

Quantitation Report (QT Reviewed)

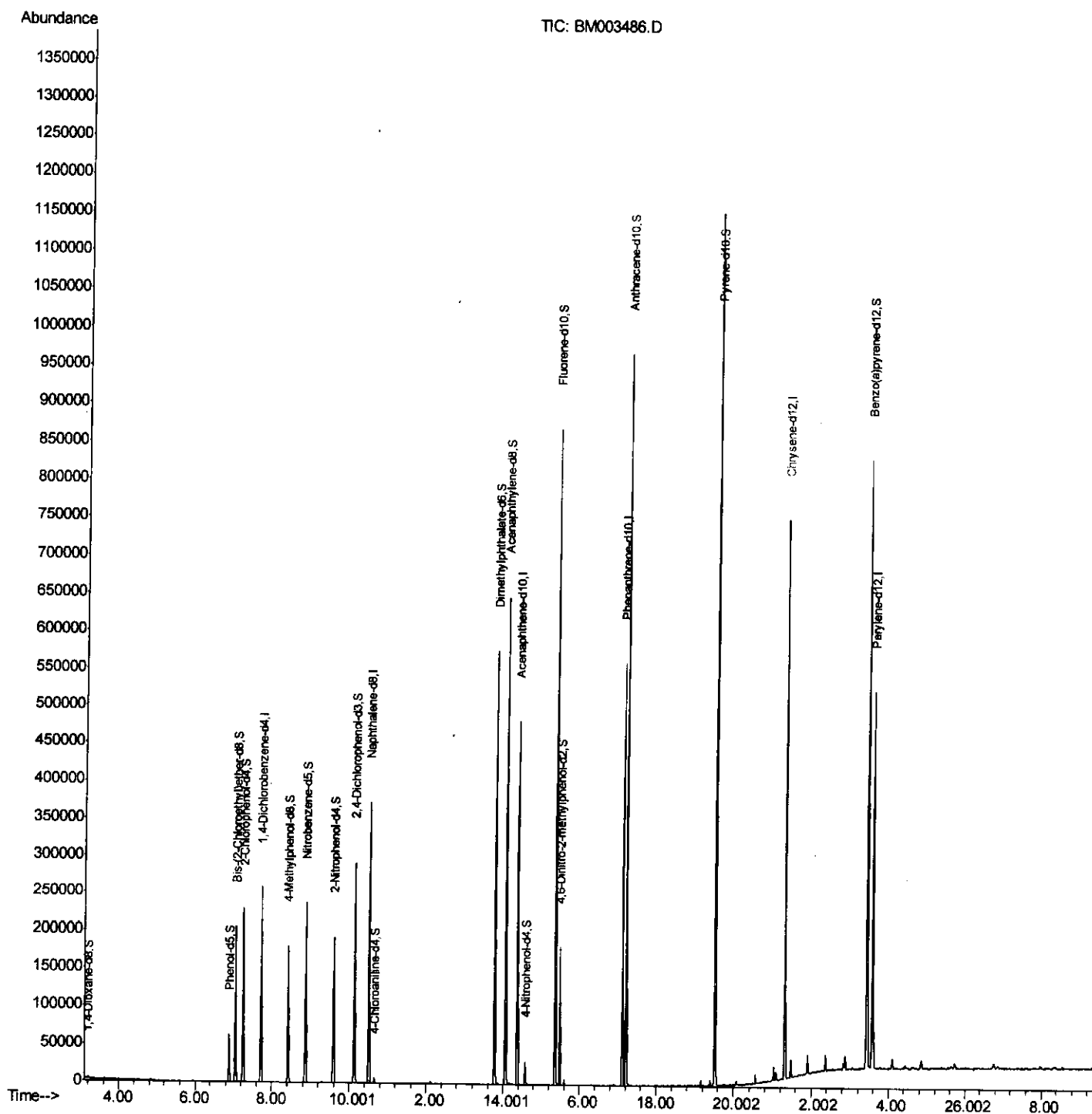
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121115\
Data File : BM003486.D
Acq On : 11 Dec 2015 17:06
Operator : UM/IZ
Sample : G4718-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AD1

Manual Integrations
APPROVED

MMdadoda
12/14/2015 6:27:46 PM

Quant Time: Dec 12 01:09:55 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM120115.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Dec 12 00:54:45 2015
Response via : Initial Calibration



Quantitation Report (Qedit)

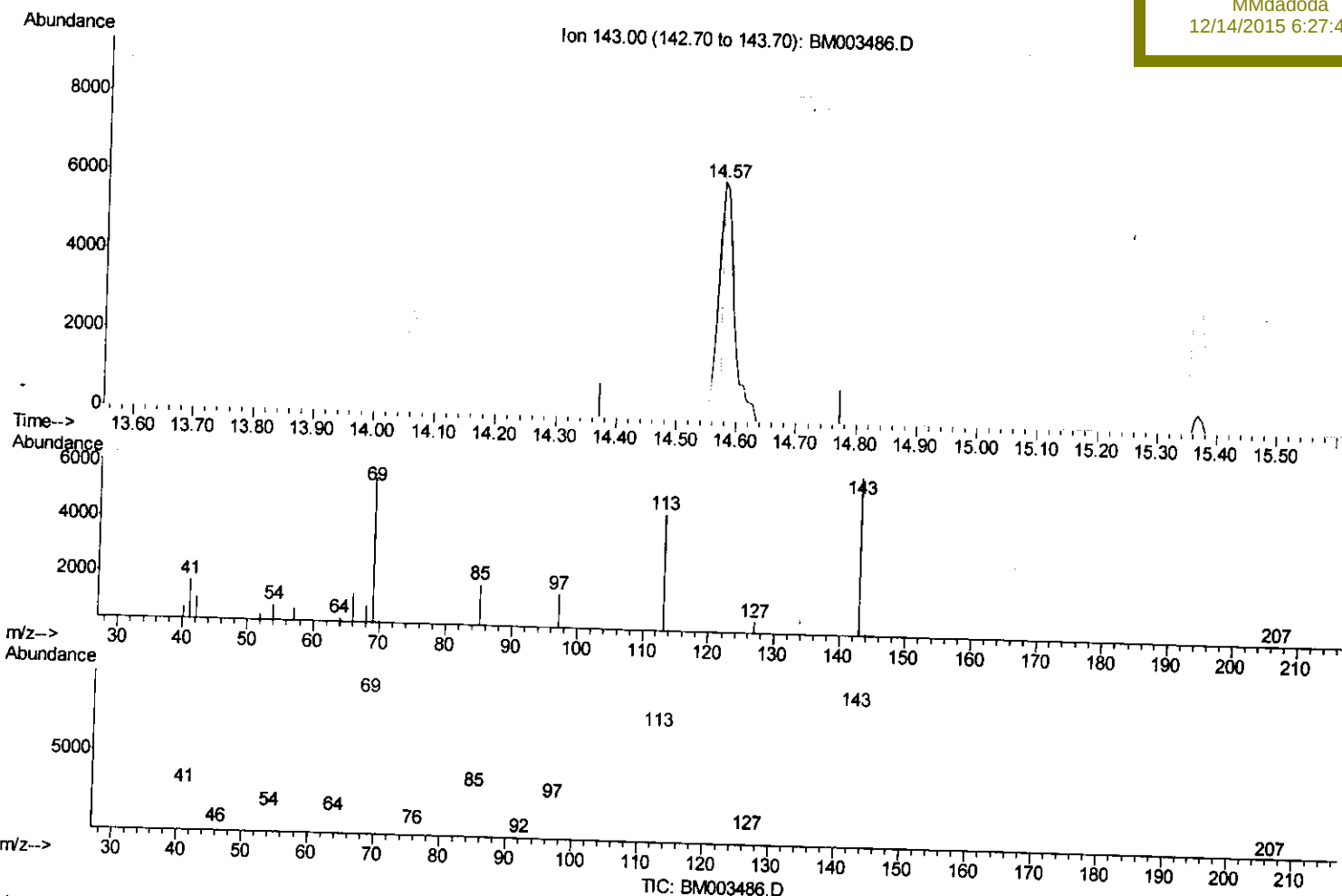
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 Response via : Initial Calibration



(52) 4-Nitrophenol-d4 (S)

14.574min (-0.000) 4.83ng/ul

response 10885

Ion	Exp%	Act%
143.00	100	100
113.00	70.40	74.90
41.00	27.00	28.67
42.00	17.60	18.75

Quantitation Report (Qedit)

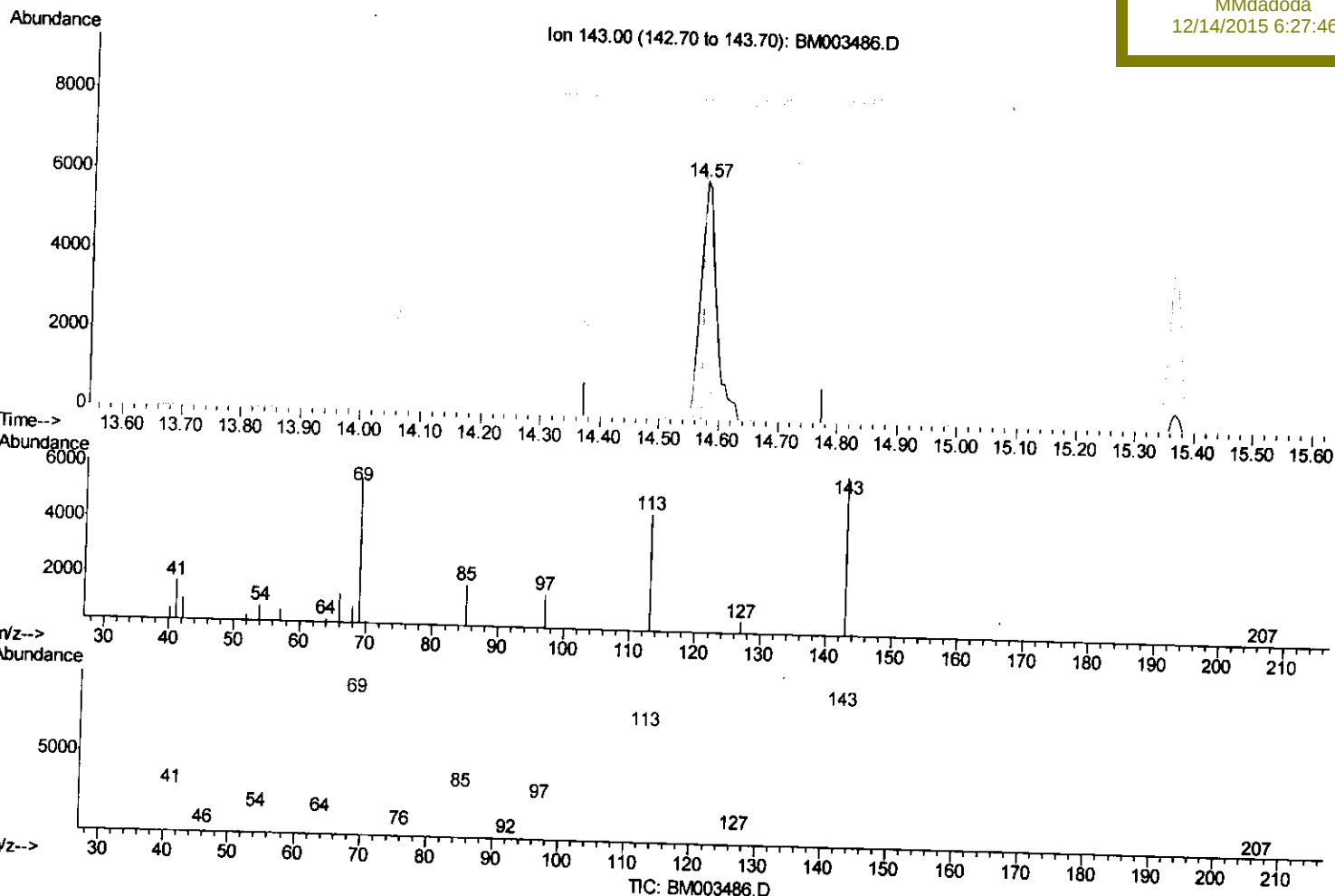
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(52) 4-Nitrophenol-d4 (S)

14.574min (-0.000) 4.90ng/ul m U.M
 response 11040 12/15/15

Ion	Exp%	Act%
143.00	100	100
113.00	70.40	74.90
41.00	27.00	28.67
42.00	17.60	18.75

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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	7.72	152	72681	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	322058	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	167275	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.12	188	334854	20.00	ng/ul	0.00
78) Chrysene-d12	21.31	240	349900	20.00	ng/ul	0.00
86) Perylene-d12	23.57	264	347495	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.19	96	2087	1.43	ng/uL	0.00
5) Phenol-d5	6.89	99	38810	6.69	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	86513	27.03	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	113289	23.61	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	71345	14.66	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	63314	28.82	ng/ul	0.00
22) 2-Nitrophenol-d4	9.60	143	67279	29.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	112776	24.76	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	6022	1.17	ng/ul	0.00
44) Dimethylphthalate-d6	13.78	166	391571	31.14	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	462898	29.95	ng/ul	0.00
52) 4-Nitrophenol-d4	14.57	143	11040m	4.90	ng/ul	0.00
58) Fluorene-d10	15.37	176	343522	31.74	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	40733	23.78	ng/ul	0.00
71) Anthracene-d10	17.22	188	546153	36.71	ng/ul	0.00
79) Pyrene-d10	19.52	212	579848	33.85	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.43	264	594178	35.44	ng/ul	0.00

U.M
 12/15/15

Target Compounds

Qvalue

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