

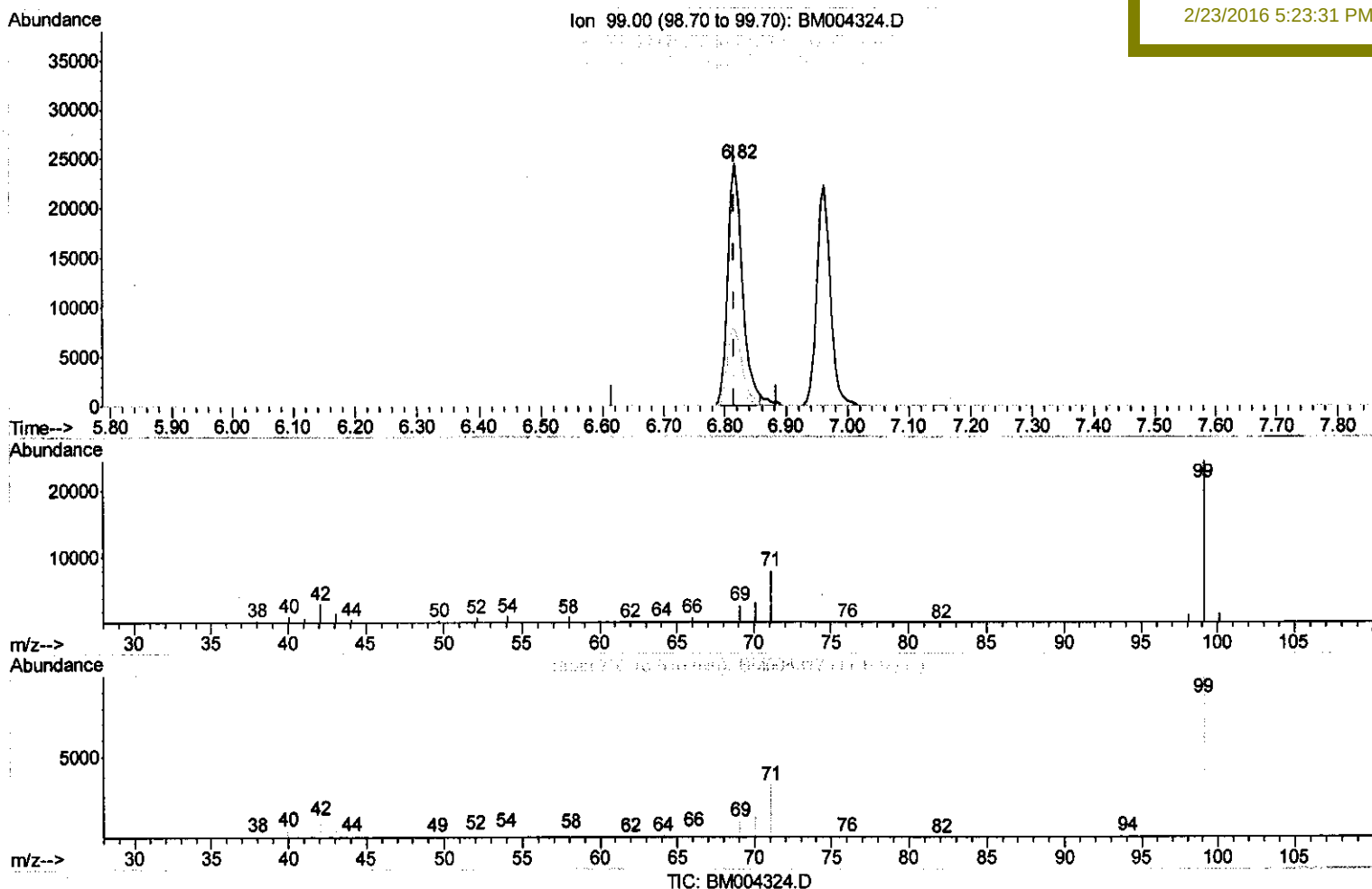
Data Path : Z:\HPCHEM1\BNA_M\Data\BM022216\
Data File : BM004324.D
Acq On : 22 Feb 2016 22:56
Operator : SJ/UM
Sample : H1499-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Quant Time: Feb 23 00:16:15 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM022216.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 23 00:11:03 2016
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
0311-S10-0312

Manual Integrations
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(5) Phenol-d5 (S)

6.816min (+0.000) 19.27ng/ul

response 40462

Ion	Exp%	Act%
99.00	100	100
71.00	35.10	32.67
42.00	14.10	12.72
0.00	0.00	0.00

Quantitation Report (Qedit)

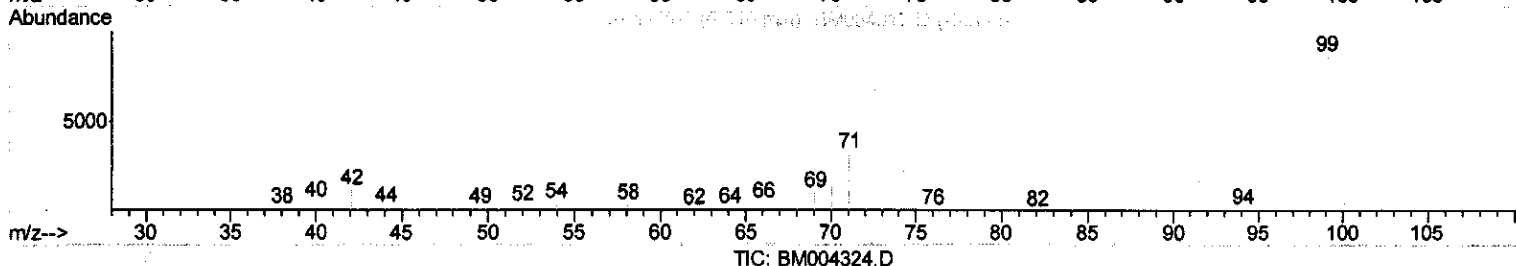
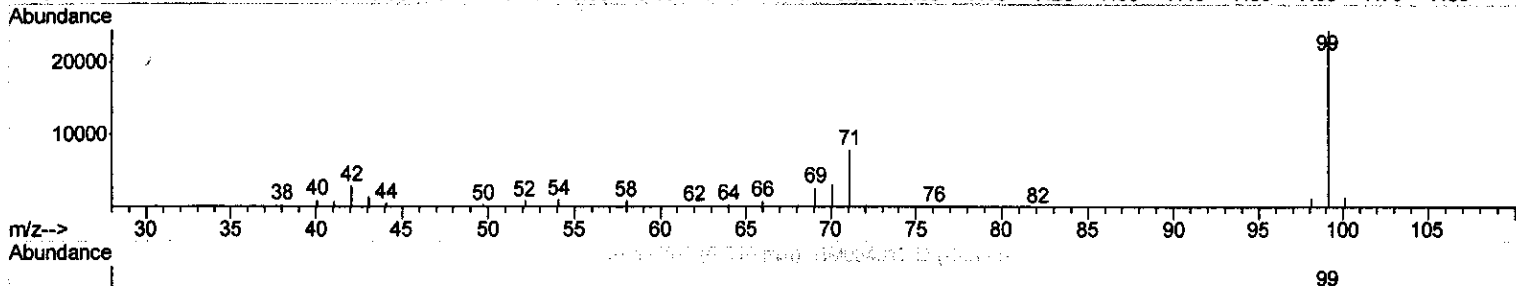
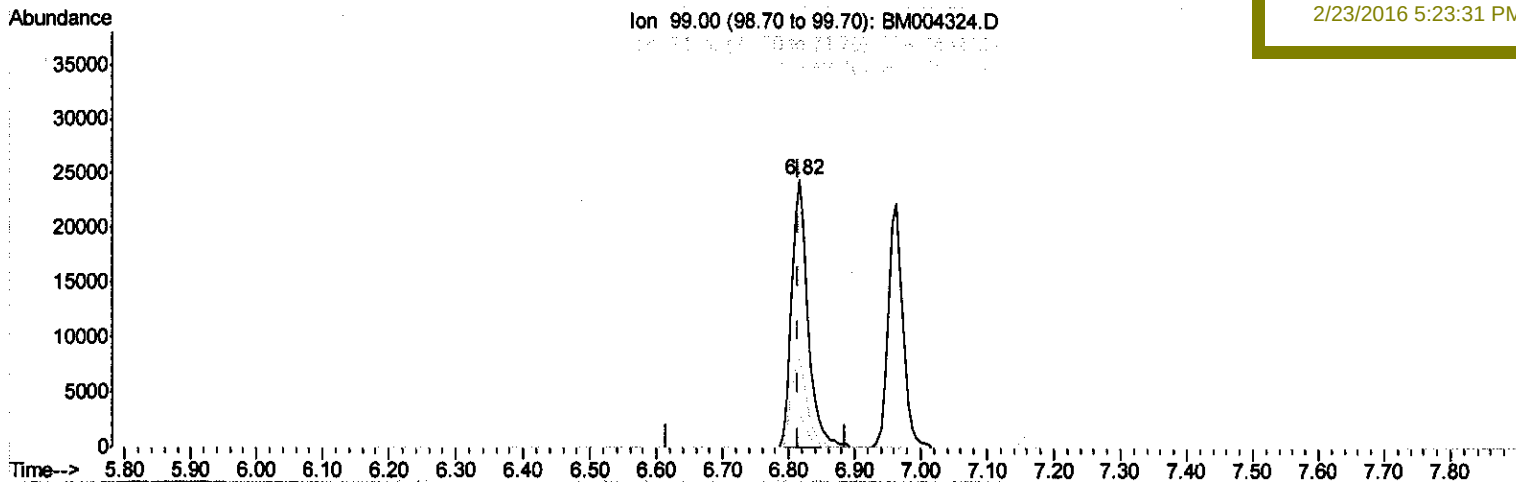
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(5) Phenol-d5 (S)

6.816min (+0.000) 19.67ng/ul m

response 41302

Ion	Exp%	Act%
99.00	100	100
71.00	35.10	32.67
42.00	14.10	12.72
0.00	0.00	0.00

> 9.5
 2/29/16

Quantitation Report (Qedit)

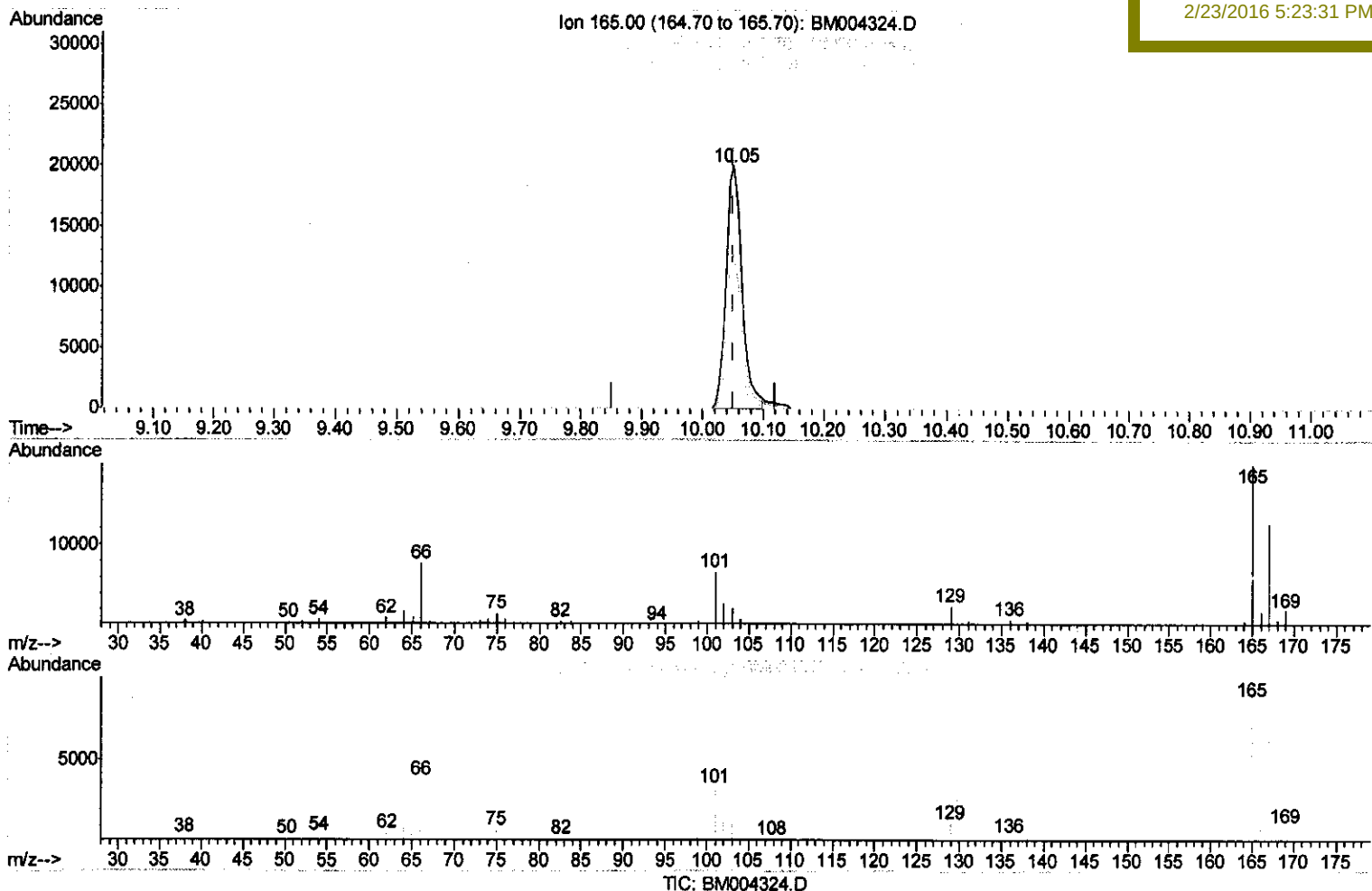
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(26) 2,4-Dichlorophenol-d3 (S)

10.051min (+0.000) 19.02ng/ul

response 36491

Ion	Exp%	Act%
165.00	100	100
167.00	65.30	63.92
101.00	34.20	33.51
0.00	0.00	0.00

Quantitation Report (Qedit)

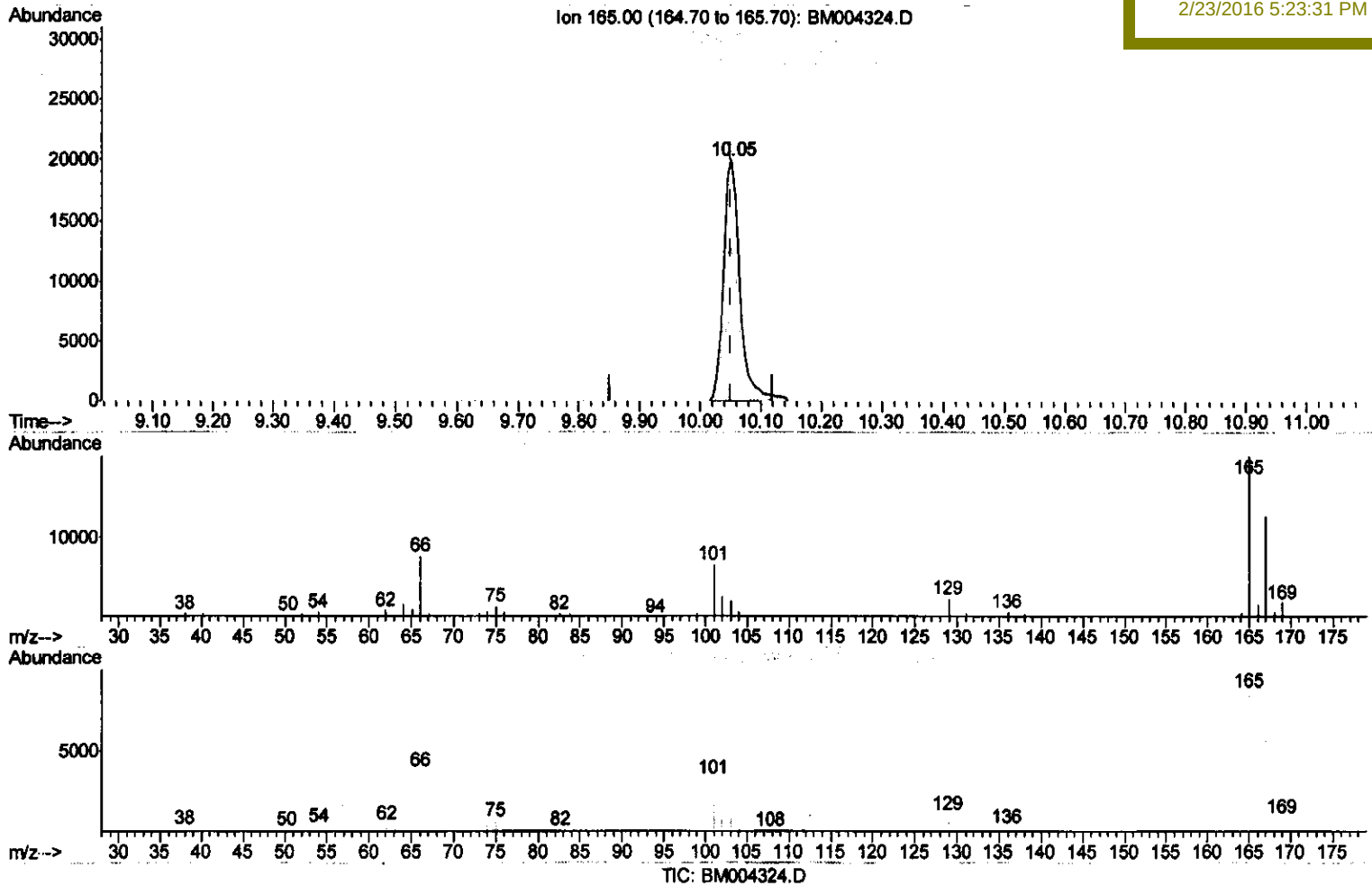
Data Path : Z:\HPCHEM1\BNA_M\Data\BM022216\
 Data File : BM004324.D
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 Operator : SJ/UM
 Sample : H1499-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

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Manual Integrations
 APPROVED

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

10.051min (+0.000) 19.63ng/ul m

response 37662

Ion	Exp%	Act%
165.00	100	100
167.00	65.30	63.92
101.00	34.20	33.51
0.00	0.00	0.00

> 9.7
 2/29/16

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Instrument :
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Quant Time: Feb 23 00:49:46 2016
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	25046	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	122382	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	80466	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	195597	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	241583	20.00	ng/ul	0.00
86) Perylene-d12	23.47	264	228980	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	1972	4.15	ng/uL	0.00
5) Phenol-d5	6.82	99	41302m	19.67	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.96	67	23867	22.35	ng/ul	0.00
9) 2-Chlorophenol-d4	7.16	132	37664	20.79	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	36486	19.44	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	19132	20.55	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	22064	20.79	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.05	165	37662m	19.63	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	57475	24.04	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	154600	22.40	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	178764	22.13	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	18689	16.85	ng/ul	0.00
58) Fluorene-d10	15.29	176	134810	22.97	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.43	200	19003	15.82	ng/ul	0.00
71) Anthracene-d10	17.14	188	216937	22.64	ng/ul	0.00
79) Pyrene-d10	19.44	212	264785	22.66	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.33	264	279429	23.25	ng/ul	0.00

Target Compounds						Qvalue
45) Dimethylphthalate	13.75	163	73241	10.18	ng/ul	99

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

5.7
 2/29/16

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