

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061217\
Quantitation Report (Qedit)

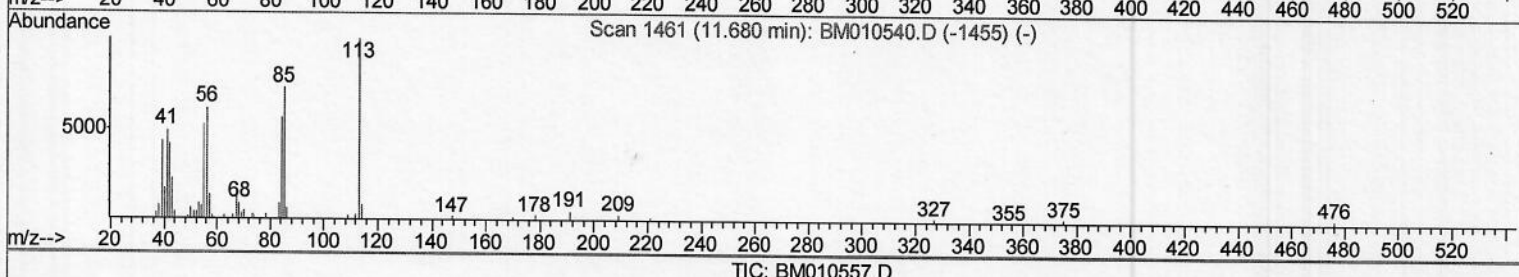
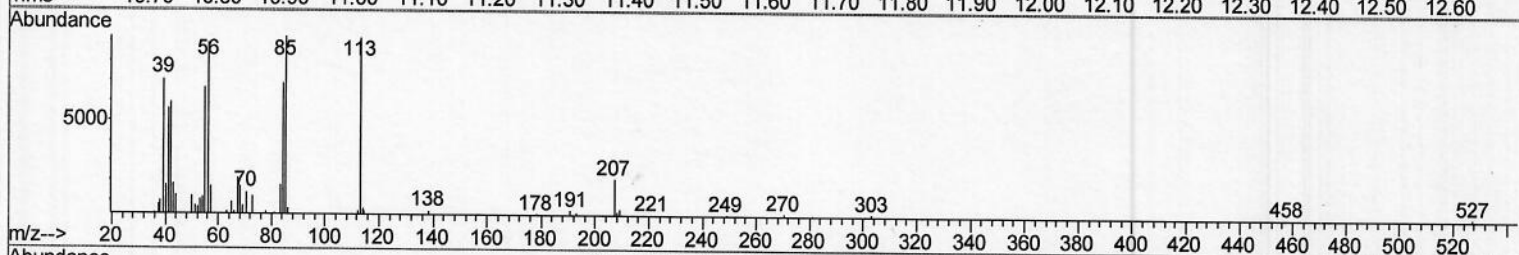
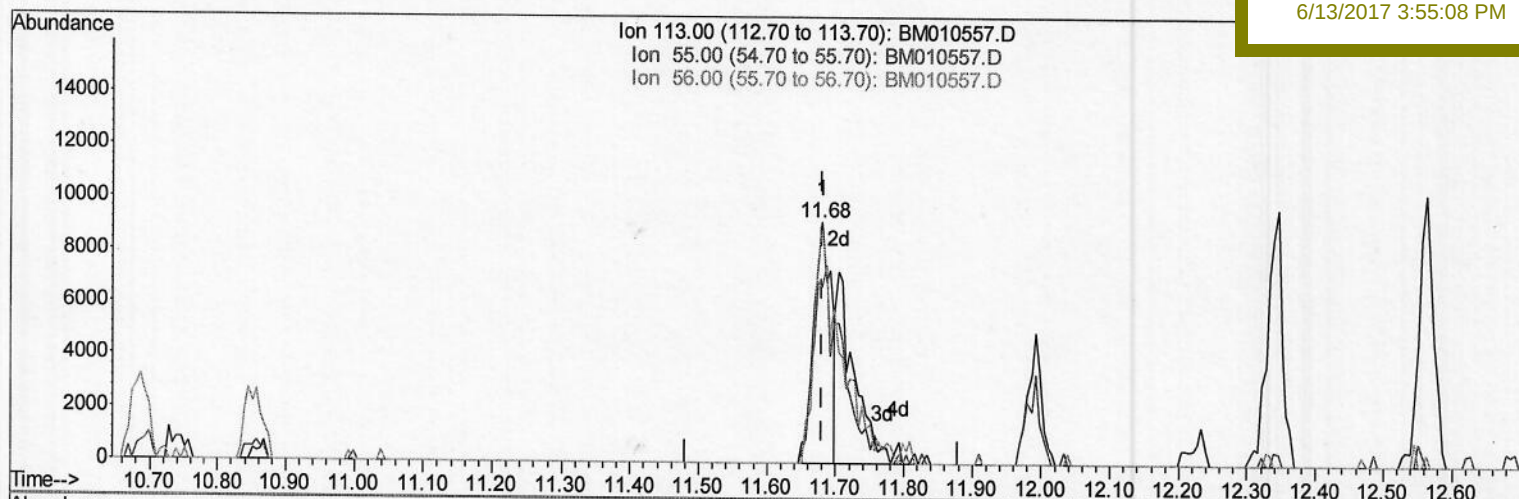
Data File : BM010557.D
Acq On : 13 Jun 2017 04:20
Operator : SJ/MA
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jun 13 06:33:14 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 13 03:33:28 2017
Response via : Initial Calibration

Instrument :
BNA_M
LabSampleId :
SSTD02028

Manual Integrations
APPROVED

mohammad
6/13/2017 3:55:08 PM



TIC: BM010557.D

(32) Caprolactam

11.680min (-0.000) 9.60ng/ul

response 16579

Ion	Exp%	Act%
113.00	100	100
55.00	118.50	72.52#
56.00	107.50	97.19
0.00	0.00	0.00

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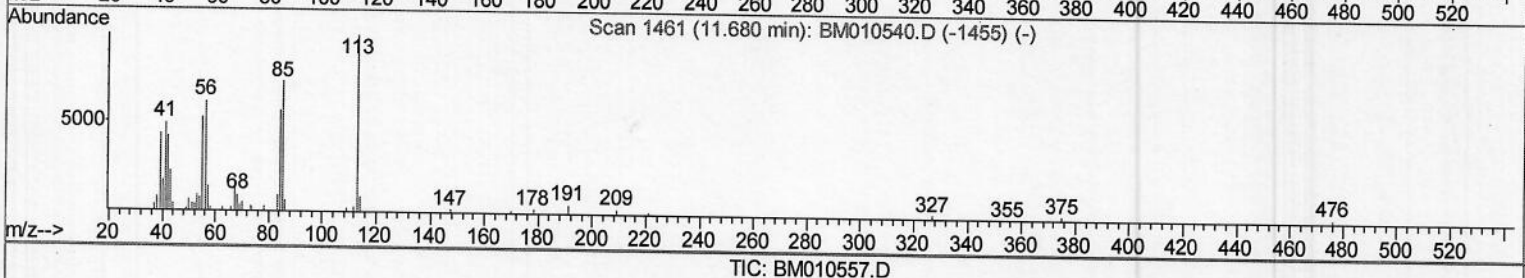
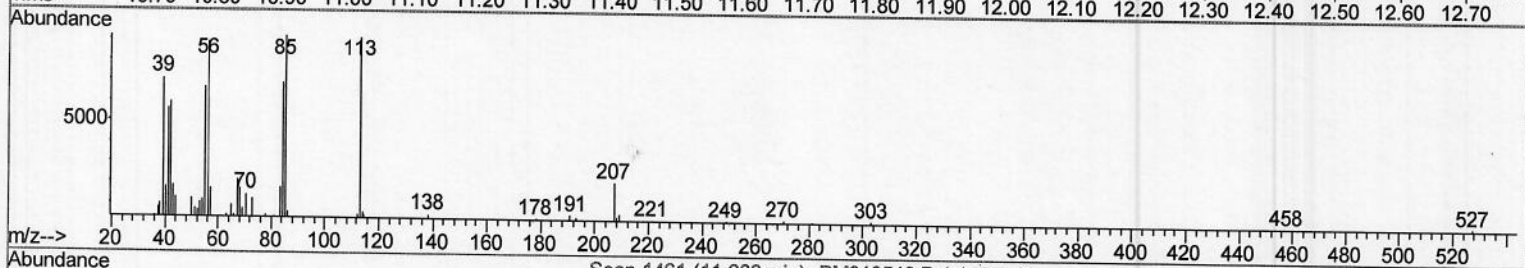
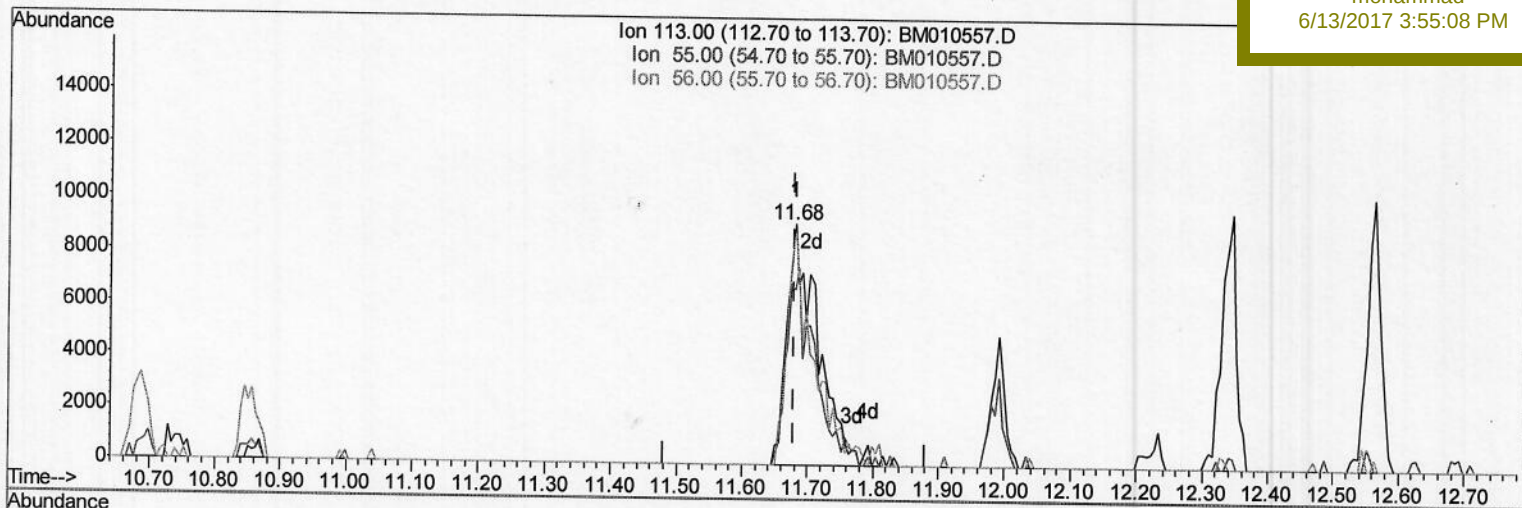
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(32) Caprolactam

11.680min (-0.000) 16.95ng/ul m

response 29278

Ion Exp% Act%

113.00 100 100

55.00 118.50 72.52#

56.00 107.50 97.19

0.00 0.00 0.00

> 5.7
06/13/17

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Acq On : 13 Jun 2017 04:20

Operator : SJ/MA

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Misc :

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Quant Time: Jun 13 06:34:36 2017

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	109797	20.00	ng/ul	0.00
18) Naphthalene-d8	10.69	136	404123	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.52	164	288482	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.27	188	741840	20.00	ng/ul	0.00
75) Chrysene-d12	21.44	240	1054026	20.00	ng/ul	0.00
83) Perylene-d12	23.77	264	1077362	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.36	96	20302	7.59	ng/uL	0.00
5) Phenol-d5	7.07	99	139565	16.77	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.23	67	83311	17.82	ng/ul	0.00
9) 2-Chlorophenol-d4	7.42	132	123131	18.57	ng/ul	0.00
13) 4-Methylphenol-d8	8.61	113	122395	18.23	ng/ul	0.00
19) Nitrobenzene-d5	9.07	128	54271	18.33	ng/ul	0.00
22) 2-Nitrophenol-d4	9.79	143	70247	20.49	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.31	165	159688	22.67	ng/ul	0.00
29) 4-Chloroaniline-d4	10.86	131	149831	22.04	ng/ul	0.00
43) Dimethylphthalate-d6	13.94	166	509689	21.26	ng/ul	0.00
46) Acenaphthylene-d8	14.22	160	554094	19.82	ng/ul	0.00
51) 4-Nitrophenol-d4	14.77	143	59709	16.65	ng/ul	0.00
57) Fluorene-d10	15.52	176	445108	21.13	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.64	200	98984	18.84	ng/ul	0.00
70) Anthracene-d10	17.37	188	726155	20.12	ng/ul	0.00
76) Pyrene-d10	19.66	212	916068	21.02	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.62	264	1015581	20.25	ng/ul	0.00

Target Compounds

					Qvalue	
2) 1,4-Dioxane	3.39	88	19073	6.69	ng/uL#	41
4) Benzaldehyde	7.05	77	115337	20.61	ng/ul	91
6) Phenol	7.10	94	140249	16.44	ng/ul#	79
8) Bis(2-Chloroethyl) ether	7.32	93	97977	16.55	ng/ul	89
10) 2-Chlorophenol	7.46	128	122317	18.38	ng/ul	94
11) 2-Methylphenol	8.34	108	112665	18.09	ng/ul	98
12) 2,2'-oxybis(1-Chloropropan	8.41	45	40657	15.23	ng/ul#	58
14) Acetophenone	8.72	105	194926	17.79	ng/ul	91
15) N-Nitroso-di-n-propylamine	8.69	70	106050	19.37	ng/ul#	89
16) 4-Methylphenol	8.67	108	115863	16.96	ng/ul	84
17) Hexachloroethane	8.95	117	65121	21.55	ng/ul	88
20) Nitrobenzene	9.12	77	175188	20.72	ng/ul#	89
21) Isophorone	9.62	82	271869	20.90	ng/ul#	90
23) 2-Nitrophenol	9.82	139	74011	20.35	ng/ul	92
24) 2,4-Dimethylphenol	9.86	107	172241	20.38	ng/ul	96
25) Bis(2-Chloroethoxy) methane	10.10	93	135854	18.58	ng/ul	91
27) 2,4-Dichlorophenol	10.34	162	155141	22.48	ng/ul#	86
28) Naphthalene	10.74	128	387252	19.10	ng/ul	99
30) 4-Chloroaniline	10.87	127	142511	21.73	ng/ul	90
31) Hexachlorobutadiene	10.98	225	158850	25.12	ng/ul	95
32) Caprolactam	11.68	113	29278m	16.95	ng/ul	
33) 4-Chloro-3-methylphenol	11.99	107	138135	20.79	ng/ul	91
34) 2-Methylnaphthalene	12.34	142	317064	20.63	ng/ul	99

5.7
06/23/17

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
36) 1,2,4,5-Tetrachlorobenzene	12.70	216	260859	21.74	ng/ul#	94
37) Hexachlorocyclopentadiene	12.66	237	141212	17.55	ng/ul#	94
38) 2,4,6-Trichlorophenol	12.96	196	146697	21.51	ng/ul	97
39) 2,4,5-Trichlorophenol	13.03	196	151340	22.00	ng/ul	98
40) 1,1'-Biphenyl	13.36	154	468167	20.00	ng/ul	98
41) 2-Chloronaphthalene	13.40	162	358912	19.52	ng/ul	94
42) 2-Nitroaniline	13.64	65	98903	21.12	ng/ul	92
44) Dimethylphthalate	13.99	163	482993	20.49	ng/ul#	96
45) 2,6-Dinitrotoluene	14.12	165	87322	20.21	ng/ul#	64
47) Acenaphthylene	14.25	152	546623	20.01	ng/ul	98
48) 3-Nitroaniline	14.47	138	70035	16.82	ng/ul	86
49) Acenaphthene	14.59	153	366687	19.13	ng/ul	97
50) 2,4-Dinitrophenol	14.66	184	61452	18.82	ng/ul	94
52) 4-Nitrophenol	14.77	109	119603	24.34	ng/ul#	73
53) Dibenzofuran	14.92	168	589250	20.79	ng/ul	93
54) 2,4-Dinitrotoluene	14.91	165	137707	21.70	ng/ul#	72
55) 2,3,4,6-Tetrachlorophenol	15.14	232	154246	23.34	ng/ul#	92
56) Diethylphthalate	15.33	149	480843	21.09	ng/ul	98
58) Fluorene	15.57	166	489217	21.05	ng/ul	96
59) 4-Chlorophenyl-phenylether	15.56	204	305566	23.00	ng/ul	95
60) 4-Nitroaniline	15.64	138	77464	17.77	ng/ul#	77
63) 4,6-Dinitro-2-methylphenol	15.66	198	104129	18.70	ng/ul	98
64) N-Nitrosodiphenylamine	15.79	169	409603	18.94	ng/ul	99
65) 4-Bromophenyl-phenylether	16.46	248	192631	21.10	ng/ul	92
66) Hexachlorobenzene	16.55	284	181371	18.97	ng/ul#	79
67) Atrazine	16.73	200	204030	21.71	ng/ul	94
68) Pentachlorophenol	16.92	266	115154	18.93	ng/ul	99
69) Phenanthrene	17.32	178	780519	19.71	ng/ul	99
71) Anthracene	17.40	178	832074	20.13	ng/ul	97
72) Carbazole	17.69	167	655921	18.74	ng/ul	98
73) Di-n-butylphthalate	18.21	149	782881	19.82	ng/ul	98
74) Fluoranthene	19.32	202	1068999	21.95	ng/ul#	95
77) Pyrene	19.69	202	1125493	20.72	ng/ul#	93
78) Butylbenzylphthalate	20.56	149	374880	19.97	ng/ul	98
79) 3,3'-Dichlorobenzidine	21.37	252	397343	19.20	ng/ul	96
80) Benzo(a)anthracene	21.42	228	1262893	21.25	ng/ul	99
81) Bis(2-ethylhexyl)phthalate	21.30	149	554782	19.86	ng/ul#	98
82) Chrysene	21.47	228	1098906	20.45	ng/ul	98
84) Di-n-octyl phthalate	22.20	149	1008062	20.33	ng/ul#	96
85) Benzo(b)fluoranthene	23.06	252	1297624	20.70	ng/ul#	98
86) Benzo(k)fluoranthene	23.10	252	1247699	20.83	ng/ul#	98
88) Benzo(a)pyrene	23.67	252	1261471	20.30	ng/ul#	98
89) Indeno(1,2,3-cd)pyrene	26.14	276	1557065	19.83	ng/ul#	98
90) Dibenzo(a,h)anthracene	26.16	278	1332126	19.70	ng/ul	98
91) Benzo(g,h,i)perylene	26.89	276	1270237	19.64	ng/ul#	97

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

(QT Reviewed)

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6/13/2017 3:55:08 PM

