

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112018\
Quantitation Report (QT Reviewed)

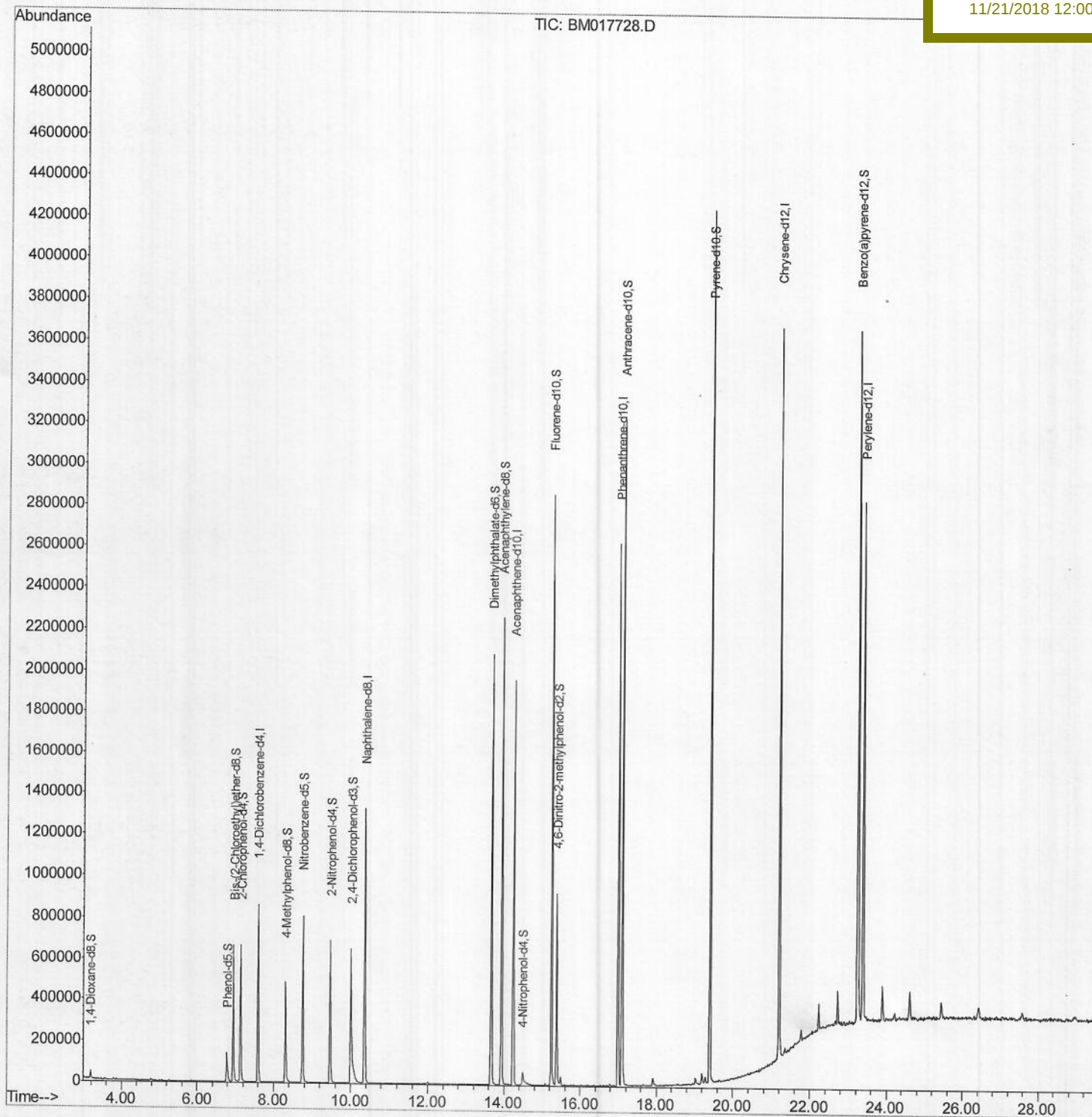
Data File : BM017728.D
Acq On : 21 Nov 2018 05:54
Operator : MJ/SJ
Sample : J5990-03
Misc :
ALS Vial : 29 Sample Multiplier: 1

Quant Time: Nov 21 07:58:57 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 21 07:44:17 2018
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
ETO78

Manual Integrations
APPROVED

Sohil
11/21/2018 12:00:54 PM



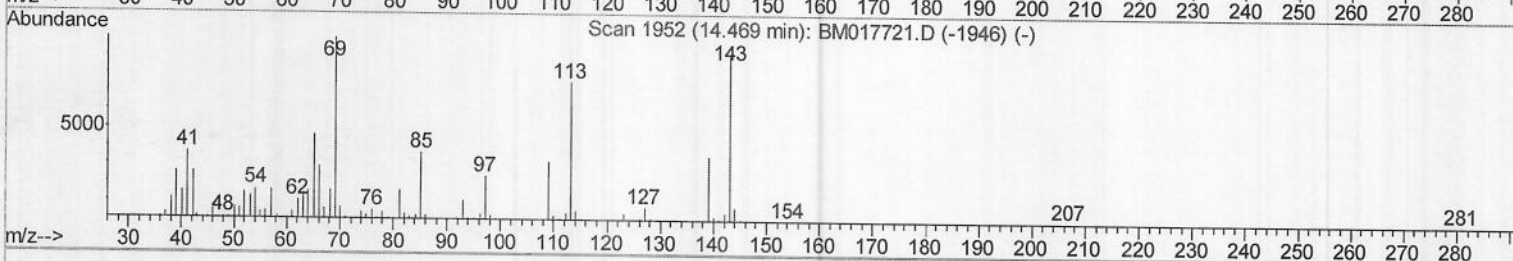
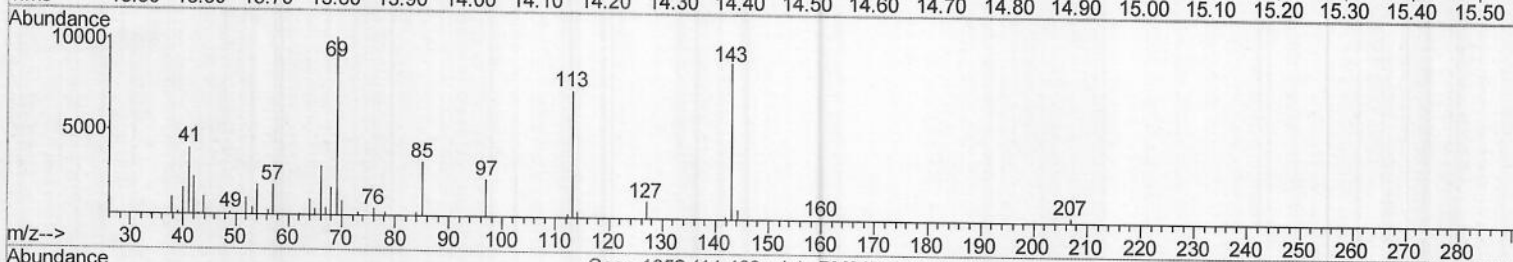
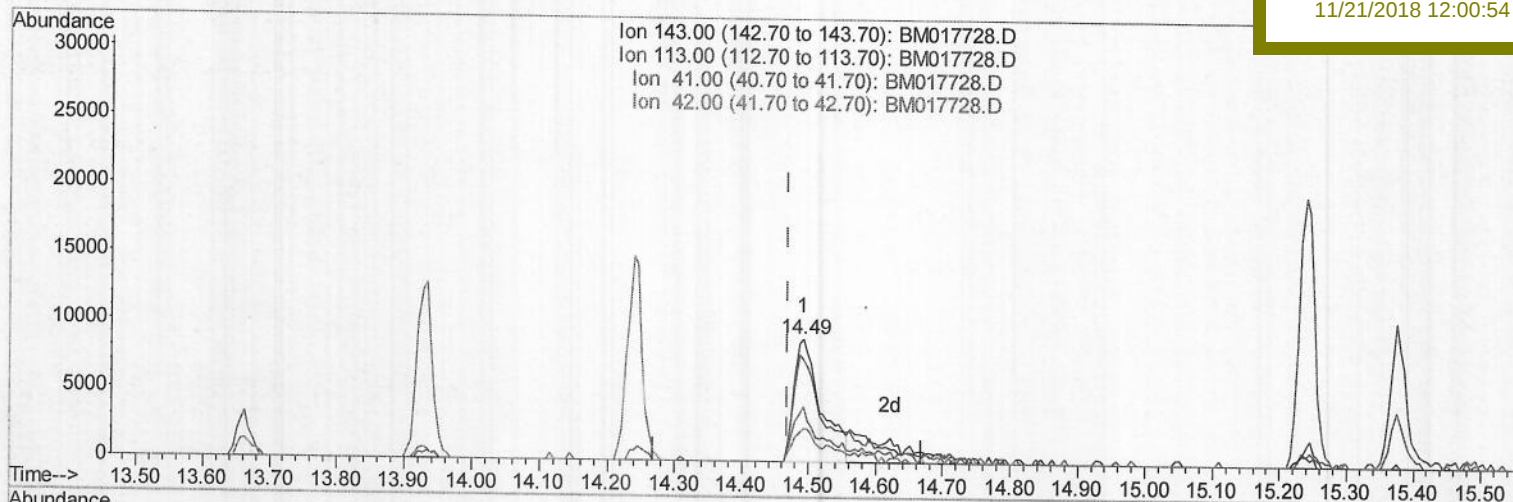
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Manual Integrations
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11/21/2018 12:00:54 PM



TIC: BM017728.D

(51) 4-Nitrophenol-d4 (S)
14.492min (+0.023) 2.48ng/ul
response 25084

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	82.28
41.00	40.10	44.98
42.00	26.80	26.60

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112018\
Quantitation Report (Qedit)

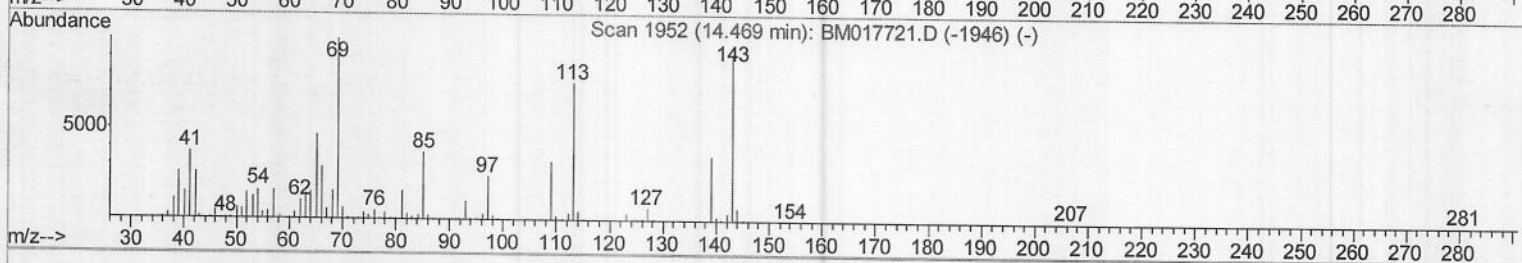
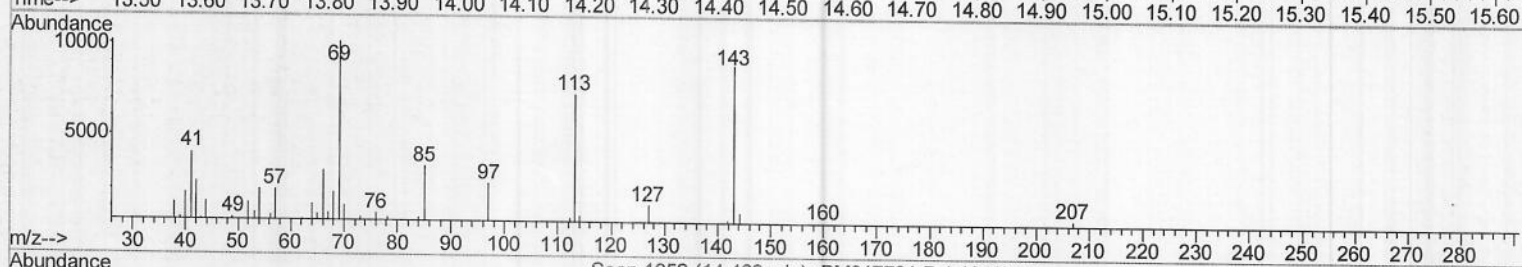
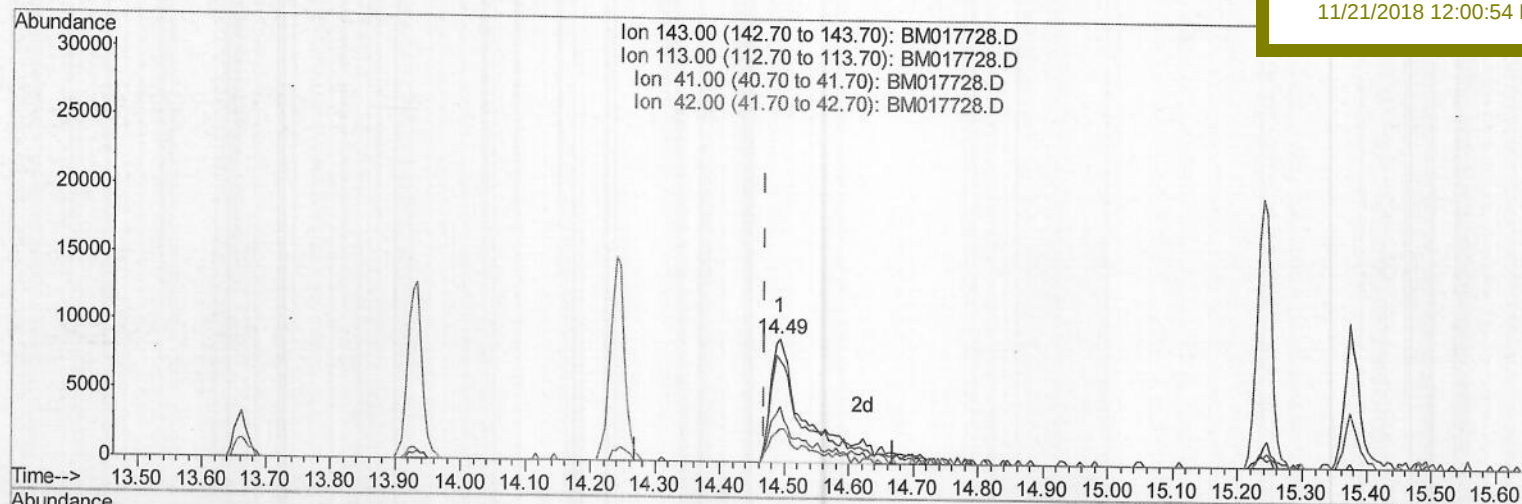
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Manual Integrations
APPROVED

Sohil
11/21/2018 12:00:54 PM



TIC: BM017728.D

(51) 4-Nitrophenol-d4 (S)

14.492min (+0.023) 3.25ng/ul m

response 32921

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	82.28
41.00	40.10	44.98
42.00	26.80	26.60

SJ
11/20/18

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112018\

Quantitation Report (QT Reviewed)

Data File : BM017728.D

Acq On : 21 Nov 2018 05:54

Operator : MJ/SJ

Sample : J5990-03

Misc :

ALS Vial : 29 Sample Multiplier: 1

Quant Time: Nov 21 07:58:57 2018

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Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 21 07:44:17 2018

Response via : Initial Calibration

Instrument :

BNA_M

ClientSampleId :

ETO78

Manual Integrations

APPROVED

Sohil

11/21/2018 12:00:54 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	236945	20.00	ng/ul	0.00
18) Naphthalene-d8	10.36	136	1071214	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.24	164	642348	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.00	188	1482451	20.00	ng/ul	0.00
77) Chrysene-d12	21.20	240	1696990	20.00	ng/ul	0.00
85) Perylene-d12	23.39	264	1743729	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	21914	4.23	ng/uL	0.00
5) Phenol-d5	6.79	99	101741	4.70	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	298467	22.87	ng/ul	0.00
9) 2-Chlorophenol-d4	7.13	132	318640	18.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.30	113	198929	11.25	ng/ul	0.00
19) Nitrobenzene-d5	8.75	128	192134	23.35	ng/ul	0.00
22) 2-Nitrophenol-d4	9.46	143	213583	22.65	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.00	165	356076	19.84	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	0.00
43) Dimethylphthalate-d6	13.66	166	1372500	24.63	ng/ul	0.00
46) Acenaphthylene-d8	13.93	160	1620914	24.03	ng/ul	0.00
51) 4-Nitrophenol-d4	14.49	143	32921m	3.25	ng/ul	0.02
57) Fluorene-d10	15.24	176	1178115	24.48	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.37	200	205270	19.37	ng/ul	0.00
70) Anthracene-d10	17.10	188	1841761	24.95	ng/ul	0.00
78) Pyrene-d10	19.40	212	2111442	25.71	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.24	264	2365922	25.02	ng/ul	0.00

Target Compounds

Qvalue

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