

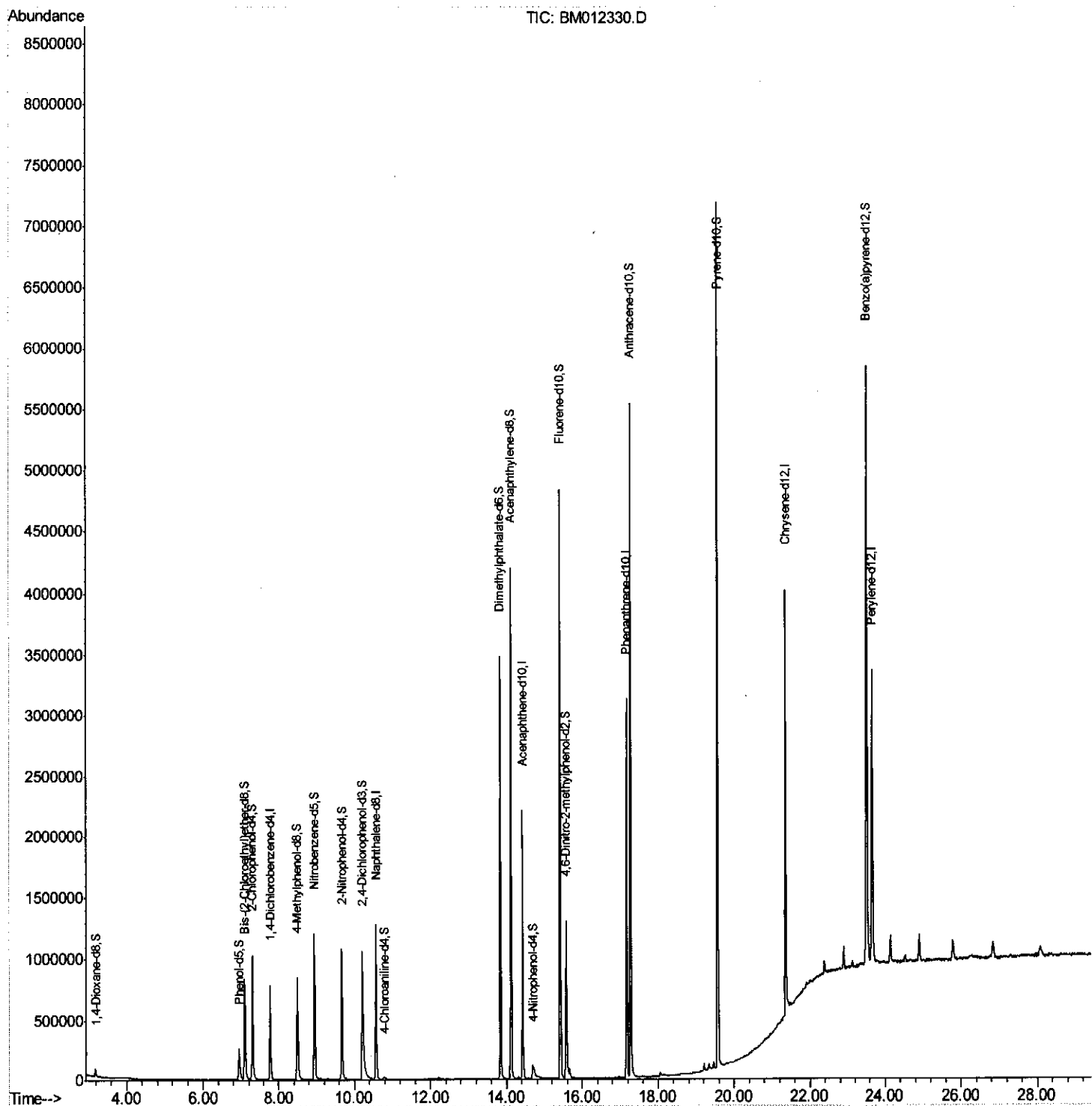
Data Path : Z:\HPCHEM1\BNA M\DATA\BM102017\
Data File : BM012330.D
Acq On : 20 Oct 2017 21:34
Operator : SJ/JU
Sample : I5846-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BD9T2

Manual Integrations
APPROVED

Sohil
10/23/2017 3:53:29 PM

Quant Time: Oct 23 02:25:16 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 23 01:17:35 2017
Response via : Initial Calibration



Quantitation Report (Qedit)

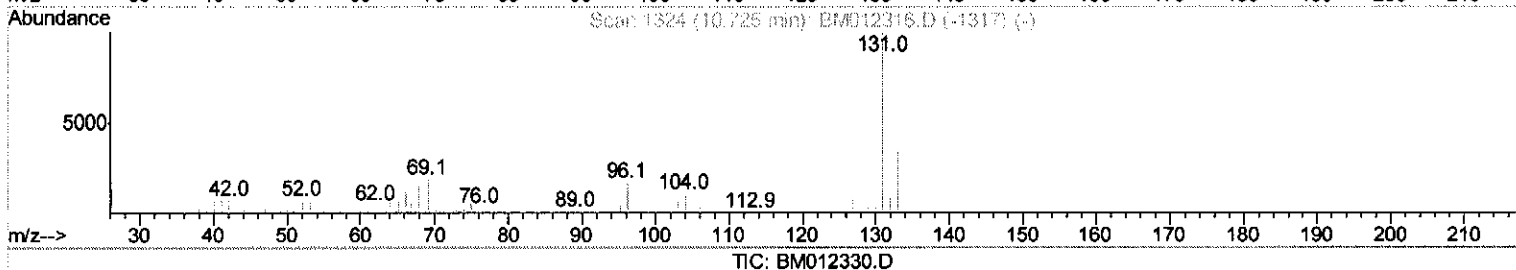
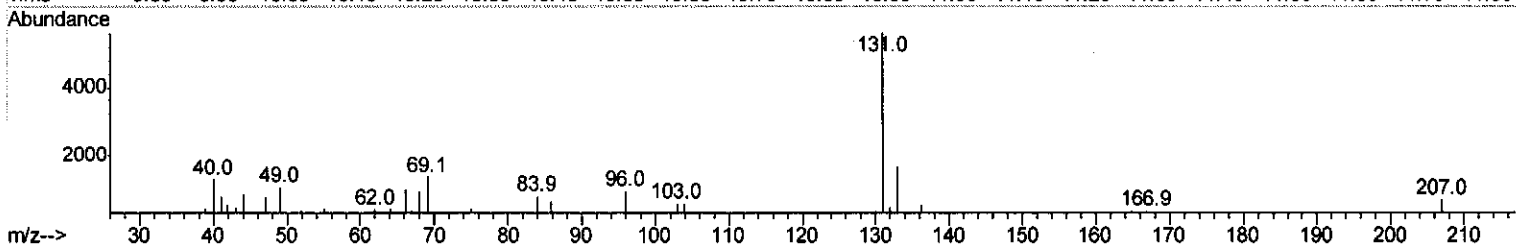
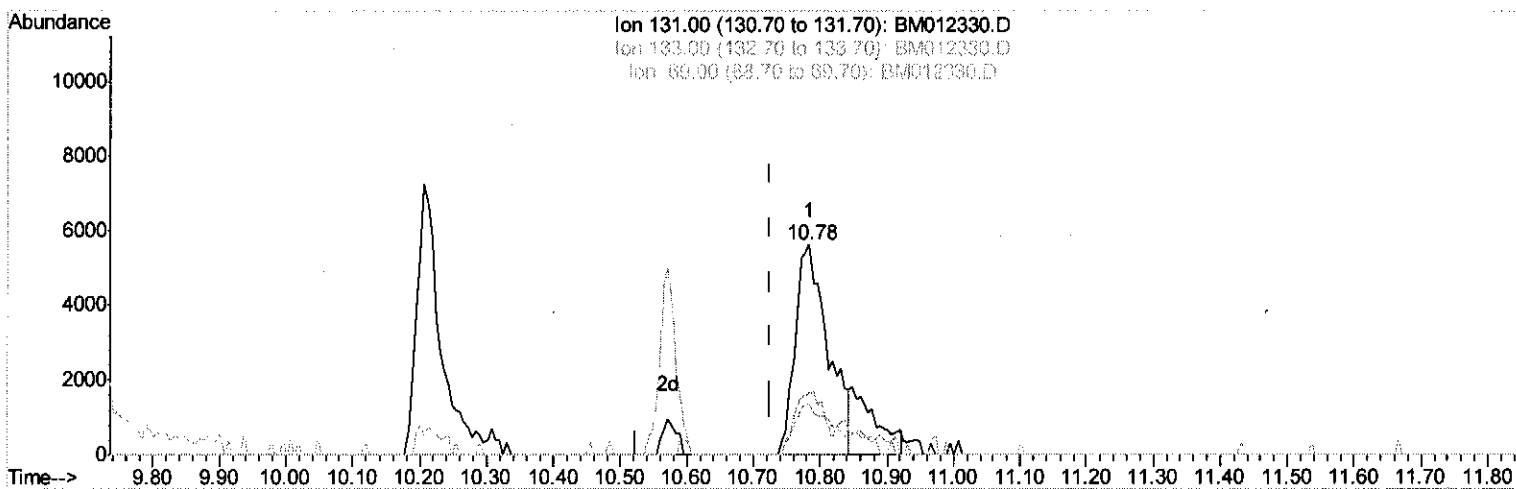
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(29) 4-Chloroaniline-d4 (S)

10.784min (+0.059) 1.16ng/ul

response 19289

Ion	Exp%	Act%
131.00	100	100
133.00	31.40	29.57
69.00	17.40	24.02#
0.00	0.00	0.00

Quantitation Report (Qedit)

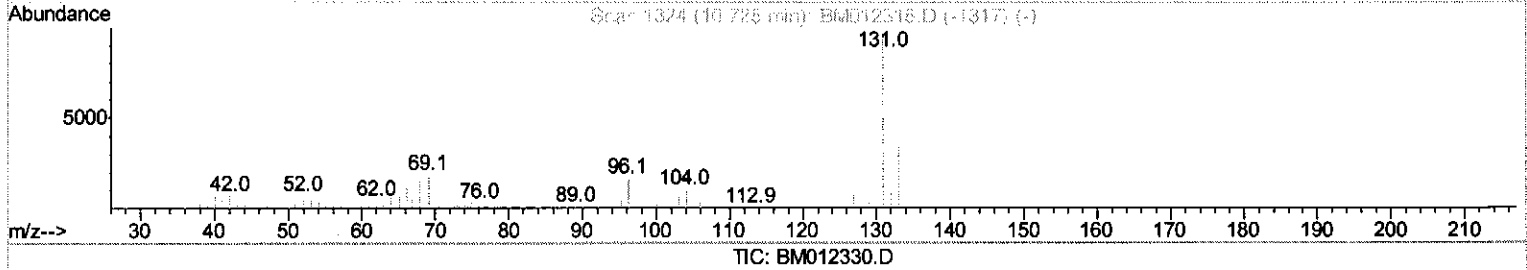
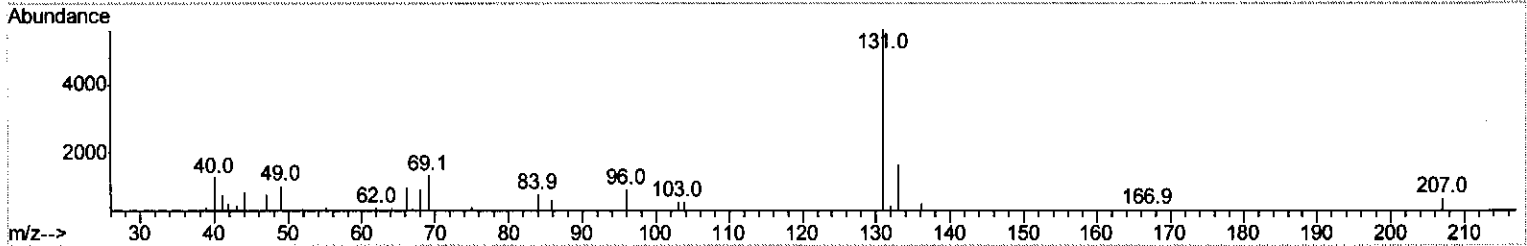
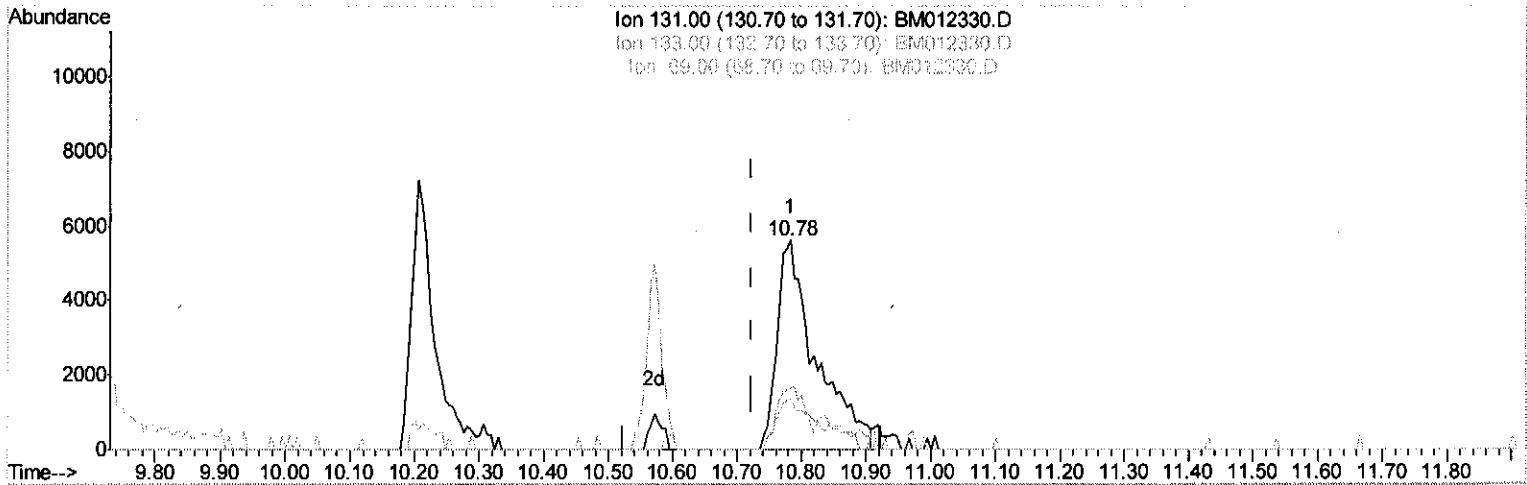
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(29) 4-Chloroaniline-d4 (S)

10.784min (+0.059) 1.41ng/ul m

JU 10/24/17

response 23491

Ion	Exp%	Act%
131.00	100	100
133.00	31.40	29.57
69.00	17.40	24.02#
0.00	0.00	0.00

Quantitation Report (Qedit)

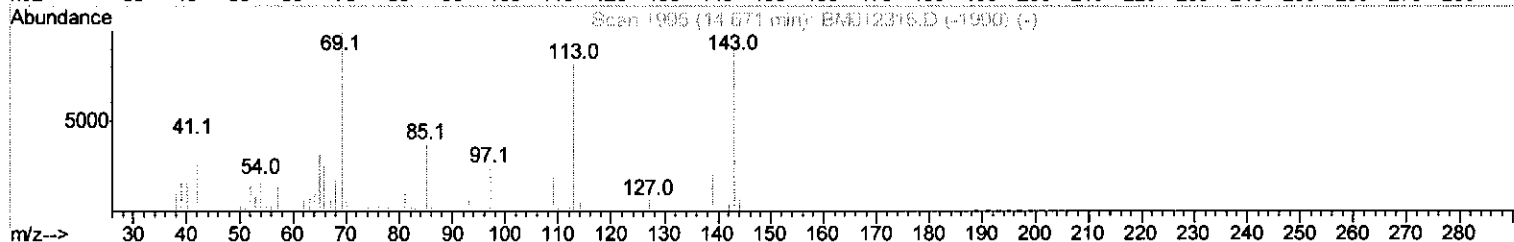
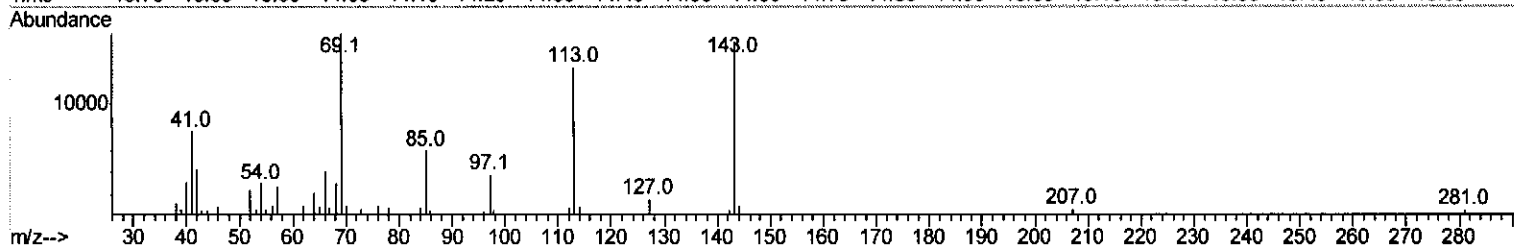
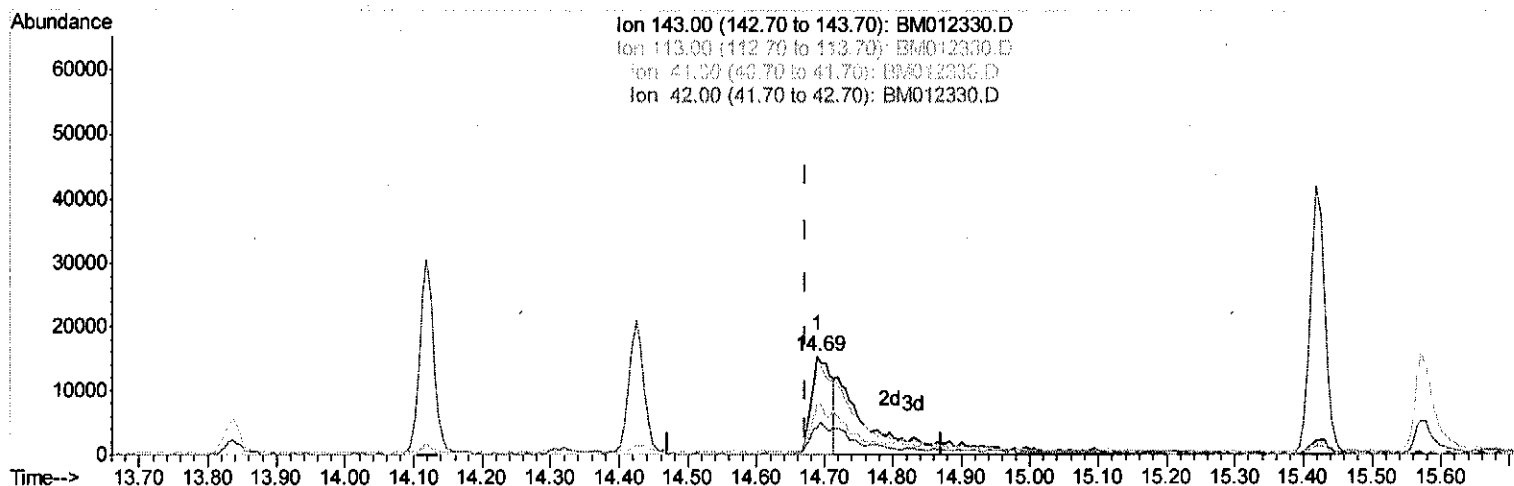
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TIC: BM012330.D

(51) 4-Nitrophenol-d4 (S)

14.689min (+0.018) 3.33ng/ul

response 31206

Ion	Exp%	Act%
143.00	100	100
113.00	81.80	85.83
41.00	40.00	49.37#
42.00	25.70	27.30

Quantitation Report (Qedit)

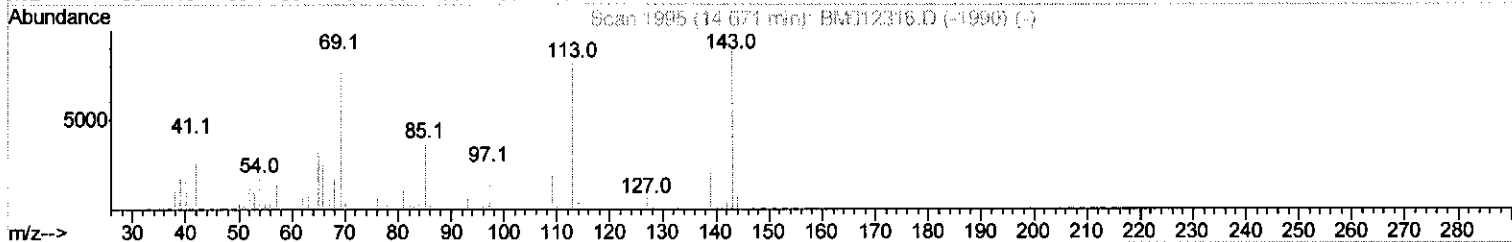
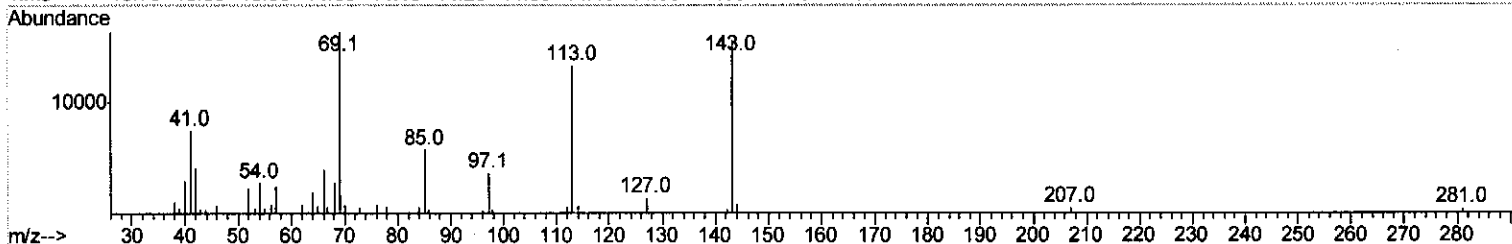
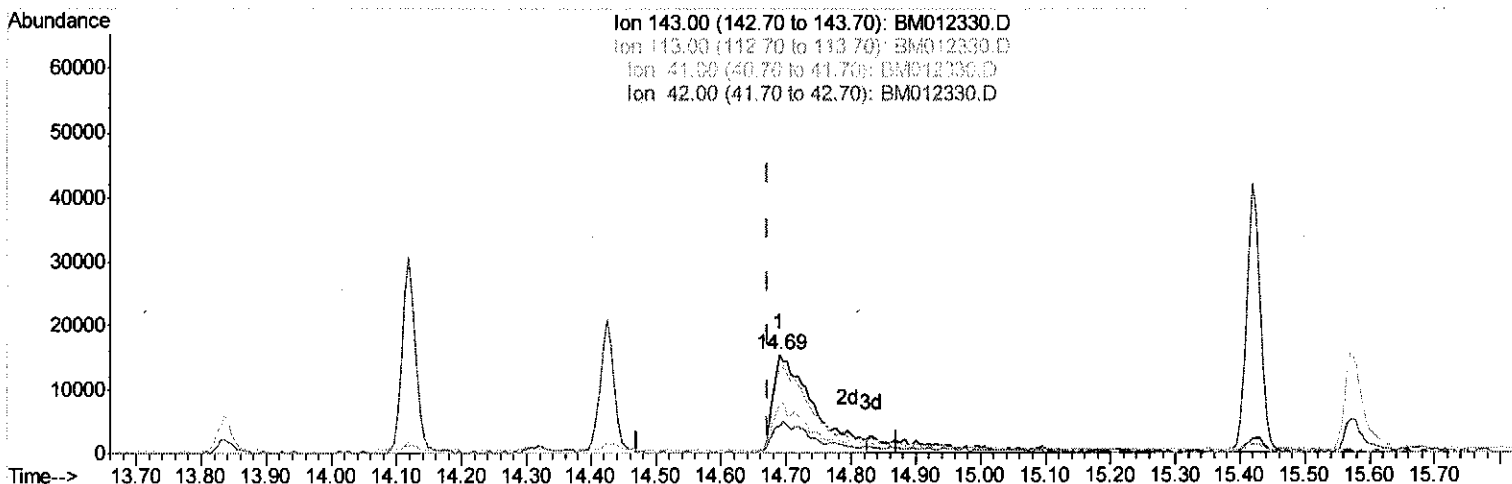
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Manual Integrations
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TIC: BM012330.D

(51) 4-Nitrophenol-d4 (S)

14.689min (+0.018) 6.97ng/ul m

JU 10/24/17

response 65365

Ion	Exp%	Act%
143.00	100	100
113.00	81.80	85.83
41.00	40.00	49.37#
42.00	25.70	27.30

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	232201	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	1044689	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.42	164	705888	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.18	188	1683849	20.00	ng/ul	0.00
75) Chrysene-d12	21.37	240	1881771	20.00	ng/ul	0.00
83) Perylene-d12	23.65	264	1840564	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	32237	6.60	ng/uL	0.00
5) Phenol-d5	6.95	99	197054	10.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	409616	36.25	ng/ul	0.00
9) 2-Chlorophenol-d4	7.31	132	485557	30.38	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	359175	22.14	ng/ul	0.00
19) Nitrobenzene-d5	8.94	128	280626	35.20	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	344965	37.03	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.21	165	564782	31.76	ng/ul	0.00
29) 4-Chloroaniline-d4	10.78	131	23491m	→ 1.41	ng/ul	0.06 → JU 10/24/17
43) Dimethylphthalate-d6	13.84	166	2427788	38.91	ng/ul	0.00
46) Acenaphthylene-d8	14.12	160	2786308	36.89	ng/ul	0.00
51) 4-Nitrophenol-d4	14.69	143	65365m	→ 6.97	ng/ul	0.02 → JU 10/24/17
57) Fluorene-d10	15.42	176	1966889	38.41	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.58	200	347596	30.08	ng/ul	0.00
70) Anthracene-d10	17.28	188	3147455	37.80	ng/ul	0.00
76) Pyrene-d10	19.57	212	3626690	37.97	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.51	264	3422758	38.57	ng/ul	0.00

Target Compounds Ovalue

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