

Instrument :
BNA_M
ClientSampleId :
SSTD160106

mohammad
6/23/2021 4:47:01 PM

TIC: BM030515.D

Abundance

Time-->

Pyrimidine.S
Benzaldehyde-68.1
Dichlorobenzene-d4.1
1,4-Dichlorobenzene-d4.1
Acetophenone
Naphthalene-d8.1
Caprolactam
Hexachlorocyclopentadiene
Nitrophenol-d4.1.S
Phenanthrene-d10.1
Pentachlorophenol.C
Carbazole
Chrysene-d12.1
Di-n-octyl phthalate
Perylene-d12.1
3,3'-Dithiobenzidine

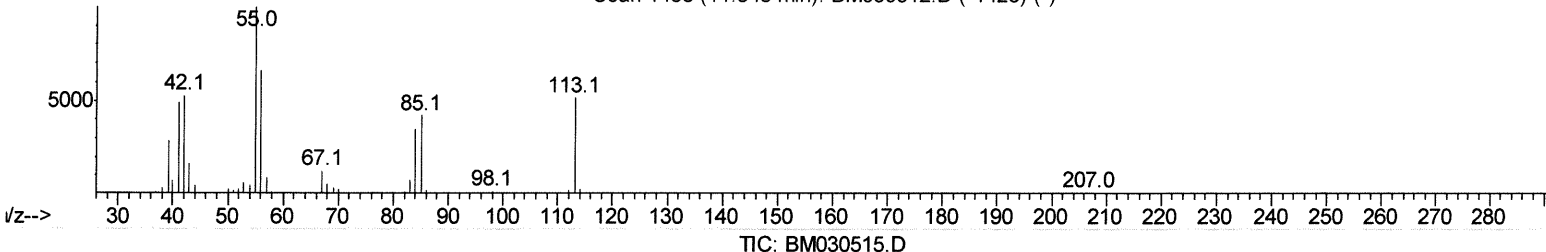
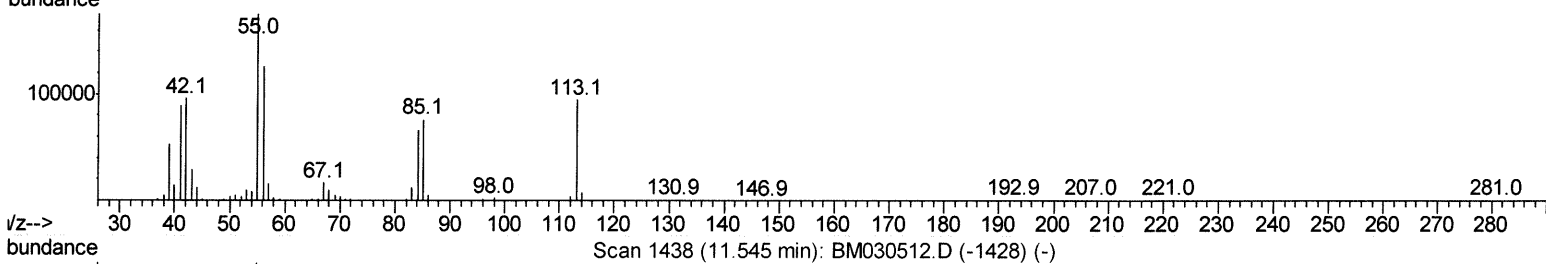
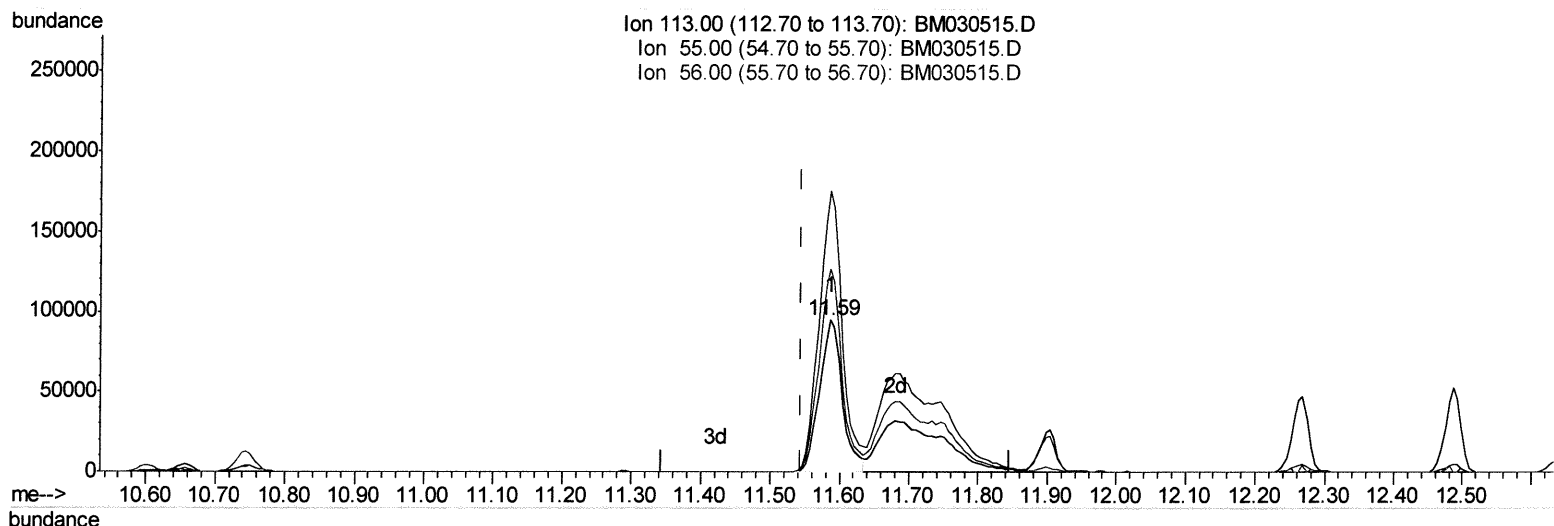
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062221\
 Data File : BM030515.D
 Acq On : 22 Jun 2021 14:29
 Operator : CG/JU
 Sample : SSTD16006
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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Manual Integrations
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Quant Time: Jun 22 15:00:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 22 13:19:30 2021
 Response via : Initial Calibration



(34) Caprolactam

11.586min (+0.041) 91.82ng/ul

response 215729

Ion	Exp%	Act%
113.00	100	100
55.00	194.00	185.08
56.00	127.40	133.58
0.00	0.00	0.00

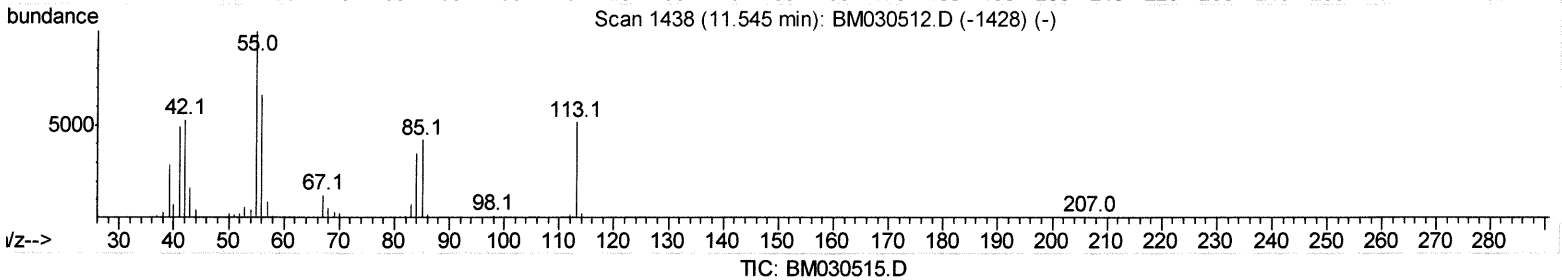
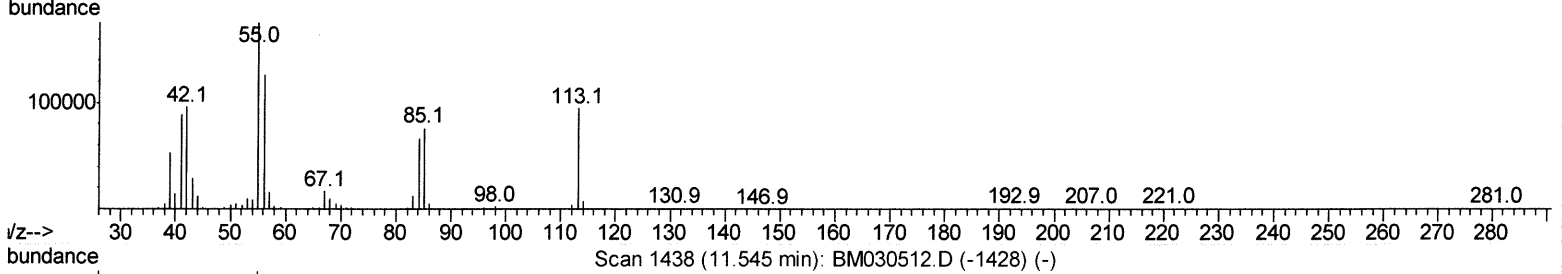
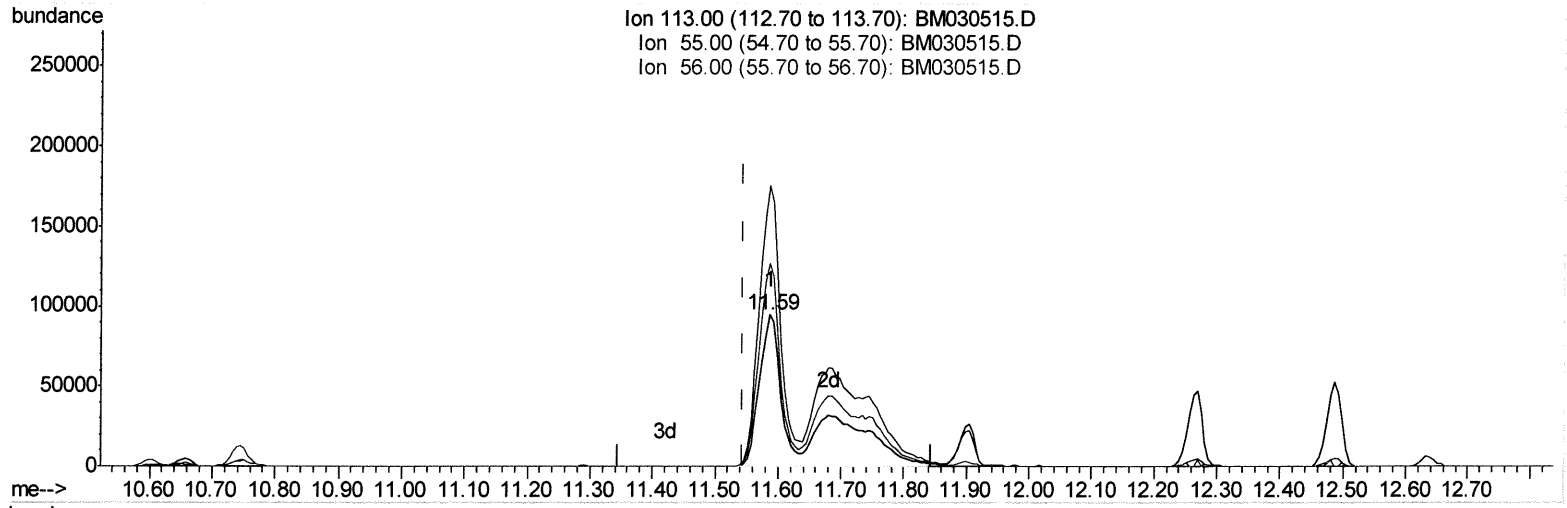
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(34) Caprolactam

11.586min (+0.041) 179.86ng/ul m

JP 6/23/21

response 422562

Ion	Exp%	Act%
113.00	100	100
55.00	194.00	185.08
56.00	127.40	133.58
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	7.81	152	127168	20.00	ng/ul	0.00
20) Naphthalene-d8	10.60	136	508812	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.44	164	313754	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.19	188	622713	20.00	ng/ul	0.00
79) Chrysene-d12	21.36	240	695245	20.00	ng/ul	0.00
88) Perylene-d12	23.63	264	725679	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
4) Pvrindine-d5	3.69	84	1526241	174.41	ng/ul	0.00
7) Phenol-d5	6.99	99	1866312	168.86	ng/ul	0.01
9) Bis-(2-Chloroethyl)ether-d	7.15	67	1110105	155.62	ng/ul	0.00
11) 2-Chlorophenol-d4	0.00	132	0d	0.00	ng/ul	
15) 4-Methylphenol-d8	8.53	113	1432356	162.27	ng/ul	0.02
21) Nitrobenzene-d5	0.00	128	0d	0.00	ng/ul	
24) 2-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
28) 2,4-Dichlorophenol-d3	0.00	165	0d	0.00	ng/ul	
31) 4-Chloroaniline-d4	10.75	131	1916307	166.36	ng/ul	0.00
46) Dimethylphthalate-d6	0.00	166	0d	0.00	ng/ul	
49) Acenaphthylene-d8	0.00	160	0d	0.00	ng/ul	
54) 4-Nitrophenol-d4	14.67	143	774457	192.16	ng/ul	0.03
60) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
65) 4,6-Dinitro-2-methylphenol	15.57	200	781339	217.37	ng/ul	0.02
73) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
81) Pyrene-d10	0.00	212	0d	0.00	ng/ul	
92) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul	

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
5) Pvrindine	3.71	79	1559649	168.821	ng/ul	98
6) Benzaldehyde	6.96	77	634234	122.441	ng/ul	96
8) Phenol	7.02	94	1881502	168.077	ng/ul	99
10) Bis(2-Chloroethyl)ether	7.25	93	1421925	157.337	ng/ul	97
13) 2-Methylphenol	8.26	108	1355021	162.781	ng/ul	97
14) 2,2'-oxybis(1-Chloropropan	8.35	45	2162416	148.548	ng/ul	99
16) Acetophenone	8.65	105	2193327	157.153	ng/ul	96
18) 4-Methylphenol	8.60	108	1460450	159.110	ng/ul	96
32) 4-Chloroaniline	10.77	127	1902964	165.050	ng/ul	100
34) Caprolactam	11.59	113	422562m	179.857	ng/ul	97
40) Hexachlorocyclopentadiene	12.62	237	1146962	198.996	ng/ul	97
51) 3-Nitroaniline	14.36	138	774281	180.170	ng/ul	94
53) 2,4-Dinitrophenol	14.57	184	578260	229.890	ng/ul	94
55) 4-Nitrophenol	14.68	109	608223	192.621	ng/ul	97
63) 4-Nitroaniline	15.53	138	717302	178.259	ng/ul	96
66) 4,6-Dinitro-2-methylphenol	15.59	198	756131	212.148	ng/ul#	93
70) Atrazine	16.66	200	1140578	173.251	ng/ul	99
71) Pentachlorophenol	16.84	266	876755	207.826	ng/ul	98
77) Carbazole	17.59	167	5304279	175.747	ng/ul	98
84) 3,3'-Dichlorobenzidine	21.27	252	2324859	184.909	ng/ul	97
89) Di-n-octyl phthalate	22.15	149	7111292	152.145	ng/ul	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed