

Quantitation Report (Qedit)

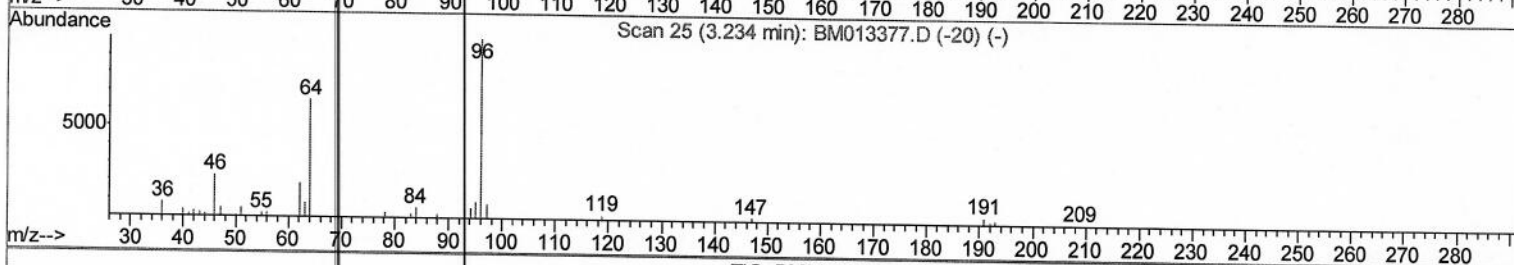
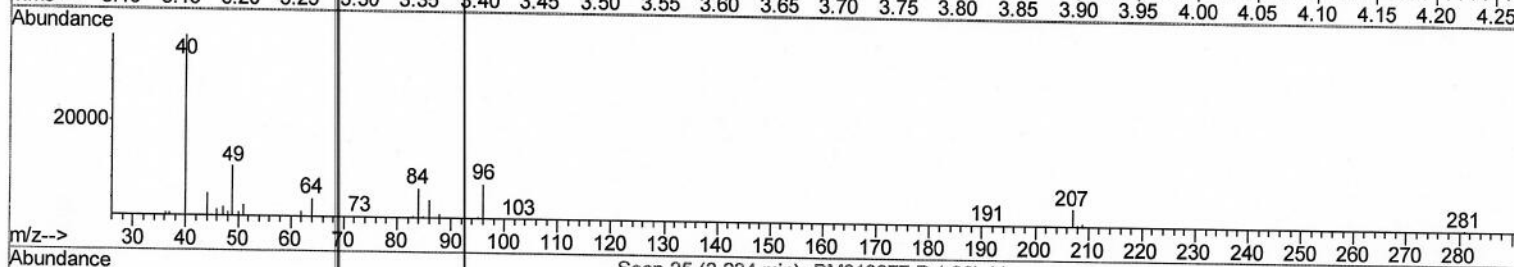
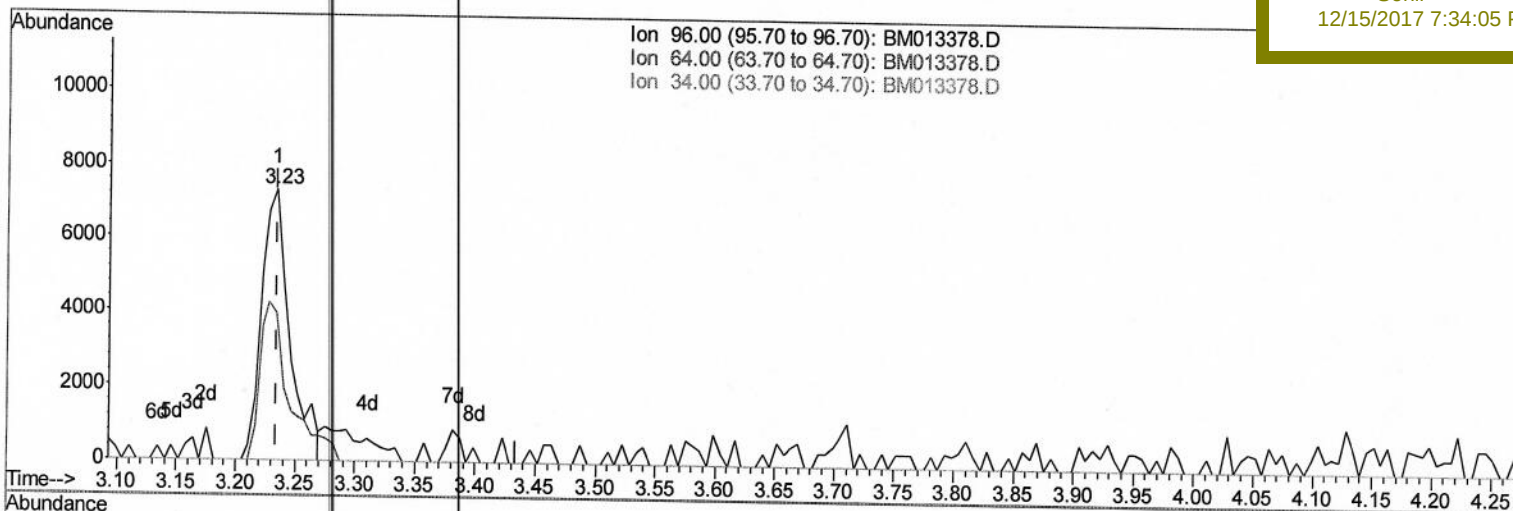
Data Path : Z:\HPCHEM1\BNA M\DATA\BM121317\
Data File : BM013378.D
Acq On : 13 Dec 2017 17:48
Operator : SJ/JU
Sample : PB104973BL
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Dec 14 01:19:55 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 14 01:15:39 2017
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
SBLK73

Manual Integrations
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TIC: BM013378.D

(3) 1,4-Dioxane-d8 (S)
3.234min (+0.000) 6.28ng/uL
response 11842

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	57.84#
34.00	0.00	0.00
0.00	0.00	0.00

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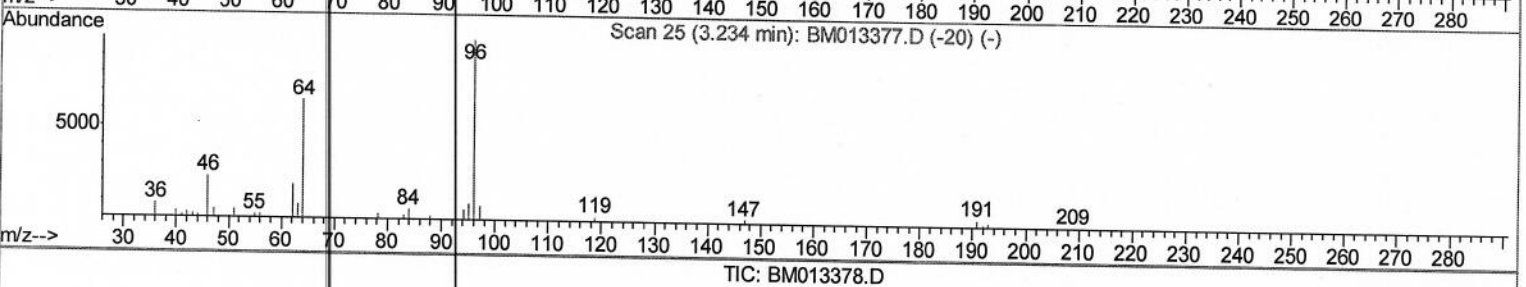
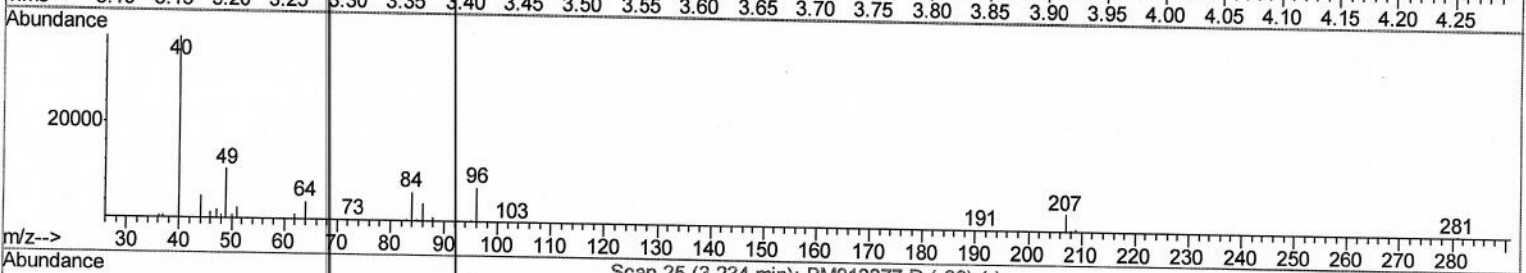
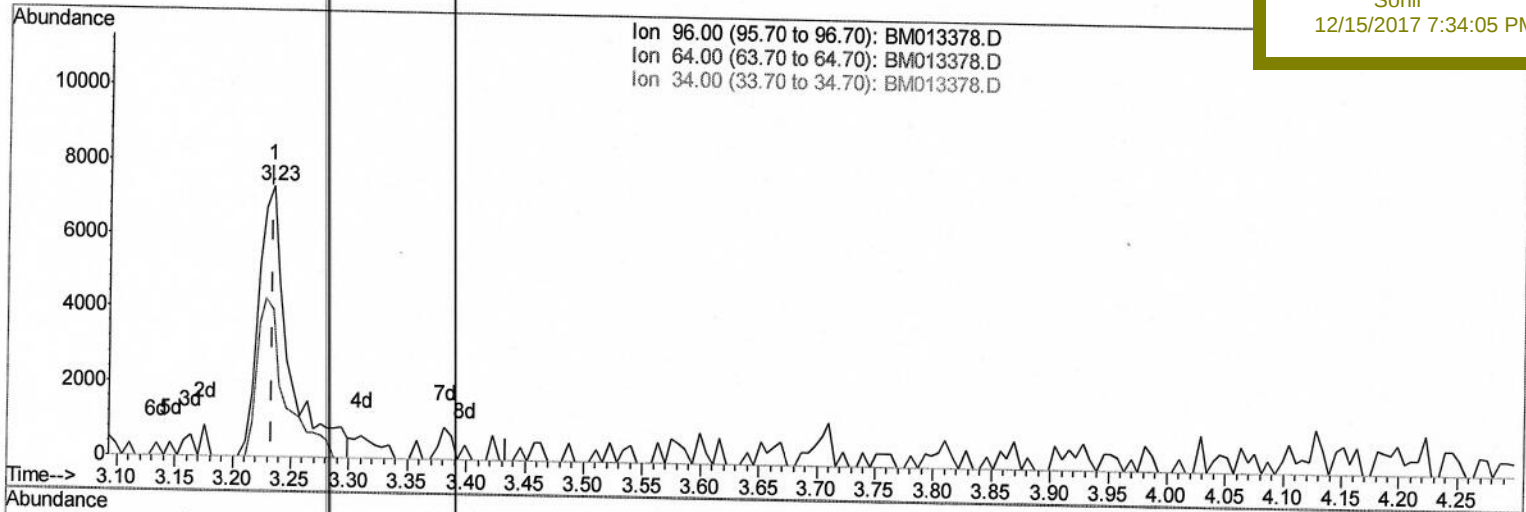
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(3) 1,4-Dioxane-d8 (S)

3.234min (+0.000) 7.00ng/uL m

response 13193

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	51.91#
34.00	0.00	0.00
0.00	0.00	0.00

SJ 12/13/17

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Quant Time: Dec 14 01:31:23 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 14 01:15:39 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	93389	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	410743	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.33	164	271300	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.07	188	756270	20.00	ng/ul	0.00
75) Chrysene-d12	21.26	240	1015046	20.00	ng/ul	0.00
83) Perylene-d12	23.49	264	936811	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.23	96	13193m	7.00	ng/uL	0.00
5) Phenol-d5	6.86	99	217531	26.92	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.02	67	128835	27.96	ng/ul	0.00
9) 2-Chlorophenol-d4	7.22	132	176926	28.02	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	181803	26.34	ng/ul	0.00
19) Nitrobenzene-d5	8.84	128	96396	29.59	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	99527	28.55	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	195235	27.18	ng/ul	0.00
29) 4-Chloroaniline-d4	10.60	131	275826	41.78	ng/ul	0.00
43) Dimethylphthalate-d6	13.74	166	722706	29.94	ng/ul	0.00
46) Acenaphthylene-d8	14.02	160	886241	29.75	ng/ul	0.00
51) 4-Nitrophenol-d4	14.54	143	115201	27.46	ng/ul	0.00
57) Fluorene-d10	15.32	176	655042	31.55	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.45	200	93693	19.69	ng/ul	0.00
70) Anthracene-d10	17.17	188	1227137	32.40	ng/ul	0.00
76) Pyrene-d10	19.47	212	1570533	31.06	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.35	264	1584200	33.20	ng/ul	0.00

Target Compounds Ovalue

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

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