

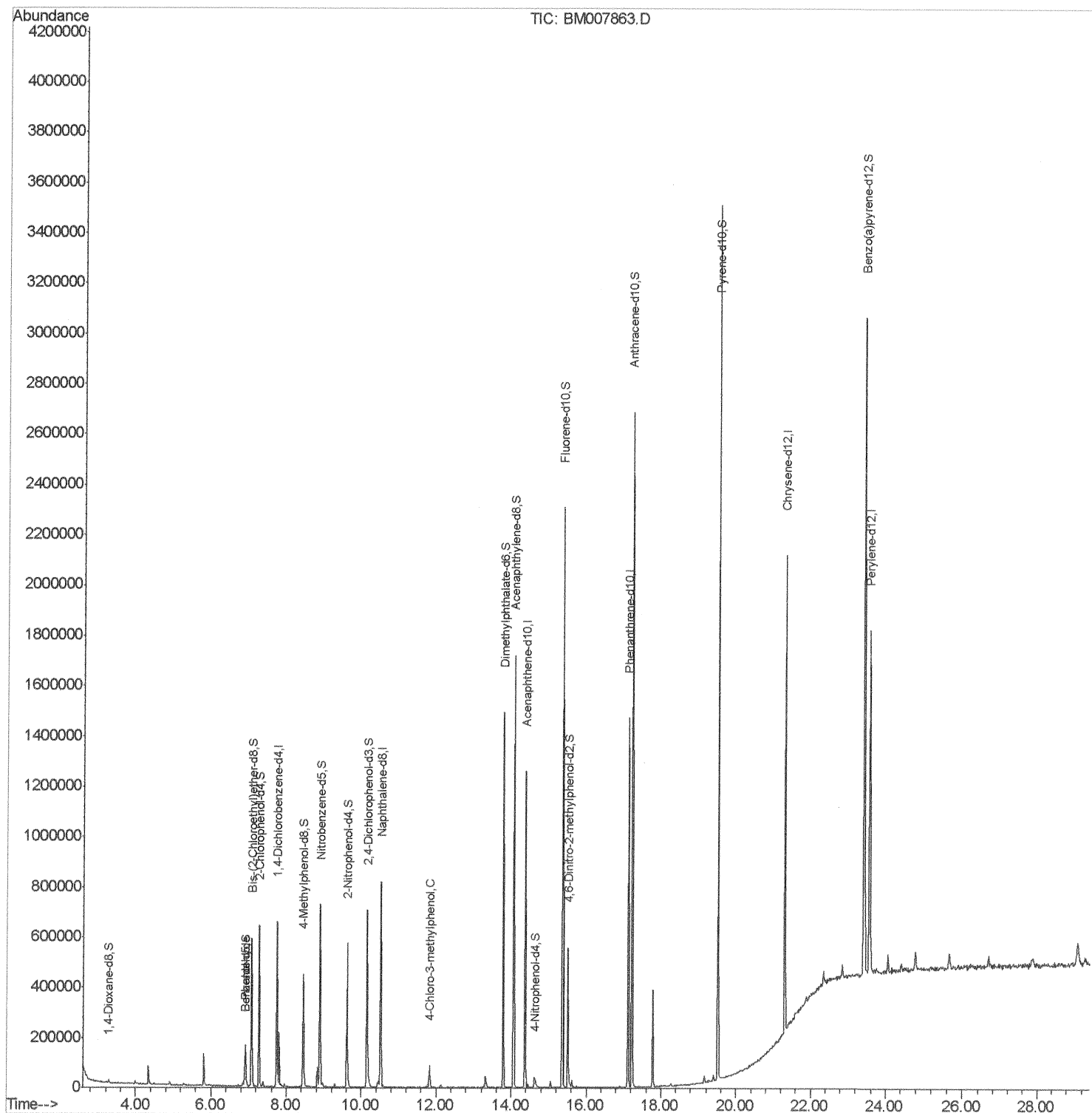
Data Path : Z:\HPCHEM1\BNA_M\Data\BM111416\
Data File : BM007863.D
Acq On : 14 Nov 2016 20:00
Operator : UM/SJ
Sample : H5623-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BD4Q2

Manual Integrations
APPROVED

umangi
11/15/2016 3:08:11 PM

Quant Time: Nov 15 03:21:18 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM102016.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 15 02:47:54 2016
Response via : Initial Calibration



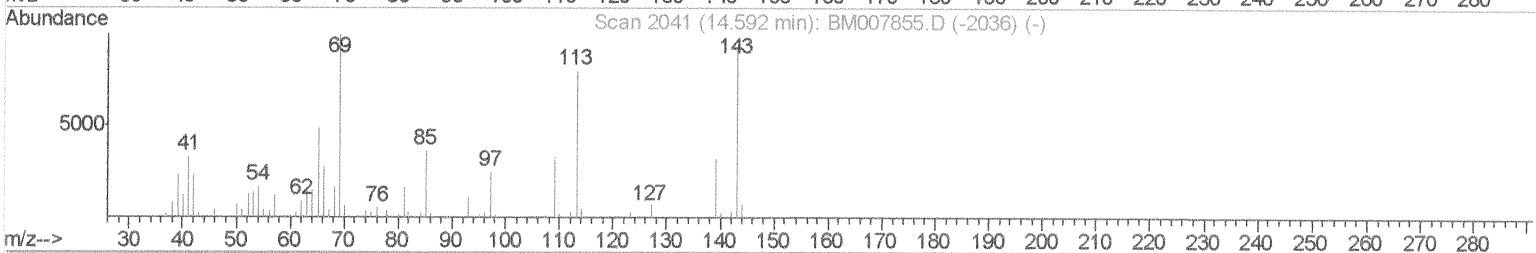
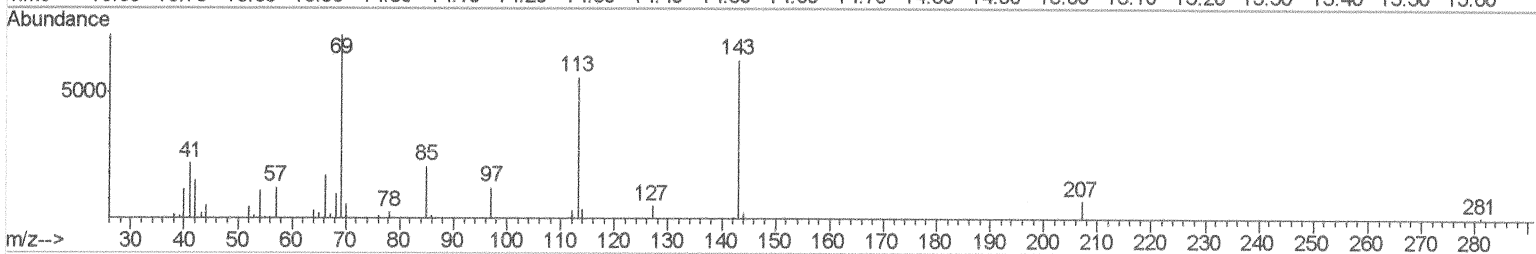
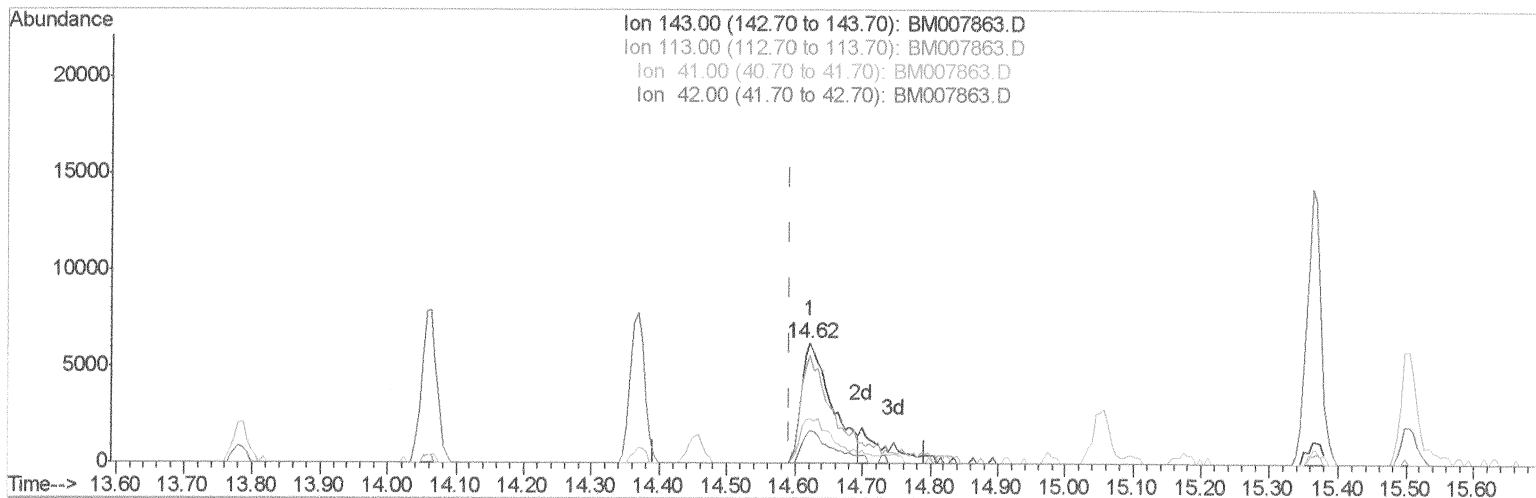
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TIC: BM007863.D

(51) 4-Nitrophenol-d4 (S)

14.621min (+0.029) 4.31ng/ul

response 19663

Ion	Exp%	Act%
143.00	100	100
113.00	93.70	89.32
41.00	36.40	37.58
42.00	22.60	27.61#

Quantitation Report (Qedit)

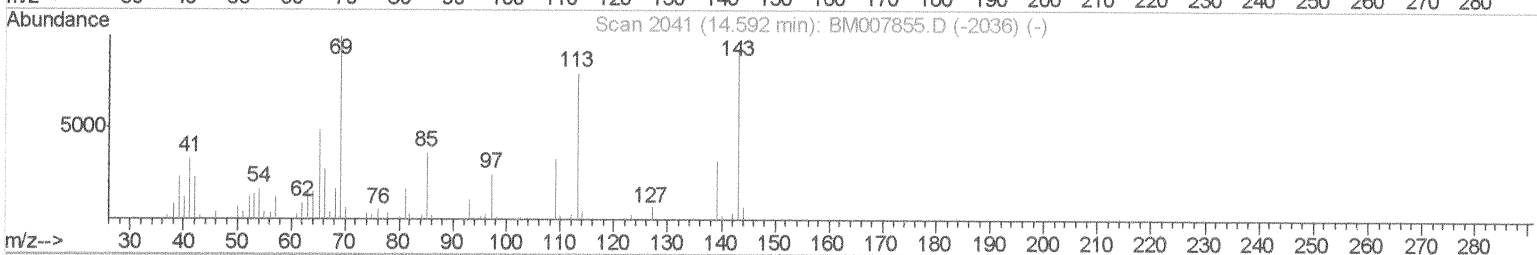
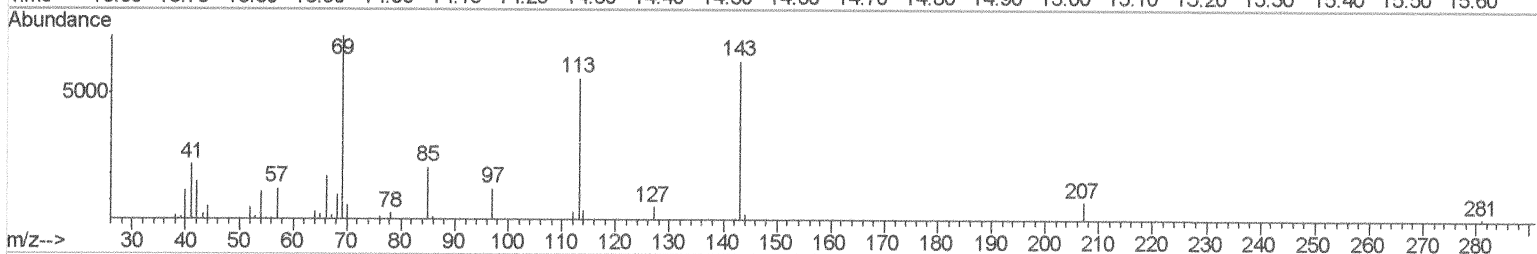
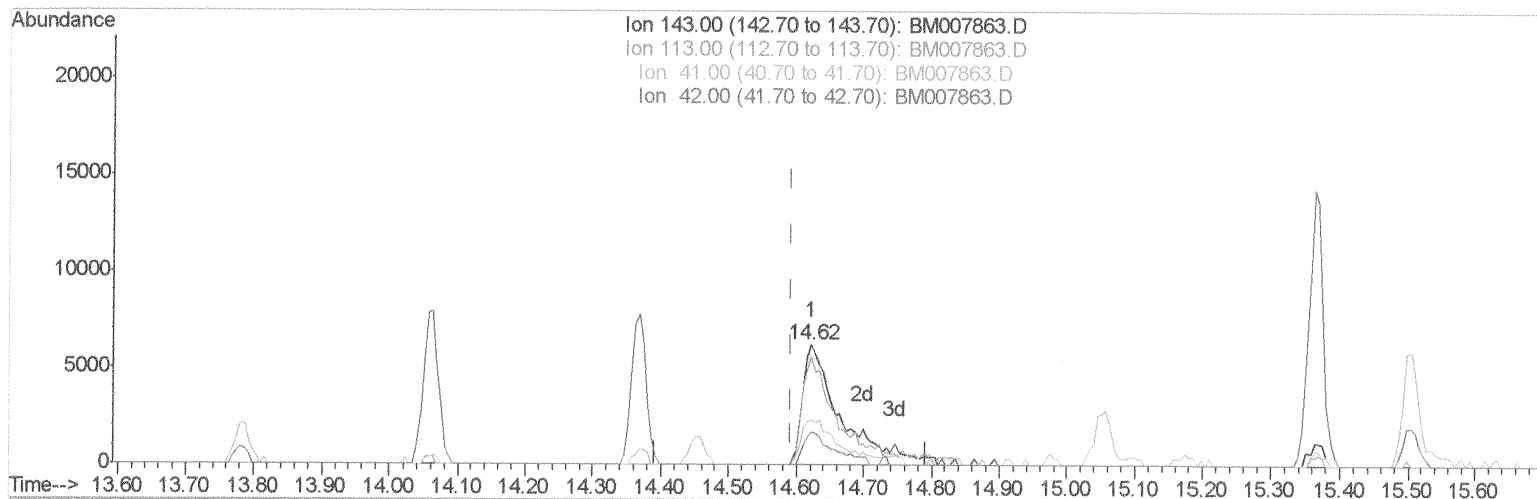
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Manual Integrations
APPROVED

umangi
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 Response via : Initial Calibration



TIC: BM007863.D

(51) 4-Nitrophenol-d4 (S)

14.621min (+0.029) 5.45ng/ul m

52 11/15/16

response 24851

Ion	Exp%	Act%
143.00	100	100
113.00	93.70	89.32
41.00	36.40	37.58
42.00	22.60	27.61#

Data Path : Z:\HPCHEM1\BNA_M\Data\BM111416\
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 BNA_M
 ClientSampleId :
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Manual Integrations
 APPROVED

umangi
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	172189	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	663164	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	423368	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.12	188	838719	20.00	ng/ul	0.00
75) Chrysene-d12	21.31	240	1047604	20.00	ng/ul	0.00
83) Perylene-d12	23.56	264	1074913	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	5711	1.42	ng/uL	0.00
5) Phenol-d5	6.92	99	93519	7.04	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.07	67	259802	32.94	ng/ul	0.00
9) 2-Chlorophenol-d4	7.27	132	283387	27.80	ng/ul	0.00
13) 4-Methylphenol-d8	8.44	113	171074	16.45	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	169323	35.48	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	180686	35.71	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	305854	30.75	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
43) Dimethylphthalate-d6	13.79	166	1031306	35.78	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	1195424	34.45	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	24851m	5.45	ng/ul	0.03
57) Fluorene-d10	15.37	176	920613	36.44	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.50	200	132294	26.55	ng/ul	0.00
70) Anthracene-d10	17.22	188	1532463	40.55	ng/ul	0.00
76) Pyrene-d10	19.52	212	1788726	41.56	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.41	264	1905786	40.61	ng/ul	0.00

Target Compounds

					Qvalue
4) Benzaldehyde	6.90	77	15123	1.80	ng/ul 89
33) 4-Chloro-3-methylphenol	11.83	107	43855	4.27	ng/ul 91

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