

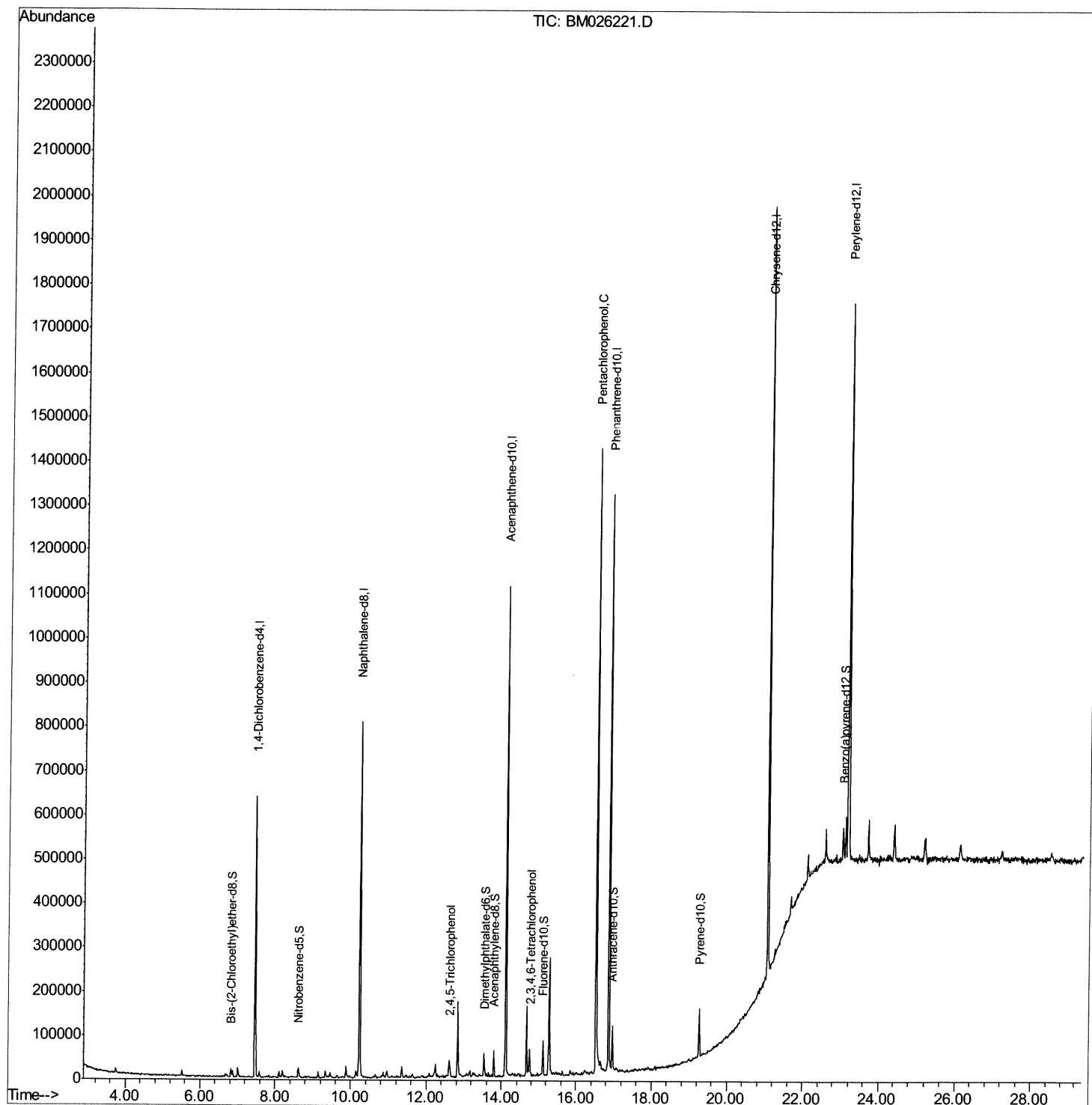
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM060120\
Data File : BM026221.D
Acq On : 02 Jun 2020 10:54
Operator : JU/CG
Sample : L2829-06DL 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C5CD0DL

Manual Integrations
APPROVED

mohammad
6/2/2020 3:22:18 PM

Quant Time: Jun 02 11:42:36 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 02 03:00:44 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

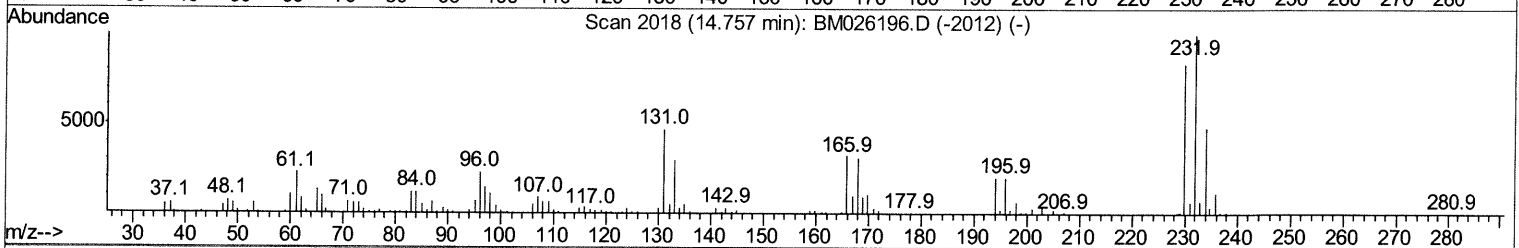
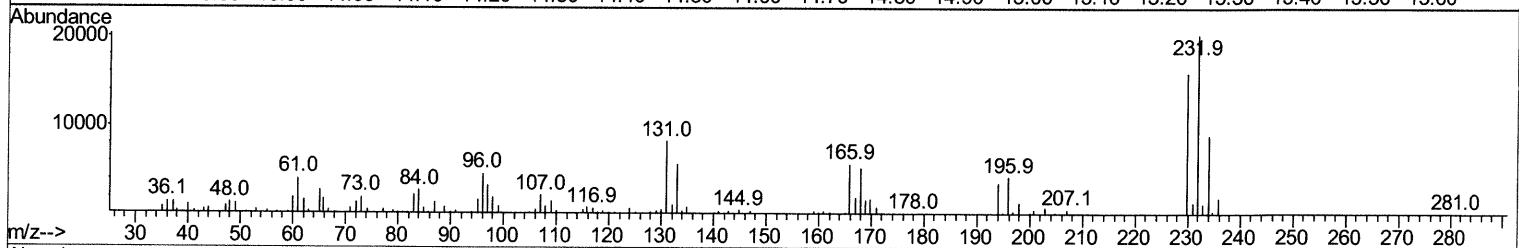
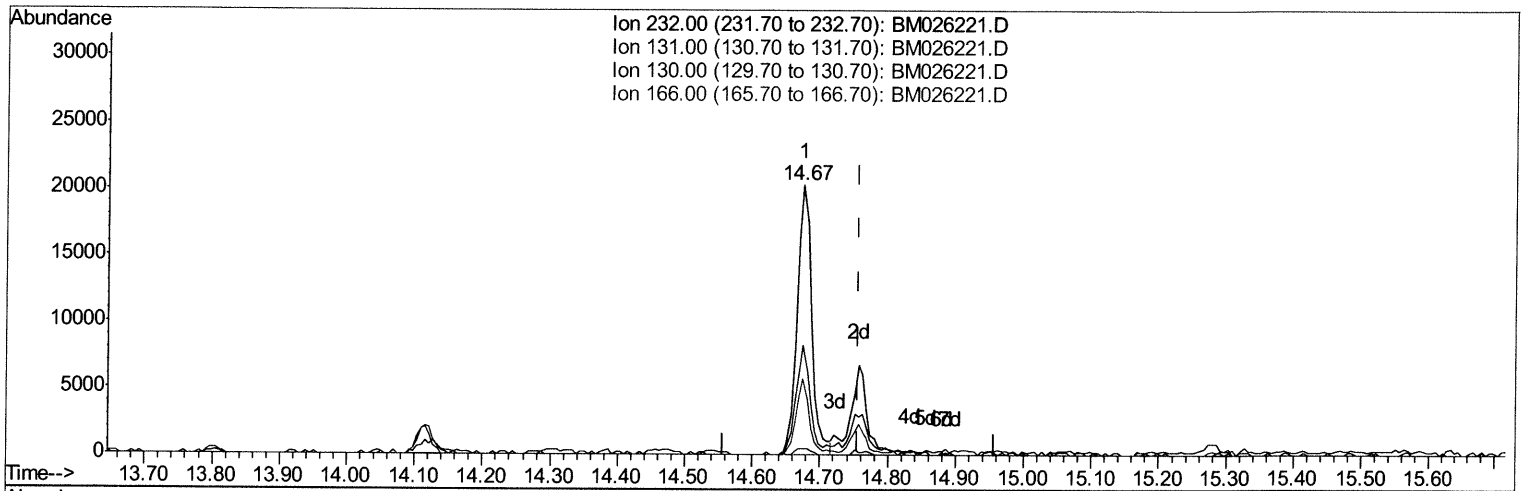
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TIC: BM026221.D

(56) 2,3,4,6-Tetrachlorophenol

14.674min (-0.082) 4.55ng/ul

response 30308

Ion	Exp%	Act%
232.00	100	100
131.00	46.20	40.62
130.00	2.30	2.47
166.00	31.70	28.10

Quantitation Report (Qedit)

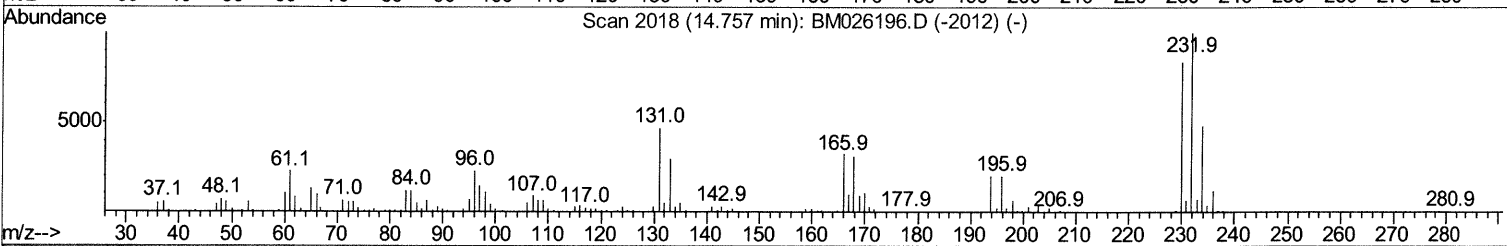
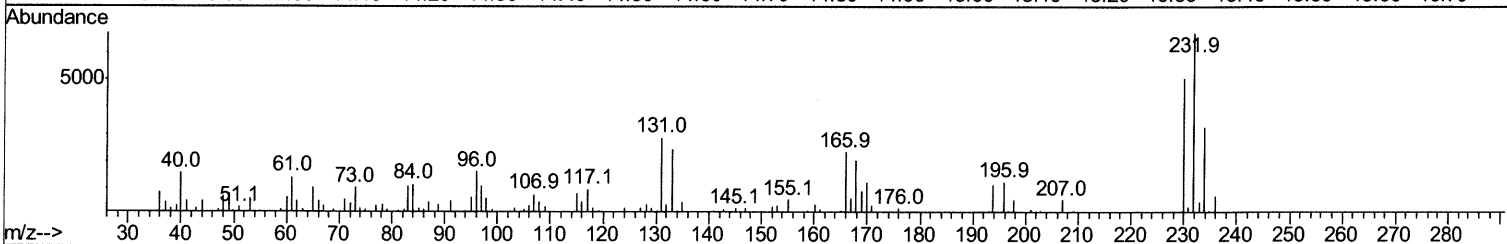
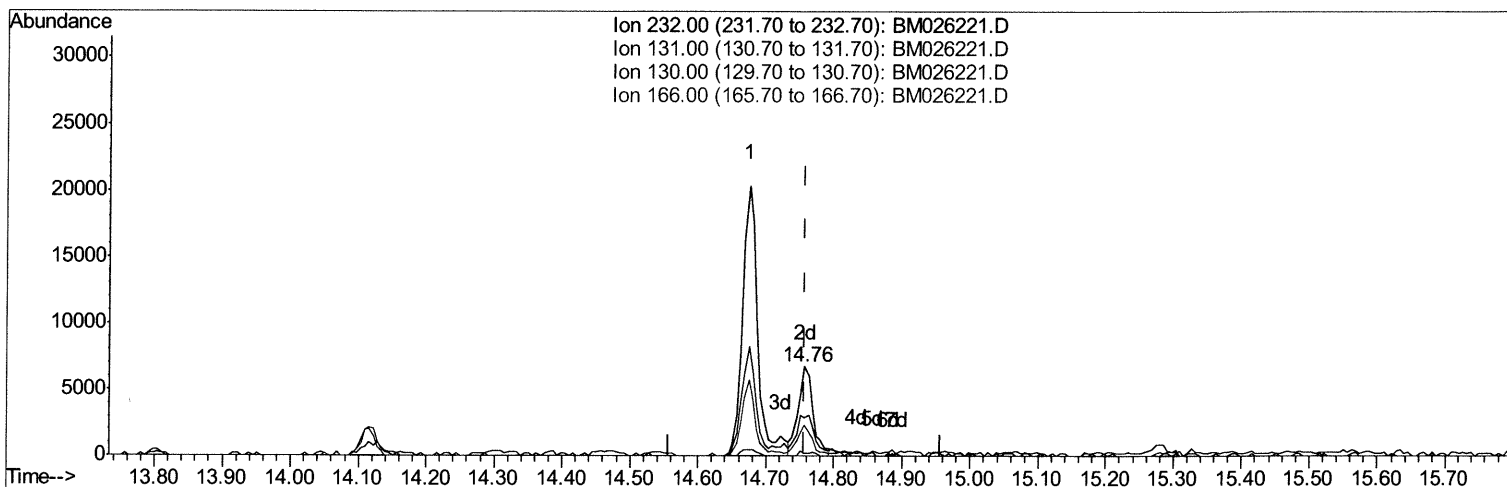
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TIC: BM026221.D

(56) 2,3,4,6-Tetrachlorophenol

14.757min (-0.000) 1.52ng/ul m *06/04/20 JU*

response 10129

Ion	Exp%	Act%
232.00	100	100
131.00	46.20	42.50
130.00	2.30	3.01#
166.00	31.70	34.80

Quantitation Report (Qedit)

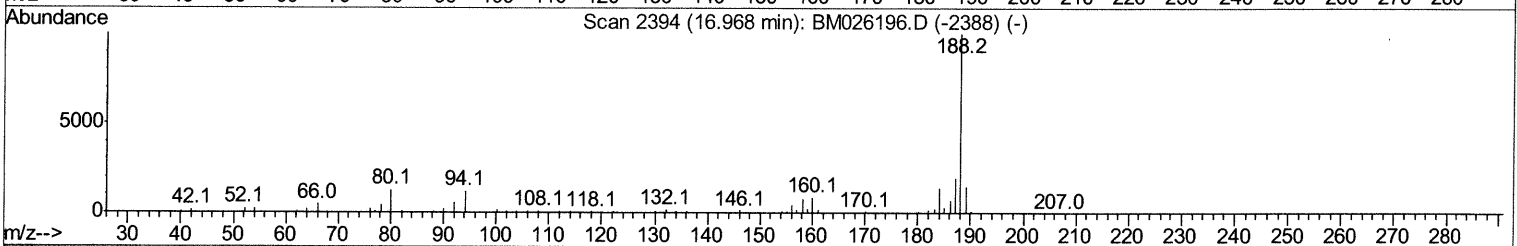
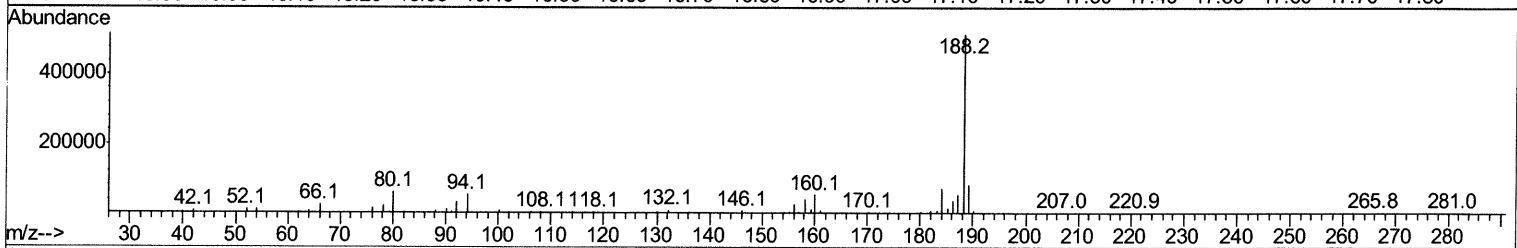
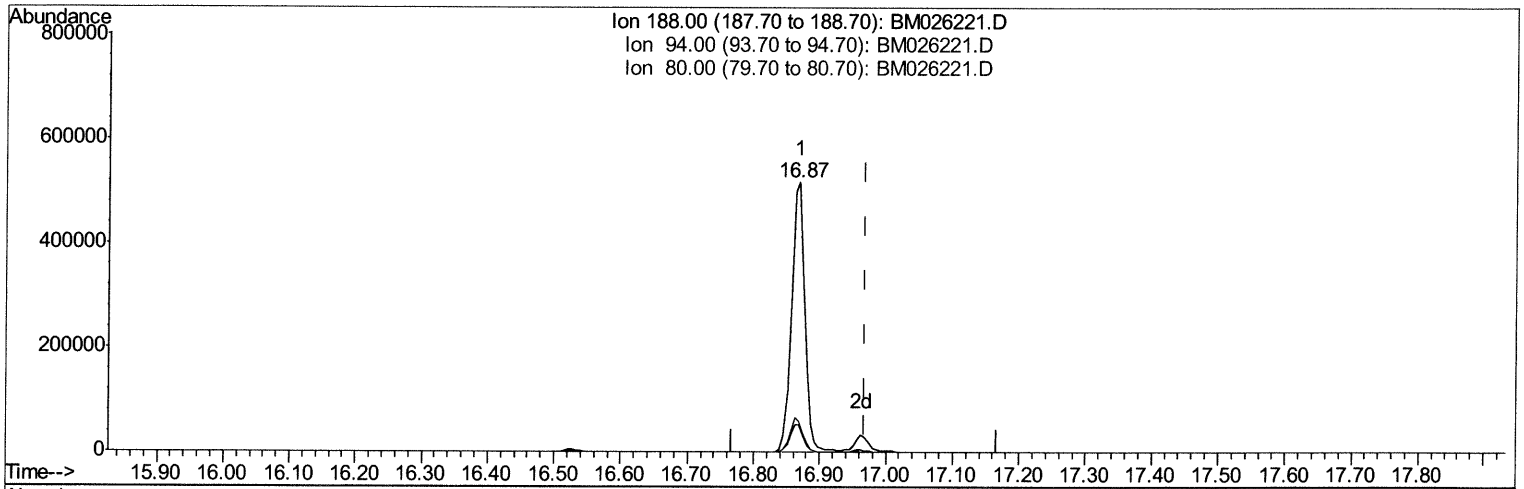
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ALS Vial : 26 Sample Multiplier: 1

Instrument :
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Manual Integrations
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TIC: BM026221.D

(71) Anthracene-d10 (S)

16.868min (-0.100) 20.29ng/ul

response 732014

Ion	Exp%	Act%
188.00	100	100
94.00	12.10	10.08
80.00	12.20	11.44
0.00	0.00	0.00

Quantitation Report (Qedit)

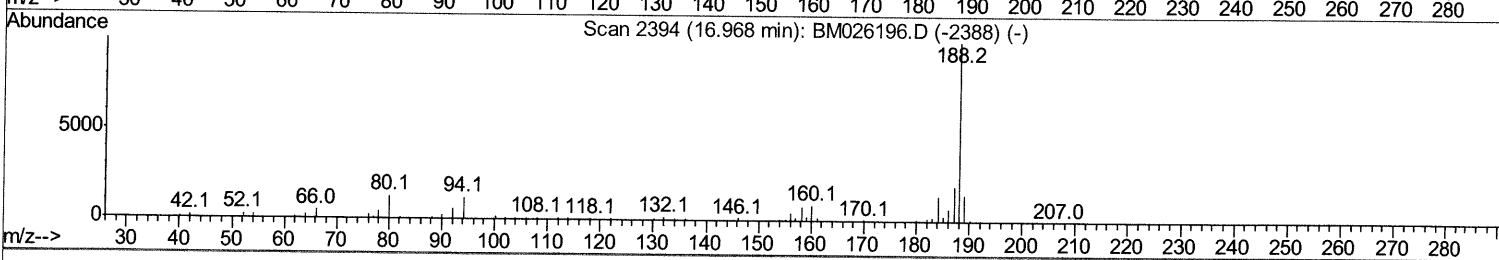
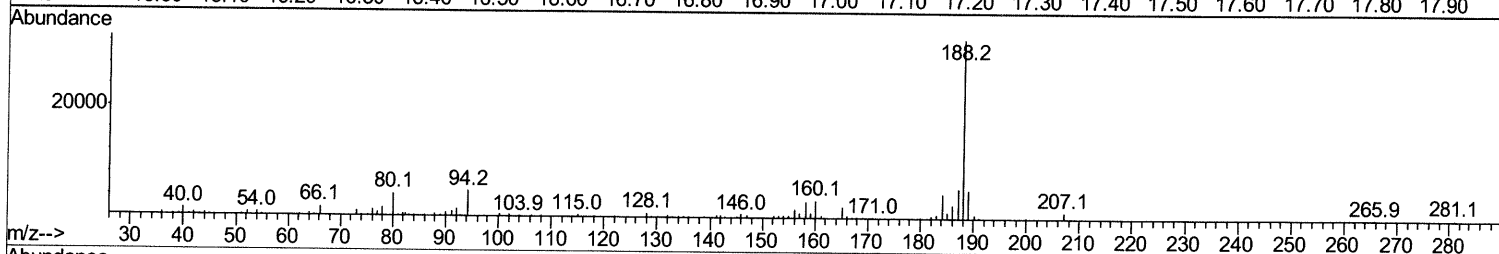
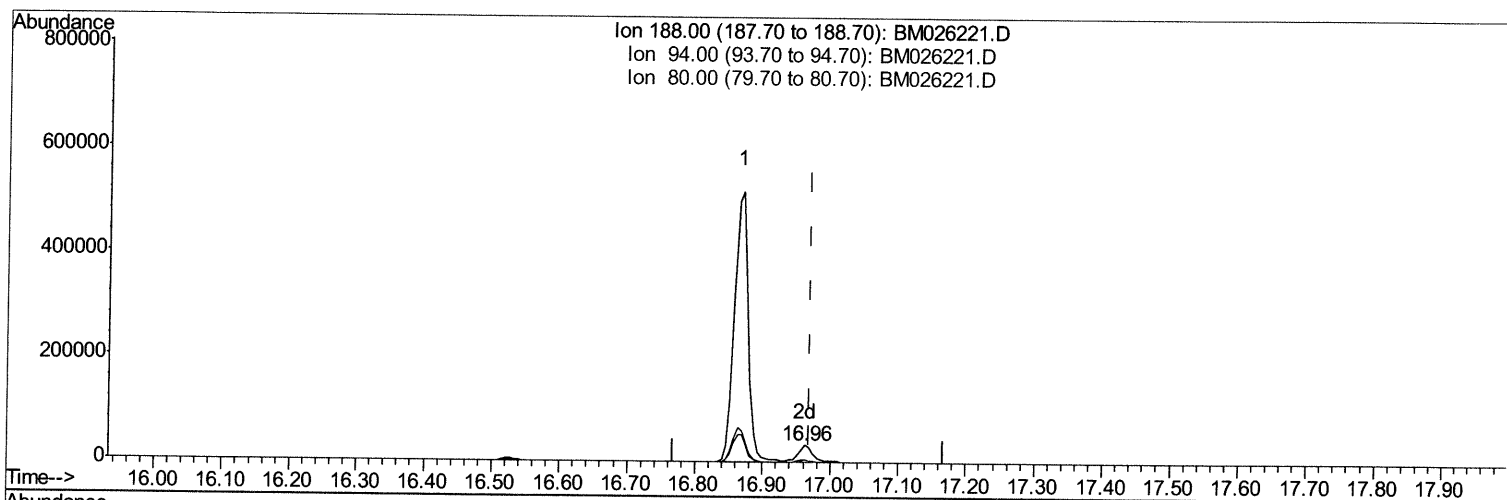
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Manual Integrations
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TIC: BM026221.D

(71) Anthracene-d10 (S)

16.962min (-0.006) 1.39ng/ul m 06/04/20 JU

response 50214

Ion	Exp%	Act%
188.00	100	100
94.00	12.10	14.54#
80.00	12.20	12.78
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	7.47	152	149118	20.00	ng/ul	0.00
18) Naphthalene-d8	10.23	136	591712	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.12	164	344493	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.87	188	732014	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	770316	20.00	ng/ul	0.00
86) Perylene-d12	23.18	264	809125	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.68	99	3554	0.26	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	6.83	67	9715	1.04	ng/ul	0.00
9) 2-Chlorophenol-d4	7.02	132	8320	0.79	ng/ul	0.00
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul	
19) Nitrobenzene-d5	8.62	128	4882	1.01	ng/ul	0.00
22) 2-Nitrophenol-d4	9.33	143	4564	0.96	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.87	165	8296	0.87	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.55	166	33515	1.21	ng/ul	0.00
47) Acenaphthylene-d8	13.80	160	35875	1.10	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	15.12	176	29341	1.22	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
71) Anthracene-d10	16.96	188	50214m>	1.39	ng/ul	> 0.00
79) Pyrene-d10	19.27	212	49718	1.26	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.04	264	47739	1.11	ng/ul	0.00

Target Compounds

				Ovalue	
40) 2,4,5-Trichlorophenol	12.62	196	8730	1.101	ng/ul# 97
56) 2,3,4,6-Tetrachlorophenol	14.76	232	10129m>	1.519	ng/ul> 97
69) Pentachlorophenol	16.53	266	241083	45.671	ng/ul 97

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