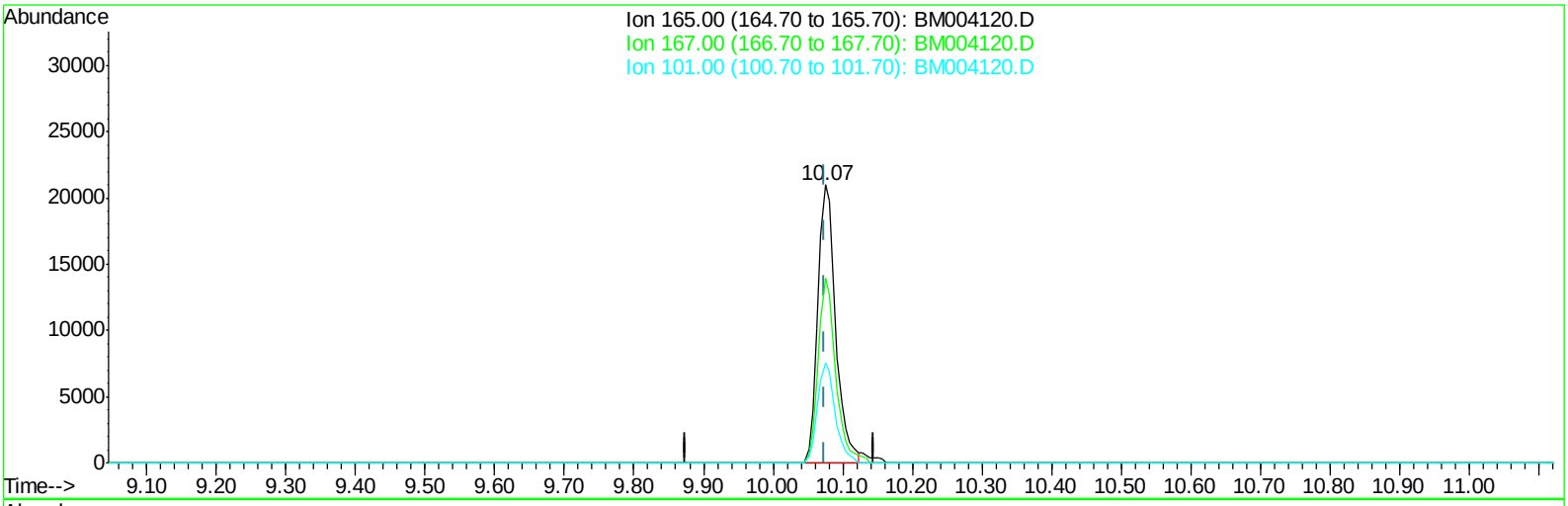
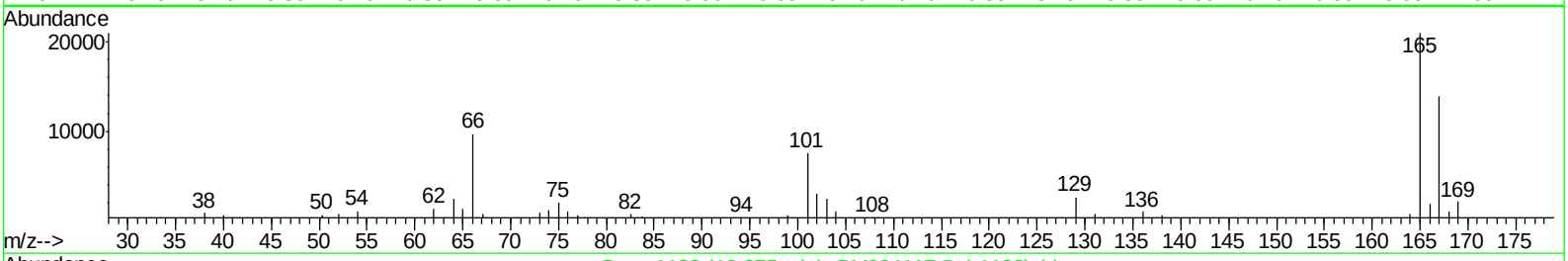
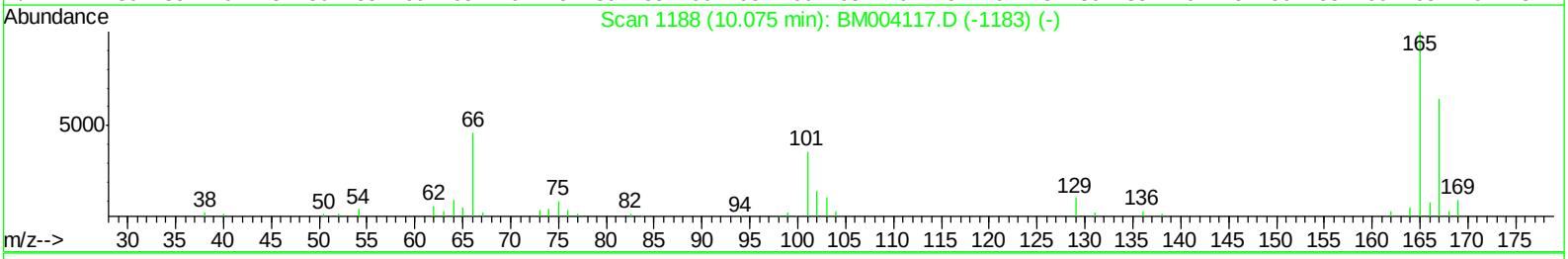
Quantitation Report (Qedit)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012316\
Data File : BM004120.D
Acq On : 22 Jan 2016 21:13
Operator : SJ/UM
Sample : H1174-15
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Jan 23 02:40:44 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: BM004120.D

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

10.075min (-0.000) 25.60ng/ul

response 37198

Ion	Exp%	Act%
165.00	100	100
167.00	65.30	66.42
101.00	34.20	36.01
0.00	0.00	0.00

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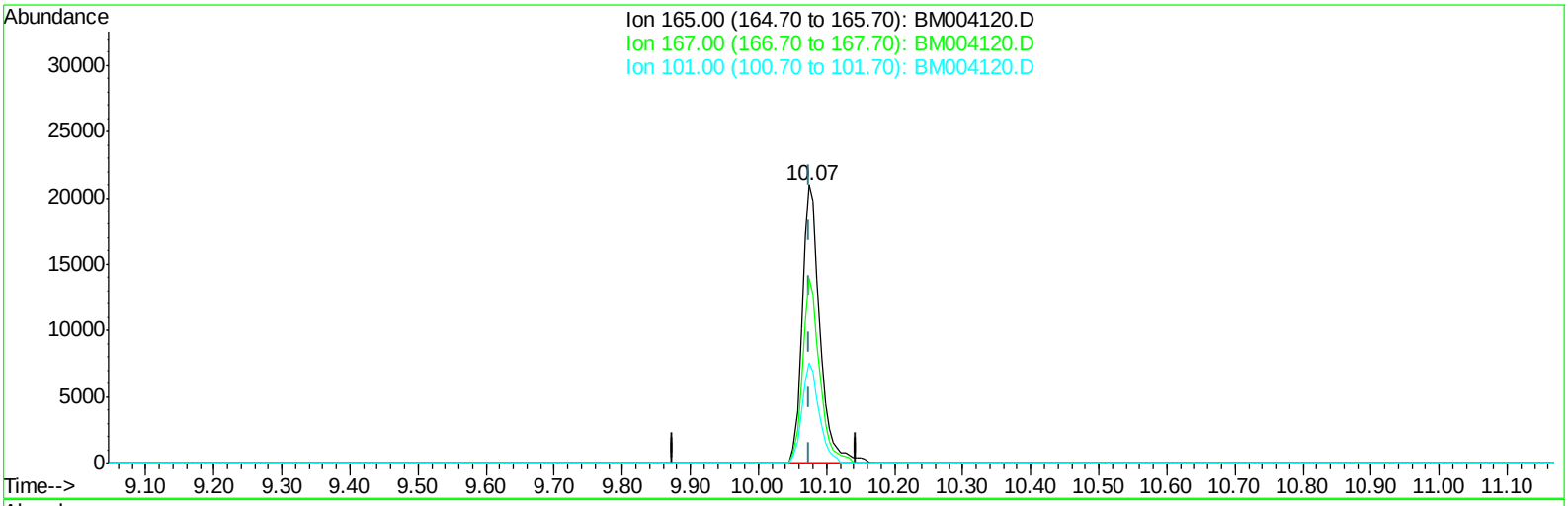

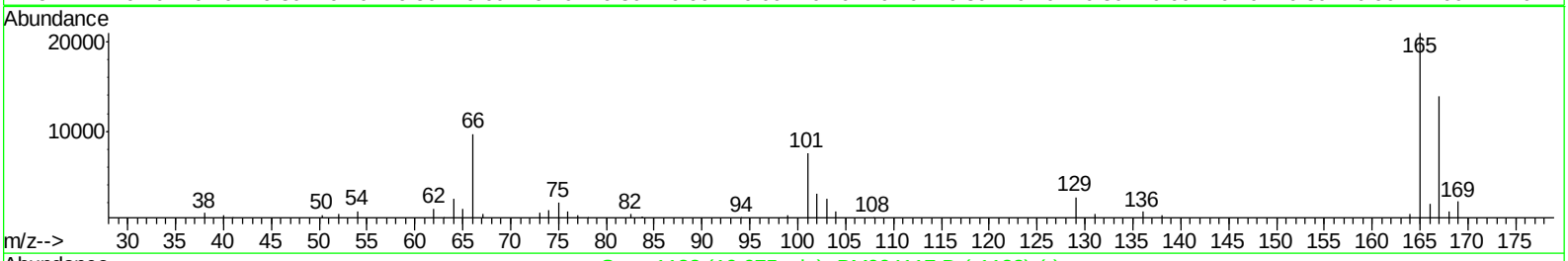

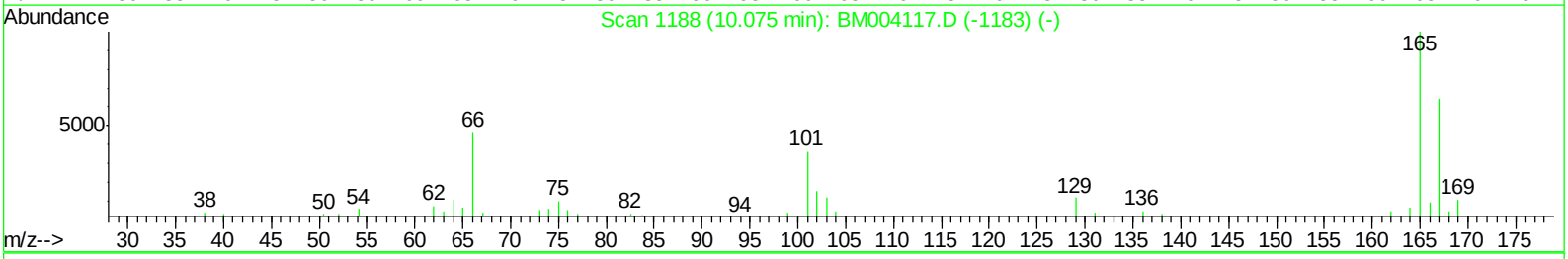

TIC: BM004120.D

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

10.075min (-0.000) 26.27ng/ul m

response 38171

Ion	Exp%	Act%
165.00	100	100
167.00	65.30	66.42
101.00	34.20	36.01
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012316\
Data File : BM004120.D
Acq On : 22 Jan 2016 21:13
Operator : SJ/UM
Sample : H1174-15
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Jan 23 02:55:27 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	19522	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	95054	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	61990	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	155459	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	212964	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	239495	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.12	96	457	1.24	ng/uL	0.00
5) Phenol-d5	6.84	99	12815	6.73	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	29155	28.76	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	35538	23.97	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	24420	14.95	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	20355	28.89	ng/ul	0.00
22) 2-Nitrophenol-d4	9.54	143	21967	28.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.07	165	38171m	26.27	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.73	166	172328	32.31	ng/ul	0.00
47) Acenaphthylene-d8	14.01	160	196146	31.59	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	4622	4.33	ng/ul	0.01
58) Fluorene-d10	15.32	176	158200	34.58	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	21887	22.84	ng/ul	0.00
71) Anthracene-d10	17.17	188	285560	37.76	ng/ul	0.00
79) Pyrene-d10	19.47	212	350821	36.77	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	461040	36.91	ng/ul	0.00

Target Compounds	Qvalue

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

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