

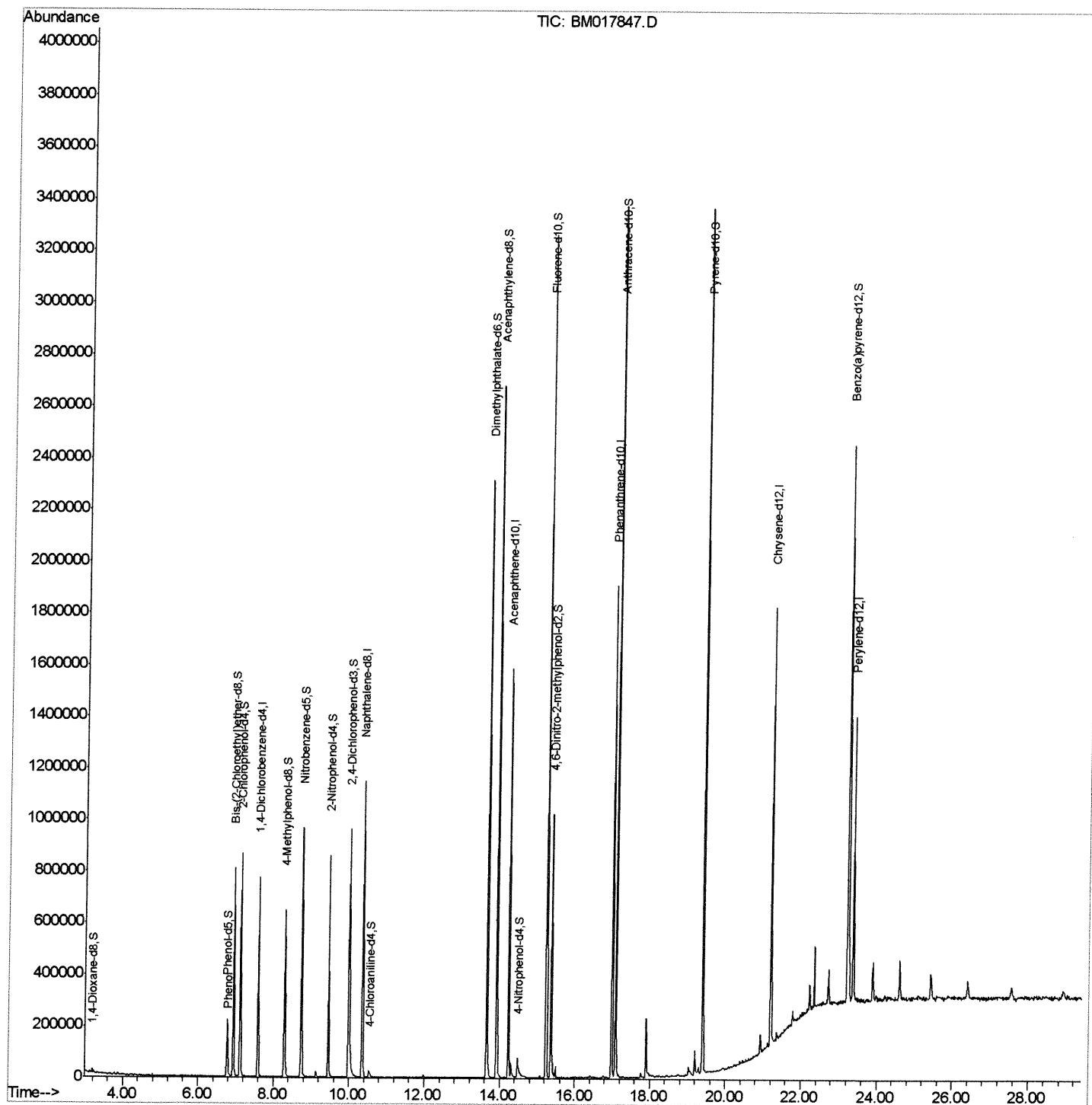
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120118\
Quantitation Report (QT/LSC Reviewed)
Data File : BM017847.D
Acq On : 01 Dec 2018 07:21
Operator : JU/SJ
Sample : J6096-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4172

Manual Integrations
APPROVED

mohammad
12/3/2018 4:57:39 PM

Quant Time: Dec 03 00:43:35 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Dec 03 00:28:52 2018
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM120118\
Quantitation Report (Qedit)

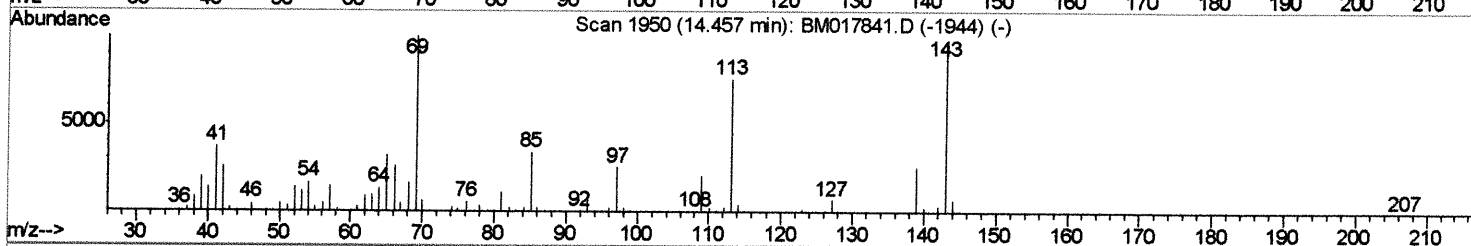
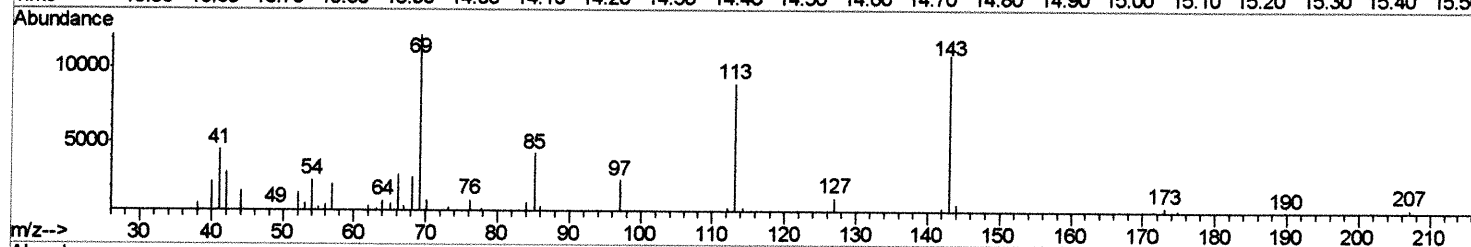
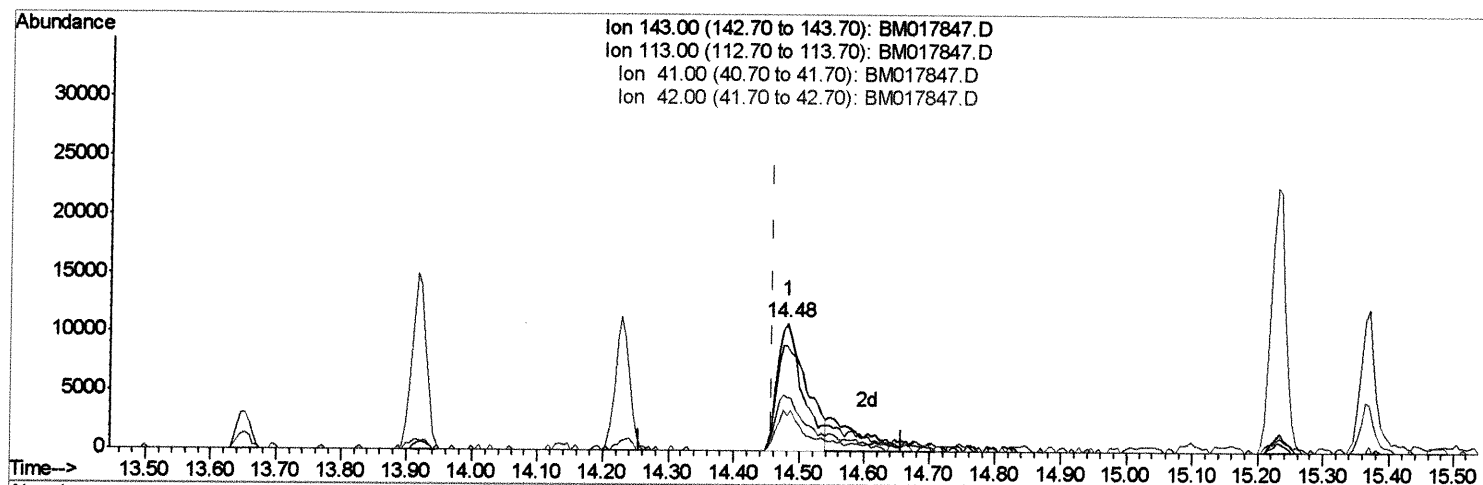
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TIC: BM017847.D

(51) 4-Nitrophenol-d4 (S)

14.481min (+0.024) 4.05ng/ul

response 33063

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	82.14
41.00	40.10	40.96
42.00	26.80	27.00

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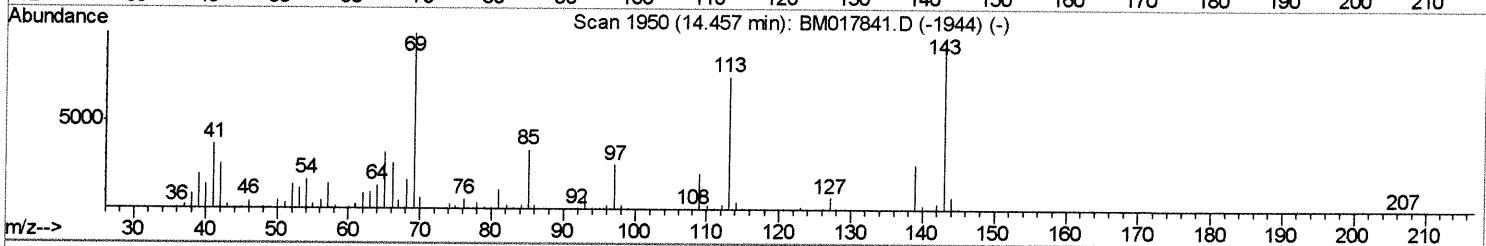
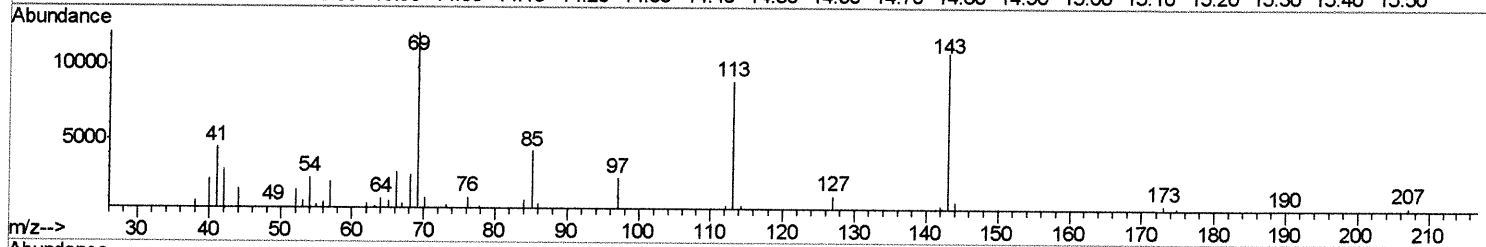
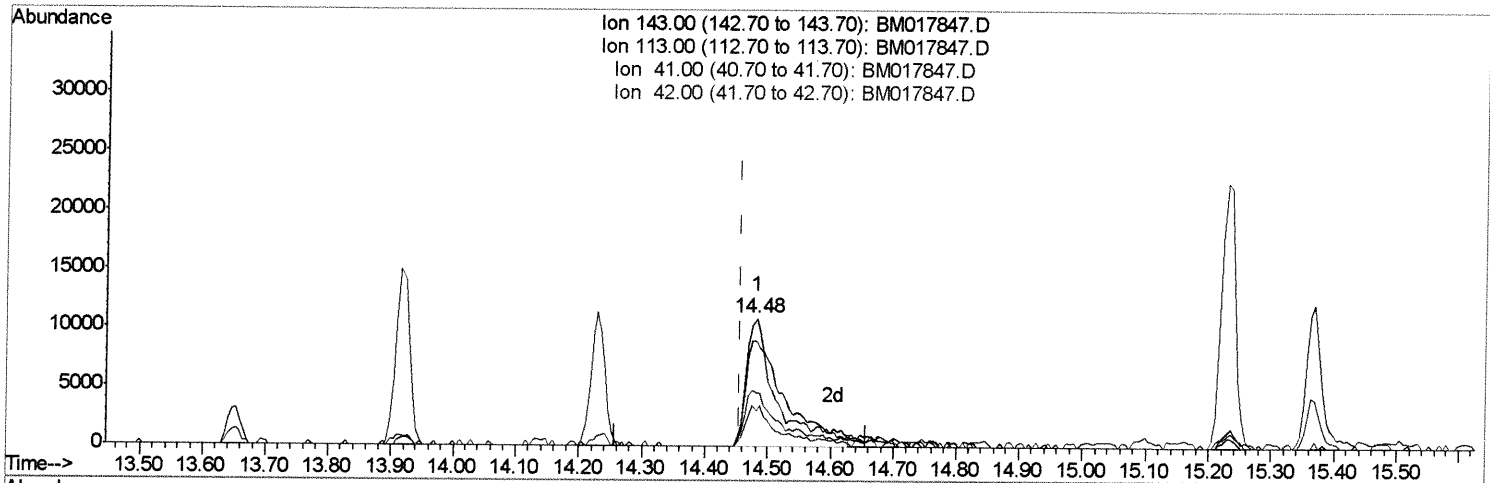
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Manual Integrations
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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BM017847.D

(51) 4-Nitrophenol-d4 (S)

14.481min (+0.024) 5.19ng/ul m > SJ 12/11/18

response 42407

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	82.14
41.00	40.10	40.96
42.00	26.80	27.00

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 Quantitation Report (QT/LSC Reviewed)
 Data File : BM017847.D
 Acq On : 01 Dec 2018 07:21
 Operator : JU/SJ
 Sample : J6096-23
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4172

Manual Integrations
 APPROVED

mohammad
 12/3/2018 4:57:39 PM

Quant Time: Dec 03 00:43:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 03 00:28:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	211903	20.00	ng/ul	0.00
18) Naphthalene-d8	10.35	136	923583	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.23	164	517940	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.99	188	1059691	20.00	ng/ul	0.00
77) Chrysene-d12	21.19	240	813348	20.00	ng/ul	0.00
85) Perylene-d12	23.37	264	779868	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	6888	1.49	ng/uL	0.00
5) Phenol-d5	6.77	99	138589	7.16	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.93	67	359279	30.79	ng/ul	0.00
9) 2-Chlorophenol-d4	7.12	132	400568	26.45	ng/ul	0.00
13) 4-Methylphenol-d8	8.29	113	255415	16.15	ng/ul	0.00
19) Nitrobenzene-d5	8.74	128	234590	33.07	ng/ul	0.00
22) 2-Nitrophenol-d4	9.45	143	269990	33.21	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.98	165	431284	27.87	ng/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	25388	1.91	ng/ul	0.03
43) Dimethylphthalate-d6	13.65	166	1505614	33.50	ng/ul	0.00
46) Acenaphthylene-d8	13.92	160	1879601	34.56	ng/ul	0.00
51) 4-Nitrophenol-d4	14.48	143	42407m	5.19	ng/ul	0.02
57) Fluorene-d10	15.23	176	1293951	33.34	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.37	200	224709	29.66	ng/ul	0.00
70) Anthracene-d10	17.09	188	1839272	34.86	ng/ul	0.00
78) Pyrene-d10	19.39	212	1702663	43.26	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.23	264	1518104	35.89	ng/ul	0.00

SJ 12/11/18

Target Compounds				Qvalue
6) Phenol	6.80	94	24221	1.254 ng/ul 95

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