

Quantitation Report (Qedit)

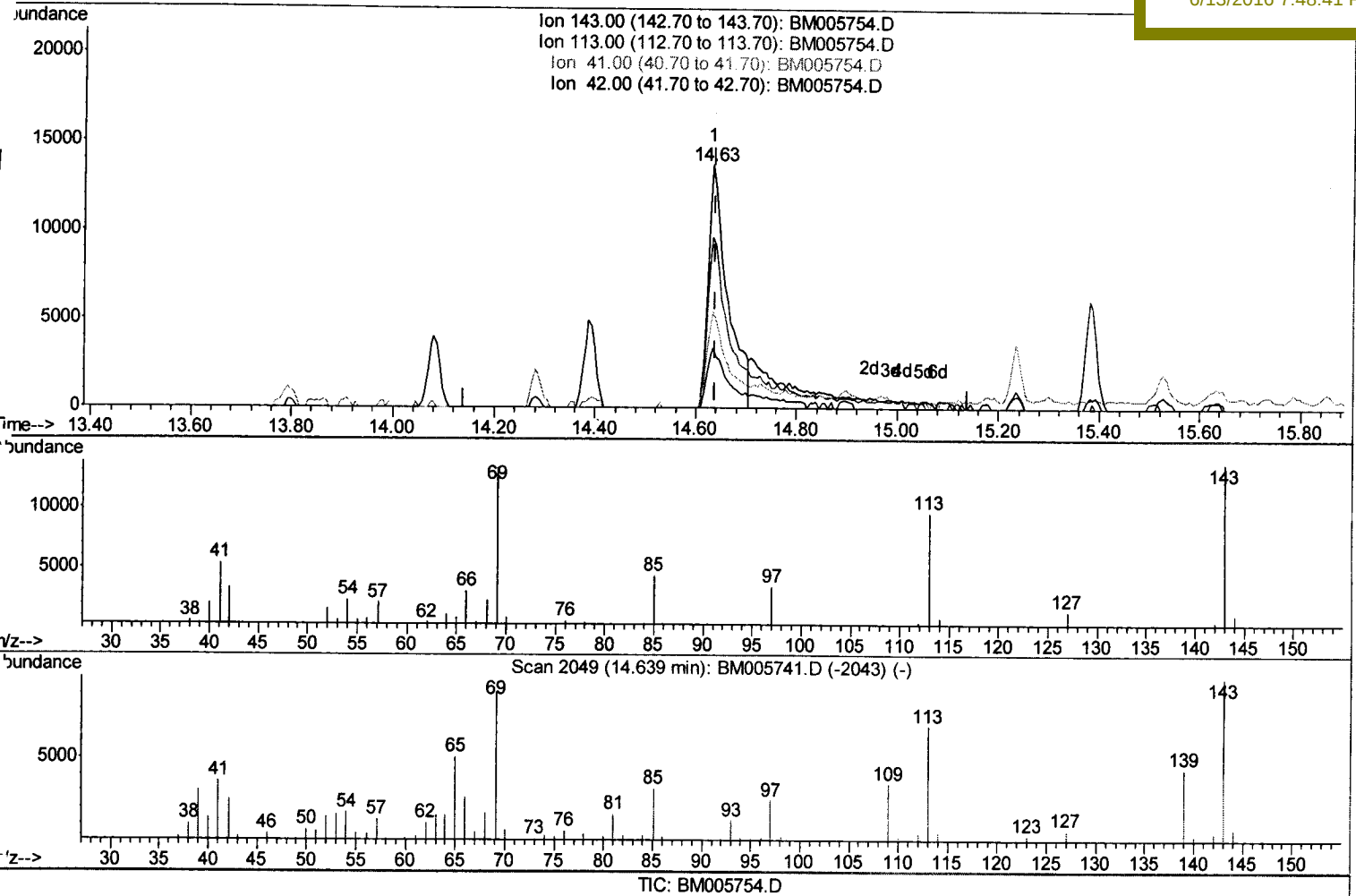
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061216\
 Data File : BM005754.D
 Acq On : 12 Jun 2016 17:51
 Operator : UM/SJ
 Sample : H3536-19
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Y29

Quant Time: Jun 13 03:31:16 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

Manual Integrations
 APPROVED

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(20) 4-Nitrophenol-d4 (S)

14.633min (-0.006) 10.89ng/ul

response 36572

Ion	Exp%	Act%
143.00	100	100
113.00	72.40	70.11
41.00	38.30	39.55
42.00	26.80	24.72

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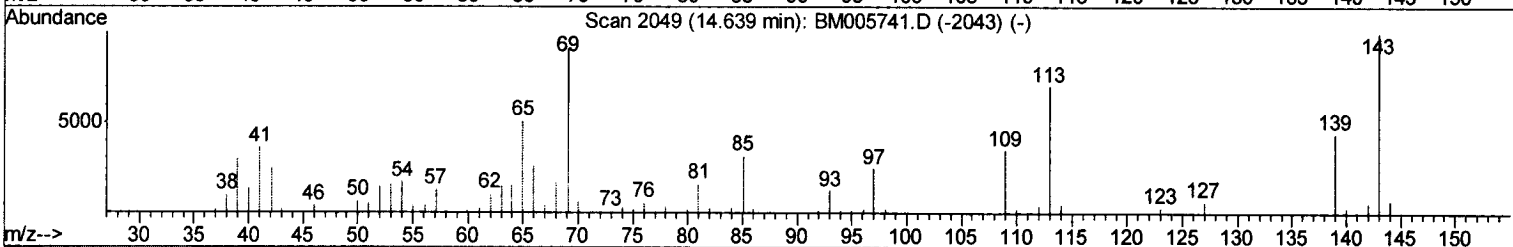
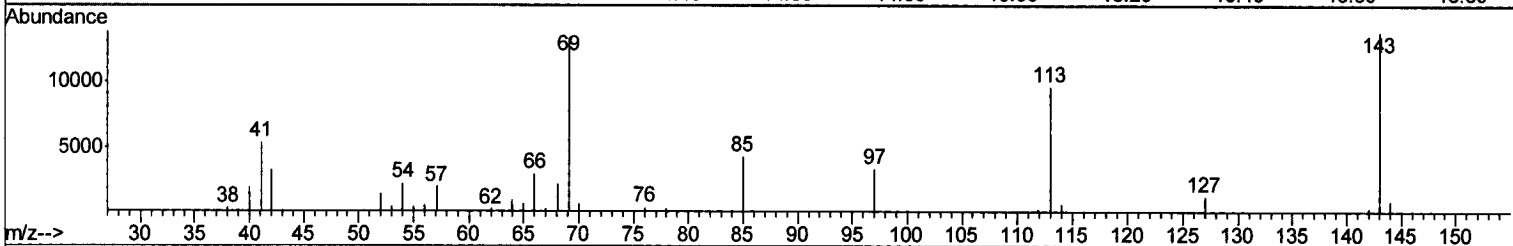
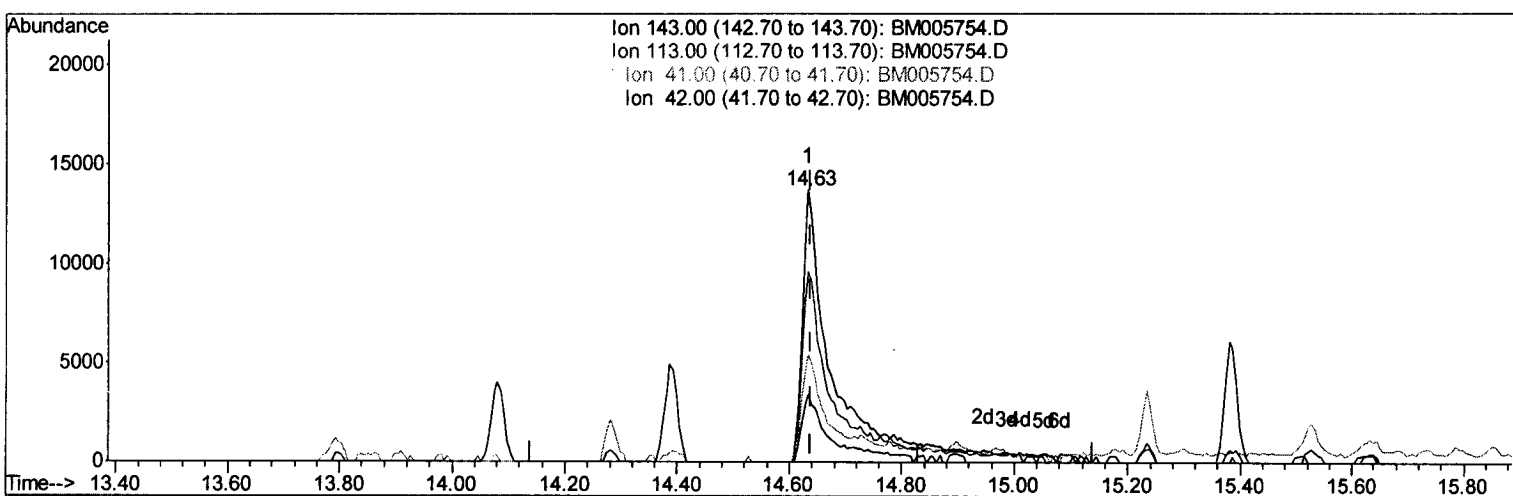
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TIC: BM005754.D

(20) 4-Nitrophenol-d4 (S)

14.633min (-0.006) 14.35ng/ul m

response 48197

Ion	Exp%	Act%
143.00	100	100
113.00	72.40	70.11
41.00	38.30	39.55
42.00	26.80	24.72

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 06/20/16

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	84556	20.00	ng/ul	0.00
7) Naphthalene-d8	10.53	136	447489	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.39	164	278352	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.13	188	621941	20.00	ng/ul	-0.01
29) Chrysene-d12	21.32	240	500812	20.00	ng/ul	0.00
34) Perylene-d12	23.58	264	430511	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.19	96	3059	1.57	ng/uL	0.00
3) Phenol-d5	6.93	99	125734	17.23	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.07	67	87514	19.96	ng/ul	0.00
5) 2-Chlorophenol-d4	7.28	132	114057	19.33	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	113800	18.46	ng/ul	0.00
8) Nitrobenzene-d5	8.90	128	60253	18.68	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	70213	19.26	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	124281	18.79	ng/ul	0.00
12) 4-Chloroaniline-d4	10.68	131	308969	38.72	ng/ul	0.00
16) Dimethylphthalate-d6	13.80	166	478632	20.85	ng/ul	0.00
17) Acenaphthylene-d8	14.08	160	600791	21.39	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	48197m	14.35	ng/ul	0.00
21) Fluorene-d10	15.39	176	415036	21.35	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	27784	8.64	ng/ul	-0.01
26) Anthracene-d10	17.23	188	610331	20.83	ng/ul	0.00
30) Pyrene-d10	19.53	212	610684	25.38	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.43	264	437550	22.28	ng/ul	-0.01

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

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