

(QT Reviewed)

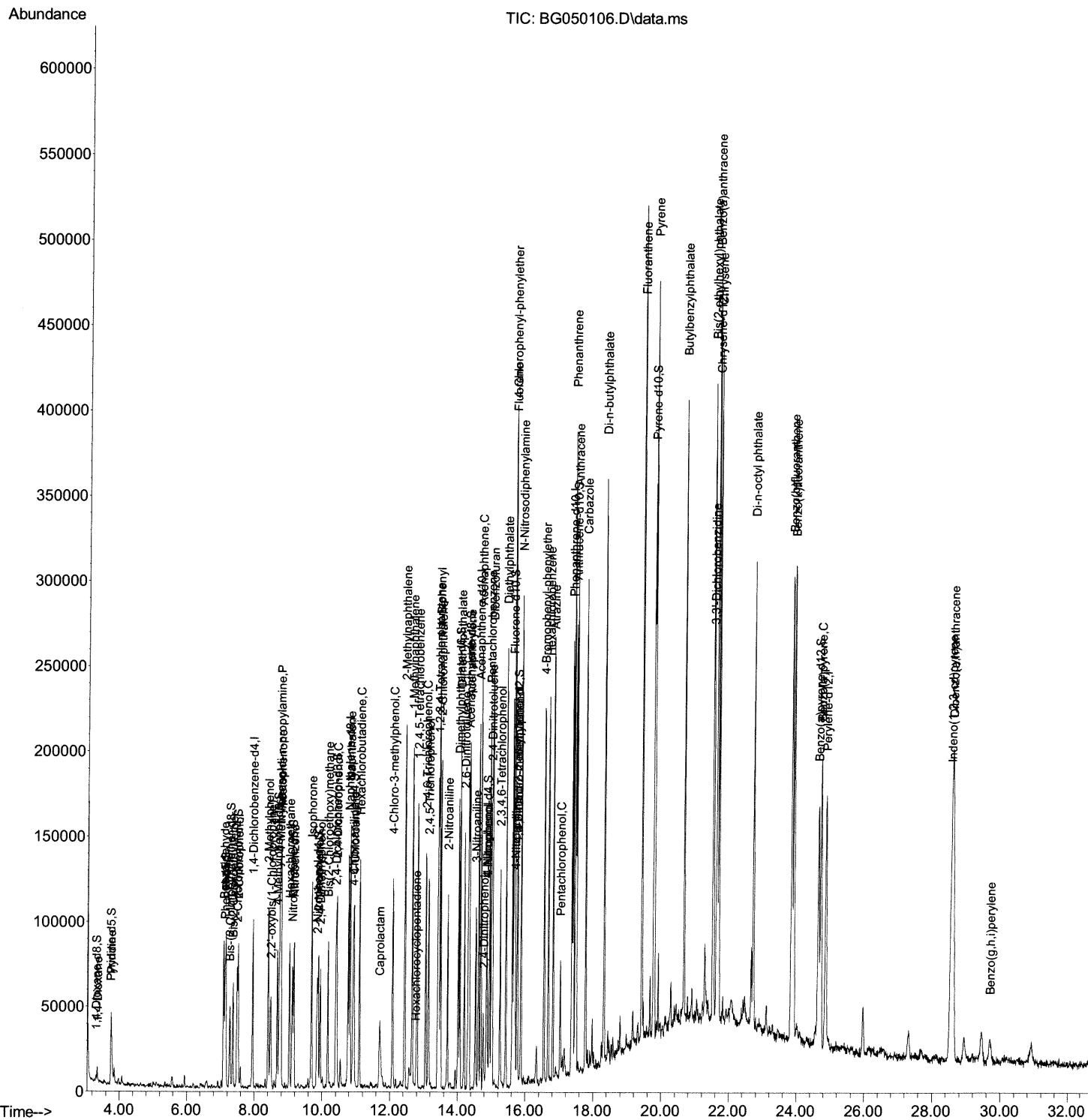
```
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\  
Data File : BG050106.D  
Acq On    : 2 Sep 2021 21:12  
Operator  : CG/JU  
Sample    : M3526-03MSD  
Misc      :  
ALS Vial  : 8 Sample Multiplier: 1
```

Instrument :
BNA_G
ClientSampleId :
EW7A9MSD

Manual Integrations APPROVED

mohammad
9/3/2021 3:31:54 PM

Quant Time: Sep 03 00:58:47 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)

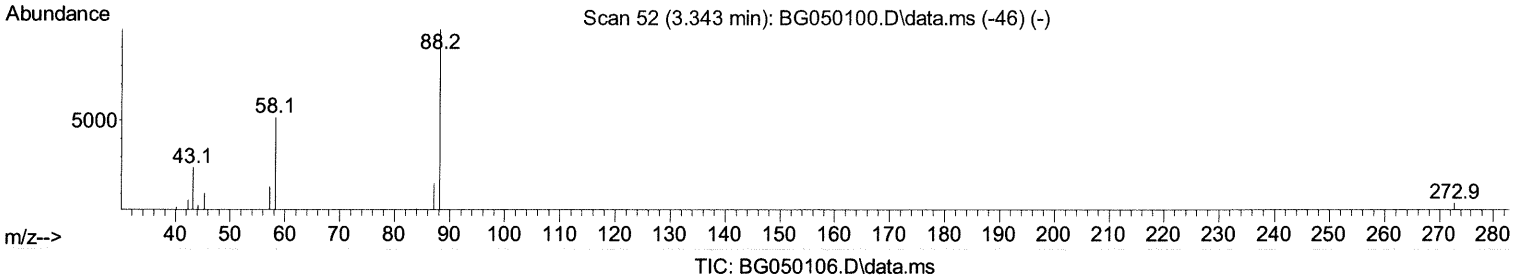
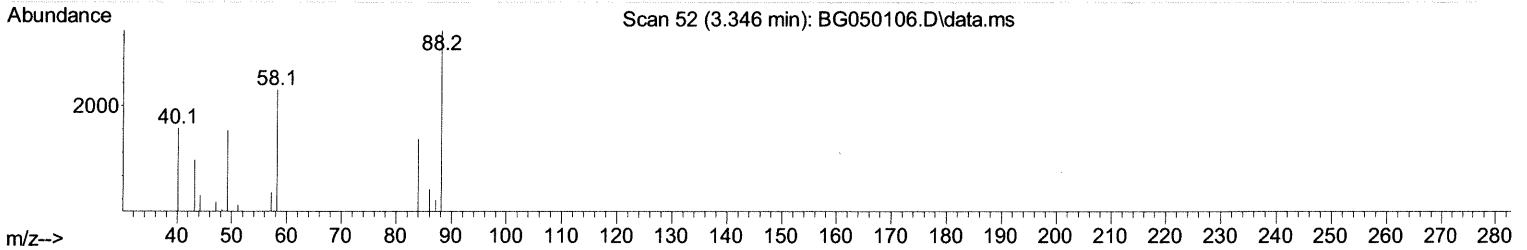
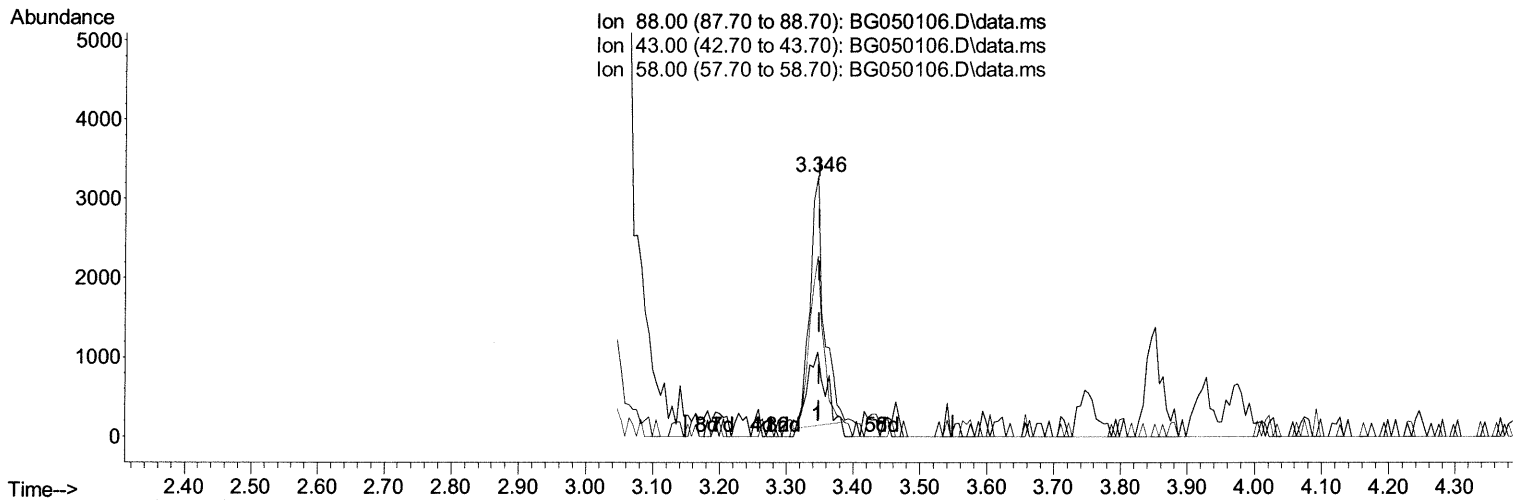
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
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(2) 1,4-Dioxane

3.346min (-0.003) 7.19 ng/uL

response 4700

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	27.00	32.52#
58.00	65.00	69.09
0.00	0.00	0.00

Quantitation Report (Qedit)

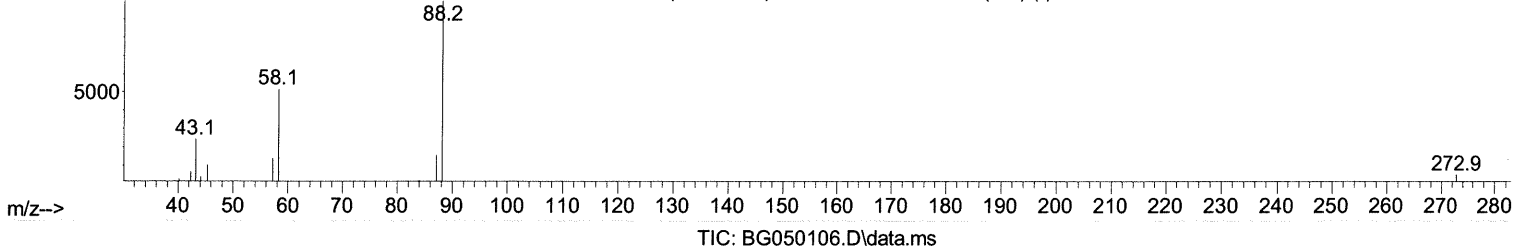
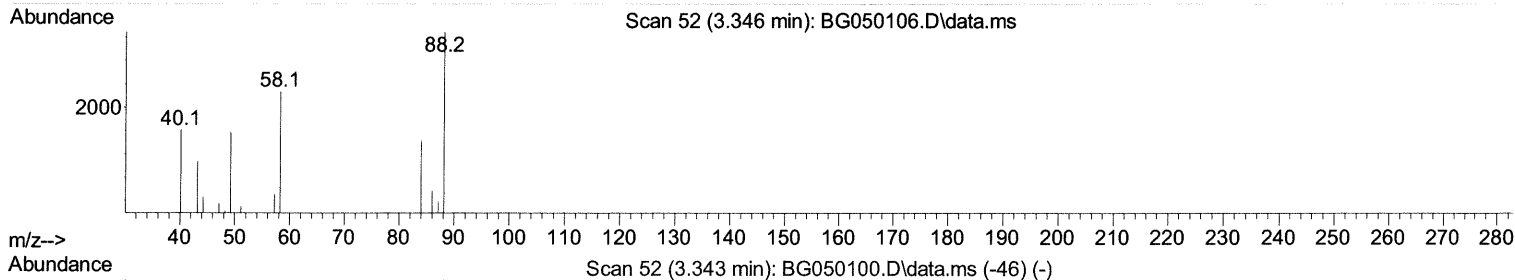
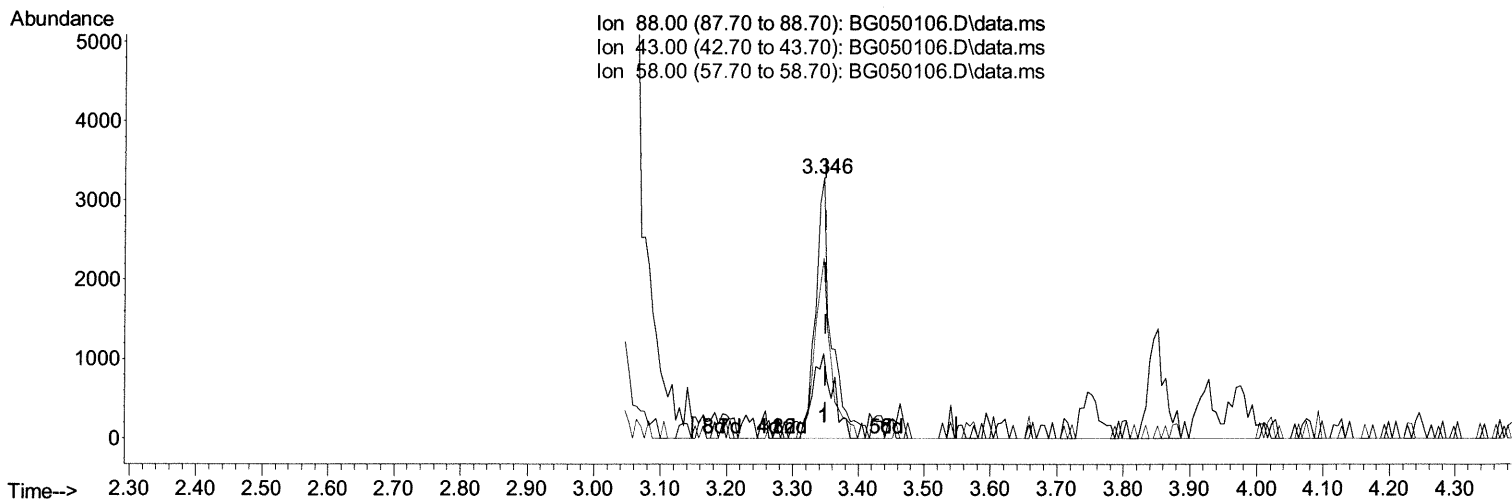
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
Data File : BG050106.D
Acq On : 2 Sep 2021 21:12
Operator : CG/JU
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(2) 1,4-Dioxane

3.346min (-0.003) 8.48 ng/uL m

response 5545

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	27.00	32.52#
58.00	65.00	69.09
0.00	0.00	0.00

Quantitation Report (Qedit)

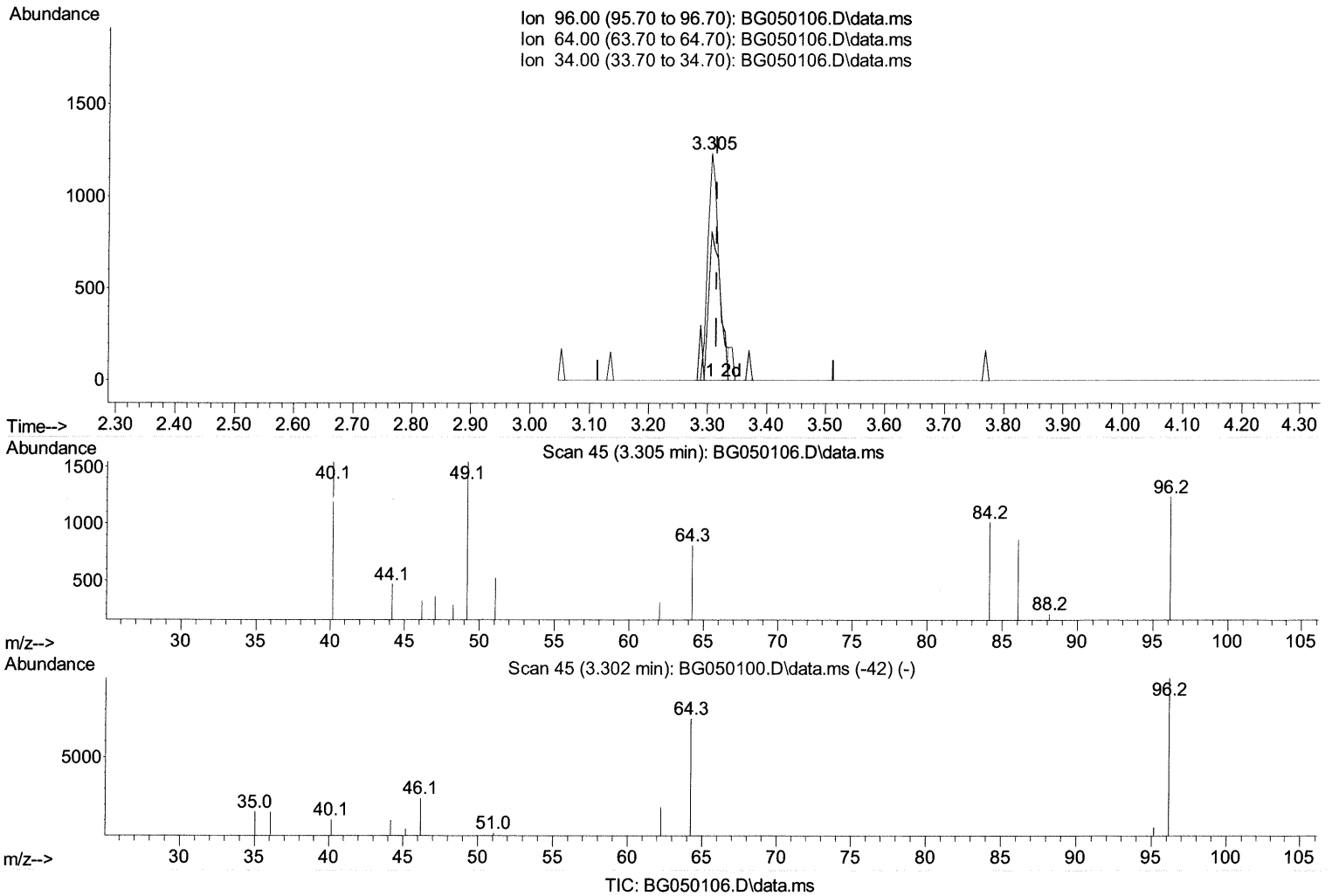
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050106.D
 Acq On : 2 Sep 2021 21:12
 Operator : CG/JU
 Sample : M3526-03MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
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(3) 1,4-Dioxane-d8 (S)

3.305min (-0.009) 3.07 ng/uL

response 1648

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	65.59
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

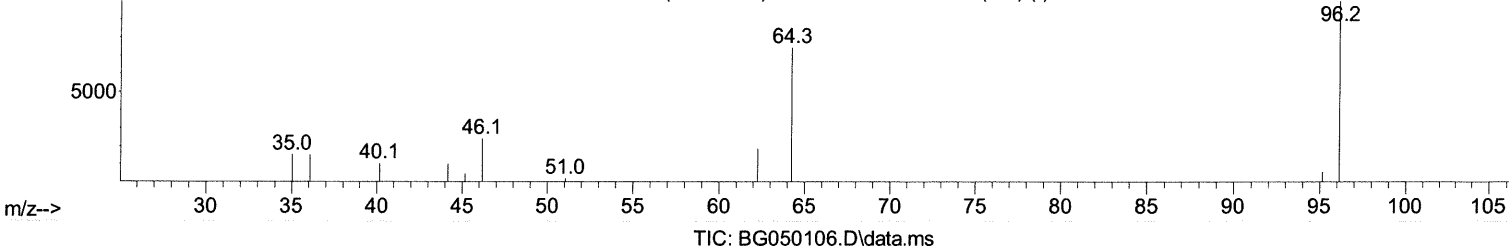
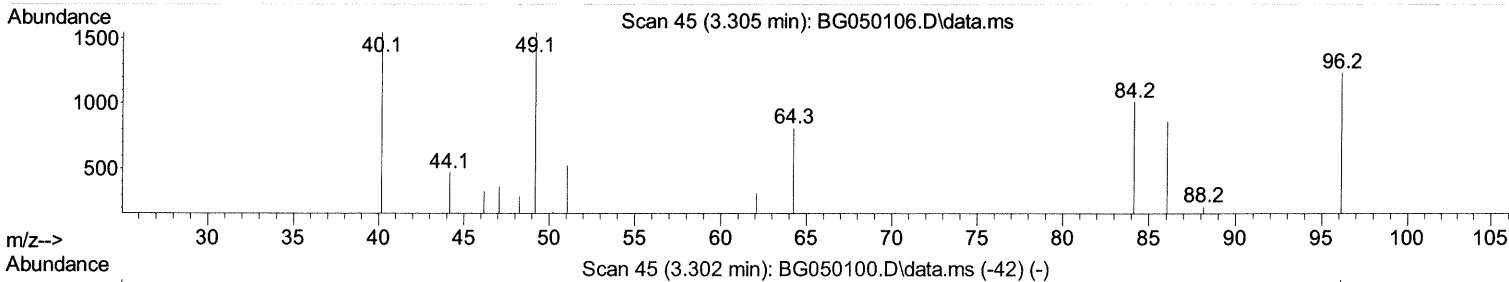
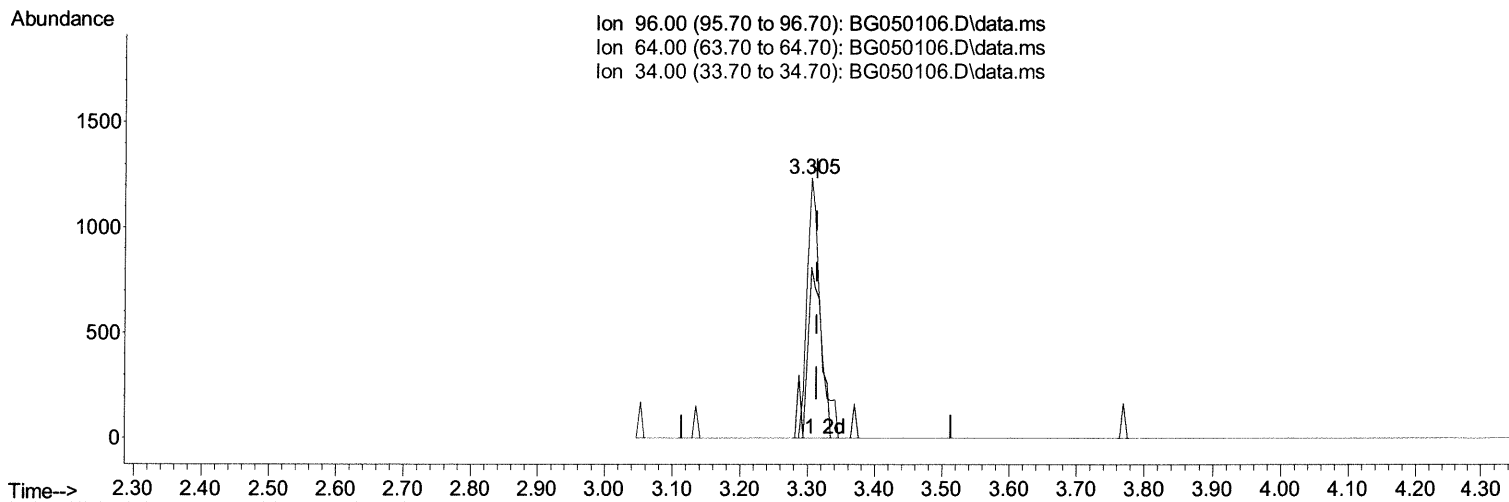
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050106.D
 Acq On : 2 Sep 2021 21:12
 Operator : CG/JU
 Sample : M3526-03MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual Integrations
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TIC: BG050106.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.305min (-0.009) 3.19 ng/uL m

response 1712

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	63.40	65.59
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

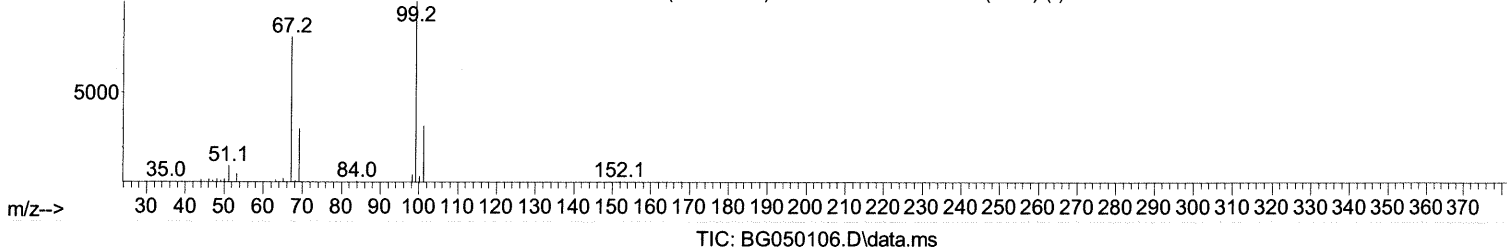
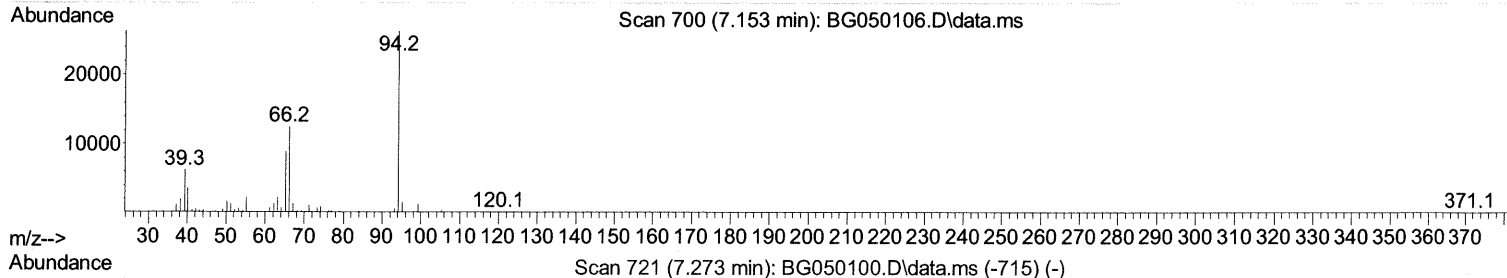
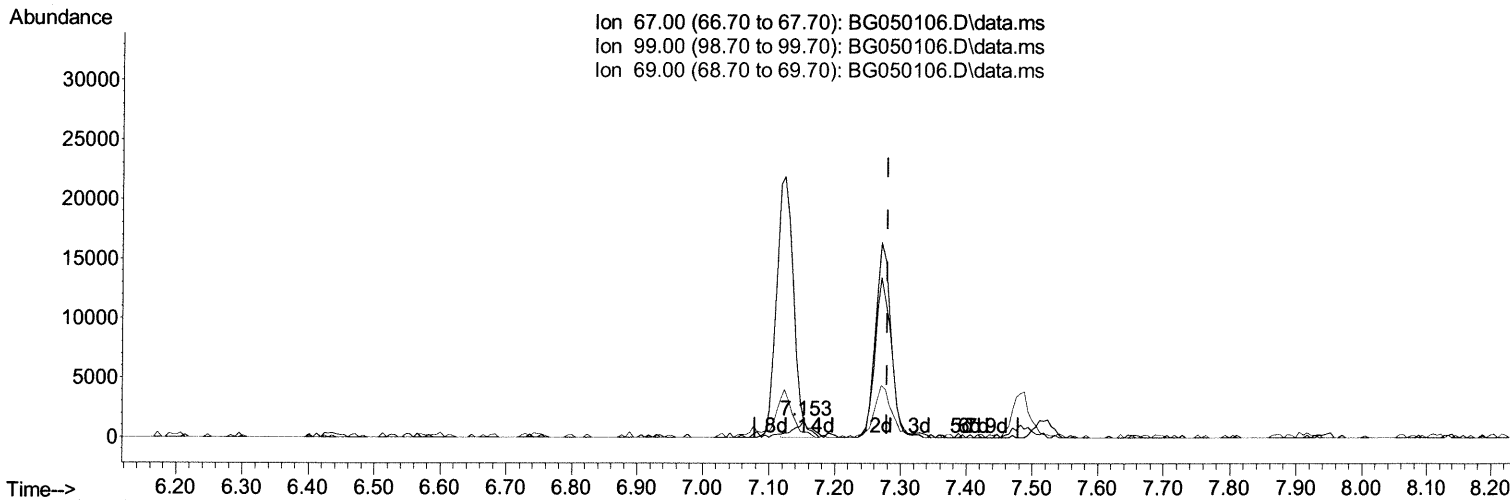
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
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(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.153min (-0.126) 1.58 ng/ul

response 2067

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	104.10	90.38
69.00	26.20	26.97
0.00	0.00	0.00

Quantitation Report (Qedit)

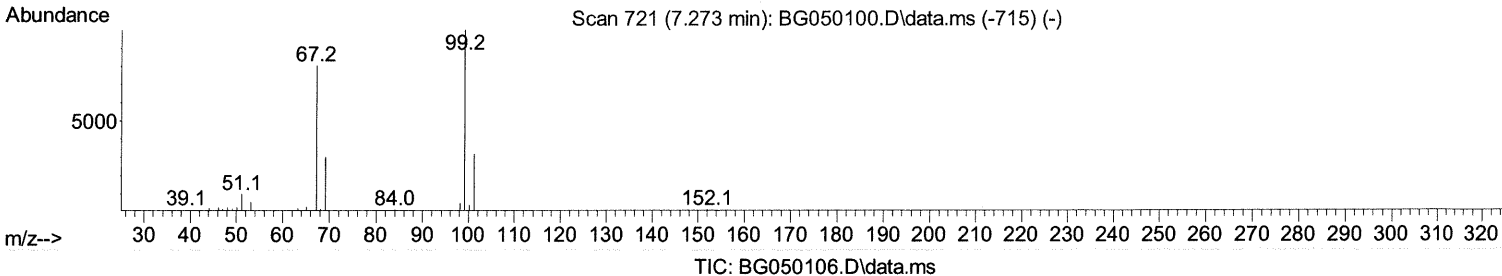
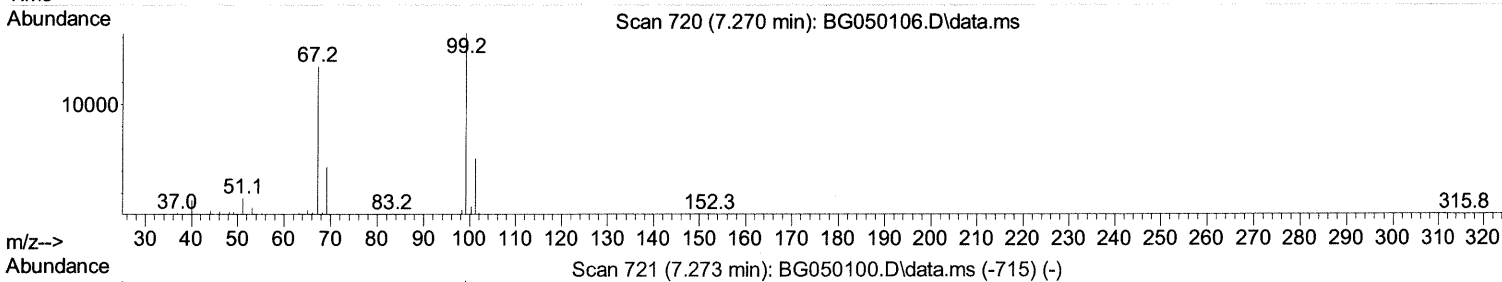
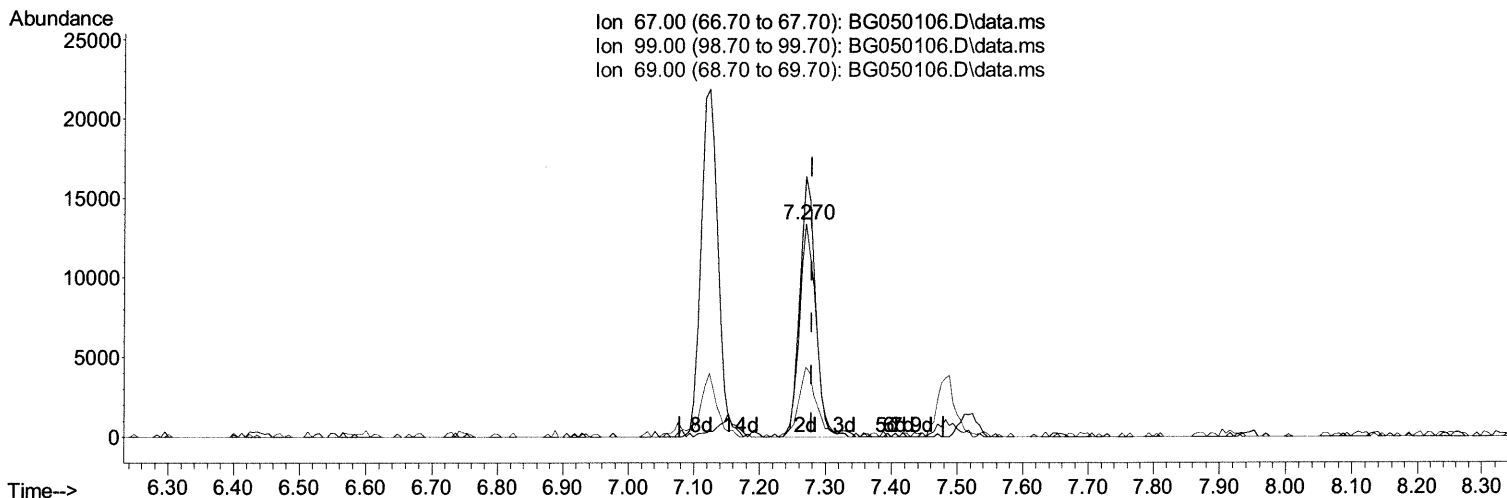
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050106.D
 Acq On : 2 Sep 2021 21:12
 Operator : CG/JU
 Sample : M3526-03MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
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TIC: BG050106.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.270min (-0.009) 18.09 ng/ul m

response 23603

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	104.10	122.09
69.00	26.20	32.56#
0.00	0.00	0.00

349/8/21

Quantitation Report (Qedit)

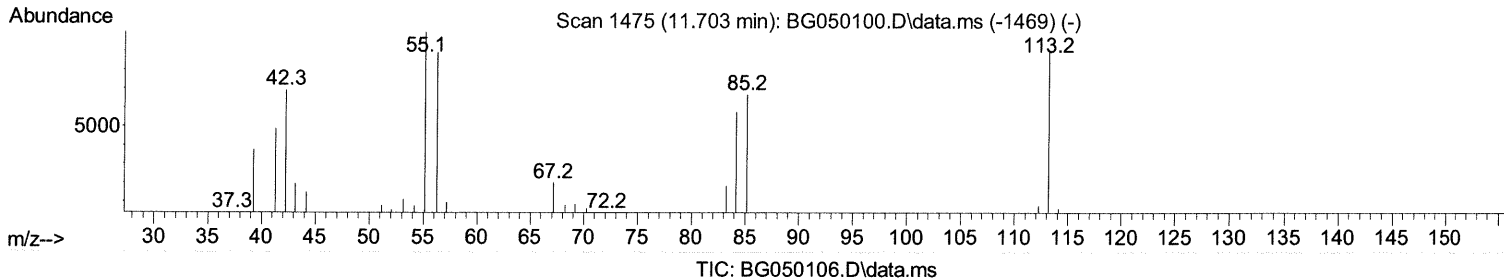
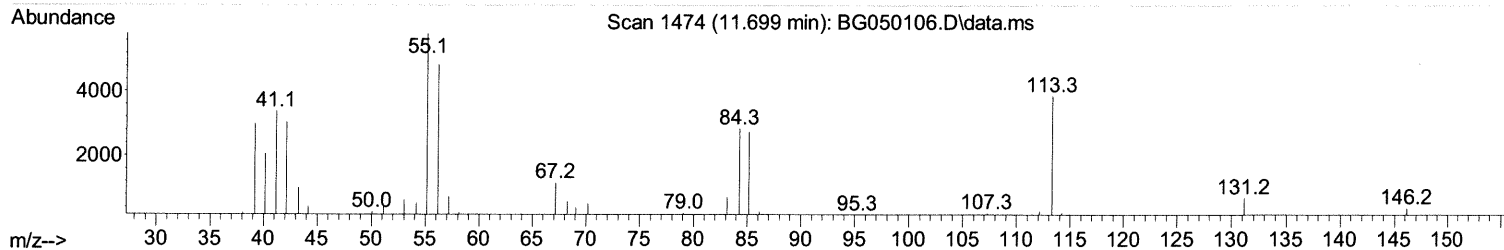
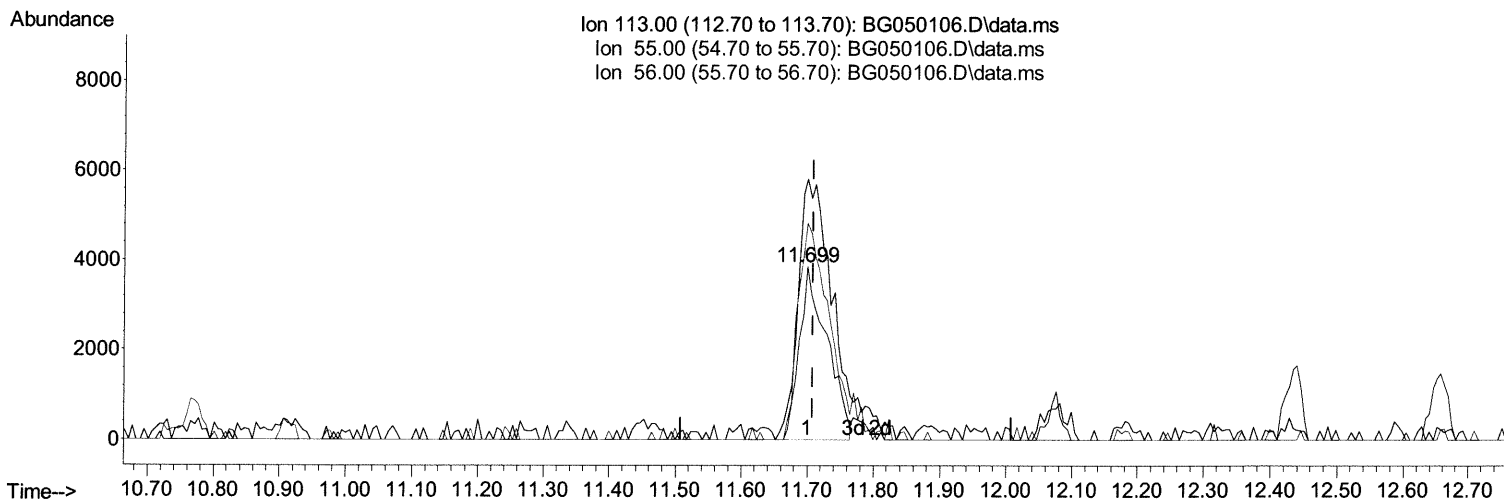
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
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 Operator : CG/JU
 Sample : M3526-03MSD
 Misc :
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Instrument :
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(34) Caprolactam

11.699min (-0.009) 18.67 ng/ul

response 11122

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	142.50	150.51
56.00	134.20	125.16
0.00	0.00	0.00

Quantitation Report (Qedit)

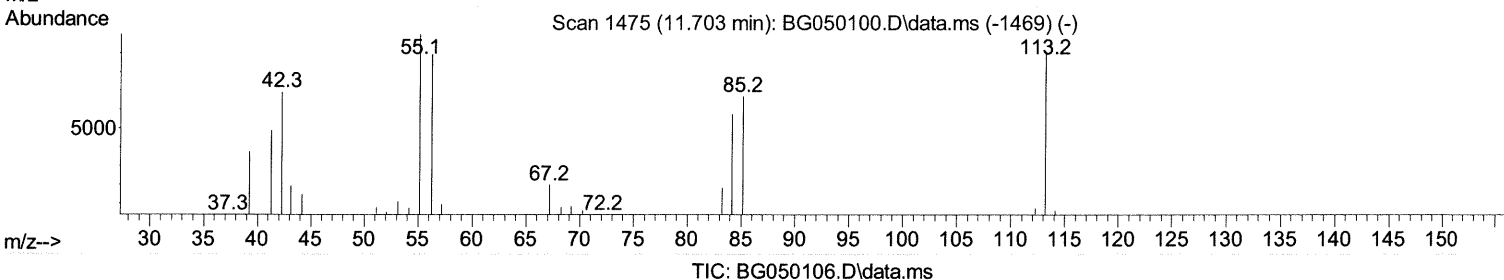
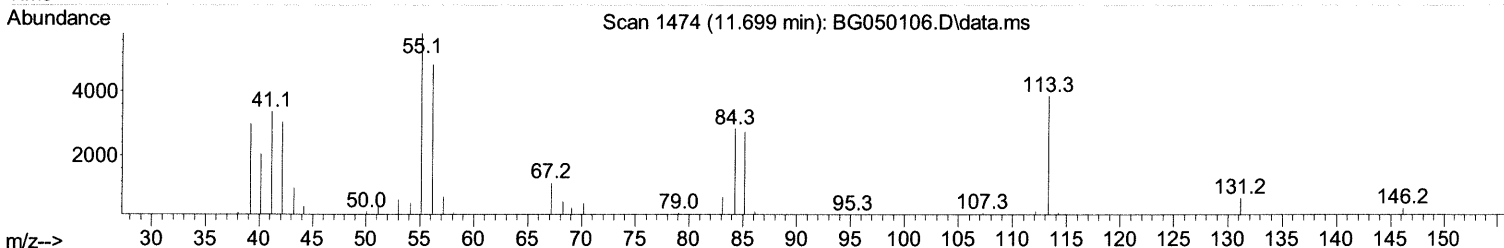
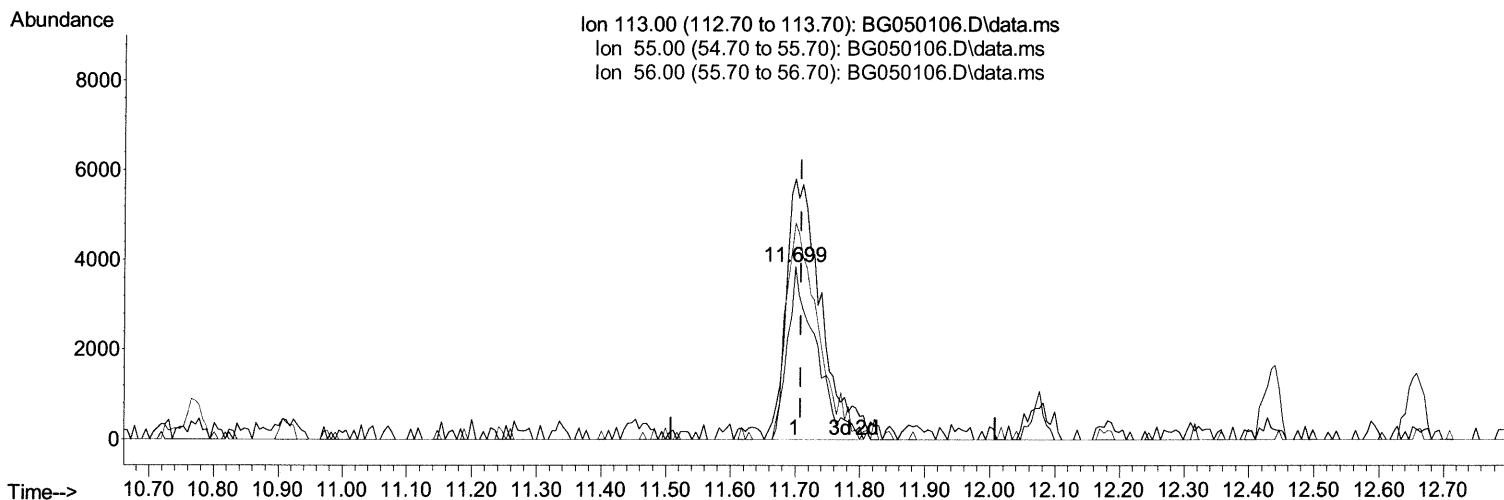
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050106.D
 Acq On : 2 Sep 2021 21:12
 Operator : CG/JU
 Sample : M3526-03MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW7A9MSD

Manual Integrations
 APPROVED

mohammad
 9/3/2021 3:31:54 PM

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(34) Caprolactam

11.699min (-0.009) 19.93 ng/ul m

response 11871

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	142.50	150.51
56.00	134.20	125.16
0.00	0.00	0.00

20 9/8/21

Quantitation Report (Qedit)

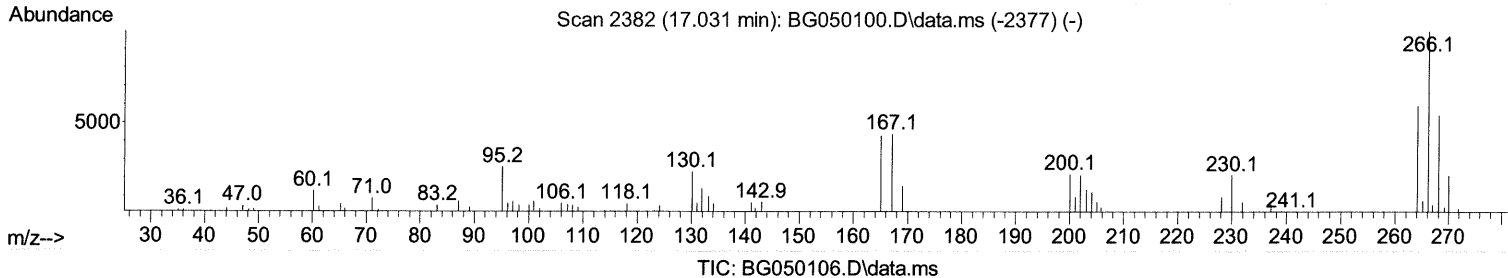
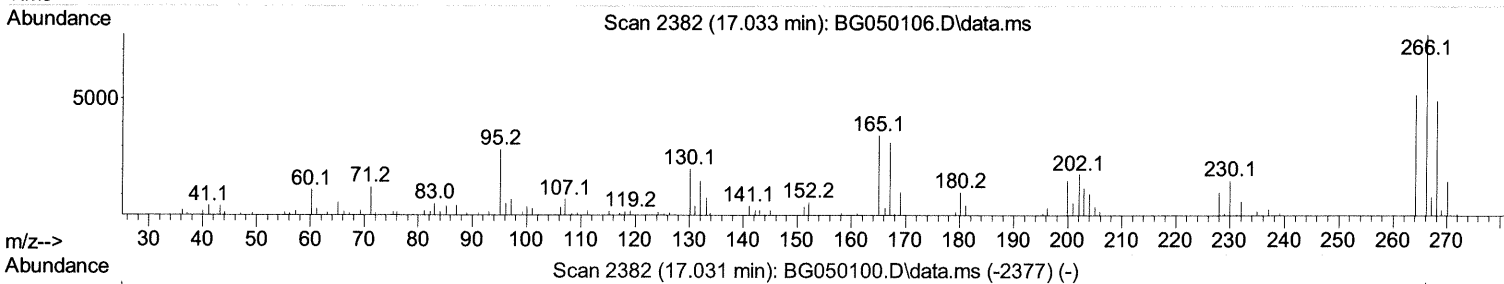
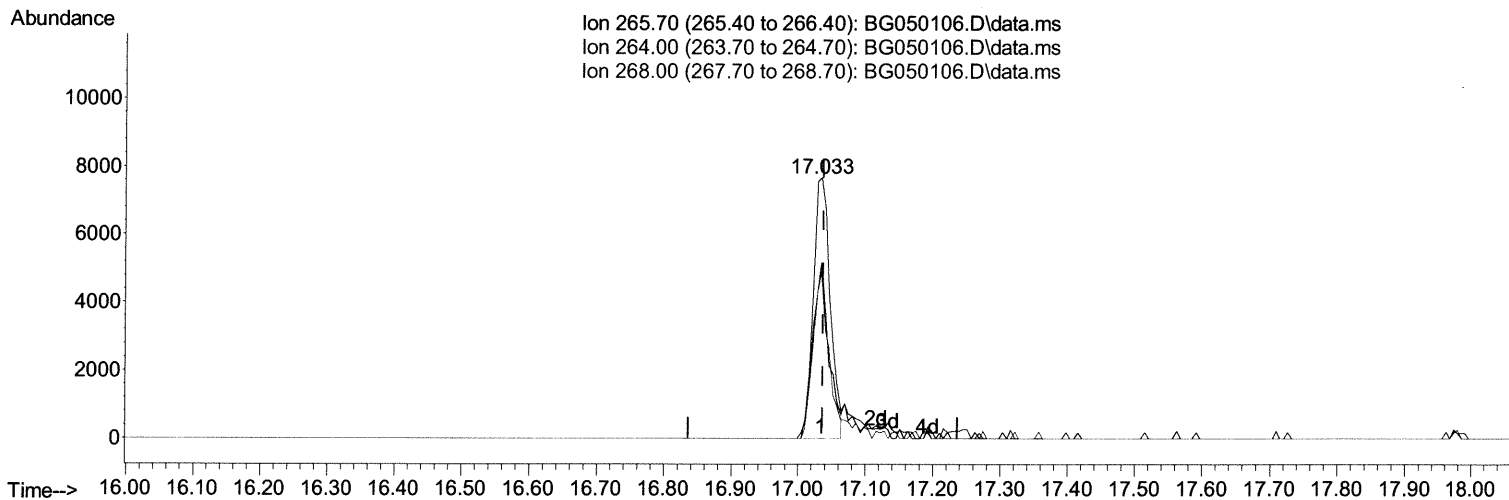
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
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TIC: BG050106.D\data.ms

(71) Pentachlorophenol (C)

17.033min (-0.003) 16.29 ng/ul

response 13423

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	67.60	67.32
268.00	65.70	64.01
0.00	0.00	0.00

Quantitation Report (Qedit)

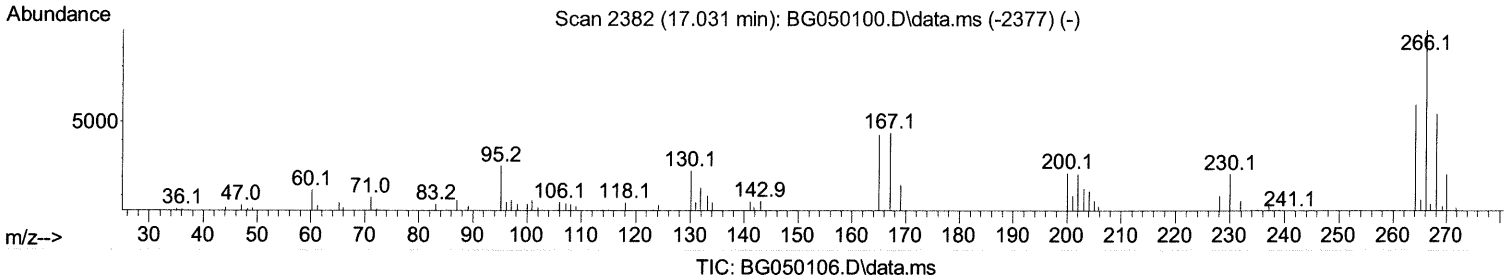
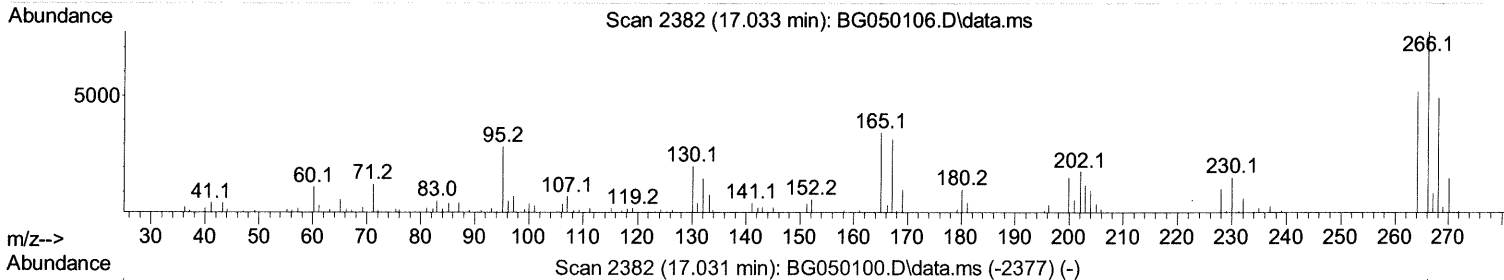
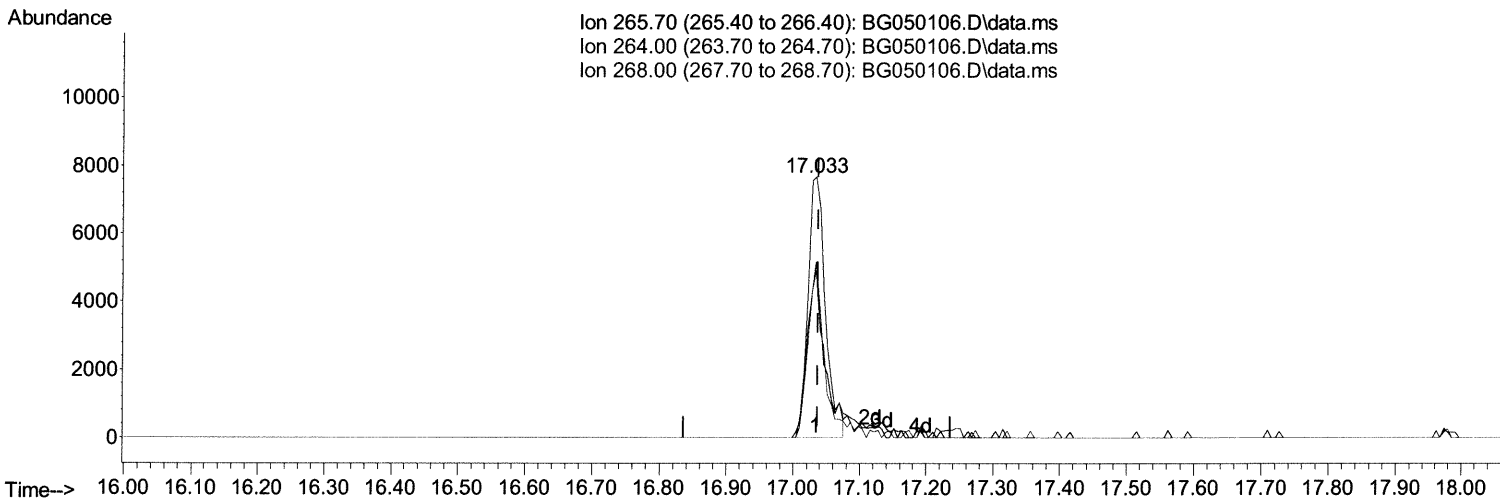
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050106.D
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 Operator : CG/JU
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 Misc :
 ALS Vial : 8 Sample Multiplier: 1

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 ClientSampleId :
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Manual Integrations
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(71) Pentachlorophenol (C)

17.033min (-0.003) 17.02 ng/ul m

response 14028

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	67.60	67.32
268.00	65.70	64.01
0.00	0.00	0.00

Quantitation Report (Qedit)

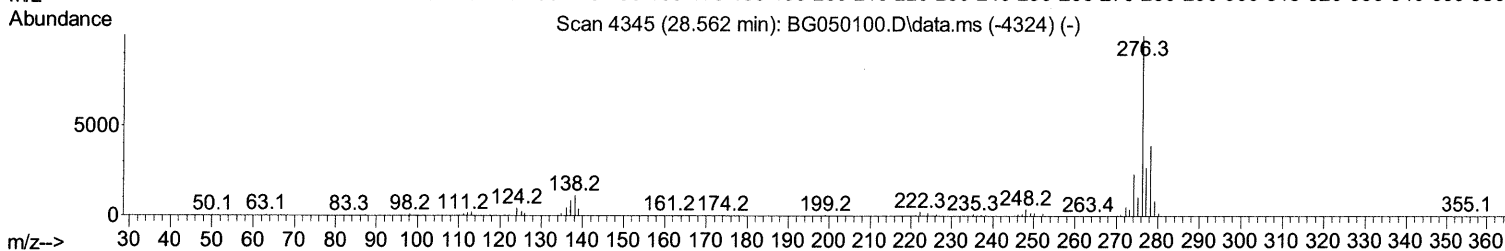
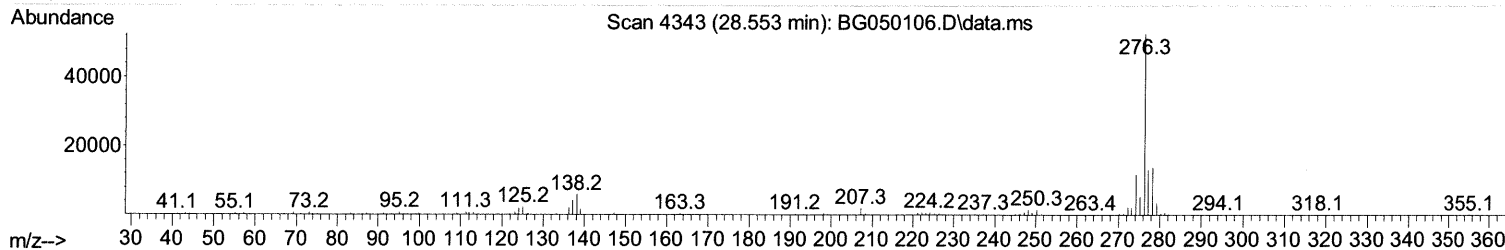
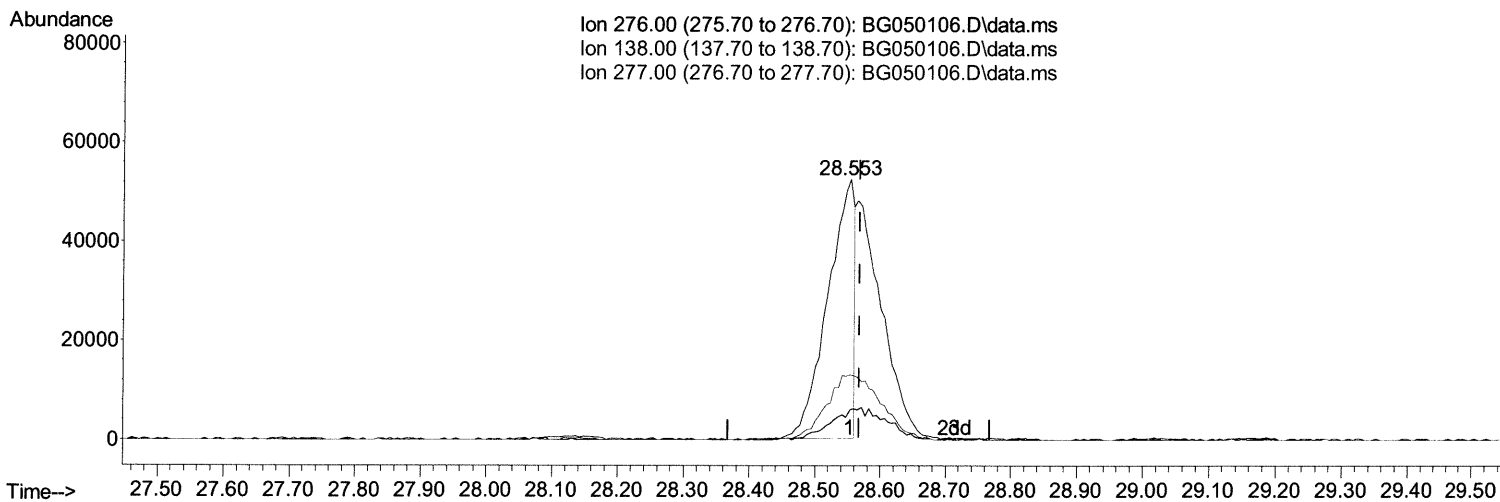
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TIC: BG050106.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

28.553min (-0.015) 13.34 ng/ul

response 148581

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	6.90	11.38#
277.00	23.20	24.84
0.00	0.00	0.00

Quantitation Report (Qedit)

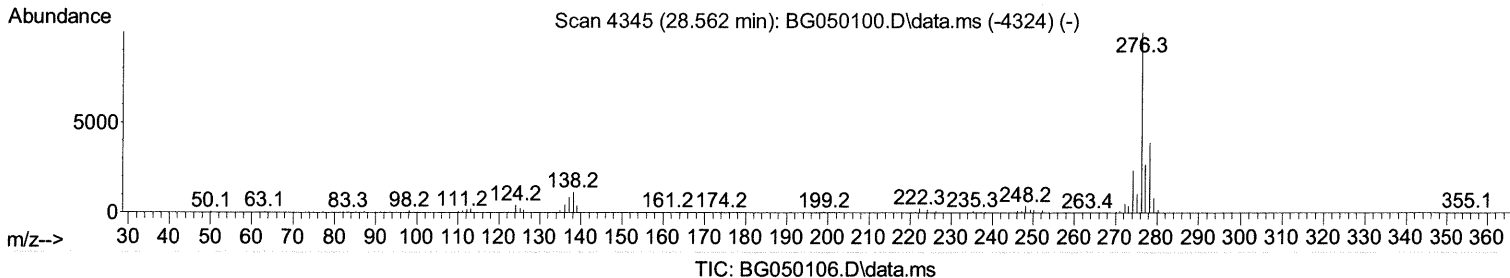
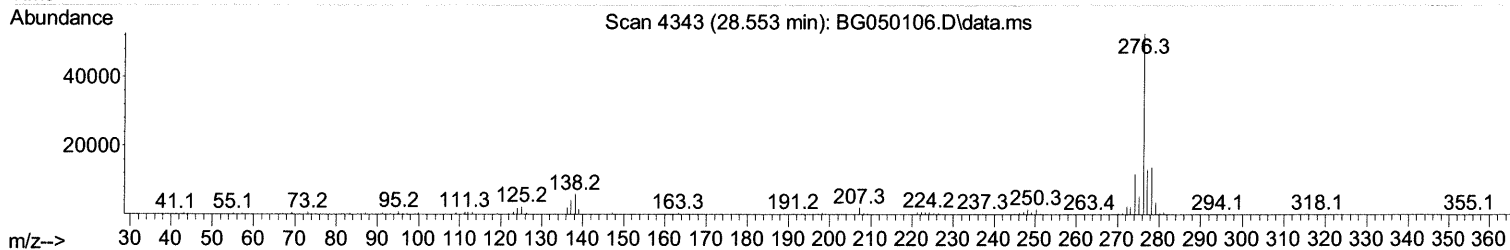
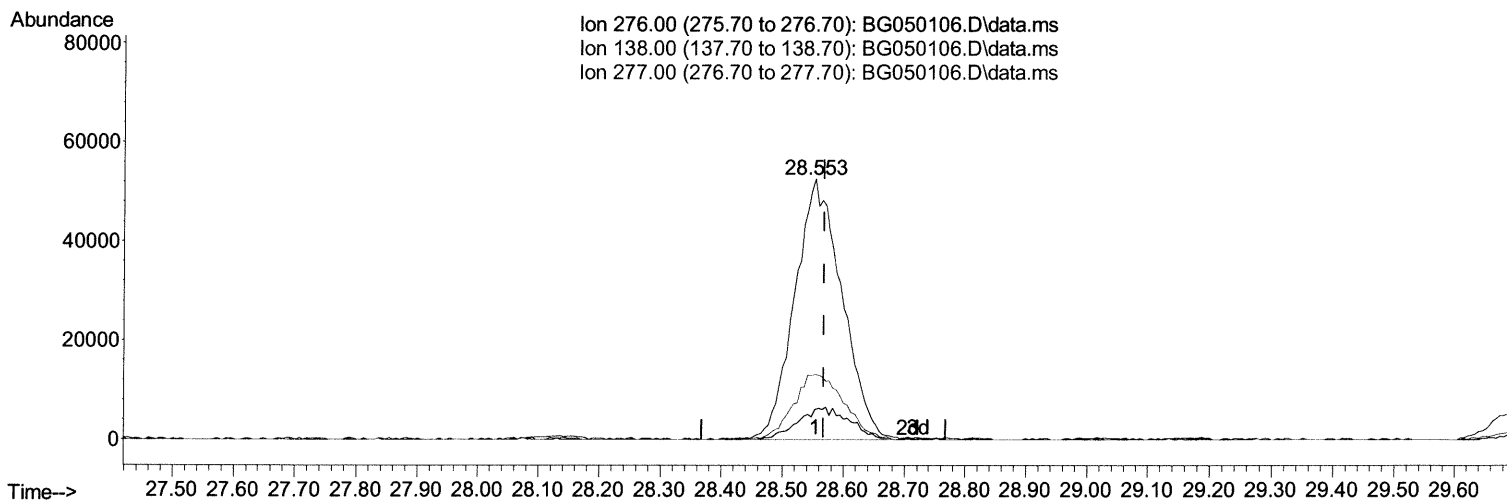
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(94) Indeno(1,2,3-cd)pyrene

28.553min (-0.015) 25.35 ng/ul m

response 282384

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	6.90	11.38#
277.00	23.20	24.84
0.00	0.00	0.00

Quantitation Report (Qedit)

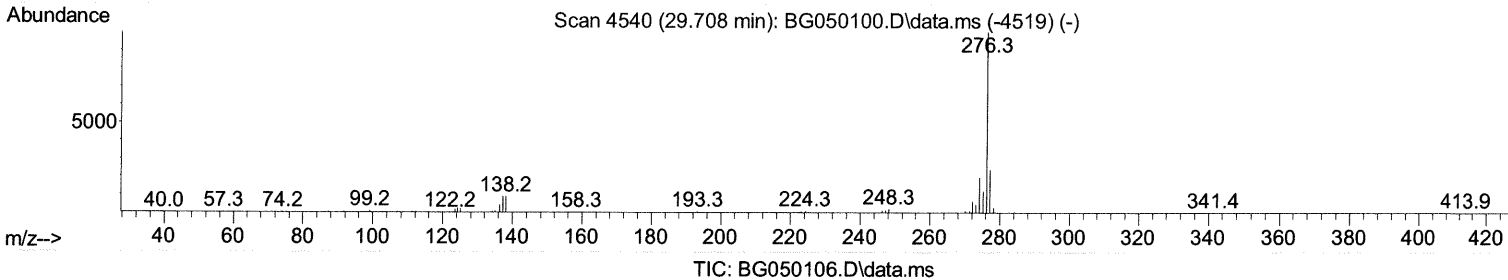
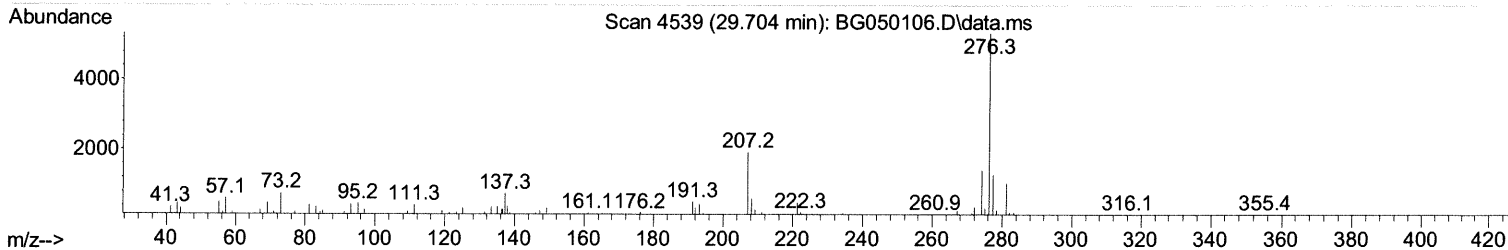
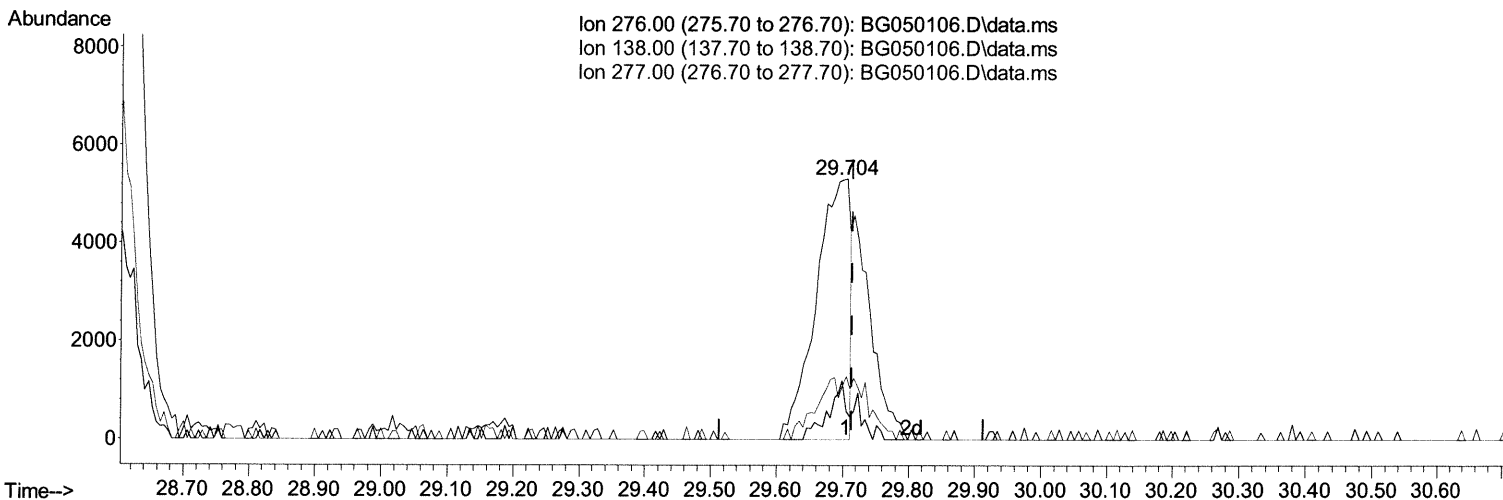
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(96) Benzo(g,h,i)perylene

29.704min (-0.009) 2.02 ng/ul

response 18875

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	10.80	10.72
277.00	21.70	24.12
0.00	0.00	0.00

Quantitation Report (Qedit)

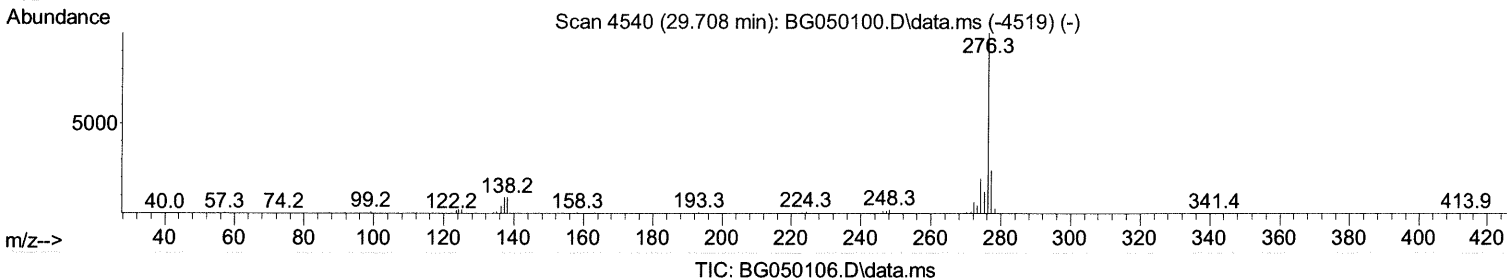
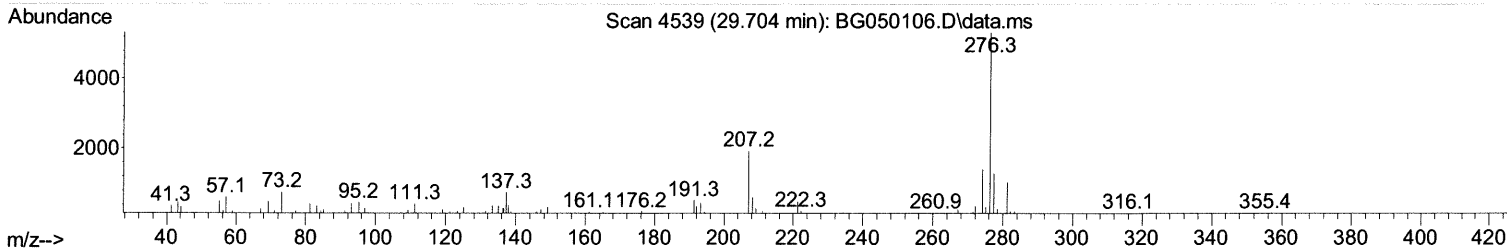
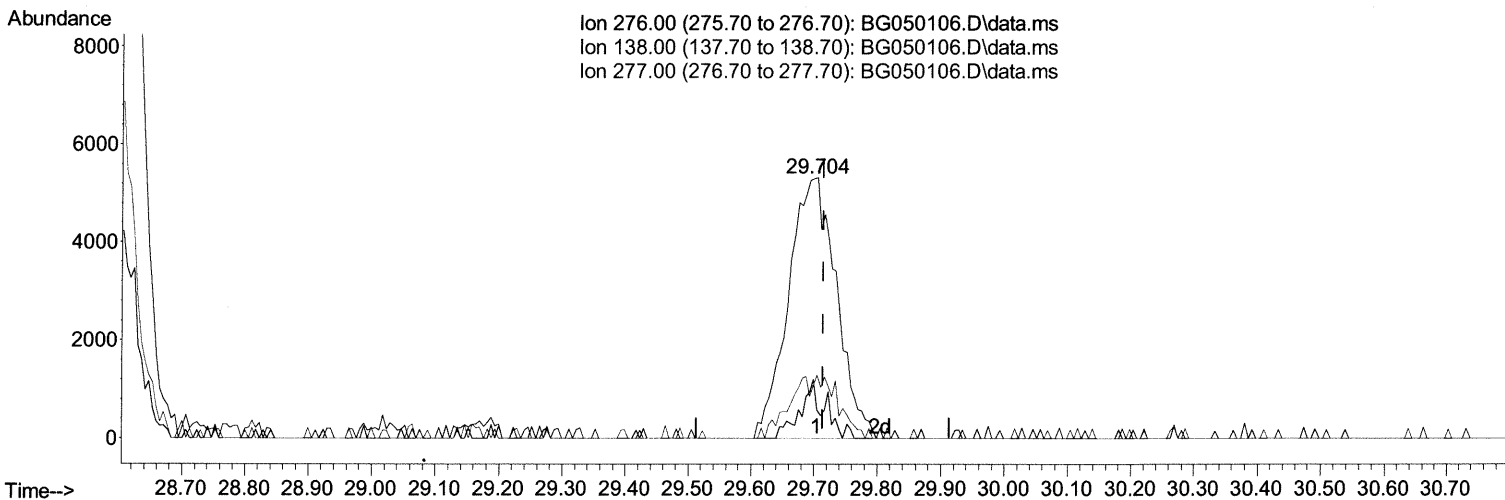
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050106.D
 Acq On : 2 Sep 2021 21:12
 Operator : CG/JU
 Sample : M3526-03MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW7A9MSD

Manual Integrations
 APPROVED

mohammad
 9/3/2021 3:31:54 PM

Quant Time: Sep 03 00:58:47 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



(96) Benzo(g,h,i)perylene

29.704min (-0.009) 2.98 ng/ul m

response 27897

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	10.80	7.13#
277.00	21.70	24.12
0.00	0.00	0.00

39/8/21

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	27654	20.000	ng/ul	0.00
20) Naphthalene-d8	10.771	136	112534	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.613	164	75975	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.368	188	152345	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.639	240	146372	20.000	ng/ul	0.00
88) Perylene-d12	24.864	264	152035	20.000	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.305	96	1712m	3.192	ng/ul	0.00	JY 8/9/21
4) Pyridine-d5	3.745	84	13565	8.398	ng/ul	0.00	
7) Phenol-d5	7.123	99	39499	16.820	ng/ul	0.00	
9) Bis-(2-Chloroethyl)eth...	7.270	67	23603m	18.092	ng/ul	0.00	JY 8/9/21
11) 2-Chlorophenol-d4	7.482	132	29372	16.716	ng/ul	0.00	
15) 4-Methylphenol-d8	8.674	113	29480	15.940	ng/ul	0.00	
21) Nitrobenzene-d5	9.121	128	16718	19.162	ng/ul	0.00	
24) 2-Nitrophenol-d4	9.855	143	18304	17.619	ng/ul	0.00	
28) 2,4-Dichlorophenol-d3	10.401	165	33777	18.229	ng/ul	0.00	
31) 4-Chloroaniline-d4	10.912	131	39964	15.846	ng/ul	0.00	
46) Dimethylphthalate-d6	14.014	166	110621	19.025	ng/ul	0.00	
49) Acenaphthylene-d8	14.308	160	132425	19.424	ng/ul	0.00	
54) 4-Nitrophenol-d4	14.848	143	12625	15.008	ng/ul	0.00	
60) Fluorene-d10	15.612	176	93644	18.995	ng/ul	0.00	
65) 4,6-Dinitro-2-methylph...	15.747	200	20004	20.410	ng/ul	0.00	
73) Anthracene-d10	17.468	188	152848	22.343	ng/ul	0.00	
81) Pyrene-d10	19.759	212	174927	22.115	ng/ul	0.00	
92) Benzo(a)pyrene-d12	24.641	264	156194	19.365	ng/ul	0.00	

Target Compounds

						Qvalue	
2) 1,4-Dioxane	3.346	88	5545m	8.482	ng/ul	90	JY 8/9/21
5) Pyridine	3.763	79	25232	14.453	ng/ul#	90	
6) Benzaldehyde	7.088	77	31502	23.312	ng/ul	93	
8) Phenol	7.147	94	48736	20.547	ng/ul	95	
10) Bis(2-Chloroethyl)ether	7.364	93	36047	22.015	ng/ul	97	
12) 2-Chlorophenol	7.517	128	38459	22.151	ng/ul#	92	
13) 2-Methylphenol	8.404	108	35708	19.513	ng/ul	95	
14) 2,2'-oxybis(1-Chloropr...	8.480	45	36085	21.016	ng/ul#	95	
16) Acetophenone	8.780	105	65902	21.850	ng/ul	95	
17) N-Nitroso-di-n-propyla...	8.762	70	32369	21.374	ng/ul	94	
18) 4-Methylphenol	8.739	108	38582	19.598	ng/ul#	74	
19) Hexachloroethane	9.032	117	16626	21.881	ng/ul	89	
22) Nitrobenzene	9.168	77	52629	22.262	ng/ul	93	
23) Isophorone	9.685	82	90821	21.562	ng/ul	98	
25) 2-Nitrophenol	9.884	139	21674	21.459	ng/ul#	92	
26) 2,4-Dimethylphenol	9.943	107	24534	10.930	ng/ul	96	
27) Bis(2-Chloroethoxy)met...	10.172	93	46979	21.599	ng/ul	98	
29) 2,4-Dichlorophenol	10.431	162	38980	22.920	ng/ul	96	
30) Naphthalene	10.818	128	128662	22.874	ng/ul	100	
32) 4-Chloroaniline	10.936	127	46685	19.221	ng/ul	93	
33) Hexachlorobutadiene	11.100	225	31270	22.729	ng/ul	92	
34) Caprolactam	11.699	113	11871m	19.927	ng/ul	92	JY 8/9/21
35) 4-Chloro-3-methylphenol	12.075	107	43629	21.465	ng/ul	98	

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 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.434	142	99676	23.839	ng/ul	98
37) 1-Methylnaphthalene	12.651	142	92946	22.023	ng/ul	91
39) 1,2,4,5-Tetrachloroben...	12.804	216	50862	22.793	ng/ul#	95
40) Hexachlorocyclopentadiene	12.774	237	3057	3.061	ng/ul#	96
41) 2,4,6-Trichlorophenol	13.051	196	31649	22.256	ng/ul	97
42) 2,4,5-Trichlorophenol	13.133	196	33414	22.419	ng/ul	91
43) 1,1'-Biphenyl	13.444	154	123553	23.277	ng/ul	99
44) 2-Chloronaphthalene	13.491	162	96889	22.725	ng/ul	95
45) 2-Nitroaniline	13.697	65	31964	22.082	ng/ul	96
47) Dimethylphthalate	14.061	163	129224	22.457	ng/ul#	99
48) 2,6-Dinitrotoluene	14.190	165	26911	22.652	ng/ul	91
50) Acenaphthylene	14.337	152	147990	23.751	ng/ul	97
51) 3-Nitroaniline	14.525	138	24940	21.755	ng/ul#	91
52) Acenaphthene	14.678	153	103976	23.004	ng/ul	93
53) 2,4-Dinitrophenol	14.754	184	10556	17.397	ng/ul#	84
55) 4-Nitrophenol	14.860	109	22674	21.258	ng/ul	95
56) Dibenzofuran	15.013	168	143793	22.972	ng/ul	99
57) 2,4-Dinitrotoluene	14.989	165	43339	25.361	ng/ul#	93
58) 2,3,4,6-Tetrachlorophenol	15.253	232	29964	24.261	ng/ul#	94
59) Diethylphthalate	15.424	149	137394	23.340	ng/ul	97
61) Fluorene	15.665	166	122105	23.122	ng/ul	96
62) 4-Chlorophenyl-phenyle...	15.653	204	63858	22.854	ng/ul	96
63) 4-Nitroaniline	15.694	138	25493	22.284	ng/ul	99
66) 4,6-Dinitro-2-methylph...	15.765	198	22204	24.994	ng/ul	98
67) N-Nitrosodiphenylamine	15.870	169	108417	27.031	ng/ul	95
68) 4-Bromophenyl-phenylether	16.552	248	45482	26.988	ng/ul	97
69) Hexachlorobenzene	16.681	284	48445	24.894	ng/ul	95
70) Atrazine	16.822	200	53462	30.004	ng/ul	99
71) Pentachlorophenol	17.033	266	14028m	17.023	ng/ul	99
72) Phenanthrene	17.409	178	243569	32.070	ng/ul	99
74) Anthracene	17.503	178	205085	26.647	ng/ul	97
75) 1,2,3,4-Tetrachloroben...	13.415	216	53788	26.316	ng/uL	95
76) Pentachlorobenzene	14.936	250	54726	25.268	ng/uL	93
77) Carbazole	17.774	167	187690	27.406	ng/ul	98
78) Di-n-butylphthalate	18.320	149	239900	27.390	ng/ul	99
80) Fluoranthene	19.424	202	310977	31.860	ng/ul	98
82) Pyrene	19.788	202	274820	28.414	ng/ul	99
83) Butylbenzylphthalate	20.664	149	101720	26.650	ng/ul	99
84) 3,3'-Dichlorobenzidine	21.533	252	68159	19.644	ng/ul#	98
85) Benzo(a)anthracene	21.621	228	269369	29.535	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.510	149	143718	27.418	ng/ul	99
87) Chrysene	21.686	228	259082	30.051	ng/ul	99
89) Di-n-octyl phthalate	22.714	149	245483	27.218	ng/ul	100
90) Benzo(b)fluoranthene	23.836	252	287291	29.541	ng/ul	100
91) Benzo(k)fluoranthene	23.906	252	260101	28.137	ng/ul	99
93) Benzo(a)pyrene	24.705	252	196458	23.261	ng/ul#	98
94) Indeno(1,2,3-cd)pyrene	28.553	276	282384m	25.346	ng/ul	99
95) Dibenzo(a,h)anthracene	28.600	278	252504	27.074	ng/ul#	97
96) Benzo(g,h,i)perylene	29.704	276	27897m	2.978	ng/ul	99

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