

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM111918\
Quantitation Report (QT Reviewed)

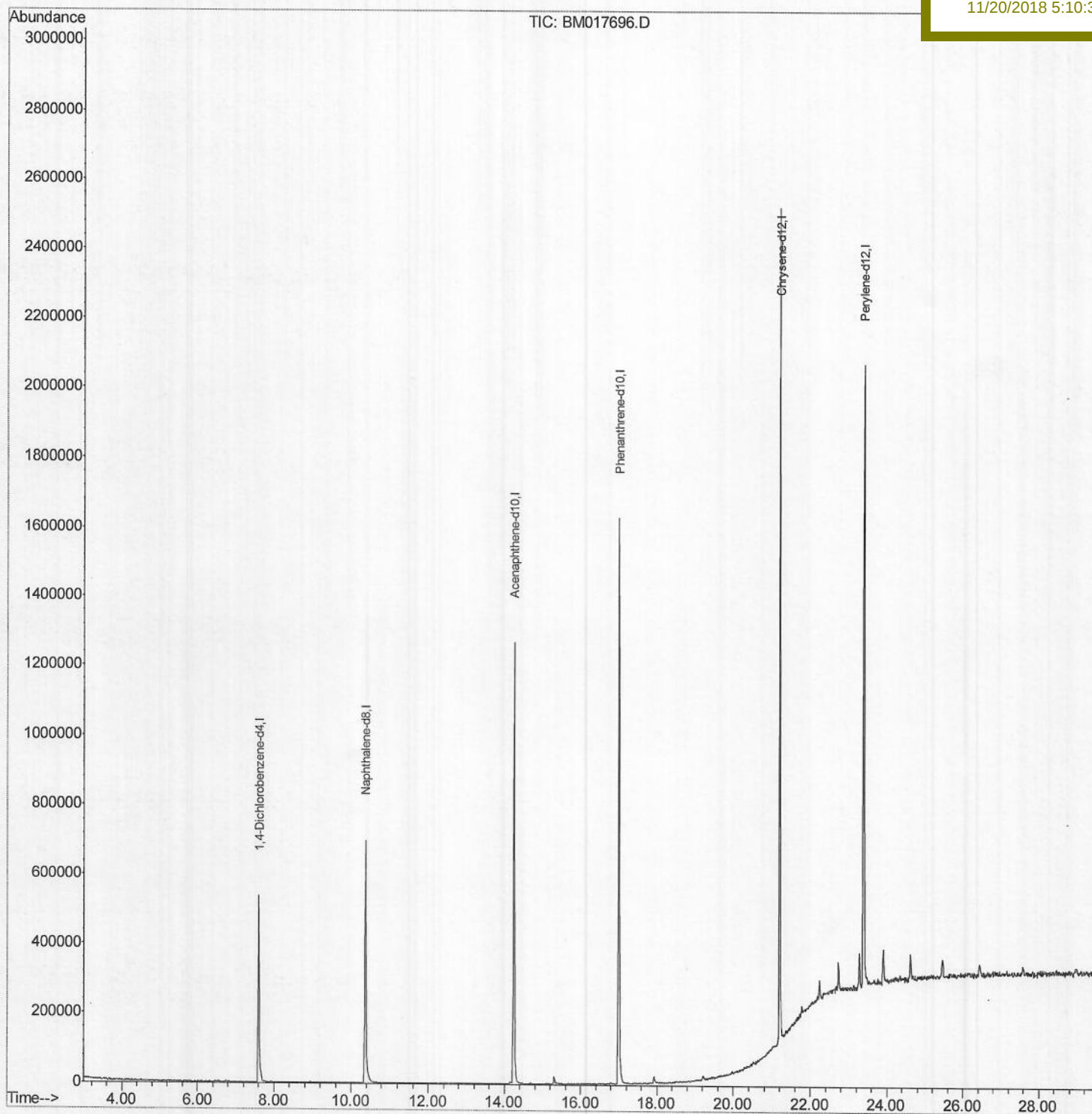
Data File : BM017696.D
Acq On : 19 Nov 2018 23:14
Operator : MJ/SJ
Sample : J5984-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Quant Time: Nov 20 01:21:59 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 20 01:04:58 2018
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
SVOC-GPC-BLANK

Manual Integrations
APPROVED

Sohil
11/20/2018 5:10:34 PM



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM111918\
Quantitation Report (Qedit)

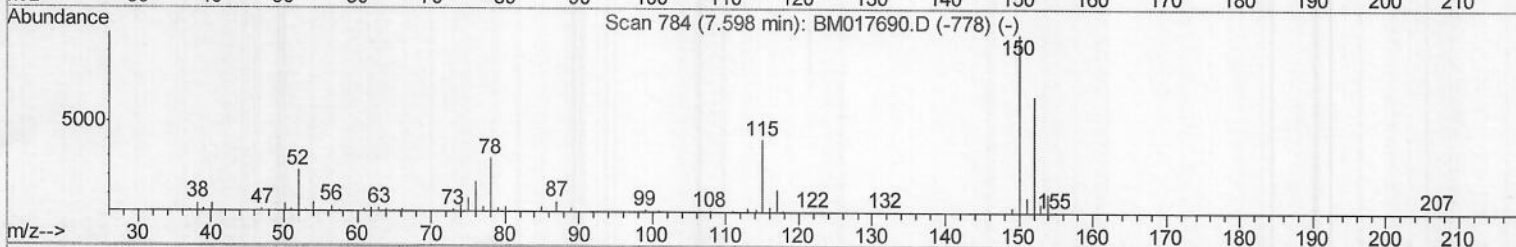
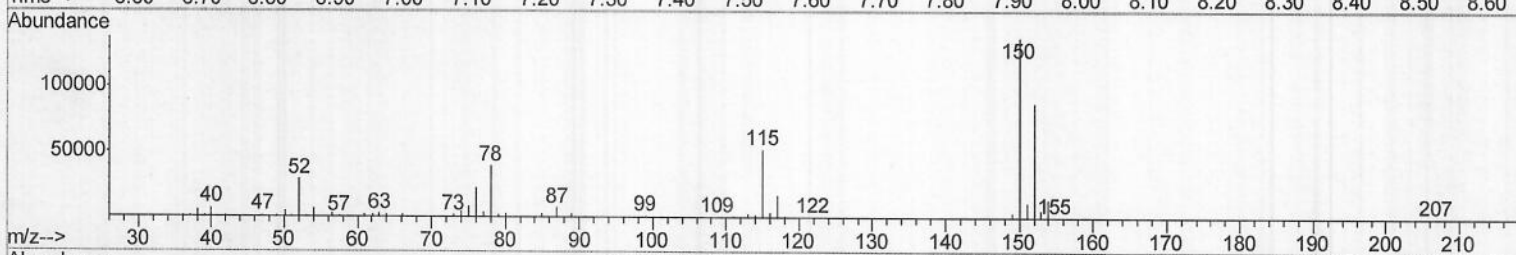
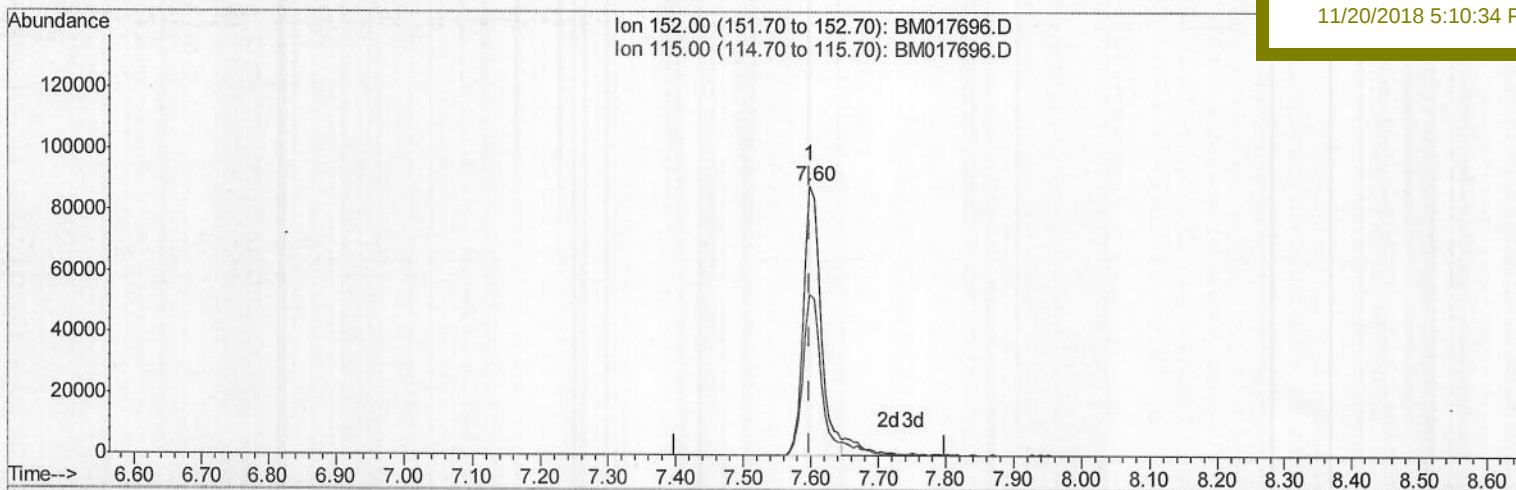
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Operator : MJ/SJ
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Quant Time: Nov 20 01:15:21 2018
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SVOC-GPC-BLANK

Manual Integrations
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TIC: BM017696.D

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

7.598min (+0.000) 20.00ng/ul

response 157958

Ion	Exp%	Act%
152.00	100	100
115.00	62.10	59.95
0.00	0.00	0.00
0.00	0.00	0.00

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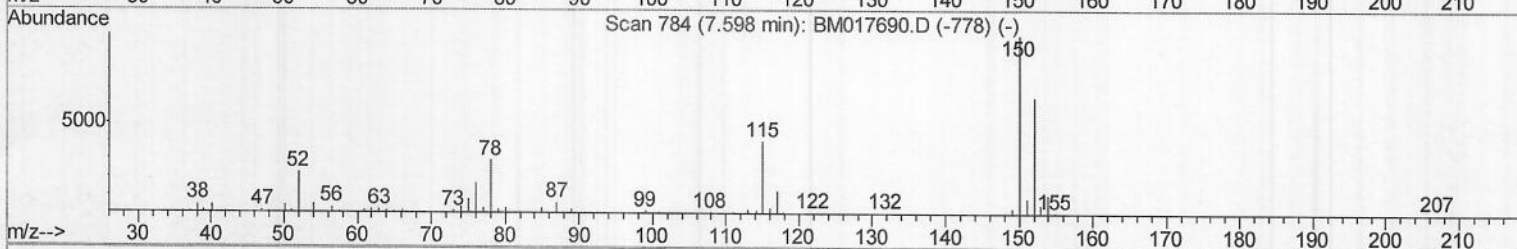
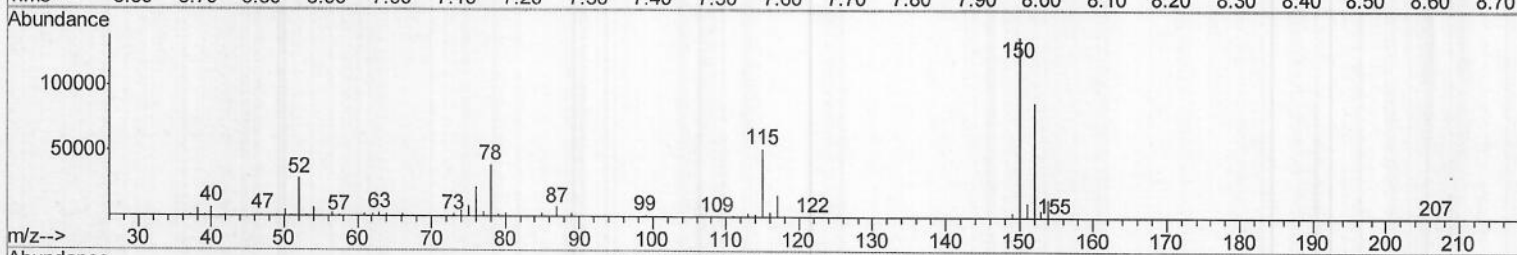
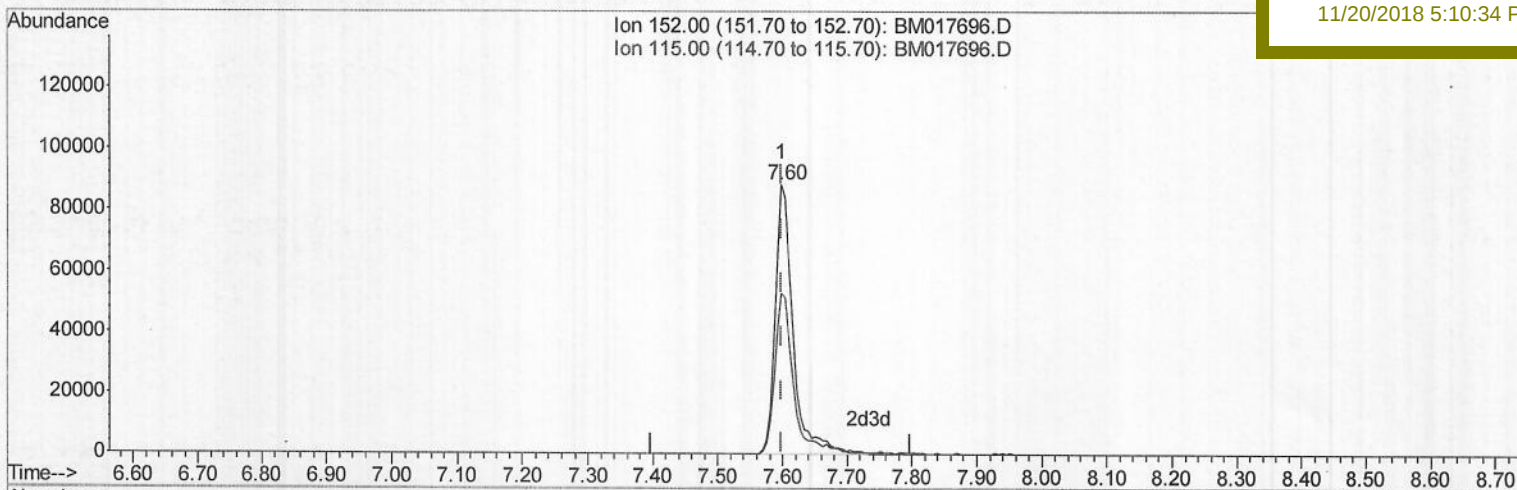
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Instrument :
BNA_M
ClientSampleId :
SVOC-GPC-BLANK

Manual Integrations
APPROVED

Sohil
11/20/2018 5:10:34 PM



TIC: BM017696.D

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

7.598min (+0.000) 20.00ng/ul m

response 169659

Ion	Exp%	Act%
-----	------	------

152.00	100	100
--------	-----	-----

115.00	62.10	59.95
--------	-------	-------

0.00	0.00	0.00
------	------	------

0.00	0.00	0.00
------	------	------

SJ
11/20/18

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Misc :

ALS Vial : 10 Sample Multiplier: 1

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Response via : Initial Calibration

Instrument :

BNA_M

ClientSampleId :

SVOC-GPC-BLANK

Manual Integrations

APPROVED

Sohil

11/20/2018 5:10:34 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	169659m	20.00	ng/ul	0.00
18) Naphthalene-d8	10.37	136	724141	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.25	164	446530	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.00	188	1067302	20.00	ng/ul	0.00
77) Chrysene-d12	21.20	240	1230088	20.00	ng/ul	0.00
85) Perylene-d12	23.39	264	1306094	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	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
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul
46) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul
57) Fluorene-d10	0.00	176	0d	0.00	ng/ul
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul
70) Anthracene-d10	0.00	188	0d	0.00	ng/ul
78) Pyrene-d10	0.00	212	0d	0.00	ng/ul
89) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul

Target Compounds

Qvalue

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