

(QT Reviewed)

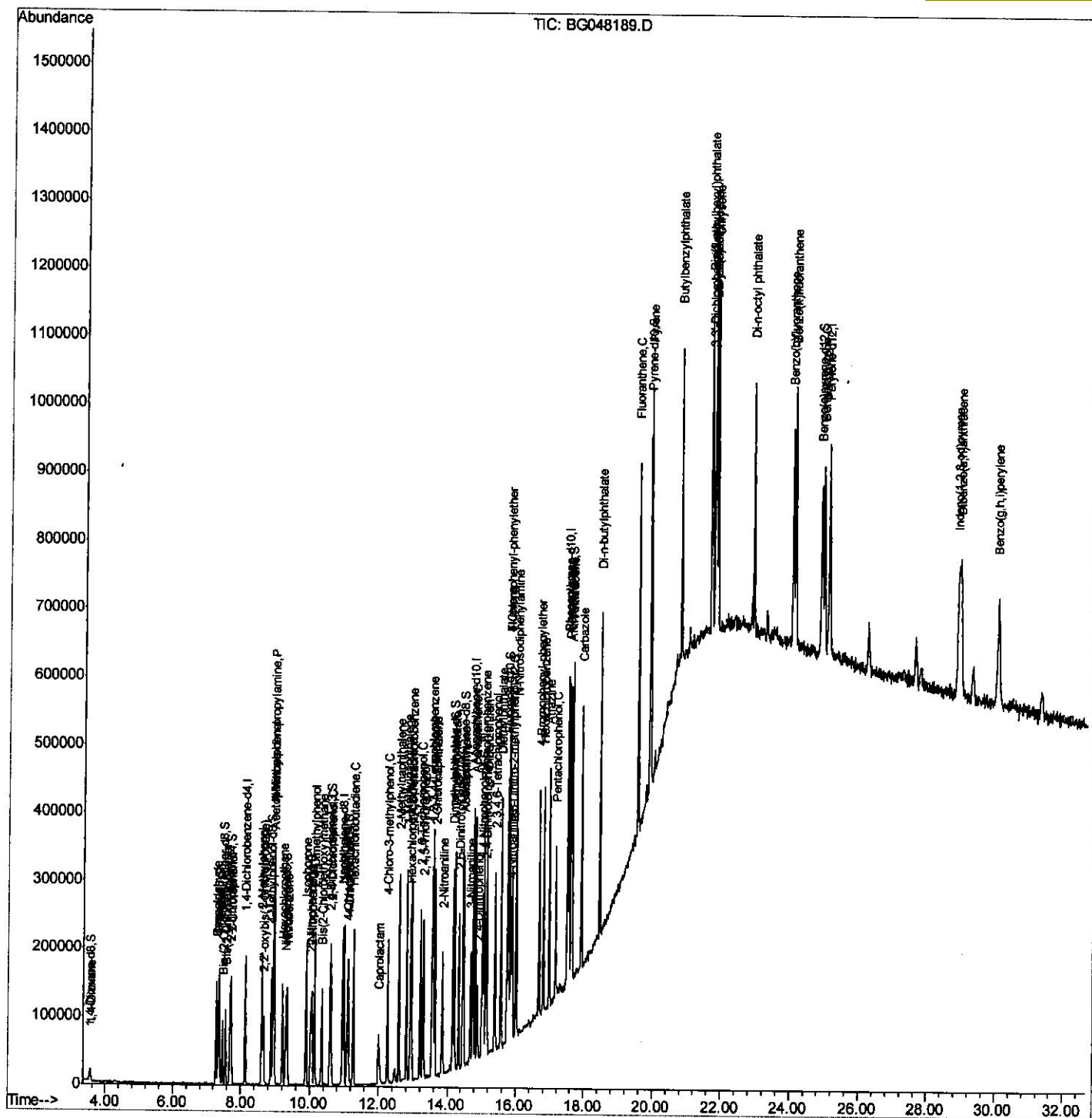
```
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021621\  
Data File : BG048189.D  
Acq On    : 17 Feb 2021   14:38  
Operator  : CG/JU  
Sample    : SSTDCCC020EC  
Misc      :  
ALS Vial  : 2      Sample Multiplier: 1
```

Instrument :
BNA_G
LabSampleId :
SSTD02035

Quant Time: Feb 17 17:37:32 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Feb 15 06:21:07 2021
Response via : Initial Calibration

Manual Integrations APPROVED

mohammad
2/18/2021 1:38:36 PM



Quantitation Report (Qedit)

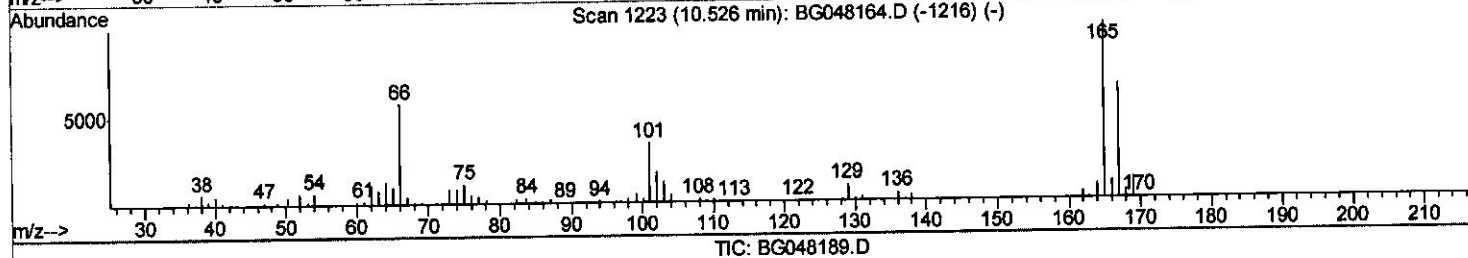
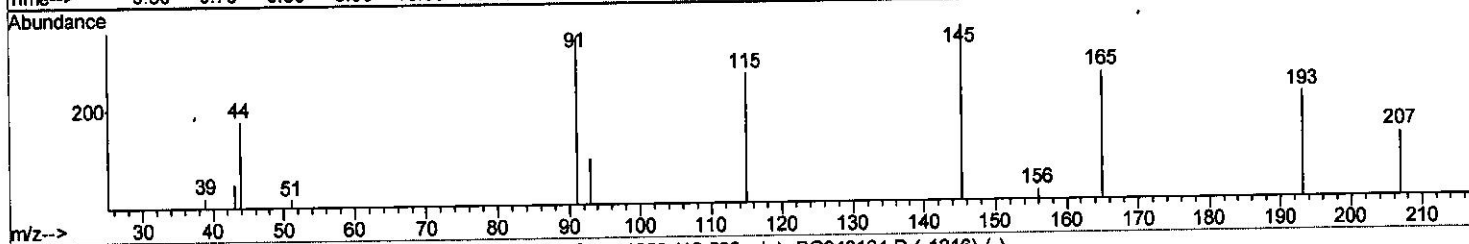
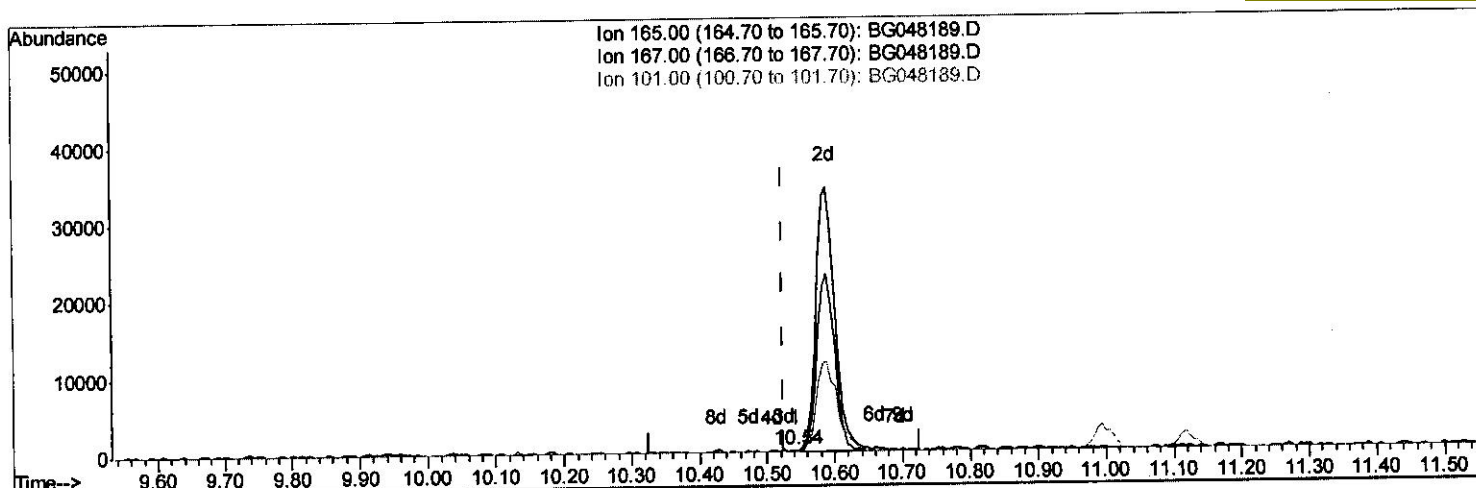
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021621\
 Data File : BG048189.D
 Acq On : 17 Feb 2021 14:38
 Operator : CG/JU
 Sample : SSTDCCC020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
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Manual Integrations
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Quant Time: Feb 17 15:52:52 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 15 06:21:07 2021
 Response via : Initial Calibration



(26) 2,4-Dichlorophenol-d3 (S)

10.542min (+0.017) 0.06ng/ul

response 188

Ion	Exp%	Act%
165.00	100	100
167.00	60.10	0.00#
101.00	26.90	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

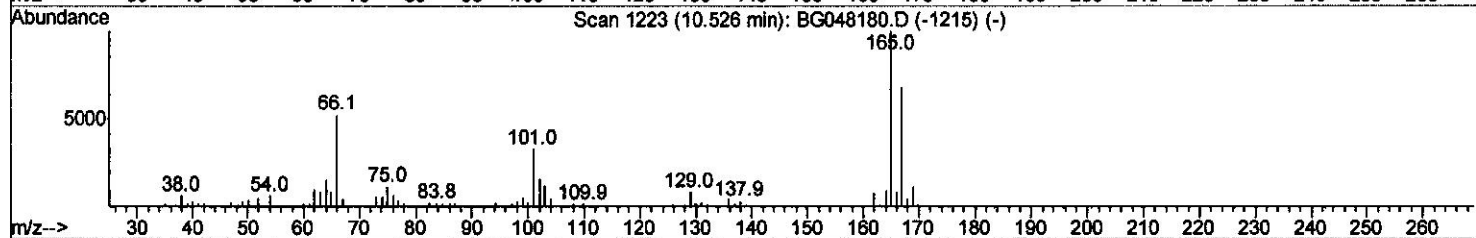
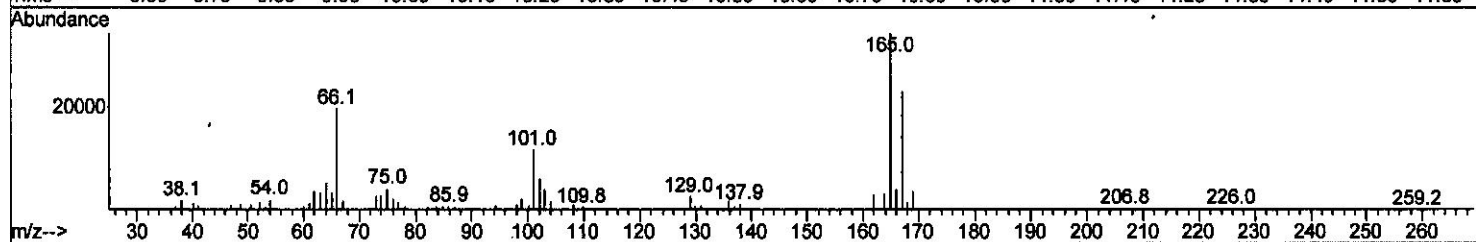
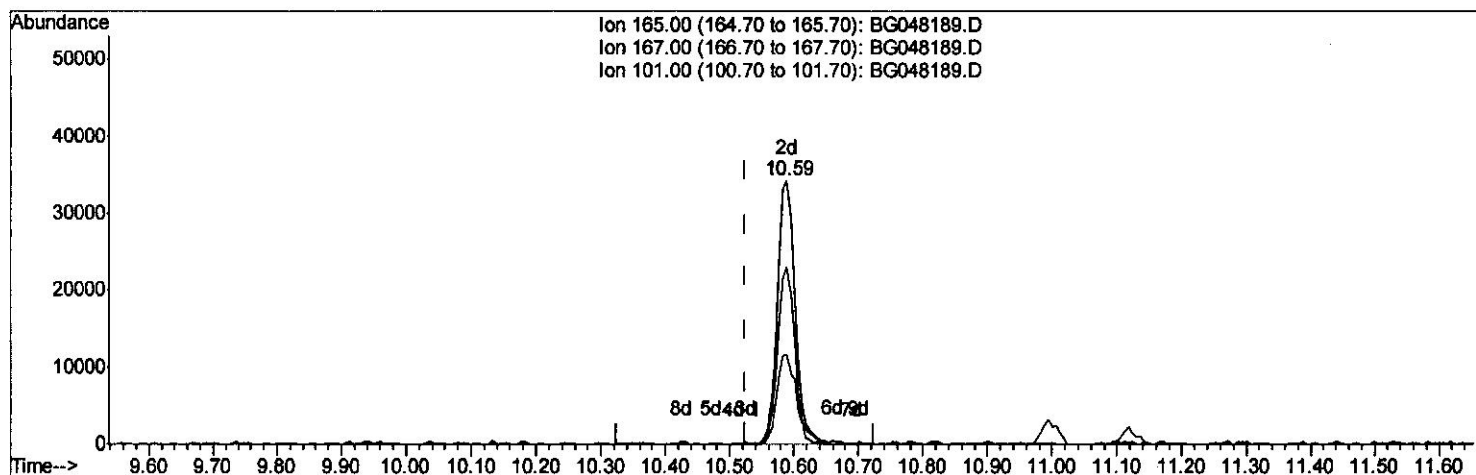
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021621\
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TIC: BG048189.D

(26) 2,4-Dichlorophenol-d3 (S)

10.589min (+0.064) 19.86ng/ul m

response 65466

Ion	Exp%	Act%
165.00	100	100
167.00	60.10	67.32
101.00	26.90	34.29#
0.00	0.00	0.00

Quantitation Report (Qedit)

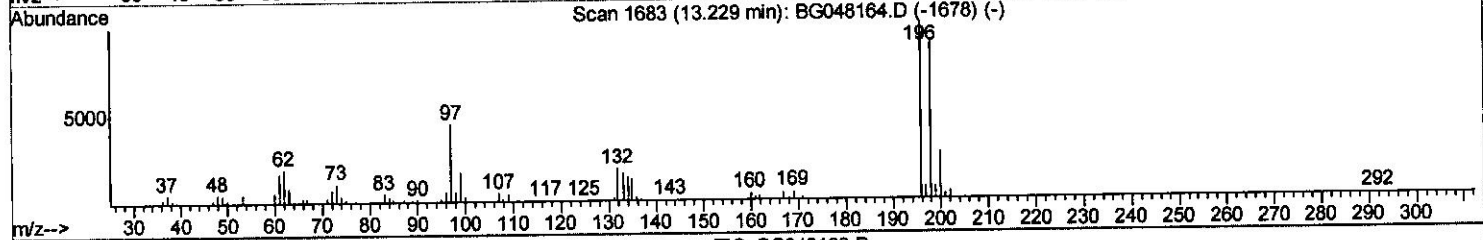
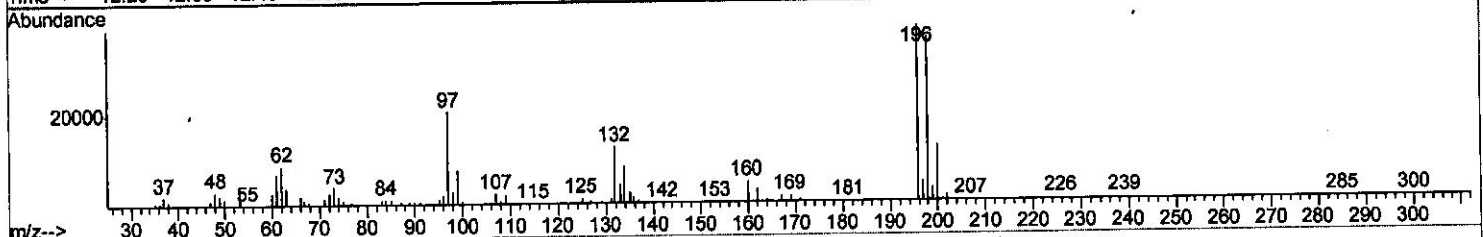
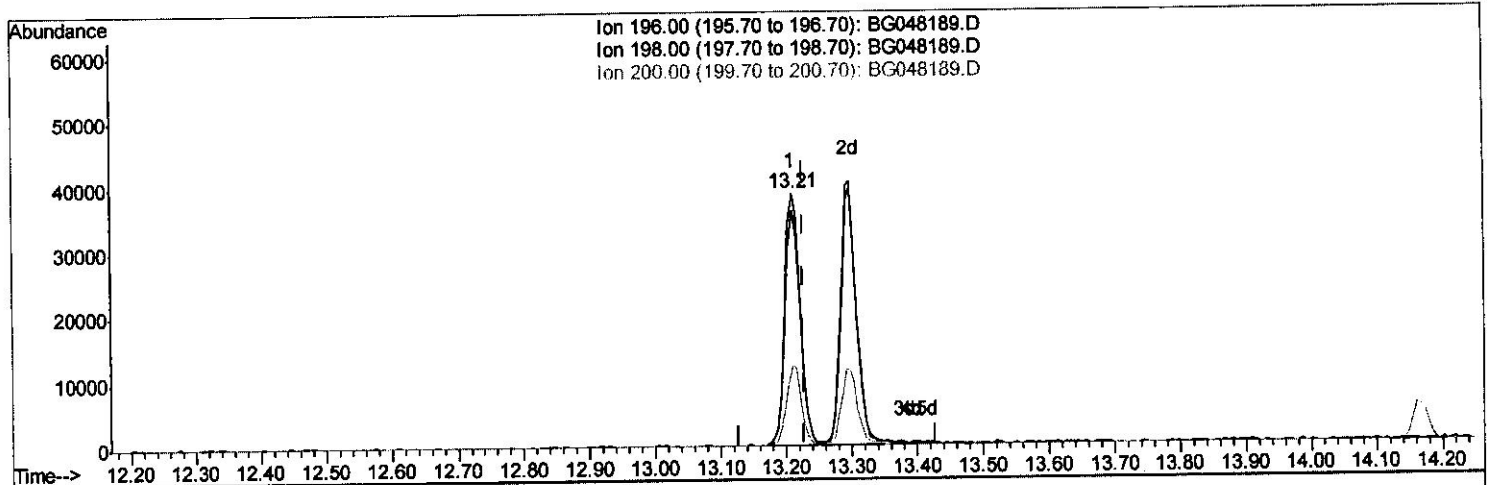
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021621\
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(40) 2,4,5-Trichlorophenol

13.210min (-0.019) 22.01ng/ul

response 63648

Ion	Exp%	Act%
196.00	100	100
198.00	96.70	93.34
200.00	36.90	31.66
0.00	0.00	0.00

Quantitation Report (Qedit)

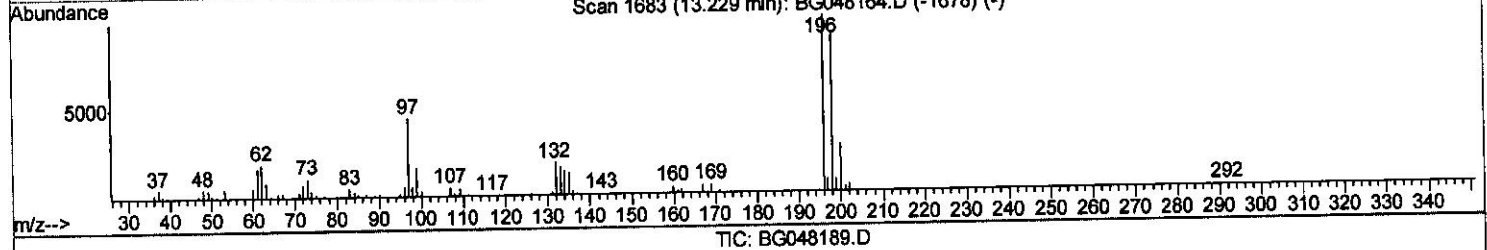
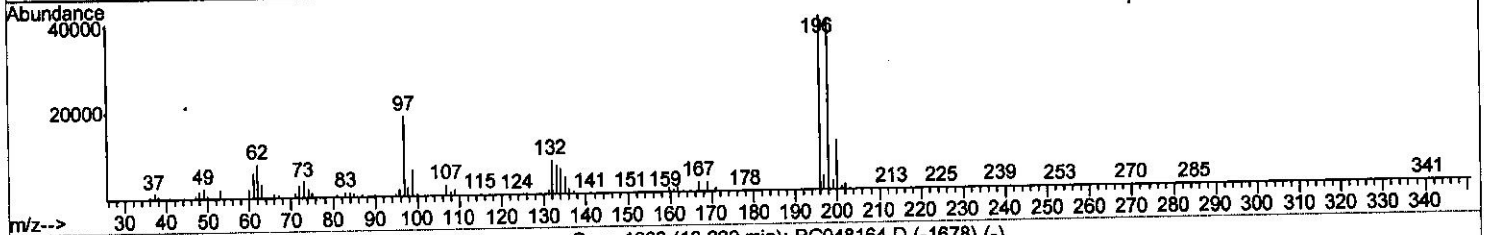
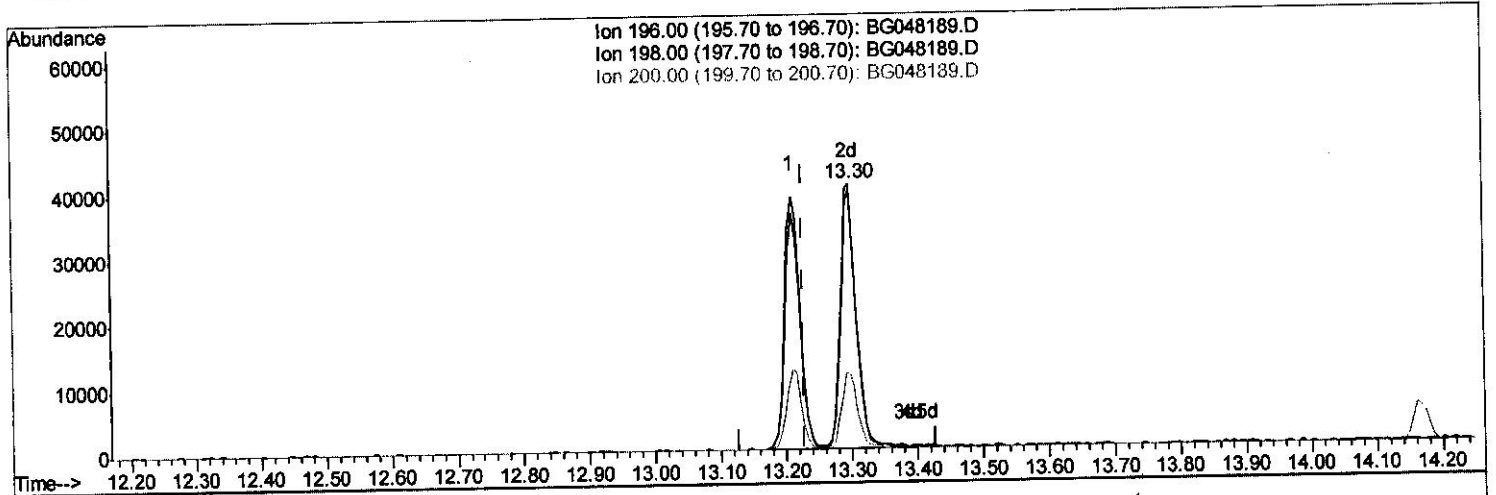
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(40) 2,4,5-Trichlorophenol

13.298min (+0.070) 23.24ng/ul m

response 67185

Ion	Exp%	Act%
196.00	100	100
198.00	96.70	97.63
200.00	36.90	28.26#
0.00	0.00	0.00

Quantitation Report (Qedit)

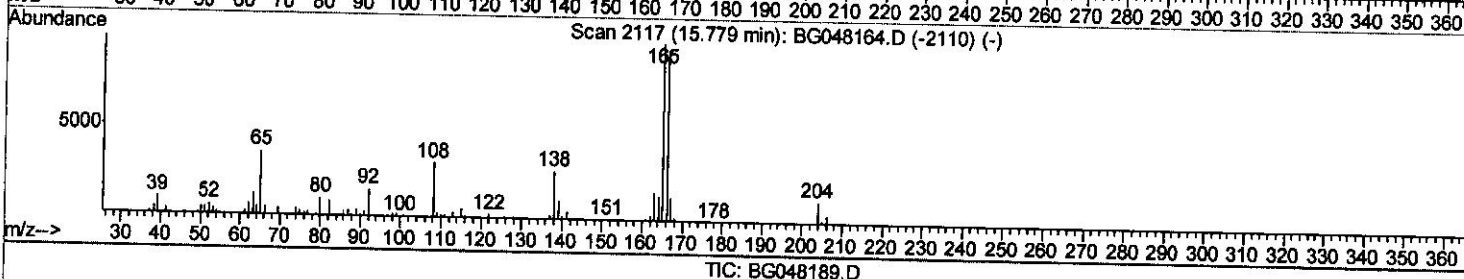
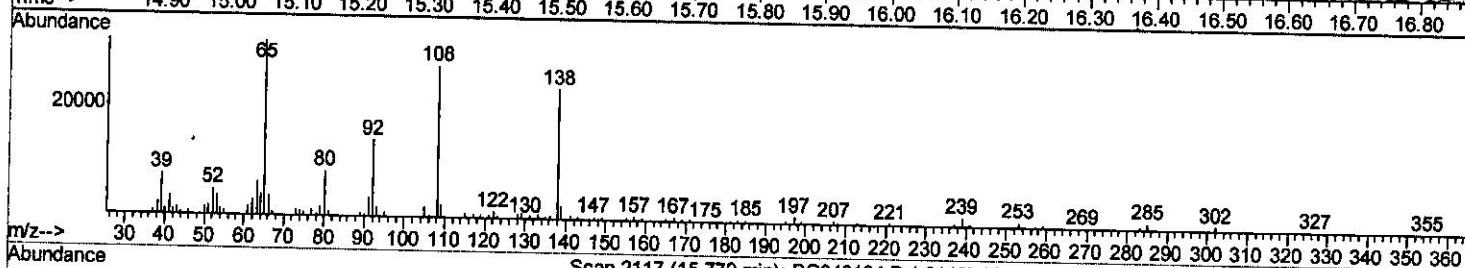
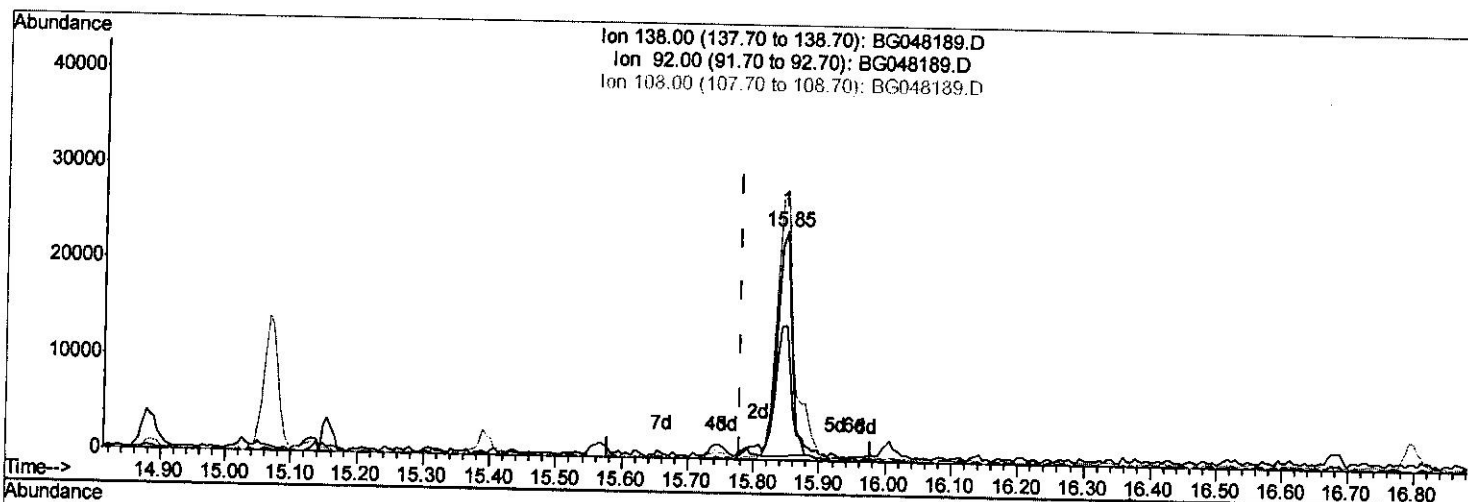
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Manual Integrations
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Quant Time: Feb 17 15:57:17 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
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(61) 4-Nitroaniline

15.848min (+0.070) 18.19ng/ui

response 35692

Ion	Exp%	Act%
138.00	100	100
92.00	57.30	59.01
108.00	108.10	115.79
0.00	0.00	0.00

Quantitation Report (Qedit)

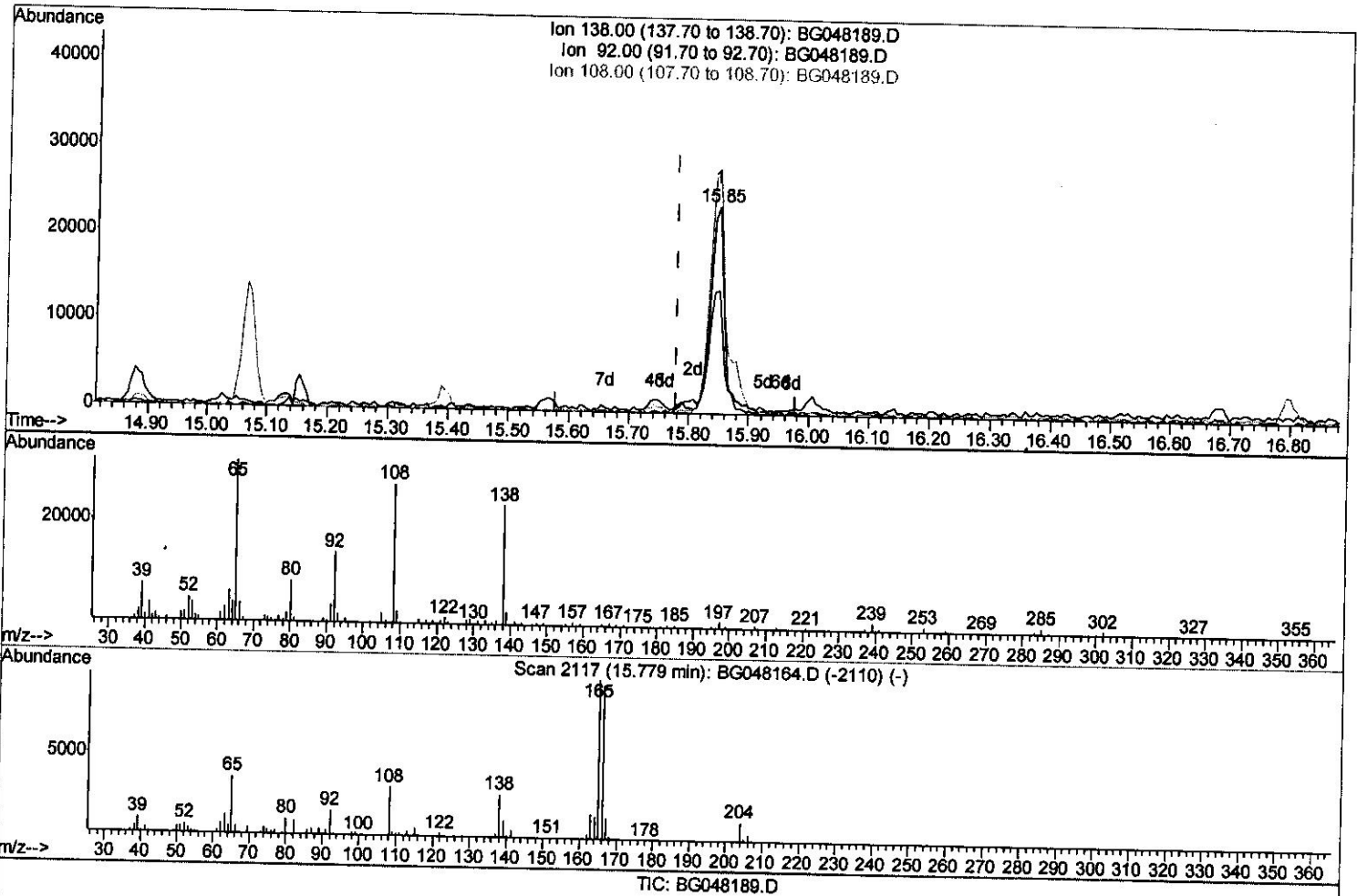
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021621\
 Data File : BG048189.D
 Acq On : 17 Feb 2021 14:38
 Operator : CG/JU
 Sample : SSTDCCC020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02035

Manual Integrations
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Quant Time: Feb 17 15:57:17 2021
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(61) 4-Nitroaniline

15.848min (+0.070) 20.83ng/ul m

response 40879

Ion	Exp%	Act%
138.00	100	100
92.00	57.30	59.01
108.00	108.10	115.79
0.00	0.00	0.00

Quantitation Report (Qedit)

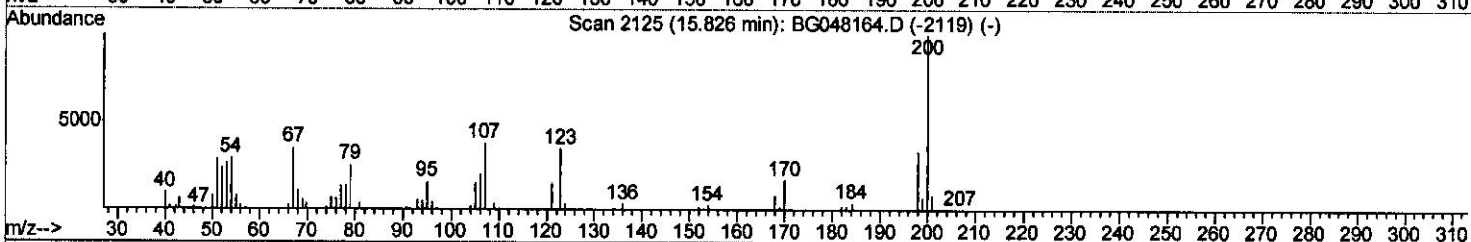
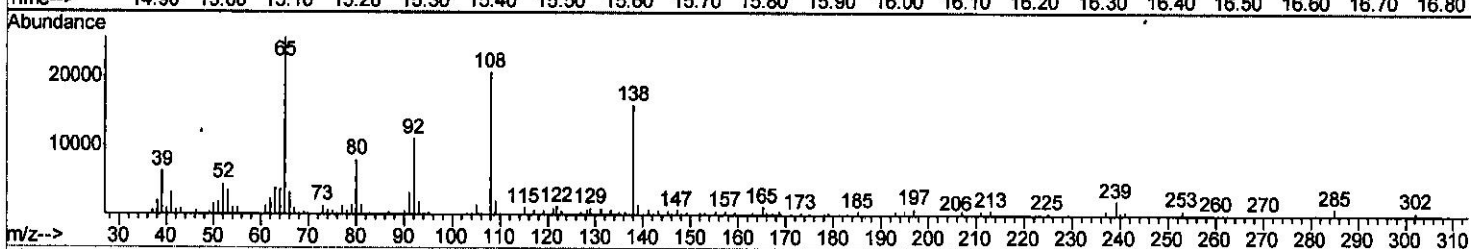
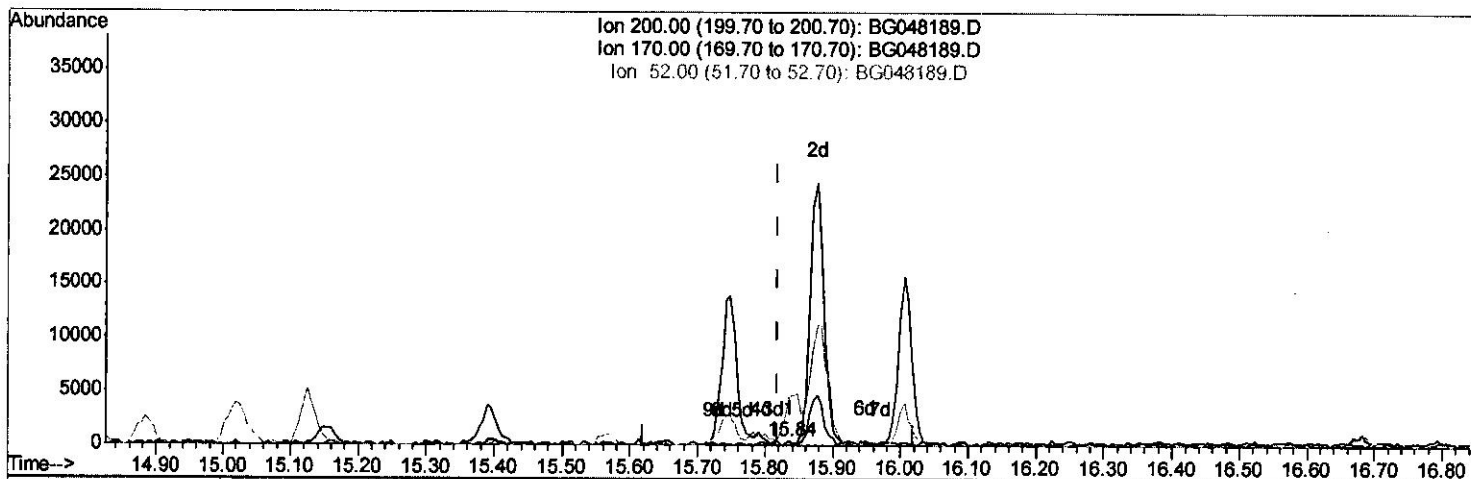
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Manual Integrations
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TIC: BG048189.D

(63) 4,6-Dinitro-2-methylphenol-d2 (S)

15.836min (+0.017) 0.11ng/ul

response 205

Ion	Exp%	Act%
200.00	100	100
170.00	18.70	0.00#
52.00	35.20	1117.40#
0.00	0.00	0.00

Quantitation Report (Qedit)

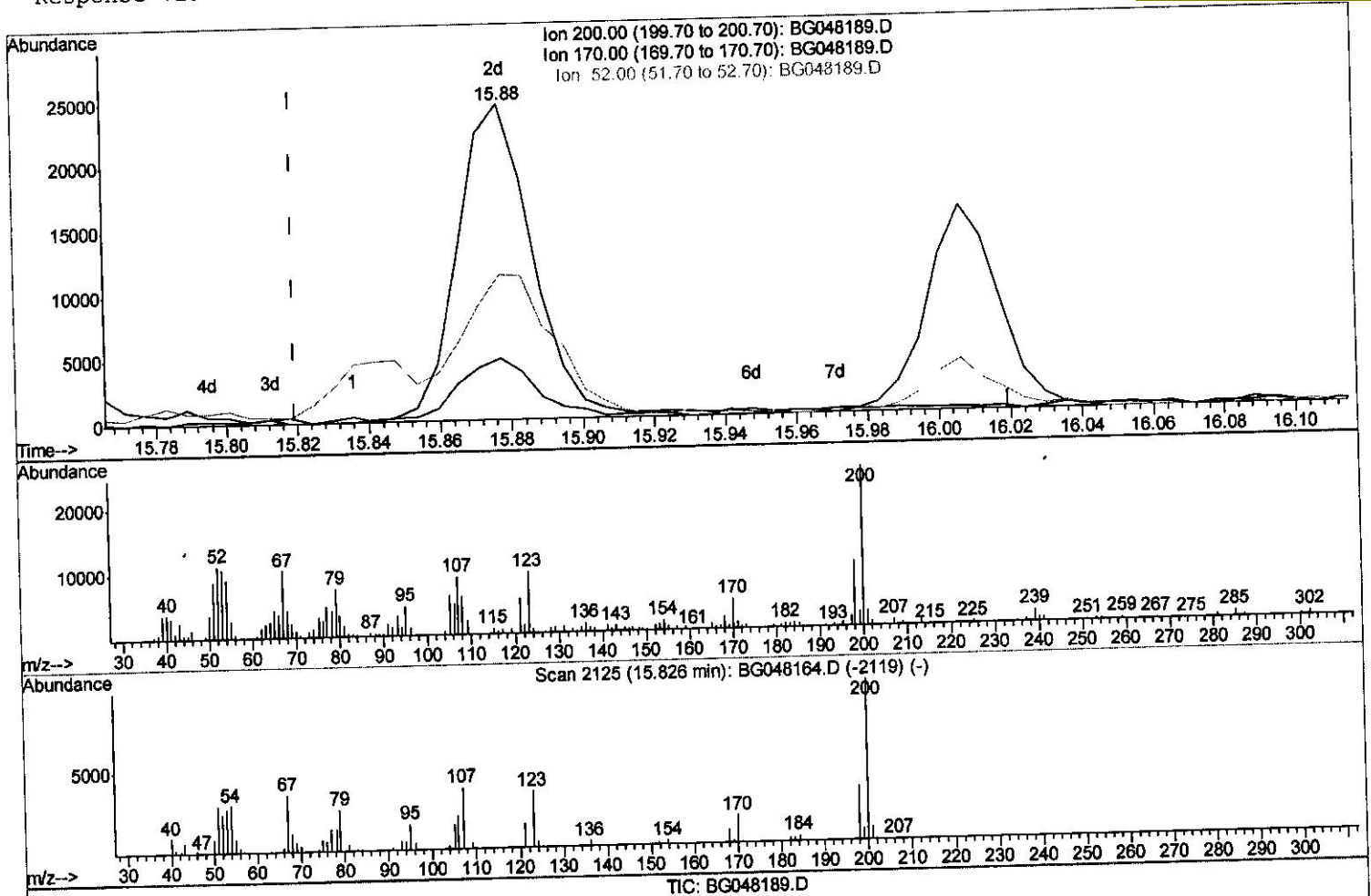
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021621\
 Data File : BG048189.D
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 Operator : CG/JU
 Sample : SSTDCCC020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
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 LabSampleId :
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Manual Integrations
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(63) 4,6-Dinitro-2-methylphenol-d2 (S)

15.877min (+0.058) 19.94ng/ul m

response 35639

Ion	Exp%	Act%
200.00	100	100
170.00	18.70	19.31
52.00	35.20	45.80#
0.00	0.00	0.00

Quantitation Report (Qedit)

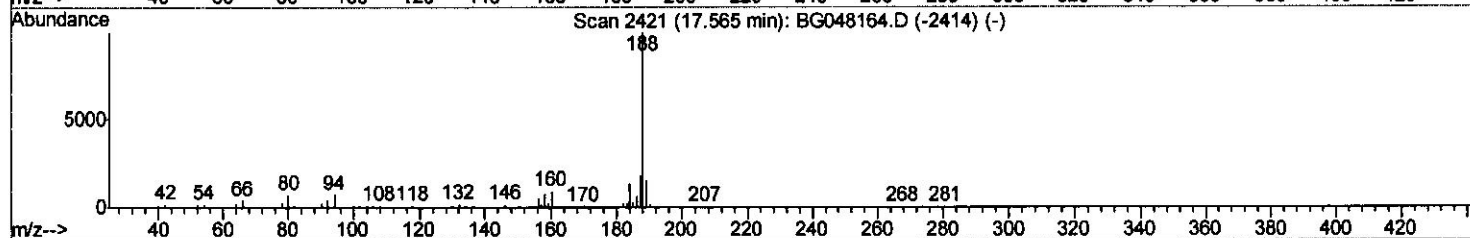
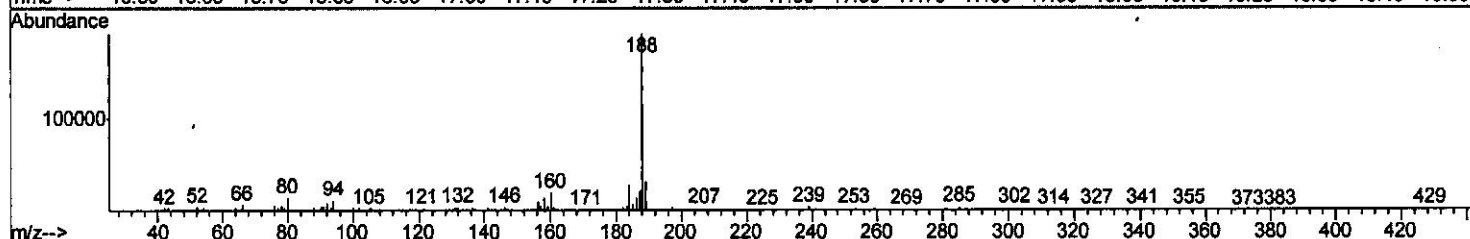
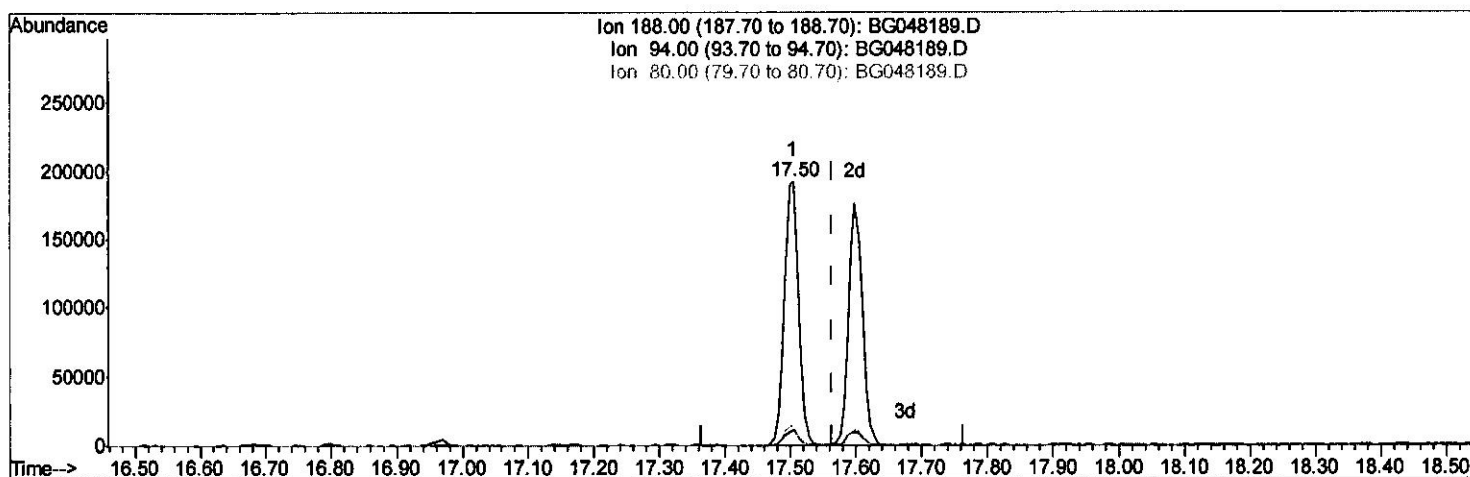
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021621\
 Data File : BG048189.D
 Acq On : 17 Feb 2021 14:38
 Operator : CG/JU
 Sample : SSTDCCC020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
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 LabSampleId :
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Manual Integrations
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TIC: BG048189.D

(71) Anthracene-d10 (S)

17.504min (-0.060) 23.37ng/ul

response 292098

Ion	Exp%	Act%
188.00	100	100
94.00	7.10	5.72
80.00	6.50	6.98
0.00	0.00	0.00

Quantitation Report (Qedit)

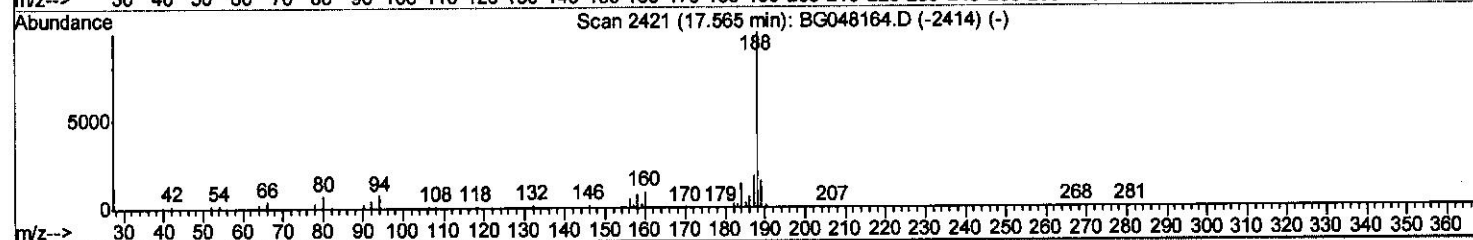
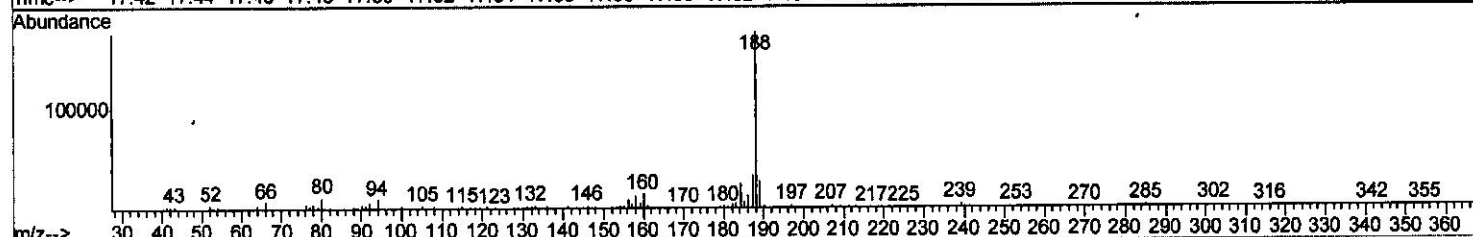
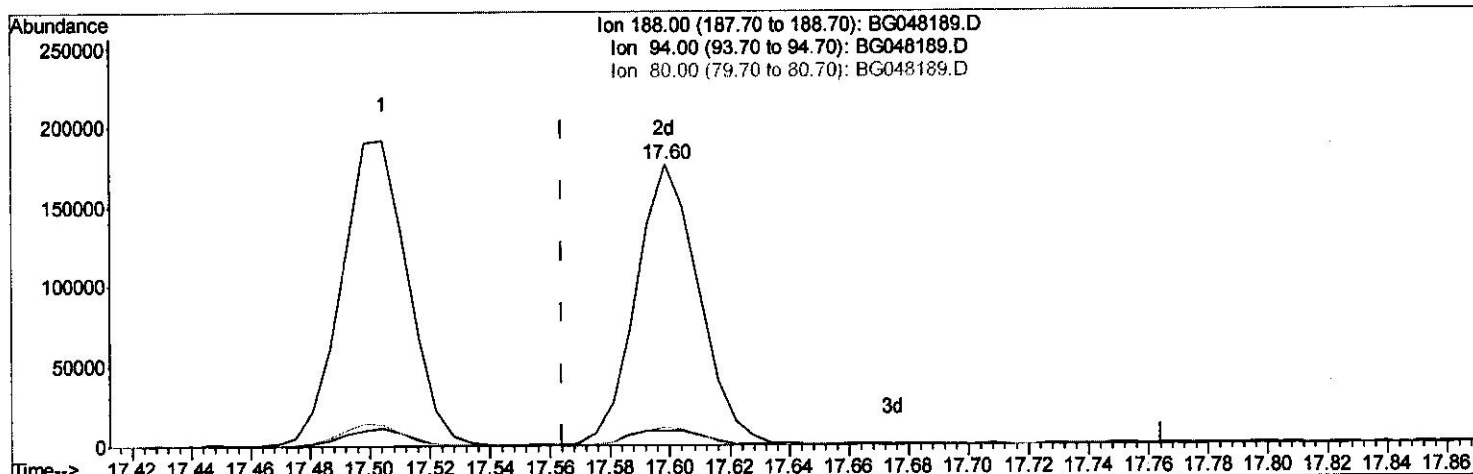
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021621\
 Data File : BG048189.D
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Instrument :
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Manual Integrations
 APPROVED

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TIC: BG048189.D

(71) Anthracene-d10 (S)

17.598min (+0.034) 20.67ng/ul m

response 258355

Ion	Exp%	Act%
188.00	100	100
94.00	7.10	5.33#
80.00	6.50	6.56
0.00	0.00	0.00

2/17/21 JV

Quantitation Report (Qedit)

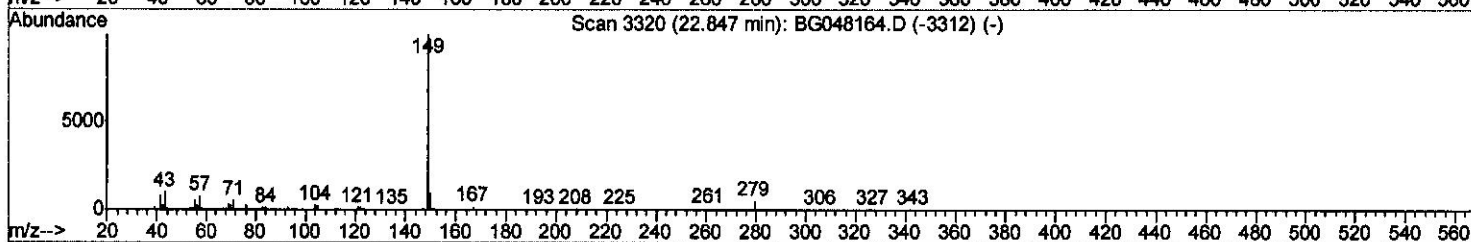
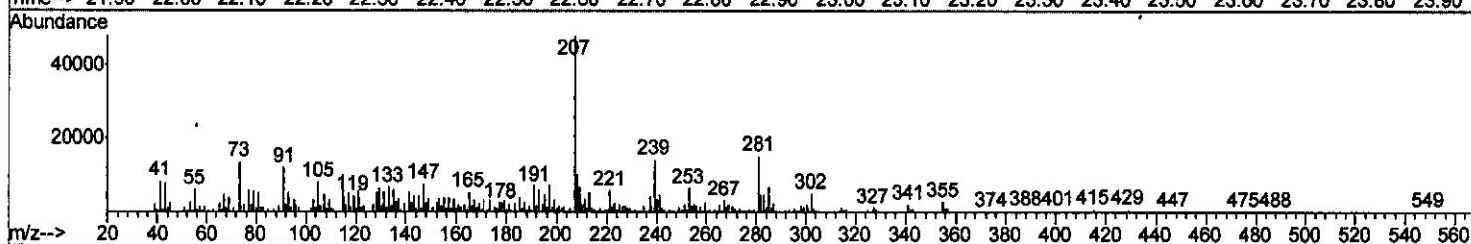
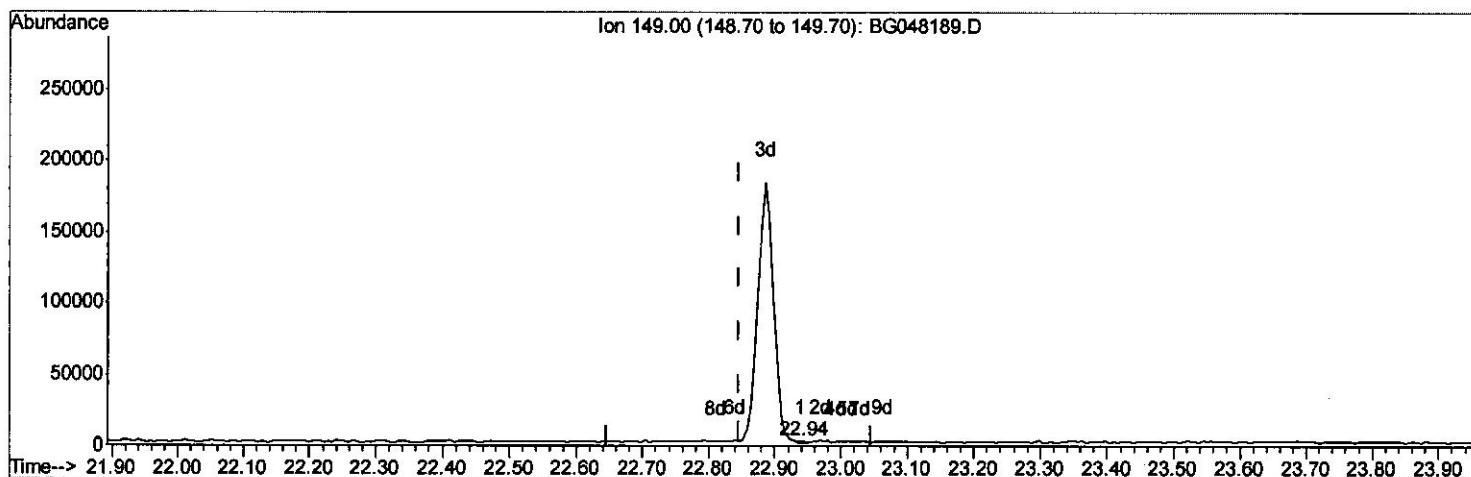
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TIC: BG048189.D

(87) Di-n-octyl phthalate

22.939min (+0.093) 0.03ng/ul

response 485

Ion	Exp%	Act%
149.00	100	100
0.00	0.00	0.00
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

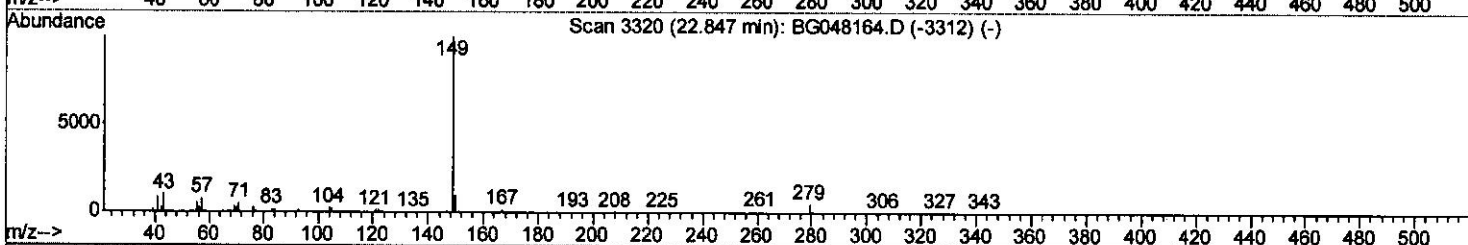
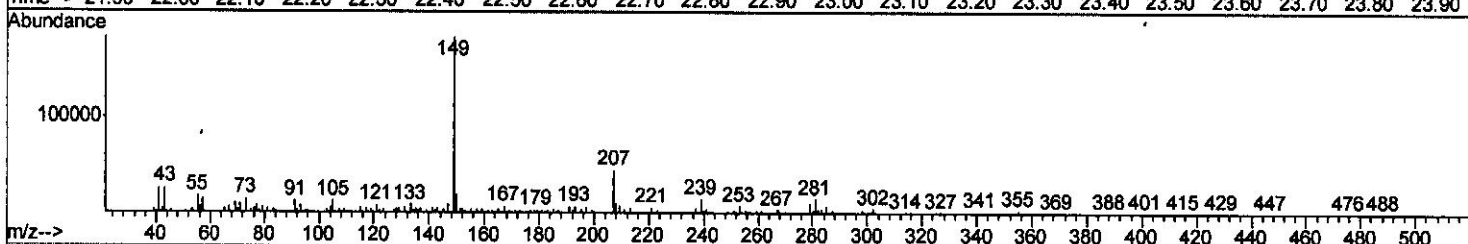
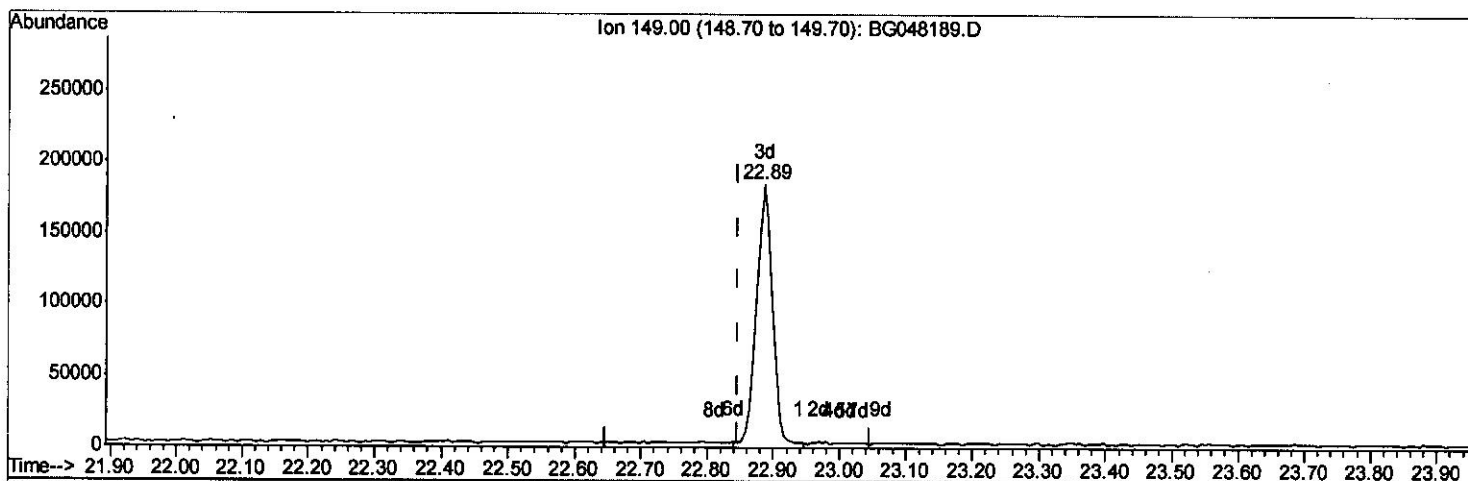
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TIC: BG048189.D

(87) Di-n-octyl phthalate

22.886min (+0.040) 21.35ng/ul m

2/19/21 JU

response 324863

Ion	Exp%	Act%
149.00	100	100
0.00	0.00	0.00
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

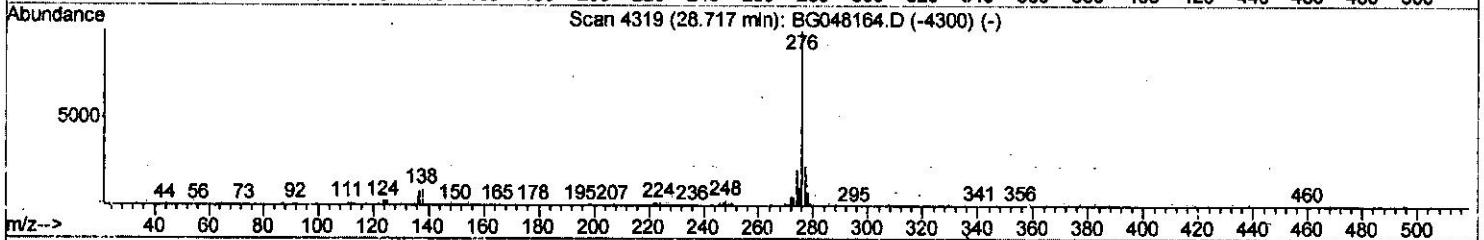
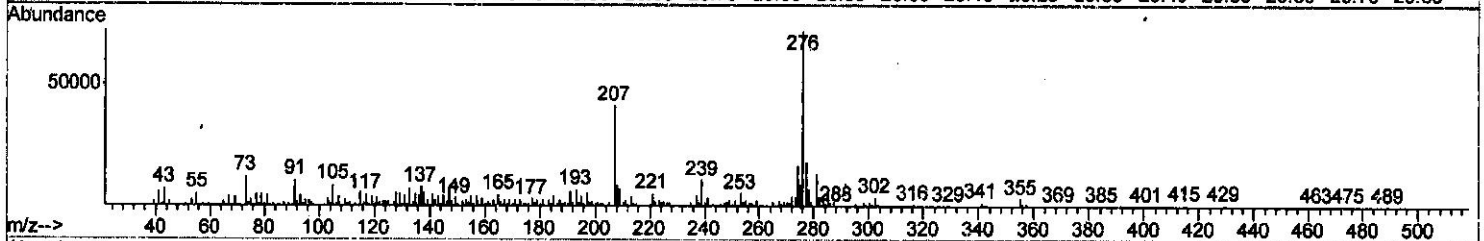
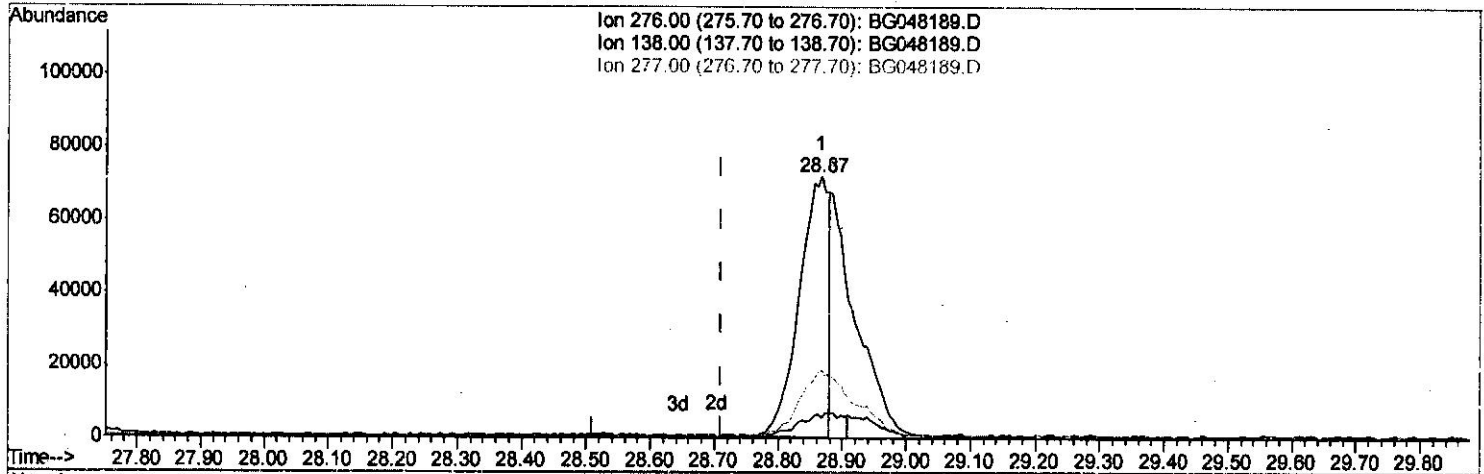
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TIC: BG048189.D

(92) Indeno(1,2,3-cd)pyrene
 28.868min (+0.150) 11.07ng/ul
 response 231784

Ion	Exp%	Act%
276.00	100	100
138.00	10.50	7.95#
277.00	24.20	25.75
0.00	0.00	0.00

Quantitation Report (Qedit)

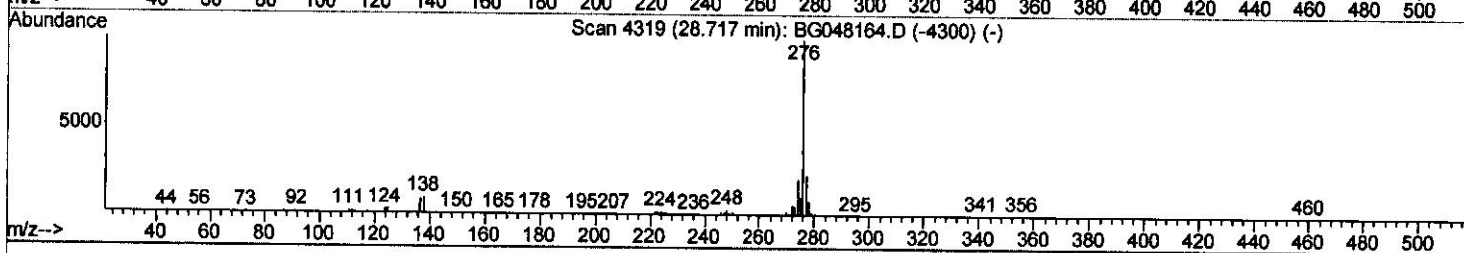
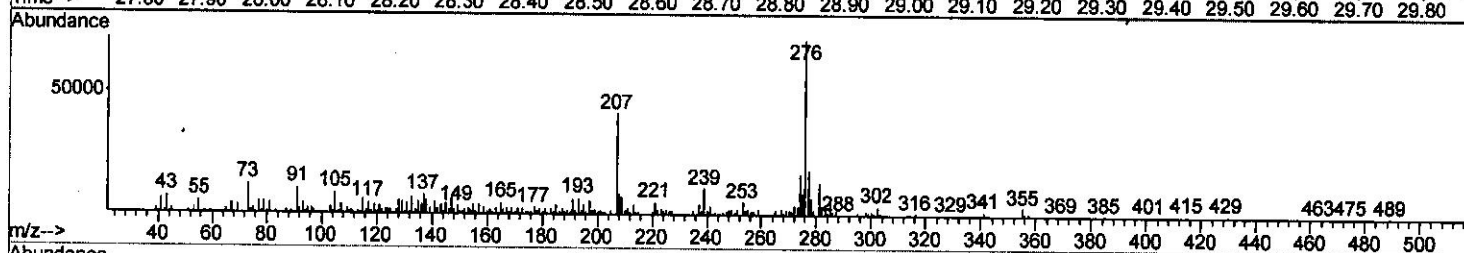
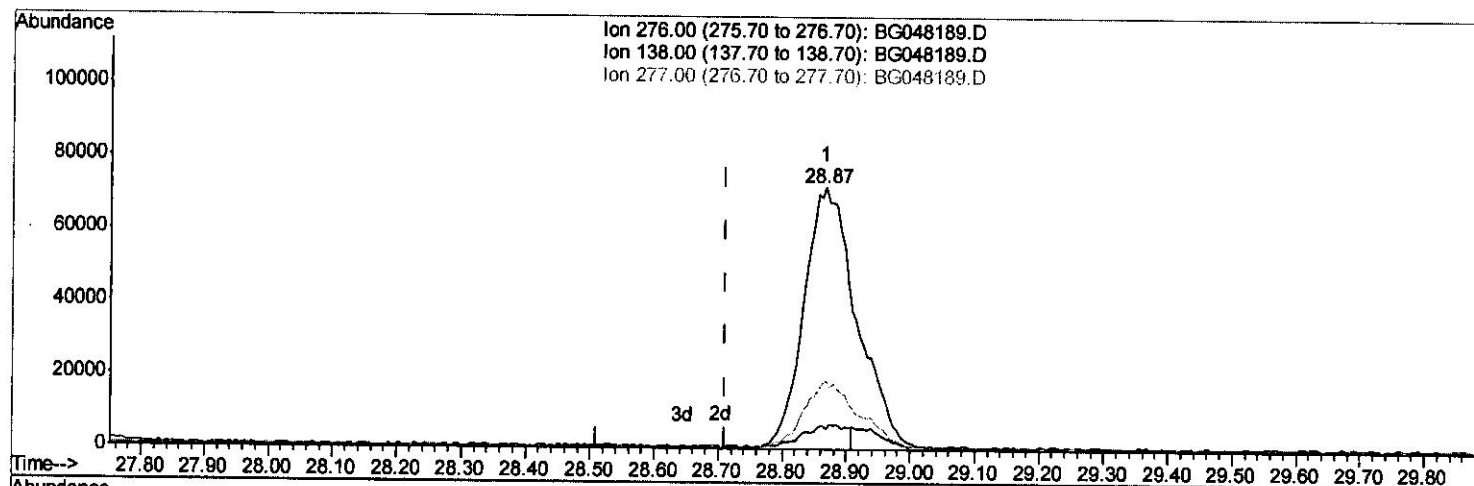
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Quant Time: Feb 17 15:52:52 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 15 06:21:07 2021
 Response via : Initial Calibration



TIC: BG048189.D

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

28.868min (+0.158) 19.86ng/ul m

response 415936

Ion	Exp%	Act%
276.00	100	100
138.00	10.50	7.95#
277.00	24.20	25.75
0.00	0.00	0.00

Quantitation Report (QT² Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021621\
 Data File : BG048189.D
 Acq On : 17 Feb 2021 14:38
 Operator : CG/JU
 Sample : SSTDCCC020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleID :
 SSTD02035

Manual Integrations
 APPROVED

mohammad
 2/18/2021 1:38:36 PM

Quant Time: Feb 17 17:37:32 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
 Quant Title : SVOA CALIBRATION
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	49340	20.00	ng/ul	0.03
18) Naphthalene-d8	10.95	136	186266	20.00	ng/ul	0.03
36) Acenaphthene-d10	14.75	164	128435	20.00	ng/ul	0.03
62) Phenanthrene-d10	17.50	188	291515	20.00	ng/ul	0.04
78) Chrysene-d12	21.78	240	291392	20.00	ng/ul	0.05
86) Perylene-d12	25.09	264	296339	20.00	ng/ul	0.09

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.53	96	8033	6.78	ng/uL	0.02
5) Phenol-d5	7.33	99	76385	18.37	ng/ul	0.09
7) Bis-(2-Chloroethyl)ether-d	7.46	67	45766	18.48	ng/ul	0.03
9) 2-Chlorophenol-d4	7.67	132	56674	19.42	ng/ul	0.05
13) 4-Methylphenol-d8	8.88	113	60033	18.41	ng/ul	0.09
19) Nitrobenzene-d5	9.30	128	28735	20.65	ng/ul	0.04
22) 2-Nitrophenol-d4	10.03	143	36199	21.51	ng/ul	0.04
26) 2,4-Dichlorophenol-d3	10.59	165	65466m	19.86	ng/ul	0.06
29) 4-Chloroaniline-d4	11.09	131	73830	22.67	ng/ul	0.05
44) Dimethylphthalate-d6	14.17	166	184687	19.41	ng/ul	0.05
47) Acenaphthylene-d8	14.46	160	229822	19.76	ng/ul	0.03
52) 4-Nitrophenol-d4	15.02	143	30915	19.12	ng/ul	0.12
58) Fluorene-d10	15.75	176	162669	19.17	ng/ul	0.03
63) 4,6-Dinitro-2-methylphenol	15.88	200	35639m	19.94	ng/ul	0.06
71) Anthracene-d10	17.60	188	258355m	20.67	ng/ul	0.03
79) Pyrene-d10	19.88	212	290597	19.97	ng/ul	0.03
90) Benzo(a)pyrene-d12	24.86	264	294607	19.49	ng/ul	0.09

2/19/21 JU

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.57	88	8908	6.554	ng/uL# 95
4) Benzaldehyde	7.28	77	55507	22.711	ng/ul 98
6) Phenol	7.36	94	81504	18.989	ng/ul 97
8) Bis(2-Chloroethyl)ether	7.55	93	59314	18.298	ng/ul 86
10) 2-Chlorophenol	7.71	128	59126	19.153	ng/ul 94
11) 2-Methylphenol	8.61	108	59117	19.037	ng/ul 97
12) 2,2'-oxybis(1-Chloropropan	8.66	45	77899	18.386	ng/ul# 92
14) Acetophenone	8.97	105	101072	19.030	ng/ul# 78
15) N-Nitroso-di-n-propylamine	8.95	70	55708	18.889	ng/ul# 92
16) 4-Methylphenol	8.94	108	64421	19.224	ng/ul 97
17) Hexachloroethane	9.21	117	26636	19.316	ng/ul 97
20) Nitrobenzene	9.35	77	83642	19.737	ng/ul 97
21) Isophorone	9.89	82	143507	19.167	ng/ul 97
23) 2-Nitrophenol	10.06	139	36735	21.406	ng/ul 97
24) 2,4-Dimethylphenol	10.14	107	76342	20.102	ng/ul 99
25) Bis(2-Chloroethoxy)methane	10.35	93	79679	18.878	ng/ul# 98
27) 2,4-Dichlorophenol	10.61	162	66587	21.117	ng/ul 95
28) Naphthalene	10.99	128	184136	19.599	ng/ul 98
30) 4-Chloroaniline	11.12	127	73627	22.037	ng/ul 95
31) Hexachlorobutadiene	11.26	225	55990	20.359	ng/ul 95
32) Caprolactam	11.99	113	21249	19.312	ng/ul 82
33) 4-Chloro-3-methylphenol	12.26	107	66447	19.034	ng/ul 97
34) 2-Methylnaphthalene	12.59	142	136062	19.053	ng/ul 94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.81	142	128175	18.555	ng/uL	97
37) 1,2,4,5-Tetrachlorobenzene	12.95	216	97842	21.177	ng/uL	99
38) Hexachlorocyclopentadiene	12.92	237	66517	22.331	ng/uL	98
39) 2,4,6-Trichlorophenol	13.21	196	63399	22.285	ng/uL	96
40) 2,4,5-Trichlorophenol	13.30	196	67185m	23.238	ng/uL	
41) 1,1'-Biphenyl	13.59	154	177256	20.119	ng/uL	95
42) 2-Chloronaphthalene	13.64	162	144703	19.933	ng/uL	96
43) 2-Nitroaniline	13.86	65	52473	20.901	ng/uL	94
45) Dimethylphthalate	14.21	163	192381	19.562	ng/uL	97
46) 2,6-Dinitrotoluene	14.34	165	42666	20.628	ng/uL	94
48) Acenaphthylene	14.48	152	207158	19.917	ng/uL	99
49) 3-Nitroaniline	14.68	138	36852	21.139	ng/uL#	84
50) Acenaphthene	14.82	153	150683	19.303	ng/uL	96
51) 2,4-Dinitrophenol	14.88	184	22752	17.614	ng/uL#	80
53) 4-Nitrophenol	15.03	109	36865	20.200	ng/uL	92
54) Dibenzofuran	15.15	168	221247	19.420	ng/uL	95
55) 2,4-Dinitrotoluene	15.12	165	57007	19.189	ng/uL	89
56) 2,3,4,6-Tetrachlorophenol	15.40	232	61854	20.927	ng/uL	97
57) Diethylphthalate	15.57	149	190052	18.721	ng/uL	97
59) Fluorene	15.80	166	170966	18.562	ng/uL	97
60) 4-Chlorophenyl-phenylether	15.79	204	107183	19.398	ng/uL	94
61) 4-Nitroaniline	15.85	138	40879m	20.830	ng/uL	
64) 4,6-Dinitro-2-methylphenol	15.89	198	39242	22.033	ng/uL	98
65) N-Nitrosodiphenylamine	16.01	169	164448	21.979	ng/uL	96
66) 4-Bromophenyl-phenylether	16.68	248	71834	20.848	ng/uL	95
67) Hexachlorobenzene	16.81	284	75512	20.675	ng/uL#	92
68) Atrazine	16.97	200	72020	21.526	ng/uL	90
69) Pentachlorophenol	17.16	266	45287	22.206	ng/uL#	86
70) Phenanthrene	17.55	178	291969	19.934	ng/uL	97
72) Anthracene	17.63	178	299426	20.266	ng/uL	99
73) 1,2,3,4-Tetrachlorobenzene	13.56	216	100414	23.053	ng/uL	90
74) Pentachlorobenzene	15.07	250	93533	21.900	ng/uL	95
75) Carbazole	17.92	167	257418	20.833	ng/uL	97
76) Di-n-butylphthalate	18.45	149	305463	20.267	ng/uL	97
77) Fluoranthene	19.55	202	363744	20.369	ng/uL#	95
80) Pyrene	19.91	202	361533	19.970	ng/uL	98
81) Butylbenzylphthalate	20.78	149	127010	20.049	ng/uL	90
82) 3,3'-Dichlorobenzidine	21.68	252	128386	21.535	ng/uL#	98
83) Benzo(a)anthracene	21.76	228	356952	19.608	ng/uL	95
84) Bis(2-ethylhexyl)phthalate	21.65	149	176172	19.152	ng/uL	97
85) Chrysene	21.83	228	347164	19.954	ng/uL	99
87) Di-n-octyl phthalate	22.89	149	324863m	21.349	ng/uL	
88) Benzo(b)fluoranthene	24.03	252	355556	19.107	ng/uL#	86
89) Benzo(k)fluoranthene	24.10	252	353314	19.386	ng/uL	94
91) Benzo(a)pyrene	24.93	252	313349	19.138	ng/uL#	90
92) Indeno(1,2,3-cd)pyrene	28.87	276	415936m	19.857	ng/uL	
93) Dibenzo(a,h)anthracene	28.94	278	342049	19.528	ng/uL#	95
94) Benzo(a,h,i)perylene	30.05	276	331496	19.031	ng/uL	95

2/19/21 JV

2/19/21 JV

2/19/21 JV

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

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