

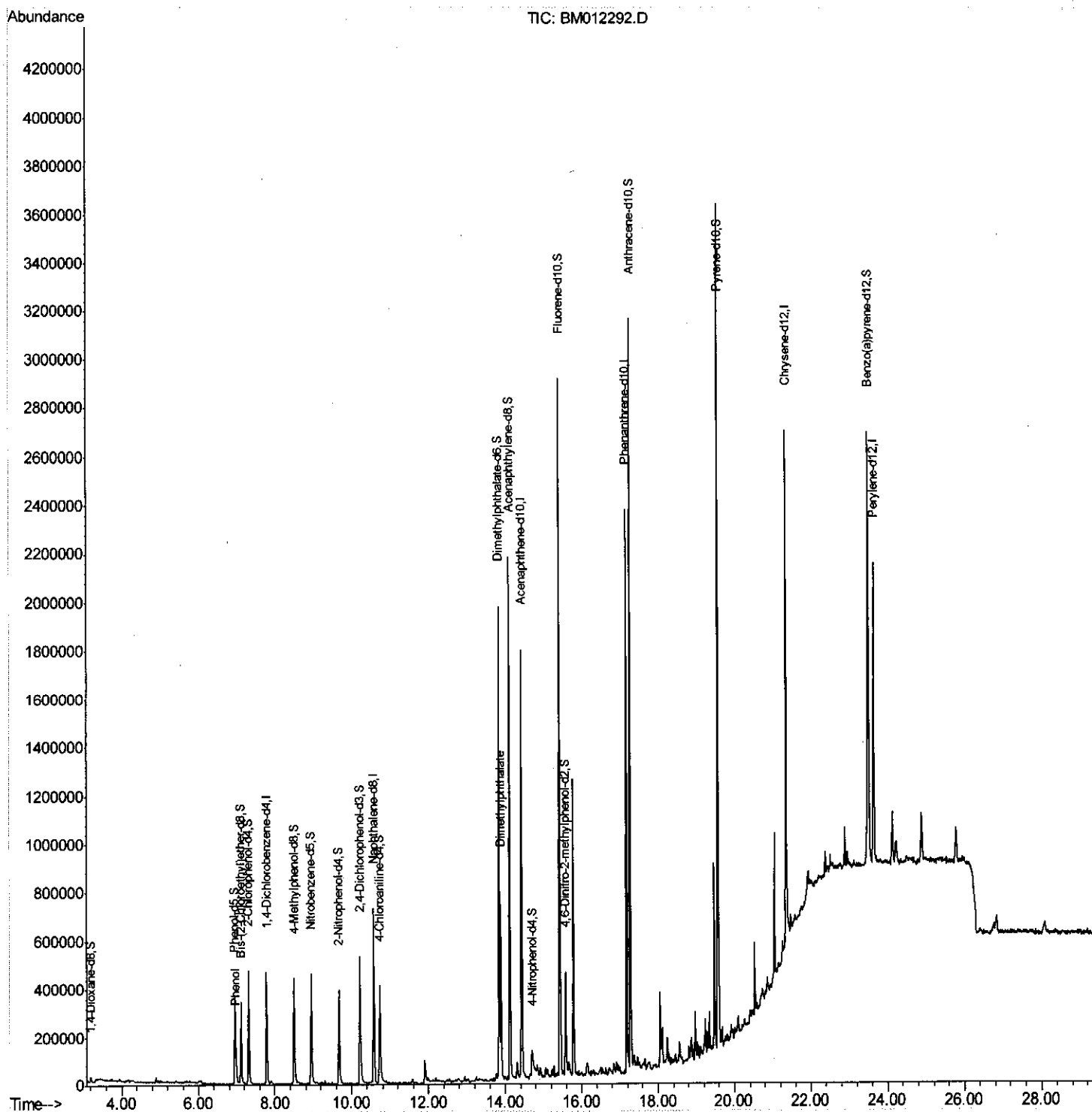
Data Path : Z:\HPCHEM1\BNA M\DATA\BM101817\
Data File : BM012292.D
Acq On : 18 Oct 2017 23:01
Operator : SJ/JU
Sample : I5656-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
Client Sampled :
B0C97

Manual Integrations
APPROVED

Sohil
10/23/2017 3:49:41 PM

Quant Time: Oct 18 23:50:24 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 18 23:09:30 2017
Response via : Initial Calibration



Quantitation Report (Qedit)

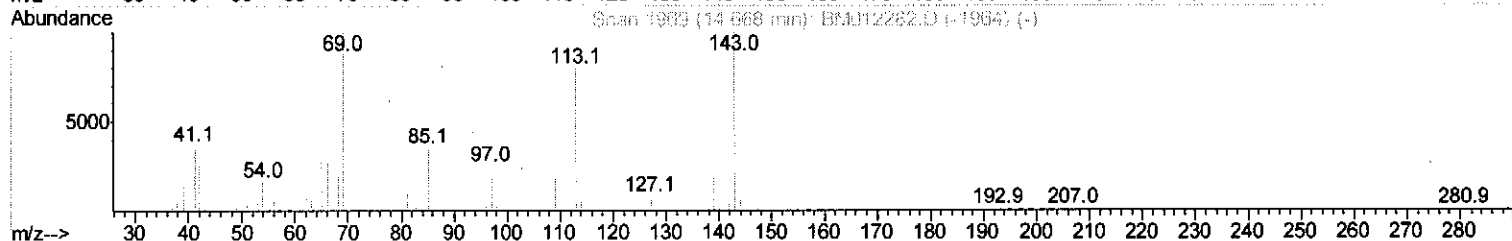
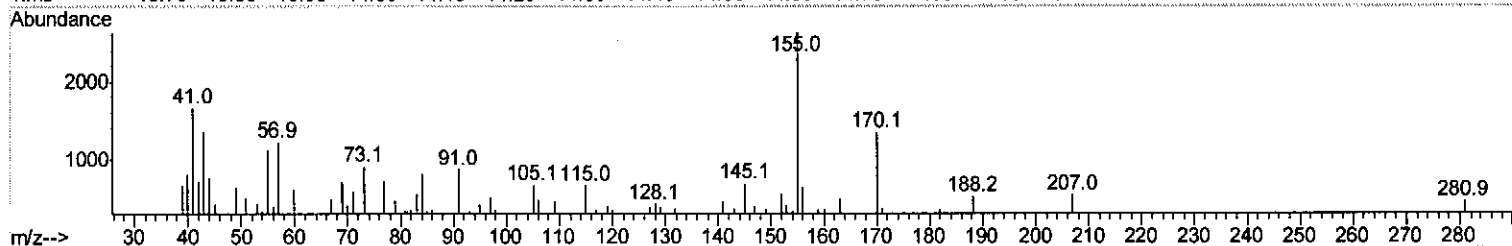
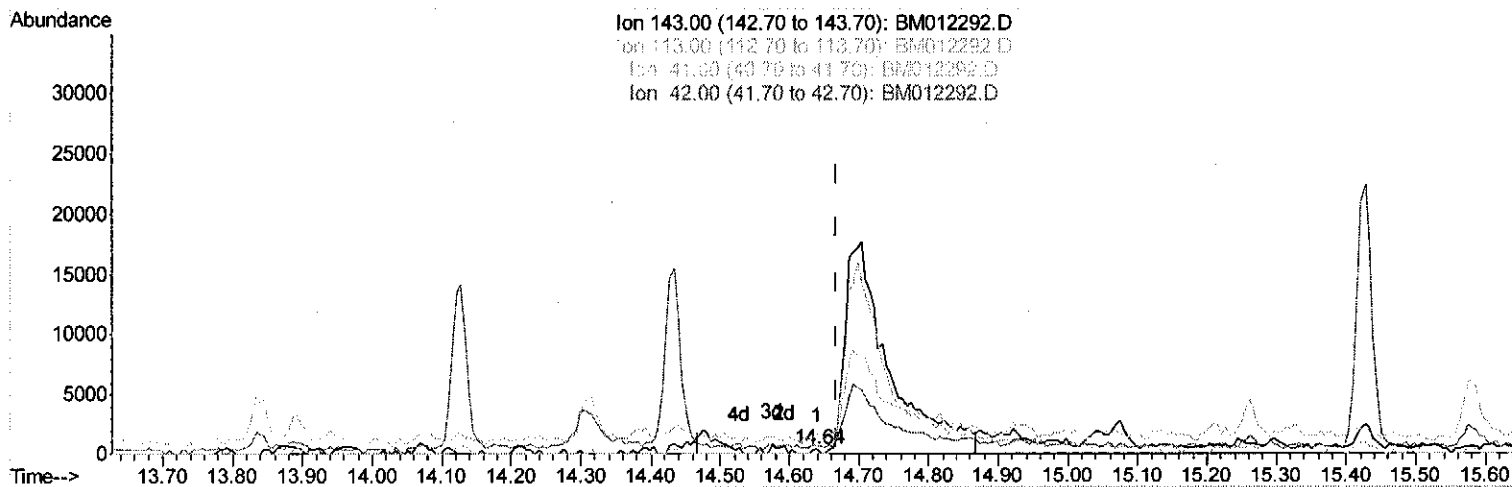
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TIC: BM012292.D

(51) 4-Nitrophenol-d4 (S)

14.639min (-0.029) 0.03ng/ul

response 242

Ion	Exp%	Act%
143.00	100	100
113.00	81.80	0.00#
41.00	40.00	449.73#
42.00	25.70	196.74#

Quantitation Report (Qedit)

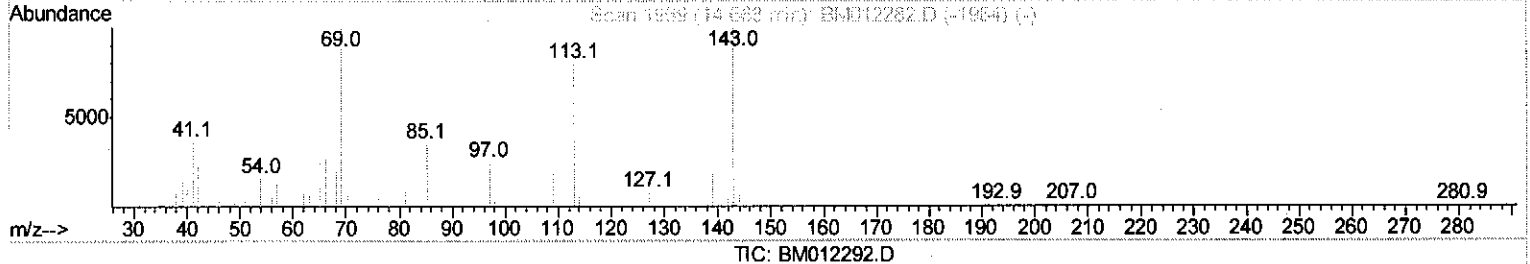
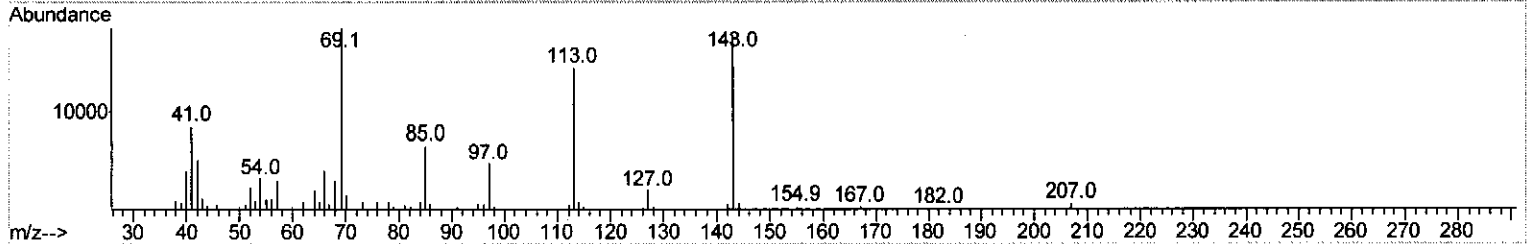
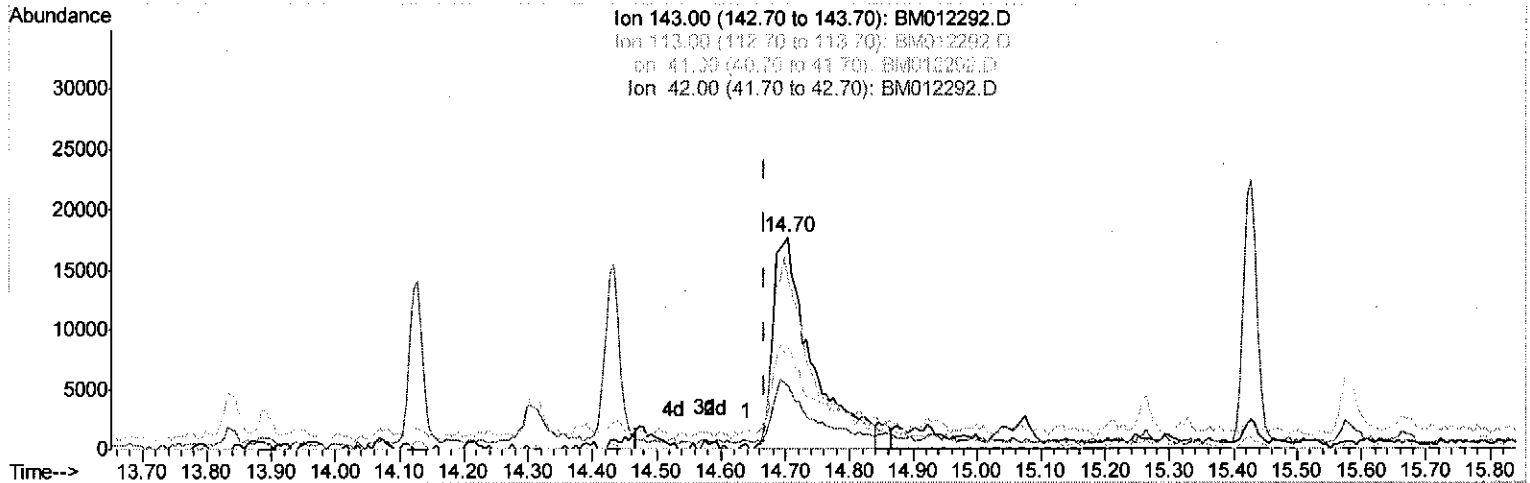
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 BNA_M
 ClientSampled :
 B0C97

Manual Integrations
 APPROVED

Sohil
 10/23/2017 3:49:41 PM

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 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 18 23:09:30 2017
 Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

14.704min (+0.035) 10.01ng/ul m

JU 10/24/17

response 75629

Ion	Exp%	Act%
143.00	100	100
113.00	81.80	81.02
41.00	40.00	47.97
42.00	25.70	30.05

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101817\
 Data File : BM012292.D
 Acq On : 18 Oct 2017 23:01
 Operator : SJ/JU
 Sample : I5656-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B0C97

Manual Integrations
 APPROVED

Sohil
 10/23/2017 3:49:41 PM

Quant Time: Oct 18 23:50:24 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM101217.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 18 23:09:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	140213	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	614425	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.43	164	568374	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.18	188	1265113	20.00	ng/ul	0.00
75) Chrysene-d12	21.36	240	1009136	20.00	ng/ul	0.00
83) Perylene-d12	23.66	264	938271	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.20	96	12535	4.25	ng/uL	0.00
5) Phenol-d5	6.95	99	234482	20.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.11	67	153832	22.55	ng/ul	0.00
9) 2-Chlorophenol-d4	7.31	132	226017	23.42	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	196821	20.09	ng/ul	0.00
19) Nitrobenzene-d5	8.95	128	117822	25.13	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	140625	25.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.22	165	253791	24.26	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	279963	28.60	ng/ul	0.00
43) Dimethylphthalate-d6	13.84	166	1281908	25.52	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	1513985	24.90	ng/ul	0.00
51) 4-Nitrophenol-d4	14.70	143	75629m	10.01	ng/ul	0.04 →
57) Fluorene-d10	15.43	176	1146706	27.81	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.58	200	118406	13.64	ng/ul	0.00
70) Anthracene-d10	17.28	188	1695833	27.11	ng/ul	0.00
76) Pyrene-d10	19.57	212	1732813	33.83	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.51	264	1260788	27.87	ng/ul	0.00

Target Compounds					Ovalue
6) Phenol	6.98	94	44177	3.756	ng/ul 95
44) Dimethylphthalate	13.89	163	519604	10.324	ng/ul 99

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