

Quantitation Report (Qedit)

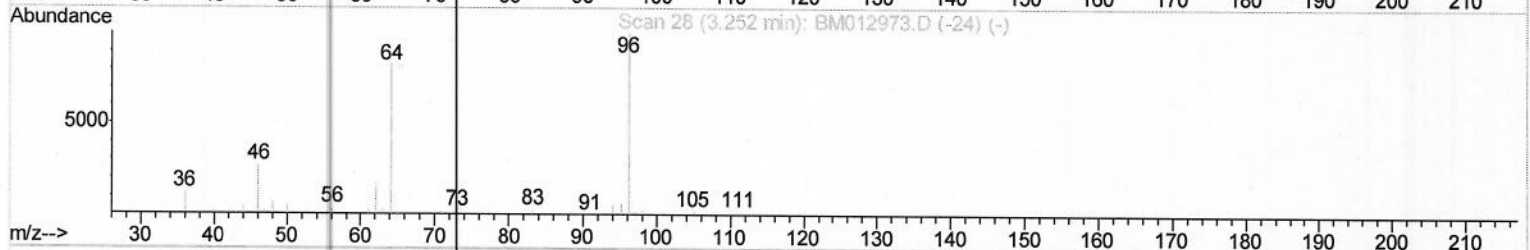
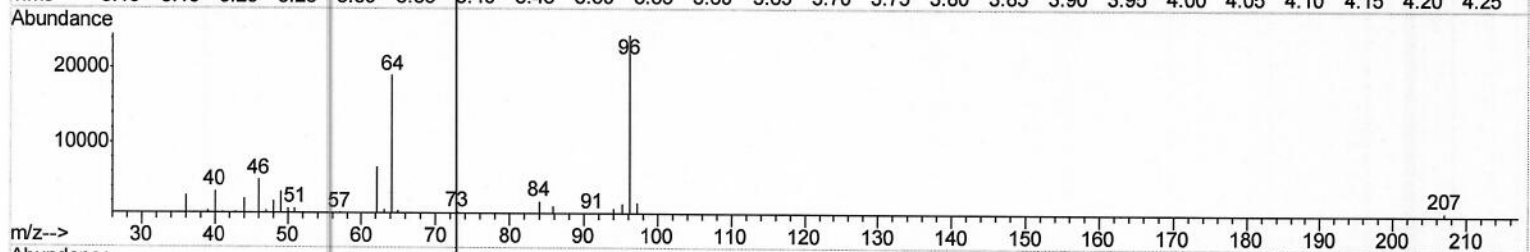
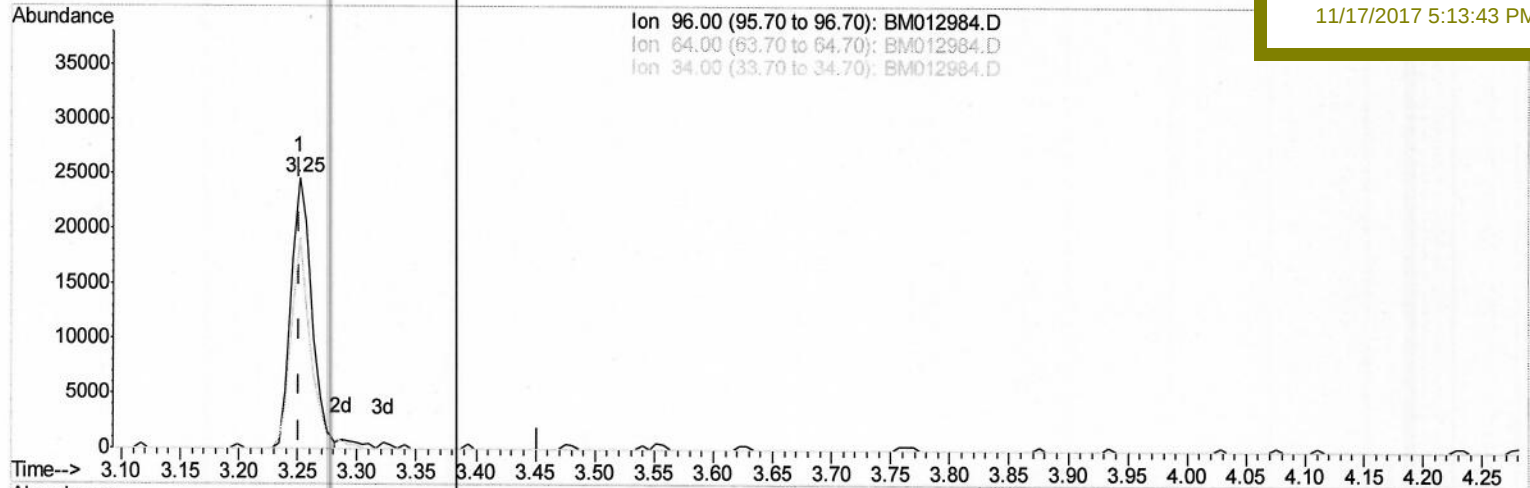
Data Path : Z:\HPCHEM1\BNA M\DATA\BM111617\
 Data File : BM012984.D
 Acq On : 17 Nov 2017 04:12
 Operator : SJ/JU
 Sample : I6361-09
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Quant Time: Nov 17 05:55:25 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111617.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 17 01:17:11 2017
 Response via : Initial Calibration

Instrument :
 BNA_M
 ClientSampleId :
 BE2R1

Manual Integrations
 APPROVED

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TIC: BM012984.D

(3) 1,4-Dioxane-d8 (S)
 3.252min (-0.000) 6.32ng/uL
 response 29904

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	74.36
34.00	0.00	0.00
0.00	0.00	0.00

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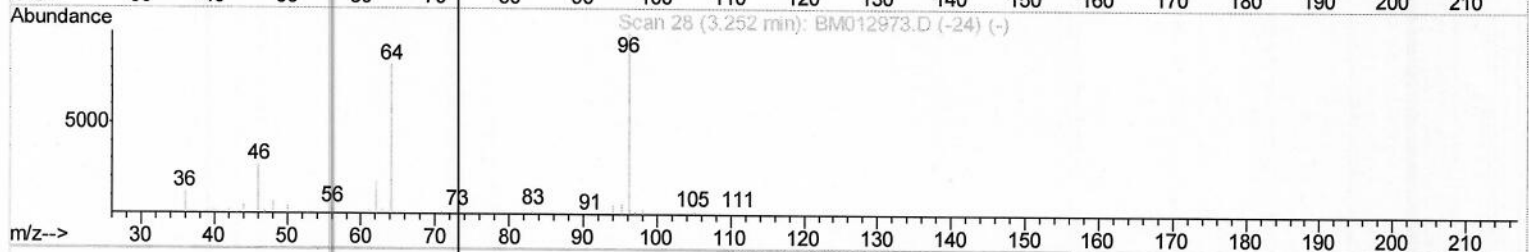
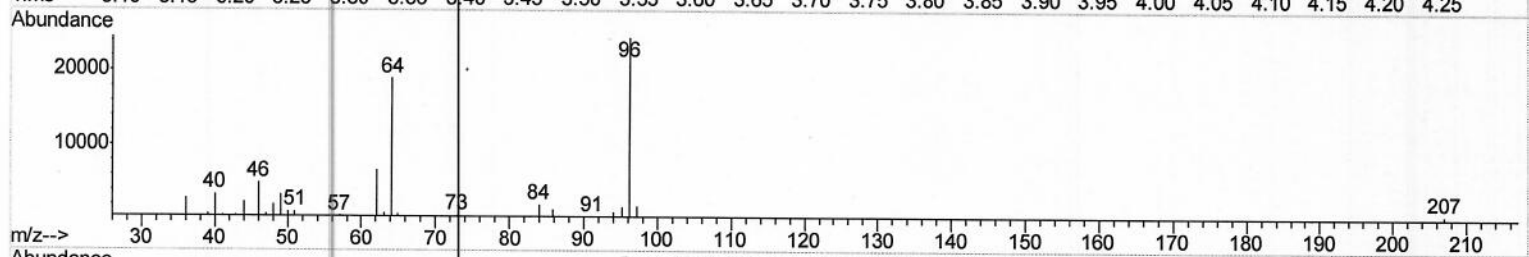
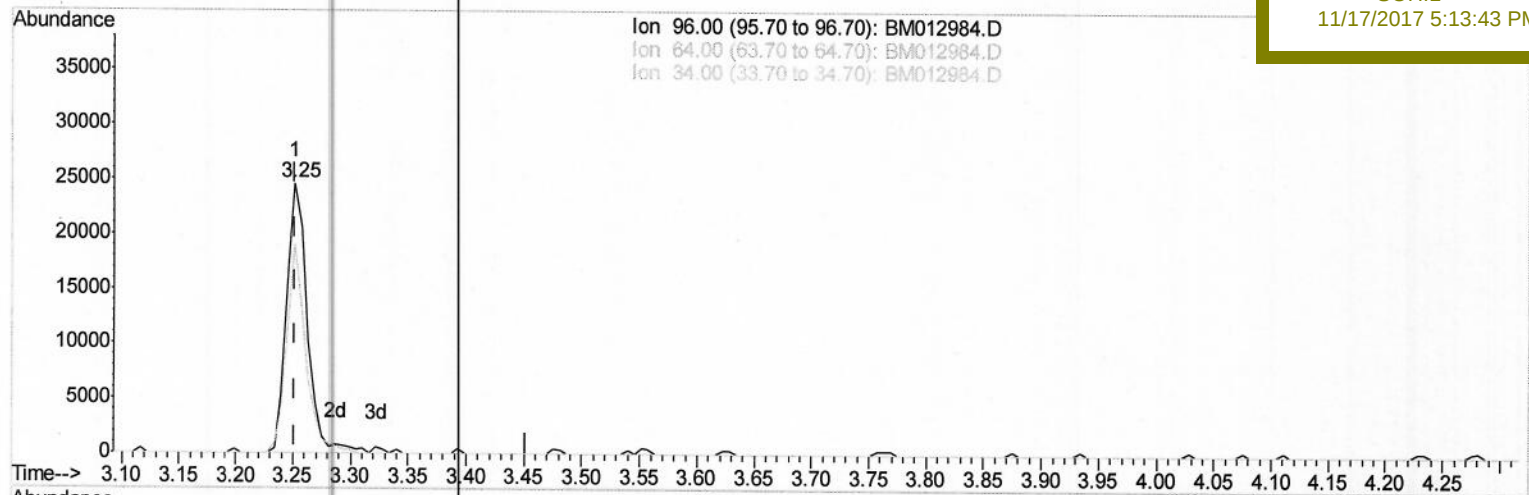
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TIC: BM012984.D

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

3.252min (-0.000) 6.54ng/uL m 25511/21117

response 30919

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	71.92
34.00	0.00	0.00
0.00	0.00	0.00

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Quant Time: Nov 17 06:11:40 2017
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	219034	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	897611	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.36	164	517340	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.11	188	1183924	20.00	ng/ul	0.00
75) Chrysene-d12	21.29	240	1324783	20.00	ng/ul	0.00
83) Perylene-d12	23.53	264	1338140	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	30919m)	6.54	ng/uL	0.00
5) Phenol-d5	6.89	99	105633	5.57	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.06	67	268483	24.94	ng/ul	0.00
9) 2-Chlorophenol-d4	7.26	132	298091	19.66	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	199113	12.81	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	175763	27.47	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	175272	28.80	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	331125	21.44	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	926	0.06	ng/ul	0.00
43) Dimethylphthalate-d6	13.78	166	1180483	26.01	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	1469620	25.42	ng/ul	0.00
51) 4-Nitrophenol-d4	14.55	143	34387	5.00	ng/ul	0.00
57) Fluorene-d10	15.36	176	1007218	26.12	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.47	200	129369	25.20	ng/ul	0.00
70) Anthracene-d10	17.21	188	1613470	26.62	ng/ul	0.00
76) Pyrene-d10	19.50	212	1779250	26.29	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.39	264	1860477	27.42	ng/ul	0.00

Target Compounds Ovalue

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

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Manual Integrations
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