

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112018\
Quantitation Report (QT Reviewed)

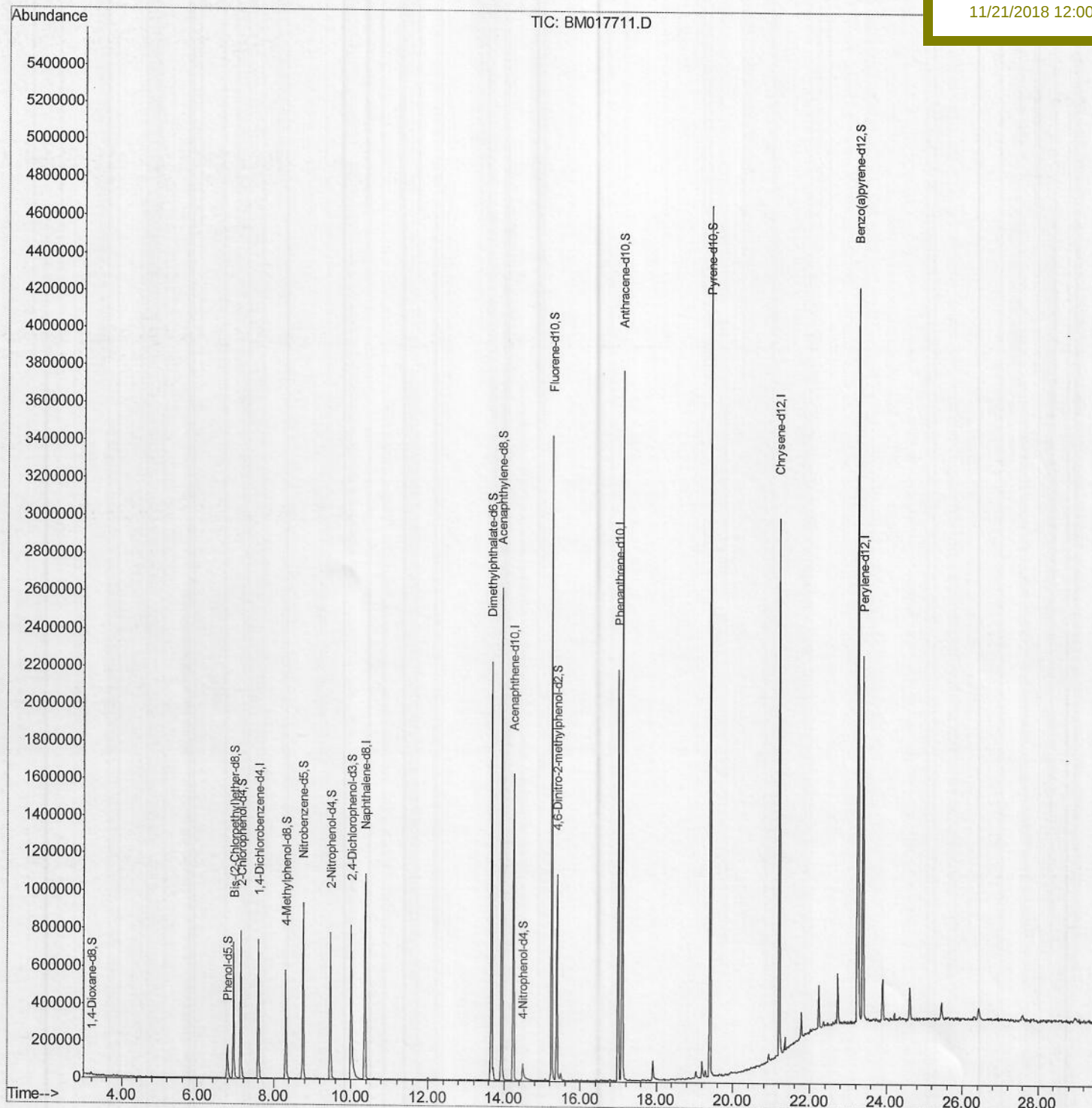
Data File : BM017711.D
Acq On : 20 Nov 2018 19:37
Operator : MJ/SJ
Sample : J6001-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 21 01:26:12 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 21 01:07:57 2018
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
A40W6

Manual Integrations
APPROVED

Sohil
11/21/2018 12:00:11 PM



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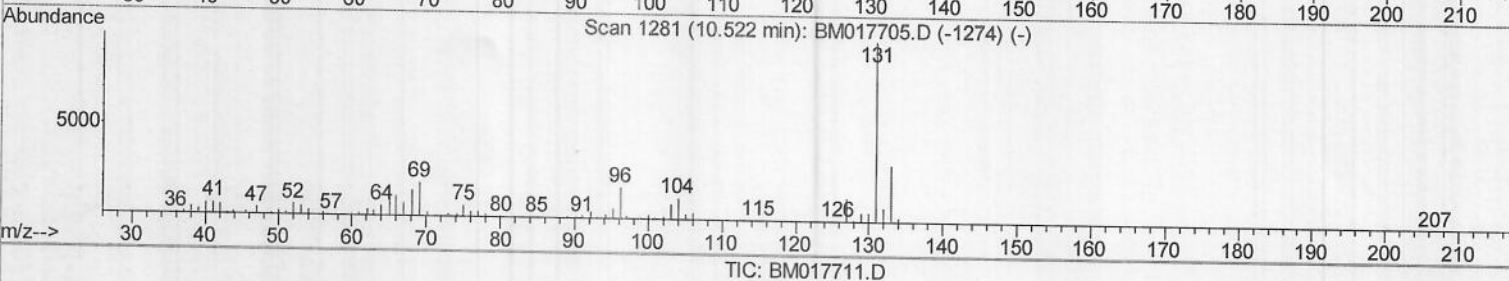
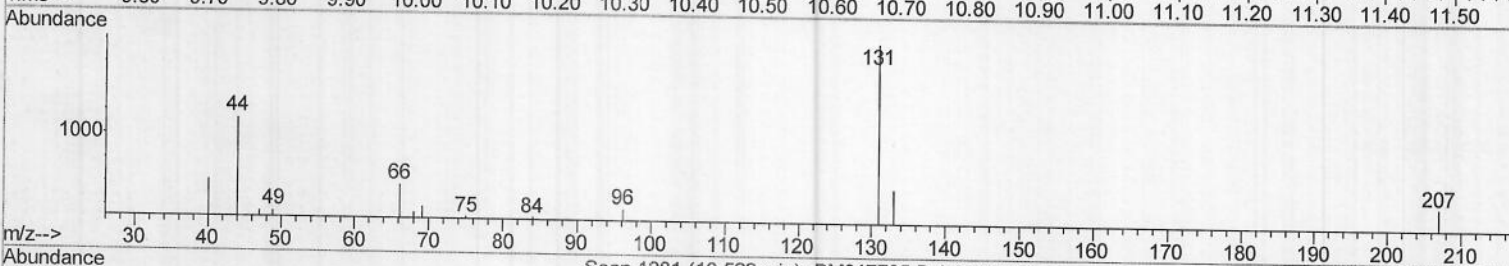
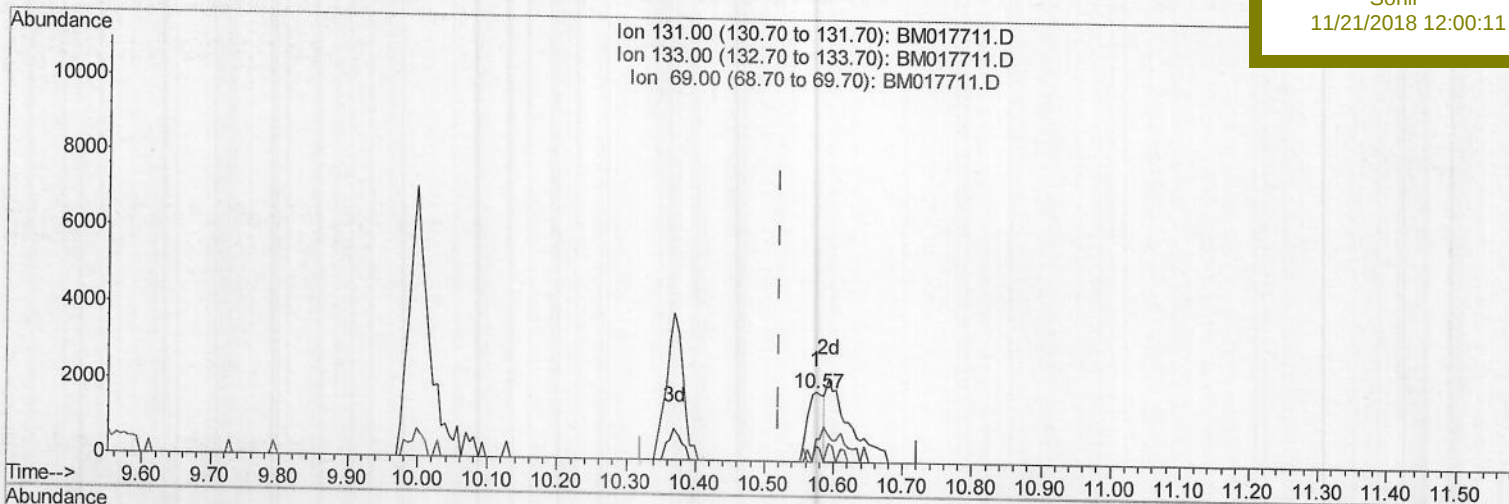
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TIC: BM017711.D

(29) 4-Chloroaniline-d4 (S)

10.575min (+0.053) 0.25ng/ul

response 3112

Ion	Exp%	Act%
131.00	100	100
133.00	32.40	32.88
69.00	18.90	22.39
0.00	0.00	0.00

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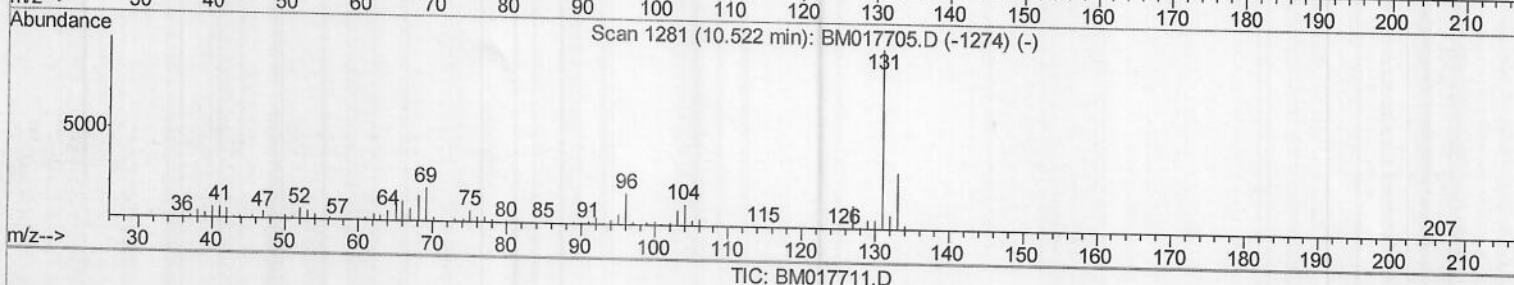
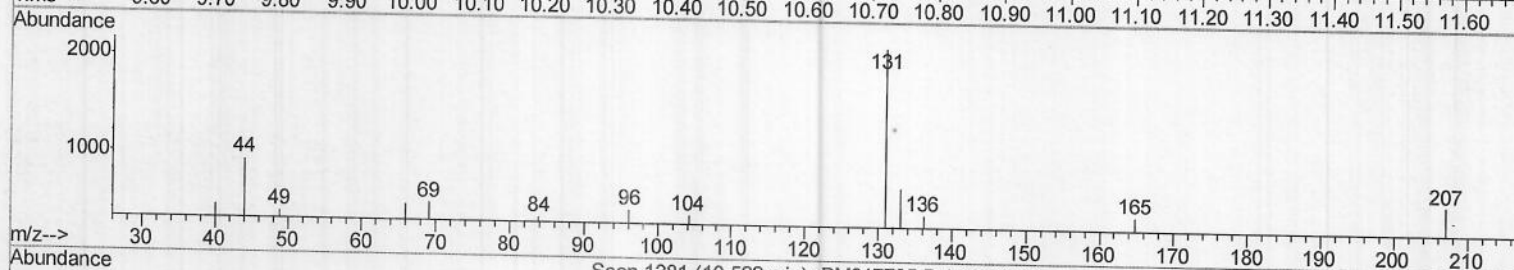
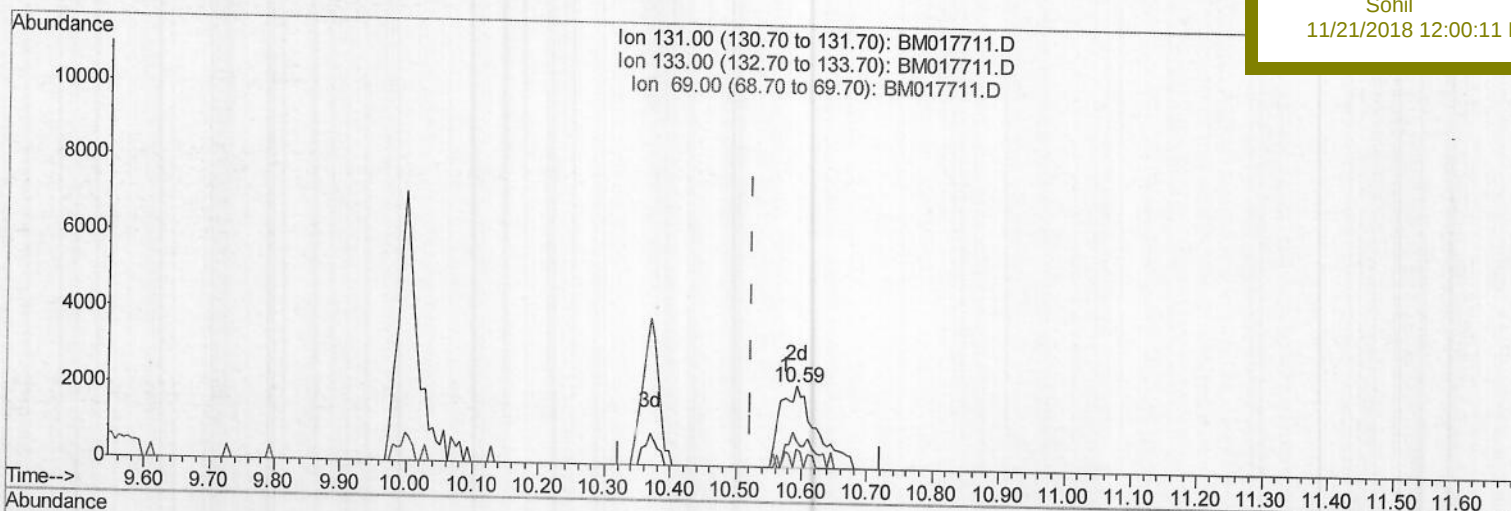
A40W6

Manual Integrations

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Sohil

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(29) 4-Chloroaniline-d4 (S)

10.592min (+0.071) 0.64ng/ul m

response 8089

Ion Exp% Act%

131.00 100 100

133.00 32.40 33.23

69.00 18.90 23.03#

0.00 0.00 0.00

SJ
11/20/18

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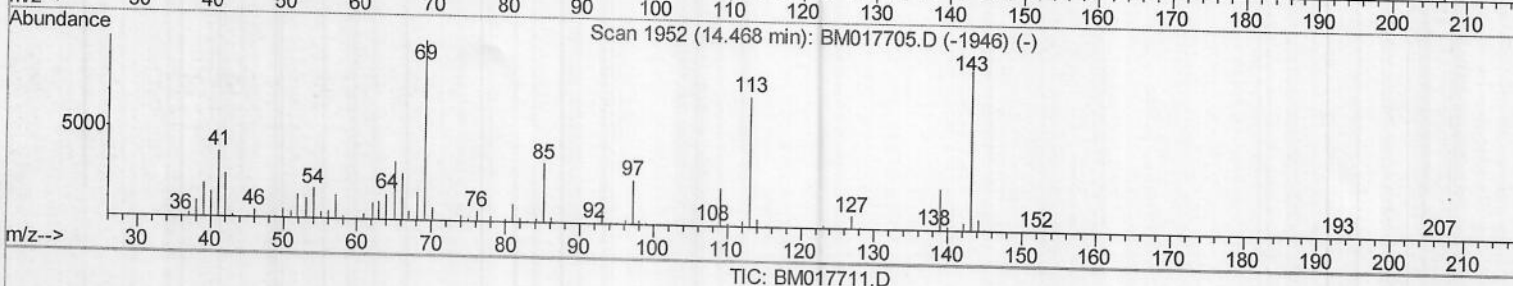
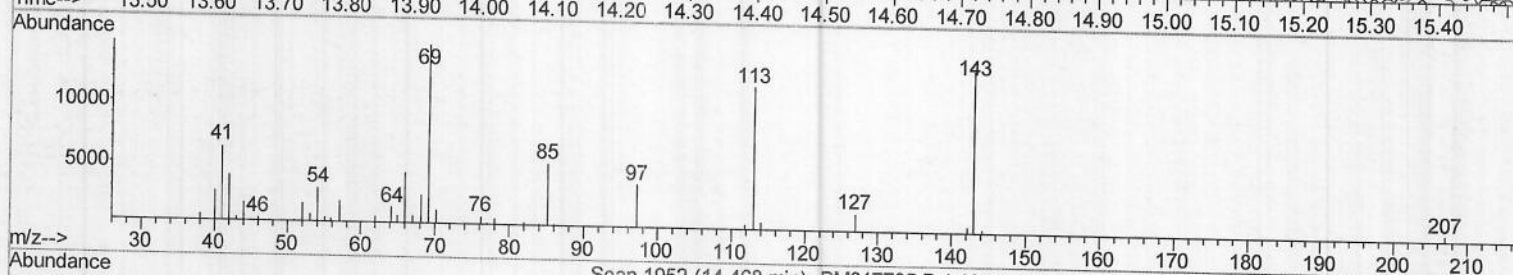
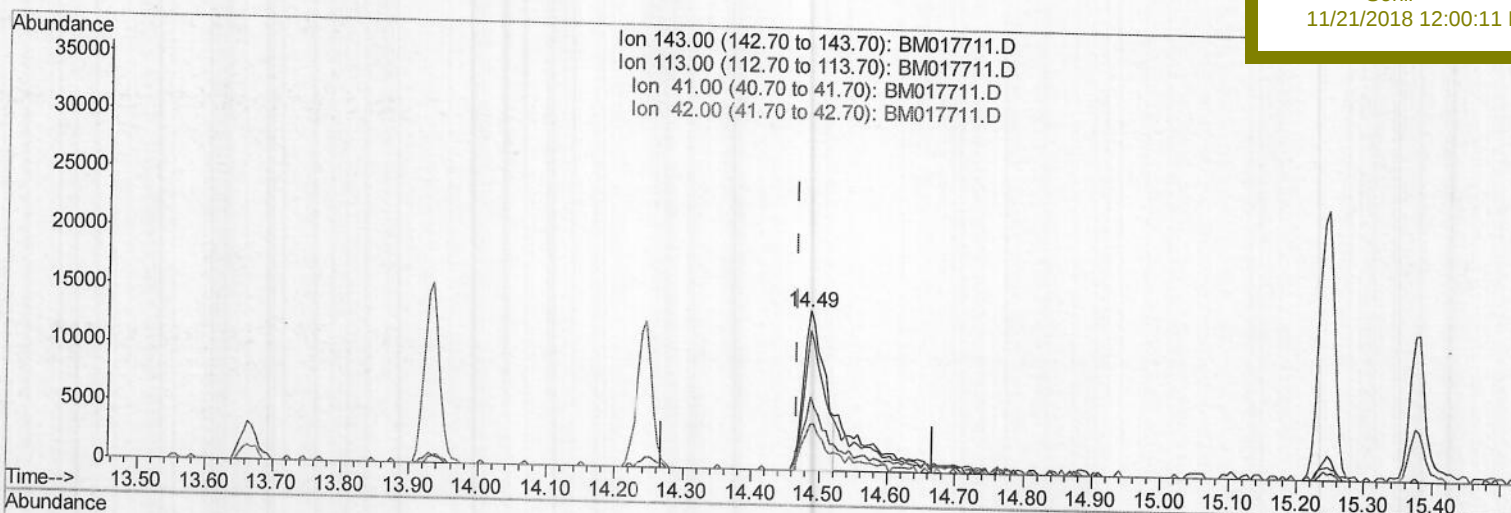
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Instrument :
BNA_M
Client Sampled :
A40W6

Manual Integrations
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(51) 4-Nitrophenol-d4 (S)
14.486min (+0.018) 3.48ng/ul
response 29396

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	87.74
41.00	40.10	46.52
42.00	26.80	29.47

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Operator : MJ/SJ

Sample : J6001-06

Misc :

ALS Vial : 13 Sample Multiplier: 1

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Instrument :

BNA_M

ClientSampleId :

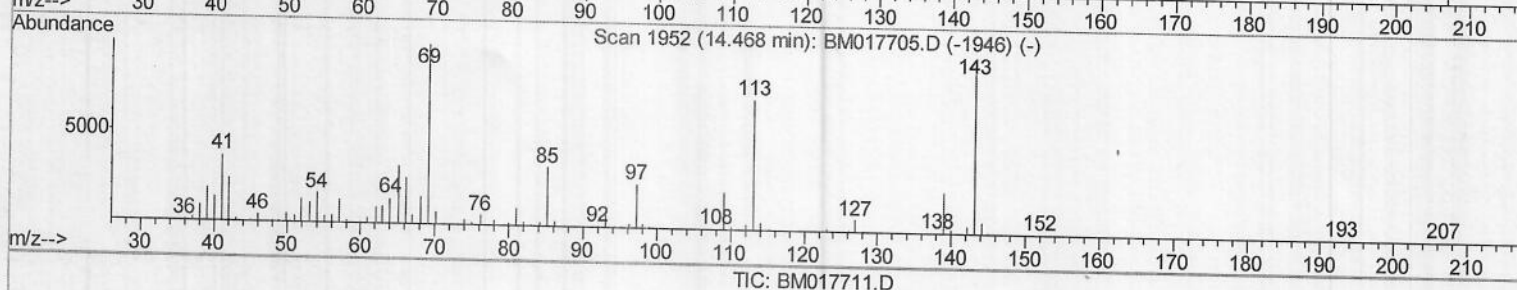
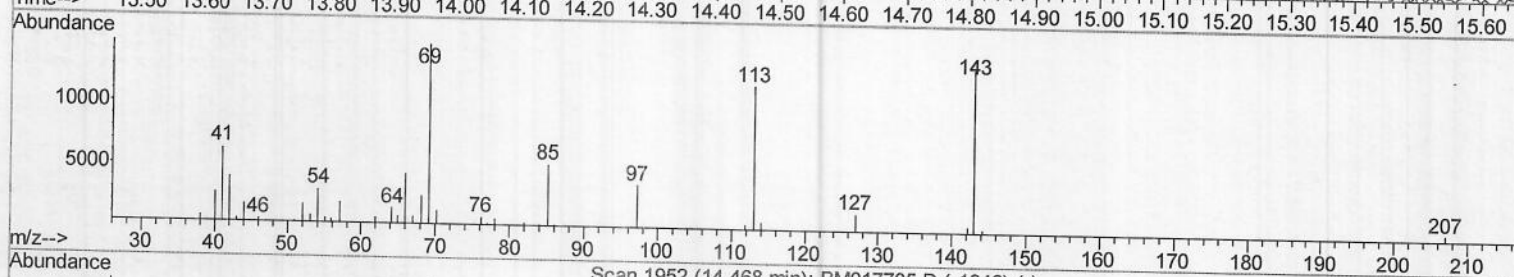
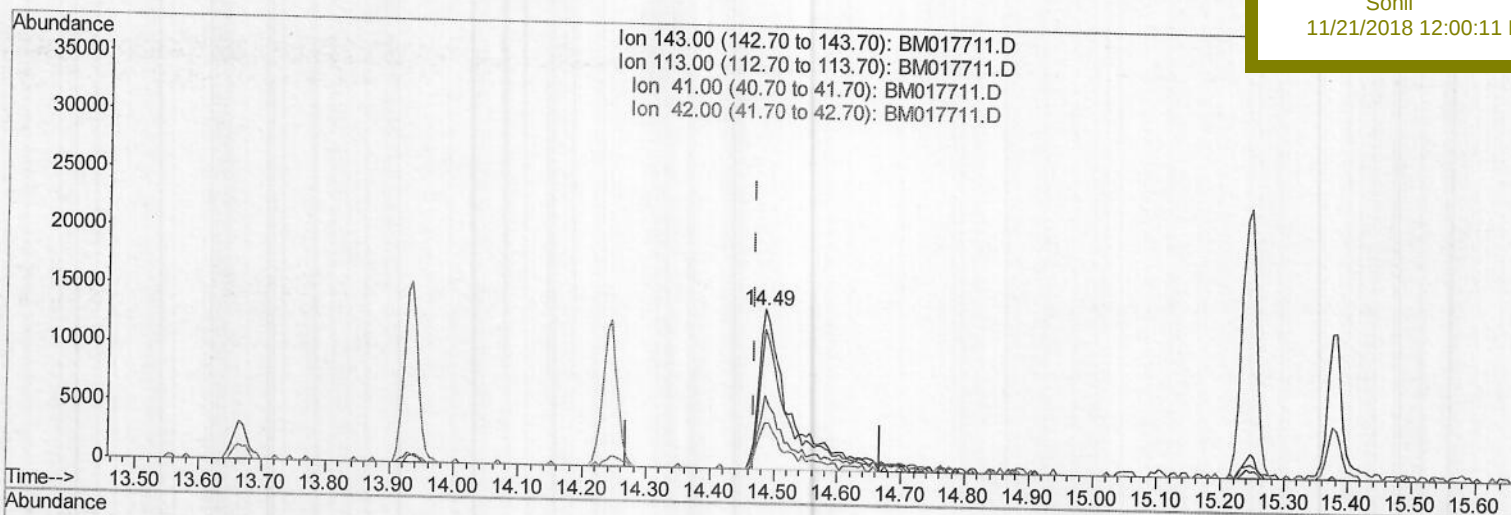
A40W6

Manual Integrations

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(51) 4-Nitrophenol-d4 (S)

14.486min (+0.018) 5.44ng/ul m

response 45910

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	87.74
41.00	40.10	46.52
42.00	26.80	29.47

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	199861	20.00	ng/ul	0.00
18) Naphthalene-d8	10.37	136	881961	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.25	164	535423	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.00	188	1250251	20.00	ng/ul	0.00
77) Chrysene-d12	21.20	240	1412263	20.00	ng/ul	0.00
85) Perylene-d12	23.39	264	1408488	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	7191	1.65	ng/uL	0.00
5) Phenol-d5	6.79	99	124919	6.84	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.95	67	330689	30.04	ng/ul	0.00
9) 2-Chlorophenol-d4	7.13	132	363176	25.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.30	113	238537	15.99	ng/ul	0.00
19) Nitrobenzene-d5	8.75	128	219044	32.34	ng/ul	0.00
22) 2-Nitrophenol-d4	9.47	143	245803	31.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.00	165	410297	27.77	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	8089m	0.64	ng/ul	0.07
43) Dimethylphthalate-d6	13.66	166	1524490	32.81	ng/ul	0.00
46) Acenaphthylene-d8	13.93	160	1830134	32.55	ng/ul	0.00
51) 4-Nitrophenol-d4	14.49	143	45910m	5.44	ng/ul	0.02
57) Fluorene-d10	15.24	176	1312473	32.72	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.38	200	235829	26.39	ng/ul	0.00
70) Anthracene-d10	17.10	188	2074889	33.33	ng/ul	0.00
78) Pyrene-d10	19.40	212	2371256	34.70	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.25	264	2600867	34.05	ng/ul	0.00

} SJ
11/20/18

Target Compounds Qvalue

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