

Instrument :
BNA_M
ClientSampleId :
SSTD010014

mohammad
6/7/2021 12:27:25 PM

Abundance

TIC: BM030298.D



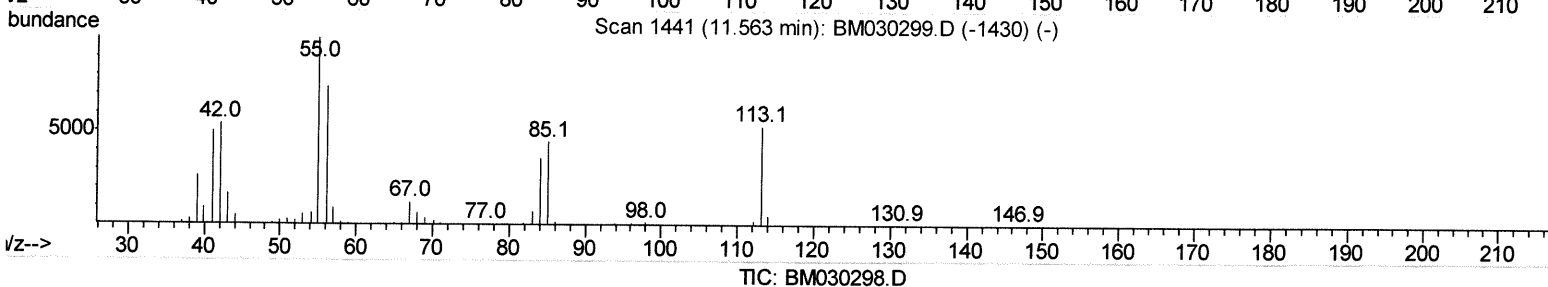
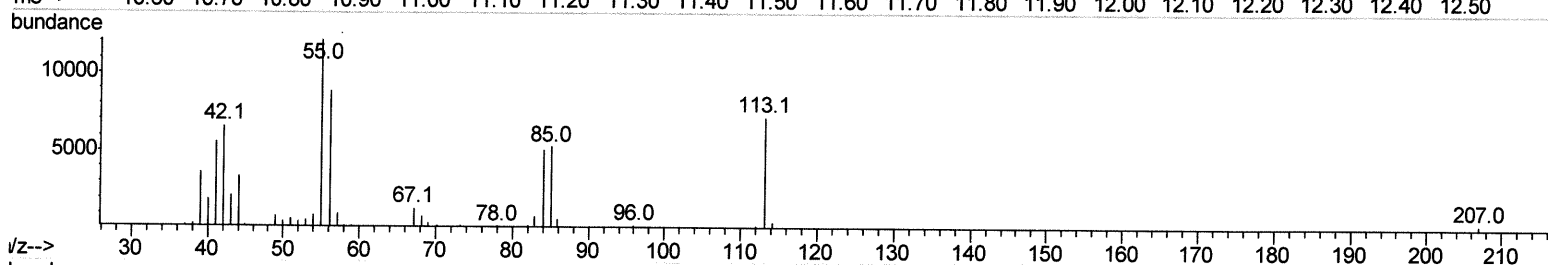
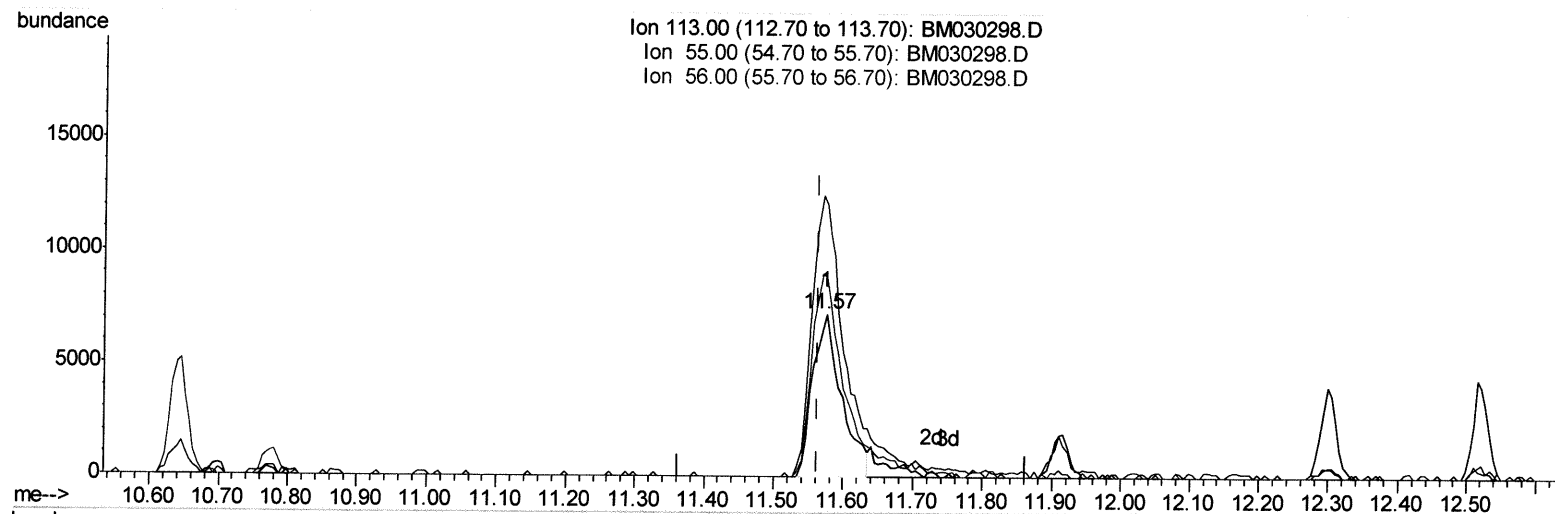
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060621\
Data File : BM030298.D
Acq On : 04 Jun 2021 16:16
Operator : CG/JU
Sample : SSTD01014
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Time: Jun 04 17:31:11 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 04 17:27:19 2021
Response via : Initial Calibration

Manual Integrations
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(34) Caprolactam

11.575min (+0.012) 6.26ng/ul

response 20858

Ion	Exp%	Act%
113.00	100	100
55.00	187.90	167.54
56.00	138.60	122.66
0.00	0.00	0.00

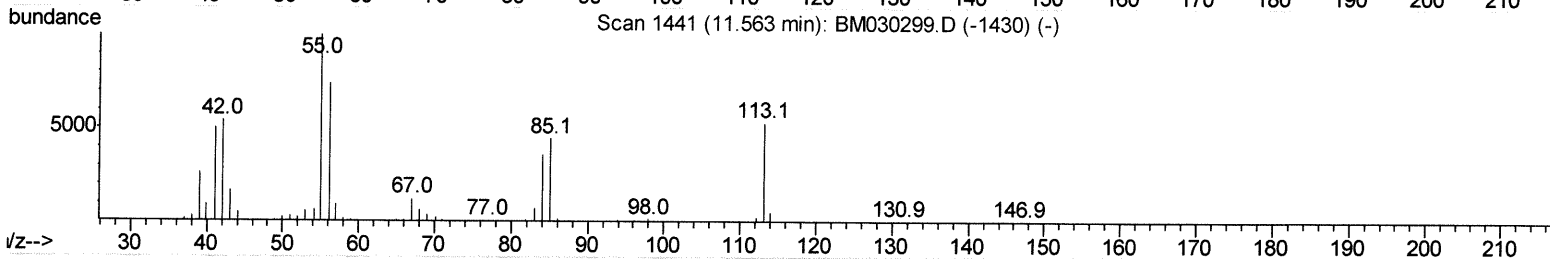
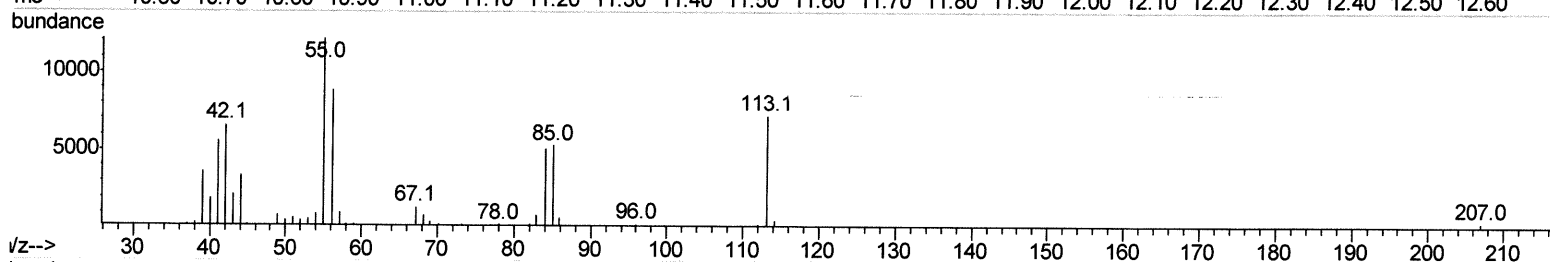
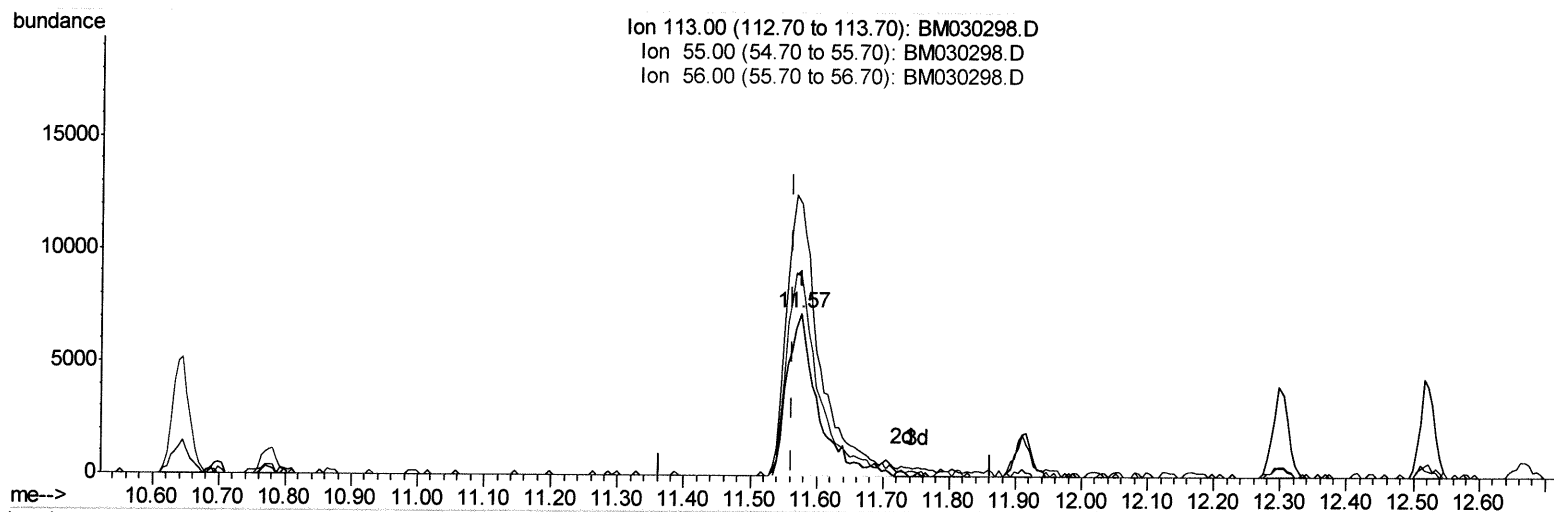
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(34) Caprolactam

11.575min (+0.012) 6.95ng/ul m

response 23172

JU 6/23/21

Ion	Exp%	Act%
113.00	100	100
55.00	187.90	167.54
56.00	138.60	122.66
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	174601	20.00	ng/ul	0.00
20) Naphthalene-d8	10.65	136	704777	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.47	164	428039	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.21	188	825218	20.00	ng/ul	0.00
79) Chrysene-d12	21.37	240	795809	20.00	ng/ul	0.00
88) Perylene-d12	23.65	264	767498	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.32	96	19871	4.66	ng/uL	0.00
4) Pividine-d5	3.74	84	96948	8.11	ng/ul	0.01
7) Phenol-d5	7.01	99	126536	8.24	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.18	67	93398	9.64	ng/ul	0.00
11) 2-Chlorophenol-d4	7.38	132	104908	9.12	ng/ul	0.00
15) 4-Methylphenol-d8	8.55	113	99262	8.14	ng/ul	0.00
21) Nitrobenzene-d5	9.00	128	45389	9.87	ng/ul	0.00
24) 2-Nitrophenol-d4	9.73	143	42464	9.69	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.26	165	99810	8.83	ng/ul	0.00
31) 4-Chloroaniline-d4	10.77	131	137837	8.21	ng/ul	0.00
46) Dimethylphthalate-d6	13.88	166	293181	8.96	ng/ul	0.00
49) Acenaphthylene-d8	14.17	160	354609	8.41	ng/ul	0.00
54) 4-Nitrophenol-d4	14.65	143	40703	7.70	ng/ul	0.00
60) Fluorene-d10	15.46	176	259567	9.09	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.57	200	34738	8.68	ng/ul	0.00
73) Anthracene-d10	17.31	188	377248	9.24	ng/ul	0.00
81) Pyrene-d10	19.59	212	400197	9.38	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.50	264	372155	8.55	ng/ul	0.00

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.36	88	20083	4.377	ng/uL 94
5) Pividine	3.76	79	110454	8.993	ng/ul 96
6) Benzaldehyde	6.99	77	62419	7.620	ng/ul 99
8) Phenol	7.03	94	133209	8.542	ng/ul 100
10) Bis(2-Chloroethyl)ether	7.28	93	116942	9.342	ng/ul 97
12) 2-Chlorophenol	7.42	128	109210	9.274	ng/ul 99
13) 2-Methylphenol	8.29	108	95745	8.119	ng/ul 98
14) 2,2'-oxybis(1-Chloropropan	8.38	45	191958	9.228	ng/ul 97
16) Acetophenone	8.67	105	167566	8.505	ng/ul 99
17) N-Nitroso-di-n-propylamine	8.65	70	84877	8.167	ng/ul 97
18) 4-Methylphenol	8.61	108	107088	8.434	ng/ul 97
19) Hexachloroethane	8.93	117	47264	9.374	ng/ul 98
22) Nitrobenzene	9.05	77	121967	9.794	ng/ul 99
23) Isophorone	9.57	82	210801	8.028	ng/ul 98
25) 2-Nitrophenol	9.76	139	48269	9.757	ng/ul 97
26) 2,4-Dimethylphenol	9.81	107	118069	8.971	ng/ul 100
27) Bis(2-Chloroethoxy)methane	10.05	93	149508	9.289	ng/ul 98
29) 2,4-Dichlorophenol	10.29	162	99941	9.160	ng/ul 98
30) Naphthalene	10.69	128	366571	9.480	ng/ul 98
32) 4-Chloroaniline	10.80	127	140823	8.102	ng/ul 99
33) Hexachlorobutadiene	10.98	225	67804	9.121	ng/ul 99
34) Caprolactam	11.57	113	23172m	6.954	ng/ul > 74 6/23/21

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35) 4-Chloro-3-methylphenol	11.91	107	94577	8.270	ng/ul	99
36) 2-Methylnaphthalene	12.30	142	249809	8.640	ng/ul	98
37) 1-Methylnaphthalene	12.52	142	254090	9.040	ng/ul	98
39) 1,2,4,5-Tetrachlorobenzene	12.67	216	132146	9.515	ng/ul	96
40) Hexachlorocyclopentadiene	12.65	237	73601	7.563	ng/ul	98
41) 2,4,6-Trichlorophenol	12.90	196	71157	9.129	ng/ul	99
42) 2,4,5-Trichlorophenol	12.97	196	79035	9.372	ng/ul	97
43) 1,1'-Biphenyl	13.31	154	321625	9.422	ng/ul	98
44) 2-Chloronaphthalene	13.35	162	247792	9.443	ng/ul	99
45) 2-Nitroaniline	13.55	65	50168	8.323	ng/ul	98
47) Dimethylphthalate	13.93	163	285453	8.831	ng/ul	100
48) 2,6-Dinitrotoluene	14.05	165	43437	8.432	ng/ul	98
50) Acenaphthylene	14.20	152	347891	8.540	ng/ul	100
51) 3-Nitroaniline	14.37	138	41860	6.972	ng/ul	96
52) Acenaphthene	14.54	153	263738	9.166	ng/ul	99
53) 2,4-Dinitrophenol	14.58	184	22720	7.025	ng/ul	90
55) 4-Nitrophenol	14.67	109	34716	5.979	ng/ul	99
56) Dibenzofuran	14.87	168	370064	9.292	ng/ul	98
57) 2,4-Dinitrotoluene	14.83	165	62337	8.343	ng/ul#	88
58) 2,3,4,6-Tetrachlorophenol	15.10	232	60328	8.472	ng/ul	96
59) Diethylphthalate	15.29	149	273977	8.368	ng/ul	98
61) Fluorene	15.52	166	300186	8.884	ng/ul	97
62) 4-Chlorophenyl-phenylether	15.52	204	147047	8.875	ng/ul	99
63) 4-Nitroaniline	15.53	138	35600	5.766	ng/ul	99
66) 4,6-Dinitro-2-methylphenol	15.59	198	35928	8.706	ng/ul	93
67) N-Nitrosodiphenylamine	15.73	169	245017	9.103	ng/ul	99
68) 4-Bromophenyl-phenylether	16.40	248	80108	8.625	ng/ul	99
69) Hexachlorobenzene	16.52	284	93101	8.769	ng/ul	100
70) Atrazine	16.67	200	74068	7.672	ng/ul	98
71) Pentachlorophenol	16.86	266	43506	6.710	ng/ul	94
72) Phenanthrene	17.25	178	460304	9.518	ng/ul	99
74) Anthracene	17.34	178	455785	9.188	ng/ul	98
75) 1,2,3,4-Tetrachlorobenzene	13.27	216	126896	9.425	ng/uL	99
76) Pentachlorobenzene	14.79	250	118954	9.525	ng/uL	98
77) Carbazole	17.61	167	376316	8.371	ng/ul	99
78) Di-n-butylphthalate	18.17	149	392974	7.870	ng/ul	99
80) Fluoranthene	19.26	202	486965	9.339	ng/ul	98
82) Pyrene	19.62	202	531486	9.752	ng/ul	100
83) Butylbenzylphthalate	20.51	149	141777	7.621	ng/ul	95
84) 3,3'-Dichlorobenzidine	21.29	252	123849	7.351	ng/ul	98
85) Benzo(a)anthracene	21.35	228	485799	9.265	ng/ul	99
86) Bis(2-ethylhexyl)phthalate	21.28	149	219318	7.347	ng/ul	98
87) Chrysene	21.40	228	482843	9.460	ng/ul	99
89) Di-n-octyl phthalate	22.17	149	359337	7.281	ng/ul	100
90) Benzo(b)fluoranthene	22.96	252	494842	9.621	ng/ul	98
91) Benzo(k)fluoranthene	23.01	252	459767	9.065	ng/ul	99
93) Benzo(a)pyrene	23.55	252	425169	9.365	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	25.97	276	537006	9.444	ng/ul	100
95) Dibenzo(a,h)anthracene	25.98	278	453714	9.544	ng/ul	98
96) Benzo(g,h,i)perylene	26.68	276	448257	9.648	ng/ul	99

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Internal Standards	R.T.	Q	Ion	Response	Conc	Units	Dev	(Min)

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