

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110518\
Quantitation Report (QT Reviewed)

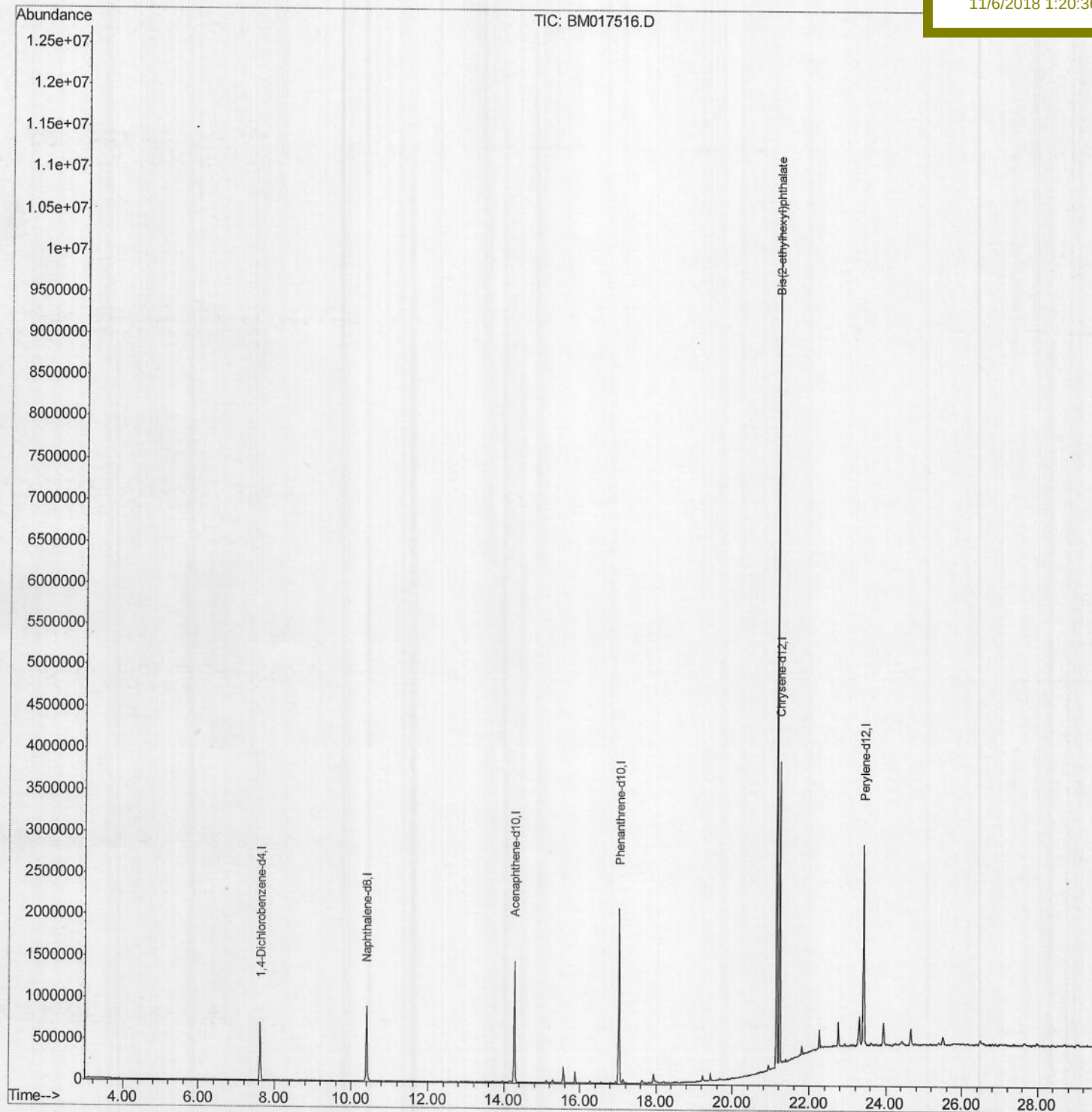
Data File : BM017516.D
Acq On : 05 Nov 2018 15:17
Operator : MJ/SJ
Sample : J5701-11DL 30X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Nov 06 01:04:43 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102418MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 06 00:34:52 2018
Response via : Initial Calibration

Instrument :
BNA_M
Client Sampled :
A40F7DL

Manual Integrations
APPROVED

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11/6/2018 1:20:36 PM



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Quantitation Report (Qedit)

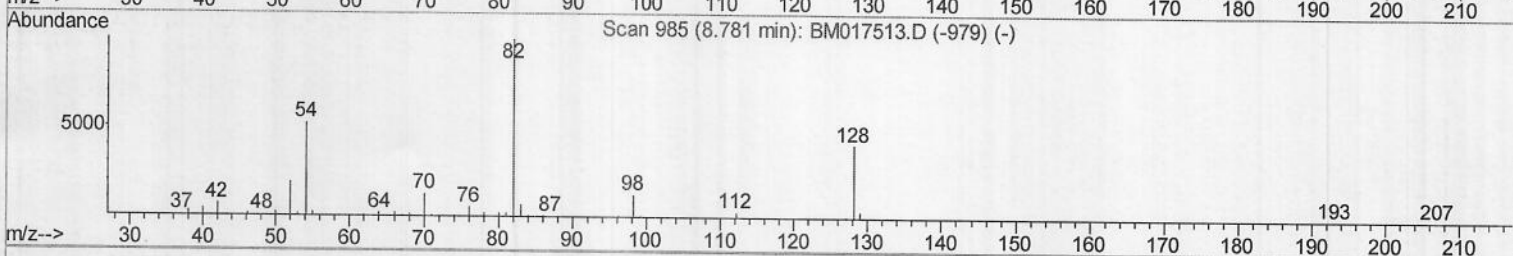
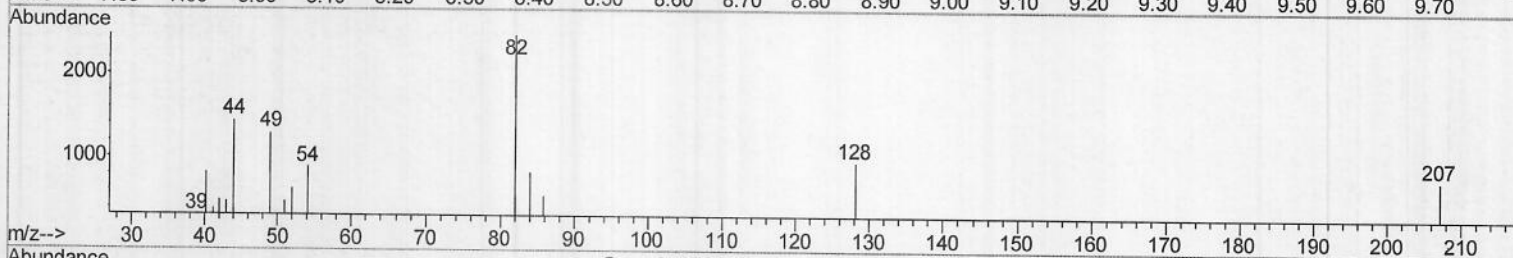
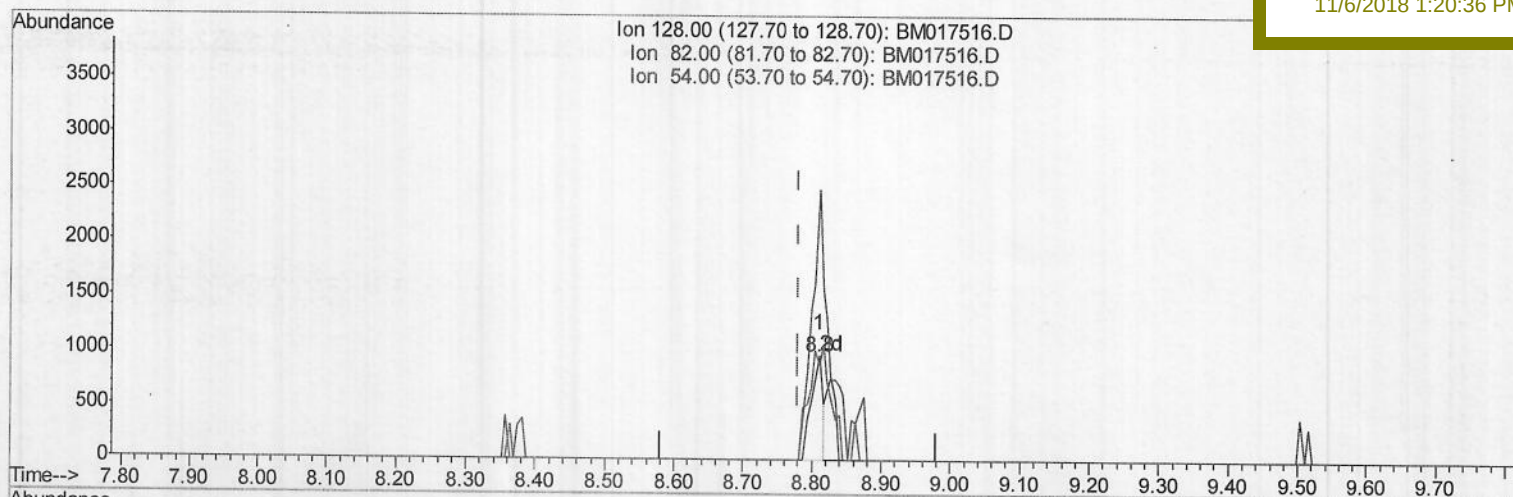
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TIC: BM017516.D

(19) Nitrobenzene-d5 (S)

8.810min (+0.029) 0.17ng/ul

response 1107

Ion	Exp%	Act%
128.00	100	100
82.00	237.30	256.52
54.00	117.80	91.41#
0.00	0.00	0.00

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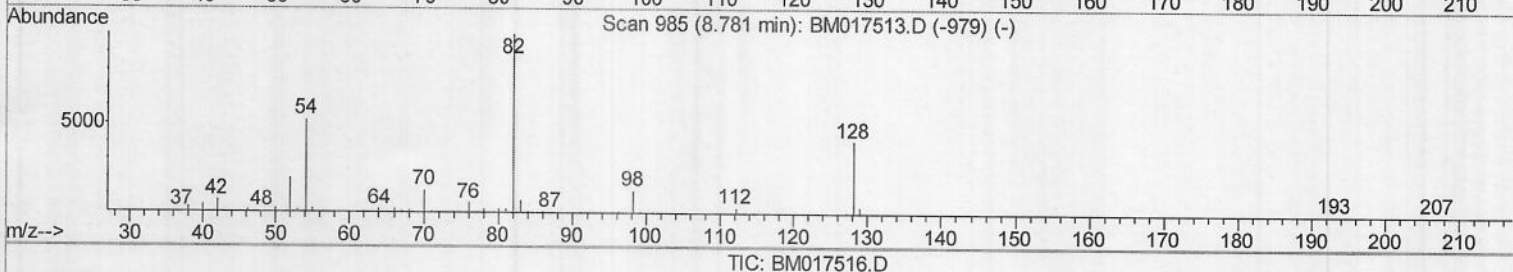
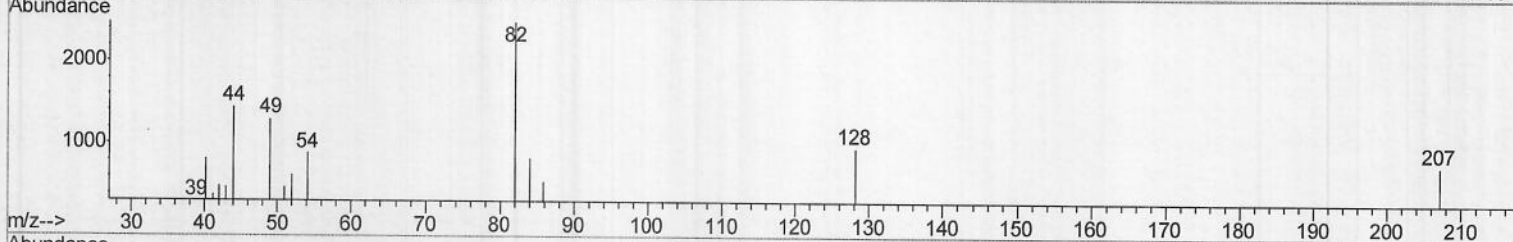
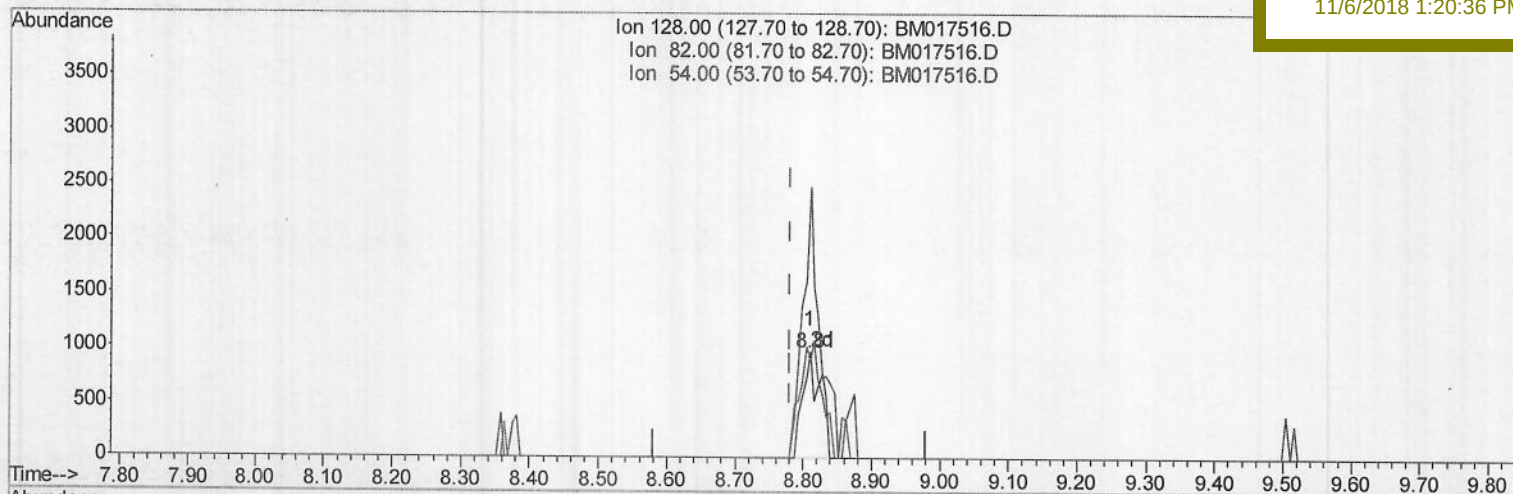
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TIC: BM017516.D

(19) Nitrobenzene-d5 (S)

8.810min (+0.029) 0.28ng/ul m

response 1803

Ion	Exp%	Act%
128.00	100	100
82.00	237.30	256.52
54.00	117.80	91.41#
0.00	0.00	0.00

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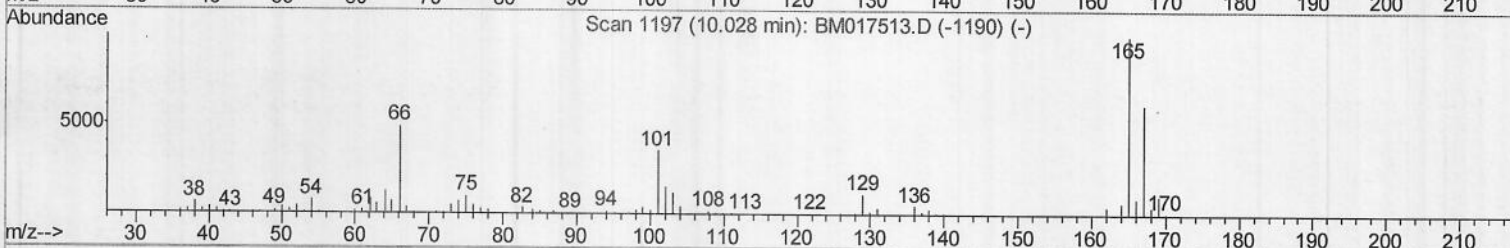
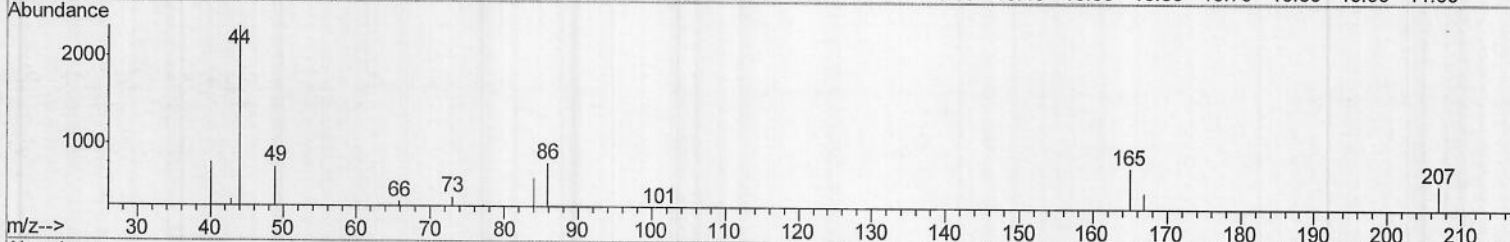
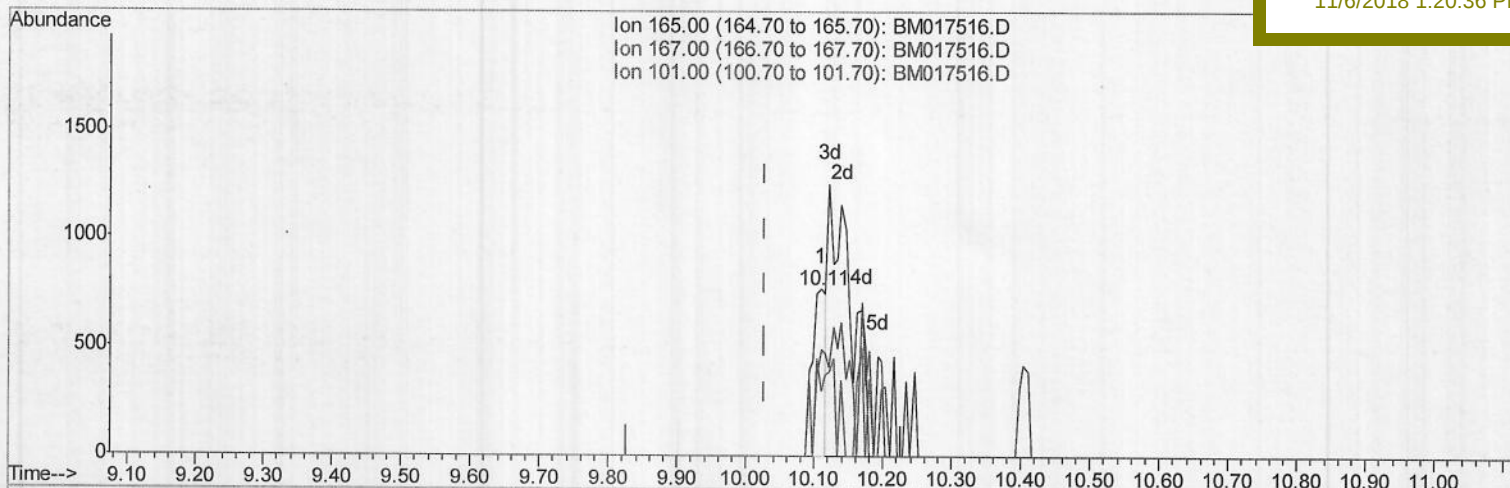
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TIC: BM017516.D

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

10.110min (+0.082) 0.08ng/ul

response 1087

Ion	Exp%	Act%
165.00	100	100
167.00	63.60	63.73
101.00	34.30	39.55
0.00	0.00	0.00

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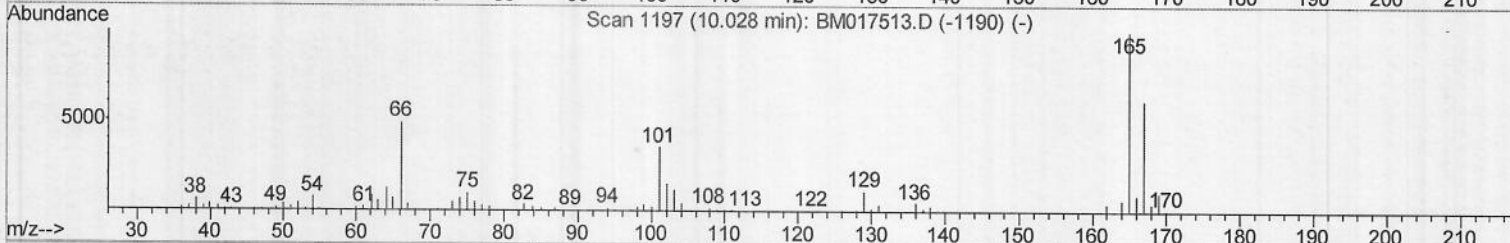
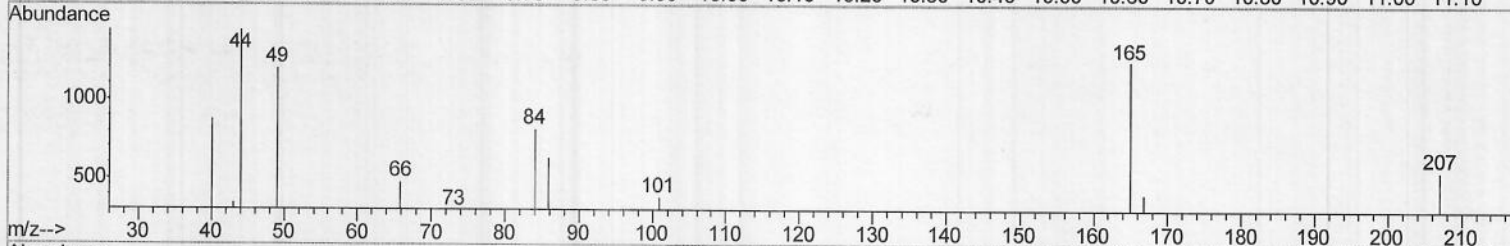
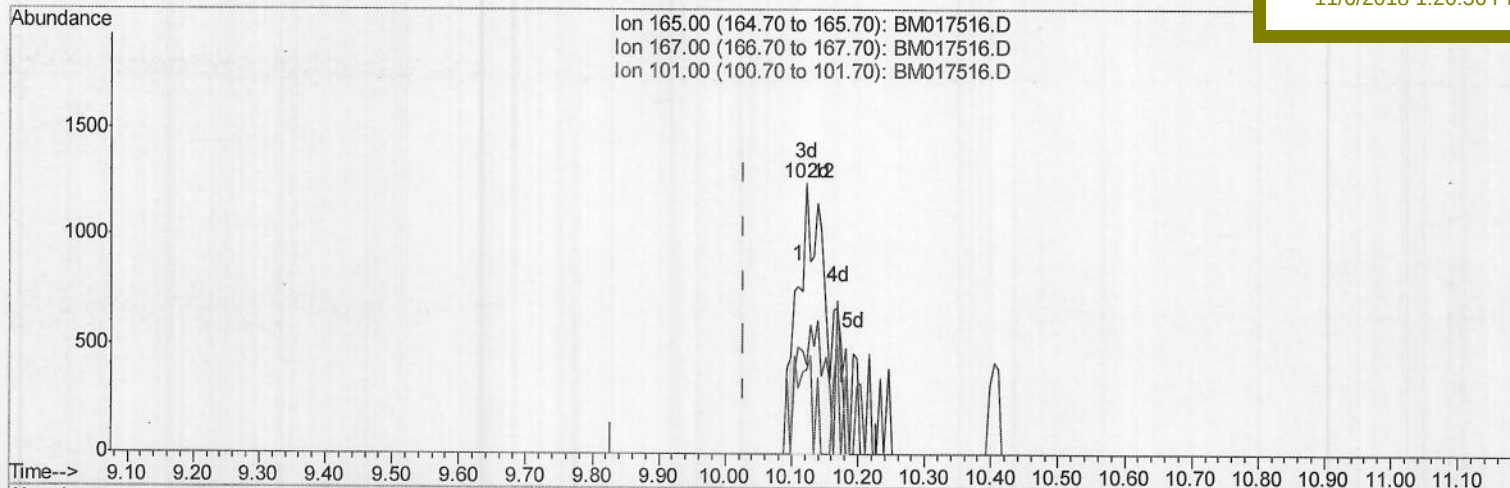
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Instrument :
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Client Sampled :
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TIC: BM017516.D

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

10.122min (+0.094) 0.27ng/ul m

response 3614

Ion	Exp%	Act%
165.00	100	100
167.00	63.60	32.64#
101.00	34.30	31.03
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	7.63	152	196122	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	818690	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	513251	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	1264857	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	1625427	20.00	ng/ul	0.00
86) Perylene-d12	23.42	264	1754883	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.22	96	409	0.10	ng/uL	0.00
5) Phenol-d5	6.83	99	3629	0.20	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	6.98	67	3569	0.34	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	4155	0.30	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	1212	0.09	ng/ul	0.02
19) Nitrobenzene-d5	8.81	128	1803m	0.28	ng/ul	0.03
22) 2-Nitrophenol-d4	9.51	143	1654	0.23	ng/ul	0.01
26) 2,4-Dichlorophenol-d3	10.12	165	3614m	0.27	ng/ul	0.09
29) 4-Chloroaniline-d4	10.62	131	5450	0.50	ng/ul	0.06
44) Dimethylphthalate-d6	13.71	166	15254	0.35	ng/ul	0.02
47) Acenaphthylene-d8	13.97	160	17405	0.33	ng/ul	0.00
52) 4-Nitrophenol-d4	14.60	143	657	0.08	ng/ul	0.11
58) Fluorene-d10	15.28	176	16204	0.42	ng/ul	0.01
63) 4,6-Dinitro-2-methylphenol	15.44	200	622	0.07	ng/ul	0.04
71) Anthracene-d10	17.13	188	31167	0.50	ng/ul	0.01
79) Pyrene-d10	19.43	212	52561	0.70	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.28	264	86689	0.94	ng/ul	0.00

Target Compounds

					Qvalue	
84) Bis(2-ethylhexyl)phthalate	21.14	149	3445852	62.579	ng/ul	97

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

} SJ
11/06/18