

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110118\
Quantitation Report (QT Reviewed)

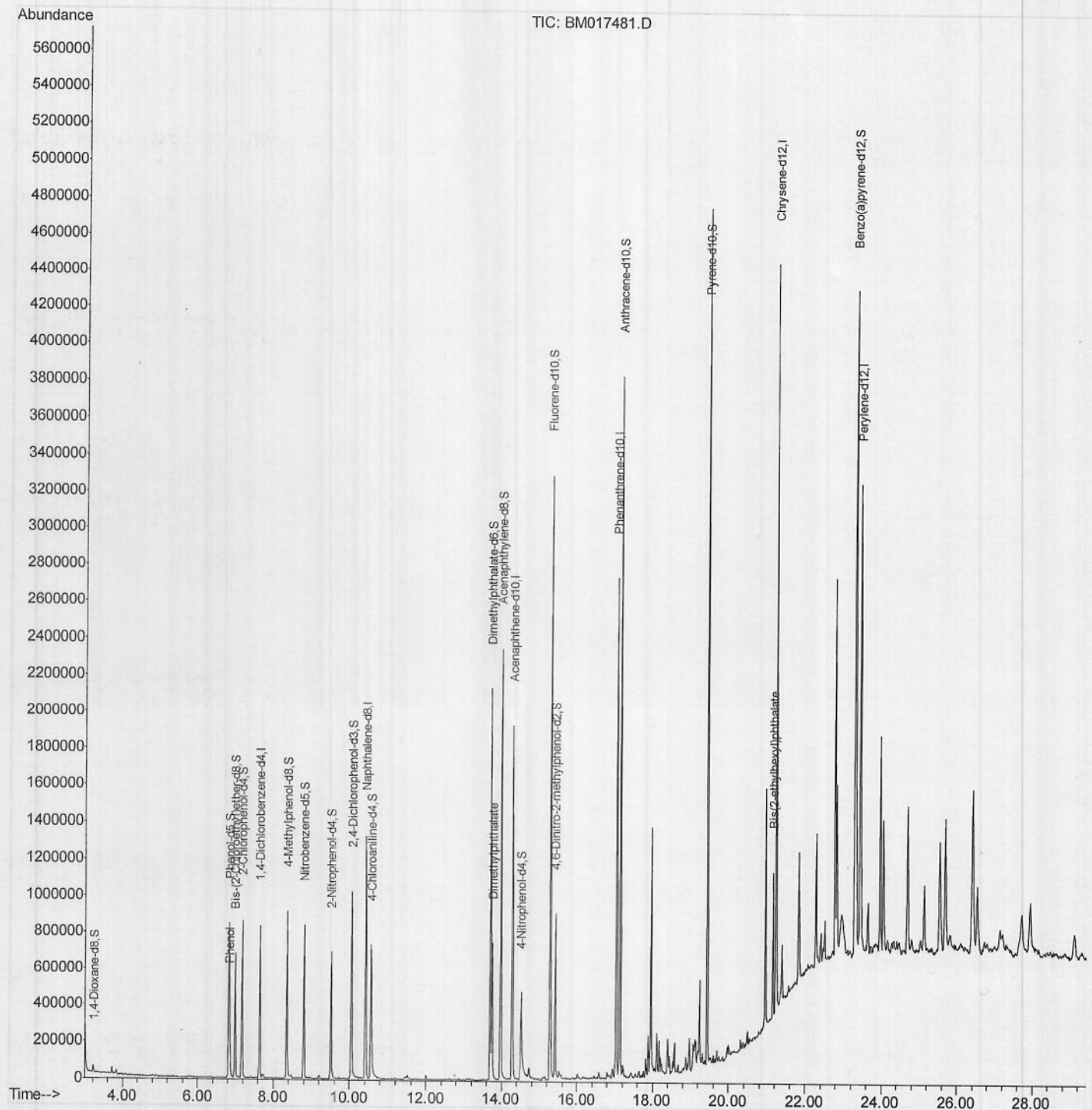
Data File : BM017481.D
Acq On : 02 Nov 2018 00:38
Operator : MJ/SJ
Sample : J5657-16
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
Client Sampled :
A40F1

Manual Integrations
APPROVED

Sohil
11/2/2018 4:12:21 PM

Quant Time: Nov 02 04:20:20 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 02 02:41:22 2018
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110118\
Quantitation Report (Qedit)

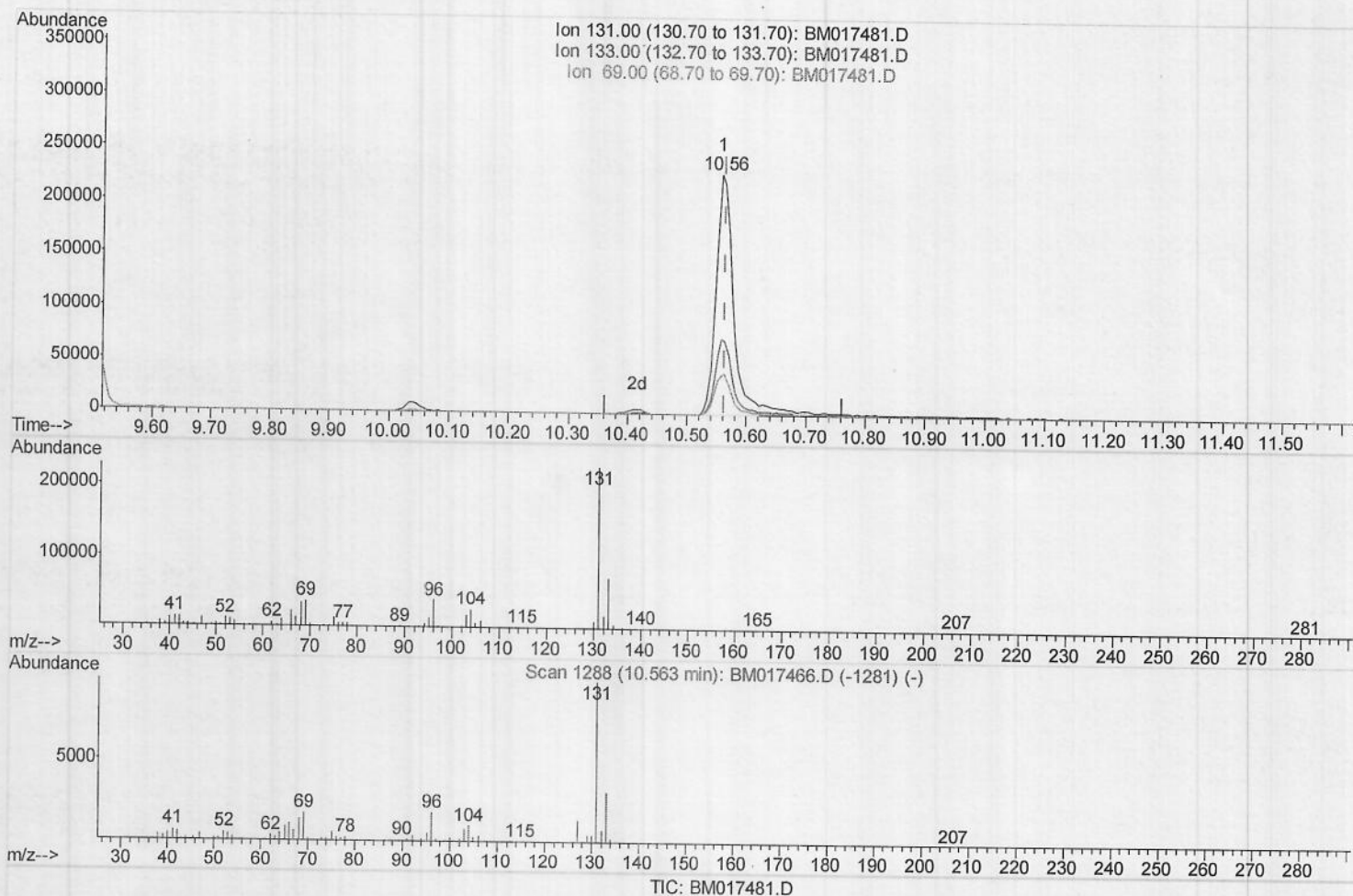
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Quant Time: Nov 02 02:46:36 2018
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(29) 4-Chloroaniline-d4 (S)

10.557min (-0.006) 31.87ng/ul

response 455193

Ion	Exp%	Act%
131.00	100	100
133.00	33.60	31.81
69.00	17.60	16.20
0.00	0.00	0.00

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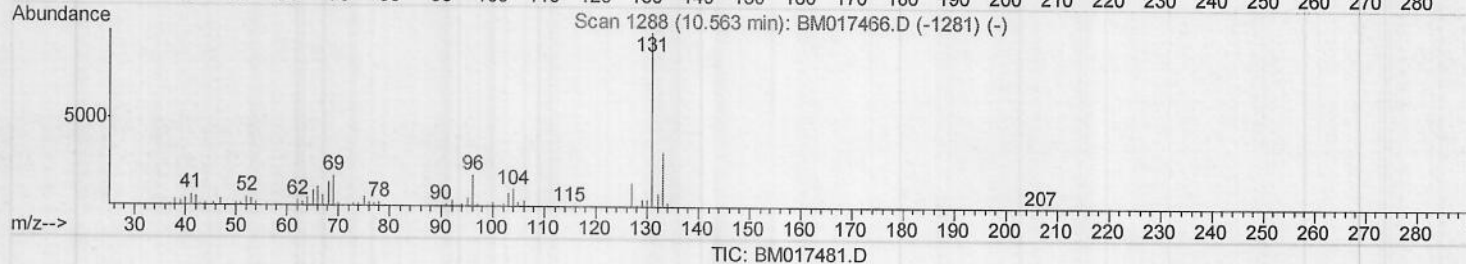
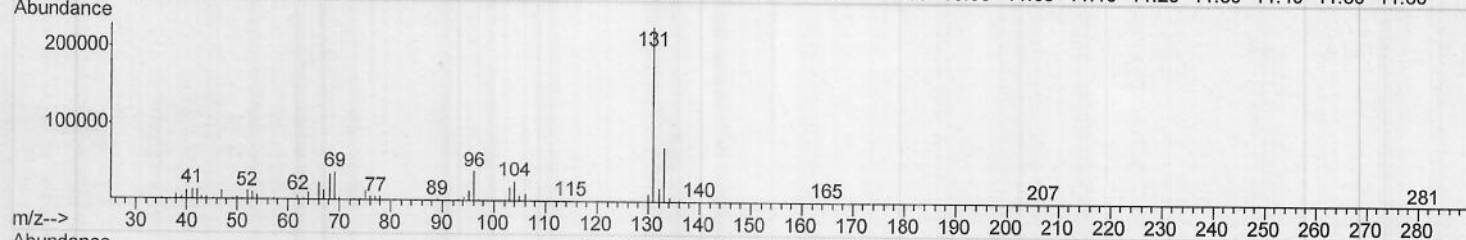
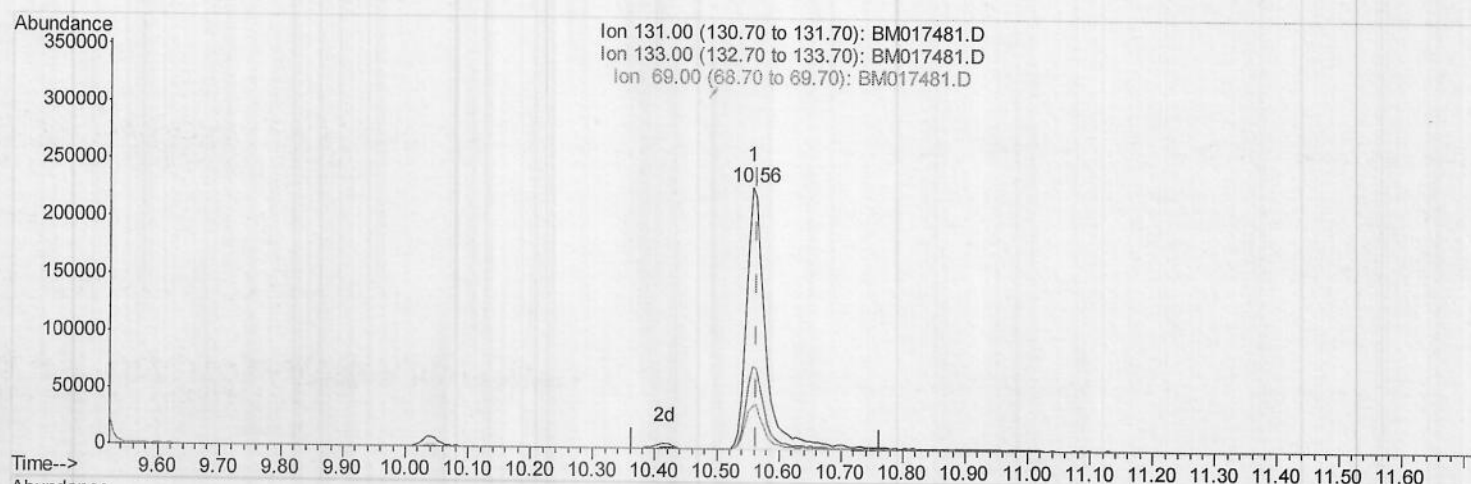
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TIC: BM017481.D

(29) 4-Chloroaniline-d4 (S)

10.557min (-0.006) 34.10ng/ul m J511/06/18

response 486947

Ion	Exp%	Act%
131.00	100	100
133.00	33.60	31.81
69.00	17.60	16.20
0.00	0.00	0.00

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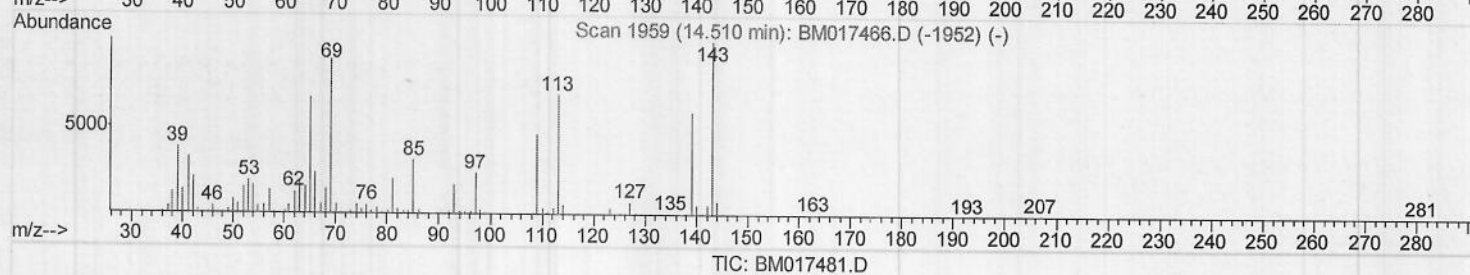
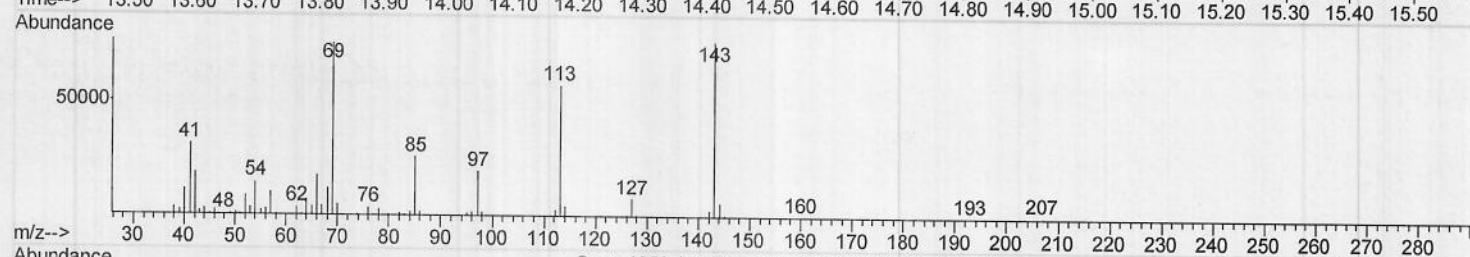
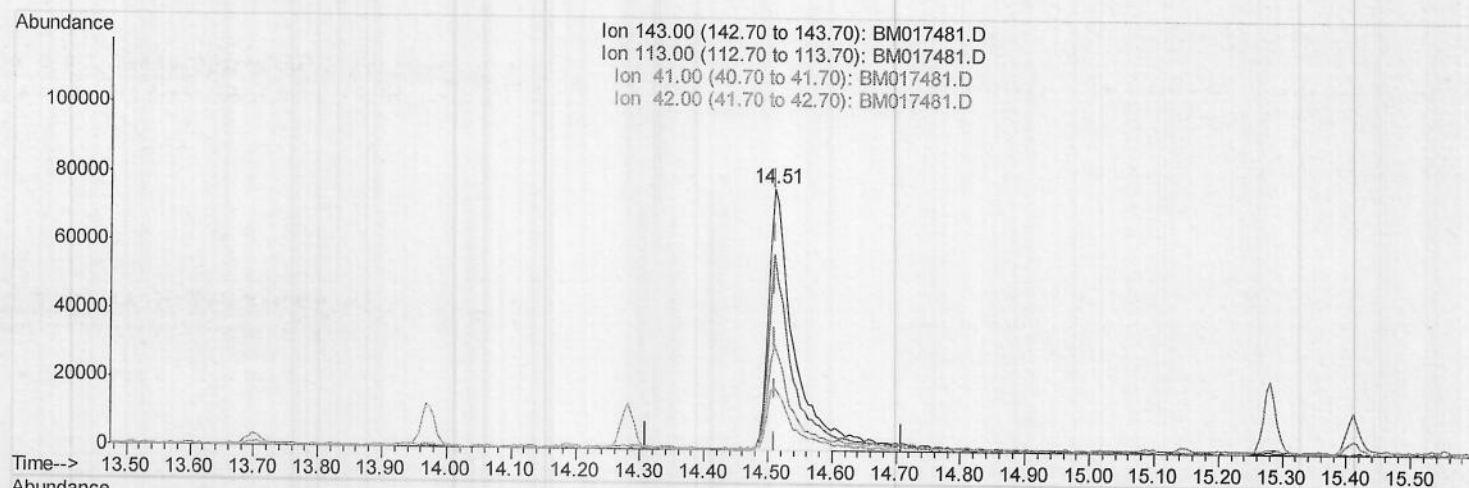
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(52) 4-Nitrophenol-d4 (S)

14.510min (+0.000) 18.52ng/ul

response 195620

Ion	Exp%	Act%
143.00	100	100
113.00	74.20	74.90
41.00	39.30	40.34
42.00	25.10	24.20

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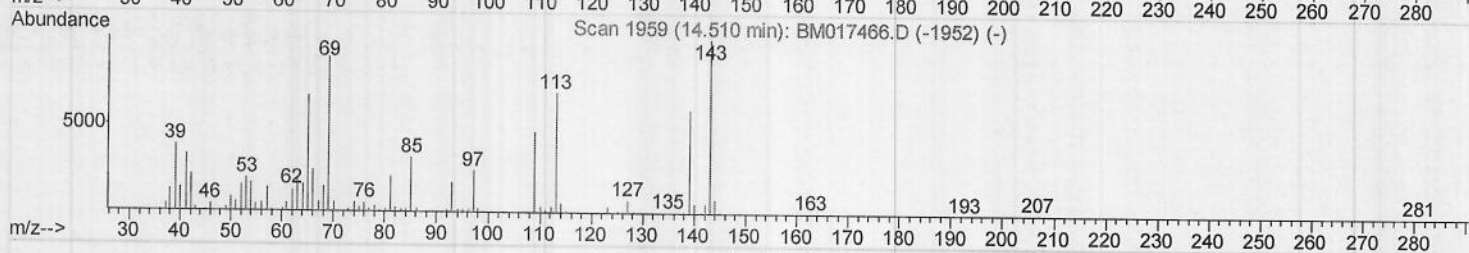
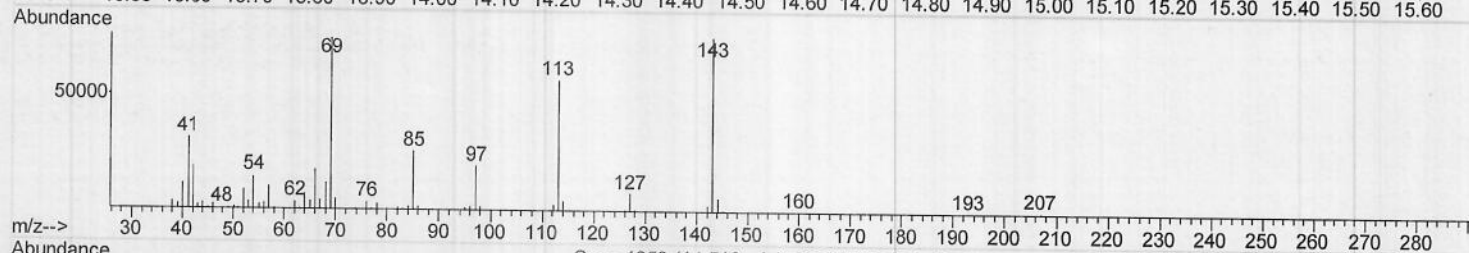
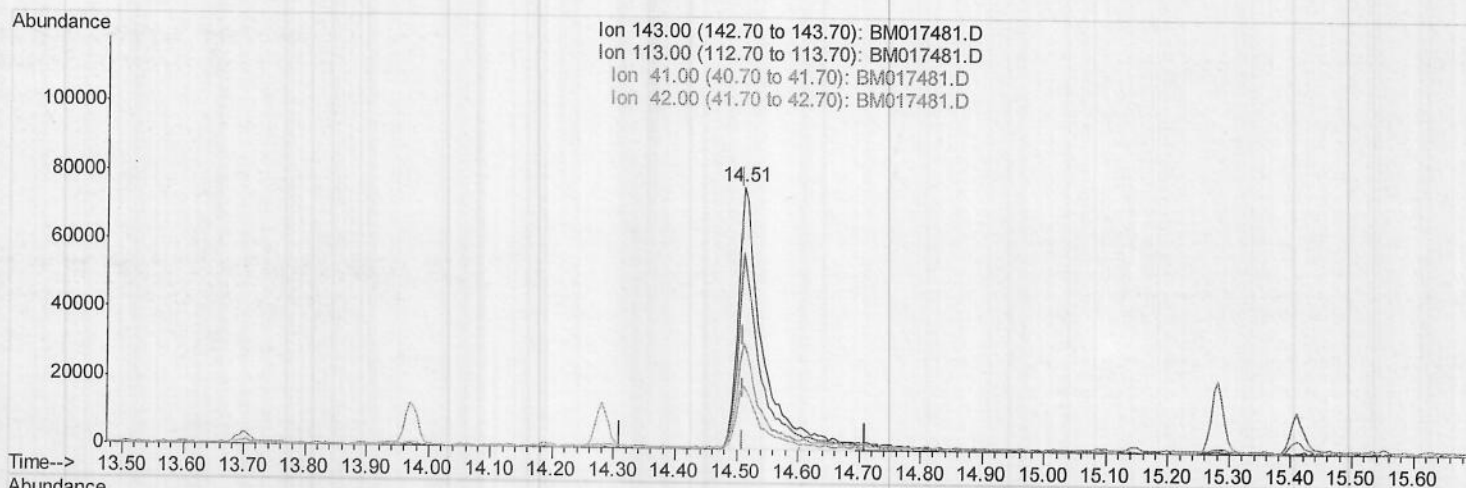
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(52) 4-Nitrophenol-d4 (S)

14.510min (+0.000) 20.53ng/ul m) 551106118

response 216948

Ion	Exp%	Act%
143.00	100	100
113.00	74.20	74.90
41.00	39.30	40.34
42.00	25.10	24.20

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Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	227389	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1074506	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.28	164	659443	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1607204	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	1894803	20.00	ng/ul	0.00
86) Perylene-d12	23.44	264	1874023	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.22	96	19860	4.02	ng/uL	0.00
5) Phenol-d5	6.82	99	492144	23.69	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	294023	24.27	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	397932	24.71	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	352848	21.48	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	207212	24.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	236718	25.13	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	407241	23.28	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	486947m	34.10	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	1456772	25.76	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	1813349	26.41	ng/ul	0.00
52) 4-Nitrophenol-d4	14.51	143	216948m	20.53	ng/ul	0.00
58) Fluorene-d10	15.28	176	1299535	26.46	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.41	200	199874	17.85	ng/ul	0.00
71) Anthracene-d10	17.13	188	2069077	26.01	ng/ul	0.00
79) Pyrene-d10	19.43	212	2429993	27.68	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.30	264	2665261	26.93	ng/ul	0.00
Target Compounds						
6) Phenol	6.85	94	131904	6.395	ng/ul	97
45) Dimethylphthalate	13.75	163	463992	8.469	ng/ul	100
84) Bis(2-ethylhexyl)phthalate	21.16	149	262235	4.085	ng/ul	96

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

{ 5/11/05/18