

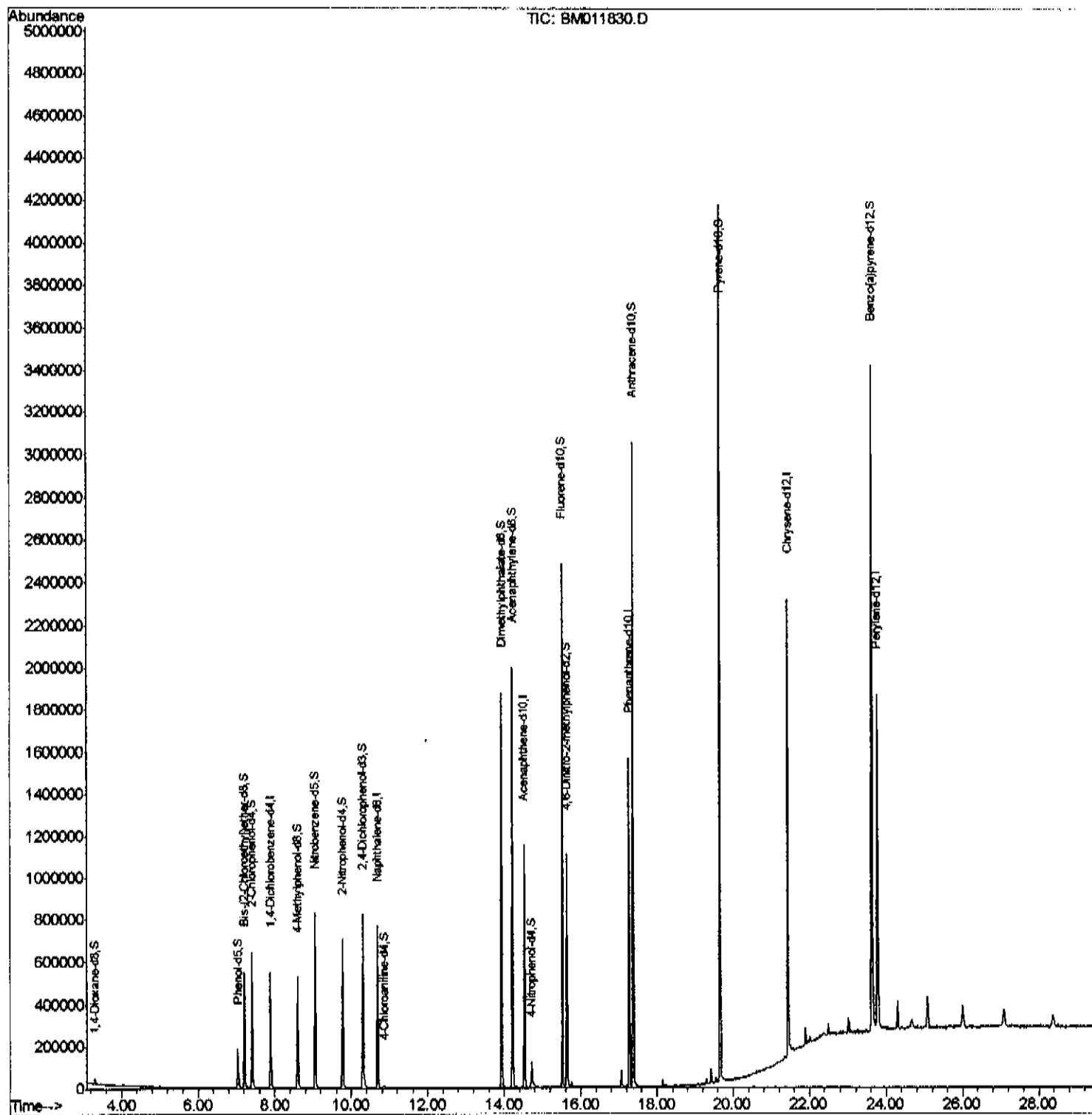
Data Path : Z:\HPCHEM1\BNA M\DATA\BM100117\
Data File : BM011830.D
Acq On : 02 Oct 2017 03:57
Operator : SJ/JU
Sample : I5458-09
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
C0AH2

Manual Integrations
APPROVED

Sohil
10/2/2017 5:52:21 PM

Quant Time: Oct 02 05:54:04 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM092717.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 02 04:55:04 2017
Response via : Initial Calibration



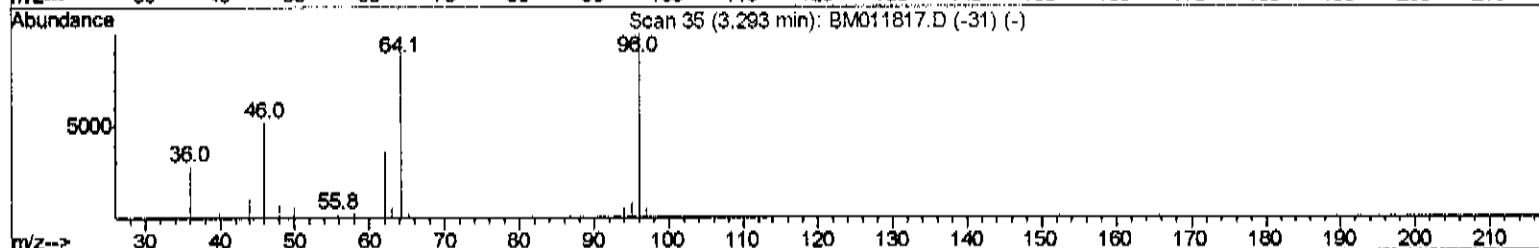
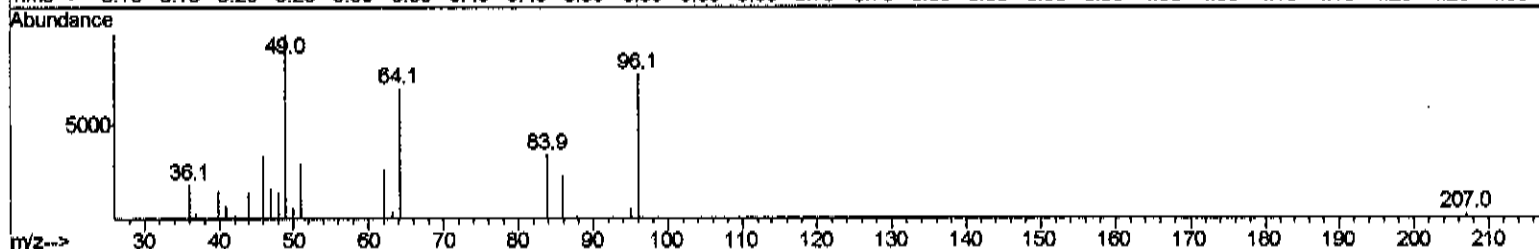
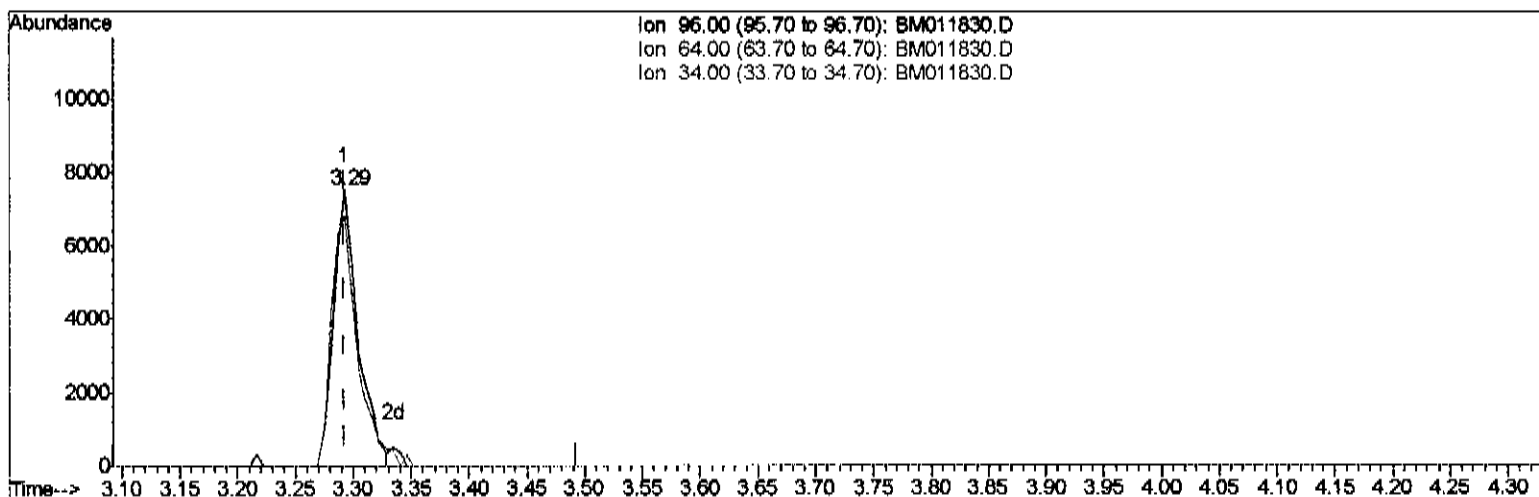
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Quant Time: Oct 02 05:13:26 2017
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TIC: BM011830.D

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

3.293min (-0.000) 4.20ng/uL

response 10987

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	97.56#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

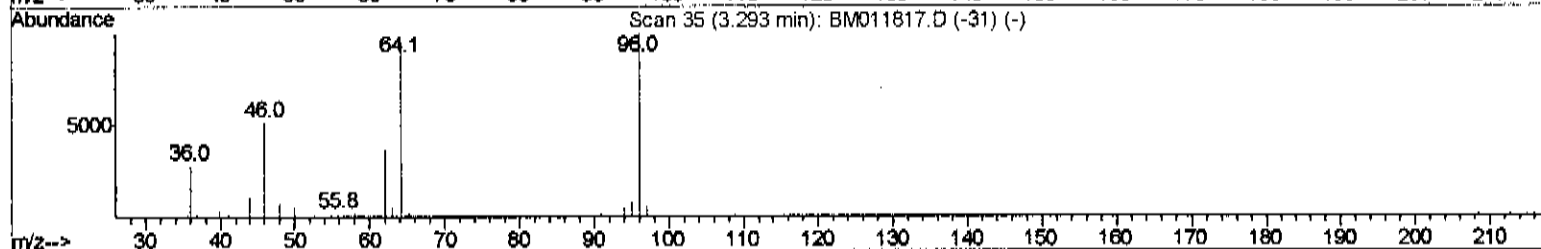
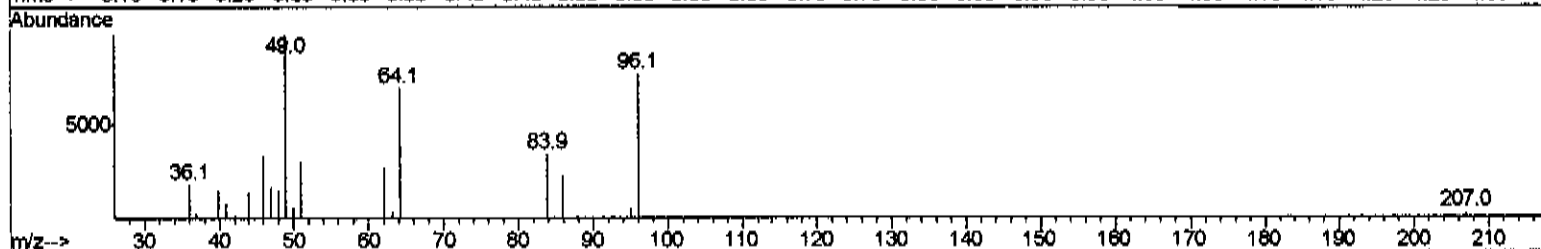
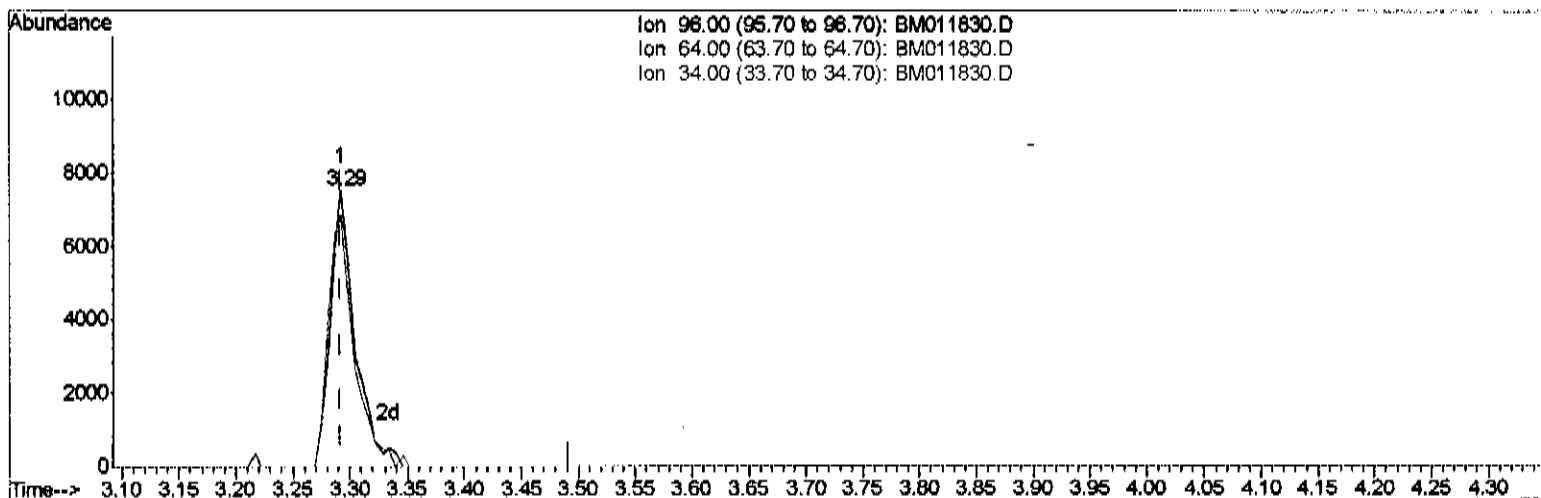
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Manual Integrations
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TIC: BM011830.D

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

3.293min (-0.000) 4.32ng/uL m

response 11299

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	94.87#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

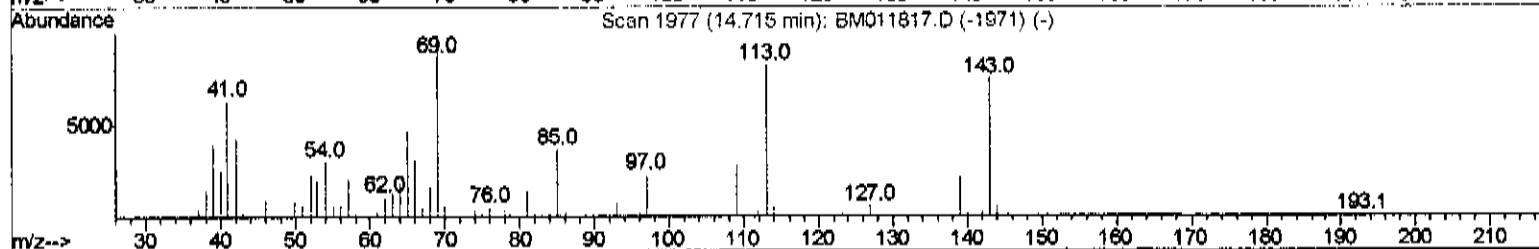
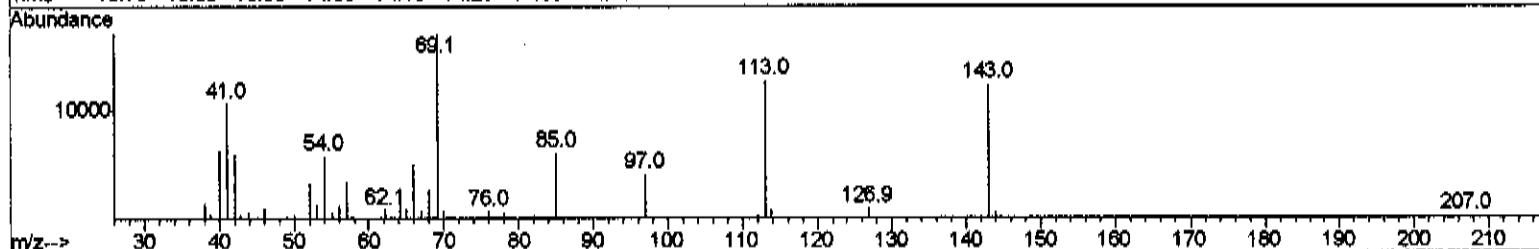
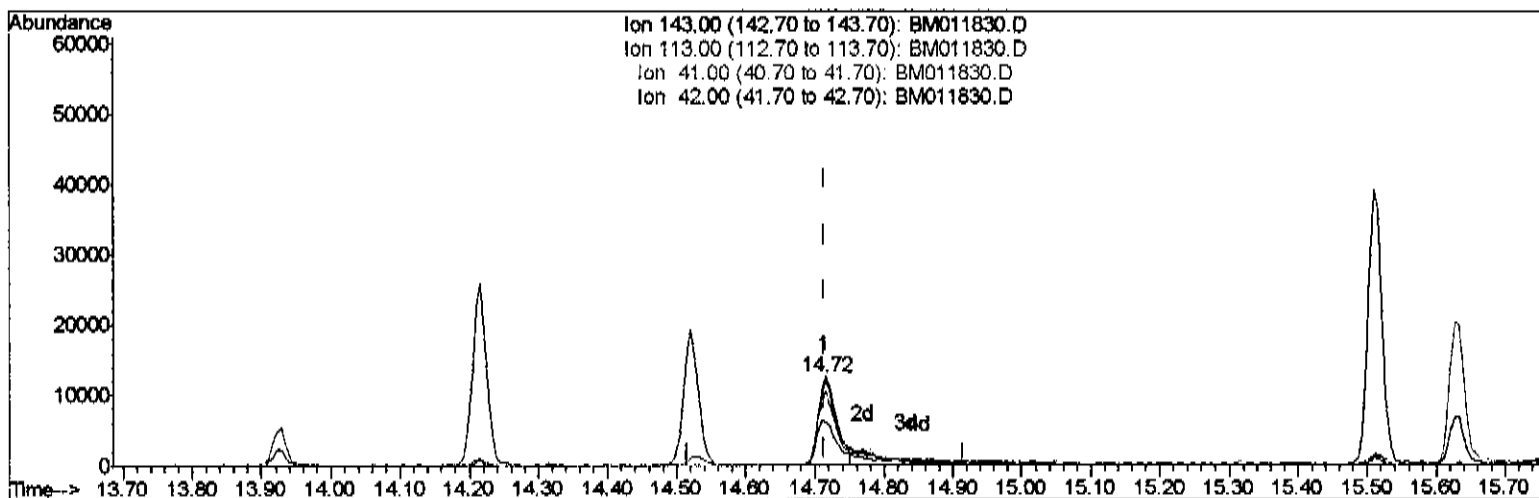
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Manual Integrations
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Sohil
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TIC: BM011830.D

(51) 4-Nitrophenol-d4 (S)

14.715min (-0.000) 4.50ng/ul

response 21674

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	102.92
41.00	34.80	86.72#
42.00	20.50	49.65#

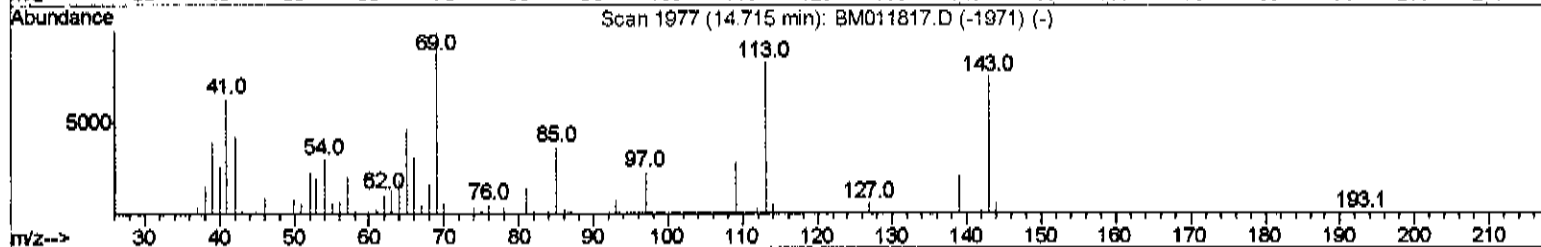
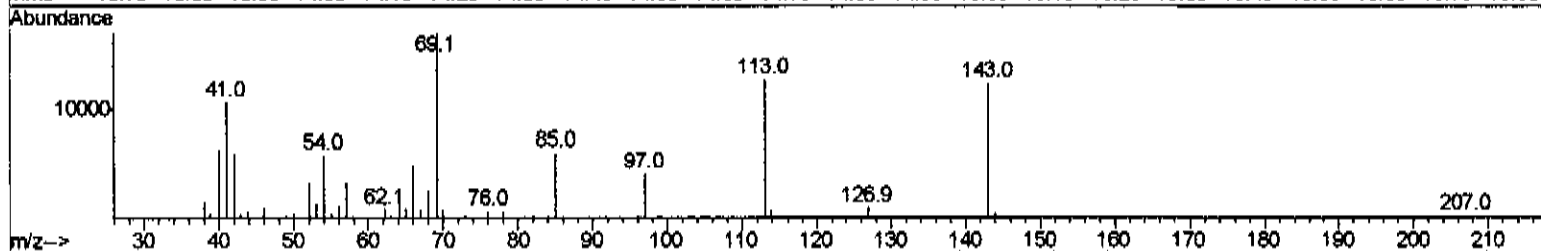
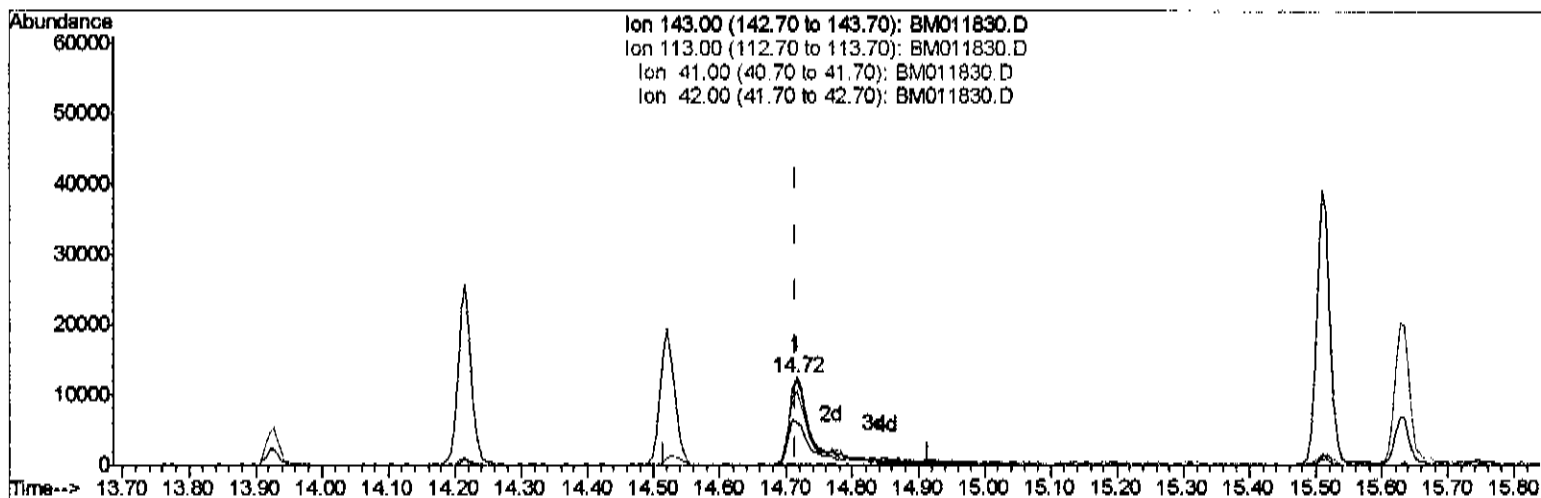
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Sample : I5458-09
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
COAH2

Manual Integrations
APPROVED

Sohil
10/2/2017 5:52:21 PM

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BM011830.D

(51) 4-Nitrophenol-d4 (S)

14.715min (-0.000) 5.87ng/ul m

response 28267

Ion	Exp%	Act%
143.00	100	100
113.00	114.60	102.92
41.00	34.80	86.72#
42.00	20.50	49.65#

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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	7.88	152	116485	20.00	ng/ul	0.00
18) Naphthalene-d8	10.69	136	536610	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.52	164	328732	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.26	188	811405	20.00	ng/ul	0.00
75) Chrysene-d12	21.43	240	1003286	20.00	ng/ul	0.00
83) Perylene-d12	23.76	264	1160460	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	11299m	4.32	ng/uL	>0.00
5) Phenol-d5	7.04	99	83303	7.27	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.20	67	263570	31.72	ng/ul	0.00
9) 2-Chlorophenol-d4	7.41	132	226942	26.63	ng/ul	0.00
13) 4-Methylphenol-d8	8.58	113	160276	17.86	ng/ul	0.00
19) Nitrobenzene-d5	9.05	128	134380	32.10	ng/ul	0.00
22) 2-Nitrophenol-d4	9.77	143	152053	34.06	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.30	165	246953	28.98	ng/ul	0.00
29) 4-Chloroaniline-d4	10.84	131	10539	1.12	ng/ul	0.02
43) Dimethylphthalate-d6	13.93	166	1029835	36.52	ng/ul	0.00
46) Acenaphthylene-d8	14.22	160	1176864	34.52	ng/ul	0.00
51) 4-Nitrophenol-d4	14.72	143	28267m	5.87	ng/ul	>0.00
57) Fluorene-d10	15.52	176	873085	35.96	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.63	200	163424	29.73	ng/ul	0.00
70) Anthracene-d10	17.36	188	1506698	36.49	ng/ul	0.00
76) Pyrene-d10	19.65	212	1936827	40.45	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.61	264	1987688	35.77	ng/ul	0.00

Target Compounds Ovalue

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