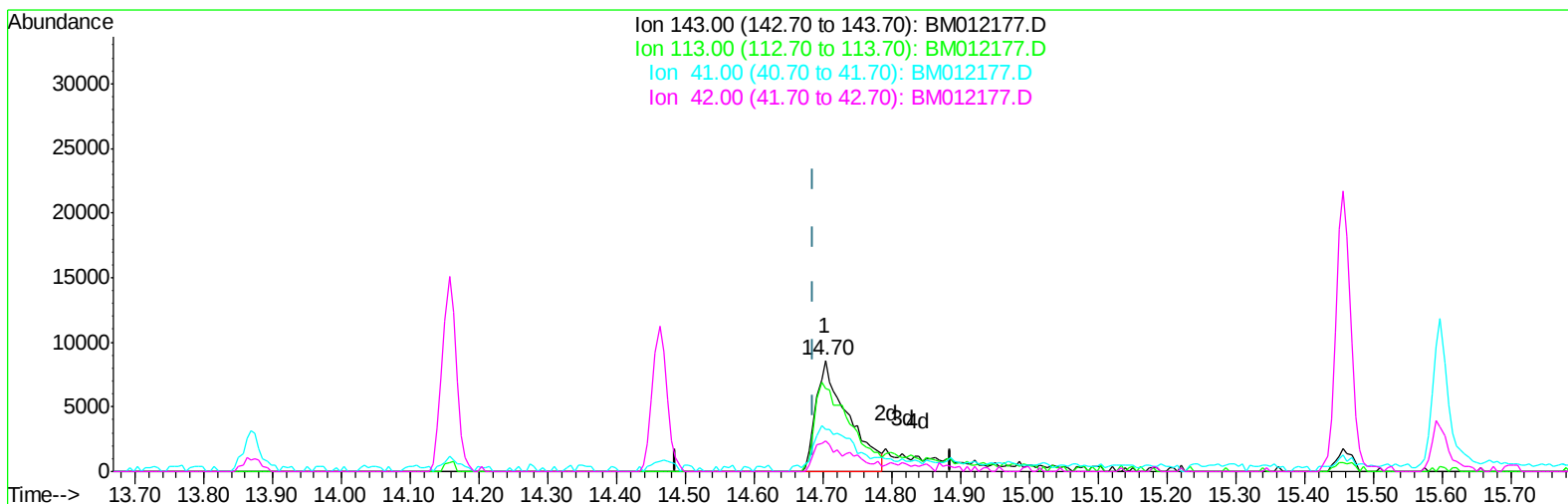
Data Path : Z:\HPCHEM1\BNA M\DATA\BM101417\
Data File : BM012177.D
Acq On : 14 Oct 2017 19:15
Operator : SJ/JU
Sample : I5840-11
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AE7

Quant Time: Oct 15 05:27:24 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Oct 15 05:22:07 2017
Response via : Initial Calibration

Manual Integrations
APPROVED

Sohil
10/16/2017 7:47:57 PM



TIC: BM012177.D

(51) 4-Nitrophenol-d4 (S)
14.704min (+0.018) 4.84ng/ul
response 27604

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	75.04#
41.00	34.80	37.86
42.00	20.50	27.38#

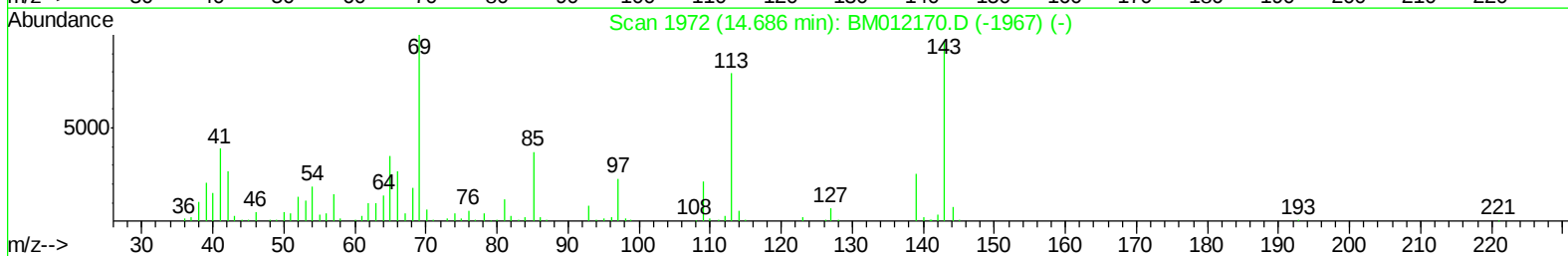
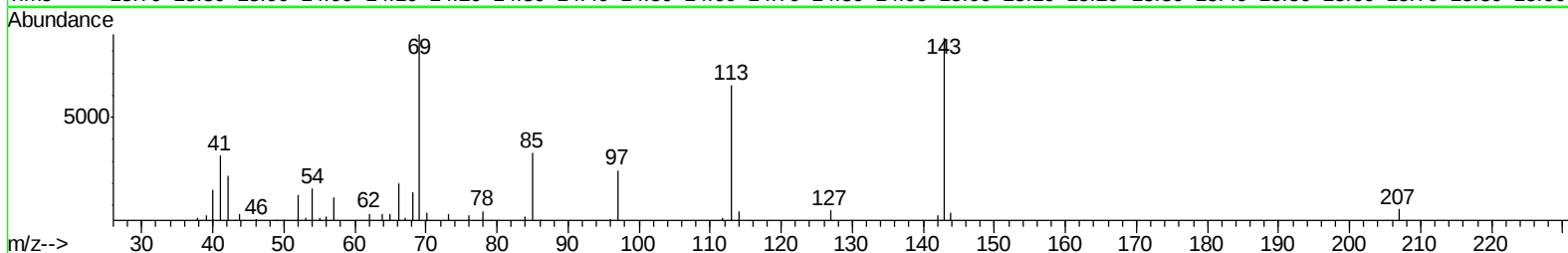
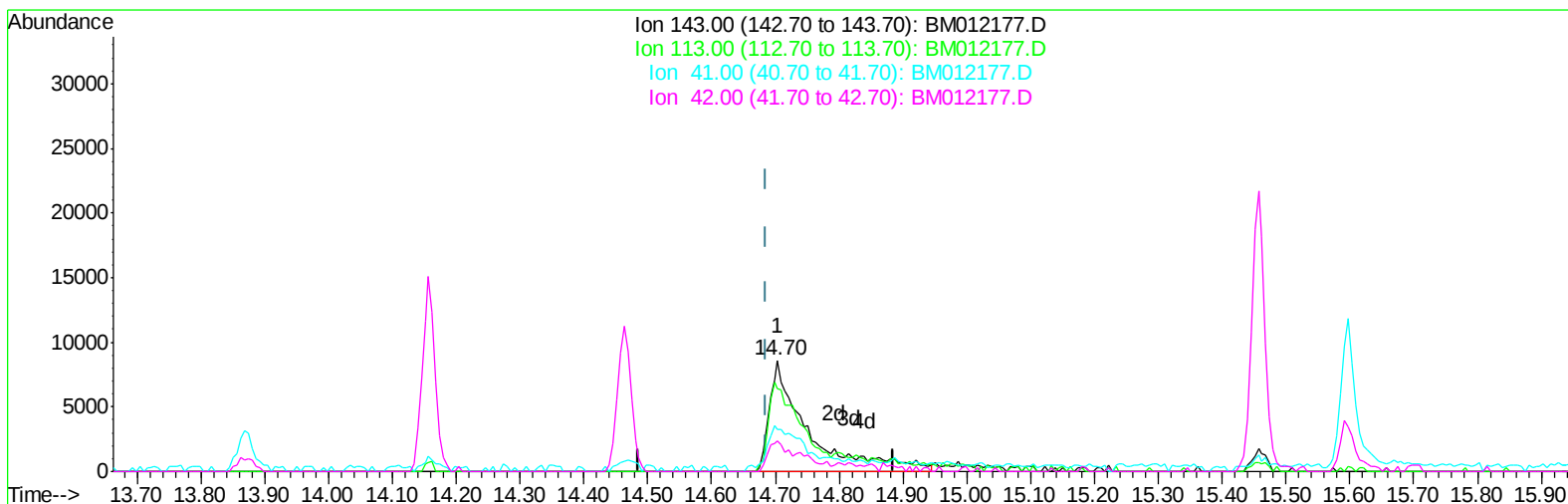
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TIC: BM012177.D

(51) 4-Nitrophenol-d4 (S)

14.704min (+0.018) 6.35ng/ul m

response 36213

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	75.04#
41.00	34.80	37.86
42.00	20.50	27.38#

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Manual Integrations
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Quant Time: Oct 15 05:53:35 2017
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 QLast Update : Sun Oct 15 05:22:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	151680	20.00	ng/ul	0.00
18) Naphthalene-d8	10.62	136	670554	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.46	164	429512	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.21	188	1046615	20.00	ng/ul	0.00
75) Chrysene-d12	21.39	240	1217503	20.00	ng/ul	0.00
83) Perylene-d12	23.70	264	1298136	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	19476	6.10	ng/uL	0.00
5) Phenol-d5	6.98	99	98616	8.01	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.14	67	244019	33.06	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	290132	27.79	ng/ul	0.00
13) 4-Methylphenol-d8	8.53	113	191410	18.06	ng/ul	0.00
19) Nitrobenzene-d5	8.98	128	176021	34.40	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	205773	34.41	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.25	165	345205	30.24	ng/ul	0.00
29) 4-Chloroaniline-d4	10.80	131	8075	0.76	ng/ul	0.04
43) Dimethylphthalate-d6	13.87	166	1434095	37.78	ng/ul	0.00
46) Acenaphthylene-d8	14.16	160	1633397	35.55	ng/ul	0.00
51) 4-Nitrophenol-d4	14.70	143	36213m	6.35	ng/ul	0.02
57) Fluorene-d10	15.46	176	1163757	37.35	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.60	200	225078	31.33	ng/ul	0.00
70) Anthracene-d10	17.31	188	1892111	36.56	ng/ul	0.00
76) Pyrene-d10	19.60	212	2286510	37.00	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.55	264	2332960	37.28	ng/ul	0.00

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

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