

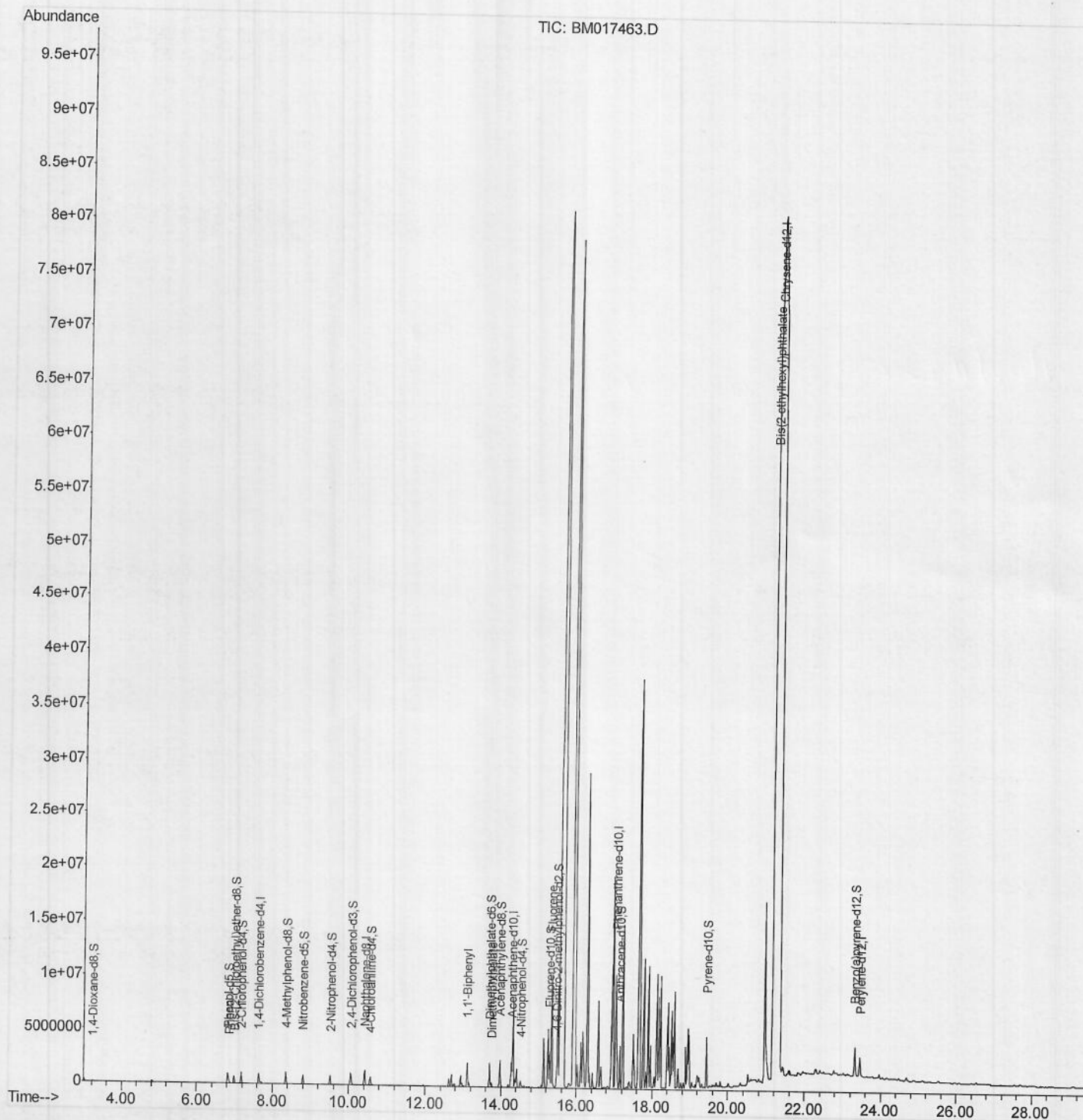
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110118\
 Quantitation Report (QT Reviewed)
 Data File : BM017463.D
 Acq On : 01 Nov 2018 13:43
 Operator : MJ/SJ
 Sample : J5704-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A40F8

Manual Integrations
 APPROVED

Sohil
 11/2/2018 4:12:06 PM

Quant Time: Nov 02 02:28:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 02 01:56:23 2018
 Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110118\
Quantitation Report (Qedit)

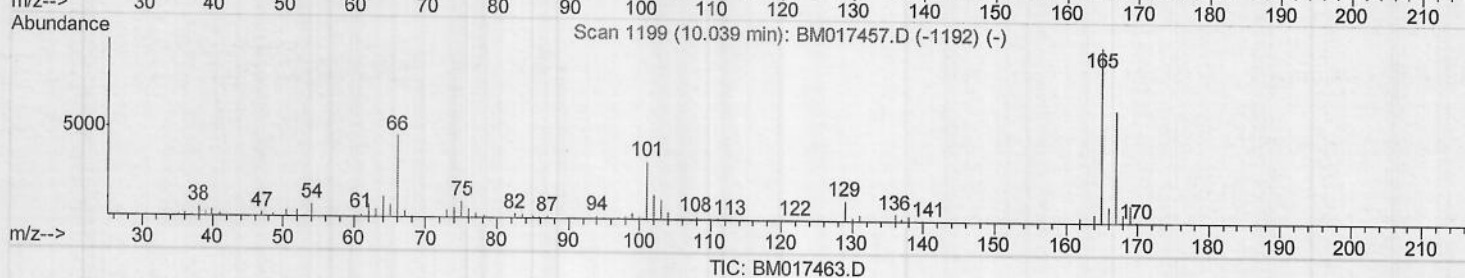
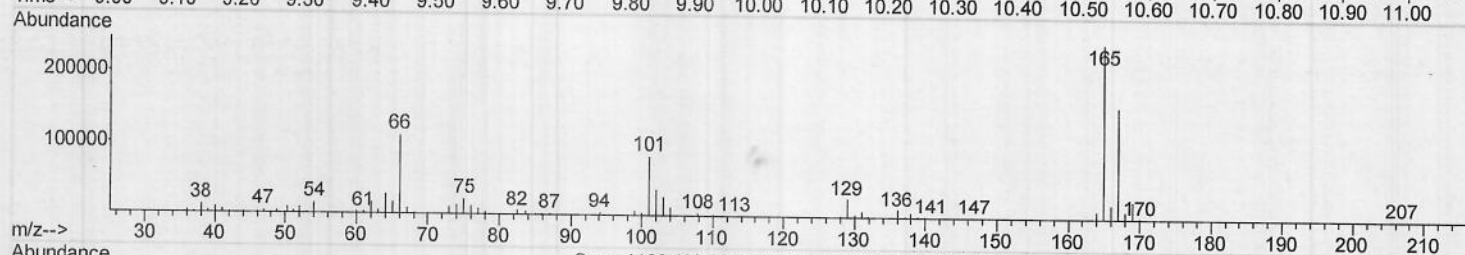
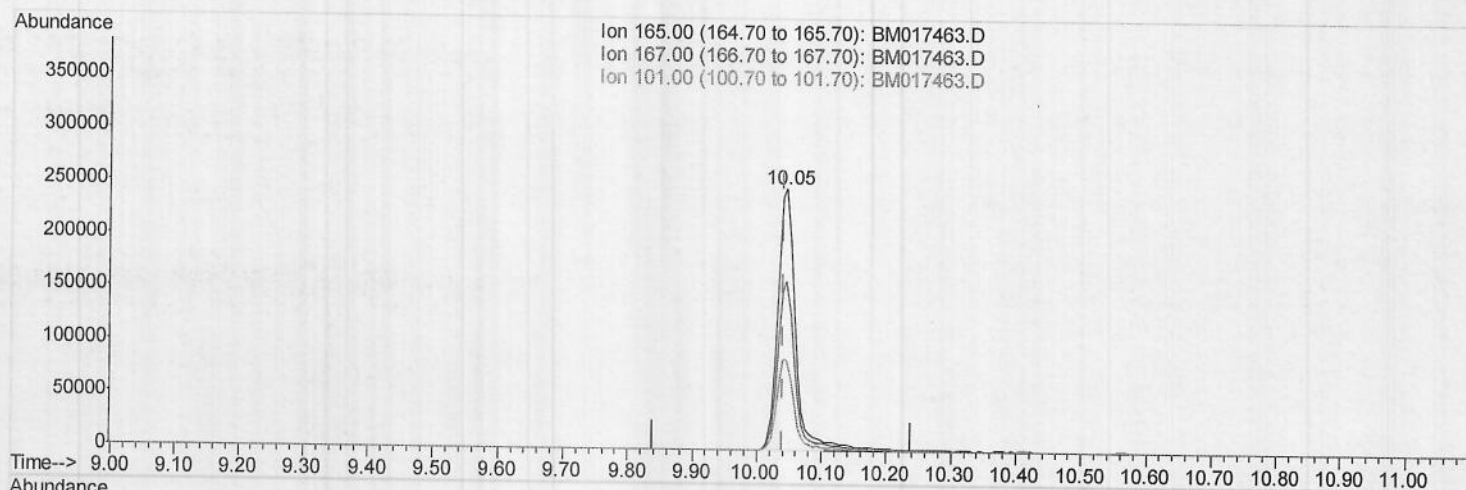
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(26) 2,4-Dichlorophenol-d3 (S)

10.045min (+0.006) 24.35ng/ul

response 452732

Ion	Exp%	Act%
165.00	100	100
167.00	63.60	64.36
101.00	34.30	34.52
0.00	0.00	0.00

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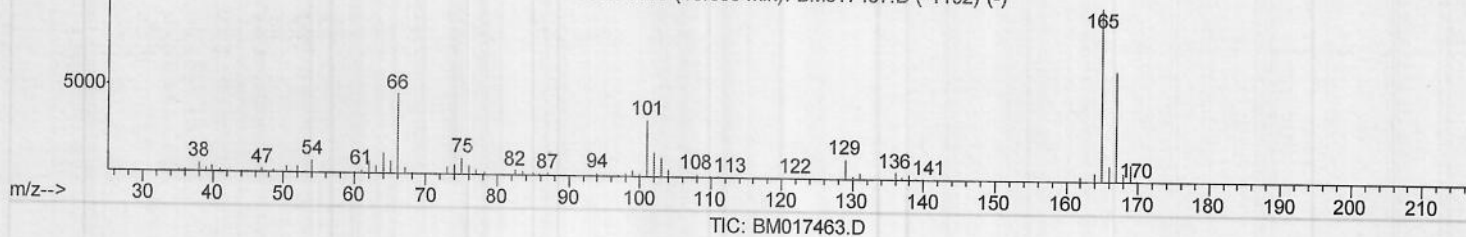
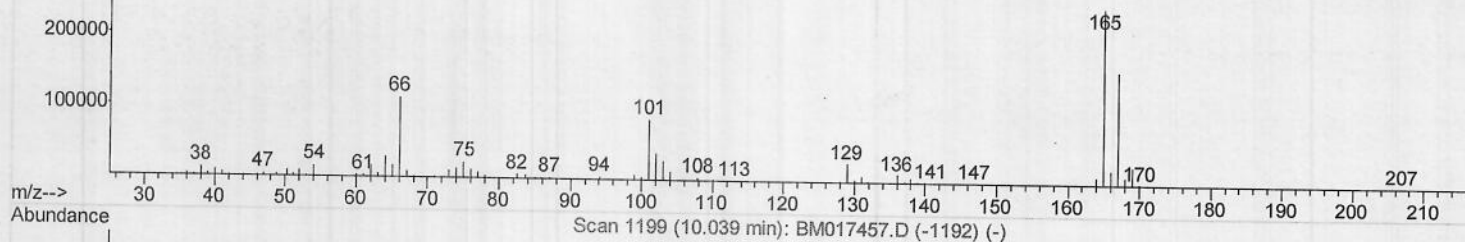
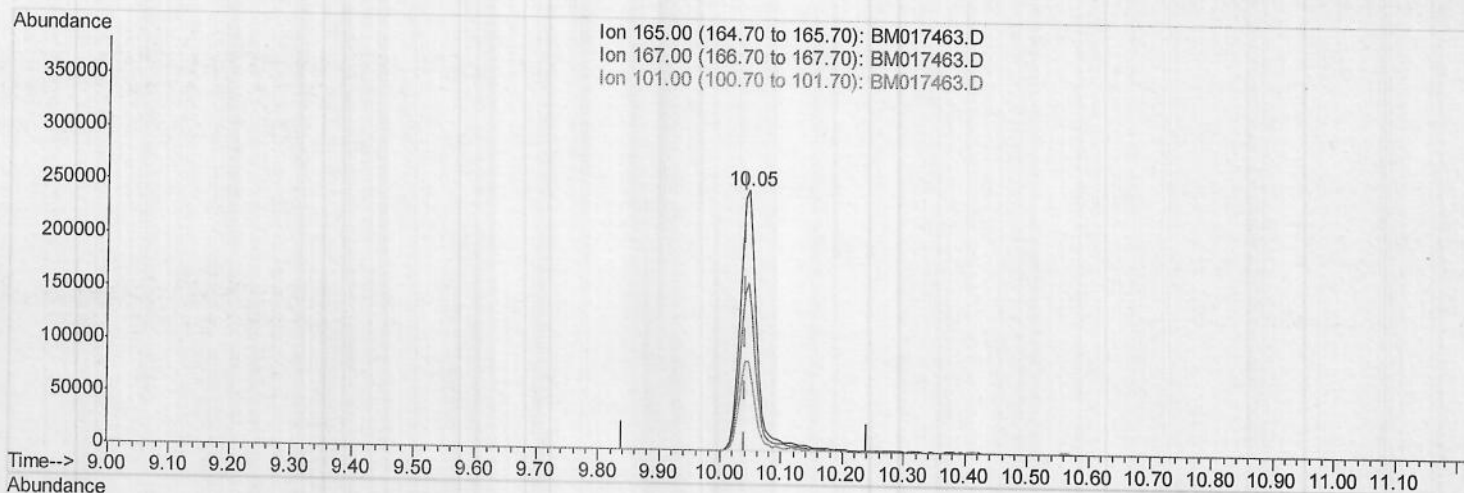
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(26) 2,4-Dichlorophenol-d3 (S)

10.045min (+0.006) 25.81ng/ul m

response 479835

Ion	Exp%	Act%
165.00	100	100
167.00	63.60	64.36
101.00	34.30	34.52
0.00	0.00	0.00

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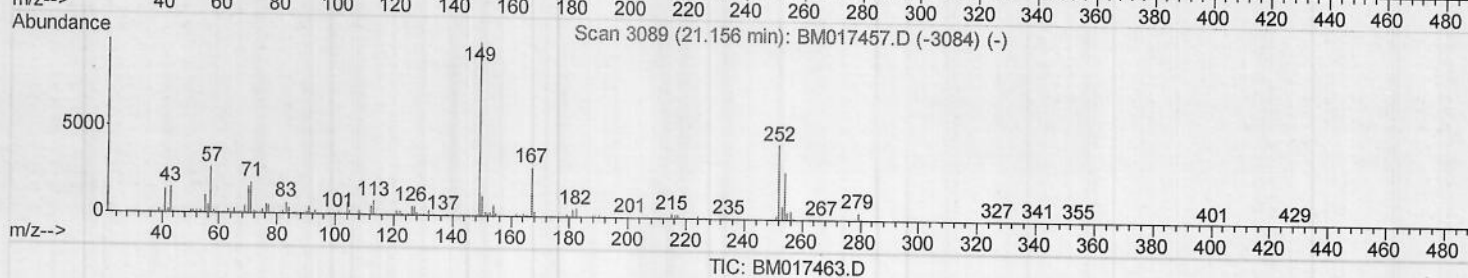
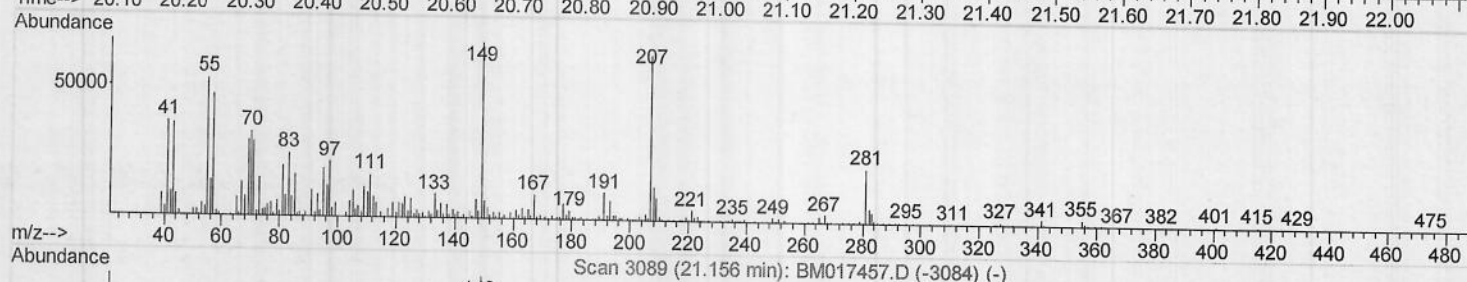
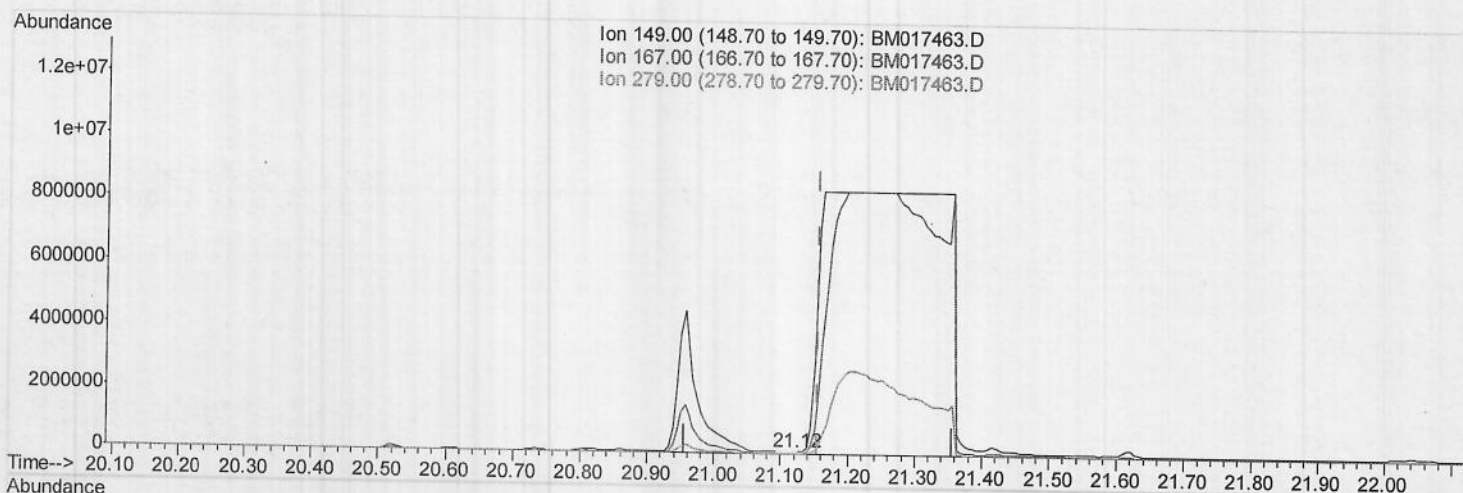
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(84) Bis(2-ethylhexyl)phthalate

21.121min (-0.035) 0.80ng/ul

response 9828

Ion	Exp%	Act%
149.00	100	100
167.00	28.80	14.23#
279.00	5.10	1.67#
0.00	0.00	0.00

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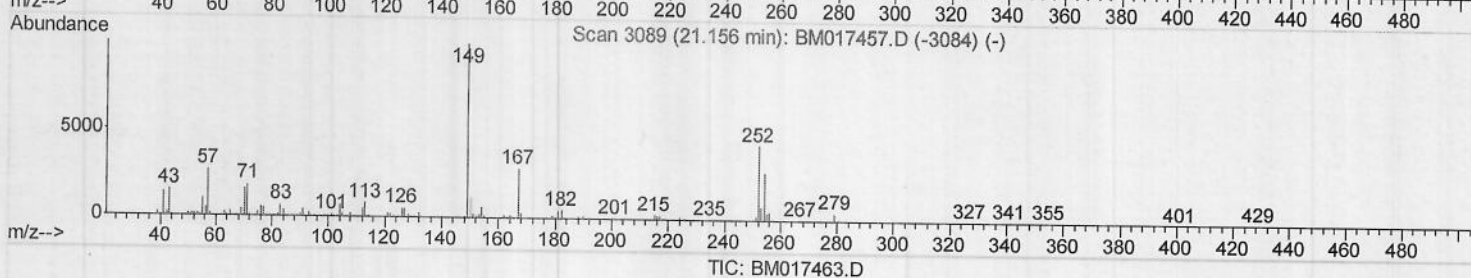
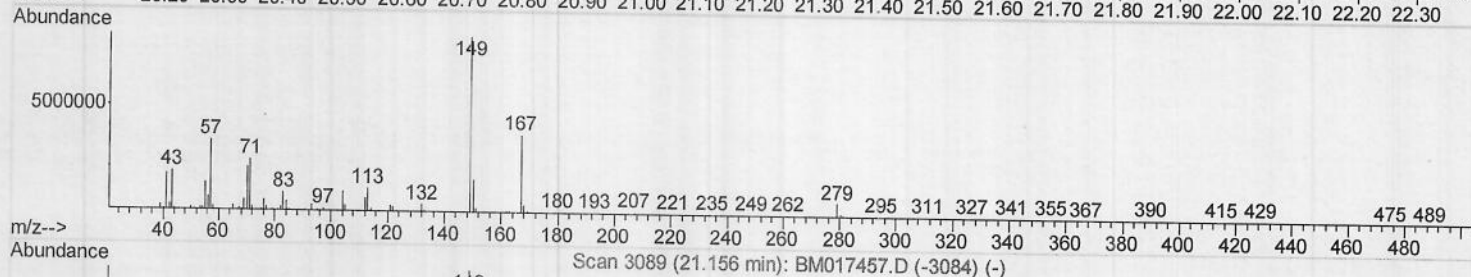
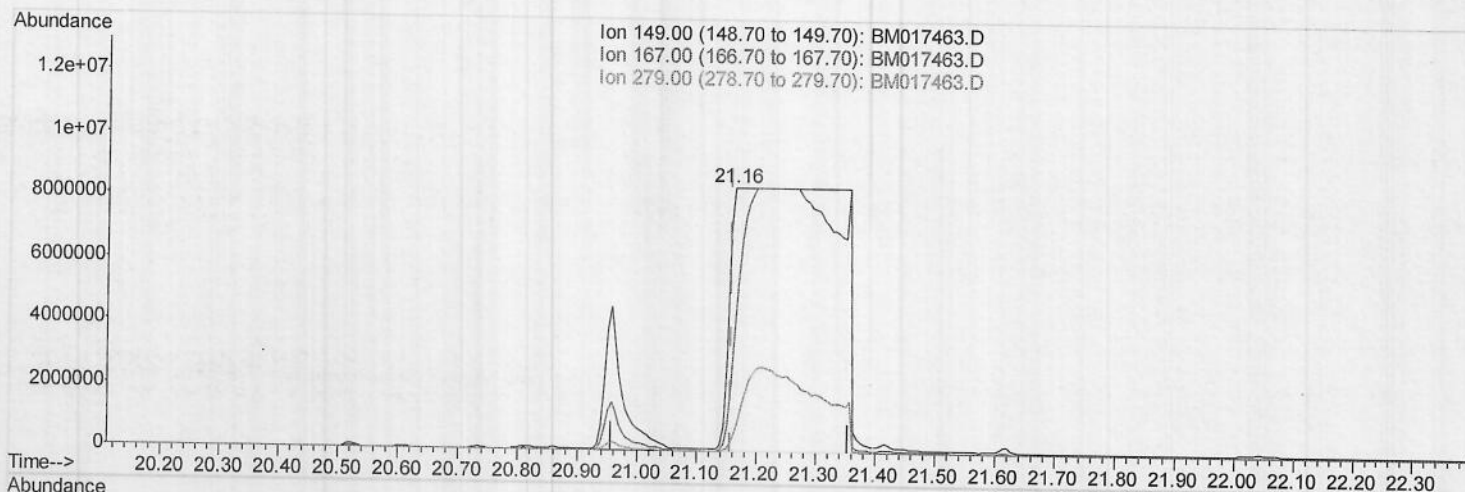
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(84) Bis(2-ethylhexyl)phthalate

21.162min (+0.006) 8629.51ng/ul m > 551106118

response 105537897

Ion	Exp%	Act%
149.00	100	100
167.00	28.80	43.90#
279.00	5.10	7.44#
0.00	0.00	0.00

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Internal Standards	R.T.	QI on	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	254307	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	1142001	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	666845	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	1896196	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	361015	20.00	ng/ul	0.00
86) Perylene-d12	23.46	264	1227888	20.00	ng/ul	0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	18915	3.42	ng/uL	0.00
5) Phenol-d5	6.82	99	565470	24.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	325118	24.00	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	455366	25.28	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	448614	24.42	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	220845	24.74	ng/ul	0.00
22) 2-Nitrophenol-d4	9.51	143	263245	26.29	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.05	165	479835m)	25.81	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	449248	29.60	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	1429583	25.00	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	1809780	26.06	ng/ul	0.00
52) 4-Nitrophenol-d4	14.52	143	192382	18.01	ng/ul	0.02
58) Fluorene-d10	15.29	176	1279030	25.75	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.47	200	189473	14.34	ng/ul	0.06
71) Anthracene-d10	17.14	188	2177816	23.20	ng/ul	0.01
79) Pyrene-d10	19.44	212	2323919	138.95	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.32	264	1763526	27.20	ng/ul	0.02

) SJ 11/06/18

Target Compounds

					Qvalue
6) Phenol	6.85	94	134186	5.817	ng/ul 98
41) 1,1'-Biphenyl	13.11	154	644798	11.688	ng/ul 99
45) Dimethylphthalate	13.75	163	264214	4.769	ng/ul 99
59) Fluorene	15.34	166	2020700	37.086	ng/ul# 29
84) Bis(2-ethylhexyl)phthalate	21.16	149	105537897m)	8629.506	ng/ul

) SJ 11/06/18

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