

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM103118\
Quantitation Report (QT Reviewed)

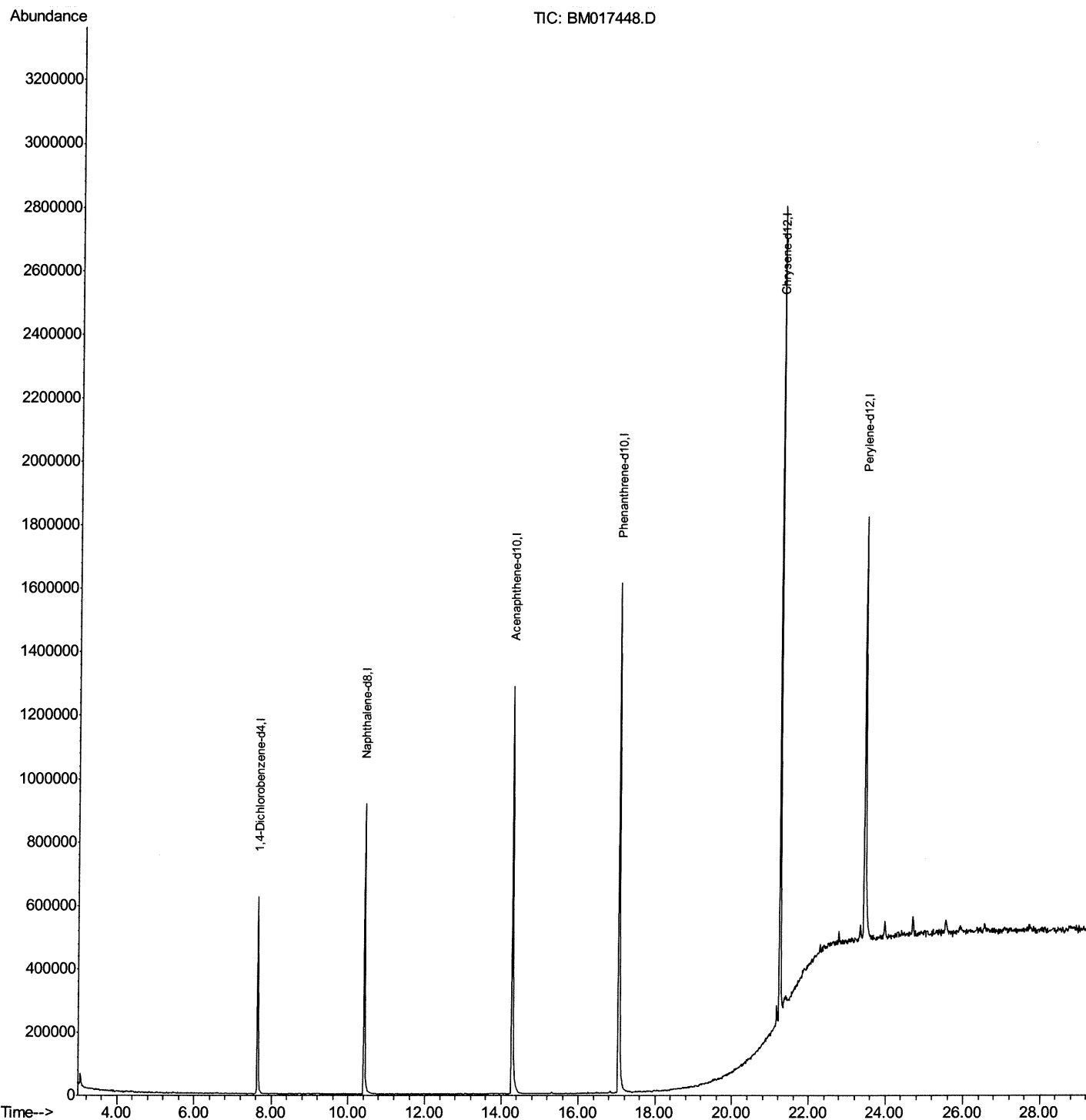
Data File : BM017448.D
Acq On : 01 Nov 2018 02:20
Operator : MJ/SJ
Sample : J5699-10
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SVOC-GPC2-BLANK

Manual Integrations
APPROVED

Sohil
11/1/2018 10:55:31 AM

Quant Time: Nov 01 07:48:12 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 01 06:46:30 2018
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM103118\
Quantitation Report (Qedit)

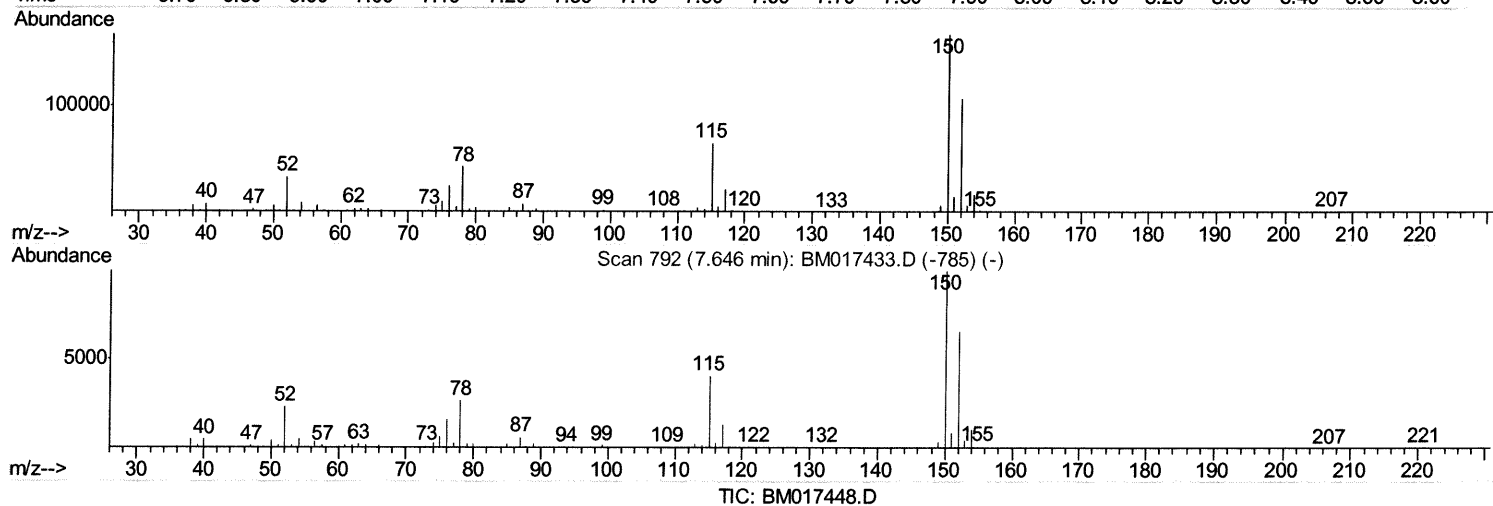
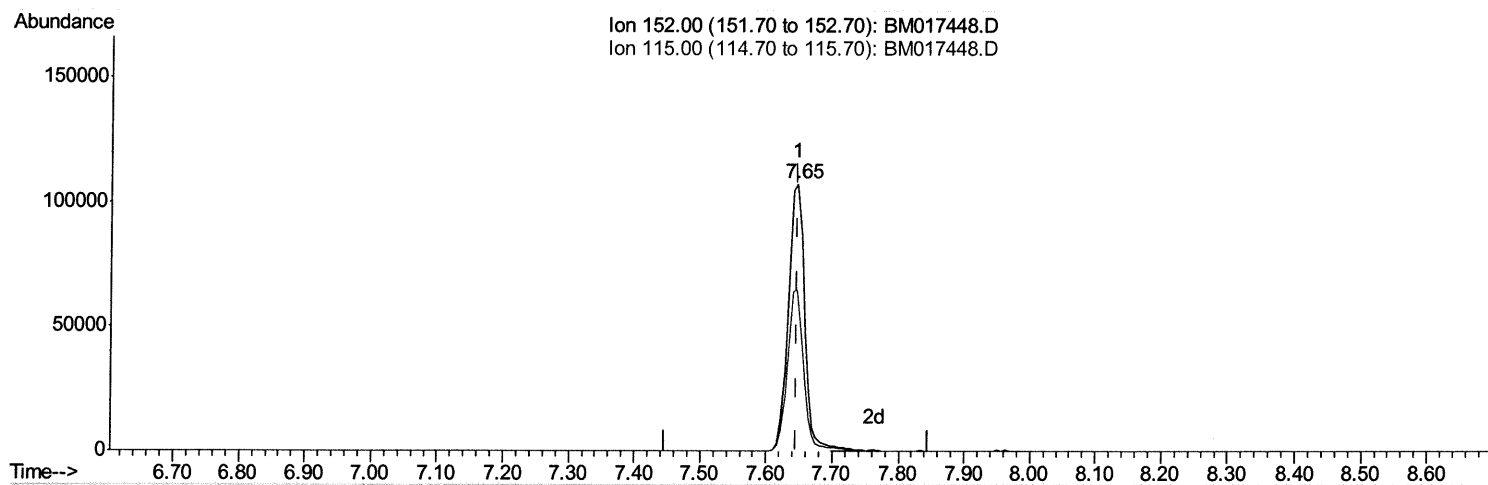
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(1) 1,4-Dichlorobenzene-d4 (I)

7.646min (+0.000) 20.00ng/ul

response 182031

Ion	Exp%	Act%
152.00	100	100
115.00	59.40	60.51
0.00	0.00	0.00
0.00	0.00	0.00

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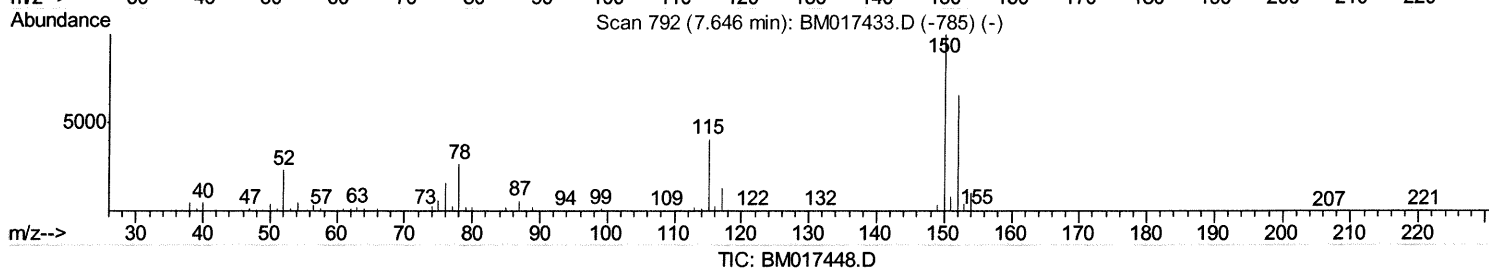
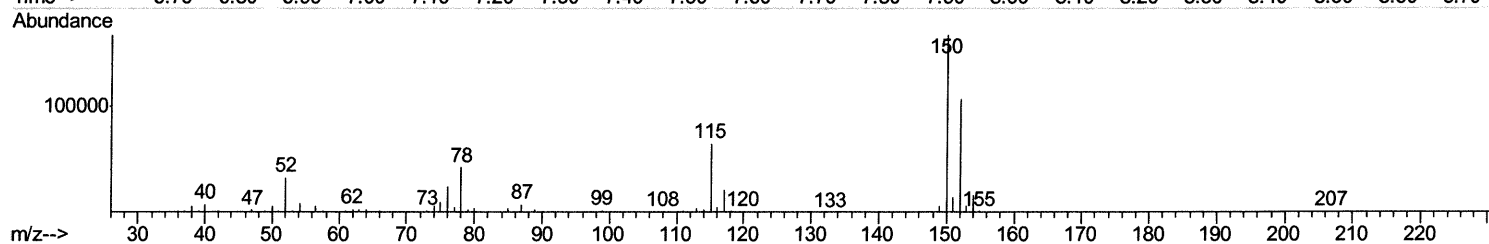
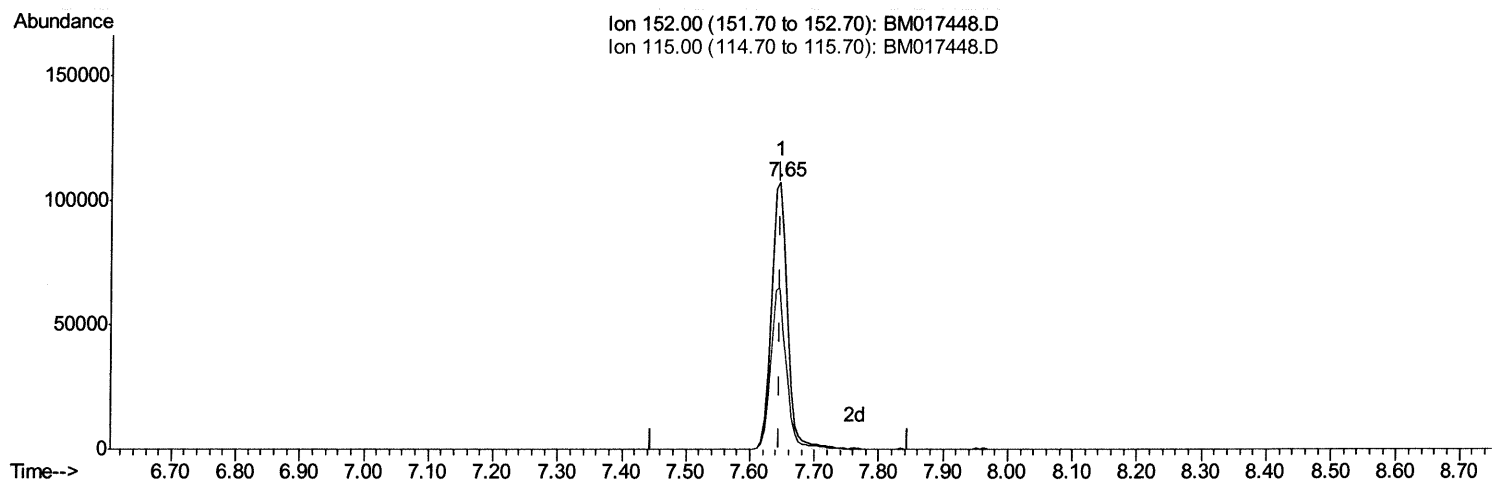
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Manual Integrations
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(1) 1,4-Dichlorobenzene-d4 (I)

7.646min (+0.000) 20.00ng/ul m

SJ11/2/18

response 185209

Ion	Exp%	Act%
152.00	100	100
115.00	59.40	60.51
0.00	0.00	0.00
0.00	0.00	0.00

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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	7.65	152	185209m	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	830541	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	551870	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.05	188	1308159	20.00	ng/ul	0.00
78) Chrysene-d12	21.24	240	1329191	20.00	ng/ul	0.00
86) Perylene-d12	23.44	264	1152806	20.00	ng/ul	0.00

5/11/2/18

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL
5) Phenol-d5	0.00	99	0	0.00	ng/ul
7) Bis-(2-Chloroethyl) ether-d	0.00	67	0	0.00	ng/ul
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul
44) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul
47) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul
58) Fluorene-d10	0.00	176	0d	0.00	ng/ul
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul
71) Anthracene-d10	0.00	188	0d	0.00	ng/ul
79) Pyrene-d10	0.00	212	0	0.00	ng/ul
90) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul

Target Compounds Qvalue

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