

Quantitation Report (Qedit)

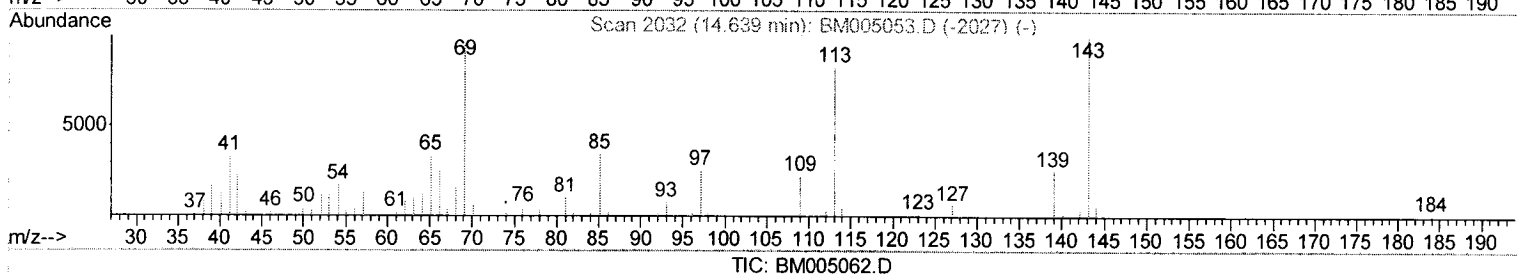
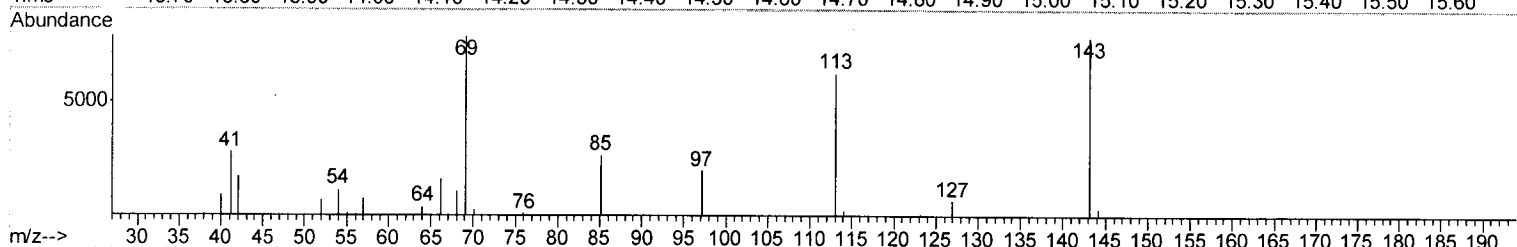
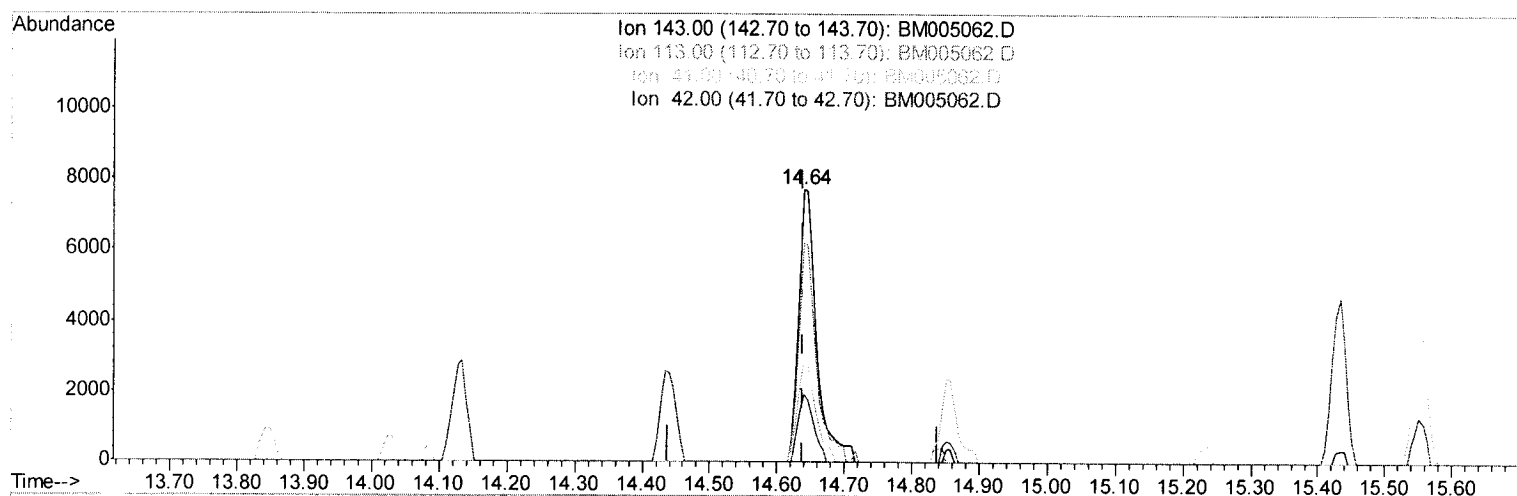
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042516\
 Data File : BM005062.D
 Acq On : 26 Apr 2016 01:18
 Operator : UM/SJ
 Sample : H2584-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA M
 ClientSampleId :
 BC7P8

Manual Integrations
 APPROVED

sohil
 4/26/2016 5:59:32 PM

Quant Time: Apr 26 04:58:55 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



(51) 4-Nitrophenol-d4 (S)

14.639min (+0.000) 6.62ng/ul m

response 14358

Ion	Exp%	Act%
143.00	100	100
113.00	75.50	80.43
41.00	35.90	37.90
42.00	22.90	25.06

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 04/27/16

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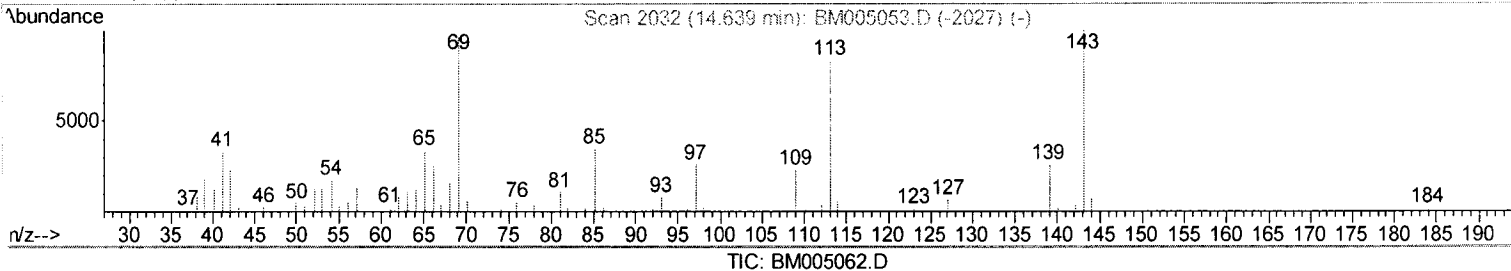
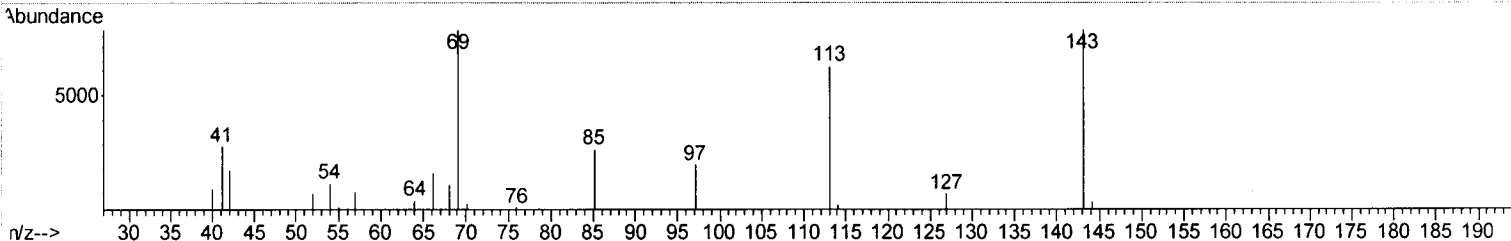
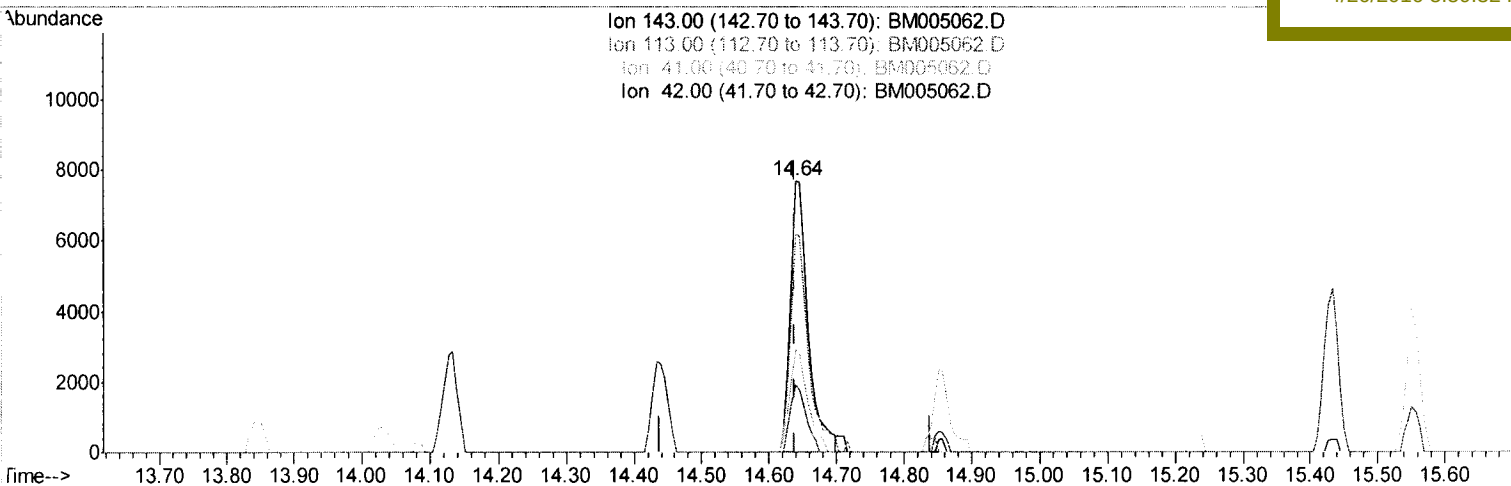
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(51) 4-Nitrophenol-d4 (S)

14.639min (+0.000) 6.48ng/ul

response 14042

Ion	Exp%	Act%
143.00	100	100
113.00	75.50	80.43
41.00	35.90	37.90
42.00	22.90	25.06

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	50411	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	241160	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	156995	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	382126	20.00	ng/ul	0.00
75) Chrysene-d12	21.37	240	451990	20.00	ng/ul	0.00
83) Perylene-d12	23.66	264	411779	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	1117	1.13	ng/uL	0.00
5) Phenol-d5	6.97	99	26190	6.28	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	63760	26.87	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	72536	21.98	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	52756	14.92	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	46838	28.04	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	54015	28.99	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.21	165	94858	26.80	ng/ul	0.00
29) 4-Chloroaniline-d4	10.74	131	583	0.14	ng/ul	0.01
43) Dimethylphthalate-d6	13.85	166	406007	32.75	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	428925	29.05	ng/ul	0.00
51) 4-Nitrophenol-d4	14.64	143	14358m	6.62	ng/ul	0.00
57) Fluorene-d10	15.43	176	338331	31.05	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.56	200	76895	35.38	ng/ul	0.00
70) Anthracene-d10	17.29	188	575539	33.59	ng/ul	0.00
76) Pyrene-d10	19.58	212	674552	32.78	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.52	264	621814	33.70	ng/ul	0.00

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Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
69) Phenanthrene	17.23	178	37658	1.85	ng/ul	100
72) Carbazole	17.59	167	18420	1.05	ng/ul	99

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

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Instrument :
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ClientSampled :
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Manual Integrations
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