

Quantitation Report (Qedit)

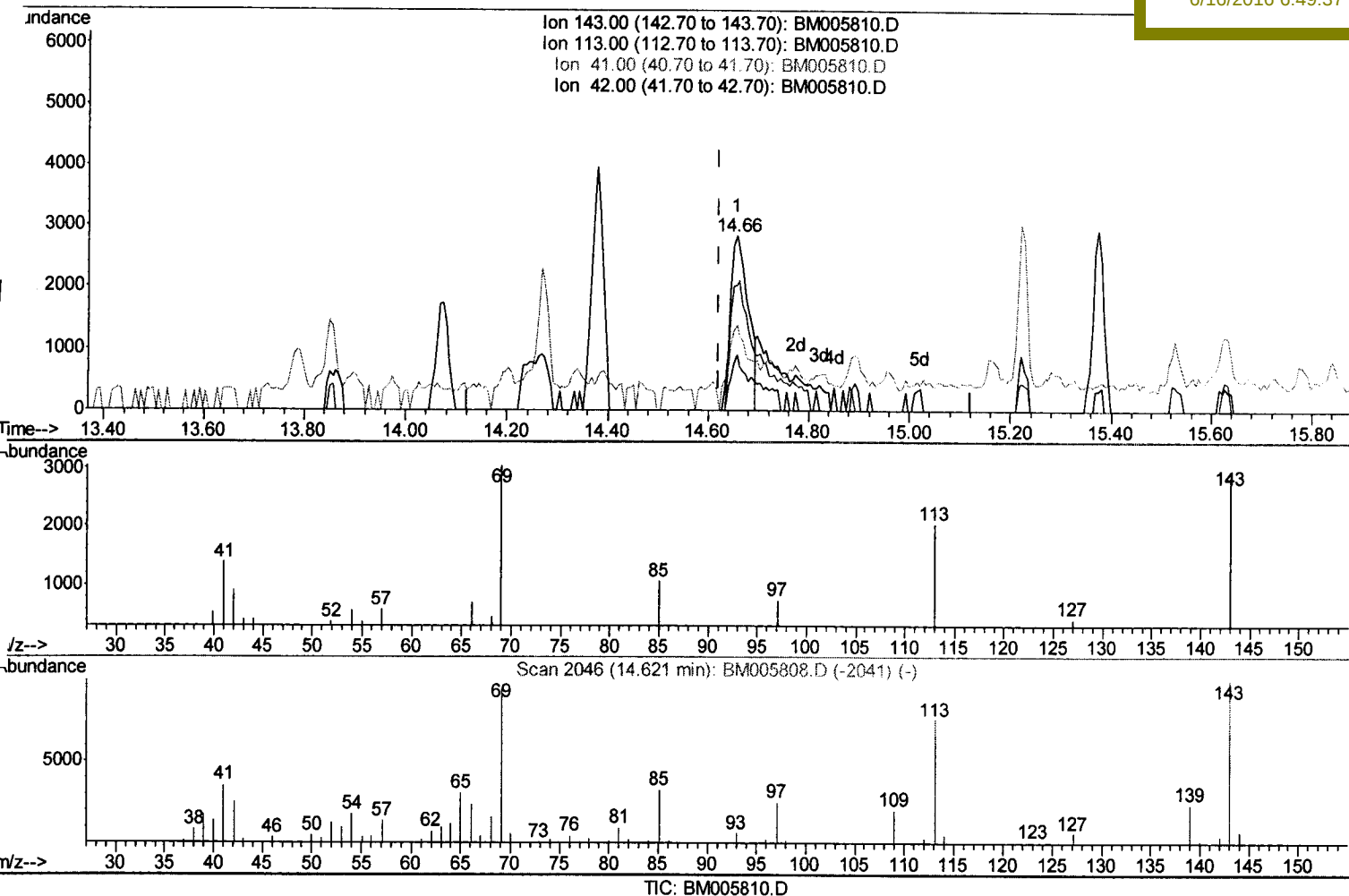
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061616\
 Data File : BM005810.D
 Acq On : 15 Jun 2016 19:02
 Operator : UM/SJ
 Sample : H3537-05
 Disc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 Client Sample ID :
 D9Y37

Quant Time: Jun 16 01:46:28 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 16 01:45:55 2016
 Response via : Initial Calibration

Manual Integrations
 APPROVED

umangi
 6/16/2016 6:49:37 PM



(20) 4-Nitrophenol-d4 (S)

14.657min (+0.035) 3.60ng/ul

response 7006

Ion	Exp%	Act%
143.00	100	100
113.00	74.90	71.86
41.00	39.60	48.70#
42.00	23.50	31.93#

Quantitation Report (Qedit)

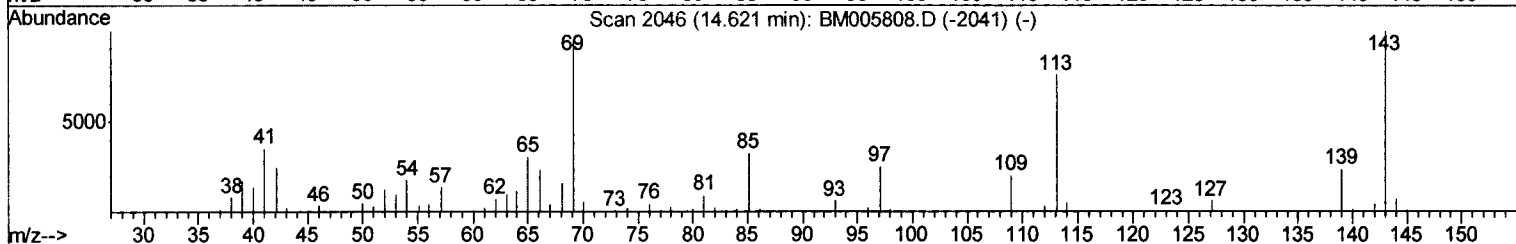
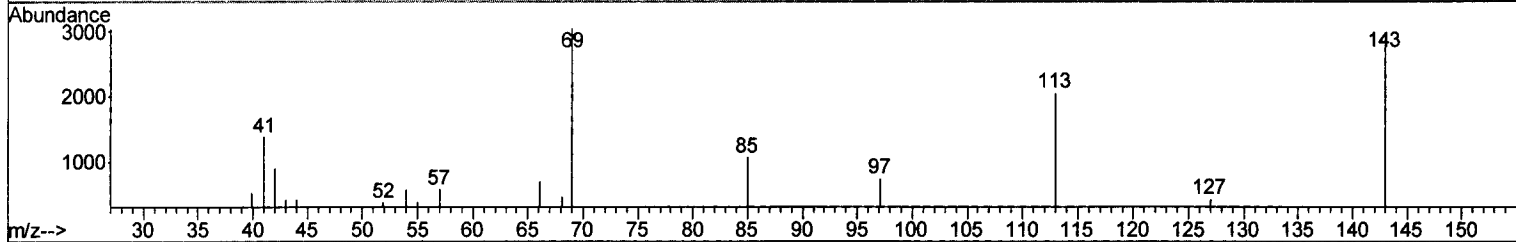
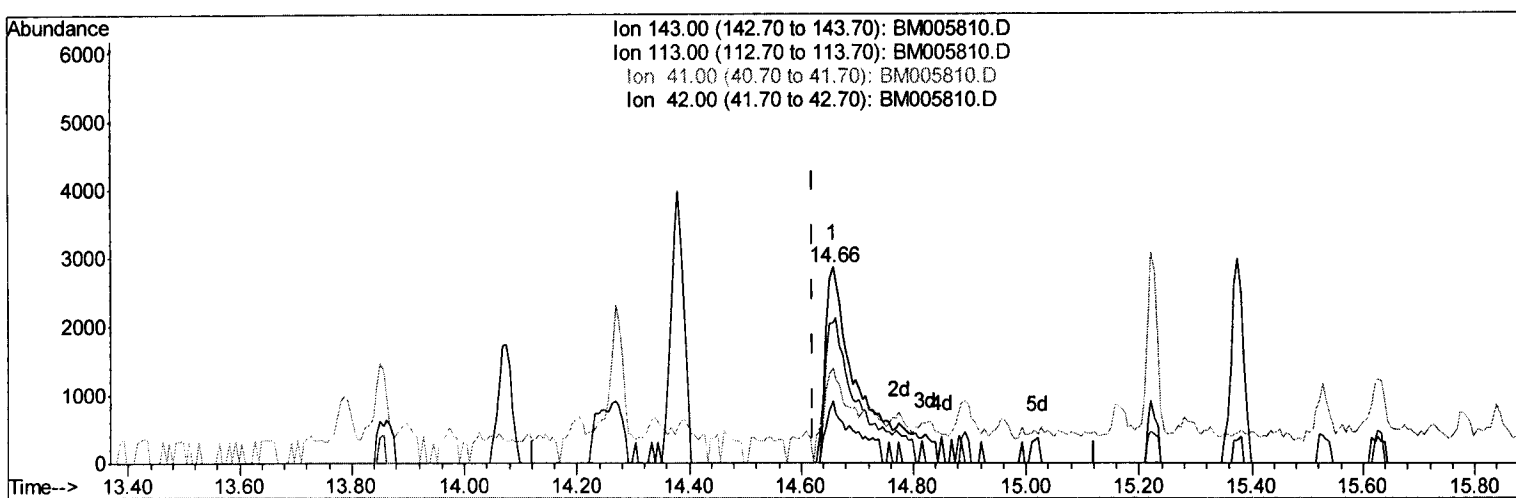
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TIC: BM005810.D

(20) 4-Nitrophenol-d4 (S)

14.657min (+0.035) 6.30ng/ul m

response 12268

Ion	Exp%	Act%
143.00	100	100
113.00	74.90	71.86
41.00	39.60	48.70#
42.00	23.50	31.93#

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 06/20/16

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 ClientSampled :
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	59851	20.00	ng/ul	0.00
7) Naphthalene-d8	10.53	136	300123	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.38	164	213286	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.13	188	524405	20.00	ng/ul	0.00
29) Chrysene-d12	21.31	240	642423	20.00	ng/ul	0.00
34) Perylene-d12	23.56	264	630015	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.19	96	1315	1.02	ng/uL	0.00
3) Phenol-d5	6.92	99	42158	8.58	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	7.07	67	35105	12.00	ng/ul	0.00
5) 2-Chlorophenol-d4	7.27	132	41279	10.08	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	21990	5.29	ng/ul	0.00
8) Nitrobenzene-d5	8.90	128	20811	9.79	ng/ul	0.00
9) 2-Nitrophenol-d4	9.62	143	25085	10.62	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	42401	9.39	ng/ul	0.00
12) 4-Chloroaniline-d4	10.68	131	43287	8.72	ng/ul	0.01
16) Dimethylphthalate-d6	13.79	166	212987	12.58	ng/ul	0.00
17) Acenaphthylene-d8	14.07	160	255202	11.86	ng/ul	0.00
20) 4-Nitrophenol-d4	14.66	143	12268m	6.30	ng/ul	0.04
21) Fluorene-d10	15.37	176	187179	13.27	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	14480	6.17	ng/ul	0.00
26) Anthracene-d10	17.23	188	297547	12.23	ng/ul	0.00
30) Pyrene-d10	19.52	212	364865	11.67	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.42	264	351551	12.08	ng/ul	0.00

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Target Compounds Qvalue

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

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