

Quantitation Report (Qedit)

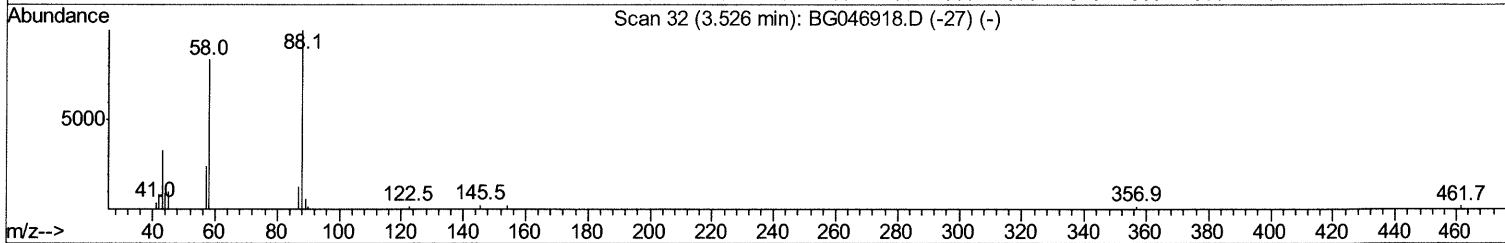
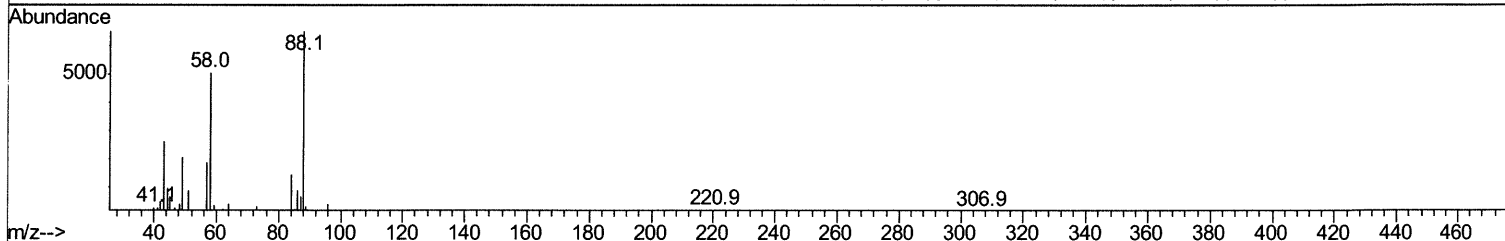
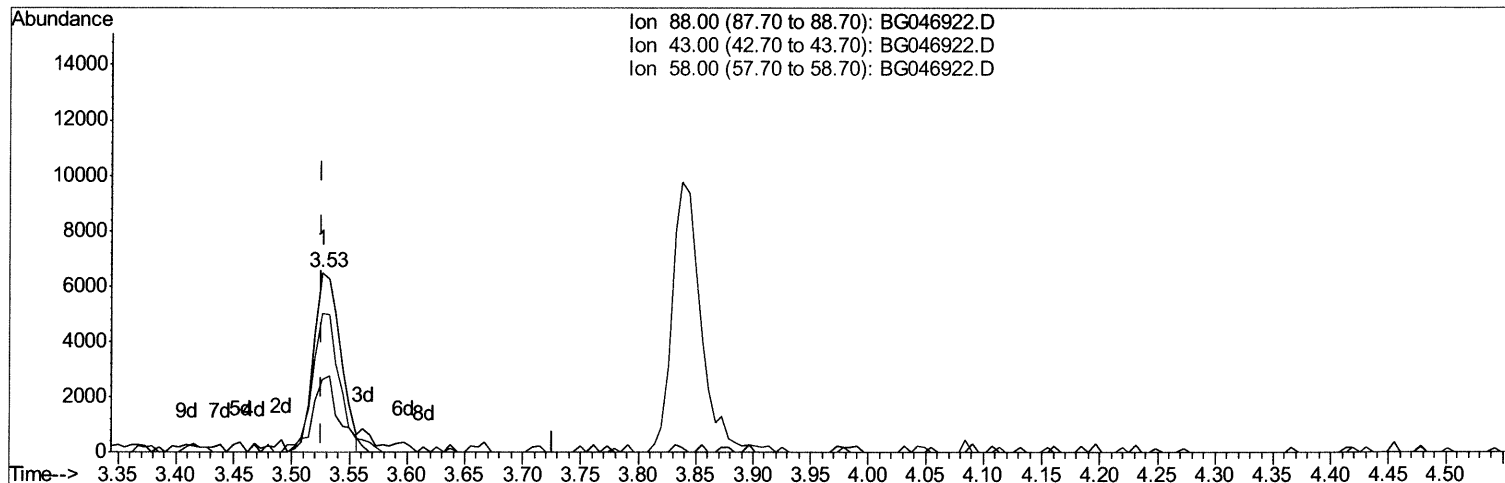
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101520\
 Data File : BG046922.D
 Acq On : 15 Oct 2020 21:09
 Operator : CG/JU
 Sample : SSTDICV020
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SICV31

Manual Integrations
 APPROVED

mohammad
 10/16/2020 12:47:11 PM

Quant Time: Oct 16 02:03:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 16 02:00:57 2020
 Response via : Initial Calibration



TIC: BG046922.D

(2) 1,4-Dioxane

3.526min (+0.000) 7.76ng/uL

response 10464

Ion	Exp%	Act%
88.00	100	100
43.00	33.50	39.99
58.00	83.50	77.35
0.00	0.00	0.00

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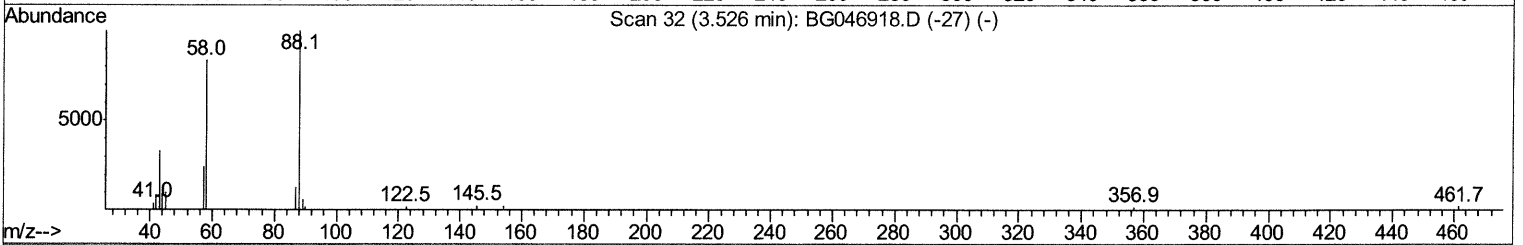
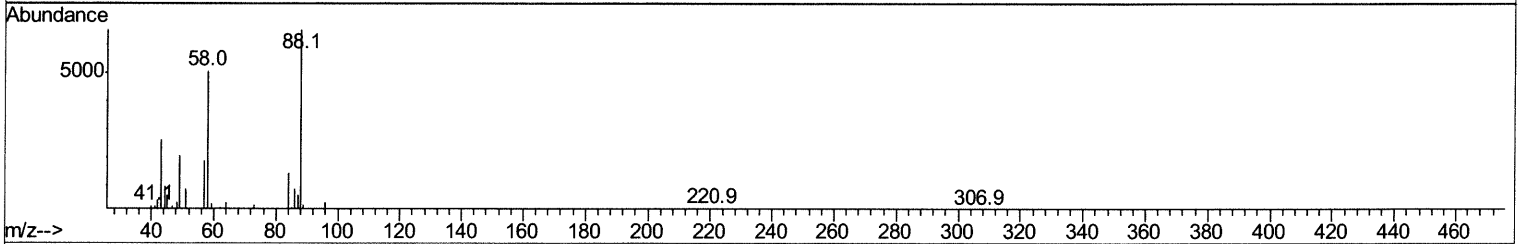
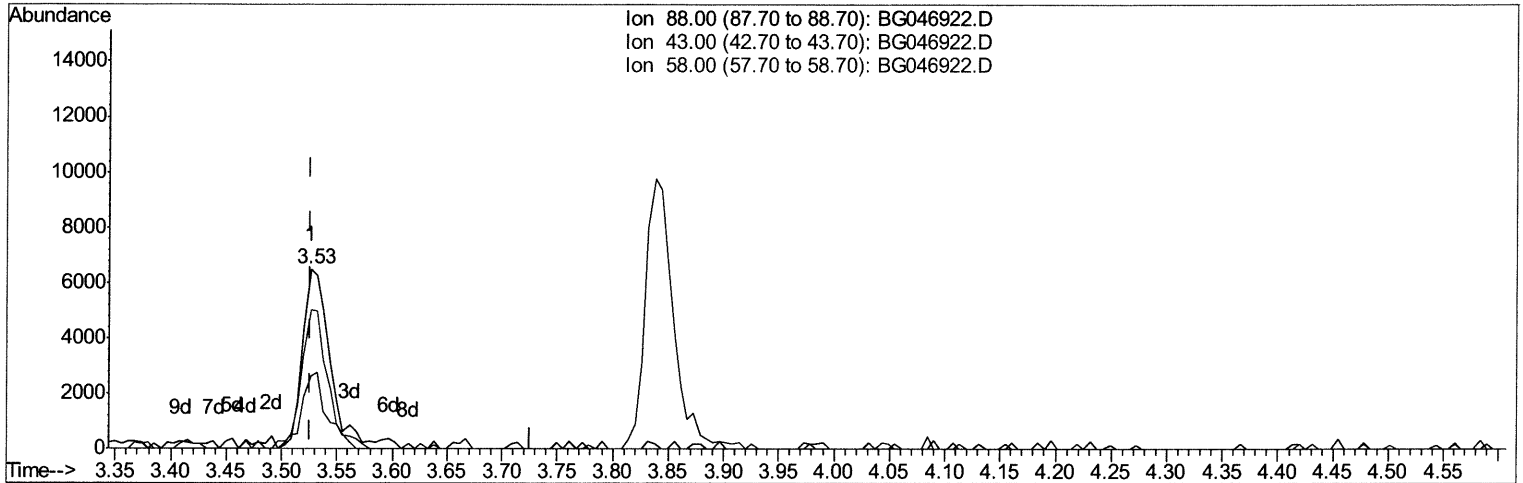
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(2) 1,4-Dioxane

3.526min (+0.000) 8.56ng/uL m 10/16/20 JU

response 11552

Ion	Exp%	Act%
88.00	100	100
43.00	33.50	39.99
58.00	83.50	77.35
0.00	0.00	0.00

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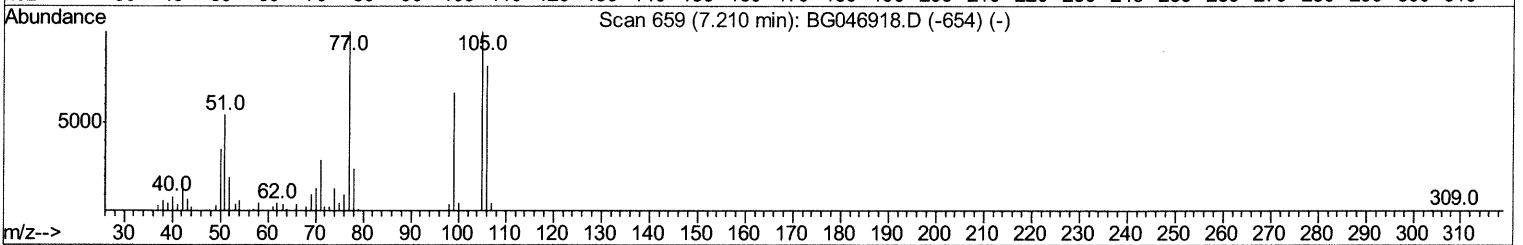
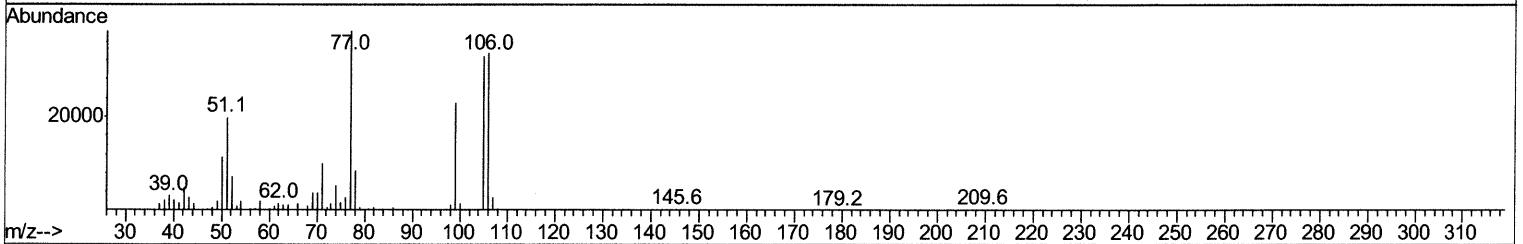
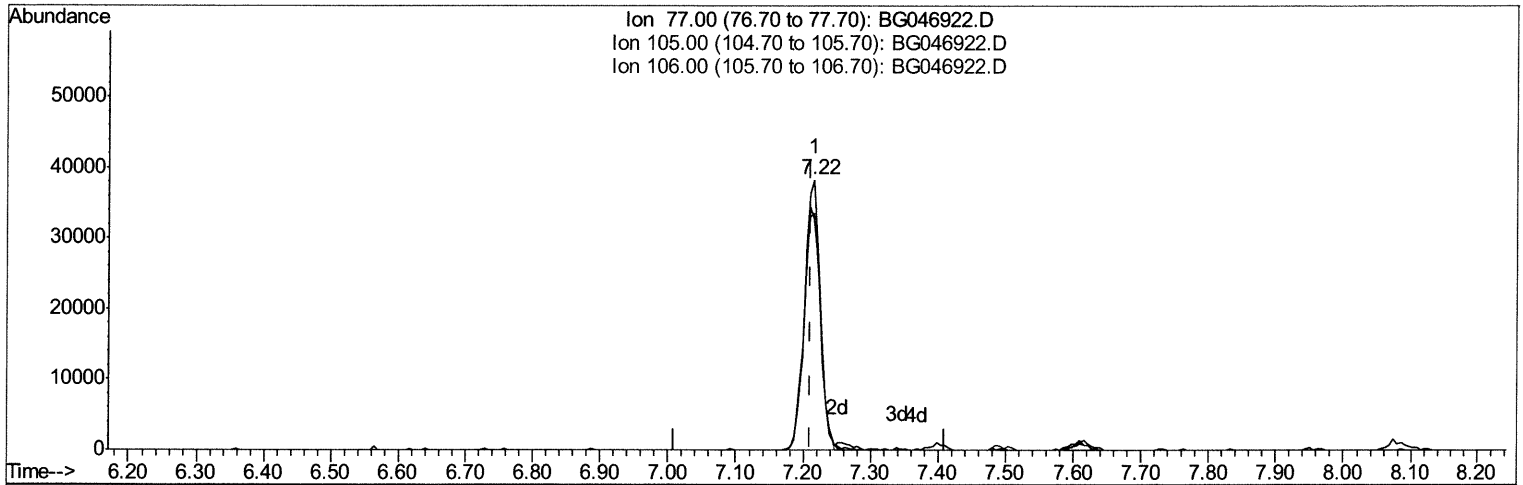
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(4) Benzaldehyde

7.216min (+0.006) 24.04ng/ul

response 61966

Ion	Exp%	Act%
77.00	100	100
105.00	100.00	85.47
106.00	82.20	87.84
0.00	0.00	0.00

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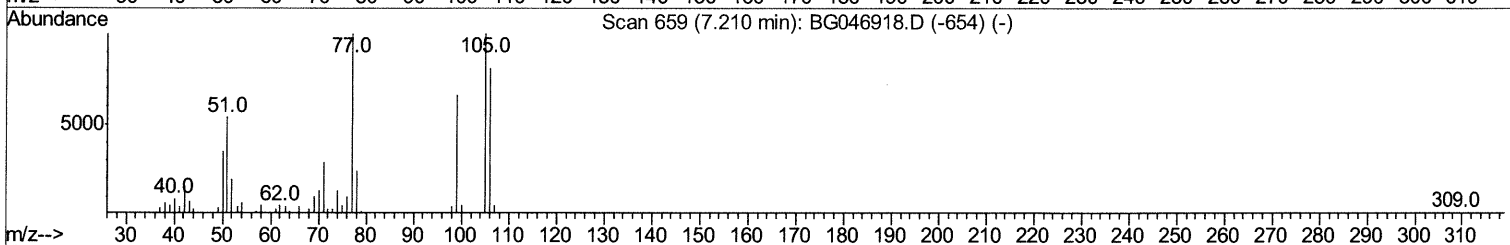
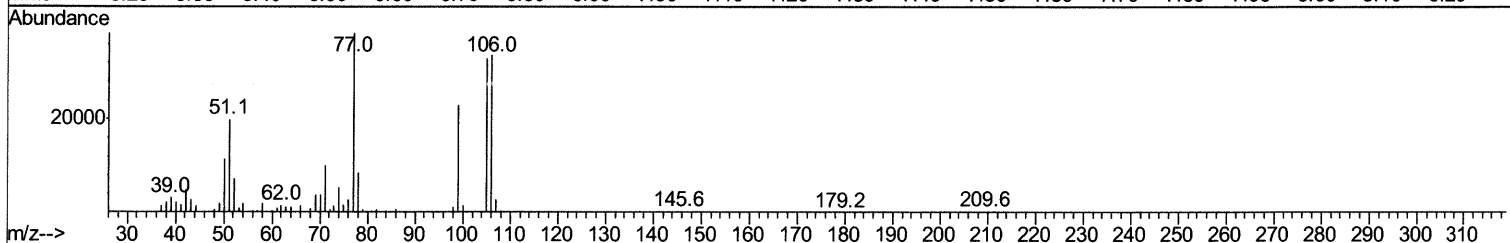
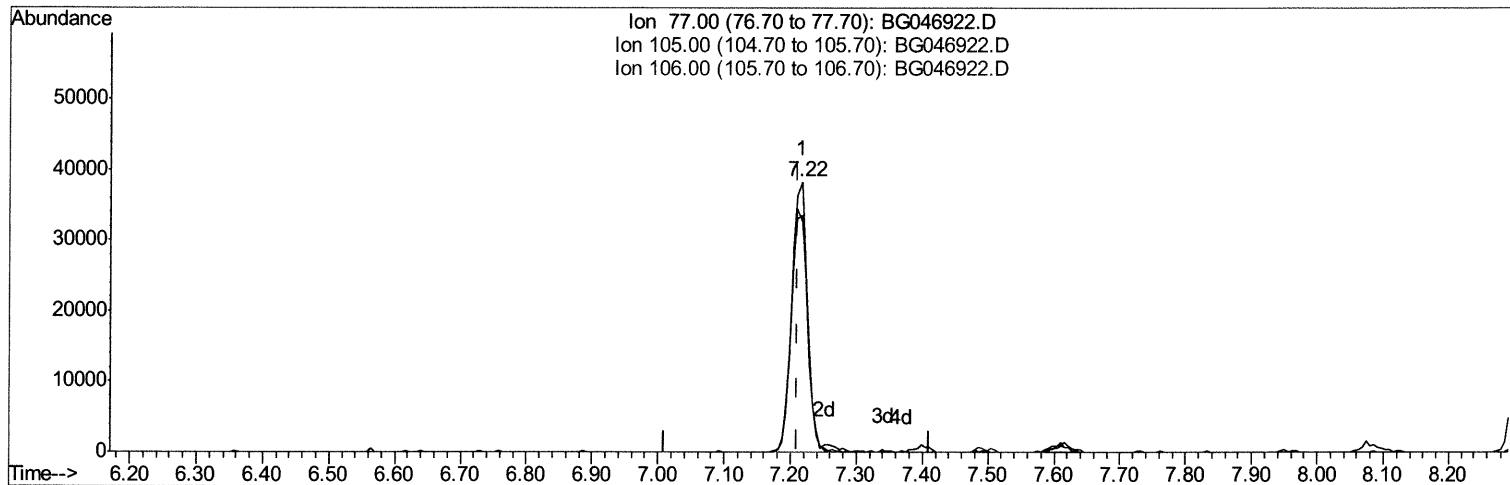
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(4) Benzaldehyde

7.216min (+0.006) 24.69ng/ul m 10/16/20 JU

response 63643

Ion	Exp%	Act%
77.00	100	100
105.00	100.00	85.47
106.00	82.20	87.84
0.00	0.00	0.00

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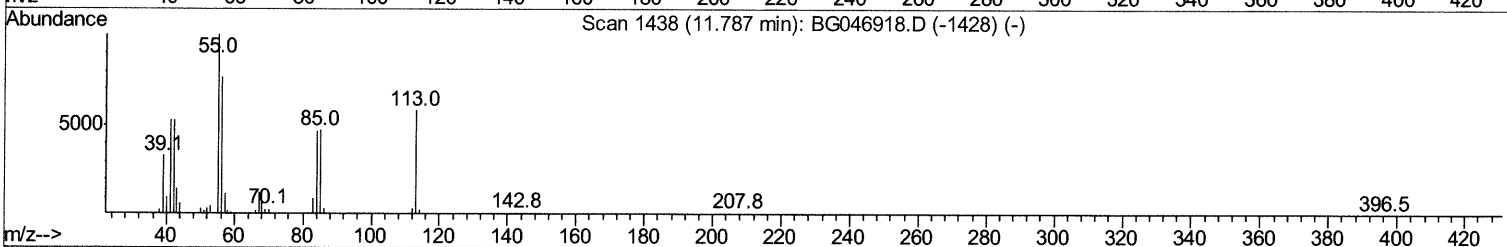
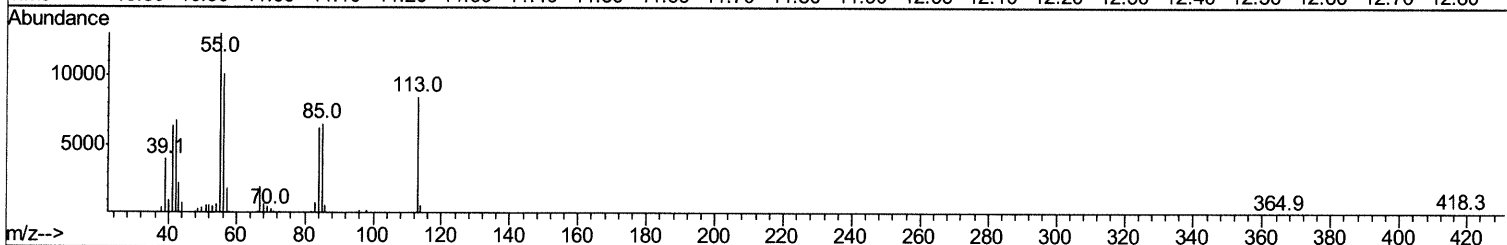
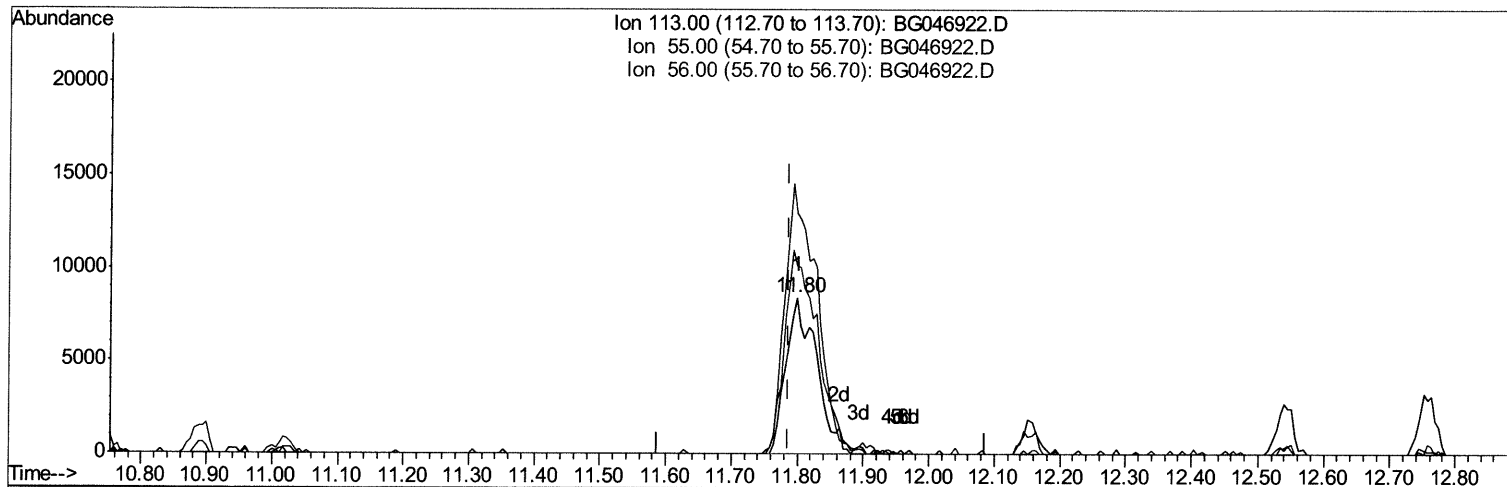
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(32) Caprolactam

11.799min (+0.012) 20.69ng/ul m 10/16/20 JU

response 27069

Ion	Exp%	Act%
113.00	100	100
55.00	172.40	154.25
56.00	131.30	120.23
0.00	0.00	0.00

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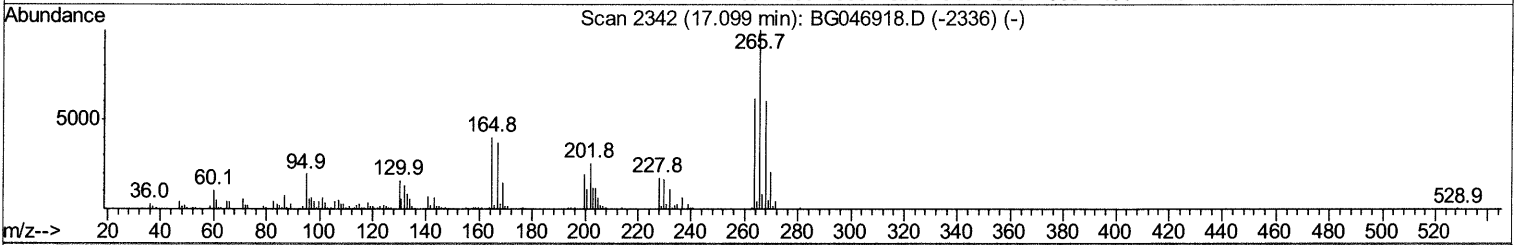
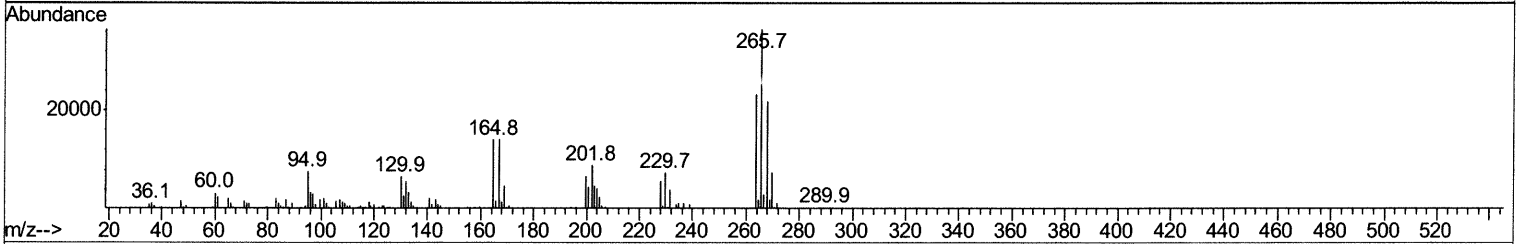
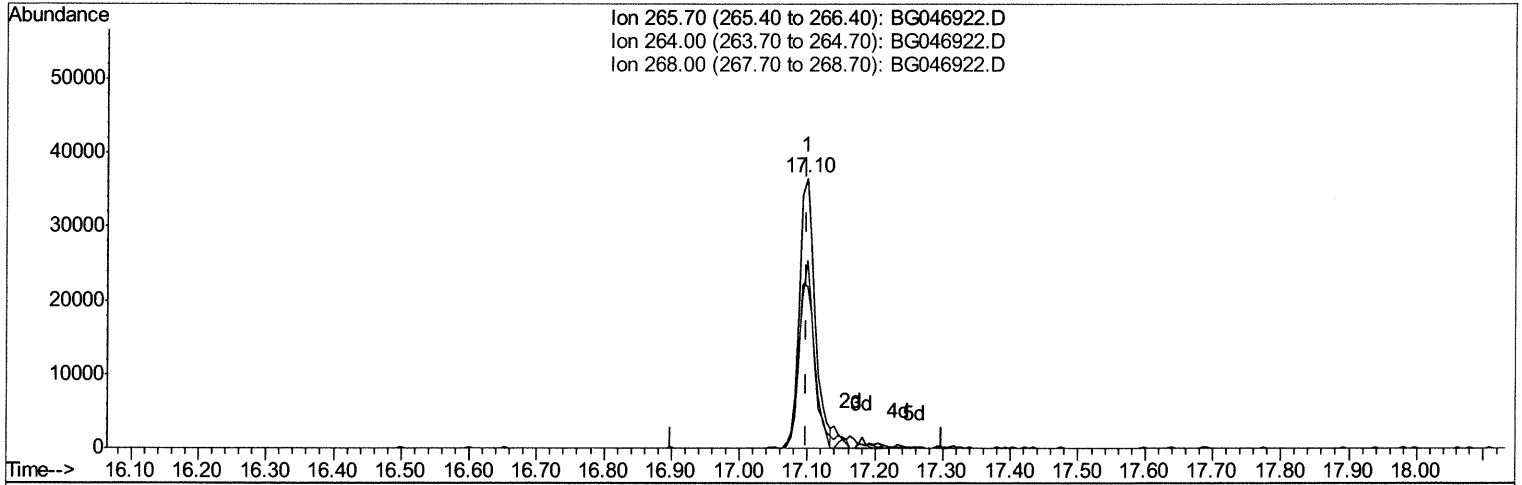
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TIC: BG046922.D

(69) Pentachlorophenol (C)

17.098min (-0.000) 17.51ng/ul

response 58550

Ion	Exp%	Act%
265.70	100	100
264.00	65.70	69.42
268.00	65.20	60.06
0.00	0.00	0.00

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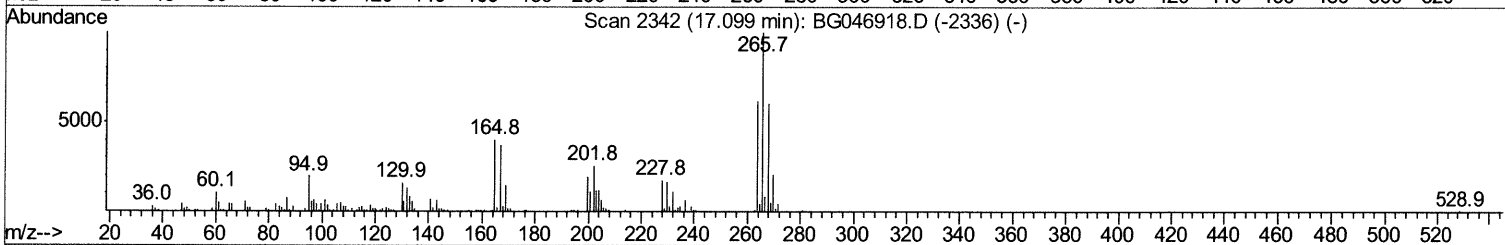
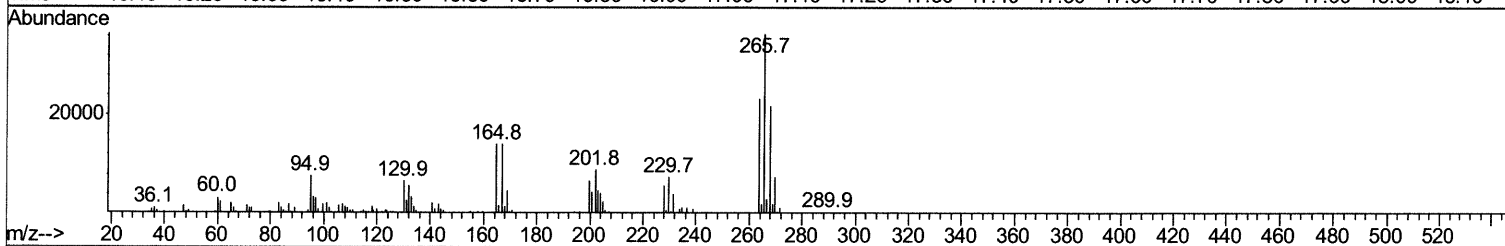
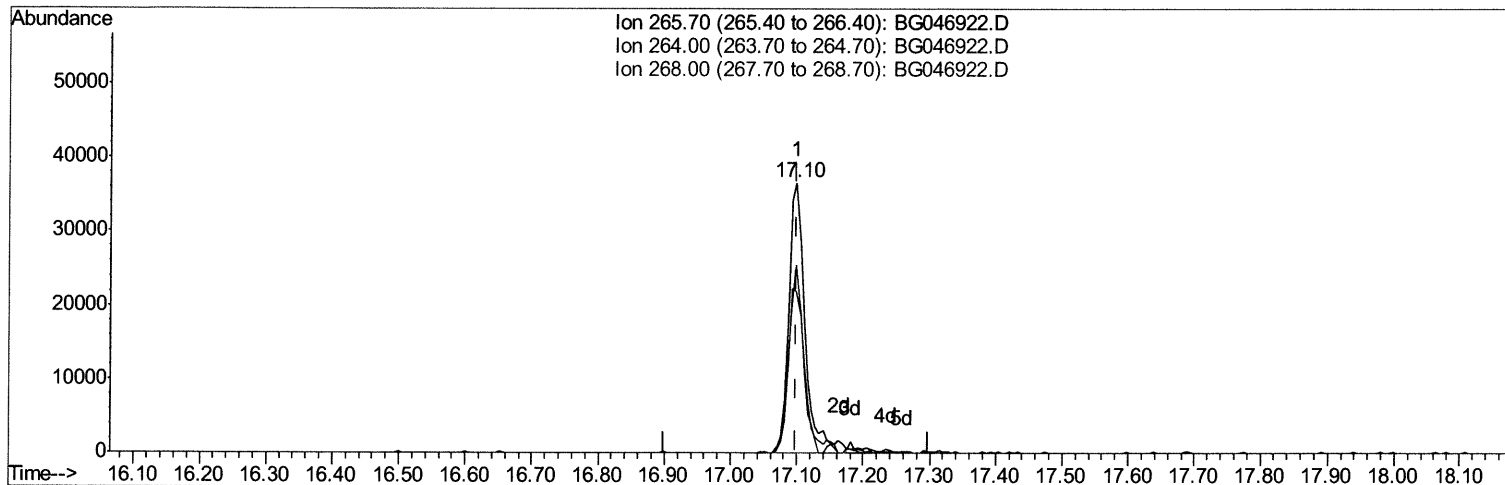
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TIC: BG046922.D

(69) Pentachlorophenol (C)

17.098min (-0.000) 18.67ng/ul m 10/16/2020

response 62415

Ion	Exp%	Act%
265.70	100	100
264.00	65.70	63.87
268.00	65.20	60.06
0.00	0.00	0.00

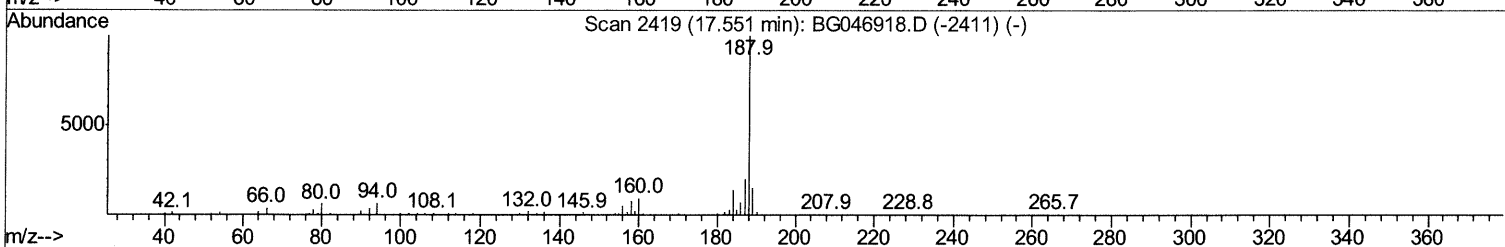
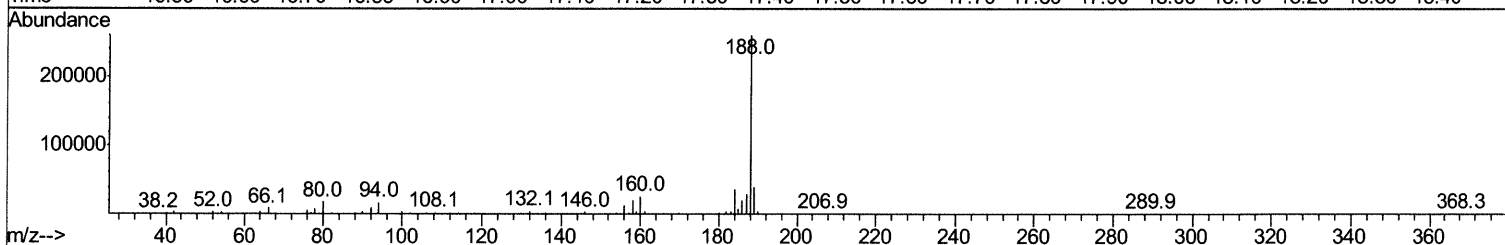
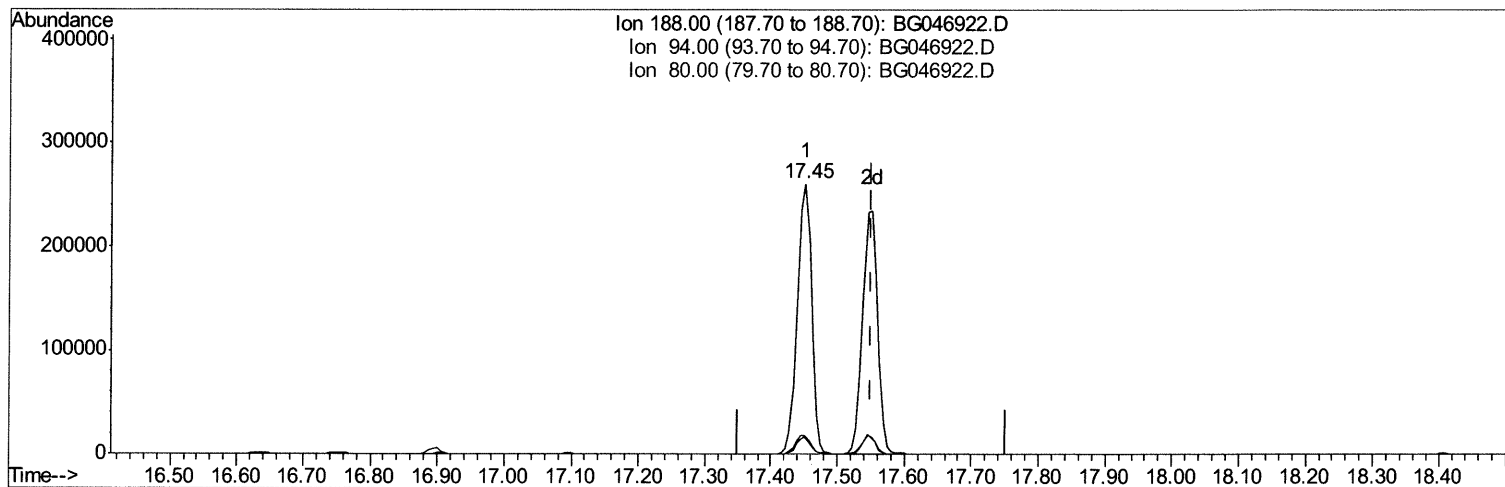
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TIC: BG046922.D

(71) Anthracene-d10 (S)

17.451min (-0.100) 20.65ng/ul

response 387712

Ion	Exp%	Act%
188.00	100	100
94.00	6.50	6.49
80.00	6.00	6.86
0.00	0.00	0.00

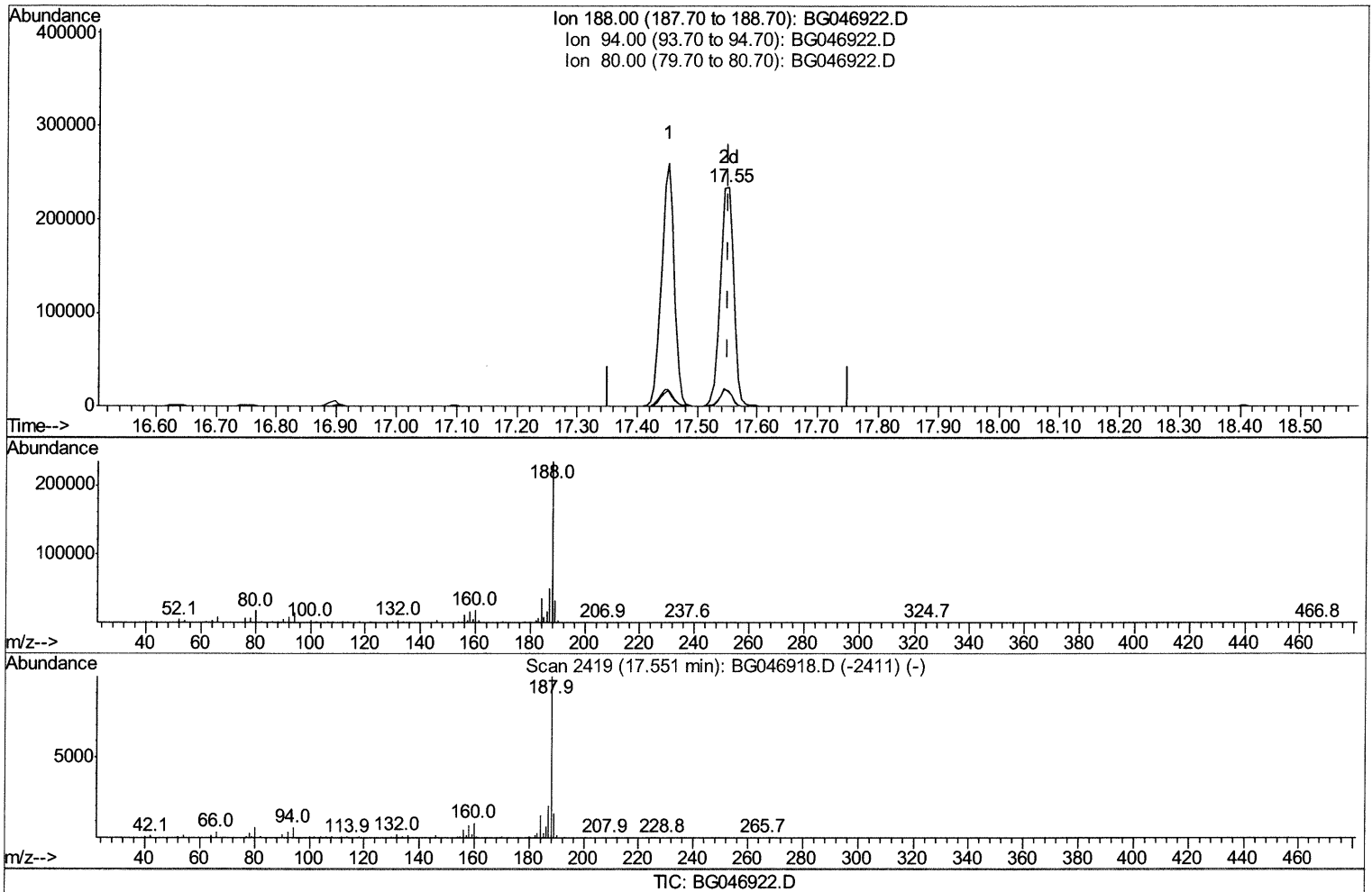
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Manual Integrations
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(71) Anthracene-d10 (S)

17.551min (-0.000) 19.33ng/ul m 10/16/2020 JU

response 362983

Ion	Exp%	Act%
188.00	100	100
94.00	6.50	6.48
80.00	6.00	7.32#
0.00	0.00	0.00

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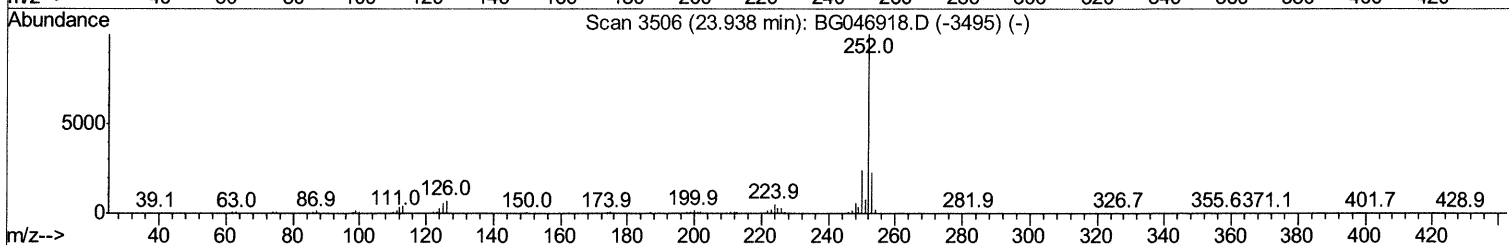
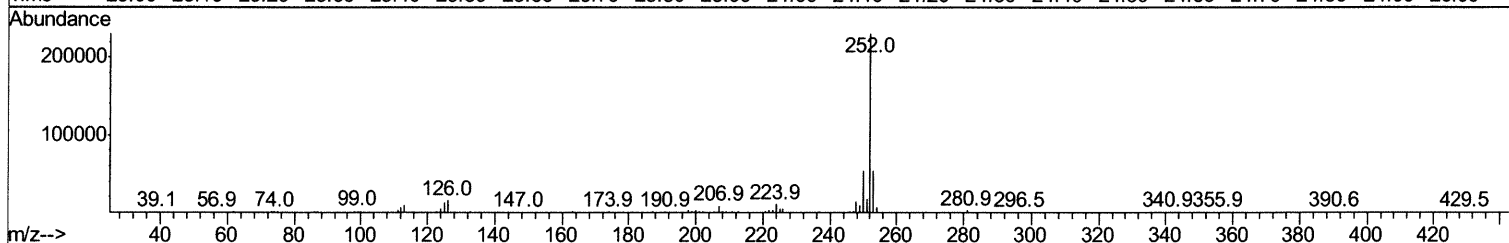
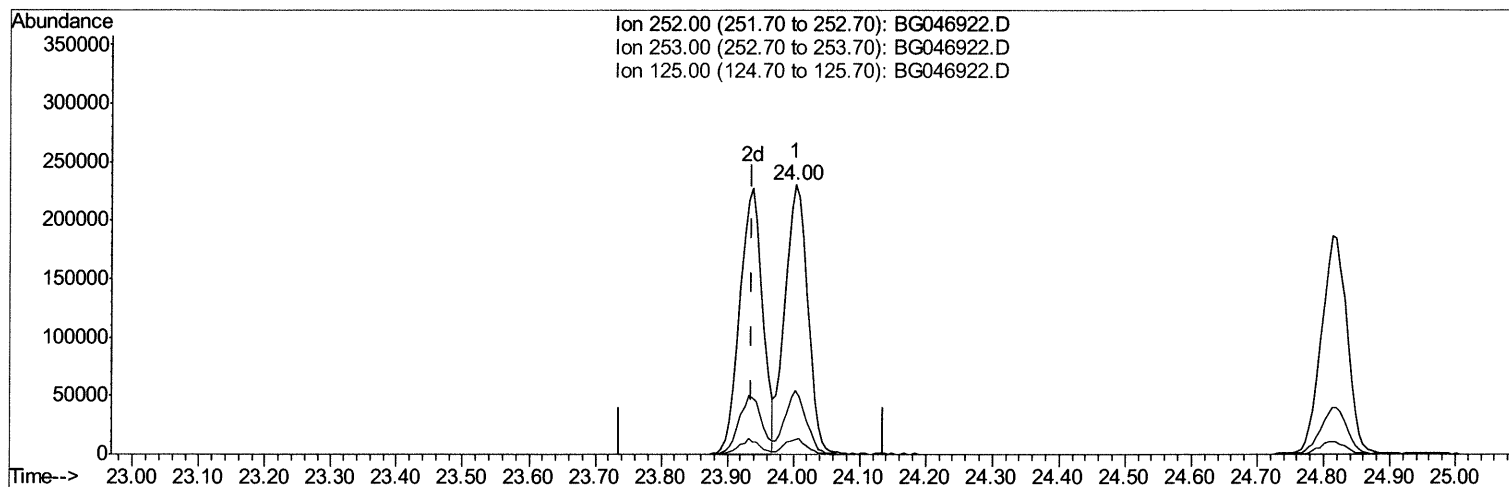
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TIC: BG046922.D

(88) Benzo(b)fluoranthene

24.002min (+0.065) 18.50ng/ul

response 547981

Ion	Exp%	Act%
252.00	100	100
253.00	22.70	23.43
125.00	5.80	5.31
0.00	0.00	0.00

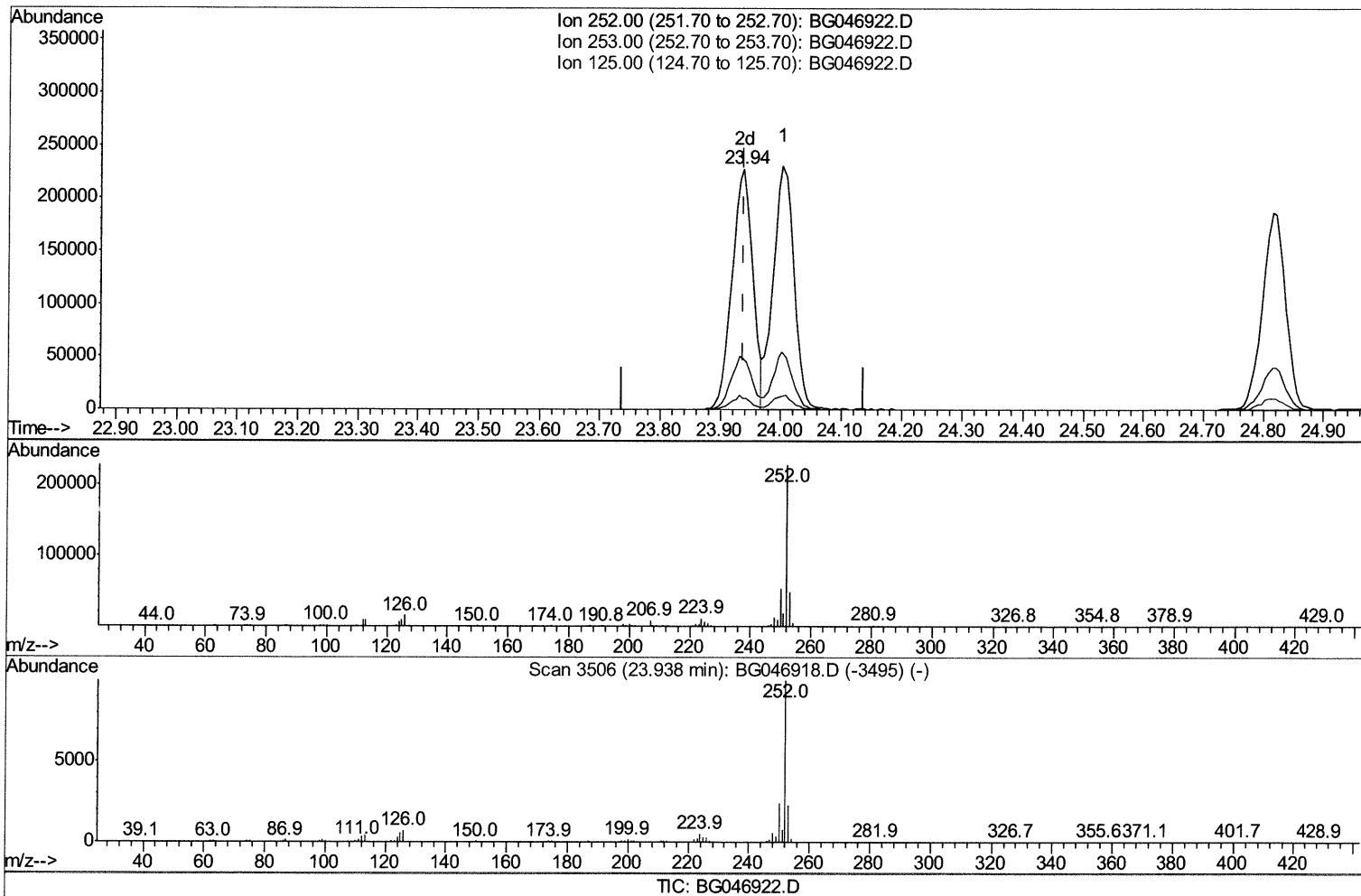
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(88) Benzo(b)fluoranthene

23.938min (-0.000) 18.41ng/ul m 10/16/2020

response 545278

Ion	Exp%	Act%
252.00	100	100
253.00	22.70	21.32
125.00	5.80	4.48#
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	8.08	152	50230	20.00	ng/ul	0.00
18) Naphthalene-d8	10.89	136	204358	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.71	164	153763	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.45	188	387712	20.00	ng/ul	0.00
78) Chrysene-d12	21.72	240	427612	20.00	ng/ul	0.00
86) Perylene-d12	24.97	264	427037	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.50	96	10590	9.36	ng/uL	0.00
5) Phenol-d5	7.23	99	88640	20.18	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.40	67	55029	21.44	ng/ul	0.00
9) 2-Chlorophenol-d4	7.61	132	69356	20.98	ng/ul	0.00
13) 4-Methylphenol-d8	8.78	113	70067	19.80	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	34624	19.93	ng/ul	0.00
22) 2-Nitrophenol-d4	9.97	143	38442	19.40	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.51	165	79140	20.36	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	95356	21.34	ng/ul	0.00
44) Dimethylphthalate-d6	14.11	166	252236	19.75	ng/ul	0.00
47) Acenaphthylene-d8	14.40	160	275056	18.89	ng/ul	0.00
52) 4-Nitrophenol-d4	14.88	143	44206	20.08	ng/ul	0.00
58) Fluorene-d10	15.69	176	217695	18.54	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.81	200	50043	18.74	ng/ul	0.00
71) Anthracene-d10	17.55	188	362983m	19.33	ng/ul	0.00
79) Pyrene-d10	19.83	212	469760	20.04	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.75	264	468346	19.47	ng/ul	0.00

Target Compounds

2) 1,4-Dioxane	3.53	88	11552m	8.562	ng/uL	
4) Benzaldehyde	7.22	77	63643m	24.688	ng/ul	
6) Phenol	7.26	94	89643	19.311	ng/ul	96
8) Bis(2-Chloroethyl)ether	7.49	93	70128	19.935	ng/ul	95
10) 2-Chlorophenol	7.64	128	69129	20.712	ng/ul	98
11) 2-Methylphenol	8.51	108	67490	19.516	ng/ul	96
12) 2,2'-oxybis(1-Chloropropan	8.60	45	99656	19.805	ng/ul#	93
14) Acetophenone	8.90	105	120722	21.269	ng/ul	96
15) N-Nitroso-di-n-propylamine	8.88	70	62452	20.979	ng/ul#	95
16) 4-Methylphenol	8.84	108	72928	19.868	ng/ul	90
17) Hexachloroethane	9.17	117	29808	19.601	ng/ul	97
20) Nitrobenzene	9.28	77	93232	19.893	ng/ul	98
21) Isophorone	9.81	82	168188	19.502	ng/ul	97
23) 2-Nitrophenol	10.00	139	40781	19.642	ng/ul	88
24) 2,4-Dimethylphenol	10.05	107	84141	19.211	ng/ul	96
25) Bis(2-Chloroethoxy)methane	10.29	93	92831	18.876	ng/ul	97
27) 2,4-Dichlorophenol	10.53	162	76792	19.865	ng/ul	97
28) Naphthalene	10.94	128	222536	18.872	ng/ul	97
30) 4-Chloroaniline	11.04	127	95675	22.376	ng/ul	100
31) Hexachlorobutadiene	11.23	225	60902	18.489	ng/ul	99
32) Caprolactam	11.80	113	27069m	20.690	ng/ul	
33) 4-Chloro-3-methylphenol	12.16	107	83387	20.616	ng/ul	94
34) 2-Methylnaphthalene	12.54	142	173351	19.995	ng/ul	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.76	142	166047	19.418	ng/uL	97
37) 1,2,4,5-Tetrachlorobenzene	12.90	216	115244	19.218	ng/uL	99
38) Hexachlorocyclopentadiene	12.89	237	74462	18.699	ng/uL	87
39) 2,4,6-Trichlorophenol	13.14	196	69405	18.807	ng/uL	96
40) 2,4,5-Trichlorophenol	13.21	196	76167	19.551	ng/uL	93
41) 1,1'-Biphenyl	13.54	154	241323	19.457	ng/uL	99
42) 2-Chloronaphthalene	13.58	162	185325	18.650	ng/uL	98
43) 2-Nitroaniline	13.78	65	59077	19.817	ng/uL	86
45) Dimethylphthalate	14.15	163	247015	18.654	ng/uL	100
46) 2,6-Dinitrotoluene	14.27	165	53591	19.798	ng/uL	97
48) Acenaphthylene	14.43	152	285544	19.402	ng/uL	98
49) 3-Nitroaniline	14.60	138	48152	21.461	ng/uL	90
50) Acenaphthene	14.77	153	191564	18.563	ng/uL	98
51) 2,4-Dinitrophenol	14.81	184	30363	17.770	ng/uL	90
53) 4-Nitrophenol	14.90	109	44733	19.360	ng/uL	93
54) Dibenzofuran	15.10	168	299381	19.522	ng/uL	98
55) 2,4-Dinitrotoluene	15.05	165	77457	19.936	ng/uL#	89
56) 2,3,4,6-Tetrachlorophenol	15.33	232	75180	19.482	ng/uL#	92
57) Diethylphthalate	15.51	149	252822	19.141	ng/uL	99
59) Fluorene	15.75	166	237016	18.834	ng/uL	97
60) 4-Chlorophenyl-phenylether	15.74	204	136441	18.603	ng/uL	95
61) 4-Nitroaniline	15.76	138	57051	22.791	ng/uL	87
64) 4,6-Dinitro-2-methylphenol	15.82	198	54339	18.924	ng/uL#	97
65) N-Nitrosodiphenylamine	15.95	169	220557	19.386	ng/uL	91
66) 4-Bromophenyl-phenylether	16.63	248	96010	18.564	ng/uL	98
67) Hexachlorobenzene	16.75	284	93886	17.429	ng/uL	95
68) Atrazine	16.90	200	92438	18.393	ng/uL	96
69) Pentachlorophenol	17.10	266	62415m>	18.668	ng/uL>	10/16/20 JU
70) Phenanthrene	17.49	178	414510	18.684	ng/uL	98
72) Anthracene	17.58	178	426484	19.032	ng/uL	98
73) 1,2,3,4-Tetrachlorobenzene	13.51	216	118424	19.218	ng/uL	95
74) Pentachlorobenzene	15.02	250	116317	18.395	ng/uL	98
75) Carbazole	17.85	167	365364	20.303	ng/uL	99
76) Di-n-butylphthalate	18.40	149	453772	19.270	ng/uL	99
77) Fluoranthene	19.50	202	557412	20.576	ng/uL	98
80) Pyrene	19.86	202	552021	19.075	ng/uL	99
81) Butylbenzylphthalate	20.74	149	203273	19.302	ng/uL	98
82) 3,3'-Dichlorobenzidine	21.62	252	216187	21.613	ng/uL	94
83) Benzo(a)anthracene	21.70	228	564968	19.143	ng/uL	99
84) Bis(2-ethylhexyl)phthalate	21.60	149	282862	18.810	ng/uL	96
85) Chrysene	21.77	228	529381	18.591	ng/uL	99
87) Di-n-octyl phthalate	22.83	149	484157	19.726	ng/uL	100
88) Benzo(b)fluoranthene	23.94	252	545278m>	18.408	ng/uL>	10/16/20 JU
89) Benzo(k)fluoranthene	24.00	252	547981	19.183	ng/uL	98
91) Benzo(a)pyrene	24.81	252	497331	18.675	ng/uL	99
92) Indeno(1,2,3-cd)pyrene	28.67	276	611187	18.734	ng/uL	99
93) Dibenzo(a,h)anthracene	28.74	278	510951	18.886	ng/uL	98
94) Benzo(a,h,i)perylene	29.82	276	512522	18.715	ng/uL	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101520\
Data File : BG046922.D
Acq On : 15 Oct 2020 21:09
Operator : CG/JU
Sample : SSTDICV020
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SICV31

Manual Integrations
APPROVED

mohammad
10/16/2020 12:47:11 PM

Quant Time: Oct 16 02:11:29 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 16 02:00:57 2020
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

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