

Quantitation Report (Qedit)

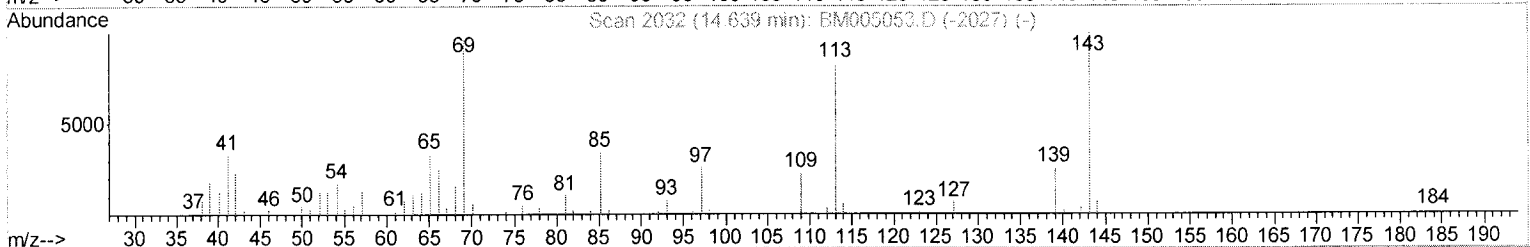
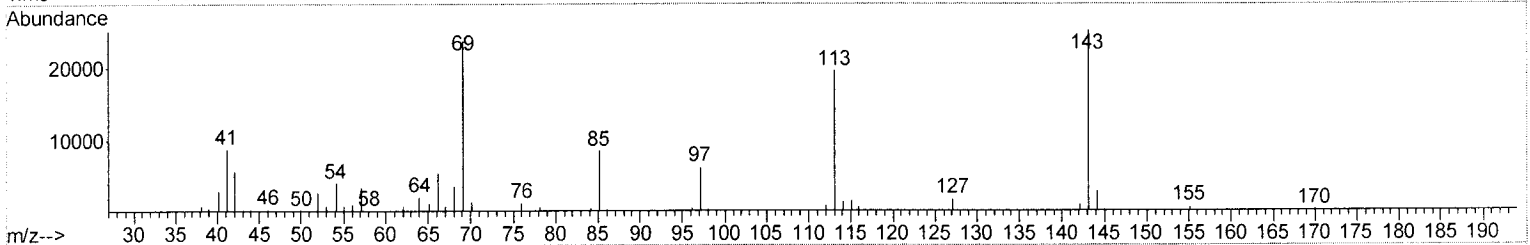
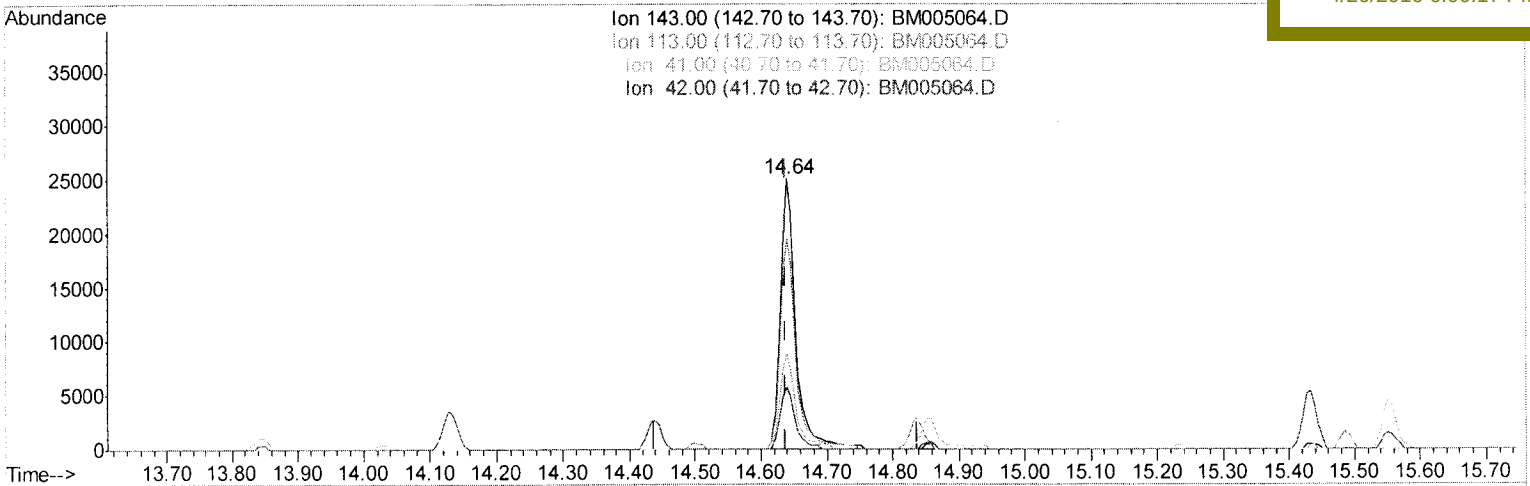
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042516\
 Data File : BM005064.D
 Acq On : 26 Apr 2016 02:31
 Operator : UM/SJ
 Sample : H2584-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
ClientSampleId :
 BC7Q0

Quant Time: Apr 26 04:59:17 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM042516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 26 00:45:11 2016
 Response via : Initial Calibration

Manual Integrations
APPROVED

sohil
 4/26/2016 6:00:17 PM



TIC: BM005064.D

(51) 4-Nitrophenol-d4 (S)

14.639min (+0.000) 18.66ng/ul

response 40889

Ion	Exp%	Act%
143.00	100	100
113.00	75.50	78.22
41.00	35.90	35.18
42.00	22.90	22.82

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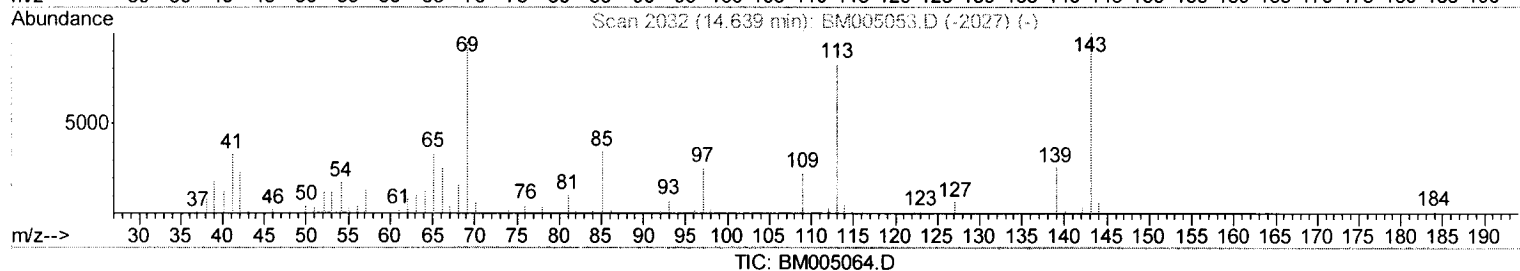
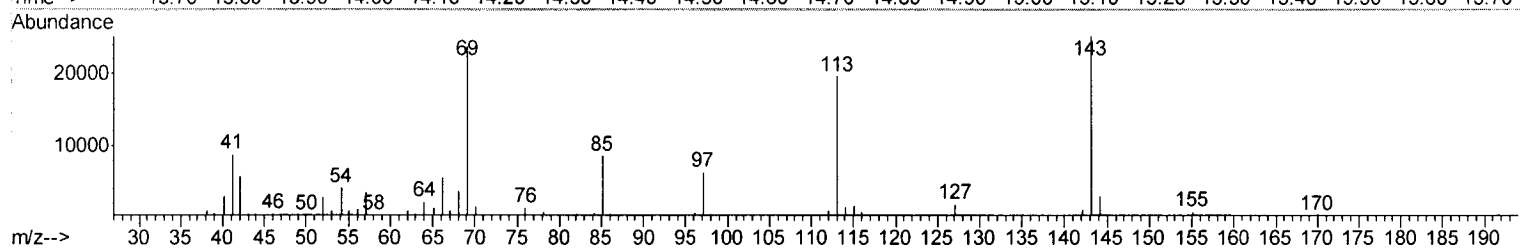
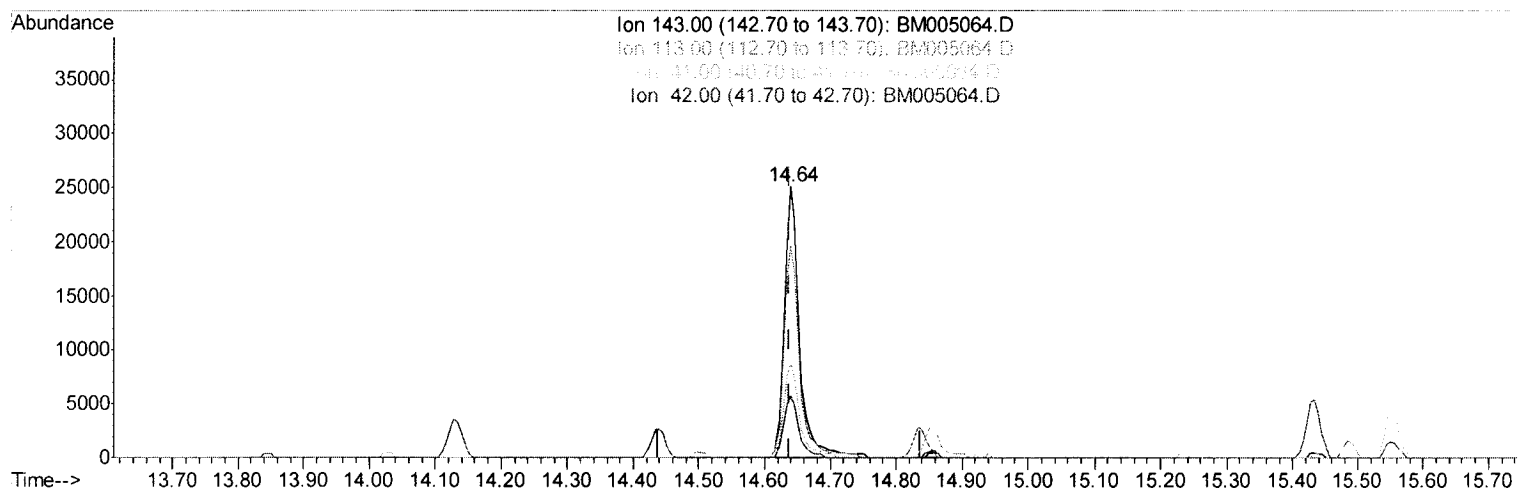
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(51) 4-Nitrophenol-d4 (S)

14.639min (+0.000) 18.72ng/ul m

response 41009

Ion	Exp%	Act%
143.00	100	100
113.00	75.50	78.22
41.00	35.90	35.18
42.00	22.90	22.82

0.00
 04/27/16

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Quant Time: Apr 26 05:19:13 2016
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	38172	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	209971	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	158705	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	382439	20.00	ng/ul	0.00
75) Chrysene-d12	21.37	240	456280	20.00	ng/ul	0.00
83) Perylene-d12	23.66	264	424746	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.30	96	3760	5.01	ng/uL	0.00
5) Phenol-d5	6.96	99	74058	23.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	97409	54.22	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	124641	49.89	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	106476	39.77	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	71451	49.13	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	83085	51.21	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.21	165	141137	45.79	ng/ul	0.00
29) 4-Chloroaniline-d4	10.75	131	666	0.18	ng/ul	0.02
43) Dimethylphthalate-d6	13.85	166	490611	39.14	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	547312	36.67	ng/ul	0.00
51) 4-Nitrophenol-d4	14.64	143	41009m	18.72	ng/ul	0.00
57) Fluorene-d10	15.43	176	410956	37.31	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.56	200	95258	43.79	ng/ul	0.00
70) Anthracene-d10	17.29	188	670537	39.10	ng/ul	0.00
76) Pyrene-d10	19.58	212	792422	38.15	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.52	264	765303	40.21	ng/ul	0.00

U.M.
 04/27/16

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
28) Naphthalene	10.66	128	2899073	273.75	ng/ul	99
34) 2-Methylnaphthalene	12.25	142	147103	18.77	ng/ul	99
40) 1,1'-Biphenyl	13.27	154	21163	1.69	ng/ul	100
47) Acenaphthylene	14.16	152	18337	1.16	ng/ul	97
49) Acenaphthene	14.50	153	102844	9.90	ng/ul	98
53) Dibenzofuran	14.84	168	75973	4.96	ng/ul	99
58) Fluorene	15.49	166	92708	7.79	ng/ul	99
69) Phenanthrene	17.23	178	177301	8.71	ng/ul	99
71) Anthracene	17.32	178	30786	1.50	ng/ul	99
72) Carbazole	17.59	167	33123	1.88	ng/ul	98
74) Fluoranthene	19.24	202	33460	1.49	ng/ul	97

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

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