

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

Data File : BM018058.D

Acq On : 11 Dec 2018 15:33

Operator : JU/SJ

Sample : J6137-13

Misc :

ALS Vial : 10 Sample Multiplier: 1

Quant Time: Dec 12 00:57:08 2018

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Dec 12 00:24:38 2018

Response via : Initial Calibration

Instrument :

BNA_M

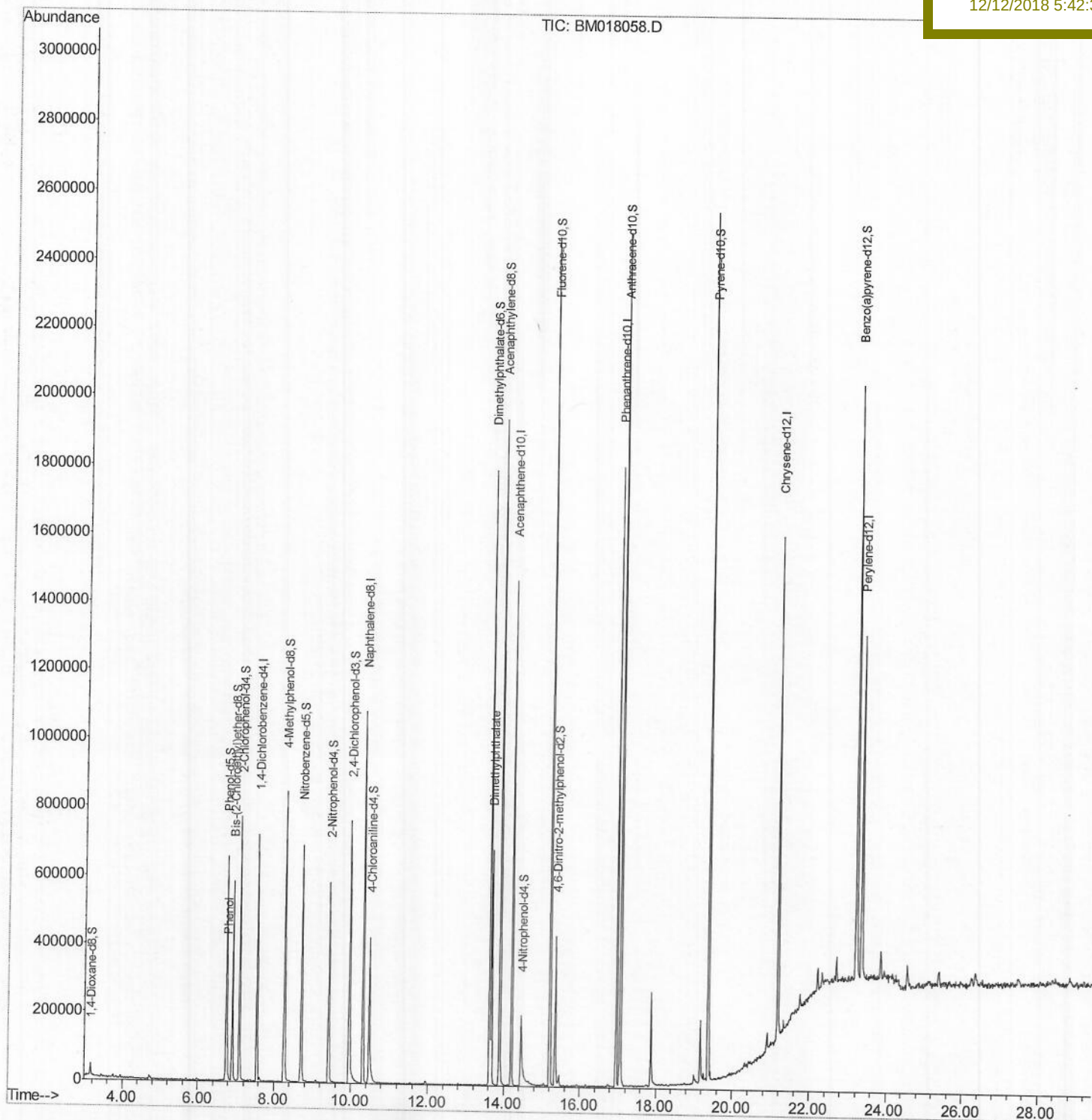
ClientSampleId :

F9J14

Manual Integrations
APPROVED

Sohil

12/12/2018 5:42:33 PM



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

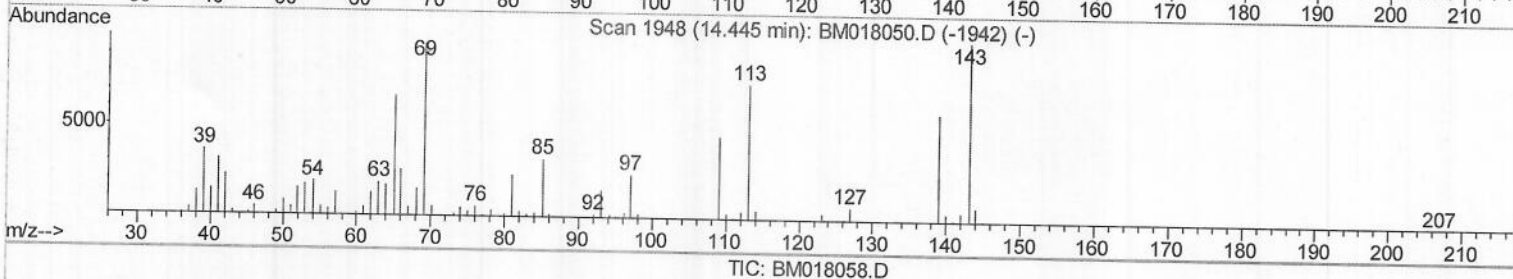
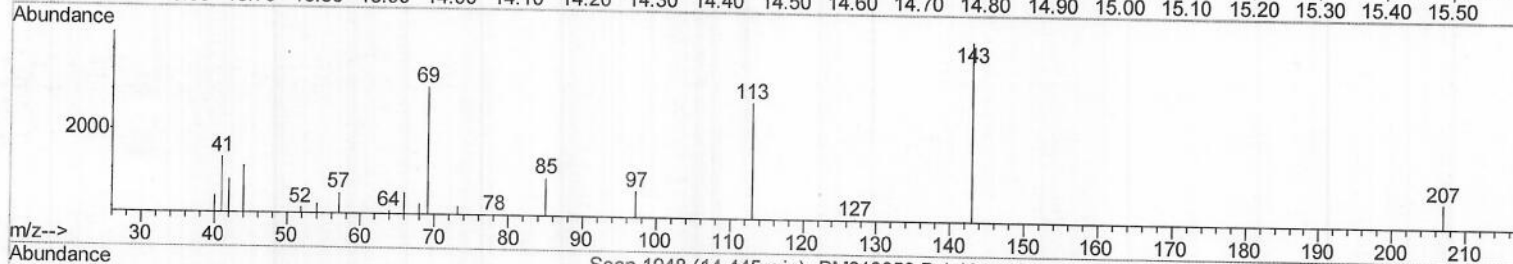
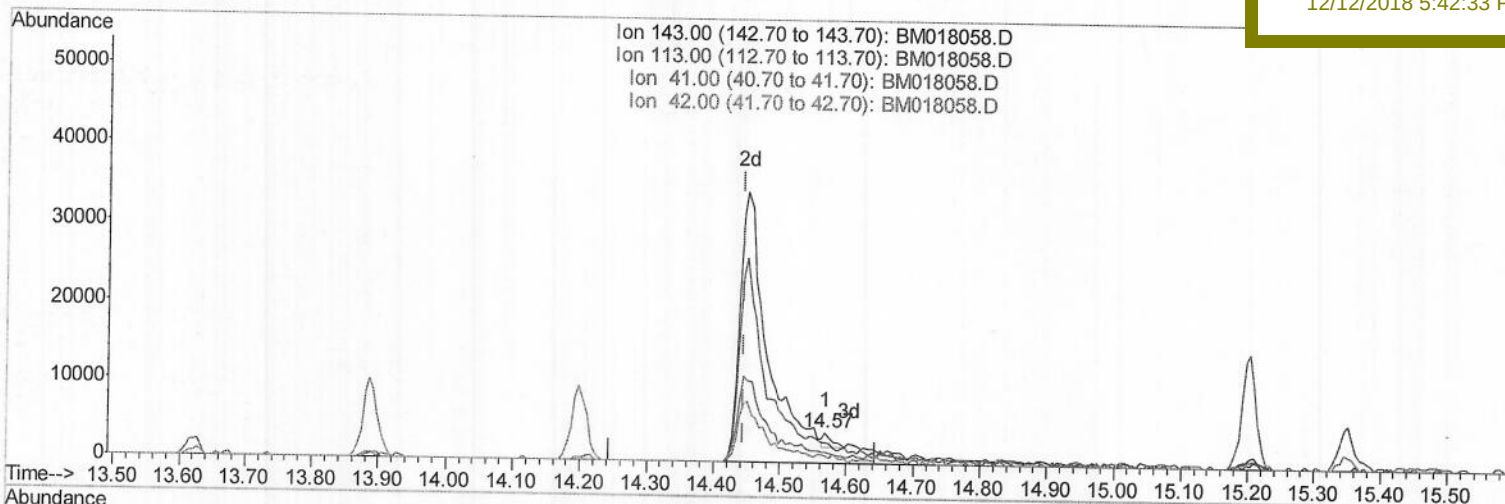
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(51) 4-Nitrophenol-d4 (S)

14.569min (+0.124) 0.23ng/ul

response 1737

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	67.68
41.00	40.10	36.79
42.00	26.80	25.33

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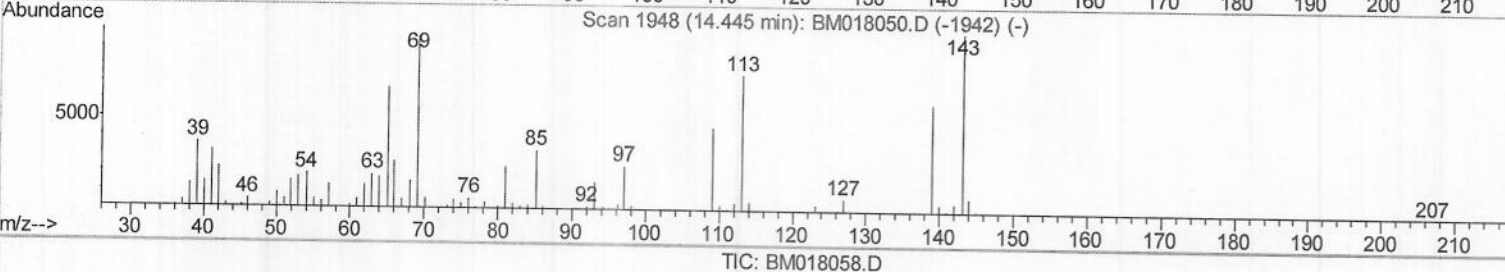
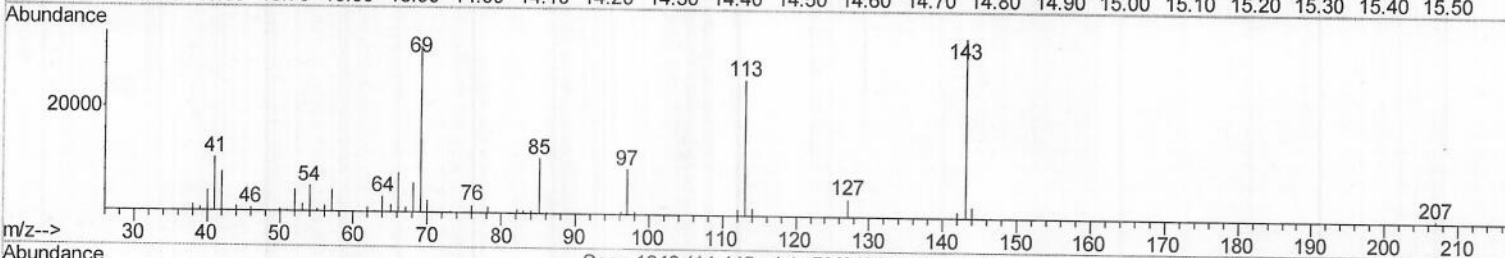
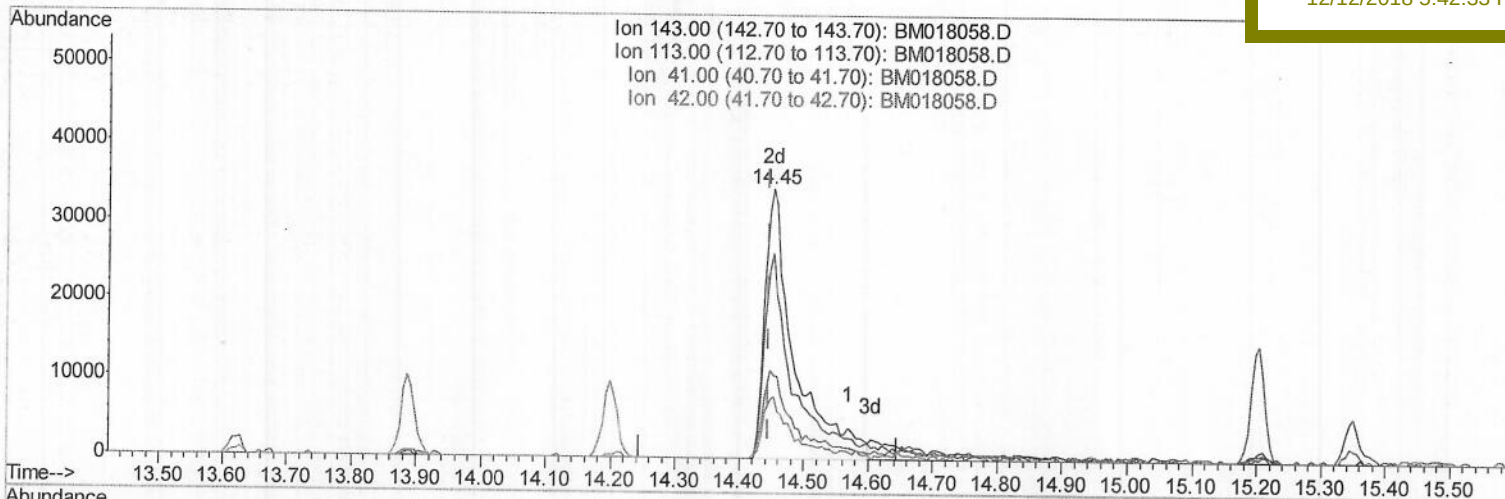
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Sohil
12/12/2018 5:42:33 PM



(51) 4-Nitrophenol-d4 (S)

14.451min (+0.006) 14.14ng/ul m

response 108942

Ion	Exp%	Act%
143.00	100	100
113.00	79.10	76.14
41.00	40.10	30.68#
42.00	26.80	22.54

15 J 12/13/18

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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	7.55	152	203584	20.00	ng/ul	0.00
18) Naphthalene-d8	10.32	136	884254	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.20	164	488535	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.96	188	981903	20.00	ng/ul	0.00
77) Chrysene-d12	21.17	240	738201	20.00	ng/ul	0.00
85) Perylene-d12	23.34	264	716939	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	19051	4.28	ng/uL	0.00
5) Phenol-d5	6.74	99	421089	22.65	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.90	67	247328	22.06	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	355566	24.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.26	113	345539	22.75	ng/ul	0.00
19) Nitrobenzene-d5	8.70	128	178941	26.35	ng/ul	0.00
22) 2-Nitrophenol-d4	9.42	143	202160	25.98	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.95	165	319952	21.60	ng/ul	0.00
29) 4-Chloroaniline-d4	10.47	131	310206	24.40	ng/ul	0.00
43) Dimethylphthalate-d6	13.62	166	1192265	28.13	ng/ul	0.00
46) Acenaphthylene-d8	13.89	160	1432876	27.93	ng/ul	0.00
51) 4-Nitrophenol-d4	14.45	143	108942m	14.14	ng/ul	0.00
57) Fluorene-d10	15.20	176	985435	26.92	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.35	200	110213	15.70	ng/ul	0.00
70) Anthracene-d10	17.06	188	1413693	28.92	ng/ul	0.00
78) Pyrene-d10	19.37	212	1346679	37.70	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.21	264	1171204	30.12	ng/ul	0.00

Target Compounds

					Qvalue
6) Phenol	6.77	94	113991	6.142	ng/ul 97
44) Dimethylphthalate	13.67	163	442752	10.781	ng/ul 99

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

SJ12/13/18