

Quantitation Report (Qedit)

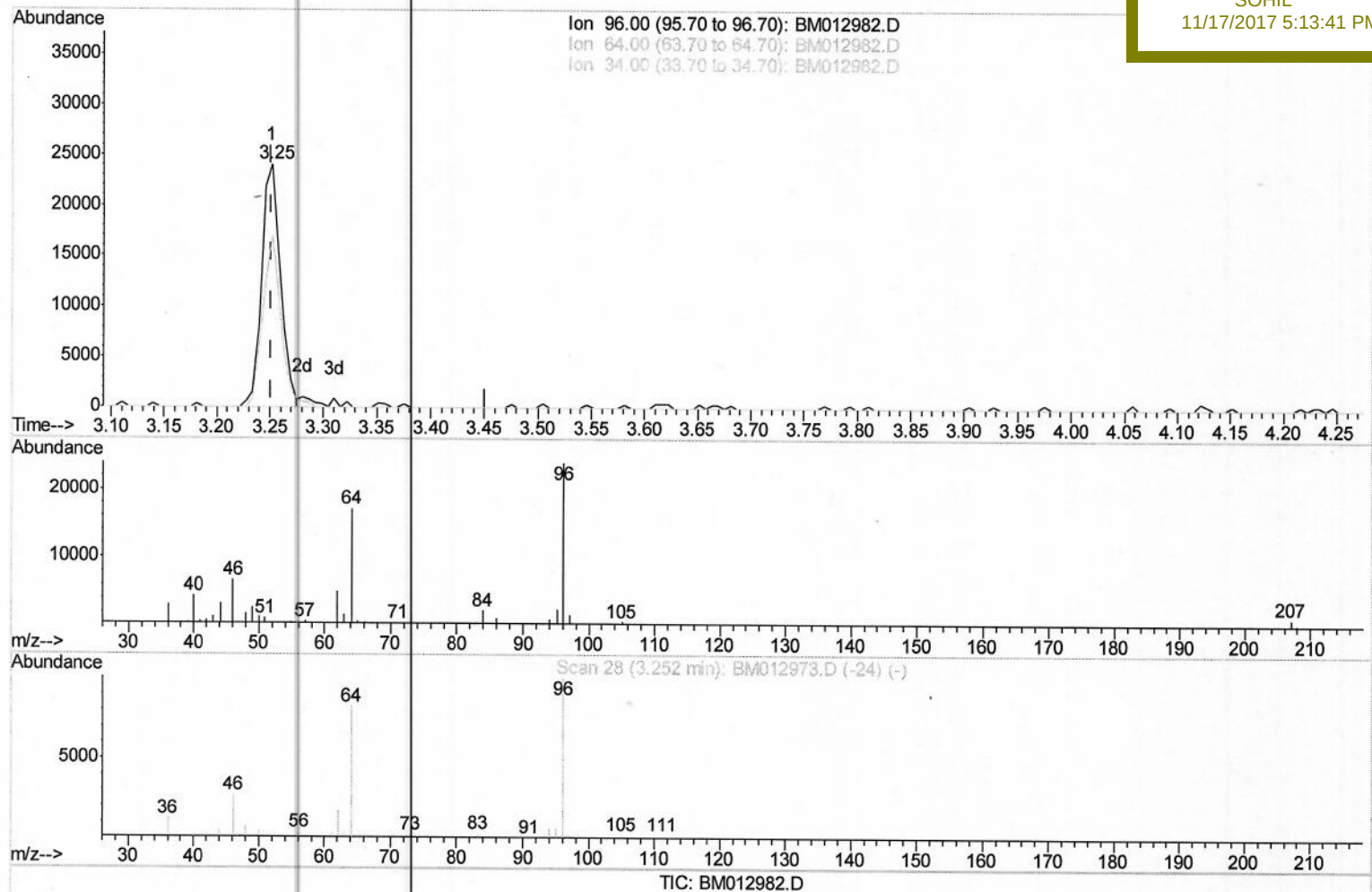
Data Path : Z:\HPCHEM1\BNA M\DATA\BM111617\
 Data File : BM012982.D
 Acq On : 17 Nov 2017 03:00
 Operator : SJ/JU
 Sample : I6361-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Nov 17 05:59:49 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111617.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 17 05:59:41 2017
 Response via : Initial Calibration

Instrument :
 BNA_M
ClientSampleId :
 BE2Q5

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

3.251min (-0.000) 7.25ng/uL

response 29416

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	68.81
34.00	0.00	0.00
0.00	0.00	0.00

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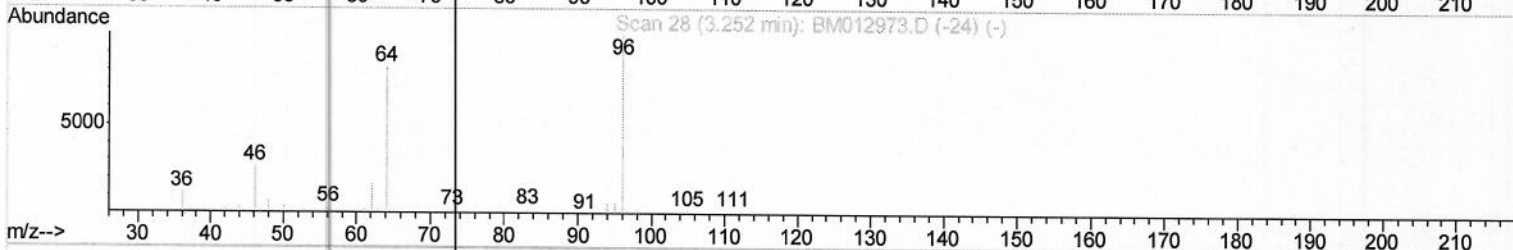
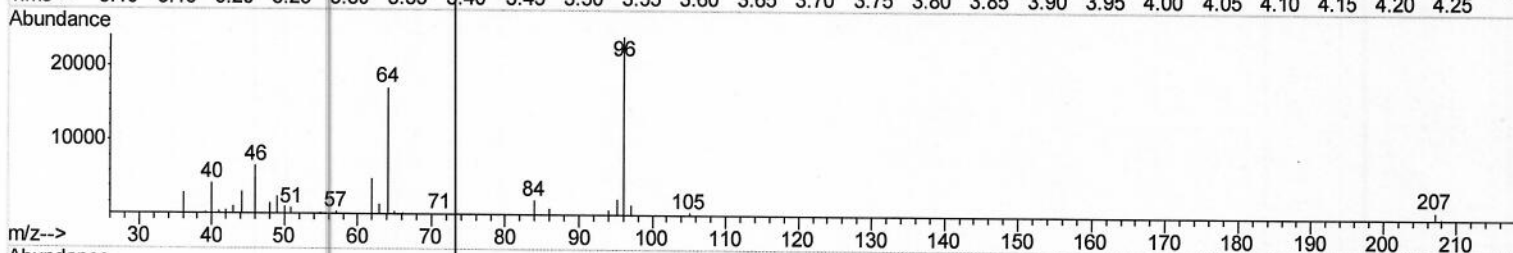
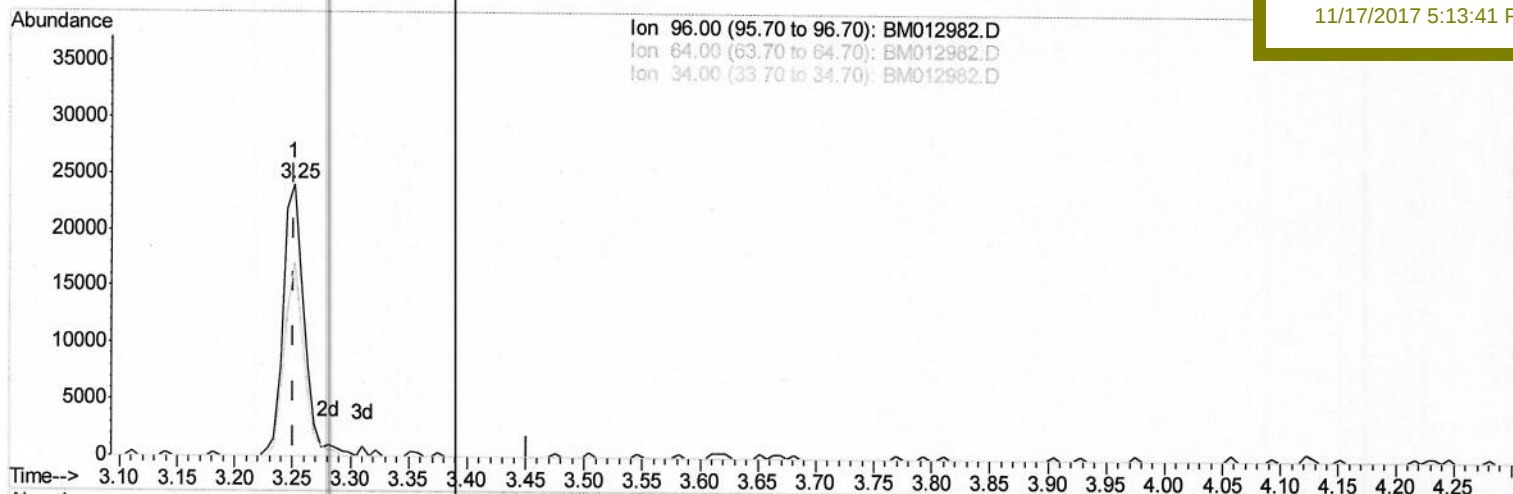
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TIC: BM012982.D

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

3.251min (-0.000) 7.48ng/uL m > 5511/21/17

response 30325

Ion	Exp%	Act%
96.00	100	100
64.00	74.90	66.74
34.00	0.00	0.00
0.00	0.00	0.00

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Quant Time: Nov 17 06:00:29 2017
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Internal Standards	R.T.	Q Ion	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	187900	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	751117	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	456509	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.11	188	1066950	20.00	ng/ul	0.00
75) Chrysene-d12	21.30	240	1213121	20.00	ng/ul	0.00
83) Perylene-d12	23.53	264	1223892	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.25	96	30325m)	7.48	ng/uL	0.00) SJ11/21/17
5) Phenol-d5	6.89	99	167325	10.28	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.06	67	131856	14.28	ng/ul	0.00
9) 2-Chlorophenol-d4	7.26	132	170903	13.14	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	150458	11.29	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	84740	15.83	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	83217	16.34	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.12	165	178826	13.83	ng/ul	0.00
29) 4-Chloroaniline-d4	10.64	131	145490	10.76	ng/ul	0.00
43) Dimethylphthalate-d6	13.78	166	621118	15.51	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	753502	14.77	ng/ul	0.00
51) 4-Nitrophenol-d4	14.55	143	54318	8.94	ng/ul	0.00
57) Fluorene-d10	15.36	176	530199	15.58	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.47	200	60247	13.02	ng/ul	0.00
70) Anthracene-d10	17.21	188	855886	15.67	ng/ul	0.00
76) Pyrene-d10	19.50	212	942386	15.21	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.39	264	982799	15.84	ng/ul	0.00
Target Compounds	Ovalue					

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

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