

Quantitation Report (Qedit)

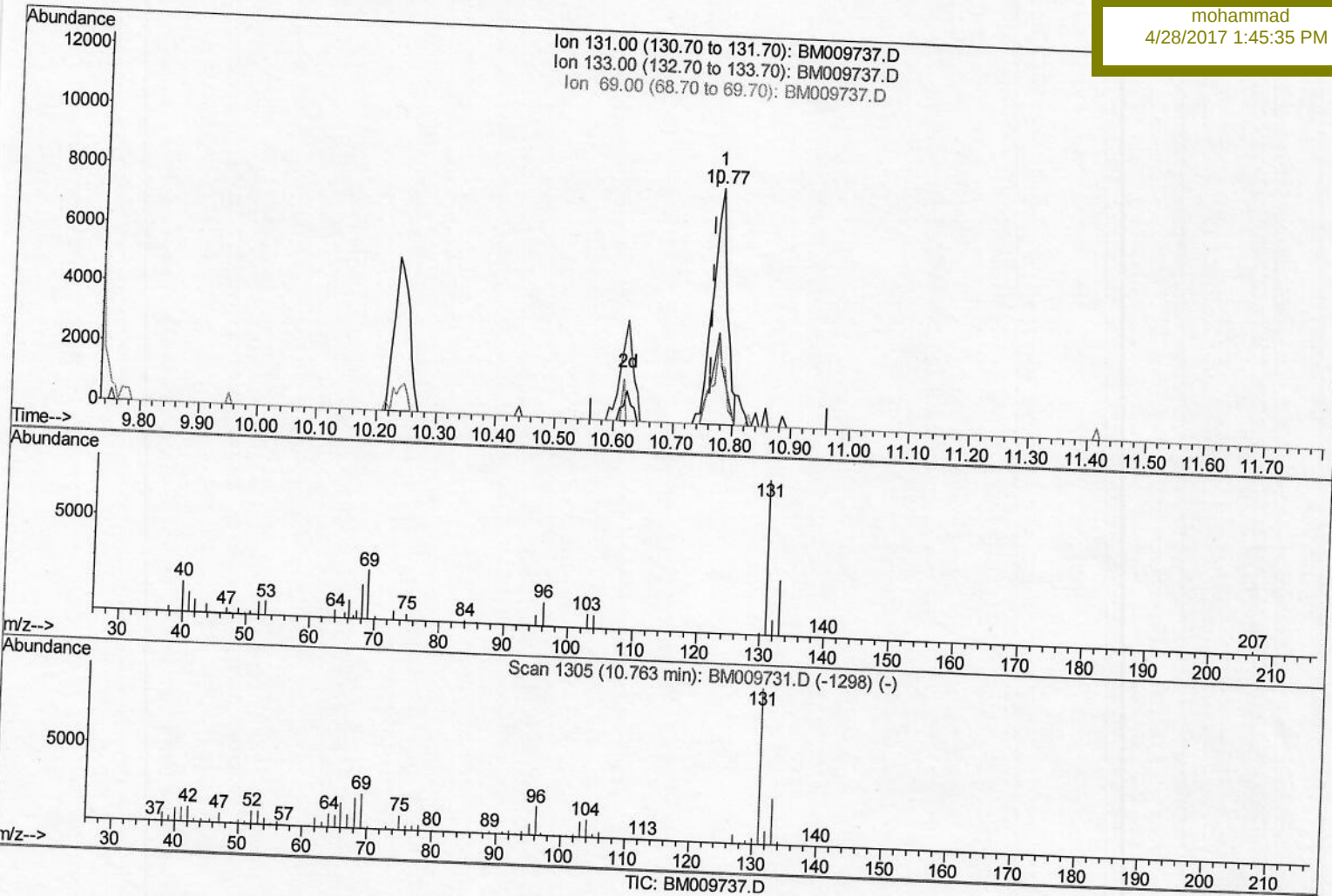
Data Path : Z:\HPCHEM1\BNA_M\Data\BM042717\
 Data File : BM009737.D
 Acq On : 27 Apr 2017 12:39
 Operator : SJ/MA
 Sample : I2809-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Apr 27 18:19:47 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 27 18:07:34 2017
 Response via : Initial Calibration

Instrument :
 BNA_M
 ClientSampleId :
 JHTC2

Manual Integrations
 APPROVED

mohammad
 4/28/2017 1:45:35 PM



(29) 4-Chloroaniline-d4 (S)
 10.775min (+0.012) 1.05ng/ul
 response 13576

Ion	Exp%	Act%
131.00	100	100
133.00	31.50	38.08#
69.00	36.50	34.16
0.00	0.00	0.00

Quantitation Report (Qedit)

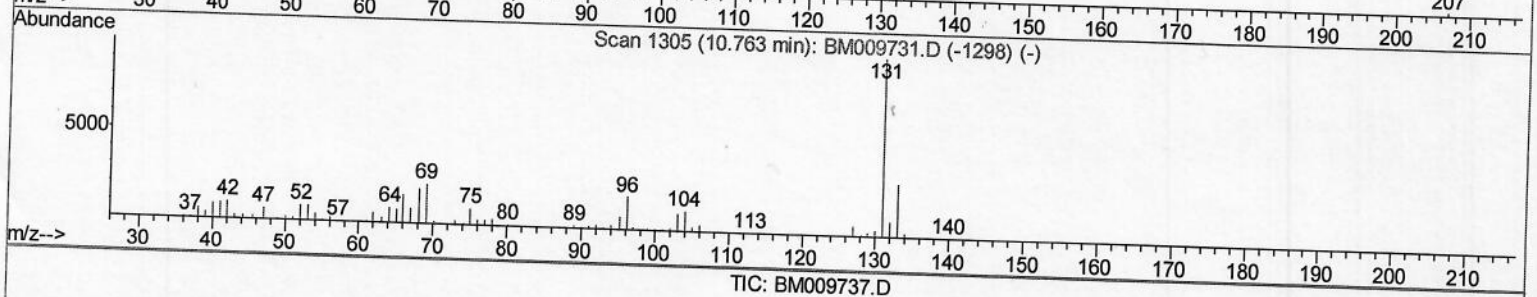
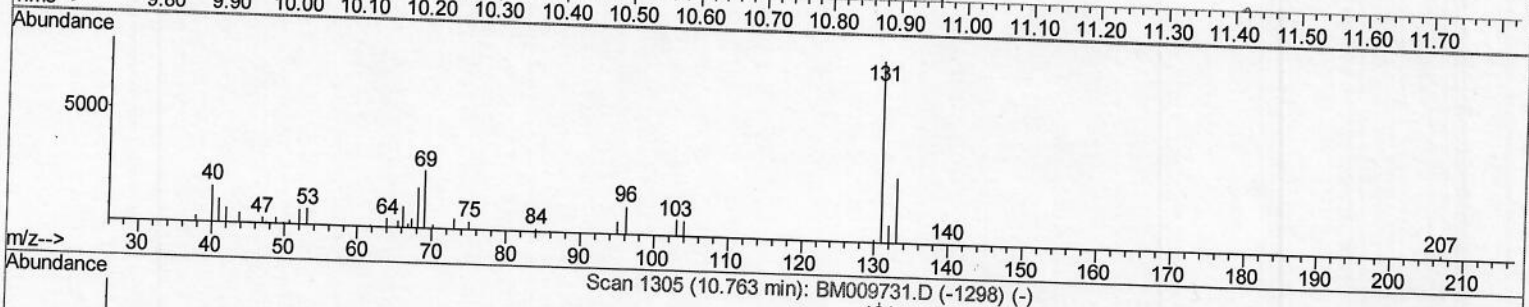
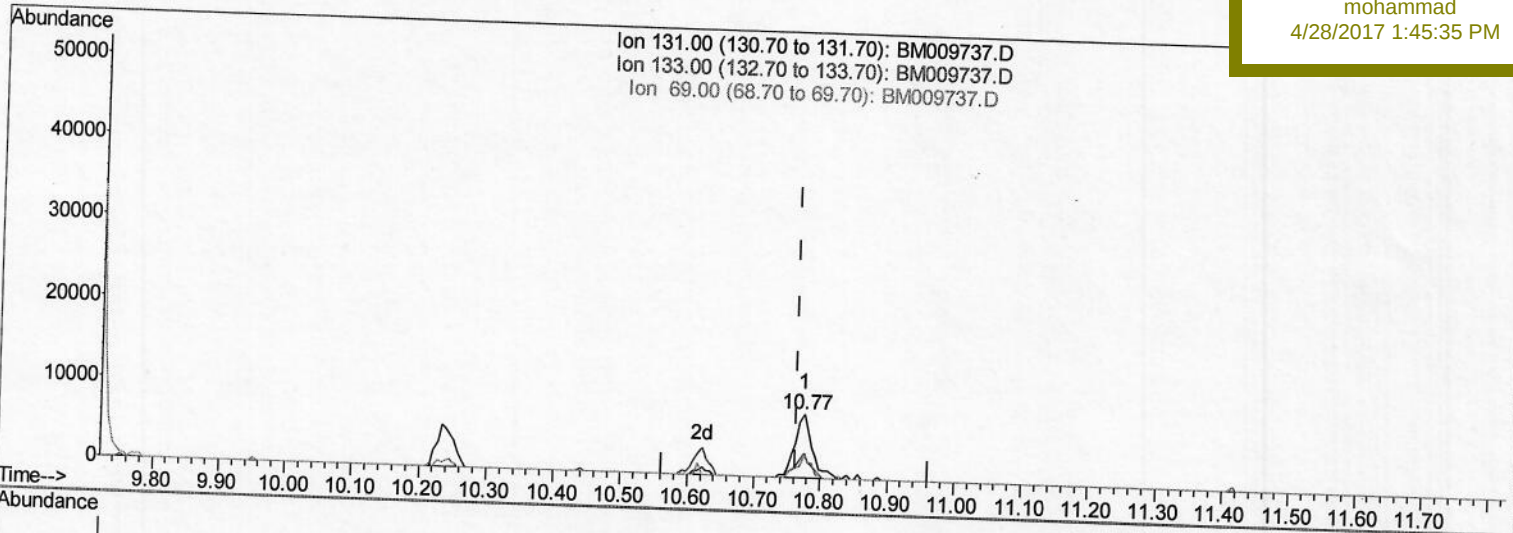
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(29) 4-Chloroaniline-d4 (S)

10.775min (+0.012) 1.11ng/ul m

response 14262

Ion Exp% Act%

131.00 100 100

133.00 31.50 38.08#

69.00 36.50 34.16

0.00 0.00 0.00

S-J
 05/04/17

Quantitation Report (QT/LSC Reviewed)

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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	7.82	152	139150	20.00	ng/ul	0.00
18) Naphthalene-d8	10.62	136	624422	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.47	164	402141	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.22	188	980495	20.00	ng/ul	0.00
75) Chrysene-d12	21.40	240	1236621	20.00	ng/ul	0.00
83) Perylene-d12	23.71	264	957782	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.30	96	15793	5.23	ng/uL	0.00
5) Phenol-d5	6.99	99	95439	6.31	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.16	67	312793	30.34	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	238586	24.81	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	180406	15.31	ng/ul	0.00
19) Nitrobenzene-d5	8.99	128	156541	30.80	ng/ul	0.00
22) 2-Nitrophenol-d4	9.71	143	176103	32.93	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.24	165	307265	28.73	ng/ul	0.00
29) 4-Chloroaniline-d4	10.77	131	14262m	1.11	ng/ul	0.01
43) Dimethylphthalate-d6	13.88	166	1129140	32.28	ng/ul	0.00
46) Acenaphthylene-d8	14.16	160	1407375	33.36	ng/ul	0.00
51) 4-Nitrophenol-d4	14.67	143	35745	5.30	ng/ul	0.00
57) Fluorene-d10	15.46	176	1037073	32.84	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.59	200	196081	28.34	ng/ul	0.00
70) Anthracene-d10	17.32	188	1732565	34.42	ng/ul	0.00
76) Pyrene-d10	19.62	212	2004916	36.74	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.57	264	1581766	33.96	ng/ul	0.00

S.J
 05/04/17

Target Compounds

Qvalue

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

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