

Quantitation Report (Qedit)

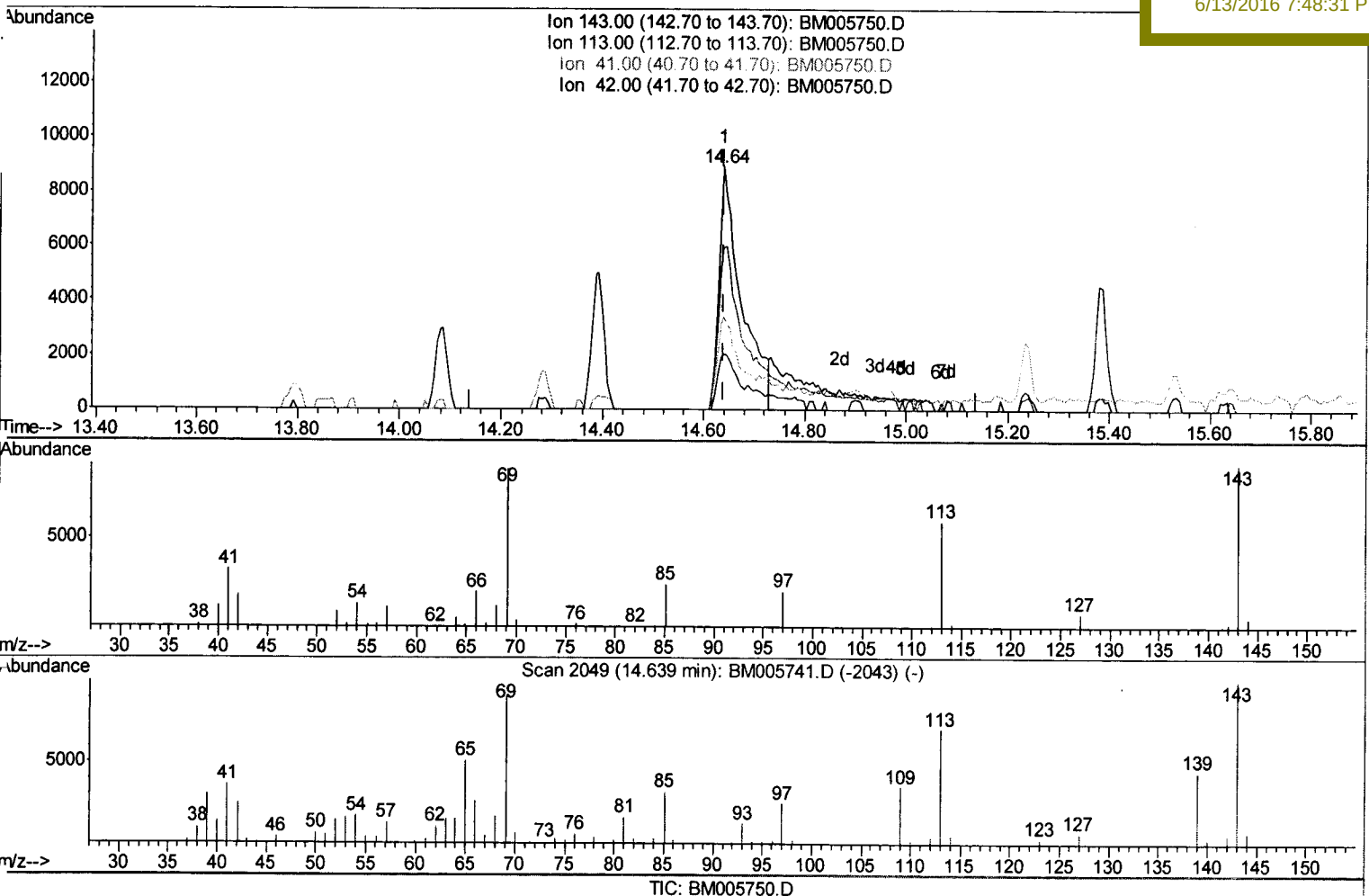
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061216\
Data File : BM005750.D
Acq On : 12 Jun 2016 15:26
Operator : UM/SJ
Sample : H3535-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Jun 13 03:30:53 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061016.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 13 03:29:45 2016
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
D9XW5

Manual Integrations
APPROVED

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6/13/2016 7:48:31 PM



(20) 4-Nitrophenol-d4 (S)

14.639min (-0.000) 7.86ng/ul

response 27404

Ion	Exp%	Act%
143.00	100	100
113.00	72.40	66.58
41.00	38.30	37.81
42.00	26.80	22.92

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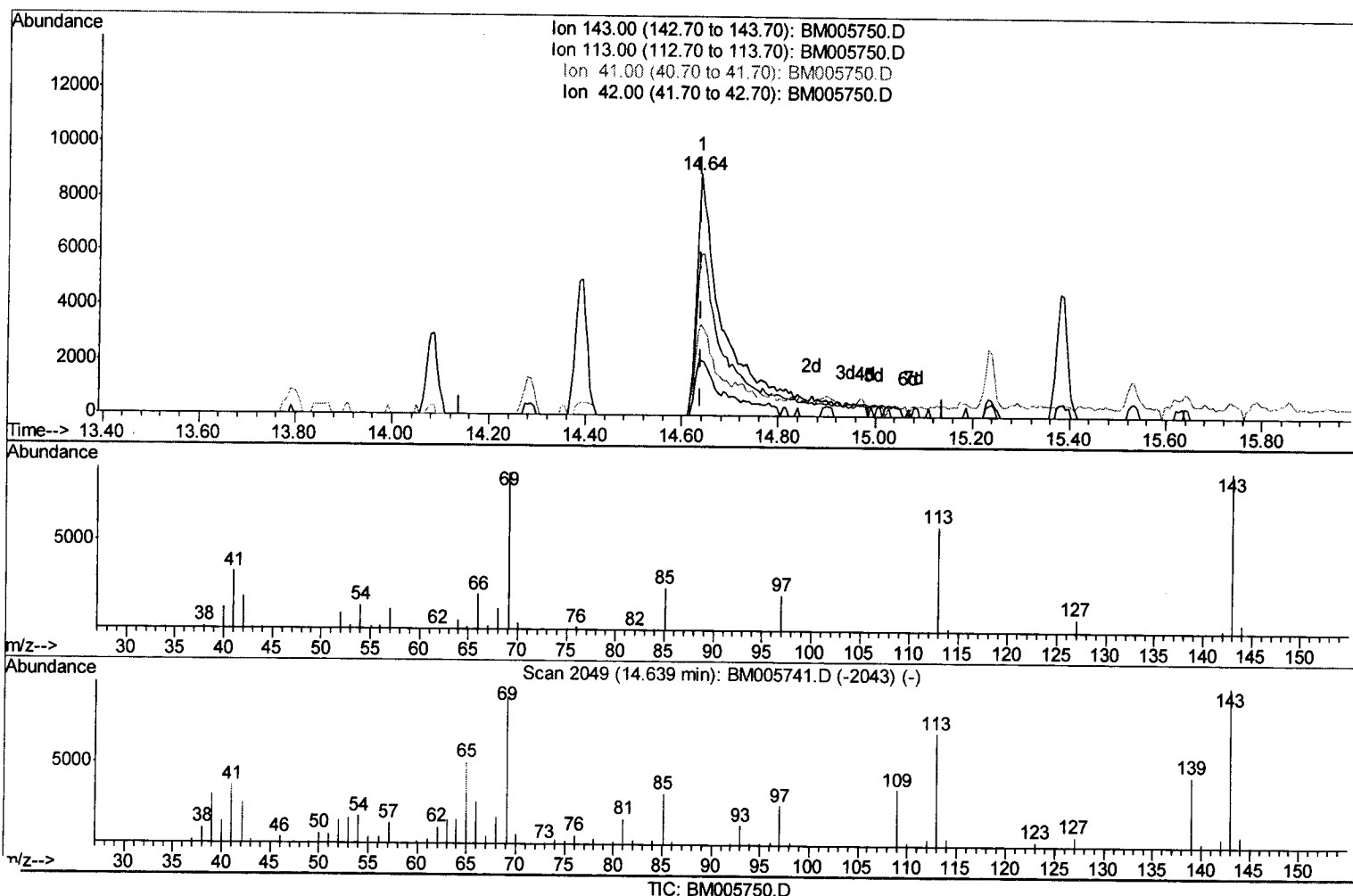
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(20) 4-Nitrophenol-d4 (S)

14.639min (-0.000) 11.43ng/ul m

response 39847

Ion	Exp%	Act%
143.00	100	100
113.00	72.40	66.58
41.00	38.30	37.81
42.00	26.80	22.92

U.M
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	86234	20.00	ng/ul	0.00
7) Naphthalene-d8	10.54	136	455907	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.39	164	288957	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	631144	20.00	ng/ul	0.00
29) Chrysene-d12	21.32	240	496630	20.00	ng/ul	0.00
34) Perylene-d12	23.58	264	436898	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.19	96	2456	1.24	ng/uL	0.00
3) Phenol-d5	6.93	99	89480	12.02	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.07	67	67113	15.01	ng/ul	0.00
5) 2-Chlorophenol-d4	7.28	132	85902	14.28	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	74231	11.81	ng/ul	0.00
8) Nitrobenzene-d5	8.90	128	45752	13.92	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	50083	13.48	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	88134	13.08	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	78342	9.64	ng/ul	0.00
16) Dimethylphthalate-d6	13.80	166	369172	15.49	ng/ul	0.00
17) Acenaphthylene-d8	14.08	160	465066	15.95	ng/ul	0.00
20) 4-Nitrophenol-d4	14.64	143	39847m	11.43	ng/ul	0.00
21) Fluorene-d10	15.39	176	321750	15.94	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	18401	5.64	ng/ul	-0.01
26) Anthracene-d10	17.23	188	475877	16.01	ng/ul	0.00
30) Pyrene-d10	19.53	212	463654	19.43	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.43	264	332240	16.67	ng/ul	-0.01

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Target Compounds Qvalue

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

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