

(QT Reviewed)

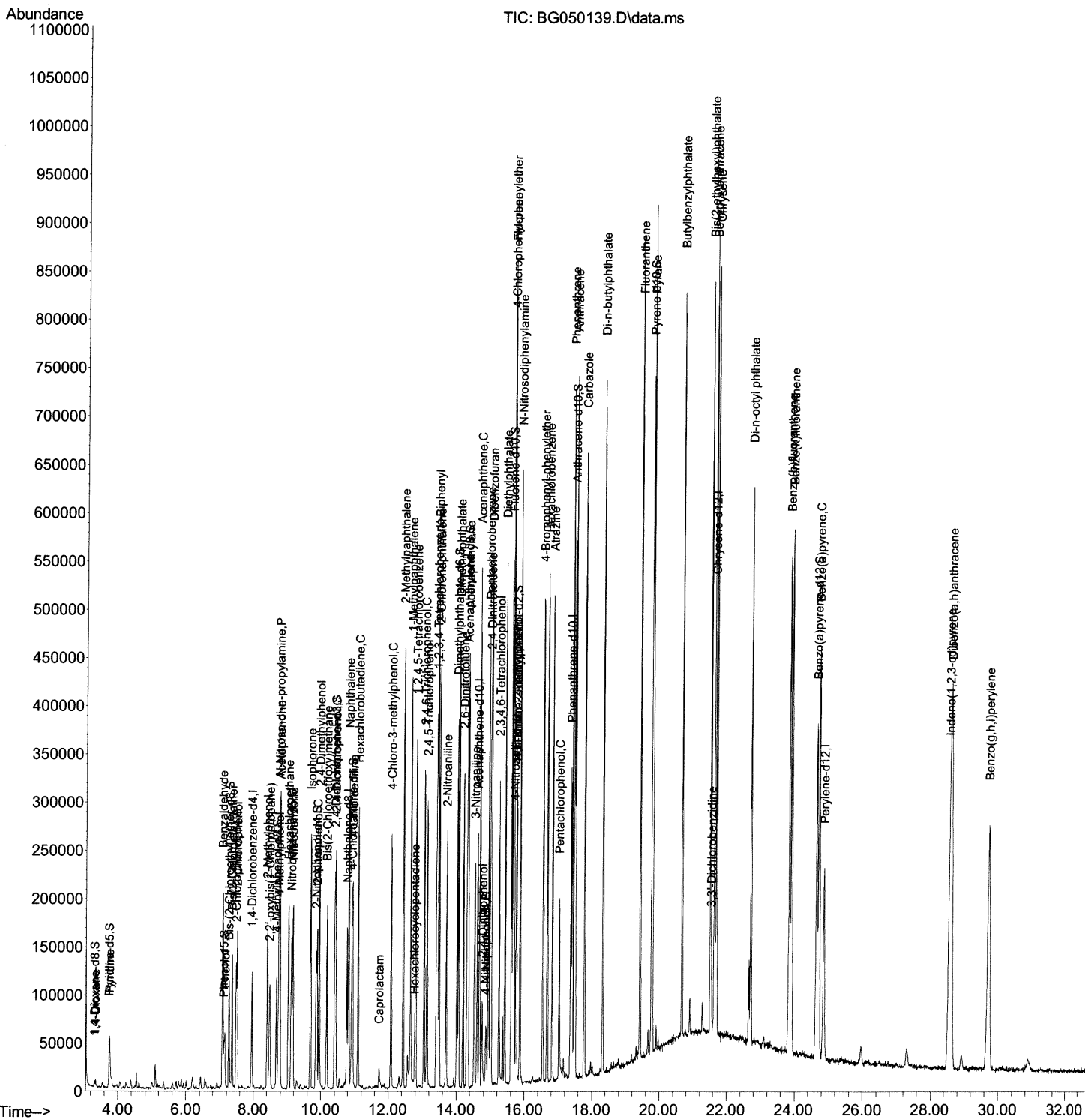
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Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\  
Data File : BG050139.D  
Acq On    : 3 Sep 2021 20:19  
Operator  : CG/JU  
Sample    : M3549-03MSD  
Misc      :  
ALS Vial  : 41 Sample Multiplier: 1
```

Instrument :
BNA_G
ClientSampleId :
EW7A1MSD

Manual Integrations

mohammad
9/7/2021 11:34:01 AM

Quant Time: Sep 03 23:47:30 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 30 14:59:55 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

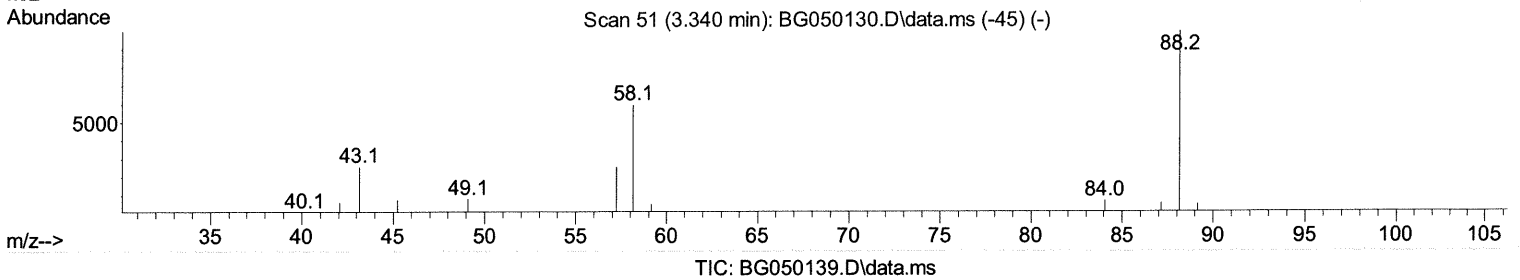
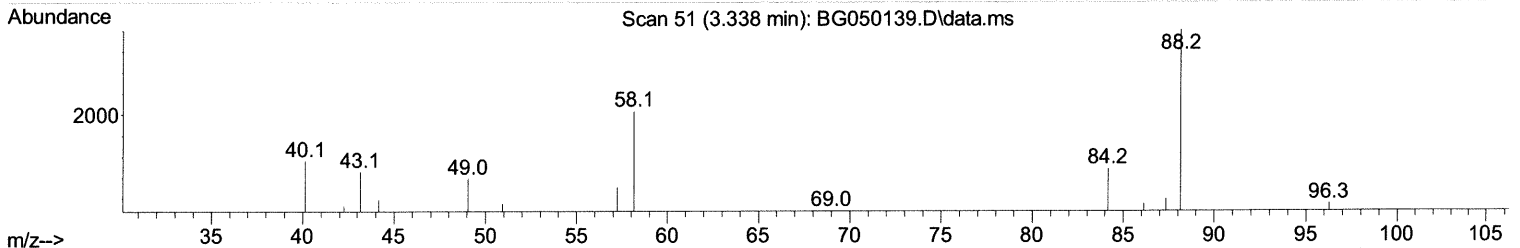
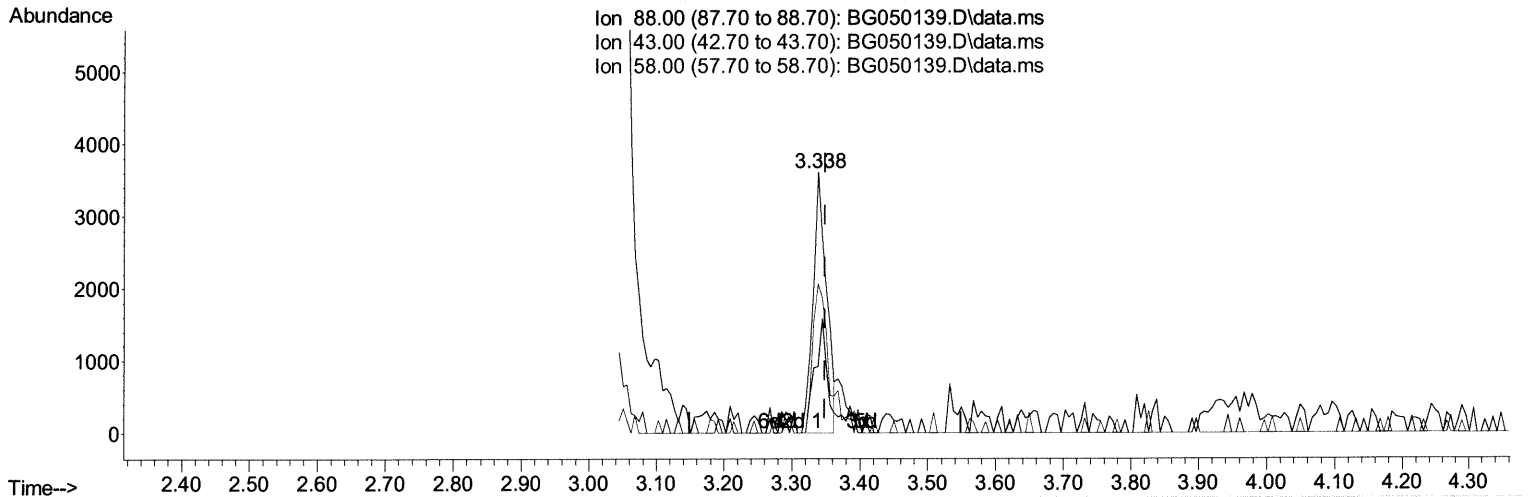
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(2) 1,4-Dioxane

3.338min (-0.011) 5.95 ng/uL

response 4936

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	27.00	25.40
58.00	65.00	57.27
0.00	0.00	0.00

Quantitation Report (Qedit)

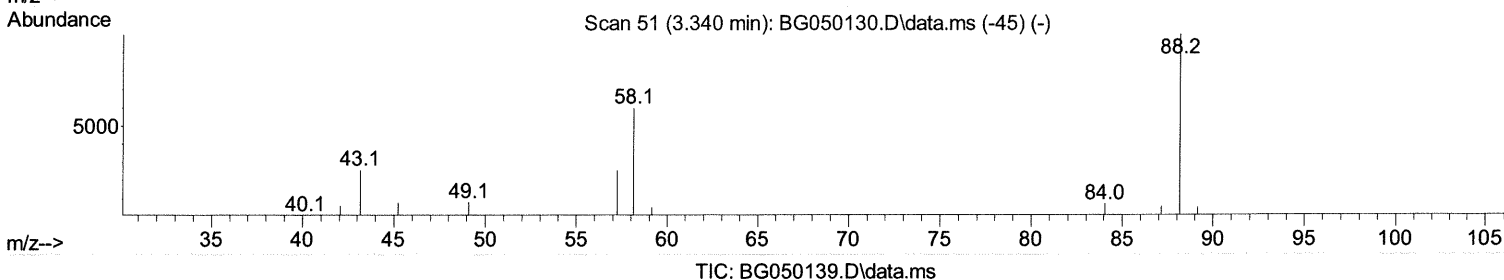
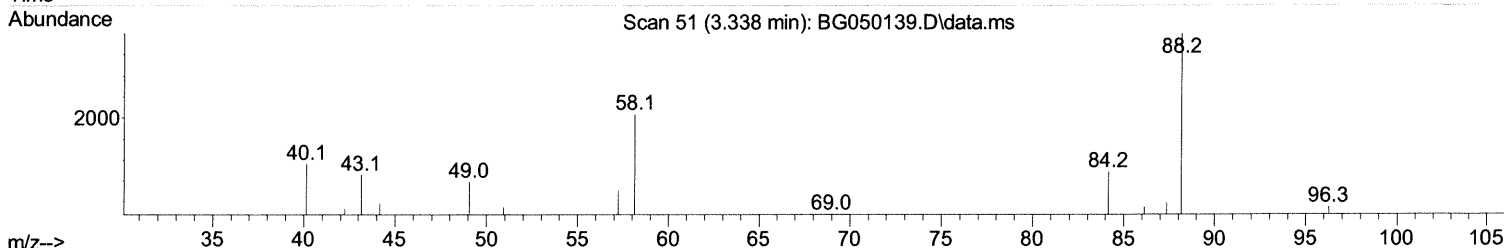
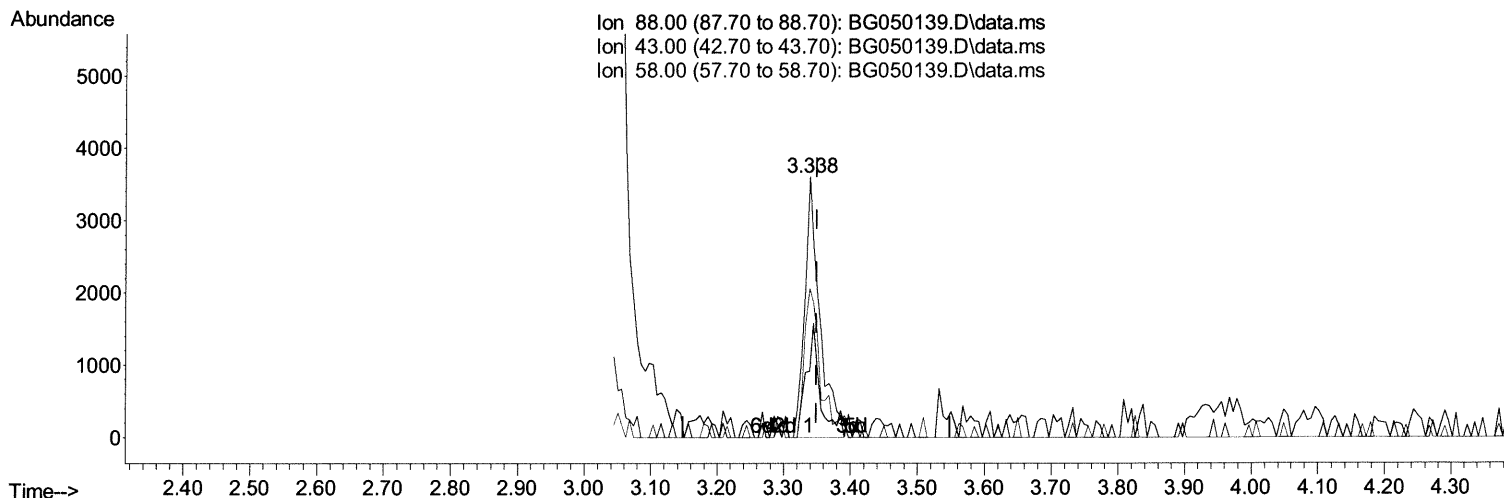
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(2) 1,4-Dioxane

3.338min (-0.011) 6.81 ng/uL m

response 5653

RP 9/8/21

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	27.00	25.40
58.00	65.00	57.27
0.00	0.00	0.00

Quantitation Report (Qedit)

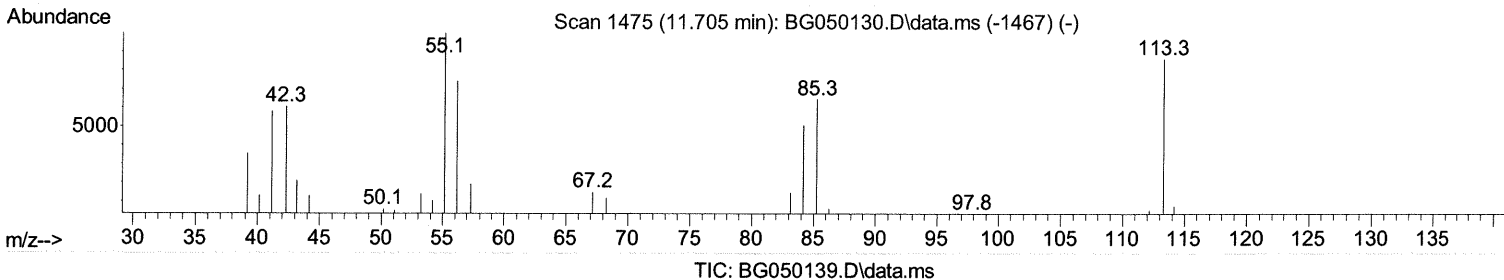
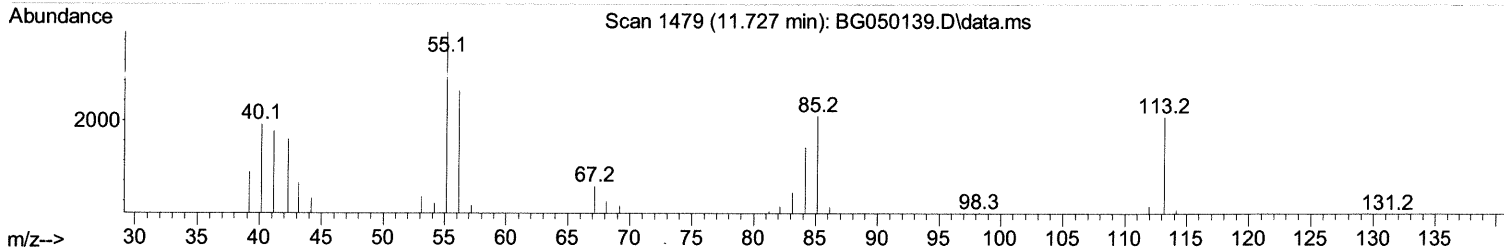
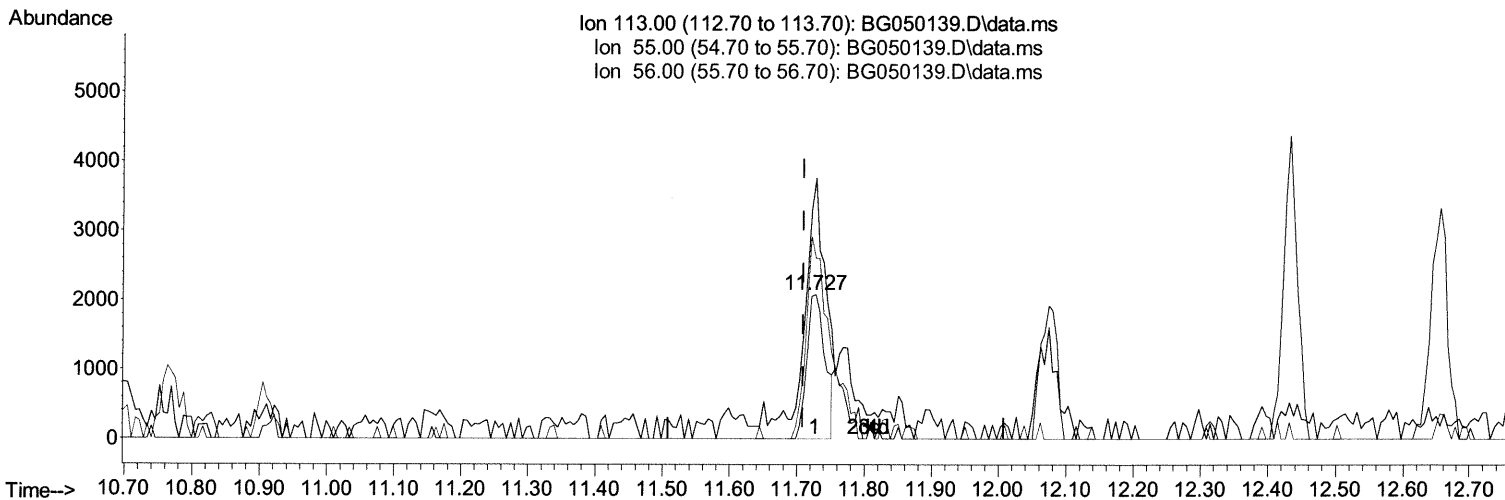
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 Operator : CG/JU
 Sample : M3549-03MSD
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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(34) Caprolactam

11.727min (+ 0.019) 5.47 ng/ul

response 3950

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	142.50	180.68#
56.00	134.20	125.04
0.00	0.00	0.00

Quantitation Report (Qedit)

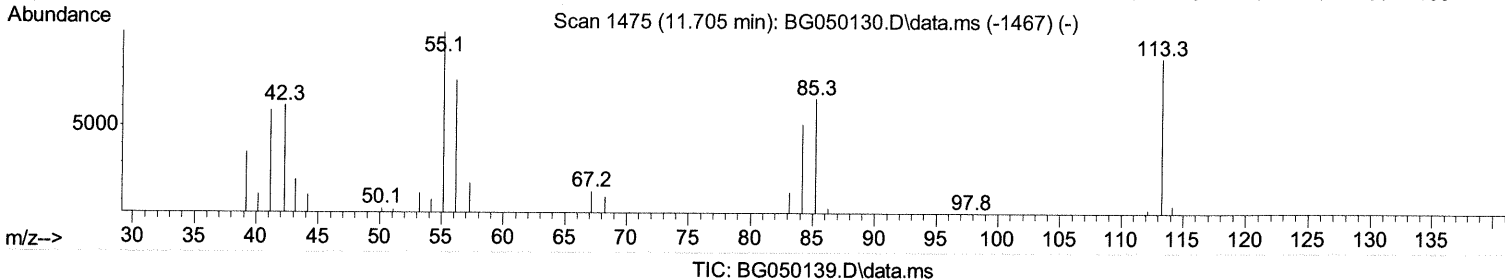
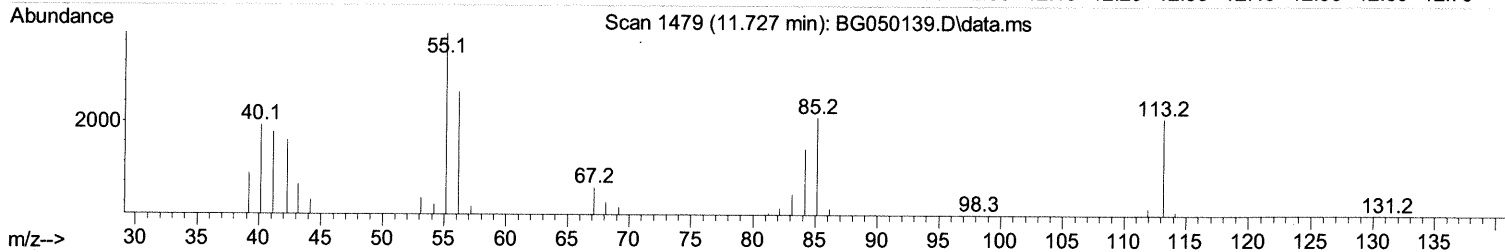
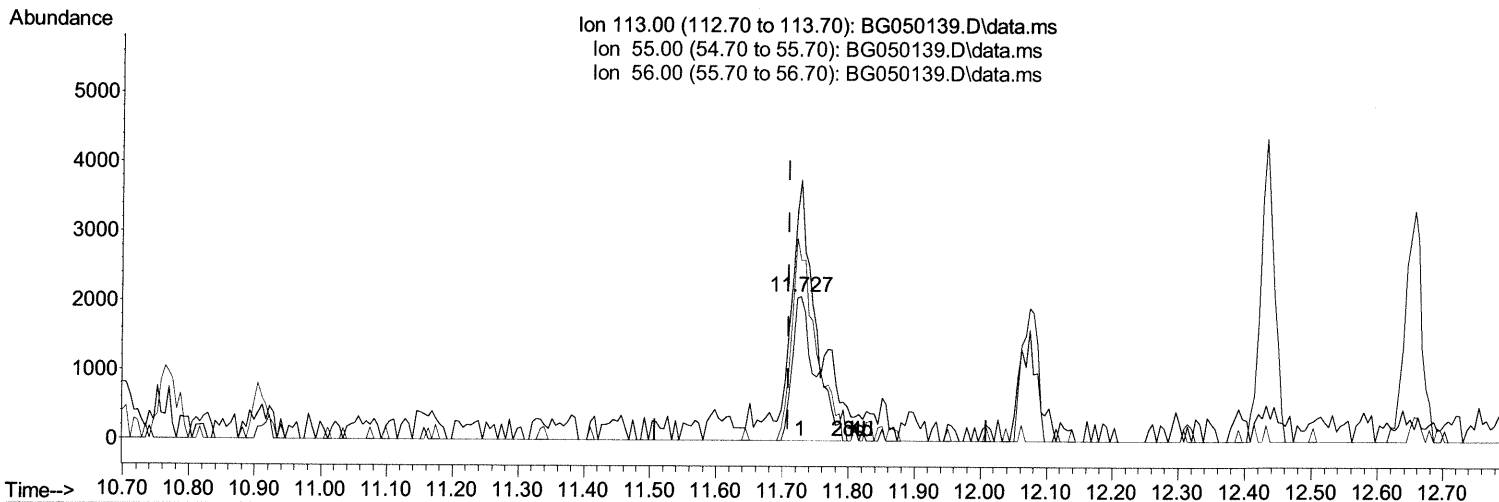
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 Data File : BG050139.D
 Acq On : 3 Sep 2021 20:19
 Operator : CG/JU
 Sample : M3549-03MSD
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW7A1MSD

Manual Integrations
 APPROVED

mohammad
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Quant Time: Sep 03 23:47:30 2021
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 Response via : Initial Calibration



(34) Caprolactam

11.727min (+ 0.019) 7.37 ng/ul m 34 9/8/21

response 5322

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	142.50	180.68#
56.00	134.20	125.04
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
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 Operator : CG/JU
 Sample : M3549-03MSD
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampled :
 EW7A1MSD

Manual Integrations
 APPROVED

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 9/7/2021 11:34:01 AM

Quant Time: Sep 03 23:47:30 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 30 14:59:55 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.950	152	35088	20.000 ng/ul	0.00
20) Naphthalene-d8	10.769	136	136363	20.000 ng/ul	-0.01
38) Acenaphthene-d10	14.611	164	90821	20.000 ng/ul	-0.01
64) Phenanthrene-d10	17.366	188	201972	20.000 ng/ul	#-0.01
79) Chrysene-d12	21.643	240	198270	20.000 ng/ul	0.00
88) Perylene-d12	24.862	264	201018	20.000 ng/ul	0.00

System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.309	96	3741	5.498 ng/uL	0.00
4) Pyridine-d5	3.738	84	24377	11.894 ng/ul	0.00
7) Phenol-d5	7.121	99	26031	8.737 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.268	67	54995	33.223 ng/ul	-0.01
11) 2-Chlorophenol-d4	7.480	132	59456	26.669 ng/ul	-0.01
15) 4-Methylphenol-d8	8.678	113	42867	18.268 ng/ul	0.00
21) Nitrobenzene-d5	9.124	128	37398	35.375 ng/ul	0.00
24) 2-Nitrophenol-d4	9.847	143	41902	33.286 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.405	165	75137	33.464 ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	86711	28.373 ng/ul	-0.01
46) Dimethylphthalate-d6	14.018	166	258787	37.232 ng/ul	0.00
49) Acenaphthylene-d8	14.306	160	298770	36.660 ng/ul	-0.01
54) 4-Nitrophenol-d4	14.852	143	9269	9.217 ng/ul	0.00
60) Fluorene-d10	15.610	176	215835	36.625 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	44423	34.188 ng/ul	0.00
73) Anthracene-d10	17.466	188	337411	37.203 ng/ul	-0.01
81) Pyrene-d10	19.757	212	403731	37.681 ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.645	264	397365	37.260 ng/ul	0.00

Target Compounds					
2) 1,4-Dioxane	3.338	88	5653m	6.815 ng/uL	Qvalue 74 9/8/21
5) Pyridine	3.755	79	30545	13.789 ng/ul	85
6) Benzaldehyde	7.080	77	71435	41.663 ng/ul	99
8) Phenol	7.151	94	30399	10.101 ng/ul	98
10) Bis(2-Chloroethyl)ether	7.362	93	77332	37.223 ng/ul	94
12) 2-Chlorophenol	7.515	128	68837	31.247 ng/ul#	91
13) 2-Methylphenol	8.408	108	58140	25.041 ng/ul	100
14) 2,2'-oxybis(1-Chloropr...	8.472	45	81717	37.510 ng/ul	96
16) Acetophenone	8.784	105	140653	36.755 ng/ul#	89
17) N-Nitroso-di-n-propyla...	8.760	70	71509	37.215 ng/ul	97
18) 4-Methylphenol	8.737	108	53999	21.618 ng/ul	86
19) Hexachloroethane	9.030	117	38080	39.497 ng/ul	94
22) Nitrobenzene	9.166	77	113283	39.546 ng/ul	99
23) Isophorone	9.688	82	200992	39.380 ng/ul	97
25) 2-Nitrophenol	9.882	139	47136	38.514 ng/ul	94
26) 2,4-Dimethylphenol	9.947	107	95890	35.254 ng/ul	92
27) Bis(2-Chloroethoxy)met...	10.170	93	107093	40.633 ng/ul	96
29) 2,4-Dichlorophenol	10.429	162	83475	40.506 ng/ul	100
30) Naphthalene	10.822	128	270075	39.625 ng/ul	98
32) 4-Chloroaniline	10.934	127	93647	31.818 ng/ul	91
33) Hexachlorobutadiene	11.098	225	64972	38.973 ng/ul	91
34) Caprolactam	11.727	113	5322m	7.372 ng/ul	74 9/8/21
35) 4-Chloro-3-methylphenol	12.073	107	93343	37.899 ng/ul	97

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Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
36) 2-Methylnaphthalene	12.432	142	200849	39.642 ng/ul	97
37) 1-Methylnaphthalene	12.655	142	196341	38.392 ng/ul	95
39) 1,2,4,5-Tetrachloroben...	12.802	216	114337	42.862 ng/ul#	92
40) Hexachlorocyclopentadiene	12.772	237	12439	10.419 ng/ul#	95
41) 2,4,6-Trichlorophenol	13.049	196	74298	43.706 ng/ul	97
42) 2,4,5-Trichlorophenol	13.131	196	80216	45.023 ng/ul	97
43) 1,1'-Biphenyl	13.442	154	263664	41.554 ng/ul	97
44) 2-Chloronaphthalene	13.489	162	206969	40.609 ng/ul	96
45) 2-Nitroaniline	13.695	65	75838	43.827 ng/ul	94
47) Dimethylphthalate	14.065	163	288819	41.987 ng/ul	99
48) 2,6-Dinitrotoluene	14.194	165	62638	44.106 ng/ul	93
50) Acenaphthylene	14.335	152	314300	42.196 ng/ul	96
51) 3-Nitroaniline	14.529	138	56454	41.195 ng/ul	97
52) Acenaphthene	14.676	153	231376	42.822 ng/ul	98
53) 2,4-Dinitrophenol	14.752	184	20421	28.153 ng/ul	91
55) 4-Nitrophenol	14.864	109	12229	9.591 ng/ul	96
56) Dibenzofuran	15.011	168	318967	42.628 ng/ul	100
57) 2,4-Dinitrotoluene	14.993	165	93051	45.550 ng/ul#	93
58) 2,3,4,6-Tetrachlorophenol	15.251	232	68892	46.662 ng/ul	95
59) Diethylphthalate	15.428	149	302898	43.044 ng/ul	96
61) Fluorene	15.663	166	273107	43.262 ng/ul	97
62) 4-Chlorophenyl-phenyle...	15.651	204	141157	42.261 ng/ul	96
63) 4-Nitroaniline	15.692	138	60428	44.186 ng/ul	97
66) 4,6-Dinitro-2-methylph...	15.763	198	46078	39.123 ng/ul	95
67) N-Nitrosodiphenylamine	15.868	169	240923	45.308 ng/ul	96
68) 4-Bromophenyl-phenylether	16.550	248	102584	45.914 ng/ul	99
69) Hexachlorobenzene	16.679	284	114570	44.407 ng/ul	93
70) Atrazine	16.826	200	105695	44.742 ng/ul	95
71) Pentachlorophenol	17.026	266	40879	37.418 ng/ul	96
72) Phenanthrene	17.413	178	450906	44.782 ng/ul	97
74) Anthracene	17.501	178	446154	43.726 ng/ul	94
75) 1,2,3,4-Tetrachloroben...	13.413	216	117046	43.194 ng/ul	92
76) Pentachlorobenzene	14.940	250	125280	43.632 ng/ul	94
77) Carbazole	17.777	167	414154	45.615 ng/ul	99
78) Di-n-butylphthalate	18.318	149	507630	43.716 ng/ul	99
80) Fluoranthene	19.428	202	547111	41.380 ng/ul#	96
82) Pyrene	19.787	202	542059	41.374 ng/ul	98
83) Butylbenzylphthalate	20.662	149	219053	42.369 ng/ul	99
84) 3,3'-Dichlorobenzidine	21.537	252	31338	6.668 ng/ul#	91
85) Benzo(a)anthracene	21.619	228	537315	43.492 ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.508	149	312565	44.021 ng/ul	100
87) Chrysene	21.690	228	509397	43.619 ng/ul	95
89) Di-n-octyl phthalate	22.712	149	517043	43.358 ng/ul	100
90) Benzo(b)fluoranthene	23.840	252	553000	43.007 ng/ul#	96
91) Benzo(k)fluoranthene	23.904	252	536380	43.885 ng/ul	99
93) Benzo(a)pyrene	24.709	252	485749	43.499 ng/ul	97
94) Indeno(1,2,3-cd)pyrene	28.557	276	654198	44.411 ng/ul#	98
95) Dibenzo(a,h)anthracene	28.610	278	541761	43.933 ng/ul#	99
96) Benzo(g,h,i)perylene	29.714	276	544328	43.954 ng/ul	98

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