

Quantitation Report (Qedit)

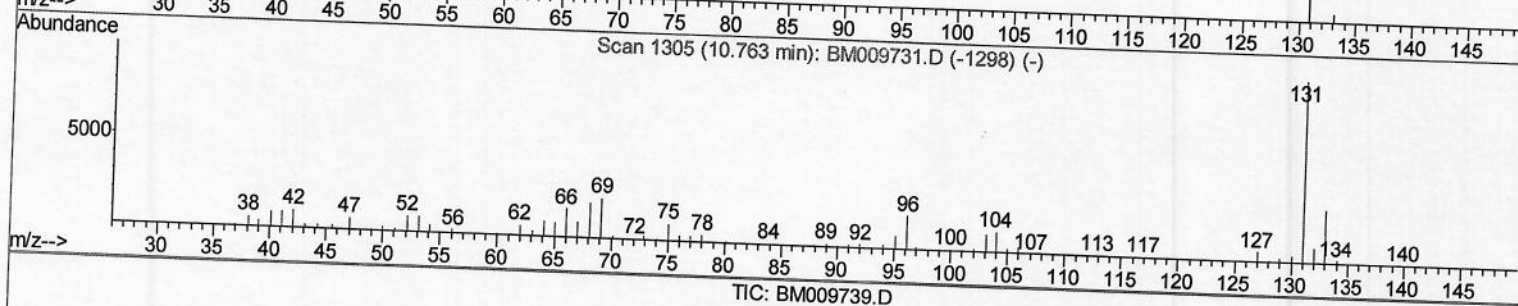
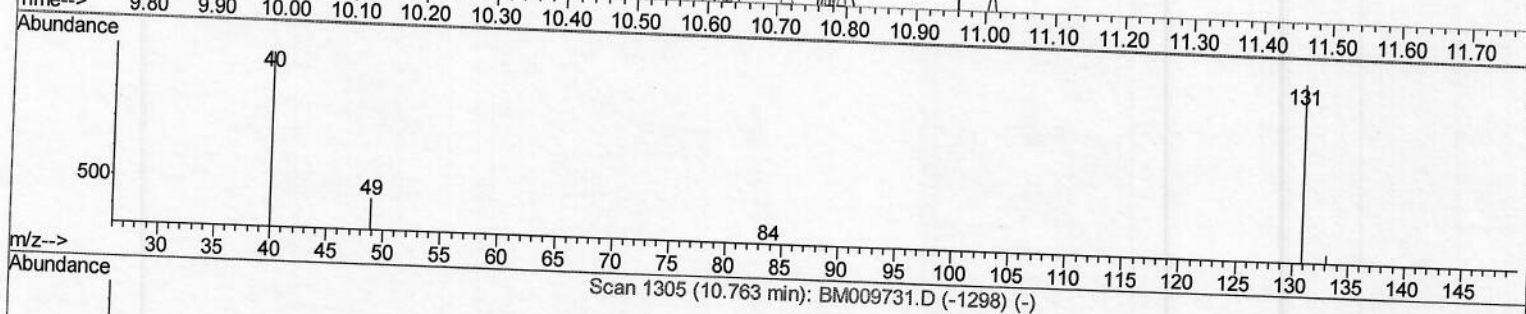
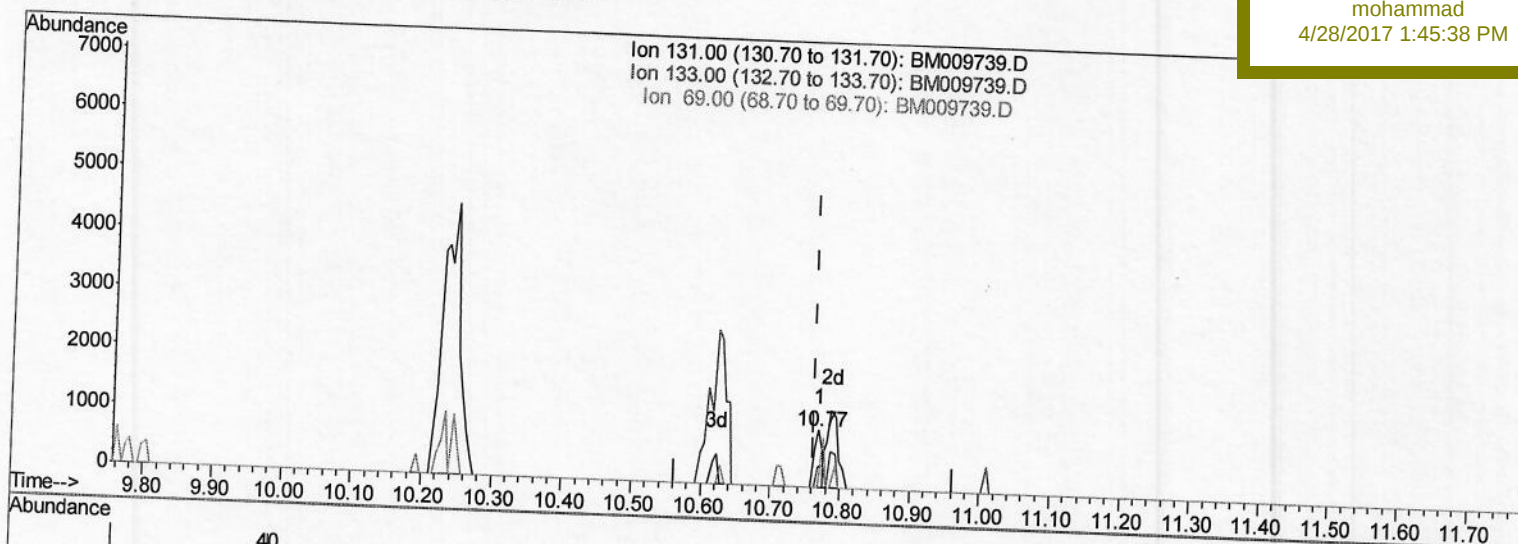
Data Path : Z:\HPCHEM1\BNA_M\Data\BM042717\
 Data File : BM009739.D
 Acq On : 27 Apr 2017 13:51
 Operator : SJ/MA
 Sample : I2809-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHTC5

Manual Integrations
 APPROVED

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Quant Time: Apr 27 18:25:46 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



(29) 4-Chloroaniline-d4 (S)
 10.769min (+0.006) 0.06ng/ul
 response 758

Ion	Exp%	Act%
131.00	100	100
133.00	31.50	38.20#
69.00	36.50	0.00#
0.00	0.00	0.00

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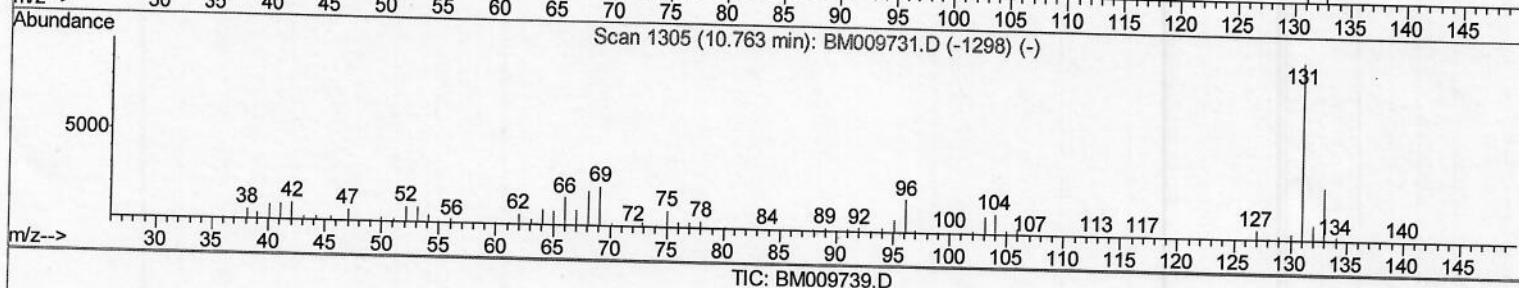
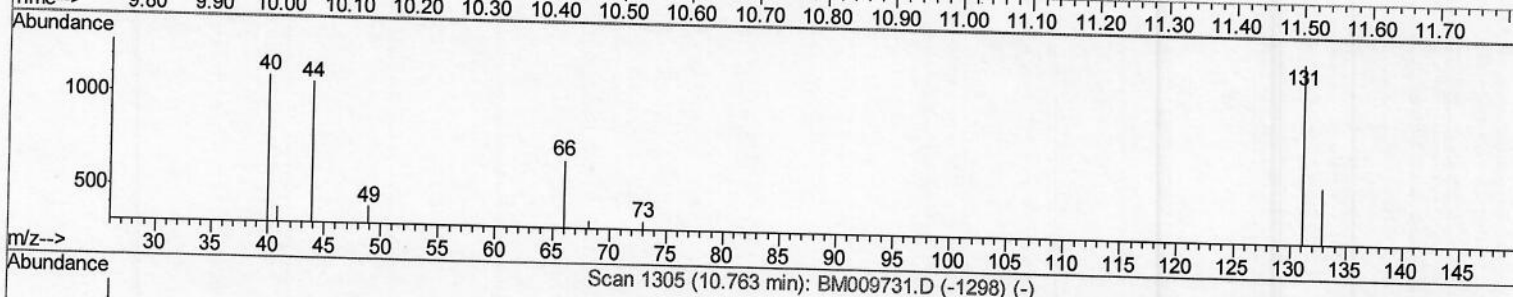
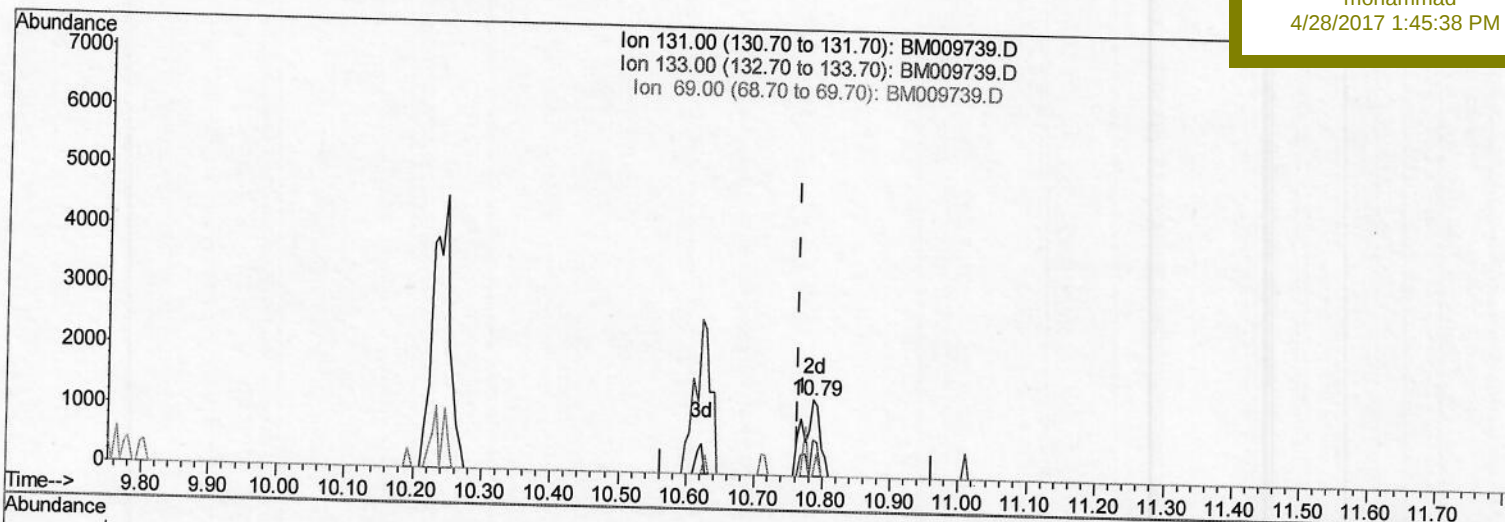
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(29) 4-Chloroaniline-d4 (S)

10.786min (+0.024) 0.18ng/ul m

response 2173

Ion	Exp%	Act%
131.00	100	100
133.00	31.50	47.03#
69.00	36.50	0.00#
0.00	0.00	0.00

> $\frac{S-J}{05/04/17}$

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	126716	20.00	ng/ul	0.00
18) Naphthalene-d8	10.62	136	577124	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.47	164	387257	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.22	188	933482	20.00	ng/ul	0.00
75) Chrysene-d12	21.40	240	1008609	20.00	ng/ul	0.00
83) Perylene-d12	23.71	264	789658	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.30	96	14919	5.42	ng/uL	0.00
5) Phenol-d5	6.99	99	83236	6.05	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.16	67	302434	32.22	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	227686	26.00	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	158520	14.77	ng/ul	0.00
19) Nitrobenzene-d5	8.99	128	157801	33.59	ng/ul	0.00
22) 2-Nitrophenol-d4	9.71	143	175411	35.49	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.24	165	283398	28.67	ng/ul	0.00
29) 4-Chloroaniline-d4	10.79	131	2173m	0.18	ng/ul	0.02
43) Dimethylphthalate-d6	13.88	166	1168229	34.68	ng/ul	0.00
46) Acenaphthylene-d8	14.16	160	1419391	34.94	ng/ul	0.00
51) 4-Nitrophenol-d4	14.67	143	30007	4.62	ng/ul	0.00
57) Fluorene-d10	15.46	176	1071215	35.23	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.59	200	197696	30.02	ng/ul	0.00
70) Anthracene-d10	17.32	188	1744443	36.40	ng/ul	0.00
76) Pyrene-d10	19.62	212	1959087	44.01	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.57	264	1437444	37.43	ng/ul	0.00

Target Compounds

Qvalue

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

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