

Quantitation Report (QT Reviewed)

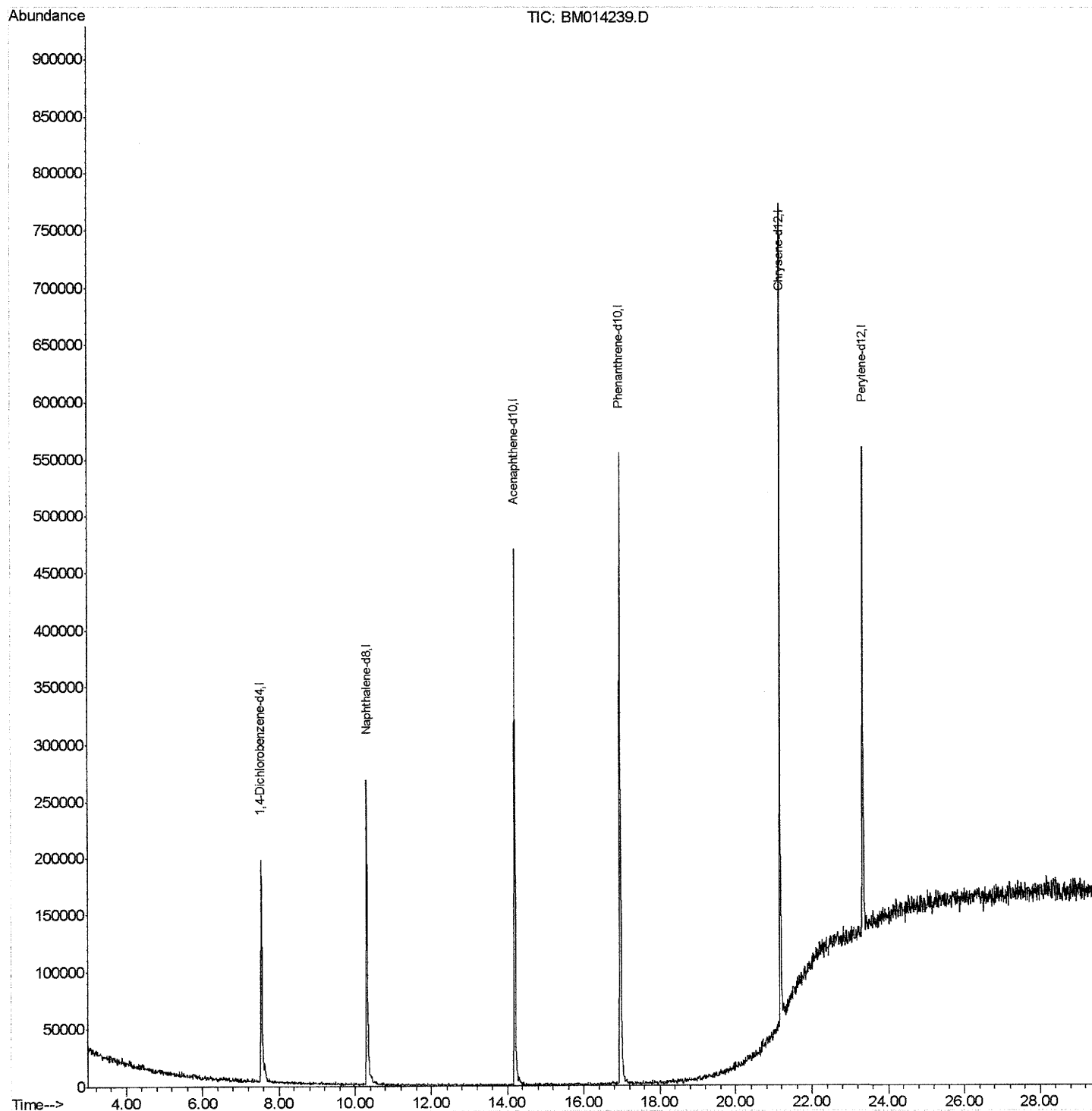
Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA_M\DATA\BM021418\
 Data File : BM014239.D
 Acq On : 14 Feb 2018 16:37
 Operator : SJ/JU
 Sample : J1524-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 Client Sampled :
 SVOC-GPC2-BLANK

Manual Integrations
 APPROVED

Sohil
 2/15/2018 11:09:03 AM

Quant Time: Feb 14 17:46:41 2018
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 14 15:20:17 2018
 Response via : Initial Calibration



Quantitation Report (Qedit)

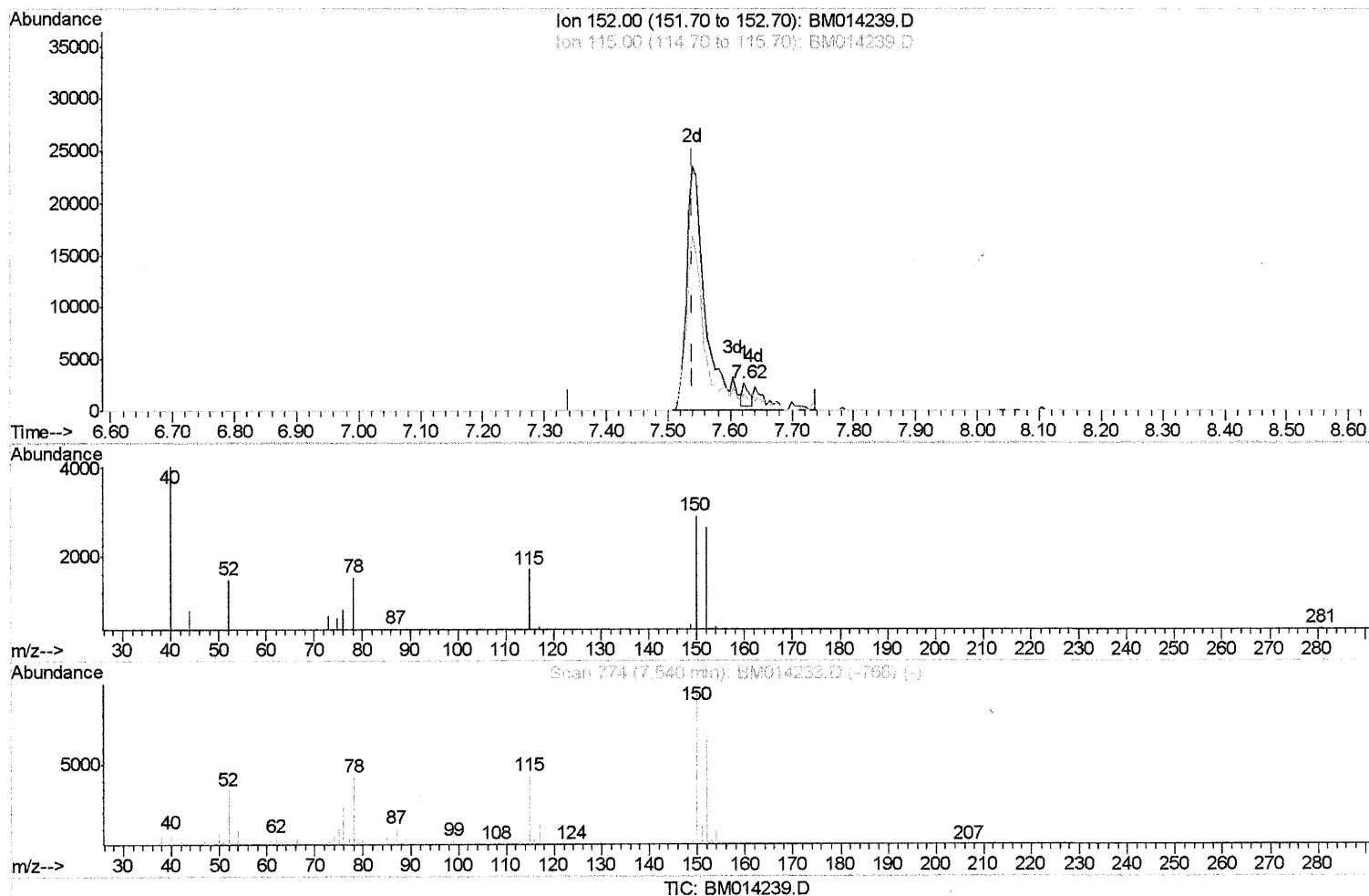
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(1) 1,4-Dichlorobenzene-d4 (I)

7.622min (+0.082) 20.00ng/ul

response 1588

Ion	Exp%	Act%
152.00	100	100
115.00	58.90	65.04
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

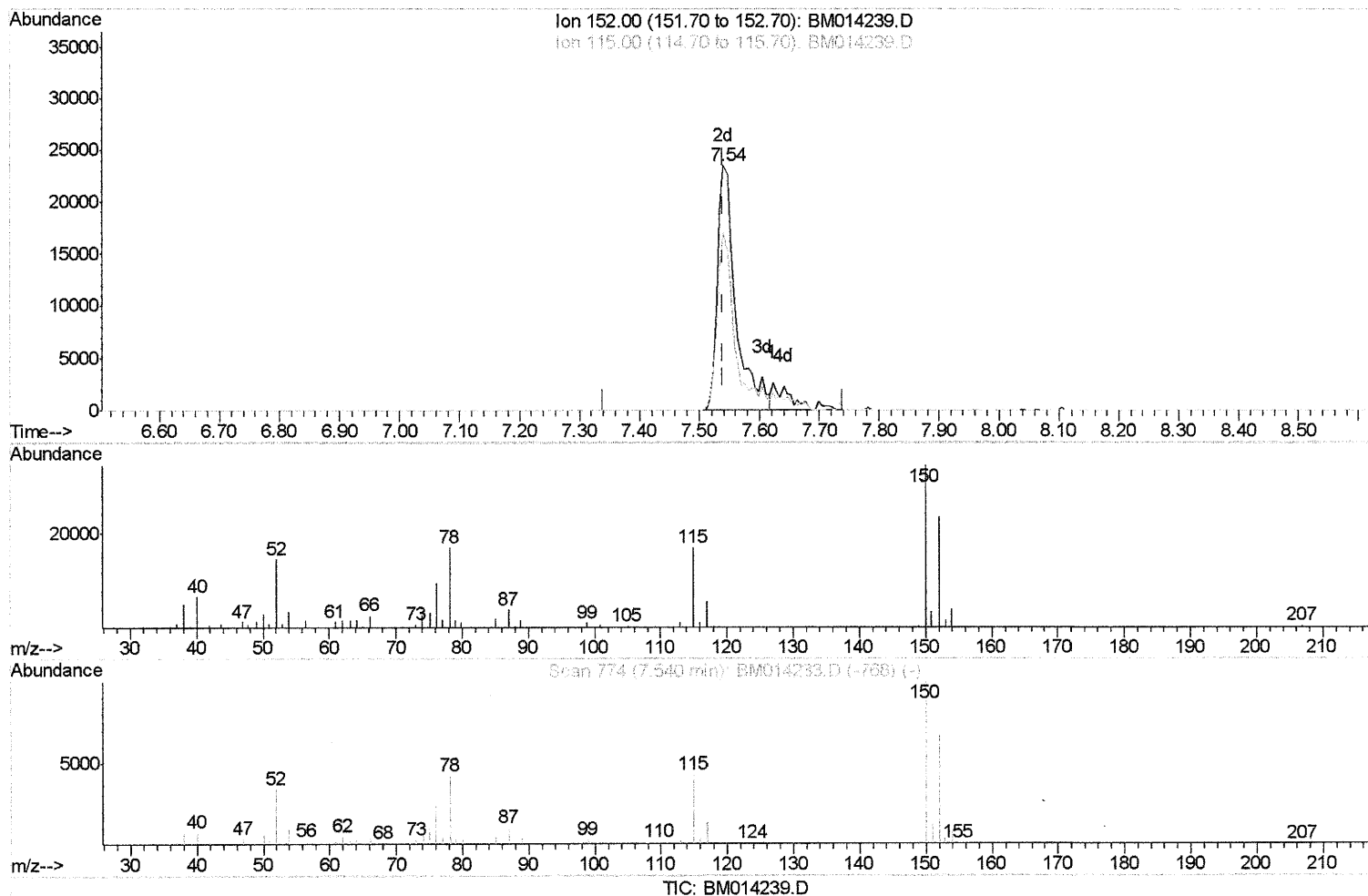
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(1) 1,4-Dichlorobenzene-d4 (I)

7.540min (+0.000) 20.00ng/ul m SJ 2/19/18

response 49815

Ion	Exp%	Act%
152.00	100	100
115.00	58.90	72.07#
0.00	0.00	0.00
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	7.54	152	49815m	20.00	ng/ul	0.00
18) Naphthalene-d8	10.31	136	205413	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.19	164	148598	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.96	188	345594	20.00	ng/ul	0.00
75) Chrysene-d12	21.16	240	341078	20.00	ng/ul	0.00
83) Perylene-d12	23.34	264	322520	20.00	ng/ul	0.00

SJ 2/19/18

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL
5) Phenol-d5	0.00	99	0	0.00	ng/ul
7) Bis-(2-Chloroethyl) ether-d	0.00	67	0	0.00	ng/ul
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul
46) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul
57) Fluorene-d10	0.00	176	0d	0.00	ng/ul
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul
70) Anthracene-d10	0.00	188	0d	0.00	ng/ul
76) Pyrene-d10	0.00	212	0	0.00	ng/ul
87) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul

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