

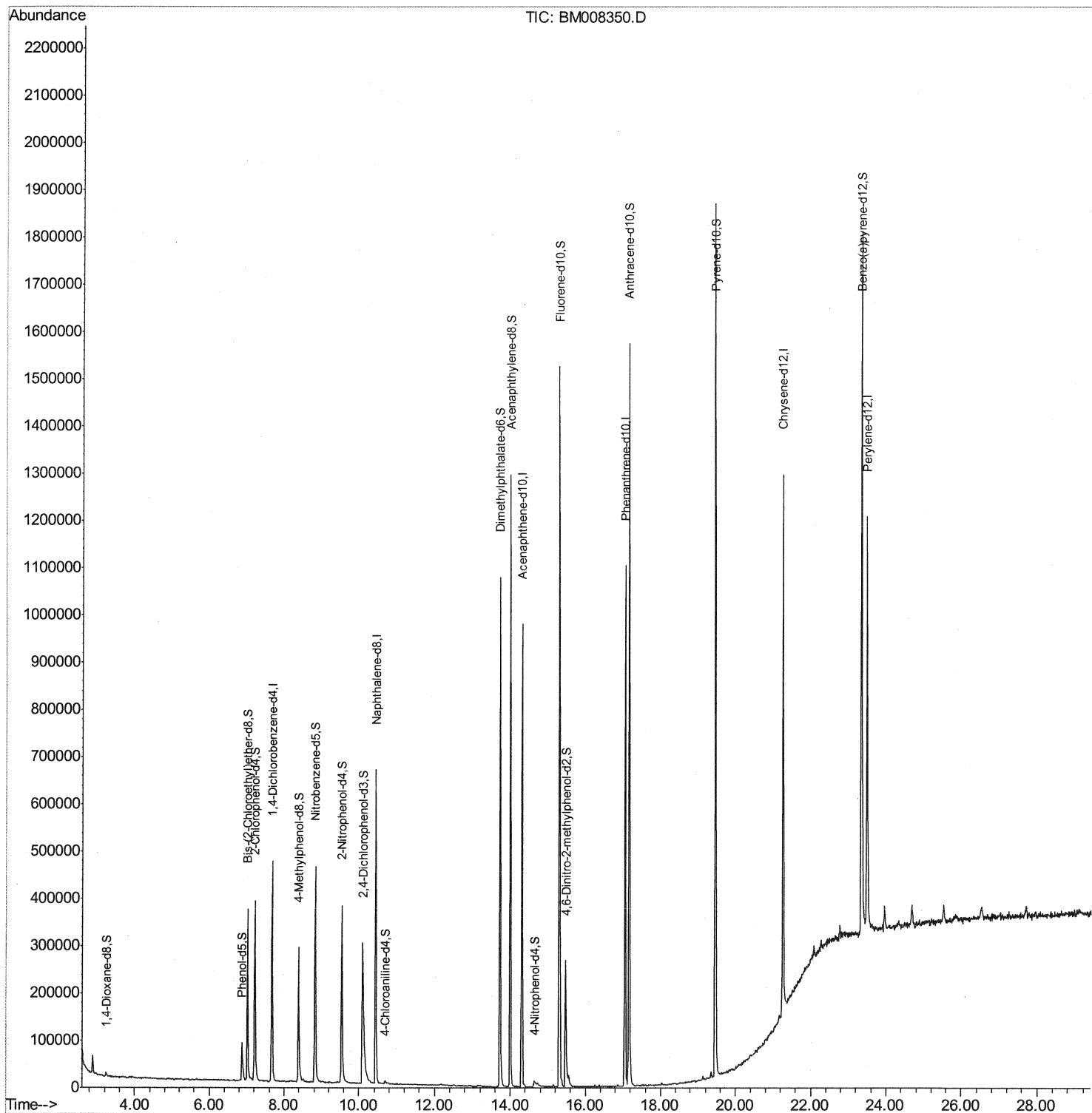
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121616\
Data File : BM008350.D
Acq On : 15 Dec 2016 19:25
Operator : UM/SJ
Sample : H6039-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DA8T5

Manual Integrations
APPROVED

sohil
12/16/2016 6:21:01 PM

Quant Time: Dec 16 03:06:11 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121416.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 16 02:33:30 2016
Response via : Initial Calibration



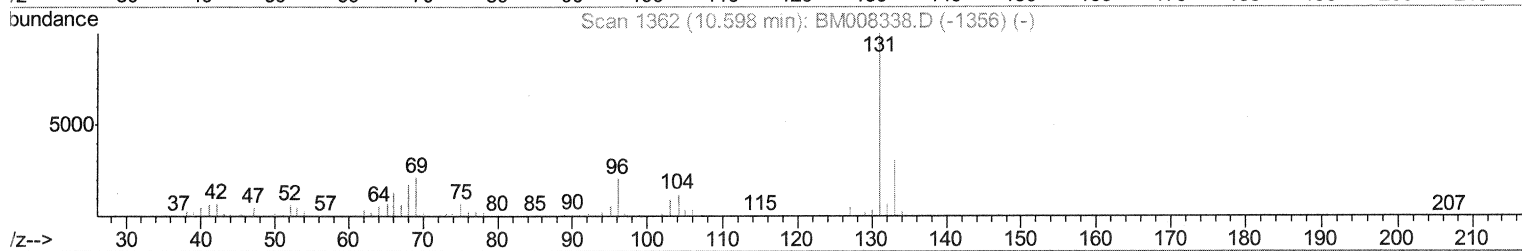
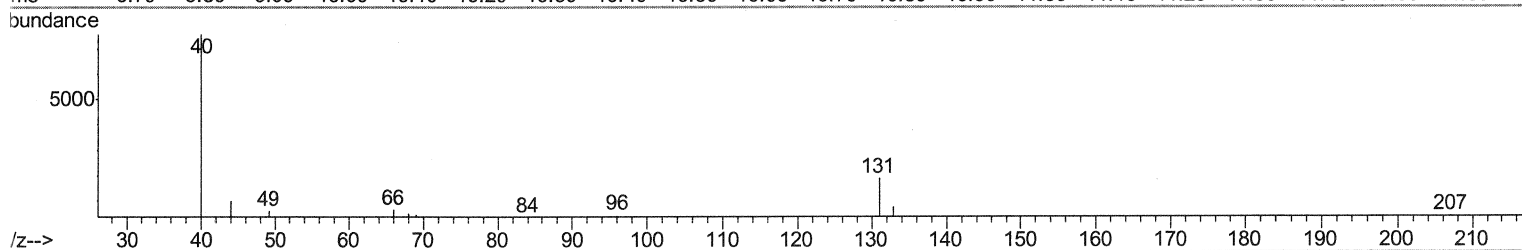
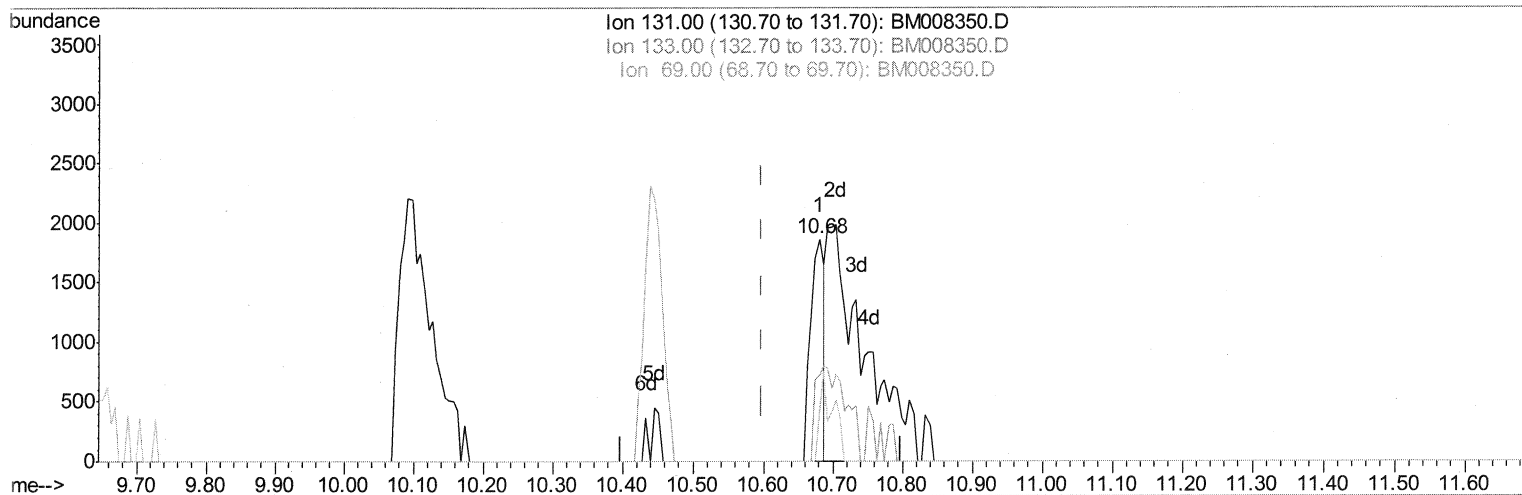
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ClientSampled :
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Manual Integrations
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Quant Time: Dec 16 02:37:53 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121416.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 16 02:33:30 2016
Response via : Initial Calibration



TIC: BM008350.D

(29) 4-Chloroaniline-d4 (S)

10.681min (+0.082) 0.28ng/ul

response 2554

Ion	Exp%	Act%
131.00	100	100
133.00	34.20	38.86
69.00	21.70	24.21
0.00	0.00	0.00

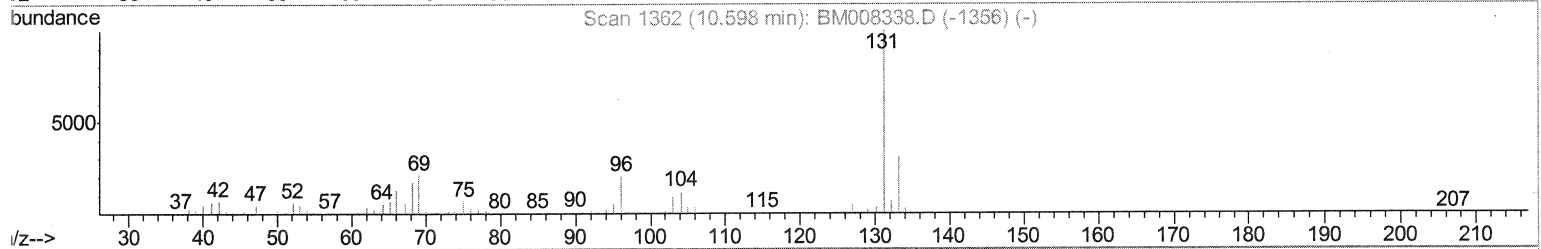
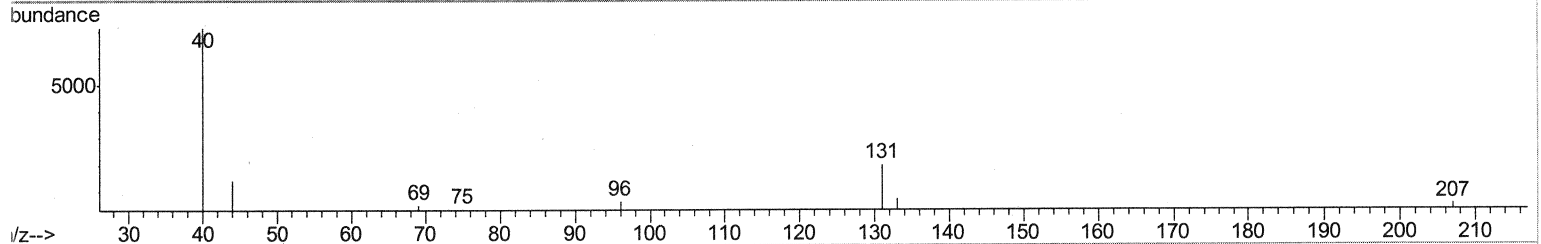
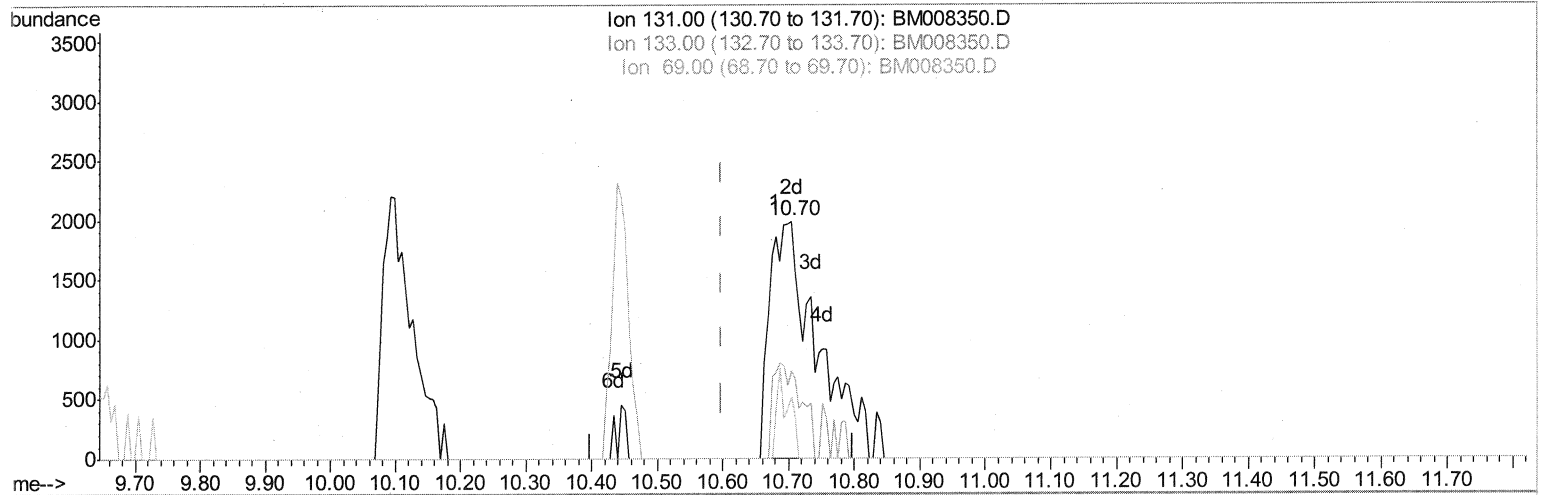
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Instrument :
BNA_M
ClientSampleId :
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Manual Integrations
APPROVED

sohil
12/16/2016 6:21:01 PM

Quant Time: Dec 16 02:37:53 2016
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Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 16 02:33:30 2016
Response via : Initial Calibration



TIC: BM008350.D

(29) 4-Chloroaniline-d4 (S)

10.704min (+0.106) 1.09ng/ul m 57 12/20/20

response 9939

Ion	Exp%	Act%
131.00	100	100
133.00	34.20	36.85
69.00	21.70	25.84
0.00	0.00	0.00

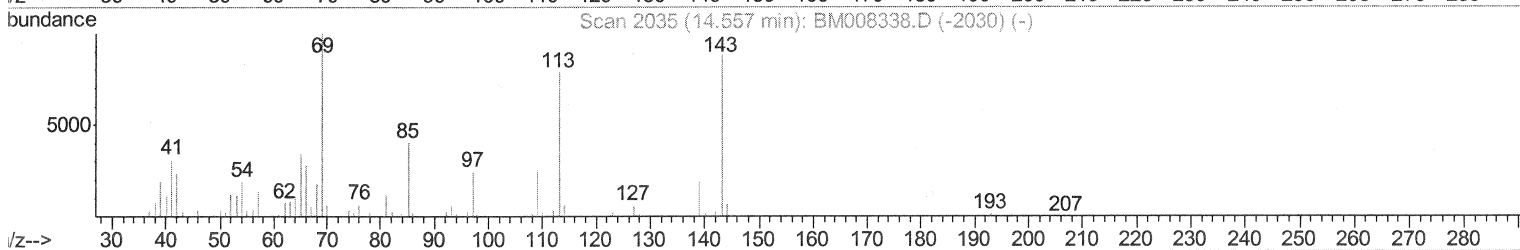
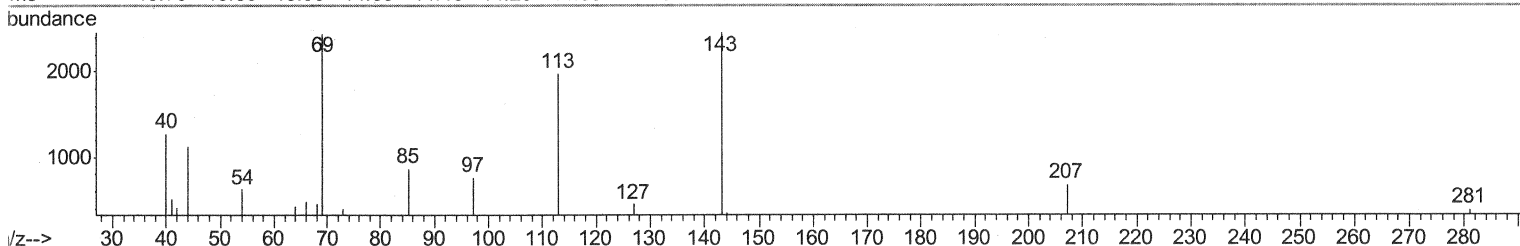
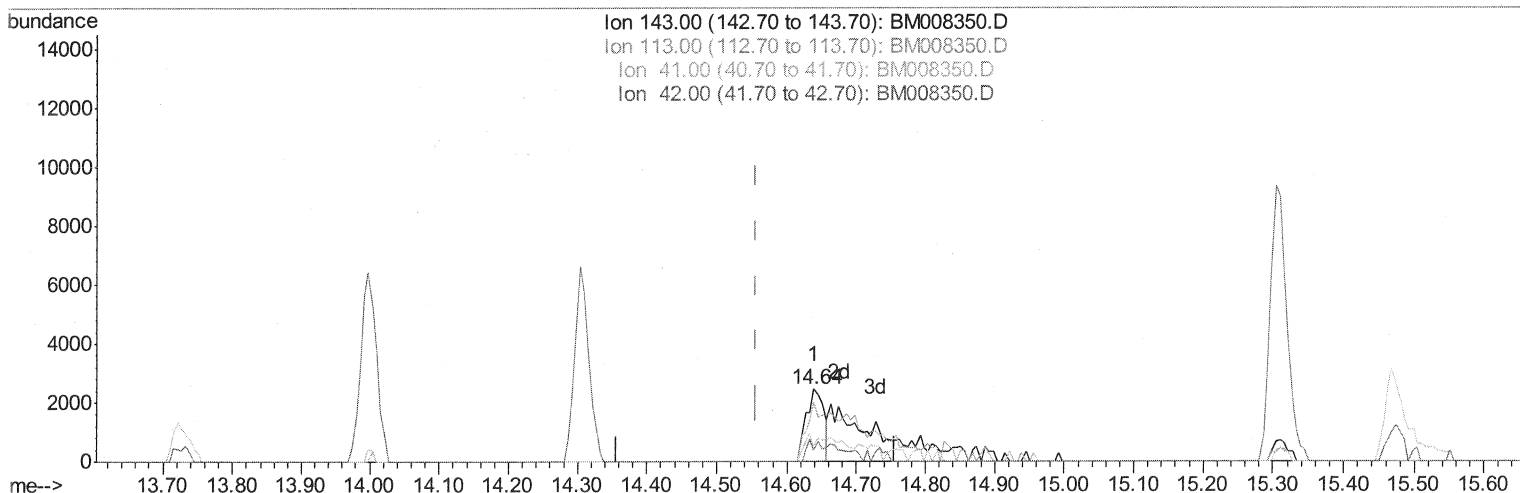
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121616\
Data File : BM008350.D
Acq On : 15 Dec 2016 19:25
Operator : UM/SJ
Sample : H6039-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DA8T5

Manual Integrations
APPROVED

sohil
12/16/2016 6:21:01 PM

Quant Time: Dec 16 02:37:53 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121416.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 16 02:33:30 2016
Response via : Initial Calibration



TIC: BM008350.D

(51) 4-Nitrophenol-d4 (S)

14.639min (+0.082) 1.29ng/ul

response 4379

Ion	Exp%	Act%
143.00	100	100
113.00	101.20	80.66#
41.00	39.40	20.97#
42.00	27.40	16.41#

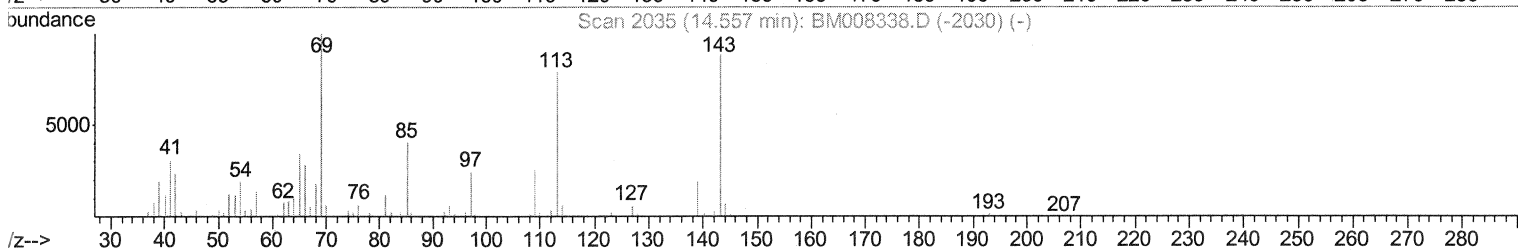
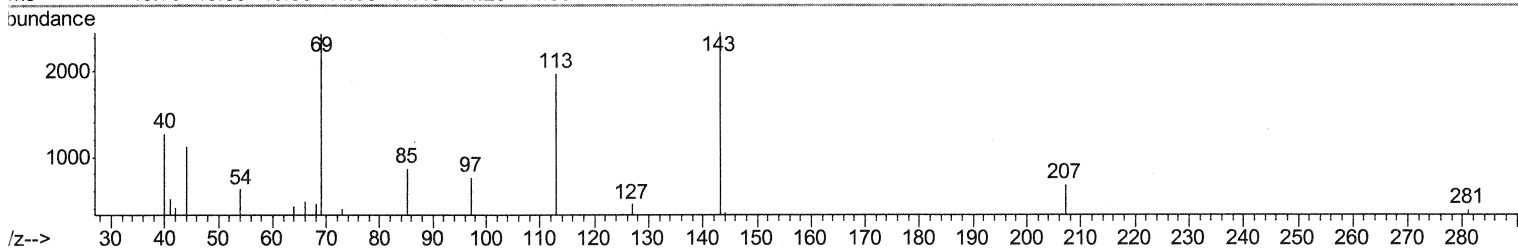
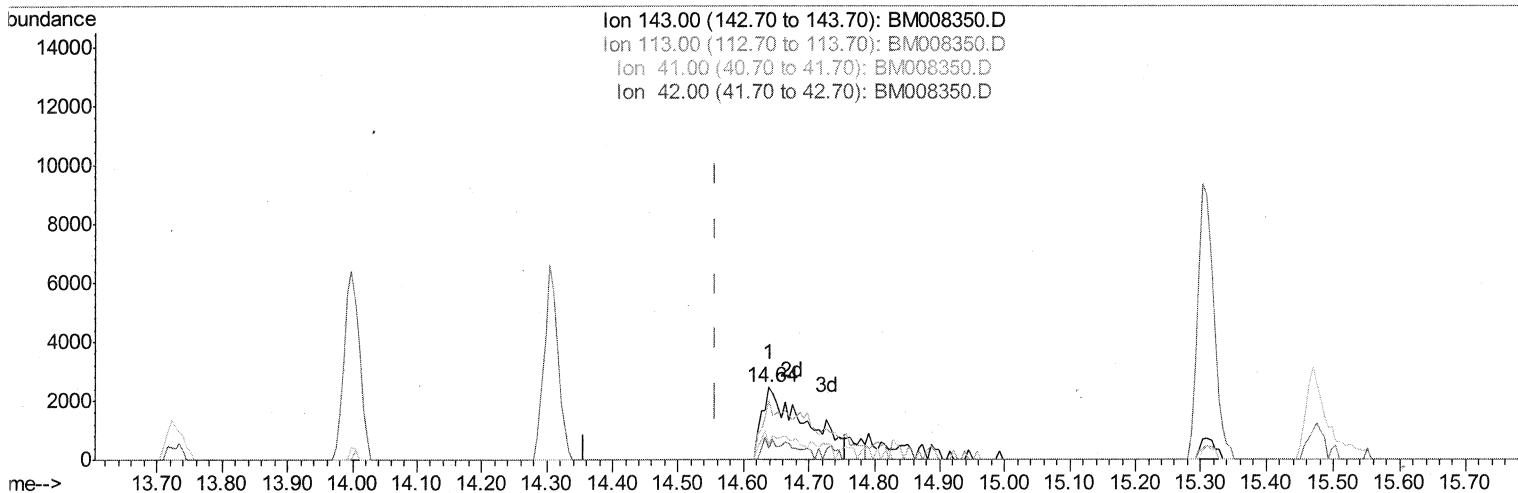
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121616\
Data File : BM008350.D
Acq On : 15 Dec 2016 19:25
Operator : UM/SJ
Sample : H6039-21
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DA8T5

Manual Integrations
APPROVED

sohil
12/16/2016 6:21:01 PM

Quant Time: Dec 16 02:37:53 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM12141
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 16 02:33:30 2016
Response via : Initial Calibration



TIC: BM008350.D

(51) 4-Nitrophenol-d4 (S)

14.639min (+0.082) 3.64ng/ul m

SJ 12/20/16

response 12347

Ion	Exp%	Act%
143.00	100	100
113.00	101.20	80.66#
41.00	39.40	20.97#
42.00	27.40	16.41#

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121616\
 Data File : BM008350.D
 Acq On : 15 Dec 2016 19:25
 Operator : UM/SJ
 Sample : H6039-21
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA8T5

Manual Integrations
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sohil
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 QLast Update : Fri Dec 16 02:33:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	132397	20.00	ng/ul	0.00
18) Naphthalene-d8	10.44	136	541417	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.31	164	345934	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	625496	20.00	ng/ul	0.00
75) Chrysene-d12	21.26	240	697360	20.00	ng/ul	0.00
83) Perylene-d12	23.49	264	717609	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.24	96	4055	1.44	ng/uL	0.00
5) Phenol-d5	6.87	99	72315	7.04	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.01	67	174602	29.58	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	189025	24.46	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	127082	15.97	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	113581	29.03	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	128078	29.54	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	203911	25.66	ng/ul	0.01
29) 4-Chloroaniline-d4	10.70	131	9939m	1.09	ng/ul	0.11
43) Dimethylphthalate-d6	13.73	166	768896	33.44	ng/ul	0.00
46) Acenaphthylene-d8	14.00	160	921126	33.89	ng/ul	0.00
51) 4-Nitrophenol-d4	14.64	143	12347m	3.64	ng/ul	0.08
57) Fluorene-d10	15.31	176	653569	32.70	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.47	200	88066	23.65	ng/ul	0.01
70) Anthracene-d10	17.16	188	990853	36.36	ng/ul	0.00
76) Pyrene-d10	19.46	212	1095252	40.67	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.34	264	1076787	35.73	ng/ul	0.00

SJ 12/20/16

SJ 12/20/16

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