

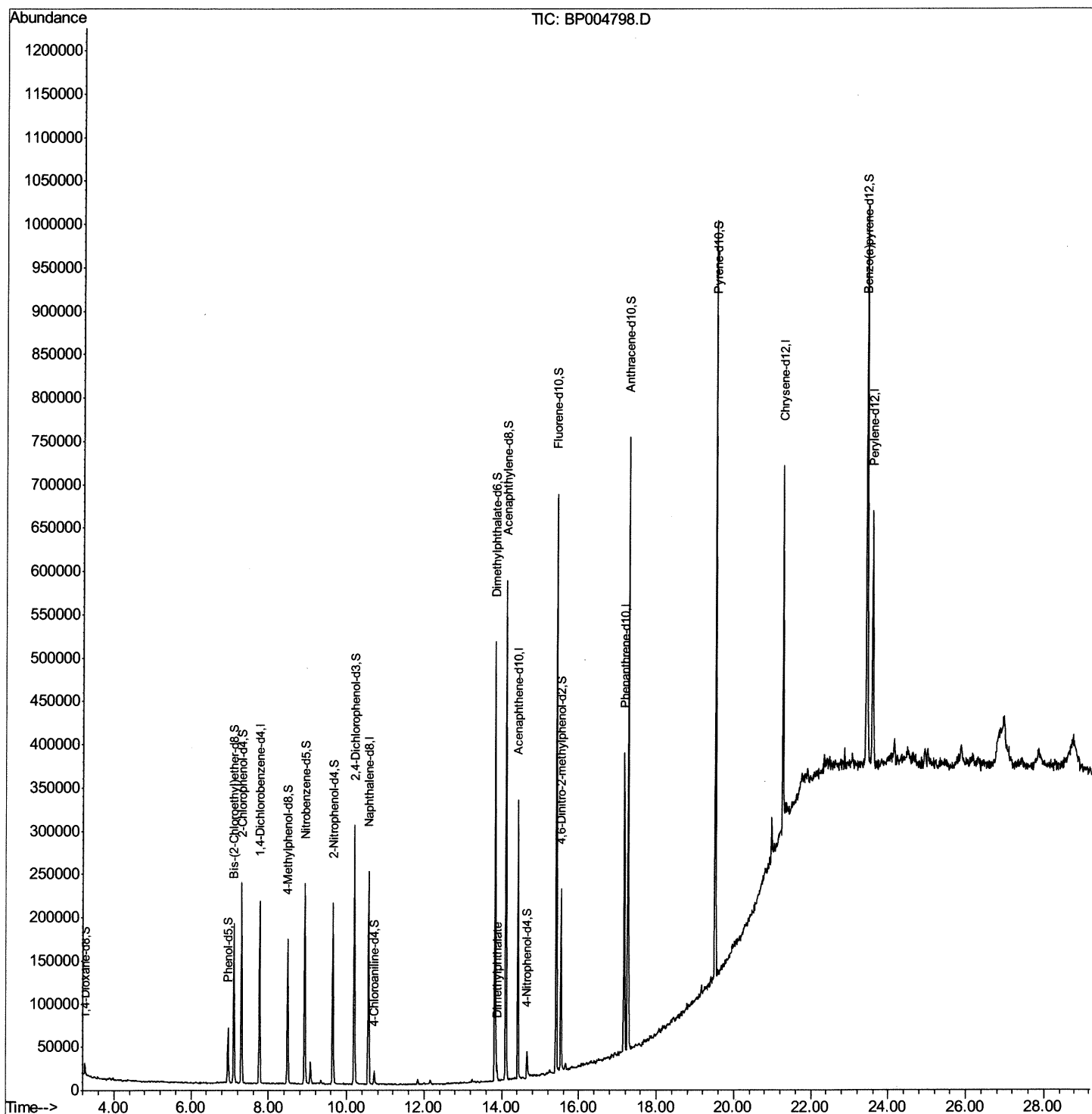
Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP021721\
Data File : BP004798.D
Acq On : 18 Feb 2021 05:48
Operator : CG/JU
Sample : M1379-17
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBGM2

Manual Integrations
APPROVED

mohammad
2/23/2021 3:42:03 PM

Quant Time: Feb 18 07:05:00 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Feb 18 03:00:36 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

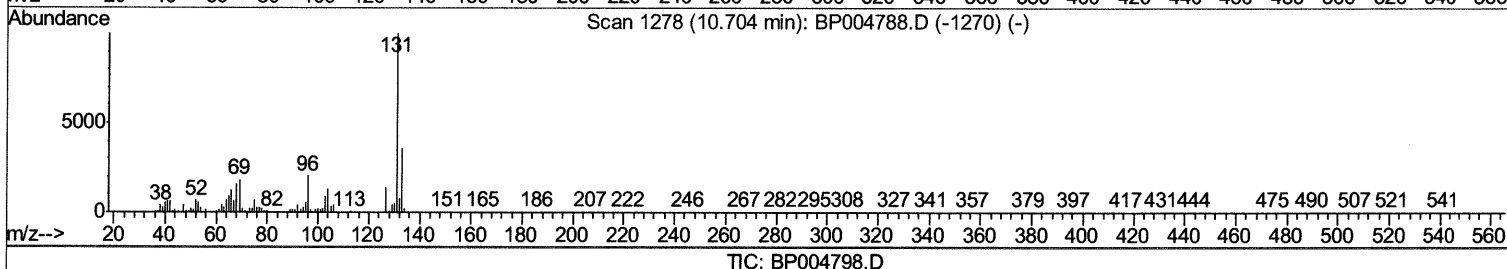
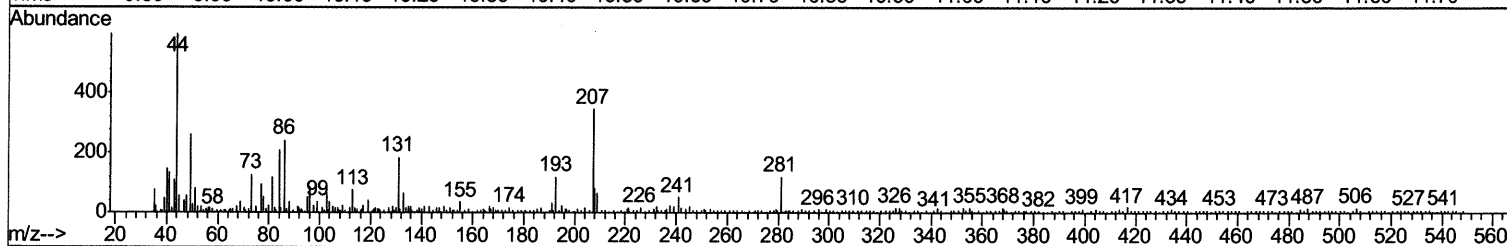
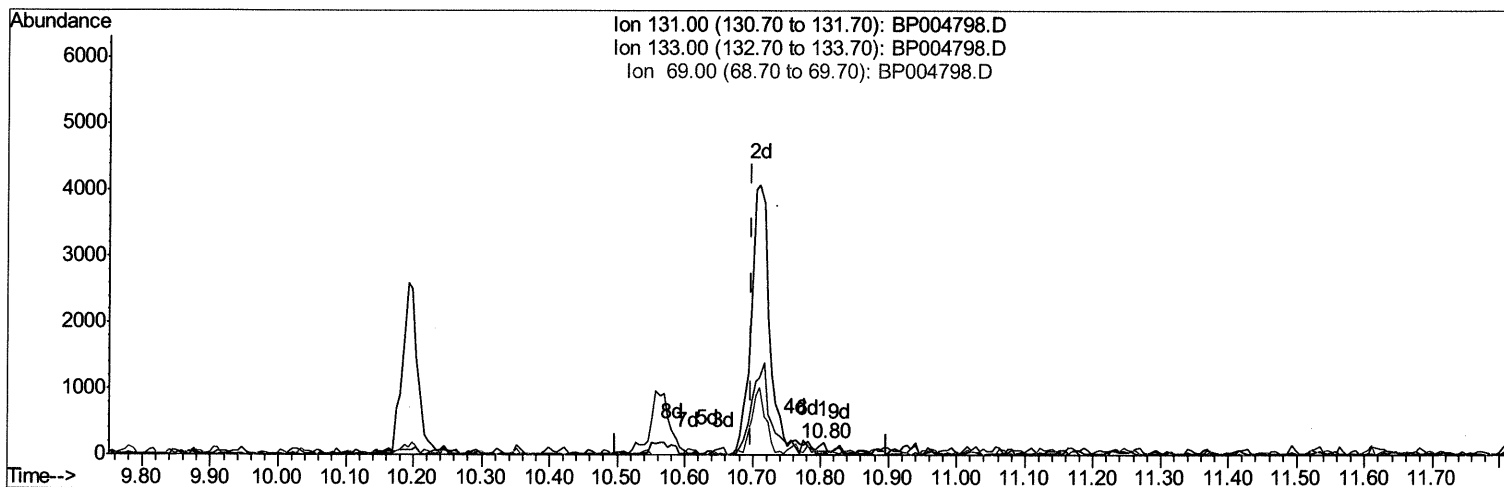
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Quant Time: Feb 18 06:56:39 2021
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(29) 4-Chloroaniline-d4 (S)
 10.804min (+0.106) 0.03ng/ul
 response 113

Ion	Exp%	Act%
131.00	100	100
133.00	31.20	32.16
69.00	20.50	18.59
0.00	0.00	0.00

Quantitation Report (Qedit)

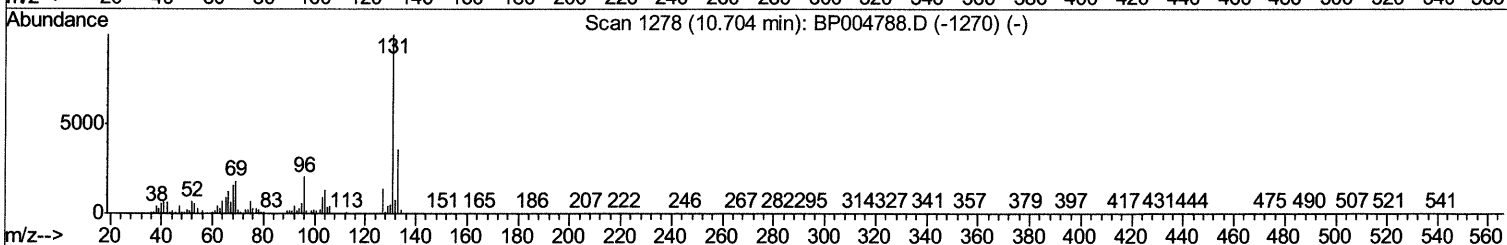
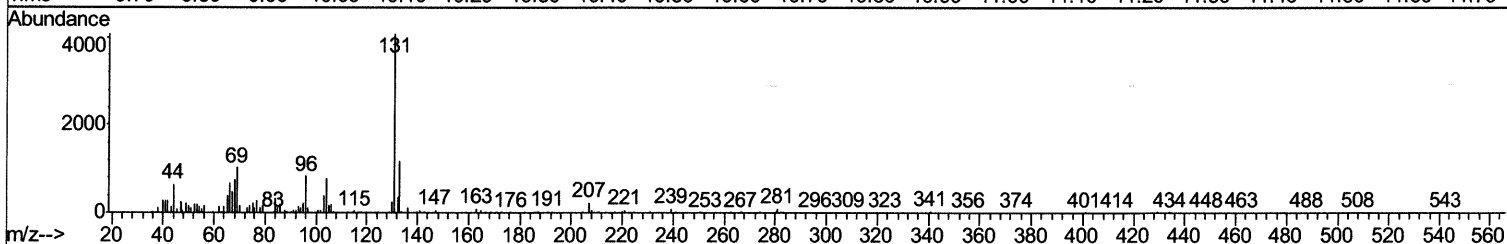
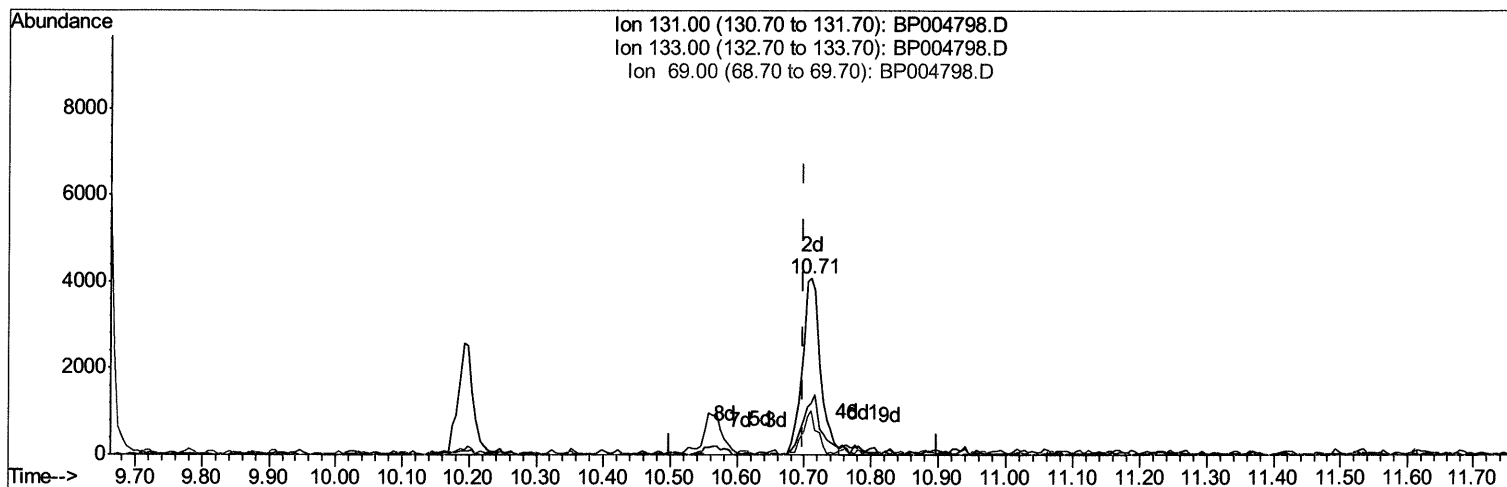
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Manual Integrations
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 2/23/2021 3:42:03 PM

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TIC: BP004798.D

(29) 4-Chloroaniline-d4 (S)

10.710min (+0.012) 2.31ng/ul m 02/22/21JU

response 7679

Ion	Exp%	Act%
131.00	100	100
133.00	31.20	28.87
69.00	20.50	25.19#
0.00	0.00	0.00

Quantitation Report (Qedit)

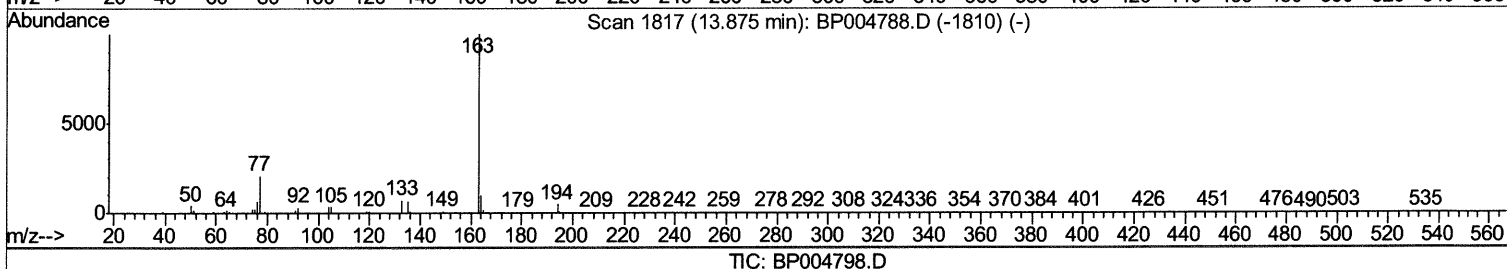
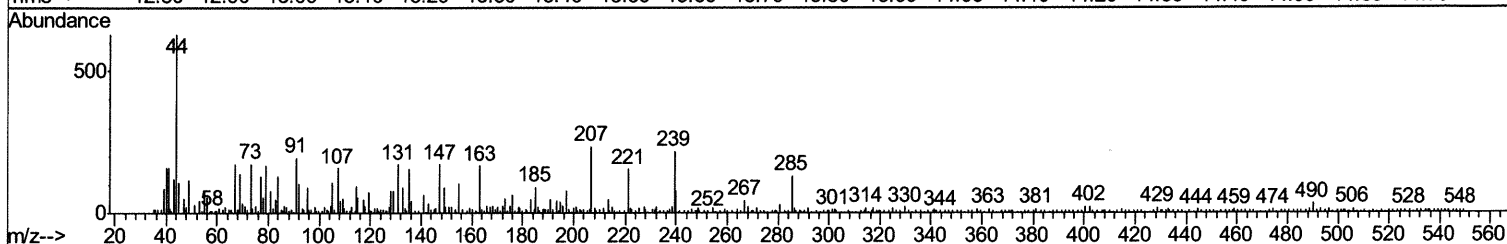
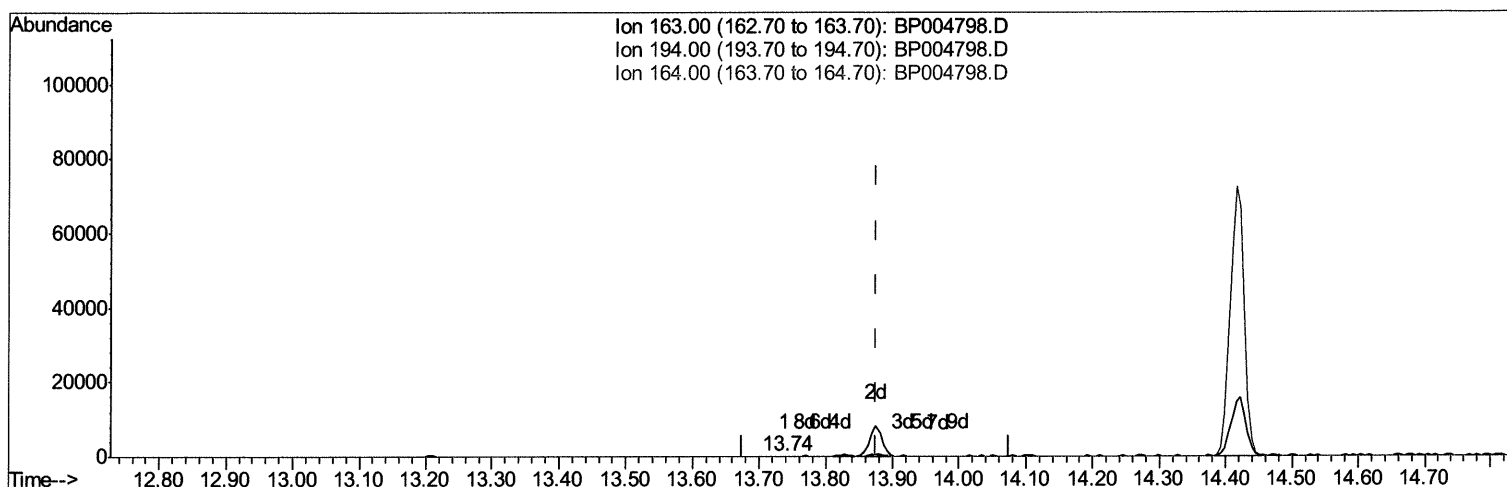
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Instrument :
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ClientSampleId :
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Manual Integrations
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(45) Dimethylphthalate

13.740min (-0.135) 0.01ng/ul

response 112

Ion	Exp%	Act%
163.00	100	100
194.00	4.70	4.19
164.00	10.00	8.98
0.00	0.00	0.00

Quantitation Report (Qedit)

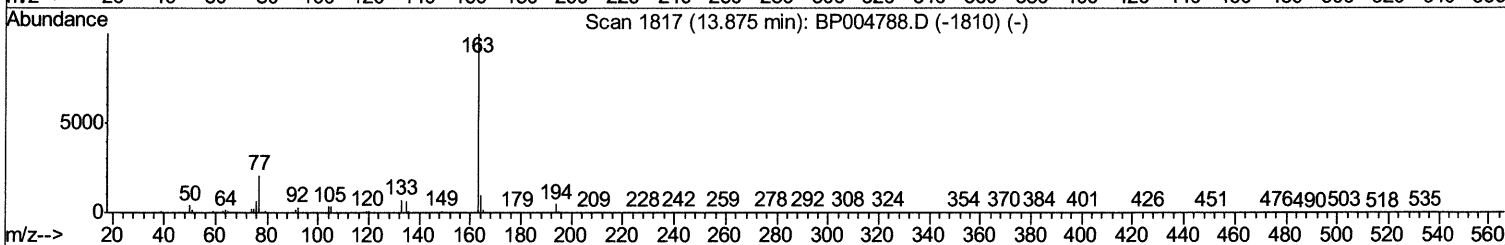
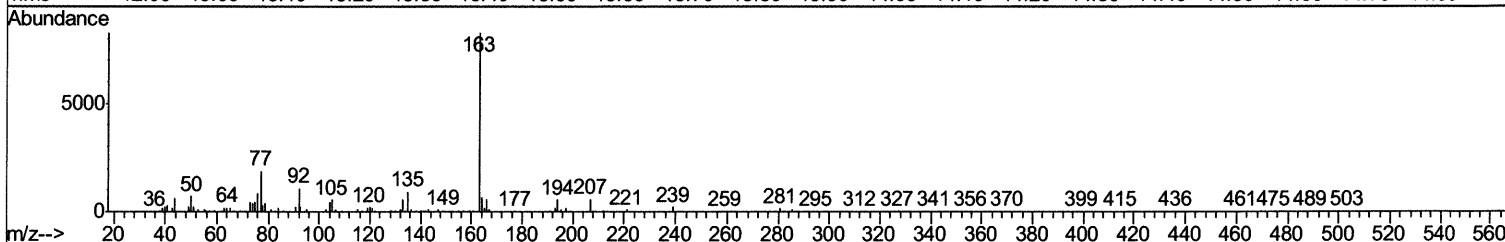
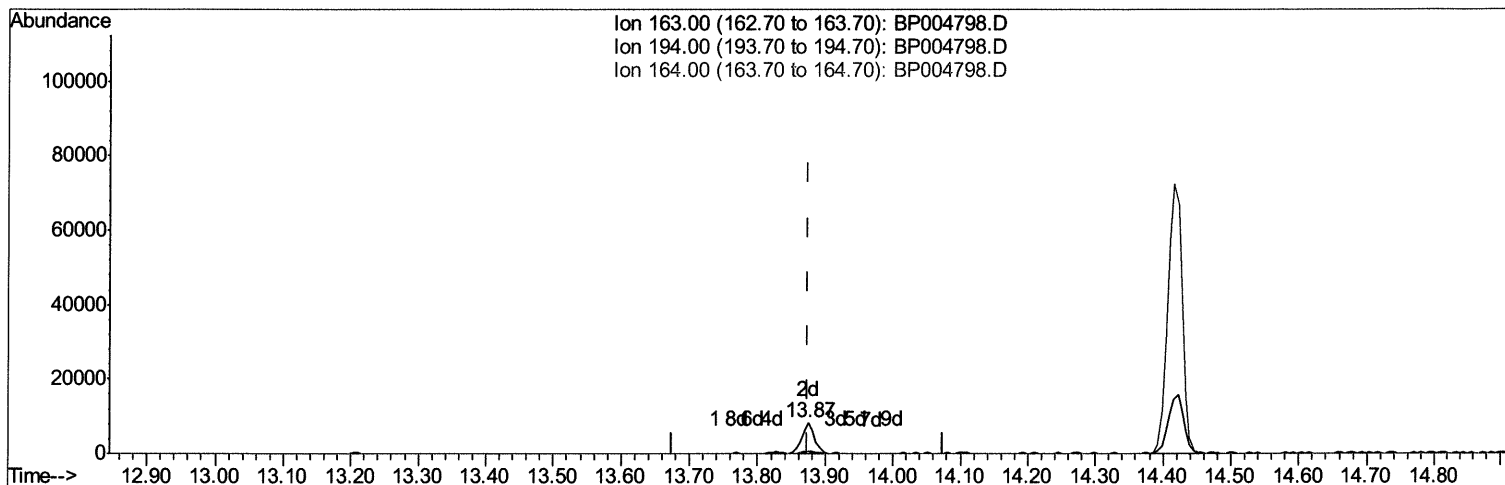
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Manual Integrations
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TIC: BP004798.D

(45) Dimethylphthalate

13.875min (-0.000) 1.31ng/ul m 02/22/2021

response 10476

Ion	Exp%	Act%
163.00	100	100
194.00	4.70	6.73#
164.00	10.00	8.34
0.00	0.00	0.00

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	52180	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	195468	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.42	164	107241	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	201105	20.00	ng/ul	0.00
78) Chrysene-d12	21.26	240	195248	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	215464	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	7192	5.33	ng/uL	0.00
5) Phenol-d5	6.96	99	33944	8.11	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	7.10	67	72671	33.41	ng/ul	0.00
9) 2-Chlorophenol-d4	7.31	132	99405	31.32	ng/ul	0.01
13) 4-Methylphenol-d8	8.49	113	59556	18.08	ng/ul	0.02
19) Nitrobenzene-d5	8.93	128	54513	40.39	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	60419	40.79	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	100897	33.85	ng/ul	0.01
29) 4-Chloroaniline-d4	10.71	131	7679m>	2.31	ng/ul>	0.01
44) Dimethylphthalate-d6	13.83	166	308005	39.57	ng/ul	0.00
47) Acenaphthylene-d8	14.11	160	383520	40.50	ng/ul	0.00
52) 4-Nitrophenol-d4	14.65	143	8681	6.38	ng/ul	0.04
58) Fluorene-d10	15.42	176	252834	37.90	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	42619	35.06	ng/ul	0.00
71) Anthracene-d10	17.27	188	390782	44.06	ng/ul	0.00
79) Pyrene-d10	19.51	212	418576	45.80	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.42	264	483818	44.39	ng/ul	0.00

02/22/21 JU

Target Compounds

				Ovalue
45) Dimethylphthalate	13.87	163	10476m>	1.305 ng/ul>

02/22/21 JU

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