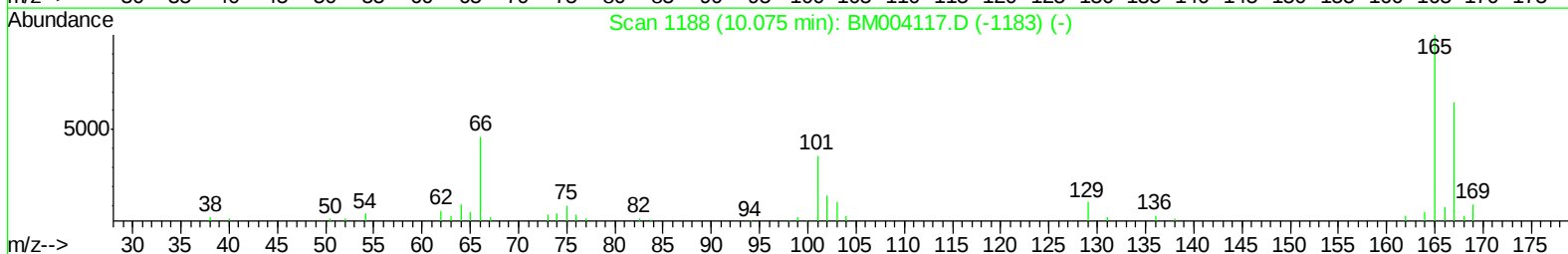
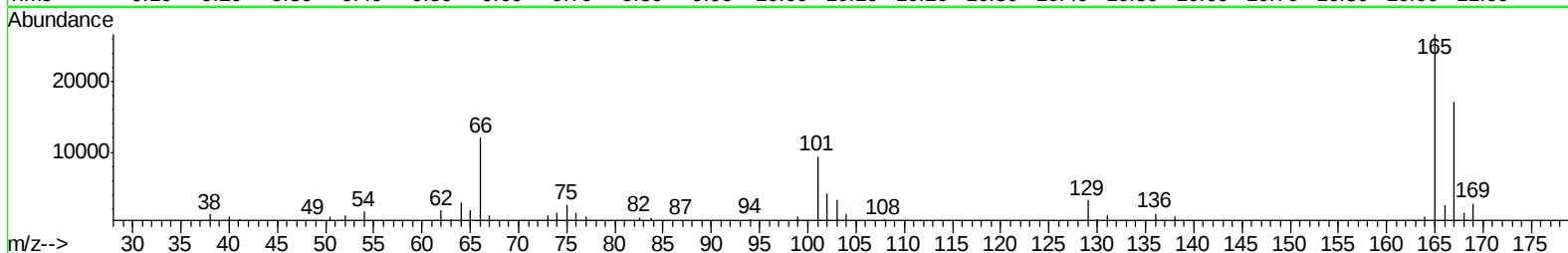
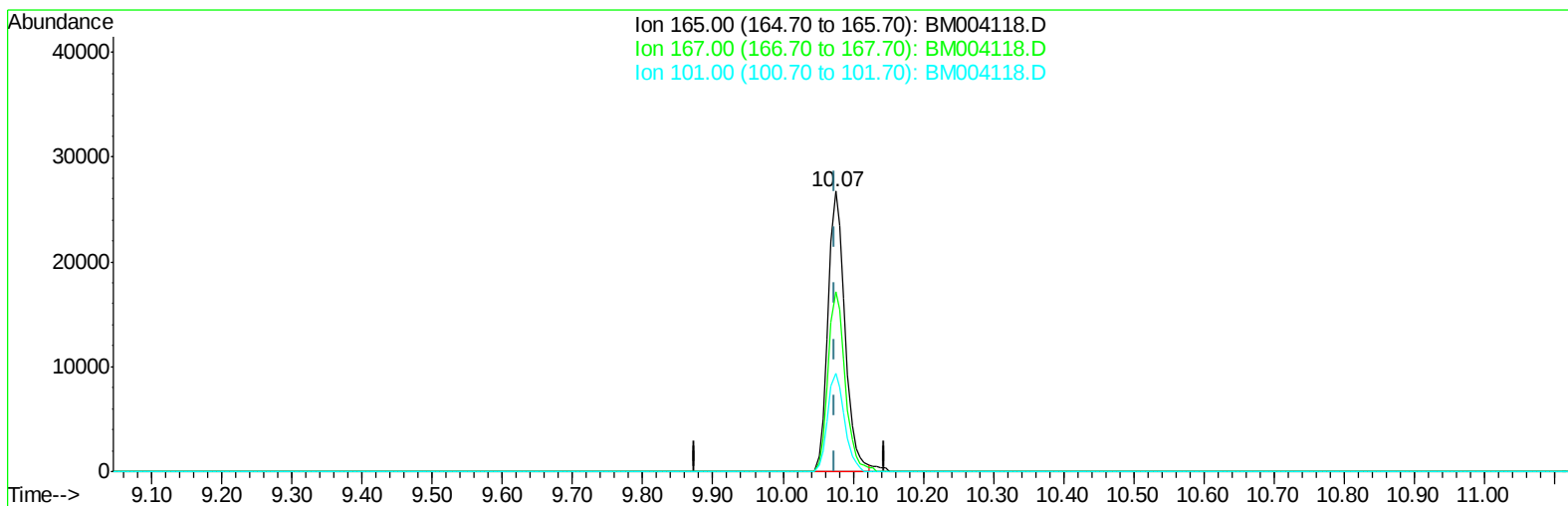


Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012316\
Data File : BM004118.D
Acq On : 22 Jan 2016 20:02
Operator : SJ/UM
Sample : PB87916BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jan 23 02:40:24 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jan 23 02:35:35 2016
Response via : Initial Calibration



TIC: BM004118.D

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

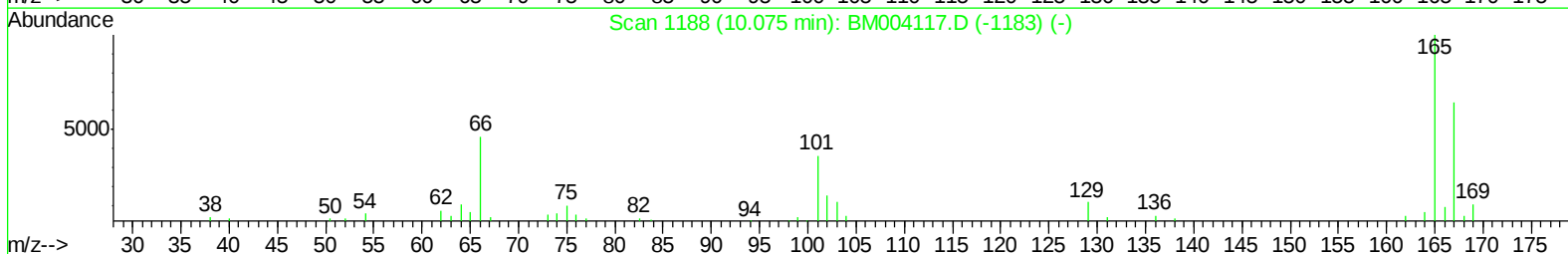
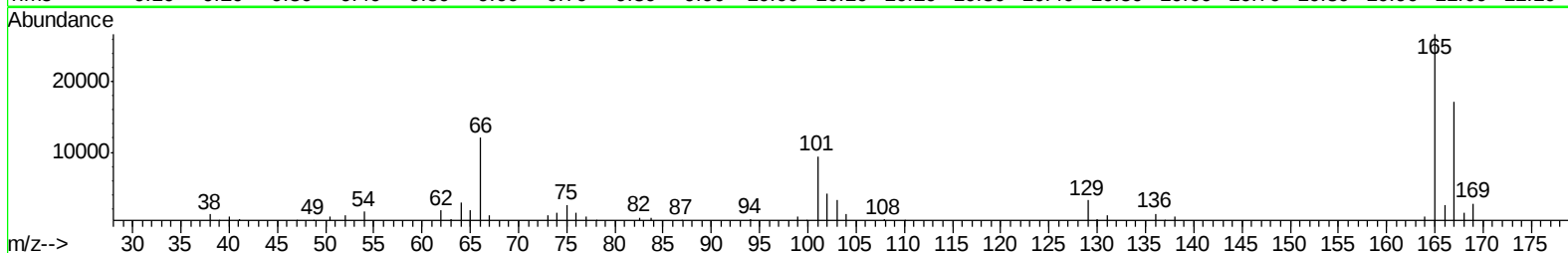
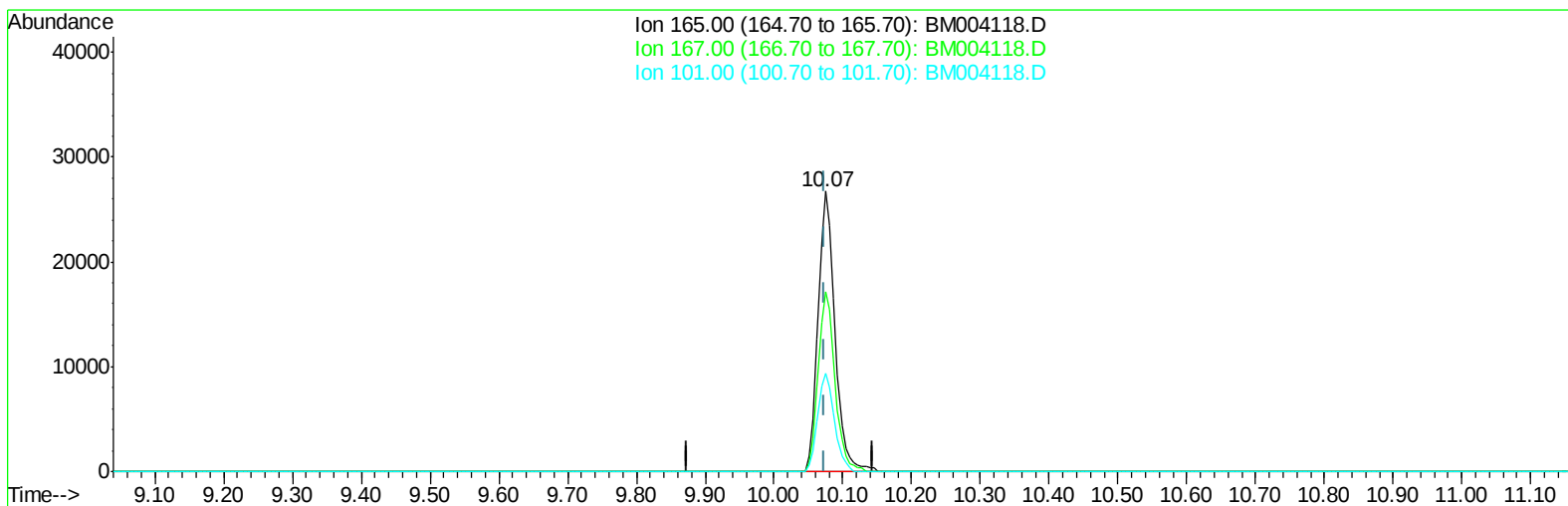
10.075min (-0.000) 30.04ng/ul

response 44654

Ion	Exp%	Act%
165.00	100	100
167.00	65.30	64.07
101.00	34.20	35.14
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012316\
Data File : BM004118.D
Acq On : 22 Jan 2016 20:02
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Sample : PB87916BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jan 23 02:40:24 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jan 23 02:35:35 2016
Response via : Initial Calibration



TIC: BM004118.D

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

10.075min (-0.000) 30.45ng/ul m

response 45256

Ion	Exp%	Act%
165.00	100	100
167.00	65.30	64.07
101.00	34.20	35.14
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012316\
 Data File : BM004118.D
 Acq On : 22 Jan 2016 20:02
 Operator : SJ/UM
 Sample : PB87916BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jan 23 02:44:51 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jan 23 02:35:35 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	20113	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	97240	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	62269	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	155812	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	211261	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	235011	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.12	96	2668	7.01	ng/uL	0.00
5) Phenol-d5	6.84	99	54931	28.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	33746	32.31	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	47378	31.02	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	48565	28.85	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	23758	32.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.54	143	25771	32.23	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.07	165	45256m	30.45	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	77145	40.77	ng/ul	0.00
44) Dimethylphthalate-d6	13.73	166	191915	35.83	ng/ul	0.00
47) Acenaphthylene-d8	14.01	160	217886	34.93	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	28176	26.28	ng/ul	0.00
58) Fluorene-d10	15.32	176	163807	35.64	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	20697	21.55	ng/ul	0.00
71) Anthracene-d10	17.17	188	278800	36.78	ng/ul	0.00
79) Pyrene-d10	19.47	212	341023	36.03	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	453920	37.03	ng/ul	0.00

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

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