

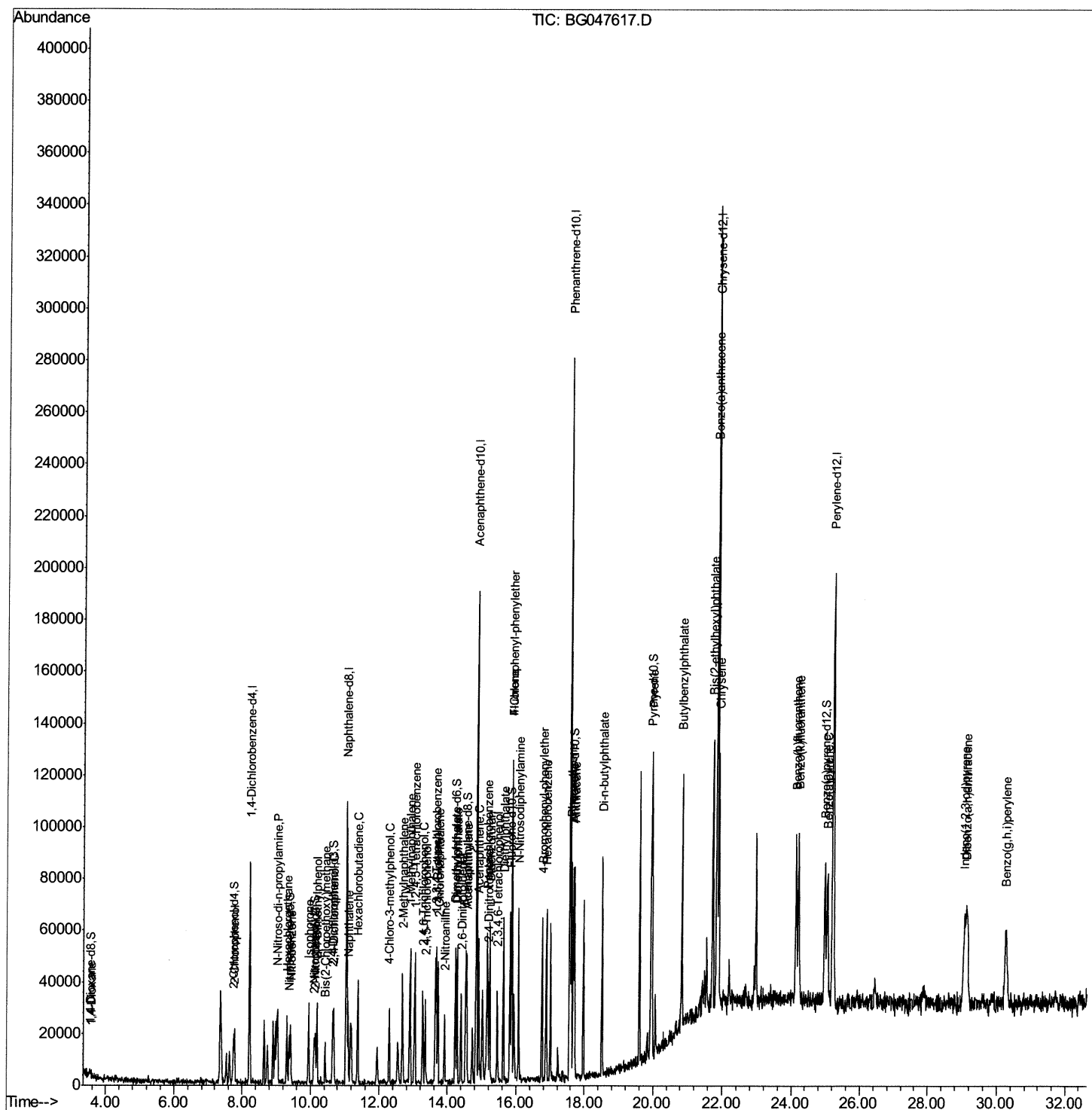
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
Data File : BG047617.D
Acq On : 7 Dec 2020 14:23
Operator : CG/JU
Sample : SSTD00525
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTD00525

Manual Integrations
APPROVED

mohammad
12/9/2020 12:18:23 PM

Quant Time: Dec 07 18:38:01 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG120720MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Dec 07 16:26:47 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

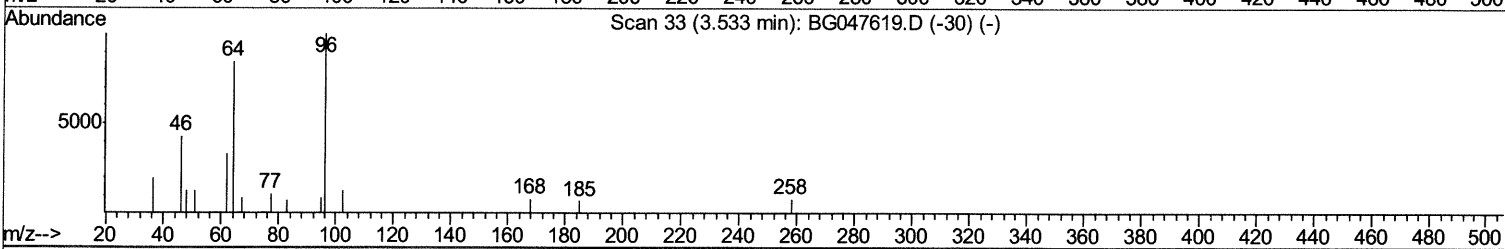
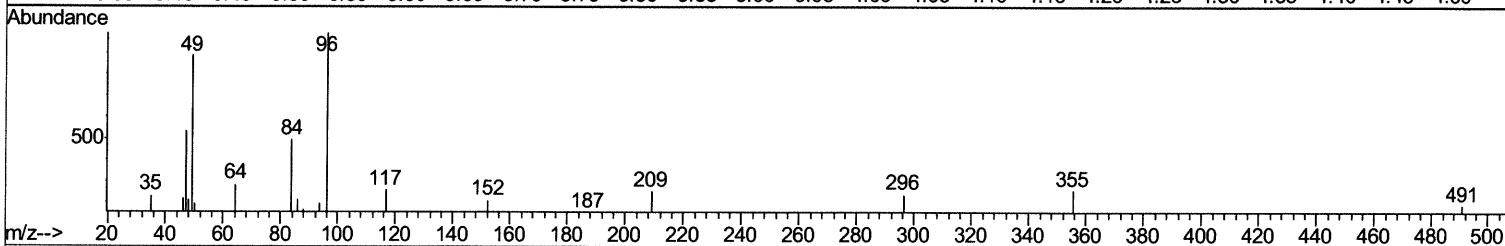
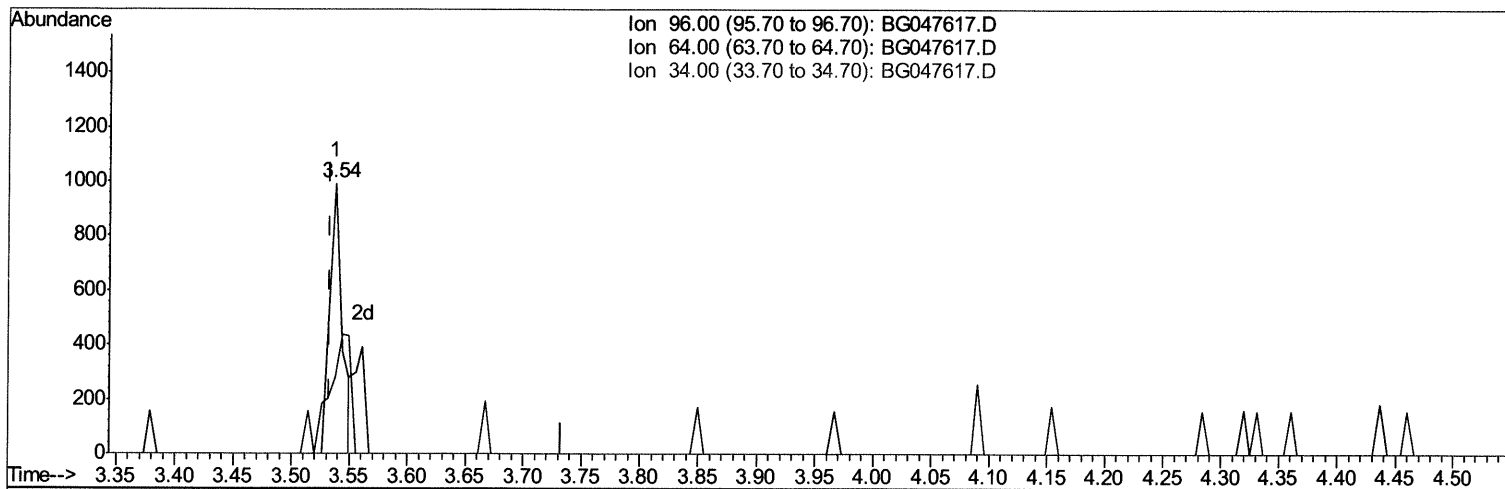
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
Data File : BG047617.D
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Operator : CG/JU
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Manual Integrations
APPROVED

mohammad
12/9/2020 12:18:23 PM

Quant Time: Dec 07 17:16:57 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG120720MA.M
Quant Title : SVOA CALIBRATION
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TIC: BG047617.D

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

3.538min (+0.005) 1.20ng/uL

response 746

Ion	Exp%	Act%
96.00	100	100
64.00	84.30	28.36#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

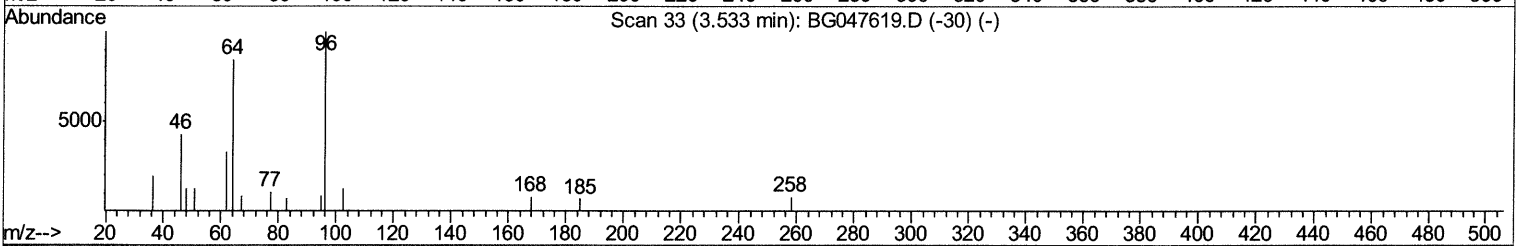
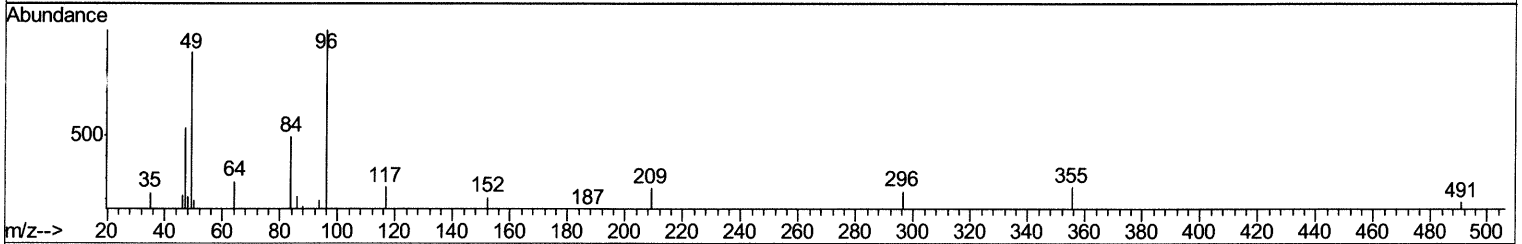
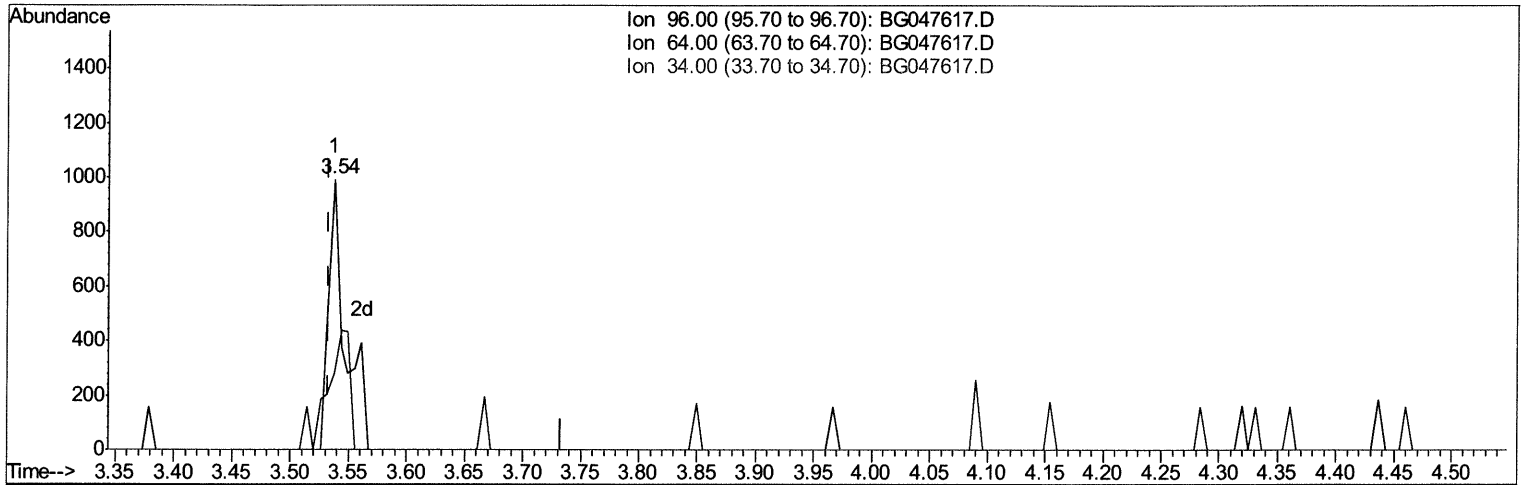
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
Data File : BG047617.D
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Manual Integrations
APPROVED

mohammad
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Quant Time: Dec 07 17:16:57 2020
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TIC: BG047617.D

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

3.538min (+0.005) 1.60ng/uL m 12/09/20 JU

response 989

Ion	Exp%	Act%
96.00	100	100
64.00	84.30	28.36#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

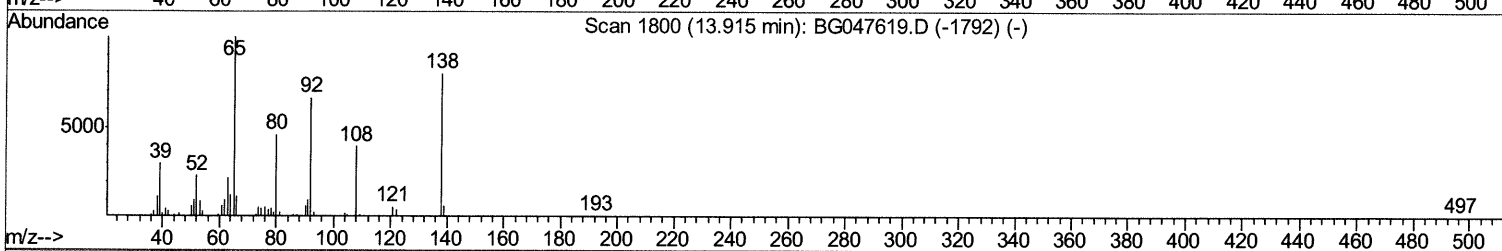
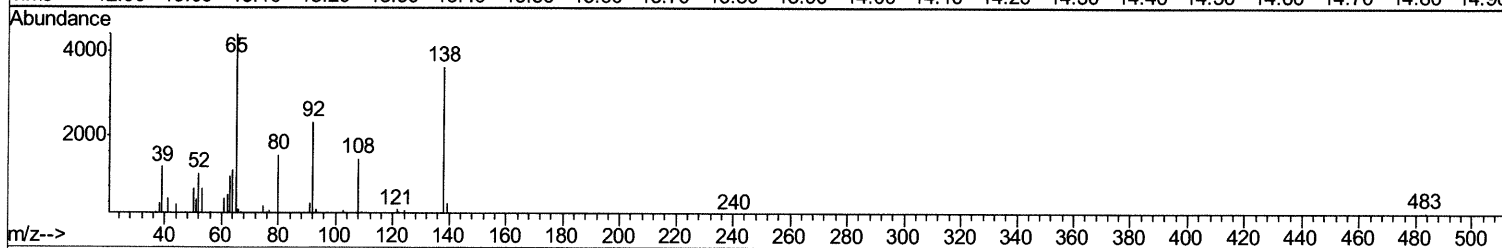
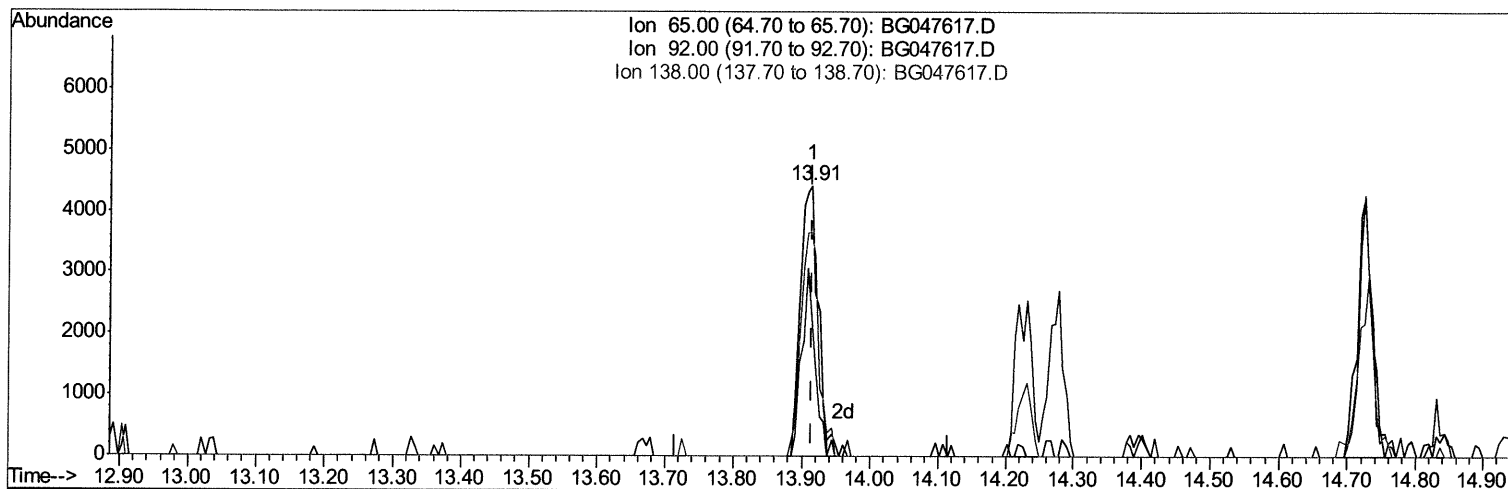
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
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Manual Integrations
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TIC: BG047617.D

(43) 2-Nitroaniline

13.914min (-0.001) 5.14ng/ul

response 7842

Ion	Exp%	Act%
65.00	100	100
92.00	66.40	52.44#
138.00	79.80	82.17
0.00	0.00	0.00

Quantitation Report (Qedit)

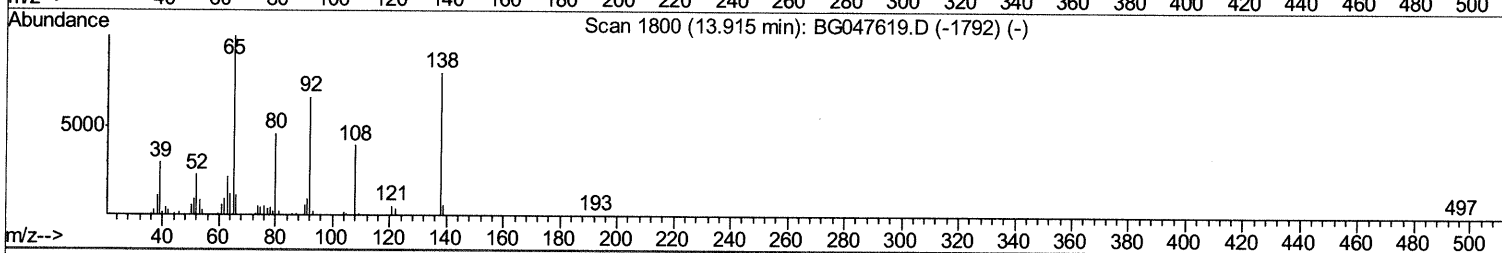
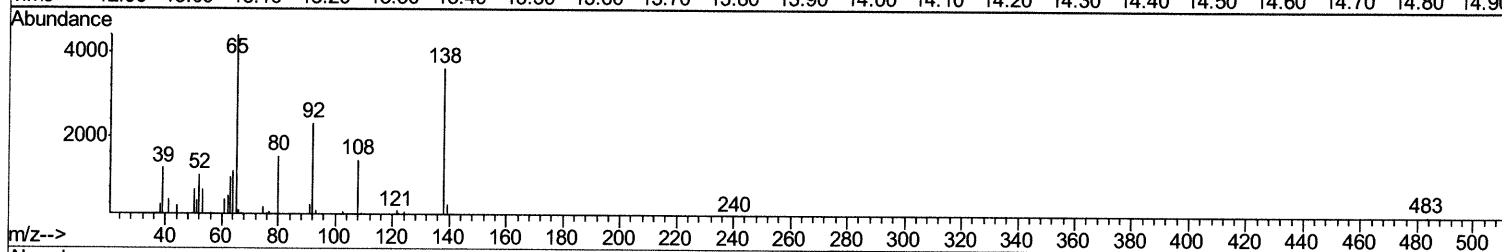
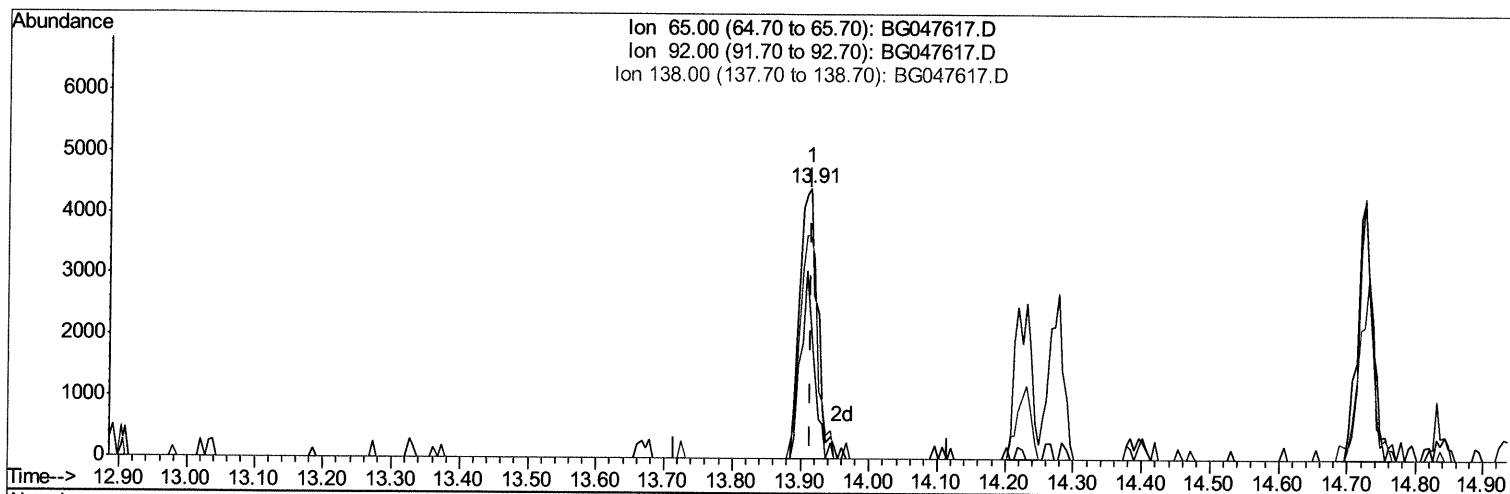
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
Data File : BG047617.D
Acq On : 7 Dec 2020 14:23
Operator : CG/JU
Sample : SST00525
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTD00525

Manual Integrations
APPROVED

mohammad
12/9/2020 12:18:23 PM

Quant Time: Dec 07 17:16:57 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG120720MA.M
Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BG047617.D

(43) 2-Nitroaniline

13.914min (-0.001) 5.26ng/ul m 1210912054

response 8040

Ion	Exp%	Act%
65.00	100	100
92.00	66.40	52.44#
138.00	79.80	82.17
0.00	0.00	0.00

Quantitation Report (Qedit)

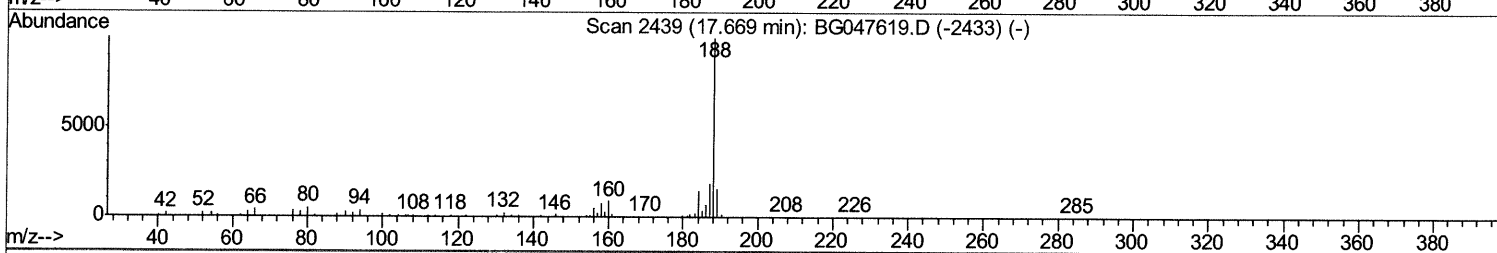
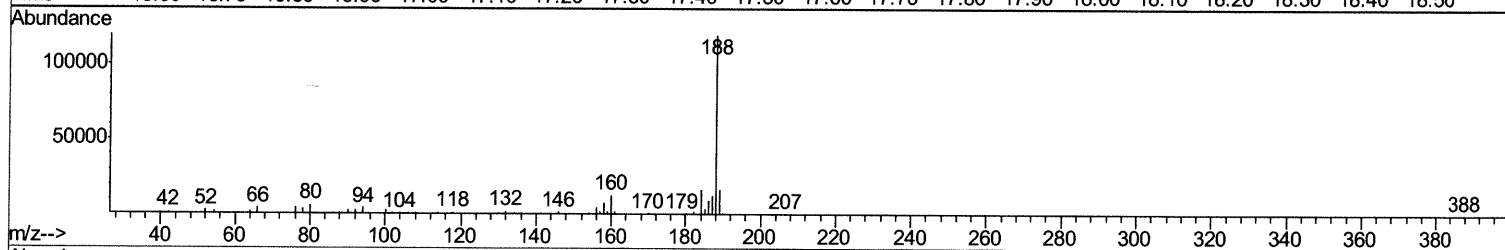
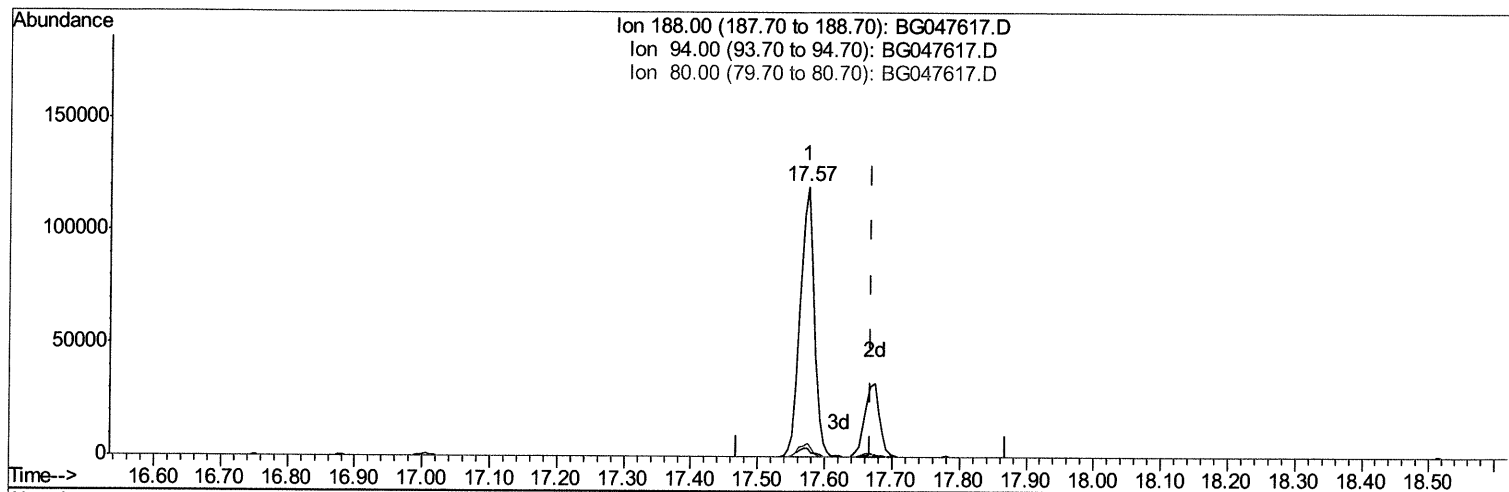
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
 Data File : BG047617.D
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 Sample : SST00525
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual Integrations
 APPROVED

mohammad
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TIC: BG047617.D

(71) Anthracene-d10 (S)

17.574min (-0.095) 20.16ng/ul

response 174553

Ion	Exp%	Act%
188.00	100	100
94.00	3.50	3.25
80.00	4.90	5.09
0.00	0.00	0.00

Quantitation Report (Qedit)

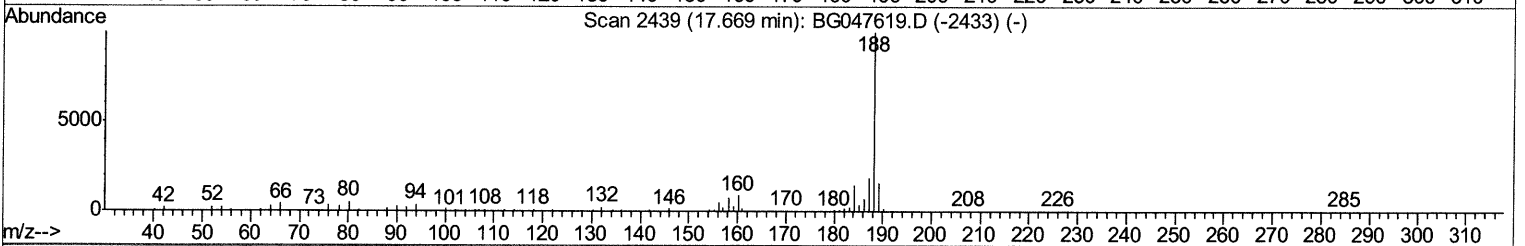
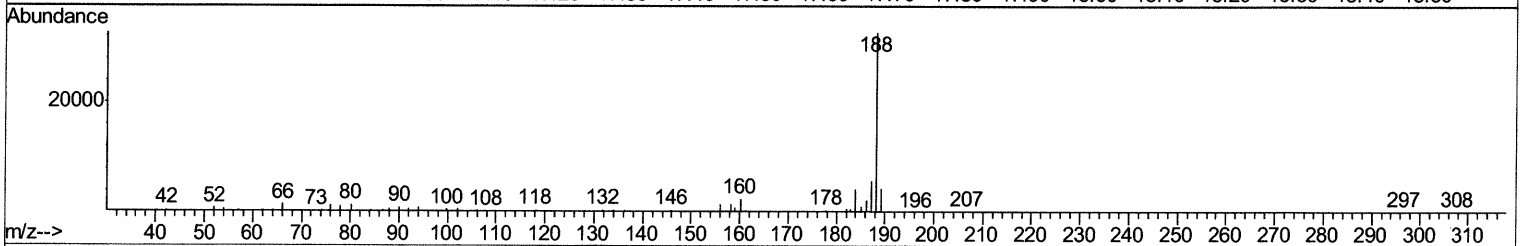
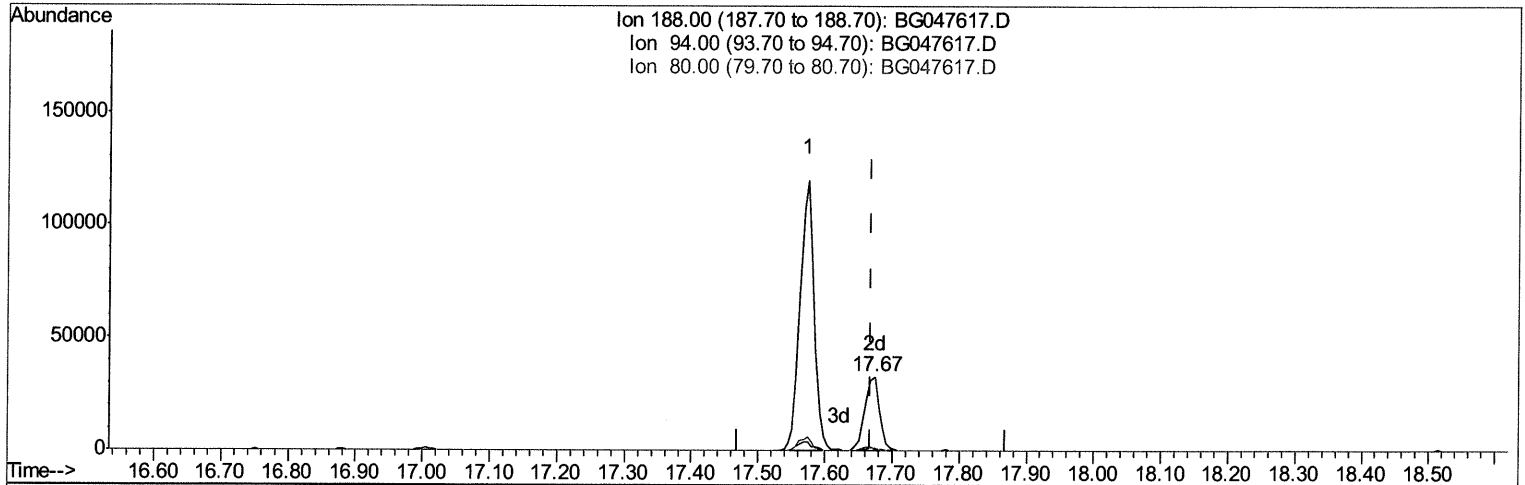
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
 Data File : BG047617.D
 Acq On : 7 Dec 2020 14:23
 Operator : CG/JU
 Sample : SST00525
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual Integrations
 APPROVED

mohammad
 12/9/2020 12:18:23 PM

Quant Time: Dec 07 17:16:57 2020
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TIC: BG047617.D

(71) Anthracene-d10 (S)

17.674min (+0.005) 5.74ng/ul m 12/09/2020

response 49692

Ion	Exp%	Act%
188.00	100	100
94.00	3.50	2.49#
80.00	4.90	4.16
0.00	0.00	0.00

Quantitation Report (Qedit)

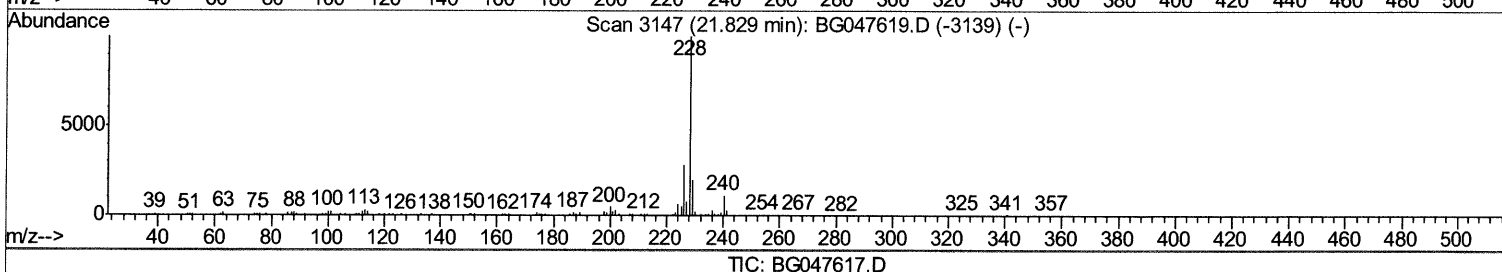
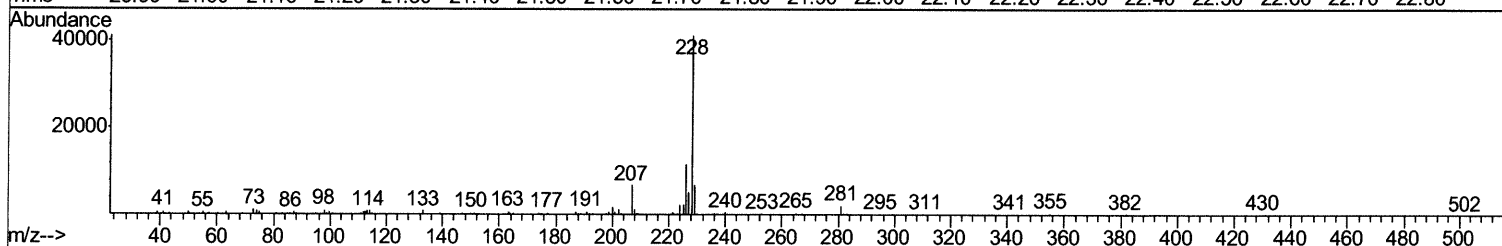
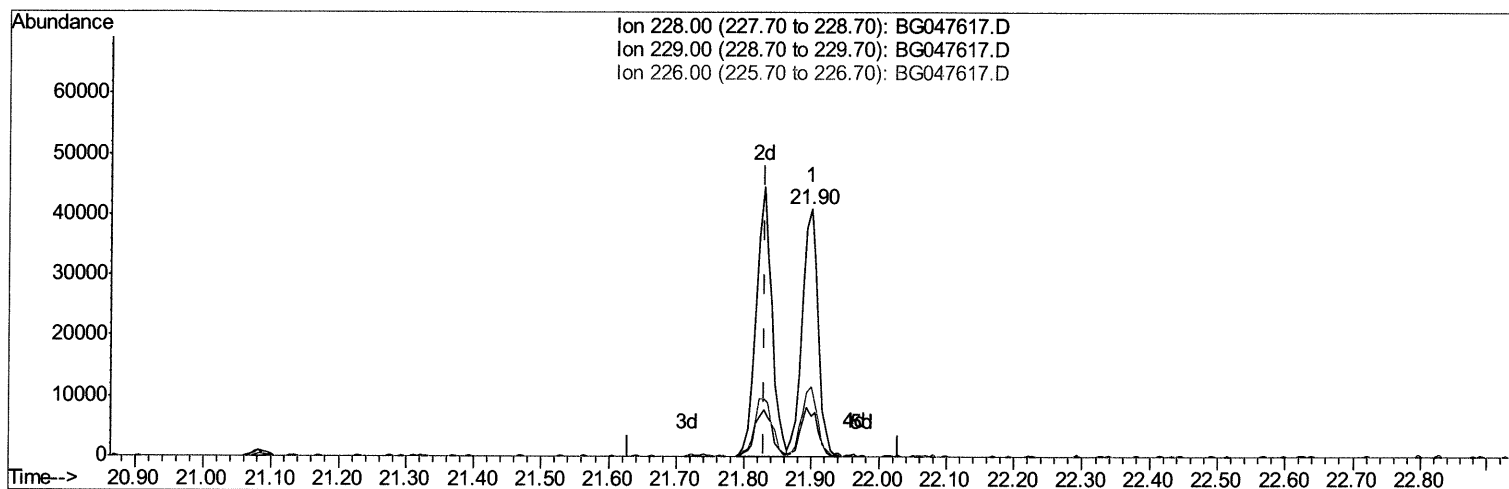
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
Data File : BG047617.D
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Instrument :
BNA_G
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SSTD00525

Manual Integrations
APPROVED

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(83) Benzo(a)anthracene

21.898min (+0.070) 5.05ng/ul

response 67615

Ion	Exp%	Act%
228.00	100	100
229.00	20.00	16.56
226.00	28.00	28.41
0.00	0.00	0.00

Quantitation Report (Qedit)

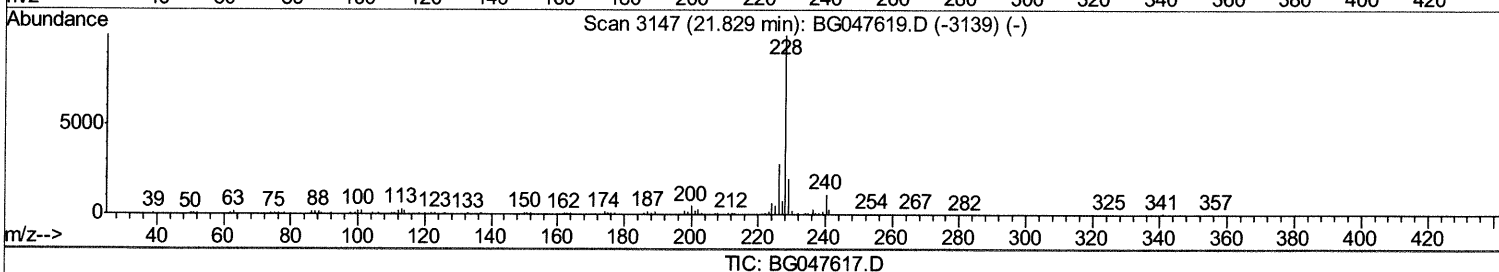
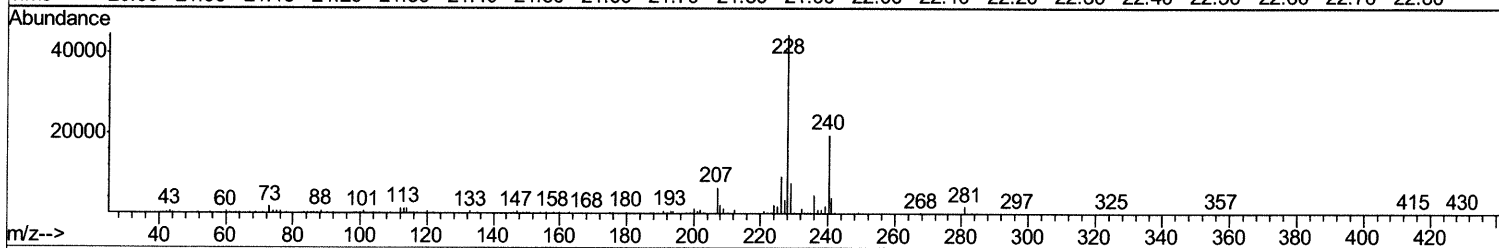
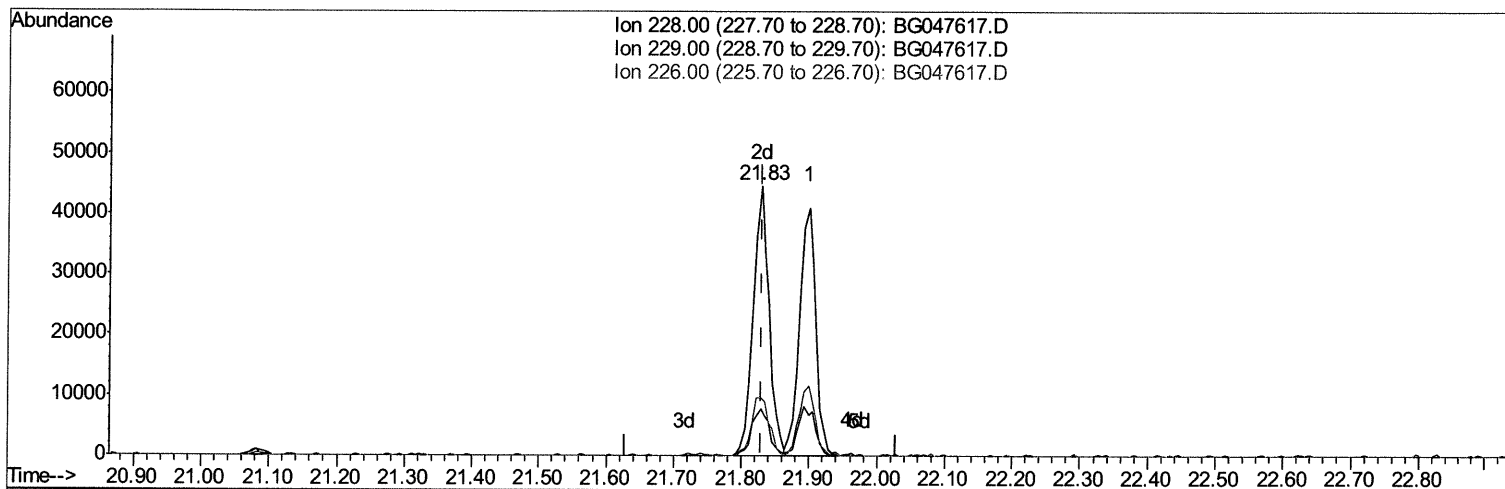
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
Data File : BG047617.D
Acq On : 7 Dec 2020 14:23
Operator : CG/JU
Sample : SSTD00525
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTD00525

Manual Integrations
APPROVED

mohammad
12/9/2020 12:18:23 PM

Quant Time: Dec 07 17:16:57 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG120720MA.M
Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BG047617.D

(83) Benzo(a)anthracene

21.828min (-0.001) 5.39ng/ul m 12/04/20 10

response 72099

Ion	Exp%	Act%
228.00	100	100
229.00	20.00	17.21
226.00	28.00	21.18#
0.00	0.00	0.00

Quantitation Report (Qedit)

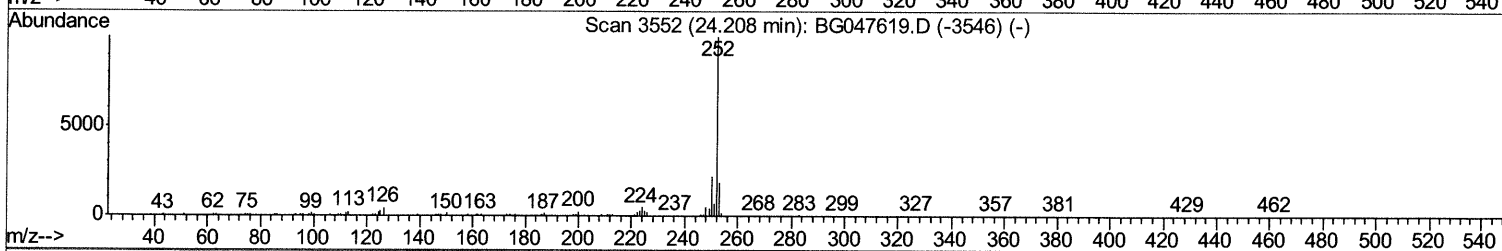
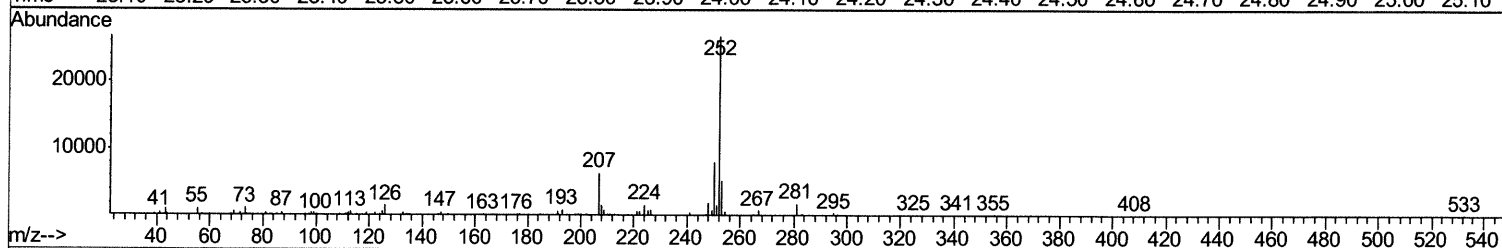
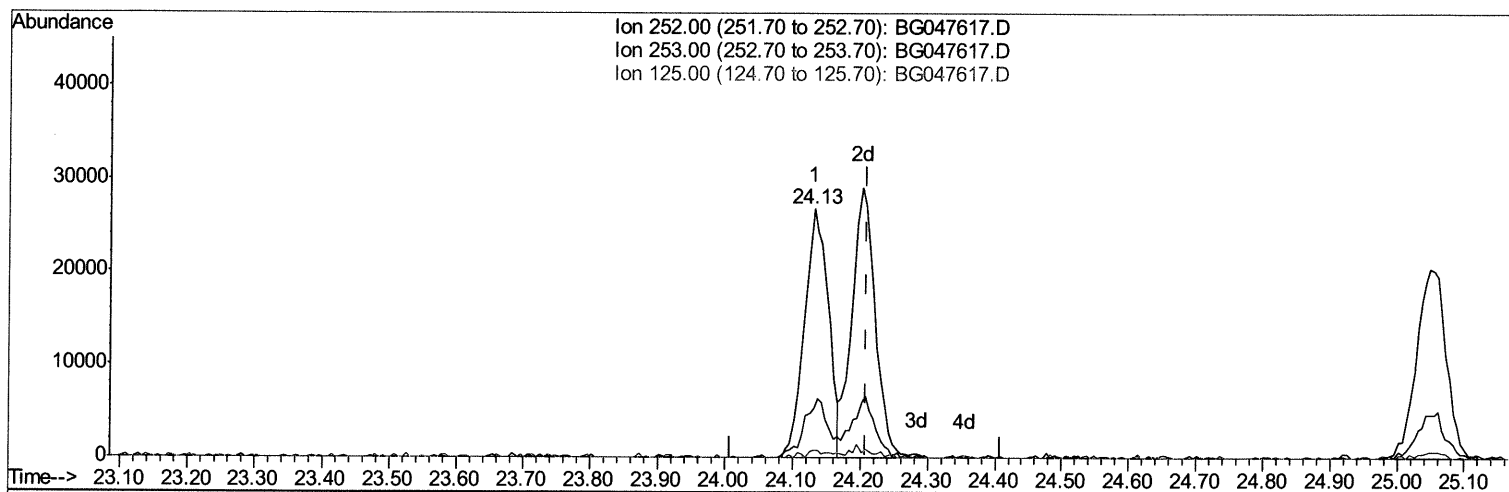
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
 Data File : BG047617.D
 Acq On : 7 Dec 2020 14:23
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Manual Integrations
 APPROVED

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TIC: BG047617.D

(89) Benzo(k)fluoranthene

24.131min (-0.077) 4.97ng/ul

response 65712

Ion	Exp%	Act%
252.00	100	100
253.00	21.10	19.57
125.00	2.80	2.81
0.00	0.00	0.00

Quantitation Report (Qedit)

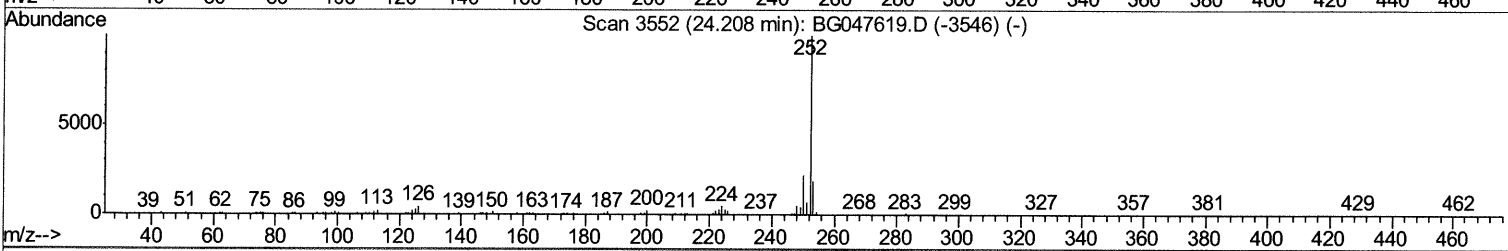
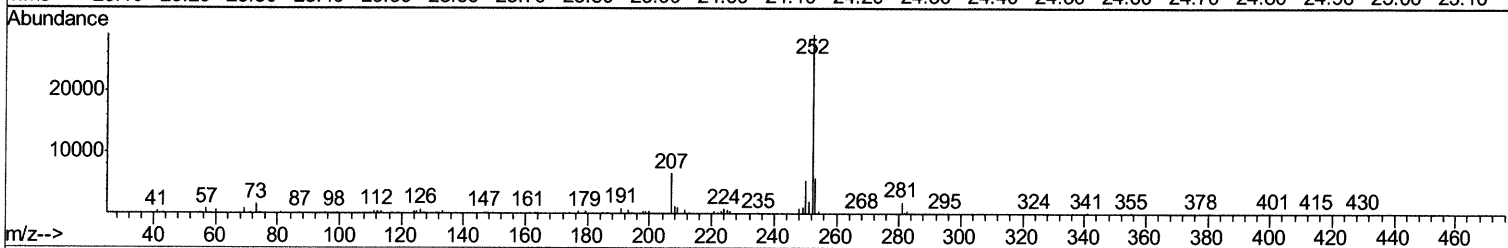
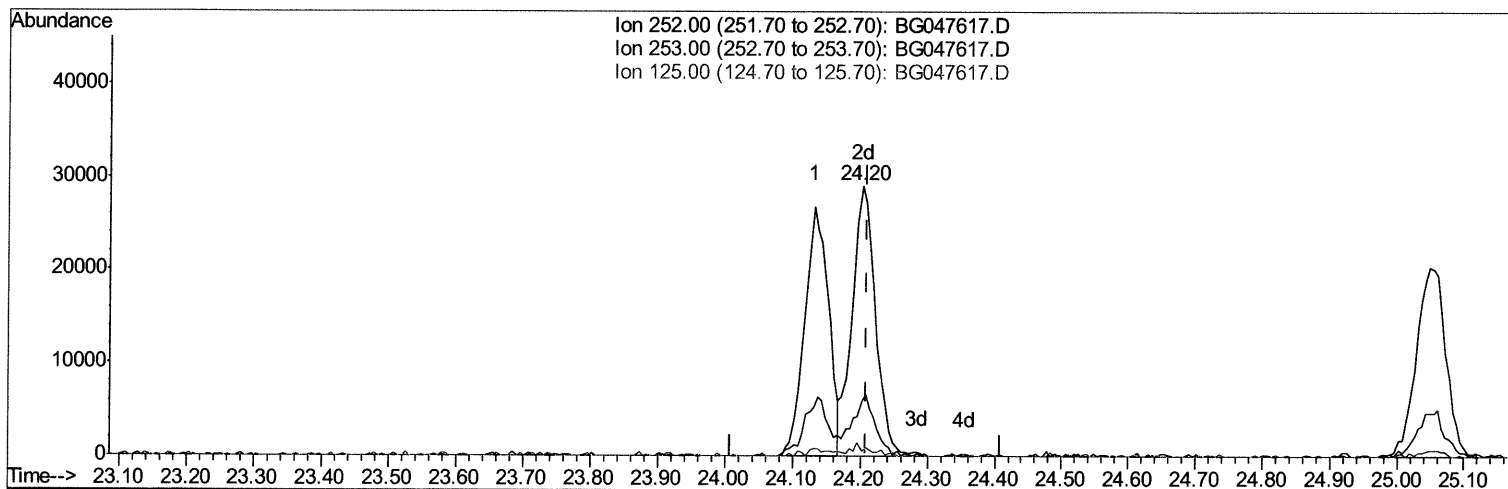
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
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Manual Integrations
 APPROVED

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TIC: BG047617.D

(89) Benzo(k)fluoranthene

24.202min (-0.007) 5.30ng/ul m 12/04/2020

response 70101

Ion	Exp%	Act%
252.00	100	100
253.00	21.10	20.44
125.00	2.80	1.90#
0.00	0.00	0.00

Quantitation Report (Qedit)

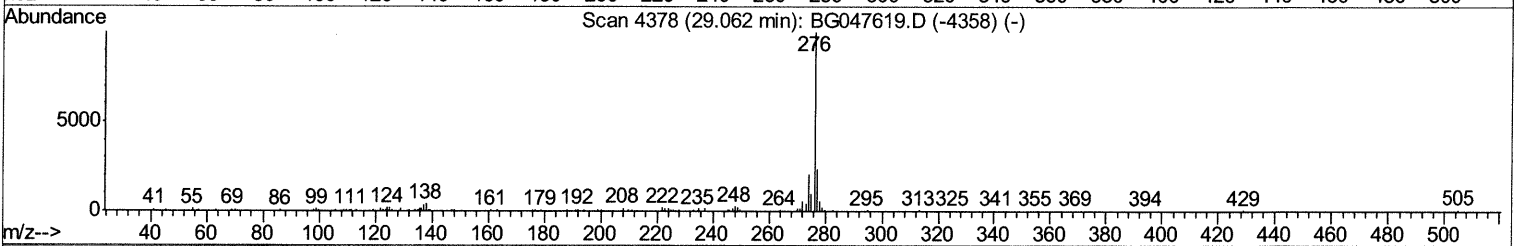
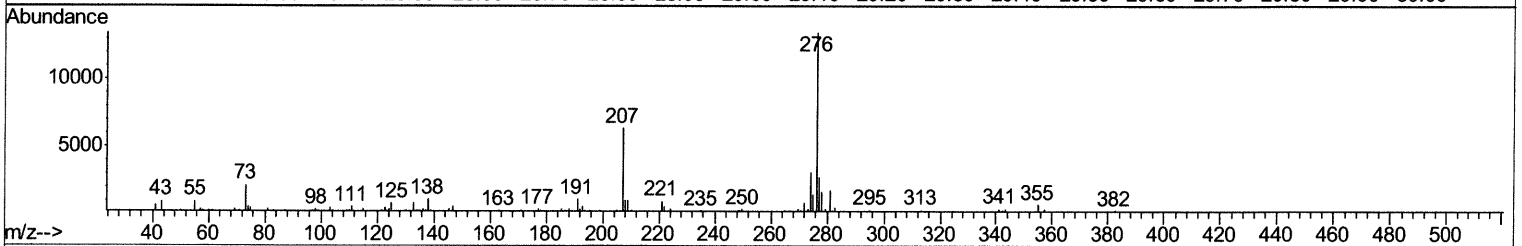
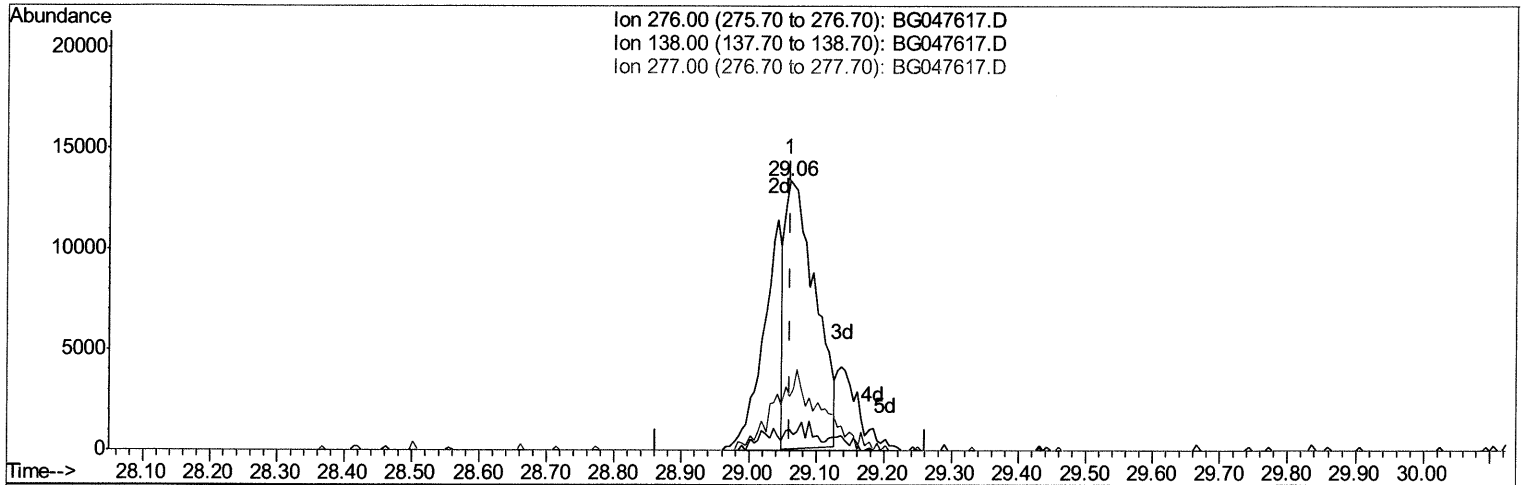
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
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Instrument :
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Manual Integrations
APPROVED

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TIC: BG047617.D

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

29.061min (-0.001) 2.79ng/ul

response 40639

Ion	Exp%	Act%
276.00	100	100
138.00	4.10	7.68#
277.00	23.20	20.19
0.00	0.00	0.00

Quantitation Report (Qedit)

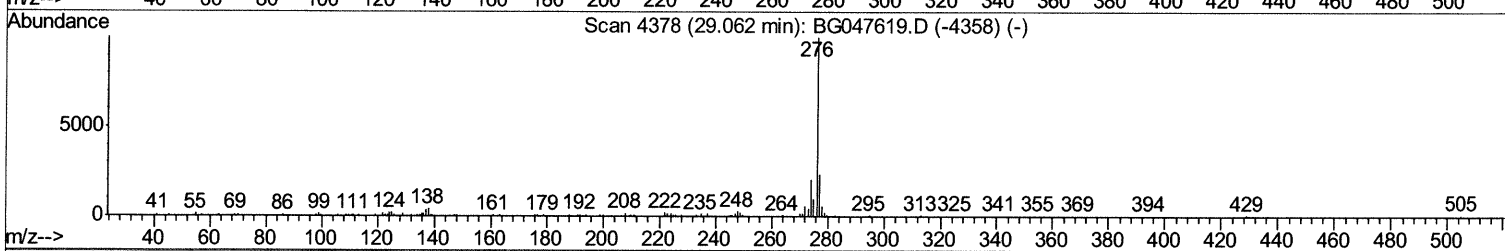
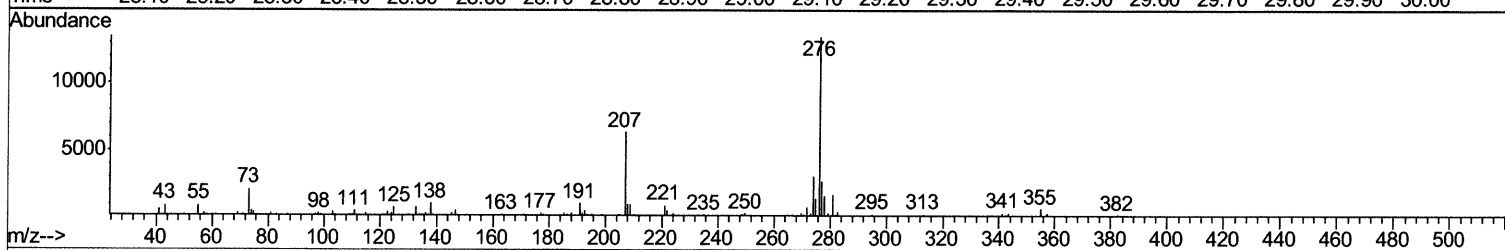
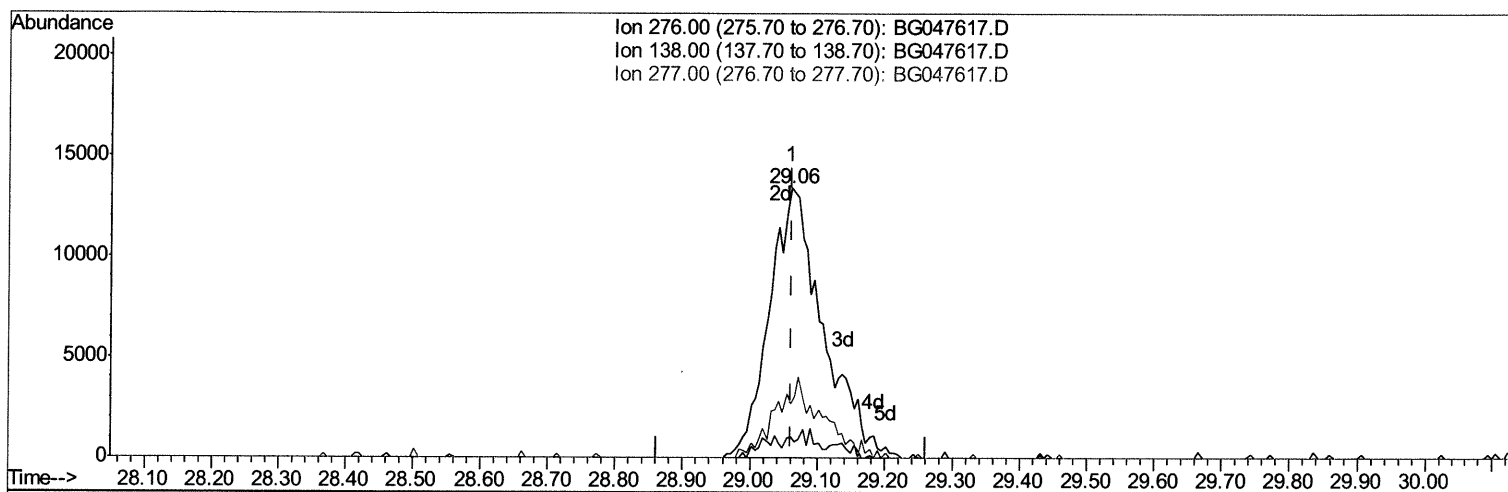
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
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Manual Integrations
APPROVED

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Response via : Initial Calibration



TIC: BG047617.D

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

29.061min (-0.001) 5.09ng/ul m 12/04/2020

response 74014

Ion	Exp%	Act%
276.00	100	100
138.00	4.10	7.68#
277.00	23.20	20.19
0.00	0.00	0.00

Quantitation Report (Qedit)

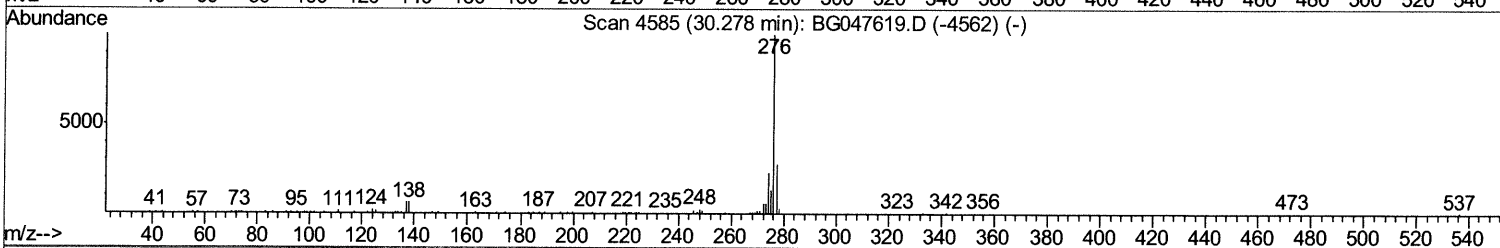
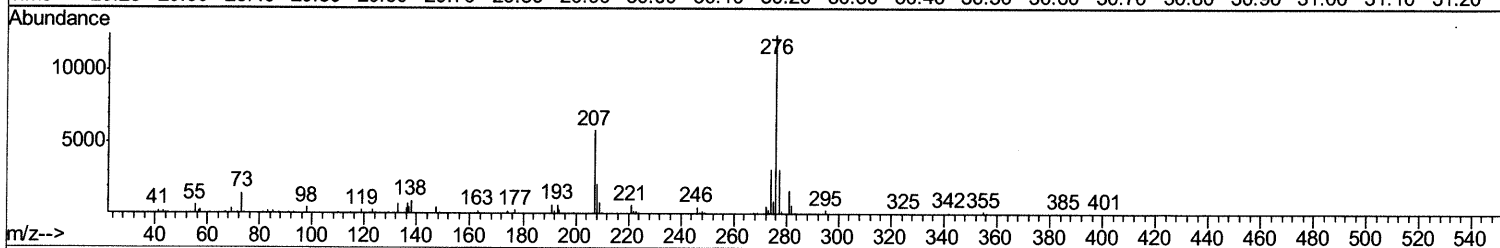
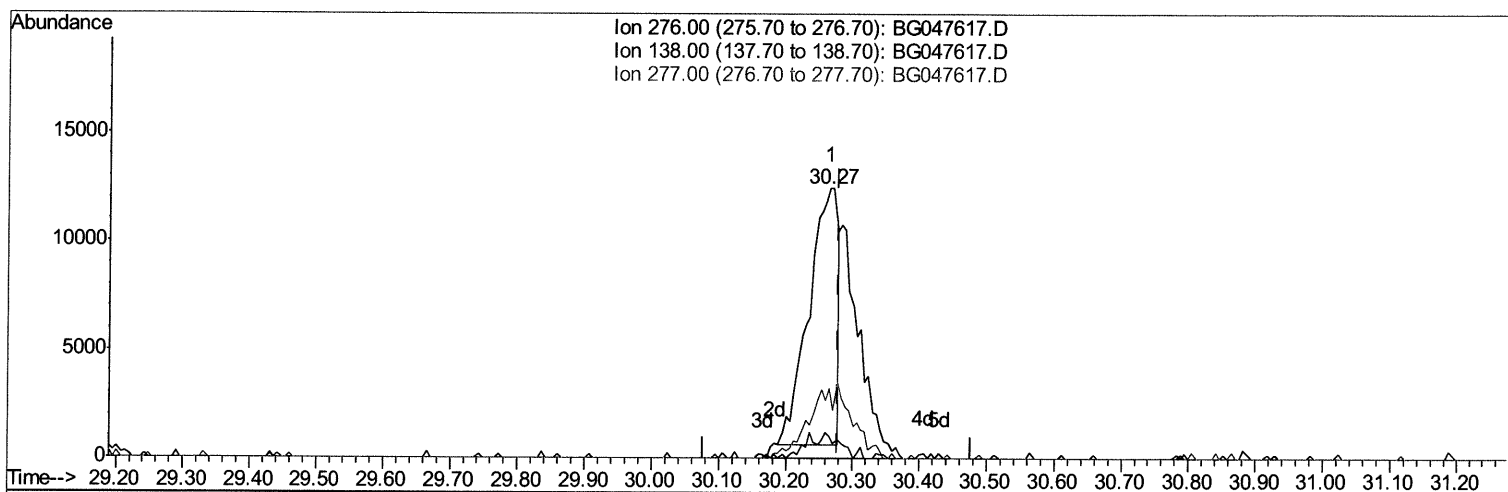
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Manual Integrations
 APPROVED

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TIC: BG047617.D

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

30.265min (-0.013) 2.94ng/ul

response 35292

Ion	Exp%	Act%
276.00	100	100
138.00	5.90	8.10#
277.00	27.40	25.85
0.00	0.00	0.00

Quantitation Report (Qedit)

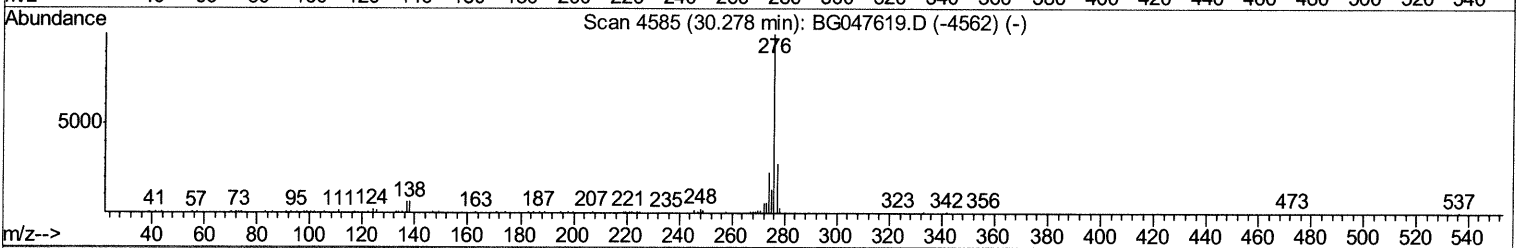
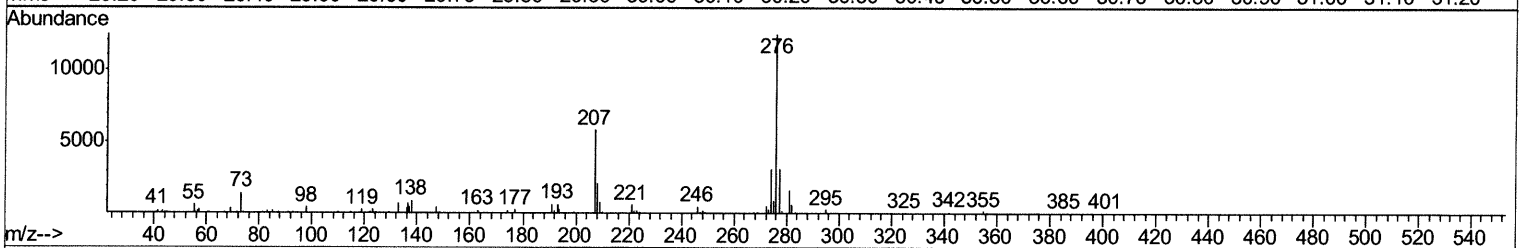
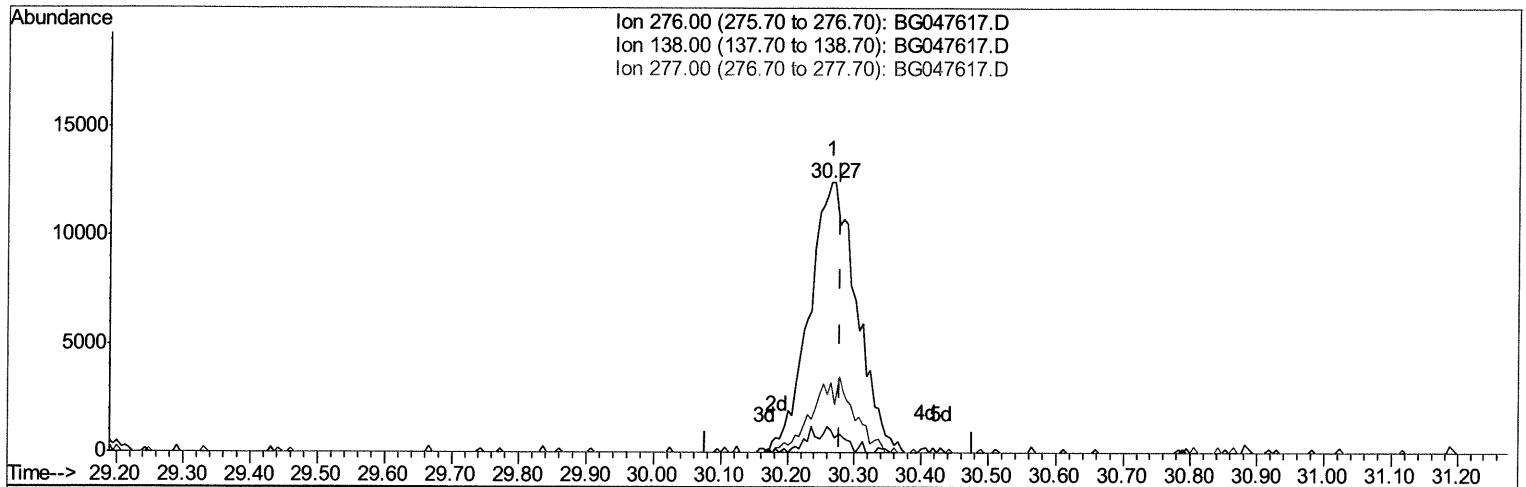
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120720\
 Data File : BG047617.D
 Acq On : 7 Dec 2020 14:23
 Operator : CG/JU
 Sample : SST00525
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
ClientSampleId :
 SST00525

Manual Integrations
APPROVED

mohammad
 12/9/2020 12:18:23 PM

Quant Time: Dec 07 17:16:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG120720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 07 16:26:47 2020
 Response via : Initial Calibration



TIC: BG047617.D

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

30.265min (-0.013) 5.12ng/ul m 12/09/2020

response 61411

Ion	Exp%	Act%
276.00	100	100
138.00	5.90	8.10#
277.00	27.40	25.85
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	8.21	152	22508	20.00	ng/ul	0.00
18) Naphthalene-d8	11.04	136	82234	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.84	164	68285	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.57	188	174553	20.00	ng/ul	0.00
78) Chrysene-d12	21.85	240	188024	20.00	ng/ul	0.00
86) Perylene-d12	25.21	264	189930	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.54	96	989m >	1.60	ng/uL >	0.00	12/09/20 34
5) Phenol-d5	0.00	99	0d	0.00	ng/ul		
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0d	0.00	ng/ul		
9) 2-Chlorophenol-d4	7.74	132	7259	4.79	ng/ul	0.00	
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul		
19) Nitrobenzene-d5	9.38	128	3334	4.91	ng/ul	0.00	
22) 2-Nitrophenol-d4	10.11	143	4003	5.40	ng/ul	0.00	
26) 2,4-Dichlorophenol-d3	10.65	165	8925	5.26	ng/ul	0.00	
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul		
44) Dimethylphthalate-d6	14.23	166	32228	5.50	ng/ul	0.00	
47) Acenaphthylene-d8	14.54	160	35658	5.26	ng/ul	0.00	
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul		
58) Fluorene-d10	15.82	176	28323	5.27	ng/ul	0.00	
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul		
71) Anthracene-d10	17.67	188	49692m >	5.74	ng/ul >	0.00	12/09/20 34
79) Pyrene-d10	19.94	212	58239	5.32	ng/ul	0.00	
90) Benzo(a)pyrene-d12	24.98	264	58630	5.36	ng/ul	0.00	

Target Compounds

					Ovalue	
2) 1,4-Dioxane	3.57	88	705	1.103	ng/uL#	74
10) 2-Chlorophenol	7.77	128	6925	4.533	ng/ul#	77
15) N-Nitroso-di-n-propylamine	9.02	70	6855	4.253	ng/ul#	74
17) Hexachloroethane	9.31	117	3699	5.152	ng/ul	80
20) Nitrobenzene	9.43	77	11808	5.375	ng/ul	95
21) Isophorone	9.95	82	19108	4.564	ng/ul#	89
23) 2-Nitrophenol	10.14	139	3734	4.634	ng/ul#	81
24) 2,4-Dimethylphenol	10.19	107	10536	5.151	ng/ul#	85
25) Bis(2-Chloroethoxy)methane	10.43	93	8698	4.319	ng/ul#	77
27) 2,4-Dichlorophenol	10.68	162	6954	4.616	ng/ul	98
28) Naphthalene	11.09	128	24291	5.215	ng/ul	95
31) Hexachlorobutadiene	11.38	225	9479	6.688	ng/ul#	83
33) 4-Chloro-3-methylphenol	12.29	107	10113	5.431	ng/ul	94
34) 2-Methylnaphthalene	12.68	142	19217	5.568	ng/ul	87
35) 1-Methylnaphthalene	12.90	142	21930	5.406	ng/uL#	99
37) 1,2,4,5-Tetrachlorobenzene	13.04	216	13445	5.011	ng/ul	91
39) 2,4,6-Trichlorophenol	13.27	196	9010	5.407	ng/ul	94
40) 2,4,5-Trichlorophenol	13.35	196	9123	5.273	ng/ul	89
41) 1,1'-Biphenyl	13.67	154	24995	4.719	ng/ul	95
42) 2-Chloronaphthalene	13.72	162	21602	5.145	ng/ul	92
43) 2-Nitroaniline	13.91	65	8040m >	5.265	ng/ul >	12/09/20 34
45) Dimethylphthalate	14.28	163	31173	5.213	ng/ul#	96
46) 2,6-Dinitrotoluene	14.40	165	7003	6.116	ng/ul#	81

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48) Acenaphthylene	14.56	152	33228	5.277	ng/ul	97
50) Acenaphthene	14.90	153	23065	5.222	ng/ul	92
54) Dibenzofuran	15.23	168	34475	5.314	ng/ul	87
55) 2,4-Dinitrotoluene	15.18	165	9208	5.668	ng/ul#	94
56) 2,3,4,6-Tetrachlorophenol	15.45	232	8139	4.913	ng/ul#	92
57) Diethylphthalate	15.63	149	31919	5.441	ng/ul	97
59) Fluorene	15.88	166	29425	5.384	ng/ul#	87
60) 4-Chlorophenyl-phenylether	15.86	204	17950	5.468	ng/ul#	94
65) N-Nitrosodiphenylamine	16.08	169	25415	5.101	ng/ul	93
66) 4-Bromophenyl-phenylether	16.76	248	11465	5.170	ng/ul	88
67) Hexachlorobenzene	16.87	284	12982	5.597	ng/ul	98
70) Phenanthrene	17.62	178	50112	5.316	ng/ul	96
72) Anthracene	17.70	178	51428	5.418	ng/ul	96
73) 1,2,3,4-Tetrachlorobenzene	13.64	216	13525	4.929	ng/uL	93
74) Pentachlorobenzene	15.15	250	14639	5.549	ng/uL	85
76) Di-n-butylphthalate	18.51	149	53985	5.535	ng/ul#	95
80) Pyrene	19.97	202	73714	5.535	ng/ul	99
81) Butylbenzylphthalate	20.84	149	23606	5.046	ng/ul	92
83) Benzo(a)anthracene	21.83	228	72099m>	5.389	ng/ul>	12/09/20 Ju
84) Bis(2-ethylhexyl)phthalate	21.72	149	33540	5.144	ng/ul	97
85) Chrysene	21.90	228	67615	5.311	ng/ul	97
88) Benzo(b)fluoranthene	24.13	252	65712	4.900	ng/ul	93
89) Benzo(k)fluoranthene	24.20	252	70101m>	5.305	ng/ul>	12/09/20 Ju
91) Benzo(a)pyrene	25.05	252	63048	5.274	ng/ul	97
92) Indeno(1,2,3-cd)pyrene	29.06	276	74014m>	5.090	ng/ul>	12/09/20 Ju
93) Dibenzo(a,h)anthracene	29.13	278	63091	5.292	ng/ul#	95
94) Benzo(a,h,i)perylene	30.27	276	61411m>	5.118	ng/ul>	12/09/20 Ju

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