

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

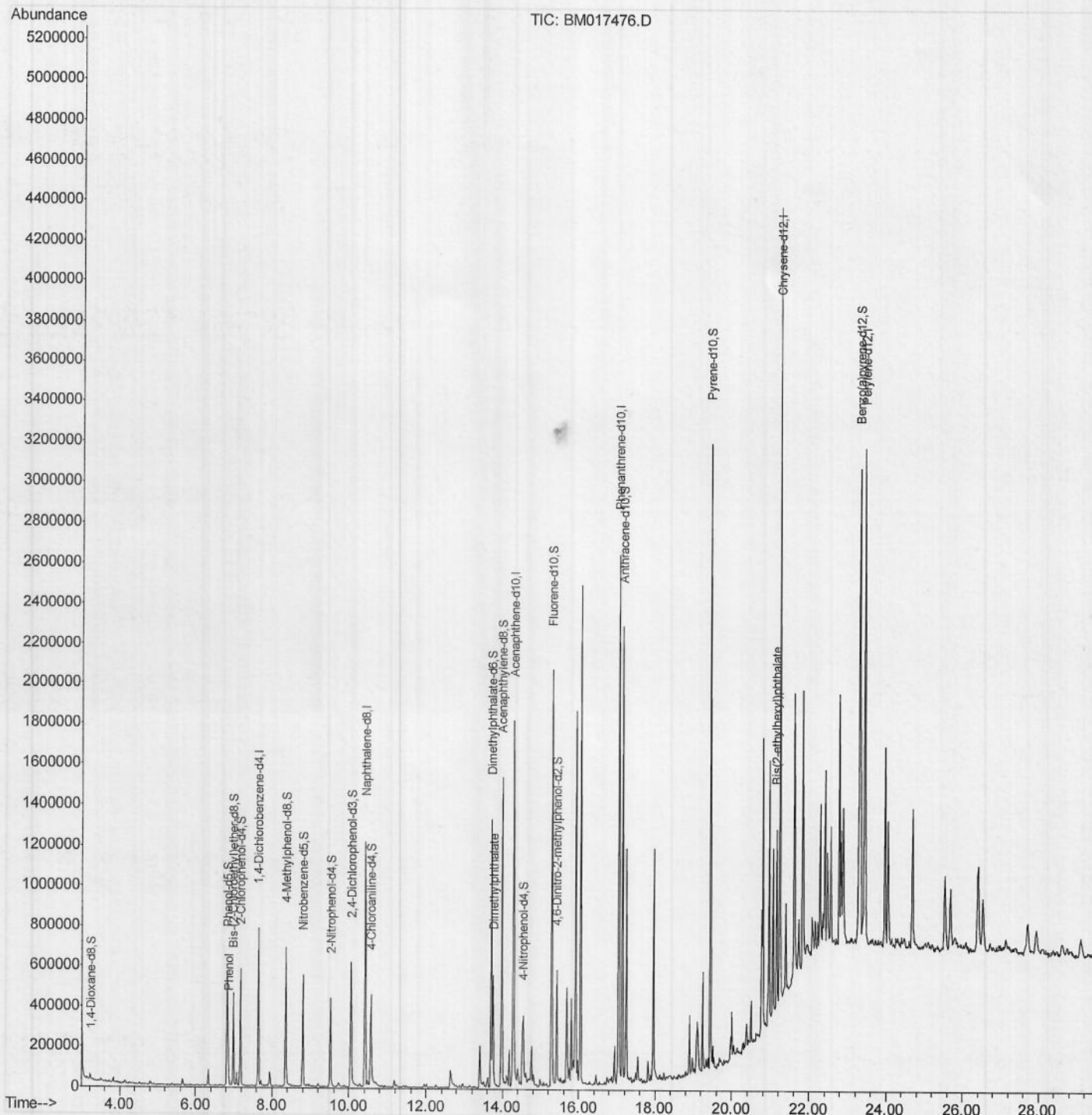
Data File : BM017476.D
Acq On : 01 Nov 2018 21:37
Operator : MJ/SJ
Sample : J5701-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
Client Sampled :
A40F0

Quant Time: Nov 02 04:06:21 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 02 02:41:22 2018
Response via : Initial Calibration

Manual Integrations
APPROVED

Sohil
11/2/2018 4:12:18 PM



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

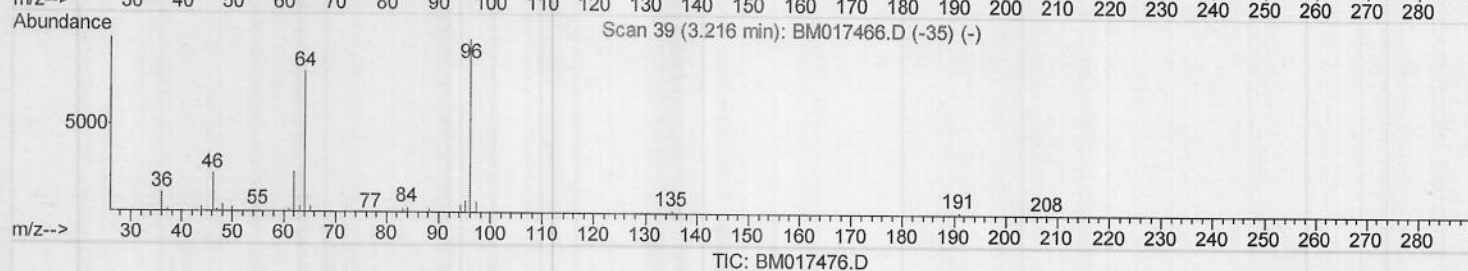
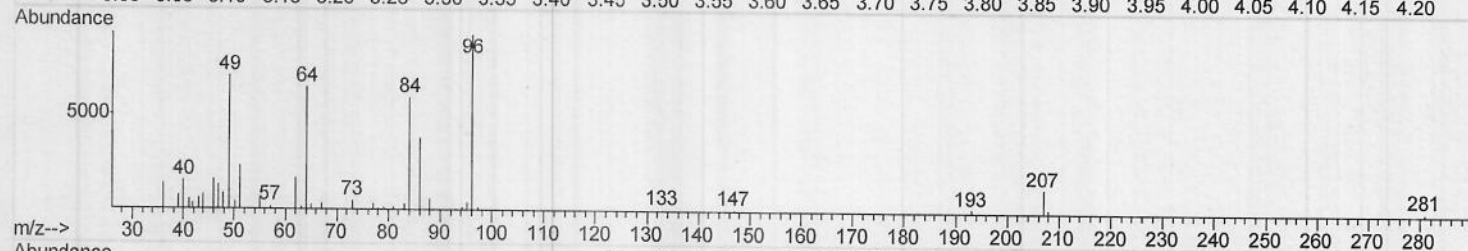
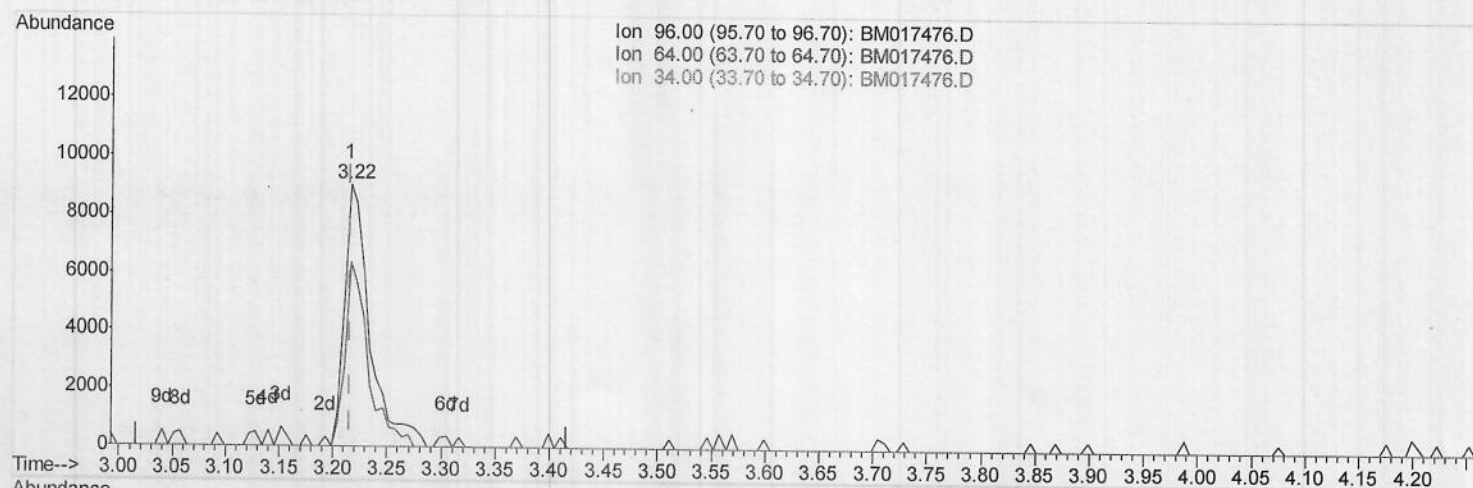
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Quant Time: Nov 02 02:45:41 2018
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(3) 1,4-Dioxane-d8 (S)

3.216min (+0.000) 3.01ng/uL

response 13711

Ion	Exp%	Act%
96.00	100	100
64.00	78.00	70.67
34.00	0.00	0.00
0.00	0.00	0.00

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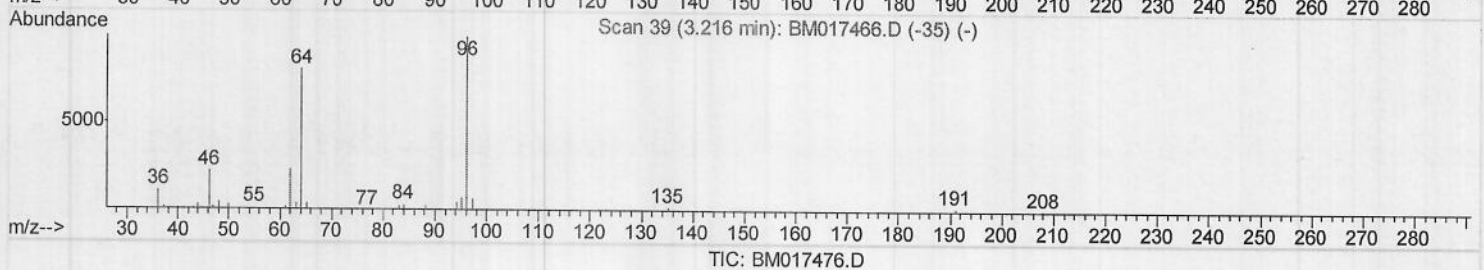
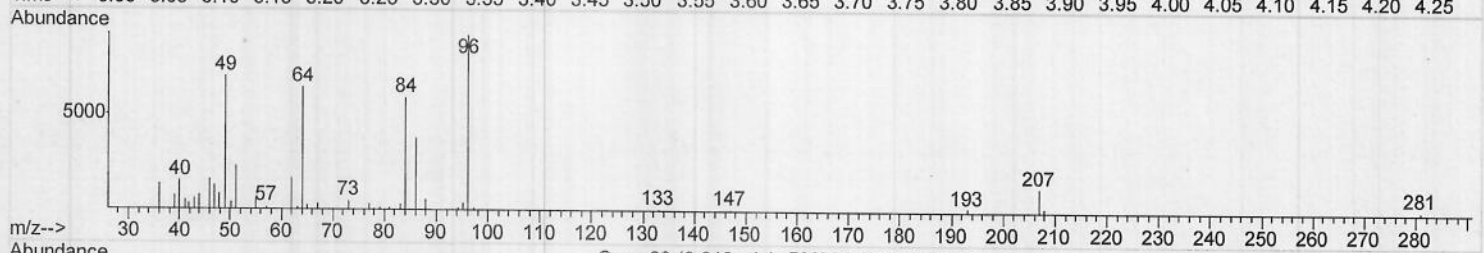
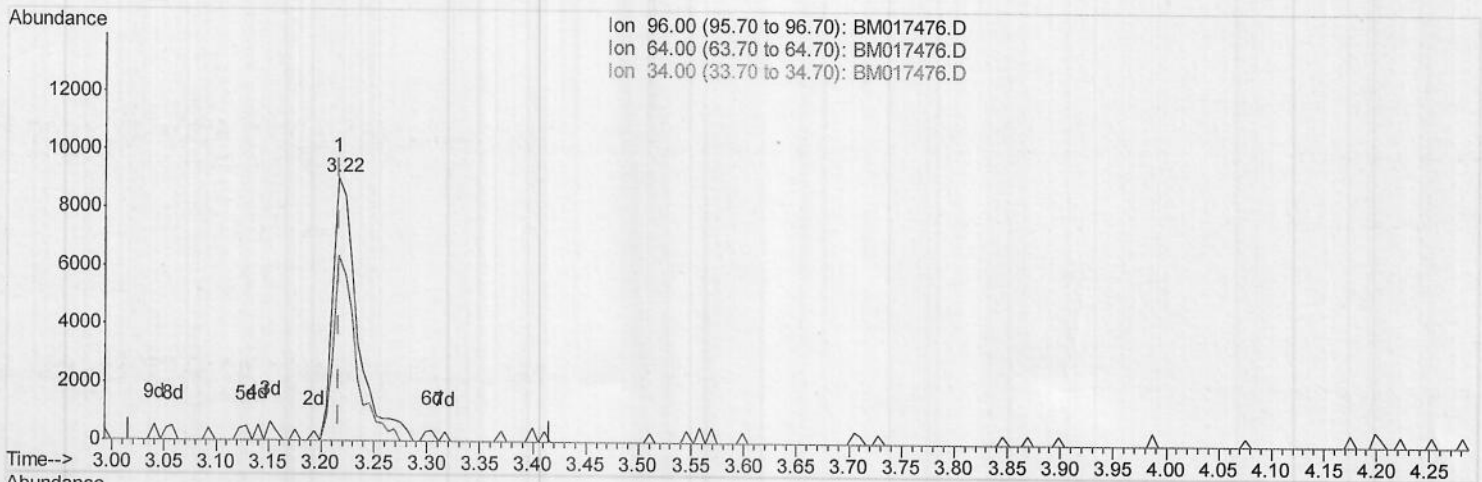
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(3) 1,4-Dioxane-d8 (S)

3.216min (+0.000) 3.20ng/uL m J5701/06/18

response 14601

Ion	Exp%	Act%
96.00	100	100
64.00	78.00	70.67
34.00	0.00	0.00
0.00	0.00	0.00

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Quantitation Report (Qedit)

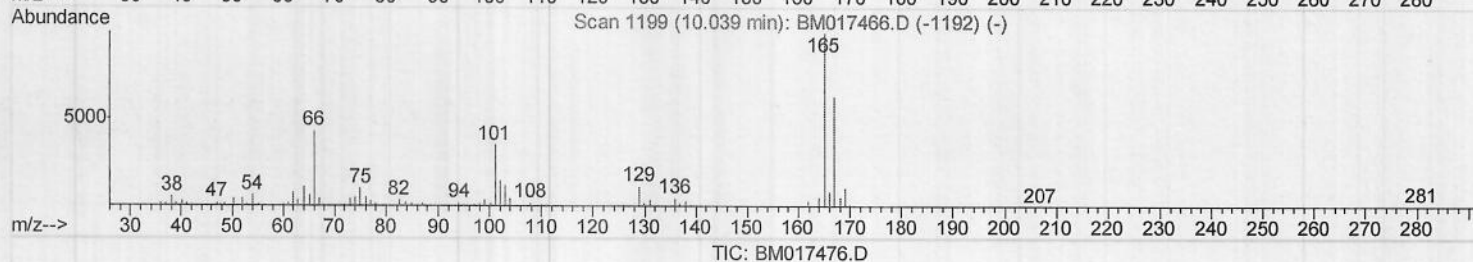
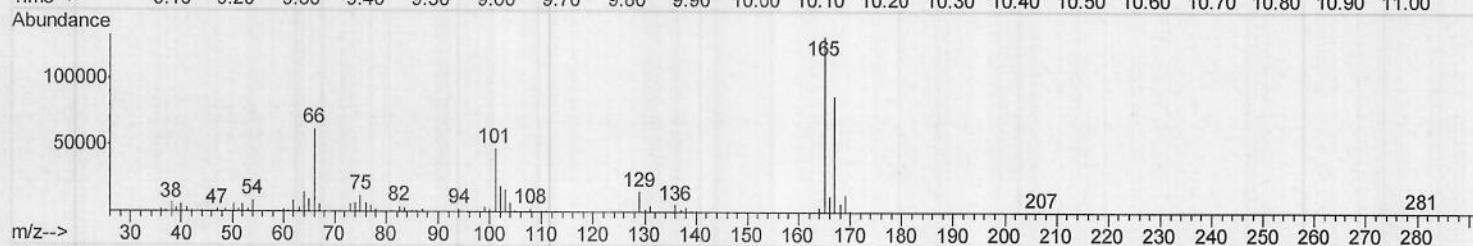
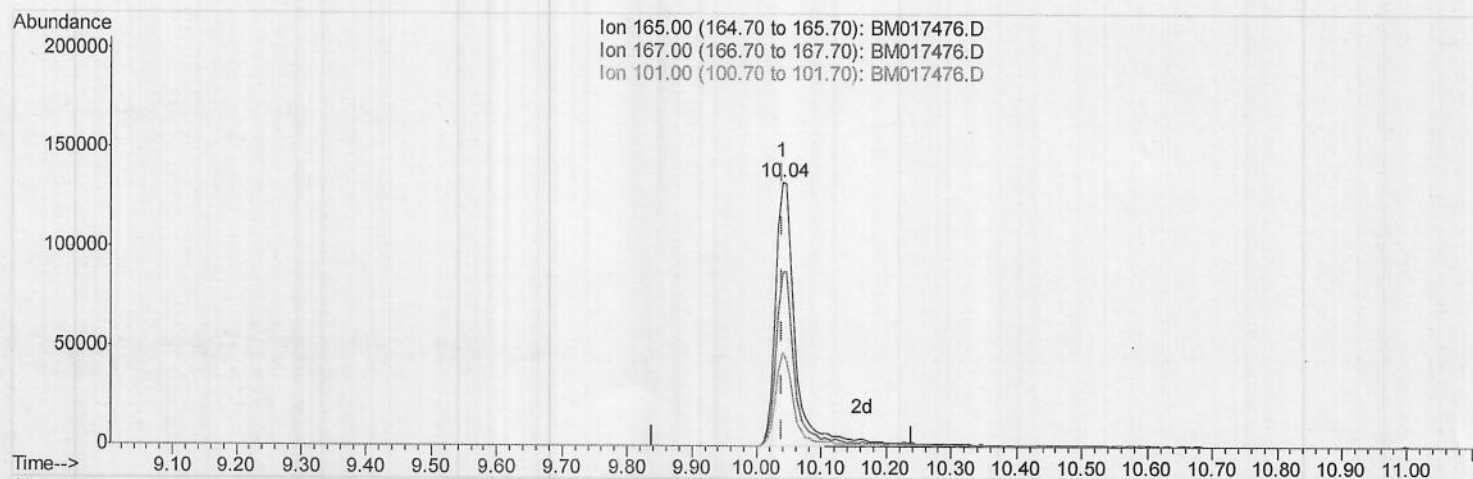
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(26) 2,4-Dichlorophenol-d3 (S)

10.039min (+0.000) 15.95ng/ul

response 261461

Ion	Exp%	Act%
165.00	100	100
167.00	63.60	65.75
101.00	34.30	36.09
0.00	0.00	0.00

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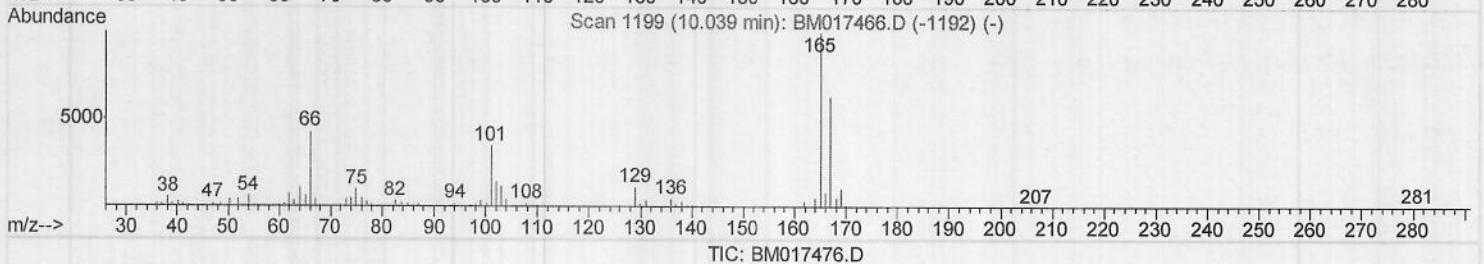
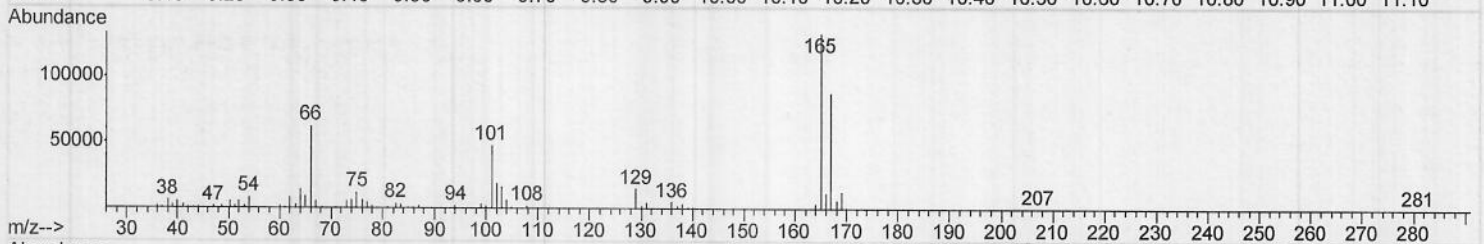
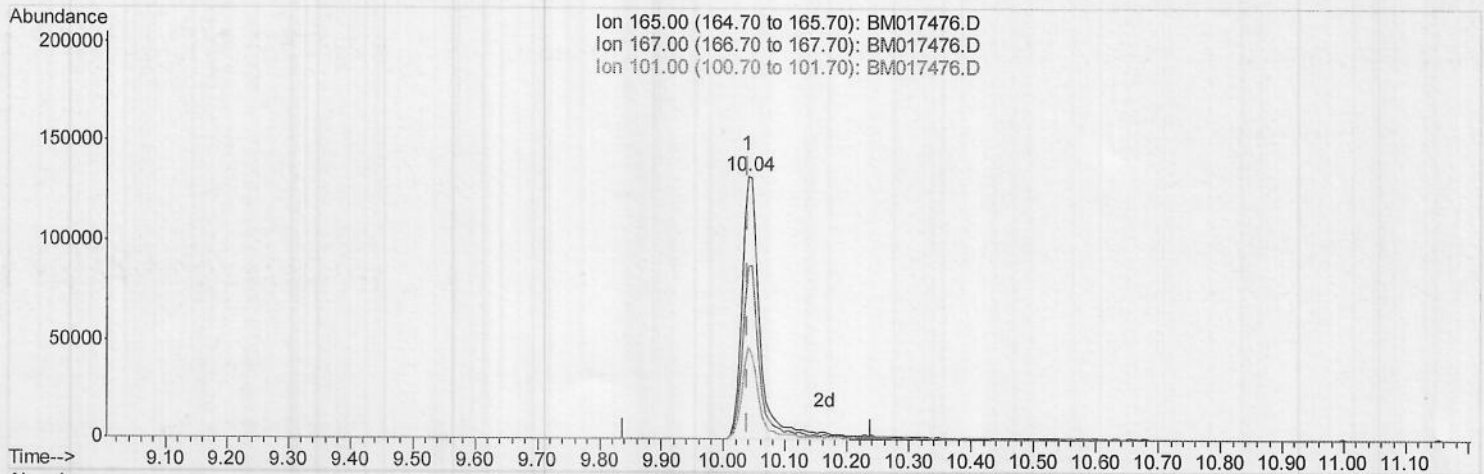
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Instrument :
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Manual Integrations
APPROVED

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

10.039min (+0.000) 17.20ng/ul m) 5511/06/18

response 281988

Ion	Exp%	Act%
165.00	100	100
167.00	63.60	65.75
101.00	34.30	36.09
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	7.64	152	209708	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1006781	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.28	164	626008	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1514210	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	1786222	20.00	ng/ul	0.00
86) Perylene-d12	23.44	264	1760189	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	14601m)	3.20	ng/uL	0.00
5) Phenol-d5	6.82	99	337950	17.64	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	195617	17.51	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	273954	18.45	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	275679	18.20	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	136684	17.37	ng/ul	0.00
22) 2-Nitrophenol-d4	9.51	143	158804	17.99	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	281988m)	17.20	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	306232	22.89	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	950106	17.70	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	1190377	18.26	ng/ul	0.00
52) 4-Nitrophenol-d4	14.52	143	144986	14.46	ng/ul	0.00
58) Fluorene-d10	15.28	176	869198	18.64	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.41	200	129527	12.28	ng/ul	0.00
71) Anthracene-d10	17.13	188	1334394	17.80	ng/ul	0.00
79) Pyrene-d10	19.43	212	1551191	18.74	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.30	264	1662535	17.88	ng/ul	0.00

SS11/06/18

Target Compounds

					Qvalue
6) Phenol	6.85	94	107435	5.648 ng/ul	94
45) Dimethylphthalate	13.75	163	372326	7.158 ng/ul	100
84) Bis(2-ethylhexyl)phthalate	21.16	149	102237	1.690 ng/ul	96

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