

Data File : BN005750.D

Operator : JU/SJ

Misc :

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M

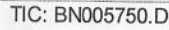
QLast Update : Sat May 18 12:10:22 2019

BNA_N

SSTD01079

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5/21/2019 7:10:03 PM



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051819\

Data File : BN005750.D

Acq On : 18 May 2019 12:59

Operator : JU/SJ

Sample : SSTD01079

Misc :

ALS Vial : 3 Sample Multiplier: 1

Quant Time: May 18 13:34:05 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Sat May 18 12:10:22 2019

Response via : Initial Calibration

Instrument :

BNA_N

Client Sampled :

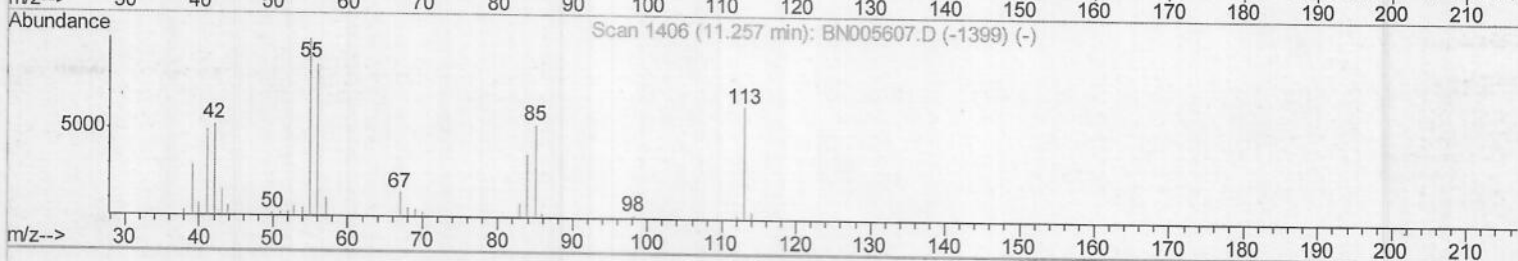
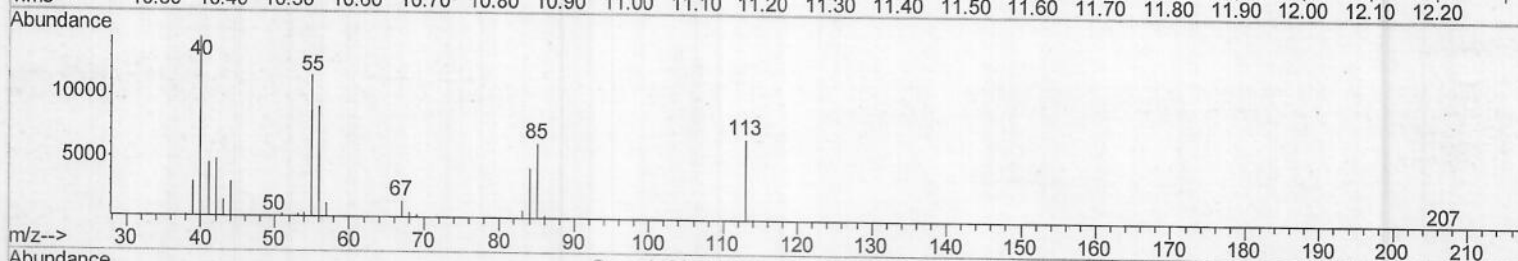
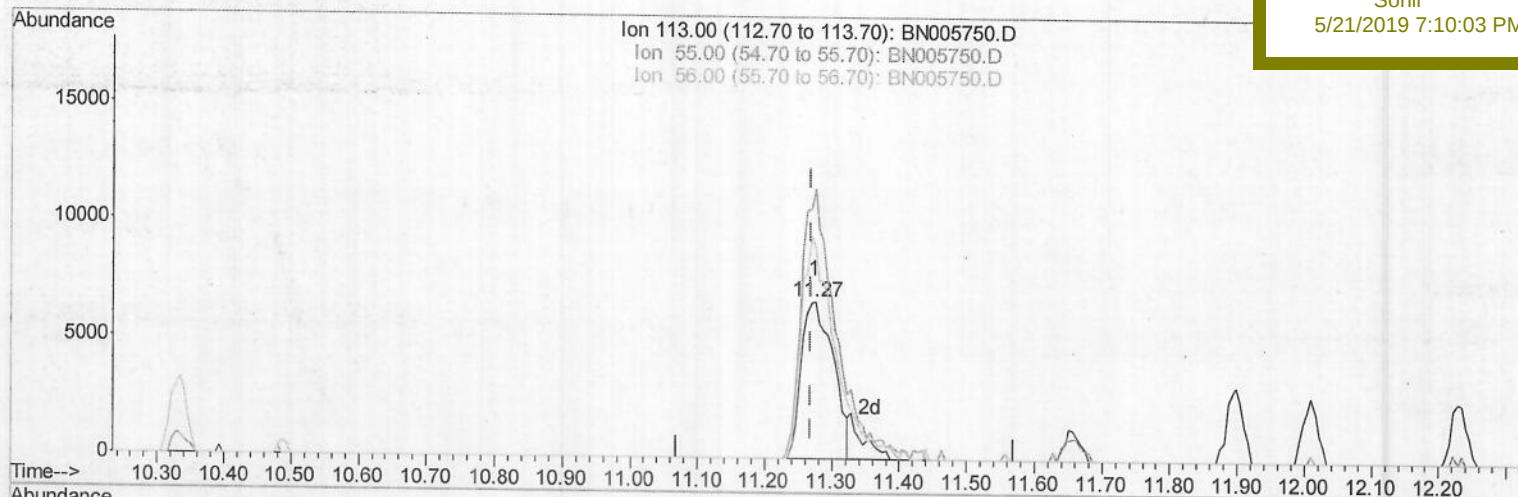
SSTD01079

Manual Integrations

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Sohil

5/21/2019 7:10:03 PM



(32) Caprolactam

11.275min (+0.006) 11.19ng/ul

response 22047

Ion	Exp%	Act%
113.00	100	100
55.00	196.20	171.83
56.00	148.90	134.35
0.00	0.00	0.00

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Misc :

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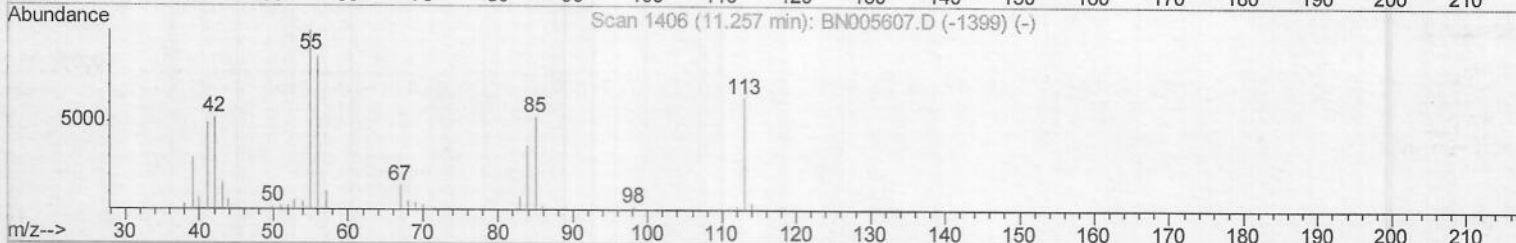
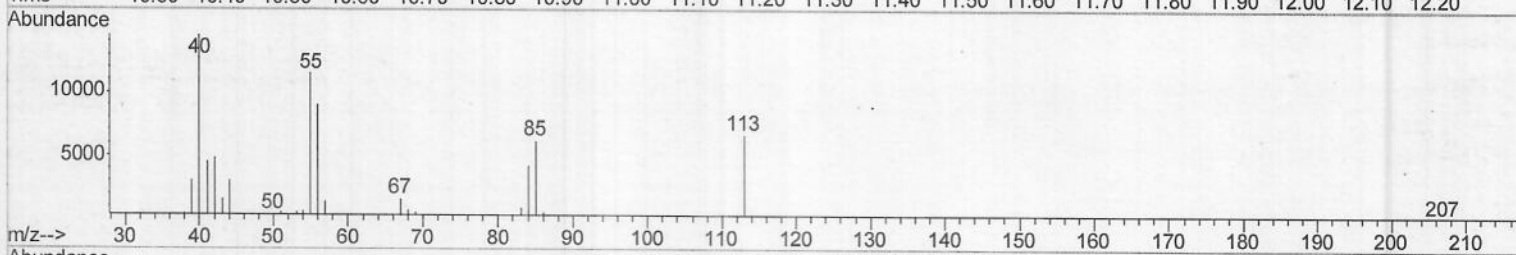
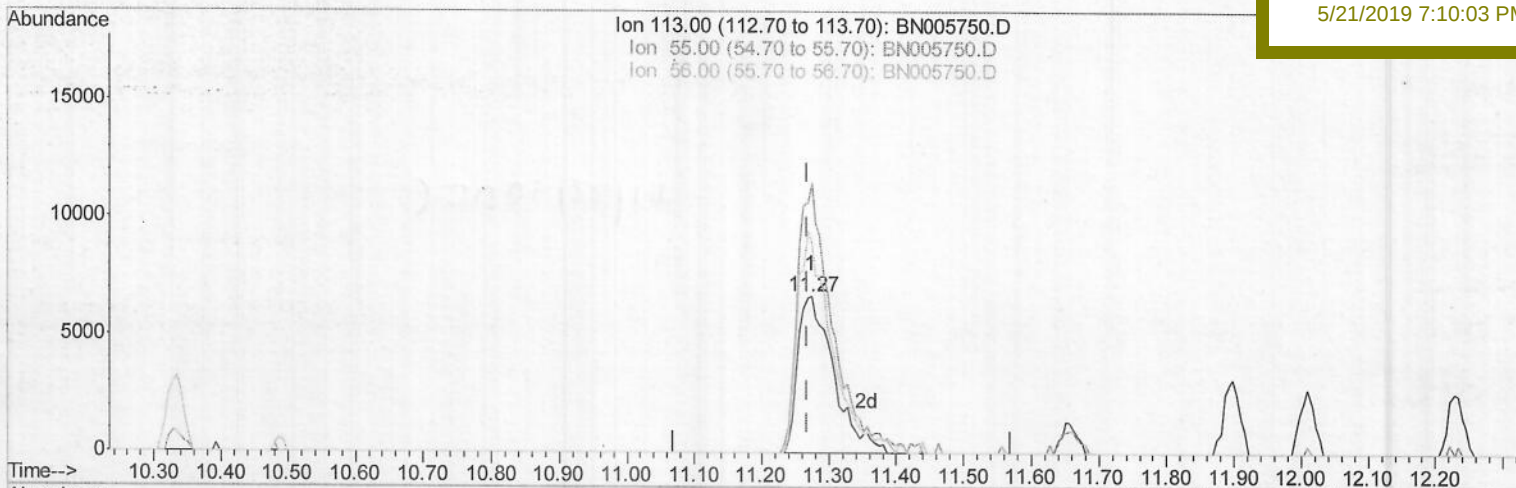
SSTD01079

Manual Integrations

APPROVED

Sohil

5/21/2019 7:10:03 PM



TIC: BN005750.D

(32) Caprolactam

11.275min (+0.006) 12.66ng/ul m > 50.05/22/19

response 24929

Ion	Exp%	Act%
113.00	100	100
55.00	196.20	171.83
56.00	148.90	134.35
0.00	0.00	0.00

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Acq On : 18 May 2019 12:59

Operator : JU/SJ

Sample : SST01079

Misc :

ALS Vial : 3 Sample Multiplier: 1

Quant Time: May 18 13:35:17 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Sat May 18 12:10:22 2019

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Client Sampled :

SSTD01079

Manual Integrations

APPROVED

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5/21/2019 7:10:03 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	102732	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	461200	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.21	164	327982	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.97	188	815885	20.00	ng/ul	0.00
77) Chrysene-d12	21.22	240	862926	20.00	ng/ul	0.00
85) Perylene-d12	23.44	264	973940	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	7400	3.48	ng/uL	0.00
5) Phenol-d5	6.76	99	71684	8.88	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.92	67	39769	7.90	ng/ul	0.00
9) 2-Chlorophenol-d4	7.11	132	60670	9.83	ng/ul	0.00
13) 4-Methylphenol-d8	8.28	113	62134	9.85	ng/ul	0.00
19) Nitrobenzene-d5	8.72	128	32121	11.32	ng/ul	0.00
22) 2-Nitrophenol-d4	9.44	143	36983	12.85	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.98	165	67649	10.26	ng/ul	0.00
29) 4-Chloroaniline-d4	10.49	131	70950	10.43	ng/ul	0.00
43) Dimethylphthalate-d6	13.63	166	245532	10.31	ng/ul	0.00
46) Acenaphthylene-d8	13.90	160	297960	10.19	ng/ul	0.00
51) 4-Nitrophenol-d4	14.46	143	25498	6.78	ng/ul	0.00
57) Fluorene-d10	15.22	176	210135	10.52	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.39	200	37736	10.28	ng/ul	0.00
70) Anthracene-d10	17.07	188	346069	10.12	ng/ul	0.00
78) Pyrene-d10	19.39	212	410714	9.30	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.30	264	430297	10.11	ng/ul	0.01

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.22	88	7877	3.442	ng/uL 97
4) Benzaldehyde	6.73	77	52622	9.203	ng/ul 96
6) Phenol	6.79	94	74005	8.917	ng/ul 96
8) Bis(2-Chloroethyl) ether	7.01	93	55730	8.386	ng/ul 92
10) 2-Chlorophenol	7.14	128	60656	9.502	ng/ul 95
11) 2-Methylphenol	8.02	108	59051	9.495	ng/ul 93
12) 2,2'-oxybis(1-Chloropropan	8.09	45	76208	6.830	ng/ul# 91
14) Acetophenone	8.39	105	103392	10.243	ng/ul 98
15) N-Nitroso-di-n-propylamine	8.37	70	50797	9.408	ng/ul 93
16) 4-Methylphenol	8.35	108	64099	9.544	ng/ul 93
17) Hexachloroethane	8.62	117	26553	9.579	ng/ul 93
20) Nitrobenzene	8.76	77	75211	9.628	ng/ul 97
21) Isophorone	9.27	82	153521	9.925	ng/ul 96
23) 2-Nitrophenol	9.47	139	38021	11.659	ng/ul 98
24) 2,4-Dimethylphenol	9.53	107	78697	9.638	ng/ul 99
25) Bis(2-Chloroethoxy)methane	9.76	93	85825	8.829	ng/ul 97
27) 2,4-Dichlorophenol	10.01	162	66035	10.143	ng/ul 96
28) Naphthalene	10.38	128	214895	10.023	ng/ul 99
30) 4-Chloroaniline	10.52	127	72117	10.408	ng/ul 100
31) Hexachlorobutadiene	10.66	225	51494	11.201	ng/ul 98
32) Caprolactam	11.27	113	24929m	12.656	ng/ul 98
33) 4-Chloro-3-methylphenol	11.65	107	76286	10.407	ng/ul 92
34) 2-Methylnaphthalene	12.01	142	166534	10.608	ng/ul 97

Sulosteezlg

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 1,2,4,5-Tetrachlorobenzene	12.39	216	102028	10.052	ng/ul	96
37) Hexachlorocyclopentadiene	12.36	237	21551	3.114	ng/ul	95
38) 2,4,6-Trichlorophenol	12.65	196	61329	10.020	ng/ul	96
39) 2,4,5-Trichlorophenol	12.73	196	61223	9.309	ng/ul	96
40) 1,1'-Biphenyl	13.04	154	230110	9.416	ng/ul	98
41) 2-Chloronaphthalene	13.08	162	181787	9.672	ng/ul	97
42) 2-Nitroaniline	13.30	65	52150	10.461	ng/ul	96
44) Dimethylphthalate	13.67	163	242685	10.230	ng/ul	99
45) 2,6-Dinitrotoluene	13.81	165	49460	12.074	ng/ul	94
47) Acenaphthylene	13.93	152	284543	9.901	ng/ul	98
48) 3-Nitroaniline	14.15	138	40850	10.727	ng/ul	96
49) Acenaphthene	14.28	153	194579	9.918	ng/ul	98
50) 2,4-Dinitrophenol	14.40	184	12418	6.223	ng/ul	94
52) 4-Nitrophenol	14.48	109	24977	7.320	ng/ul	94
53) Dibenzofuran	14.62	168	292253	10.346	ng/ul	96
54) 2,4-Dinitrotoluene	14.62	165	75346	12.759	ng/ul#	84
55) 2,3,4,6-Tetrachlorophenol	14.86	232	59180	10.816	ng/ul	91
56) Diethylphthalate	15.05	149	243618	10.241	ng/ul	97
58) Fluorene	15.27	166	235591	10.689	ng/ul	99
59) 4-Chlorophenyl-phenylether	15.27	204	130186	11.134	ng/ul	93
60) 4-Nitroaniline	15.32	138	47580	10.022	ng/ul	98
63) 4,6-Dinitro-2-methylphenol	15.40	198	38436	9.841	ng/ul#	97
64) N-Nitrosodiphenylamine	15.49	169	207658	9.299	ng/ul	98
65) 4-Bromophenyl-phenylether	16.16	248	82345	10.018	ng/ul	96
66) Hexachlorobenzene	16.29	284	93589	10.693	ng/ul	95
67) Atrazine	16.45	200	86676	10.426	ng/ul	99
68) Pentachlorophenol	16.66	266	32670	6.375	ng/ul	93
69) Phenanthrene	17.02	178	398385	10.026	ng/ul	99
71) Anthracene	17.11	178	409624	10.120	ng/ul	99
72) 1,2,3,4-Tetrachlorobenzene	13.00	216	107365	9.009	ng/ul	97
73) Pentachlorobenzene	14.55	250	105082	9.797	ng/ul	99
74) Carbazole	17.39	167	359563	10.341	ng/ul	98
75) Di-n-butylphthalate	17.96	149	424985	9.741	ng/ul	98
76) Fluoranthene	19.06	202	502383	11.399	ng/ul#	92
79) Pyrene	19.43	202	514069	9.252	ng/ul#	90
80) Butylbenzylphthalate	20.34	149	208002	9.456	ng/ul	91
81) 3,3'-Dichlorobenzidine	21.14	252	175466	10.177	ng/ul#	98
82) Benzo(a)anthracene	21.20	228	531351	9.898	ng/ul	99
83) Bis(2-ethylhexyl)phthalate	21.14	149	297607	8.786	ng/ul#	97
84) Chrysene	21.26	228	504948	10.135	ng/ul	99
86) Di-n-octyl phthalate	22.02	149	518720	8.696	ng/ul	100
87) Benzo(b)fluoranthene	22.78	252	511919	9.350	ng/ul#	95
88) Benzo(k)fluoranthene	22.82	252	521268	9.969	ng/ul#	98
90) Benzo(a)pyrene	23.34	252	510752	9.925	ng/ul#	96
91) Indeno(1,2,3-cd)pyrene	25.64	276	623844	10.965	ng/ul#	93
92) Dibenzo(a,h)anthracene	25.64	278	518897	10.747	ng/ul#	94
93) Benzo(a,h,i)perylene	26.32	276	517896	11.009	ng/ul#	91

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