

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM121118\
Quantitation Report (QT Reviewed)

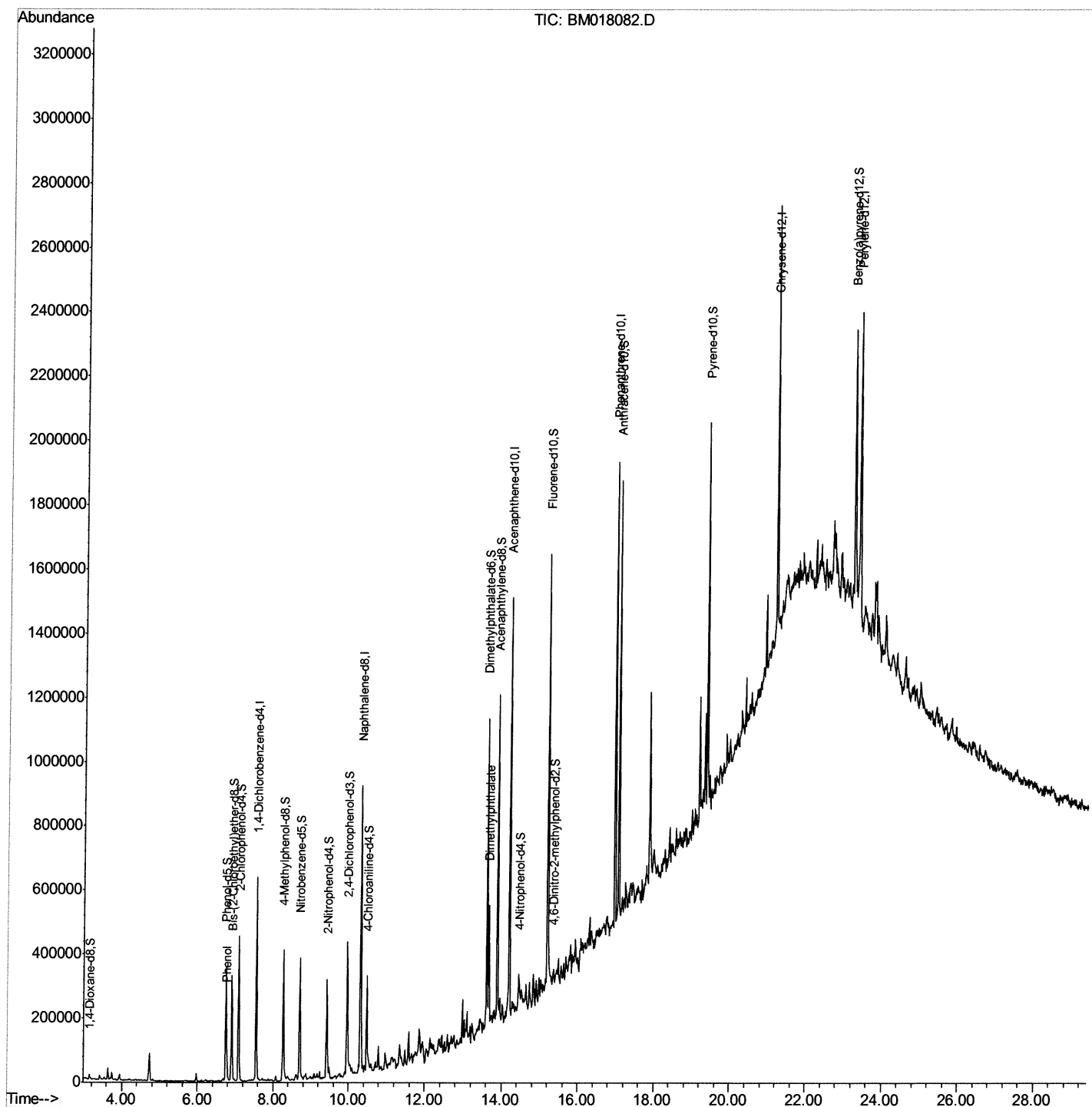
Data File : BM018082.D
Acq On : 12 Dec 2018 06:17
Operator : JU/SJ
Sample : J6171-03
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F9Y07

Manual Integrations
APPROVED

Sohil
12/12/2018 5:43:35 PM

Quant Time: Dec 12 07:15:36 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 12 06:24:35 2018
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM121118\
Quantitation Report (Qedit)

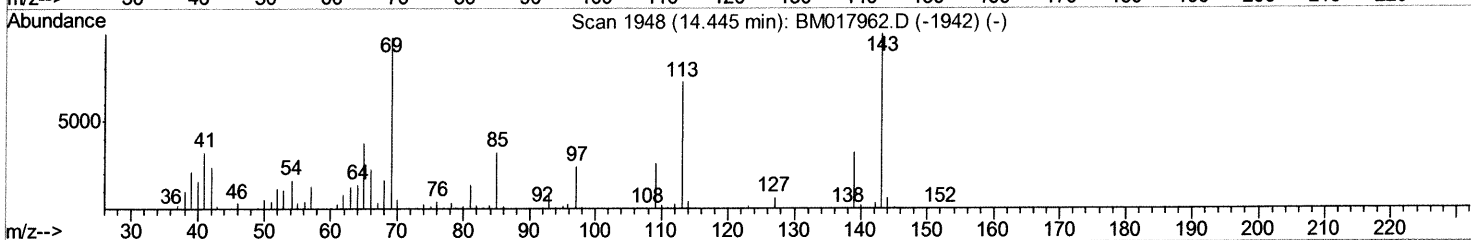
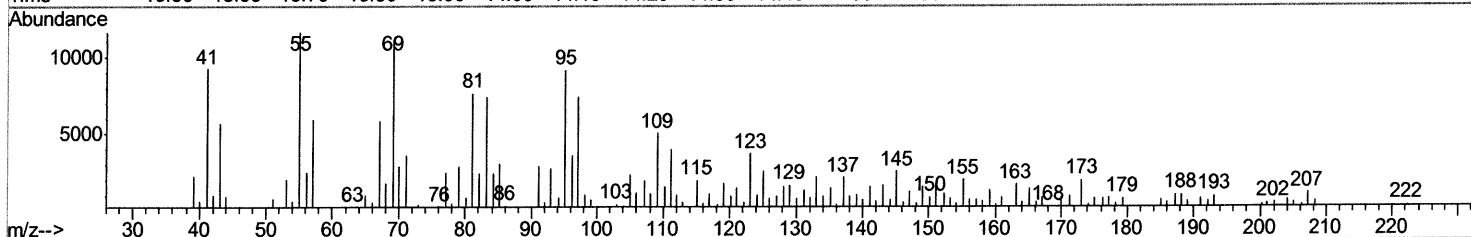
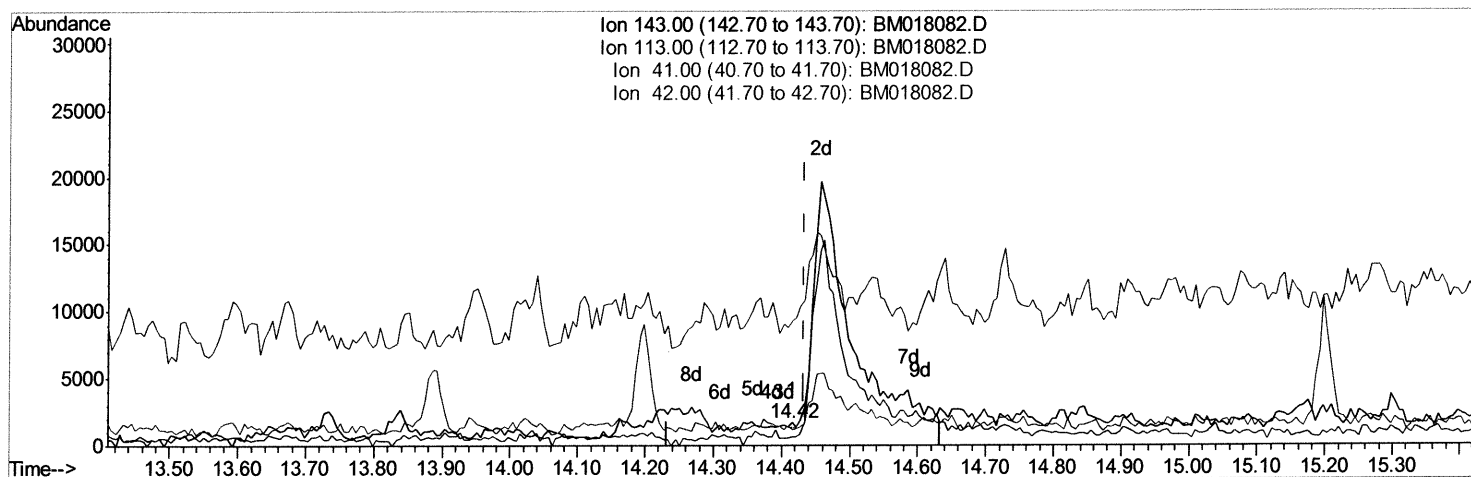
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TIC: BM018082.D

(51) 4-Nitrophenol-d4 (S)

14.416min (-0.018) 0.02ng/ul

response 123

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	36.71#
41.00	40.10	555.51#
42.00	26.80	67.31#

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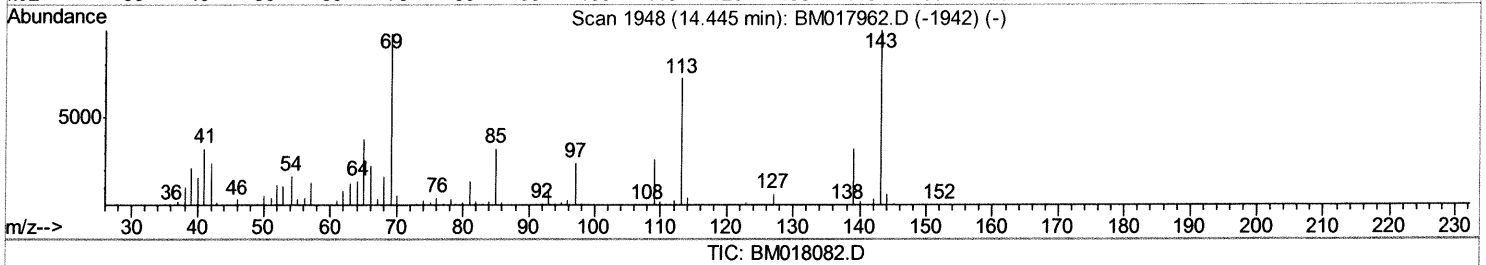
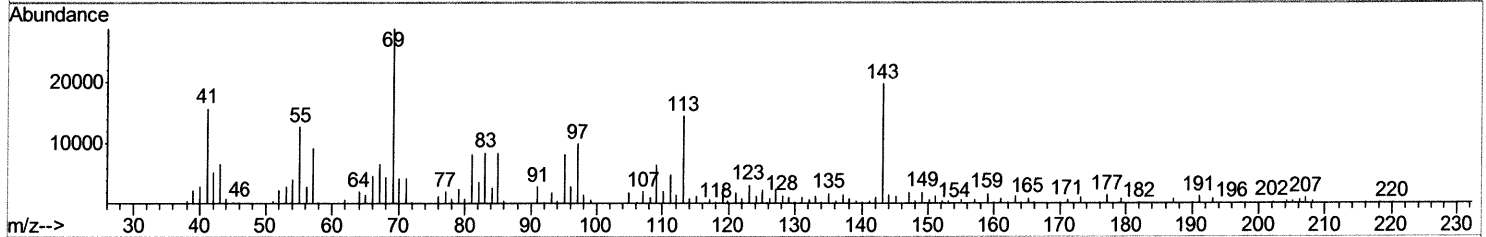
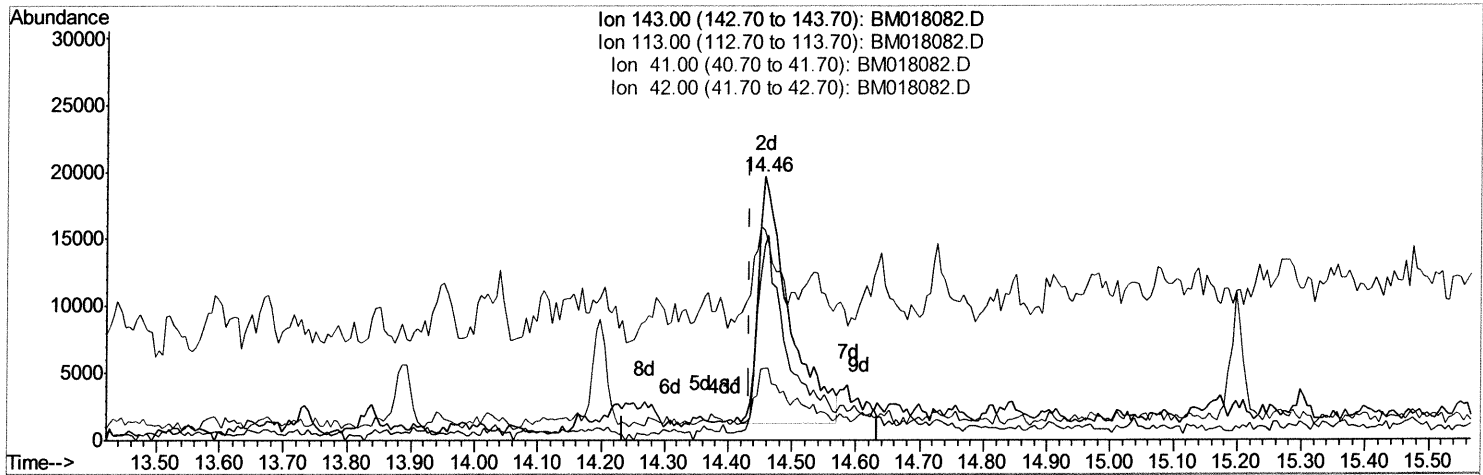
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Manual Integrations
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Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

14.457min (+0.023) 9.44ng/ul m> 50 12/12/18

response 62161

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	73.44
41.00	40.10	79.74#
42.00	26.80	27.12

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 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F9Y07

Manual Integrations
 APPROVED

Sohil
 12/12/2018 5:43:35 PM

Quant Time: Dec 12 07:15:36 2018
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 12 06:24:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	176745	20.00	ng/ul	0.00
18) Naphthalene-d8	10.32	136	757670	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.20	164	417431	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.96	188	828558	20.00	ng/ul	0.00
77) Chrysene-d12	21.17	240	608640	20.00	ng/ul	0.00
85) Perylene-d12	23.35	264	641499	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.15	96	9721	2.52	ng/uL	0.00
5) Phenol-d5	6.75	99	240044	14.87	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.90	67	144148	14.81	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	214185	16.96	ng/ul	0.00
13) 4-Methylphenol-d8	8.26	113	179815	13.63	ng/ul	0.00
19) Nitrobenzene-d5	8.70	128	102286	17.58	ng/ul	0.00
22) 2-Nitrophenol-d4	9.42	143	108624	16.29	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.95	165	189264	14.91	ng/ul	0.00
29) 4-Chloroaniline-d4	10.47	131	164417	15.09	ng/ul	0.00
43) Dimethylphthalate-d6	13.62	166	664375	18.34	ng/ul	0.00
46) Acenaphthylene-d8	13.89	160	784297	17.89	ng/ul	0.00
51) 4-Nitrophenol-d4	14.46	143	62161m	9.44	ng/ul	0.02
57) Fluorene-d10	15.20	176	553032	17.68	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.36	200	9166	1.55	ng/ul	0.01
70) Anthracene-d10	17.06	188	742072	17.99	ng/ul	0.00
78) Pyrene-d10	19.36	212	631655	21.45	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.21	264	589375	16.94	ng/ul	0.01

SJ 12/13/18

Target Compounds

					Qvalue
6) Phenol	6.77	94	56956	3.535	ng/ul 97
44) Dimethylphthalate	13.67	163	218861	6.237	ng/ul 100

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