

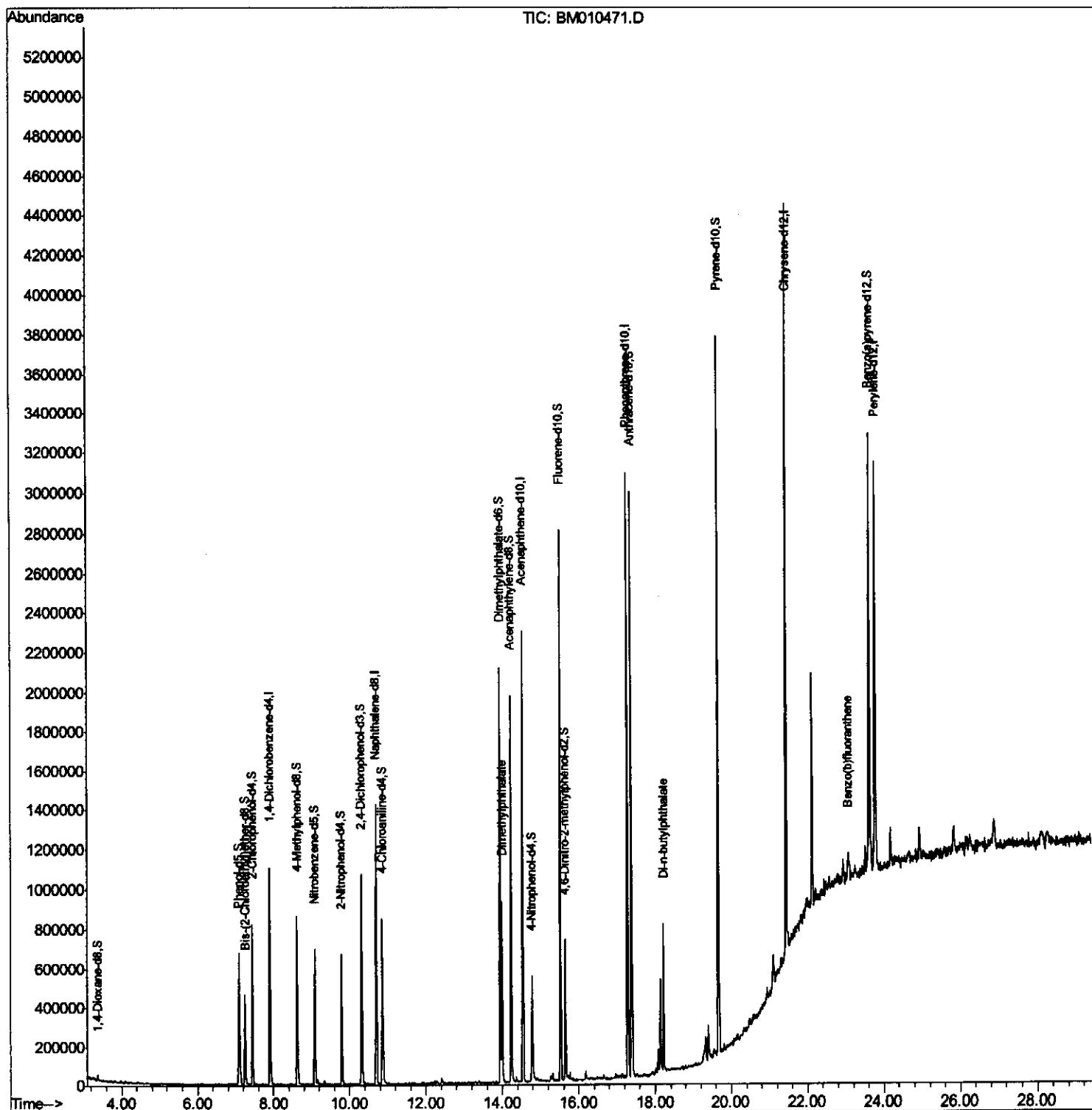
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060917\
Data File : BM010471.D
Acq On : 10 Jun 2017 01:39
Operator : SJ/MA
Sample : I3520-18
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F5B65

Manual Integrations
APPROVED

mohammad
6/12/2017 11:18:28 AM

Quant Time: Jun 10 04:08:05 2017
Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jun 10 02:32:05 2017
Response via : Initial Calibration



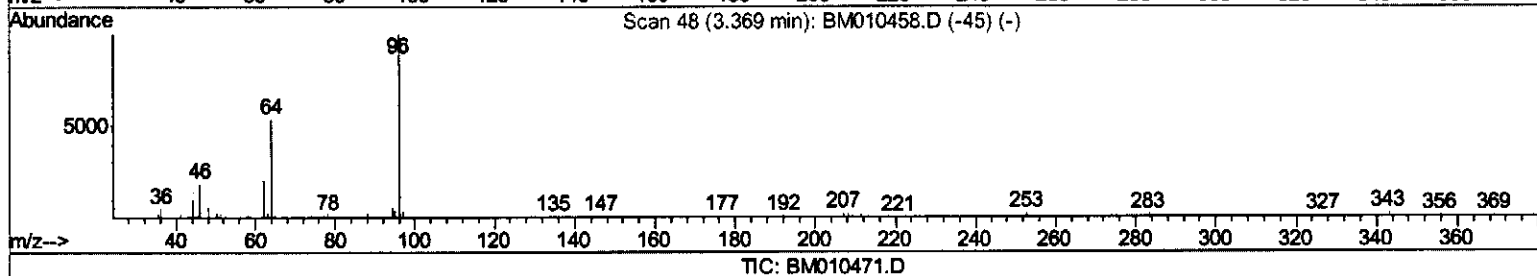
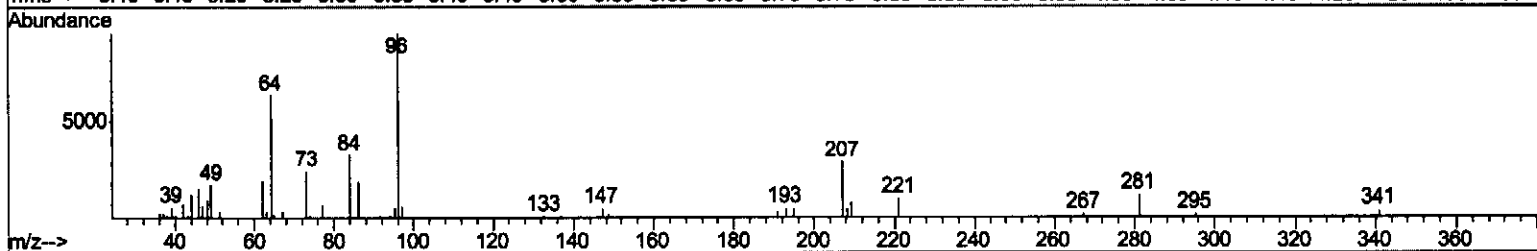
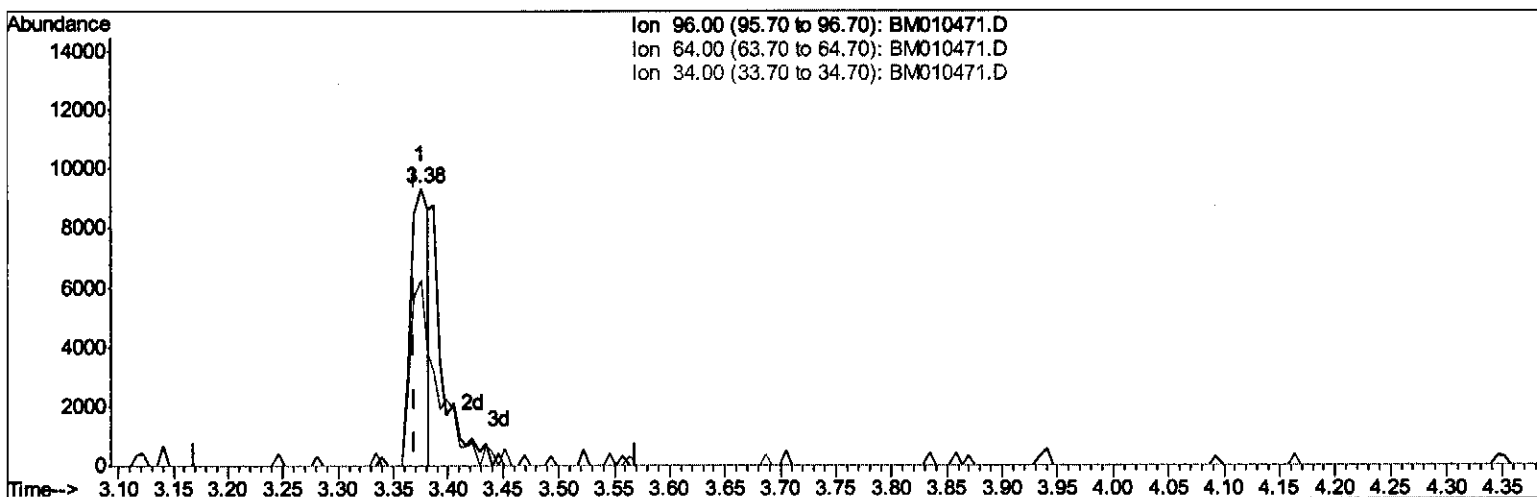
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(3) 1,4-Dioxane-d8 (S)

3.375min (+0.006) 1.65ng/uL

response 10396

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	79.53
34.00	0.00	0.00
0.00	0.00	0.00

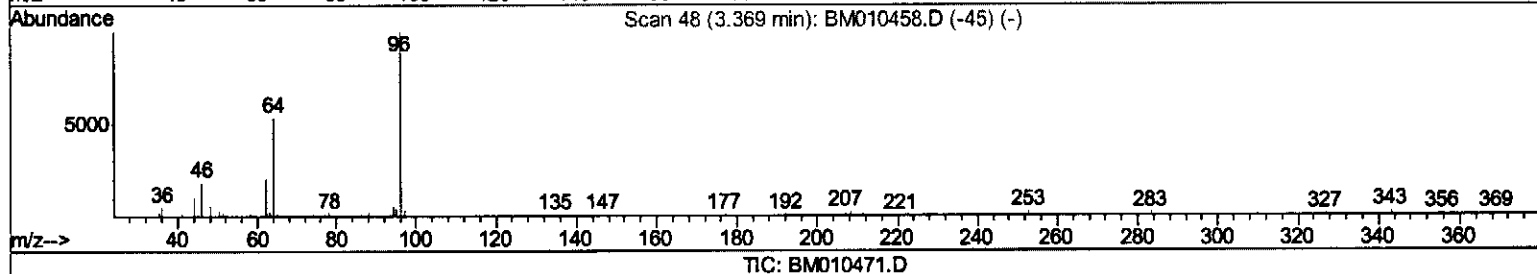
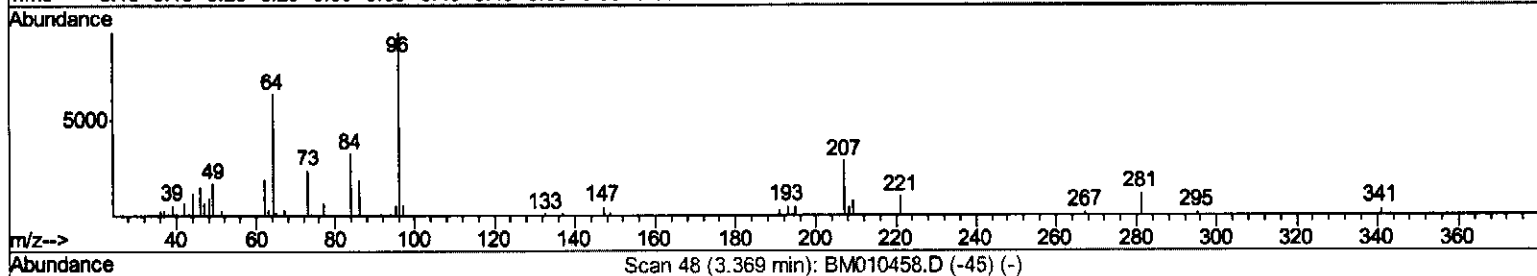
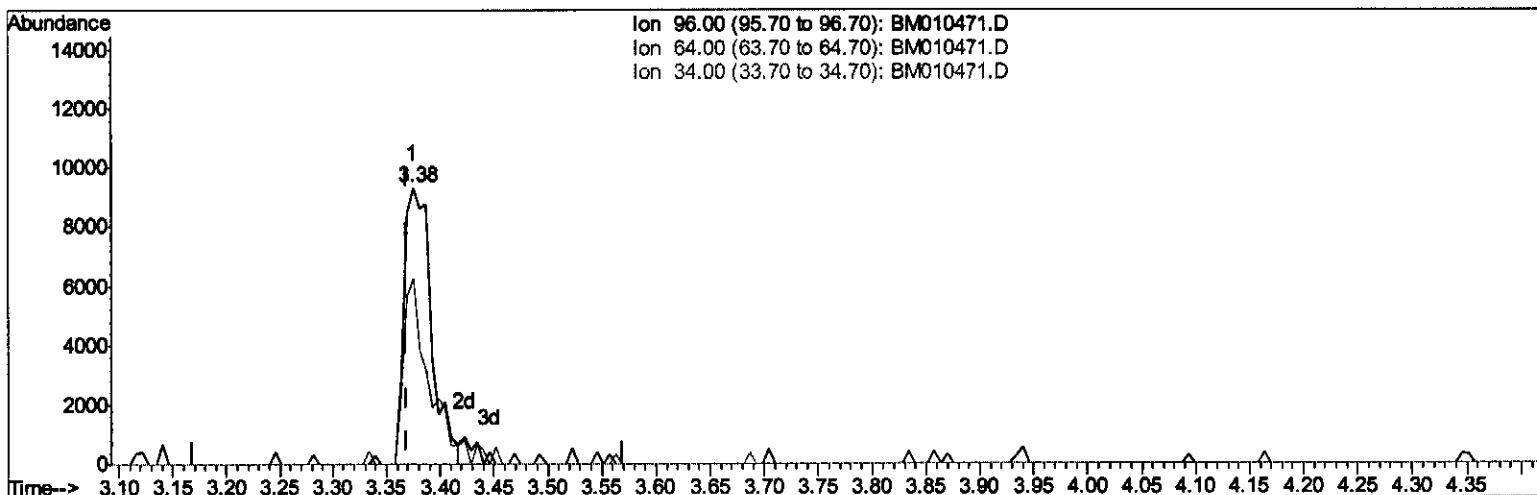
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(3) 1,4-Dioxane-d8 (S)

3.375min (+0.006) 2.64ng/uL m

response 16629

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	49.72#
34.00	0.00	0.00
0.00	0.00	0.00

05/06/19/17

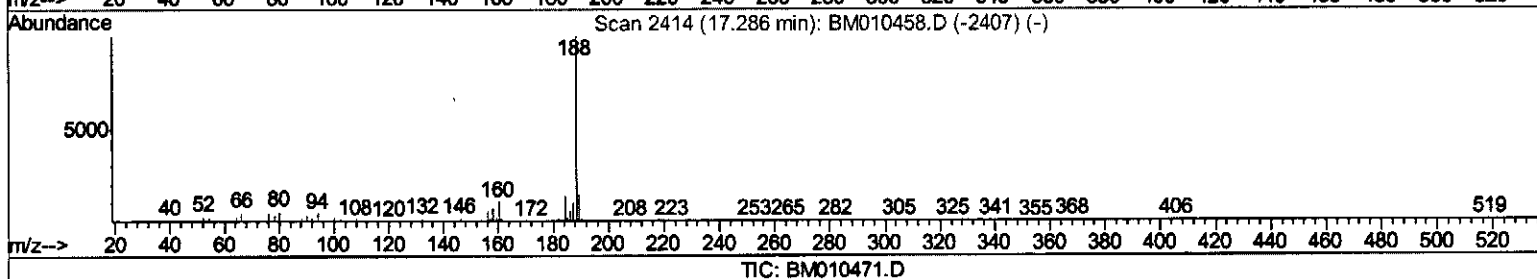
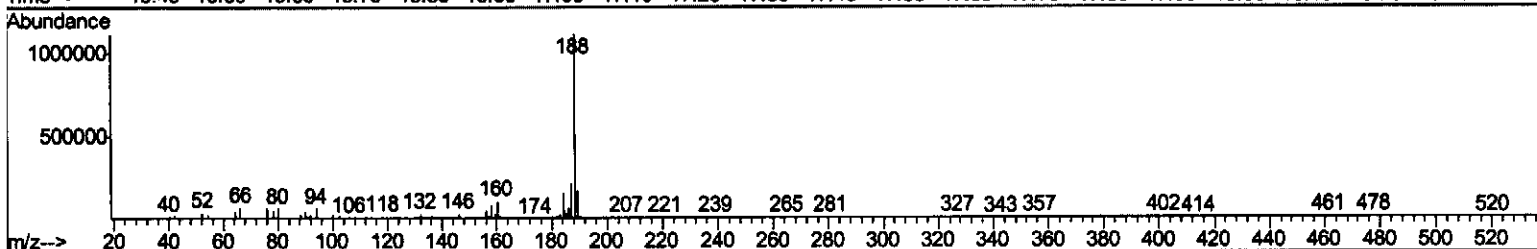
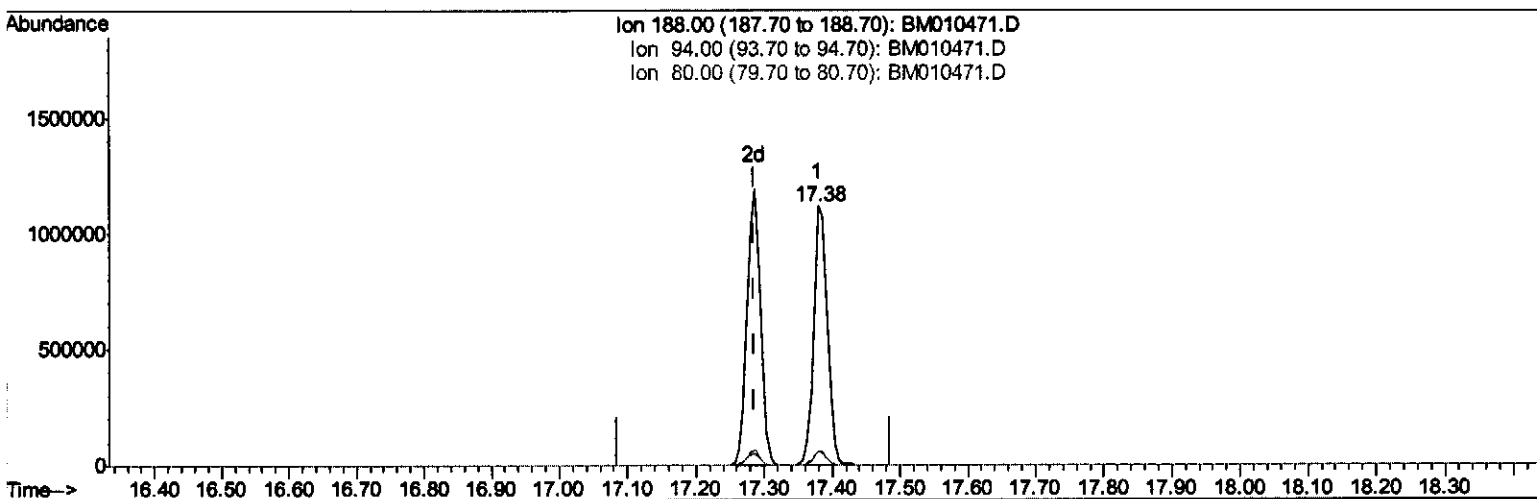
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Misc :
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Manual Integrations
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(61) Phenanthrene-d10 (I)

17.380min (+0.094) 20.00ng/ul

response 1577249

Ion	Exp%	Act%
188.00	100	100
94.00	5.50	5.23
80.00	6.60	5.36
0.00	0.00	0.00

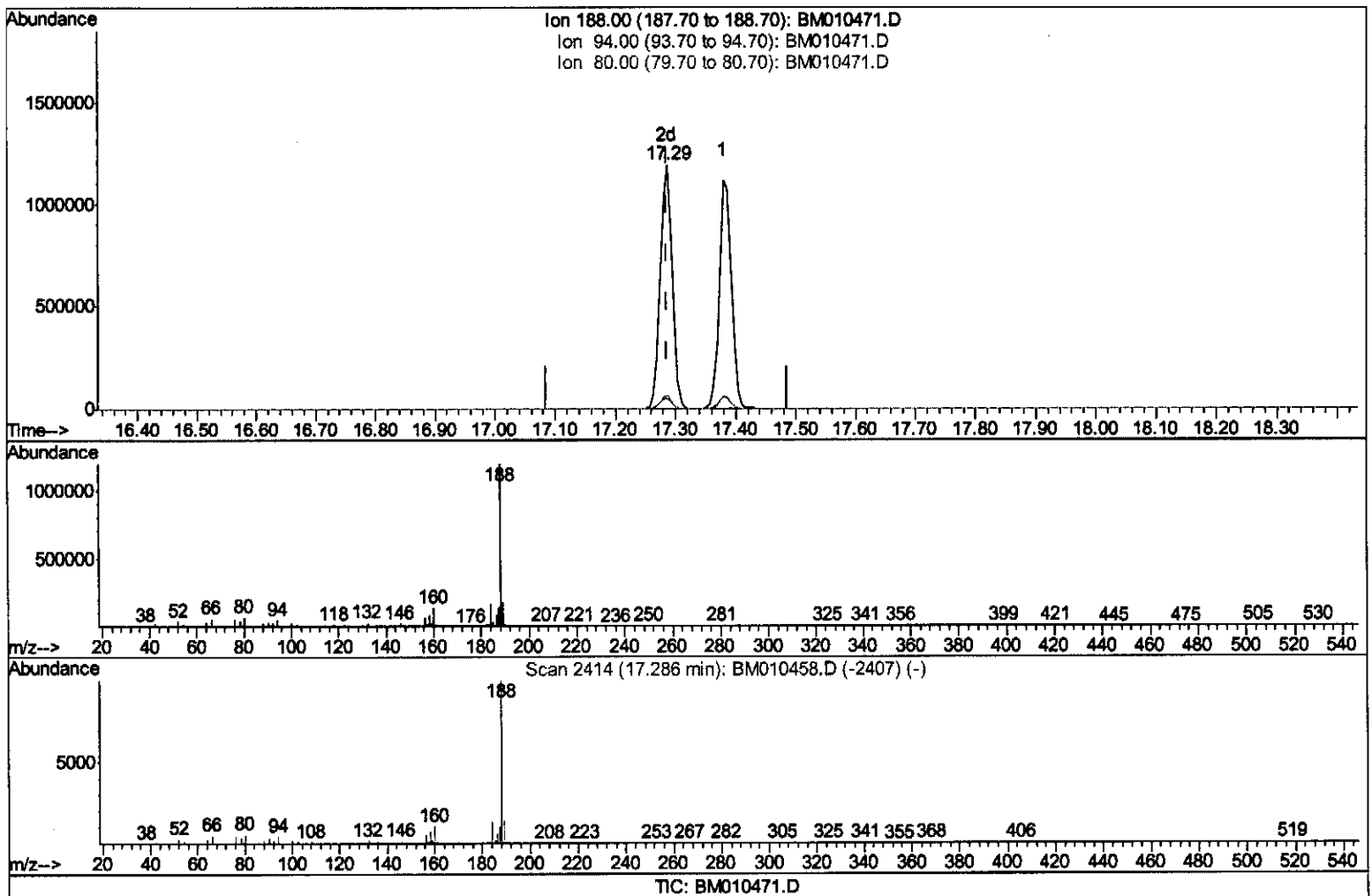
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Operator : SJ/MA
Sample : I3520-18
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
F5B65

Manual Integrations
APPROVED

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



(61) Phenanthrene-d10 (I)

17.286min (+0.000) 20.00ng/ul m

response 1637246

Ion	Exp%	Act%
188.00	100	100
94.00	5.50	3.99#
80.00	6.60	5.59
0.00	0.00	0.00

SJ 06/19/17

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 Sample : I3520-18
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
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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	7.90	152	257984	20.00	ng/ul	0.00
18) Naphthalene-d8	10.70	136	947768	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.54	164	652983	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.29	188	1637246m	20.00	ng/ul	0.00
75) Chrysene-d12	21.45	240	1735384	20.00	ng/ul	0.00
83) Perylene-d12	23.79	264	1605453	20.00	ng/ul	0.00

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System Monitoring Compounds

3) 1,4-Dioxane-d8	3.38	96	16629m	2.64	ng/uL	0.00
5) Phenol-d5	7.09	99	314829	16.10	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.25	67	175626	15.99	ng/ul	0.00
9) 2-Chlorophenol-d4	7.44	132	278439	17.87	ng/ul	0.00
13) 4-Methylphenol-d8	8.62	113	262955	16.67	ng/ul	0.00
19) Nitrobenzene-d5	9.09	128	126190	18.18	ng/ul	0.00
22) 2-Nitrophenol-d4	9.80	143	156845	19.50	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.33	165	333686	20.20	ng/ul	0.00
29) 4-Chloroaniline-d4	10.87	131	341323	21.41	ng/ul	0.00
43) Dimethylphthalate-d6	13.95	166	1143932	21.08	ng/ul	0.00
46) Acenaphthylene-d8	14.23	160	1251152	19.77	ng/ul	0.00
51) 4-Nitrophenol-d4	14.78	143	106743	13.15	ng/ul	0.00
57) Fluorene-d10	15.53	176	965578	20.25	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.66	200	166115	14.32	ng/ul	0.00
70) Anthracene-d10	17.38	188	1577249	19.80	ng/ul	0.00
76) Pyrene-d10	19.67	212	1828625	25.48	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.64	264	1511847	20.23	ng/ul	0.00

SJ 06/19/17

Target Compounds

						Qvalue
44) Dimethylphthalate	14.00	163	480817	9.012	ng/ul	99
73) Di-n-butylphthalate	18.22	149	461886	5.298	ng/ul#	94
85) Benzo(b)fluoranthene	23.08	252	122624	1.313	ng/ul#	90

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