

Quantitation Report (QT Reviewed)

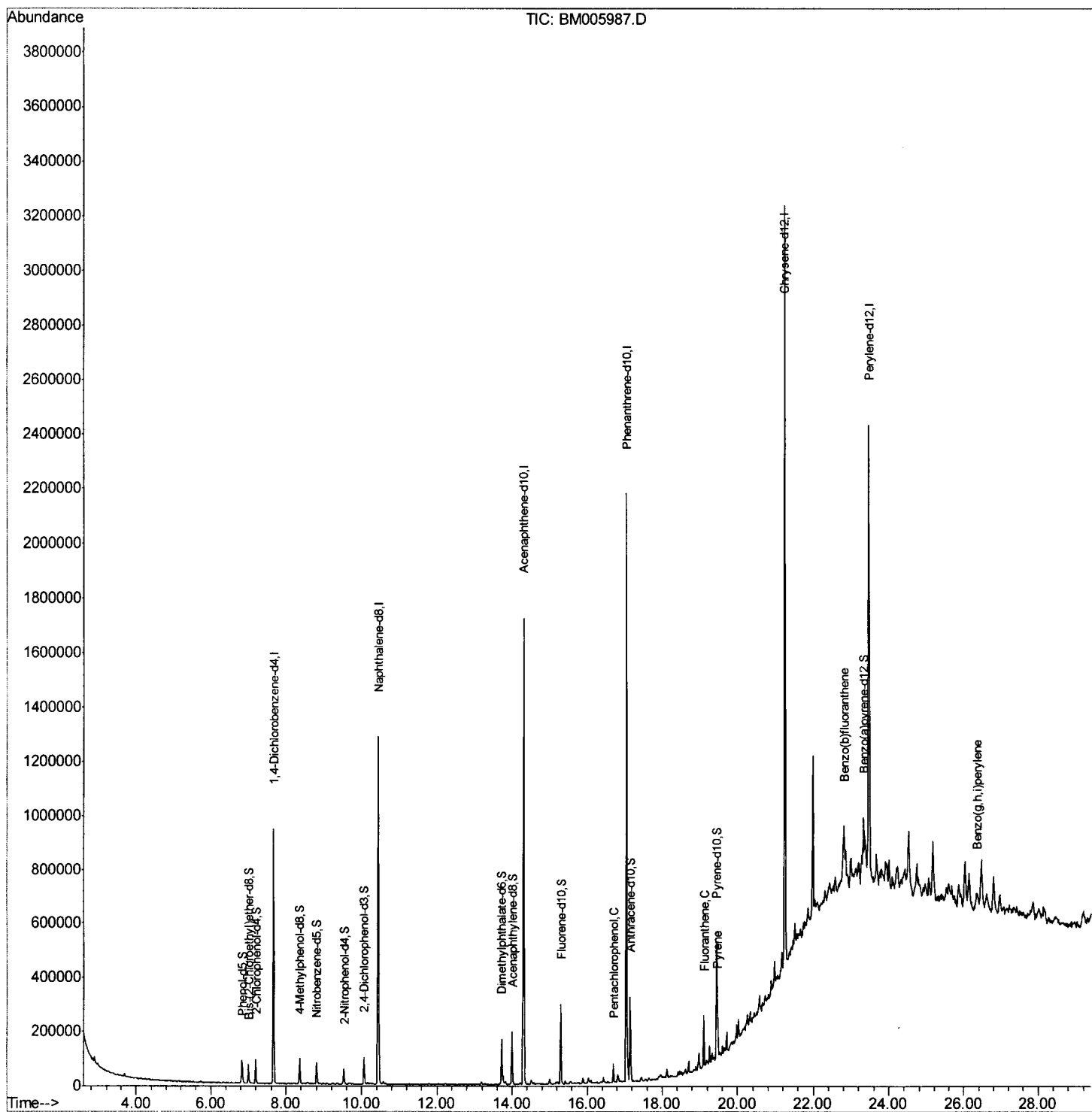
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062416\
 Data File : BM005987.D
 Acq On : 25 Jun 2016 04:01
 Operator : UM/SJ
 Sample : H3612-11 10X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Z29

Manual Integration

Quant Time: Jun 25 07:27:25 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 23:55:23 2016
 Response via : Initial Calibration

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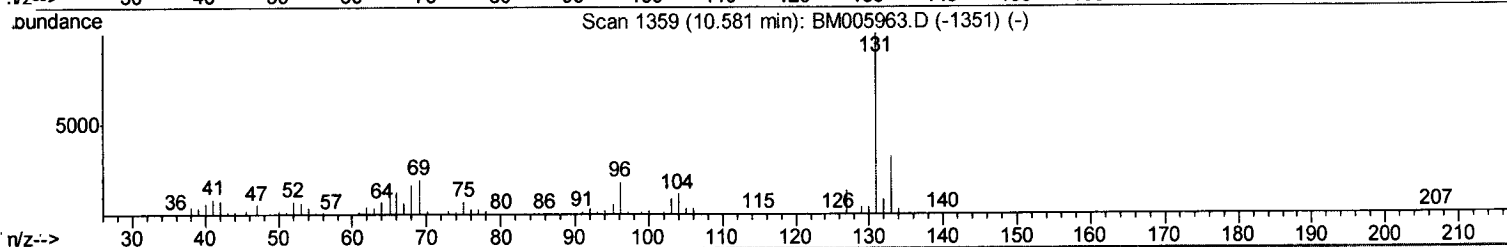
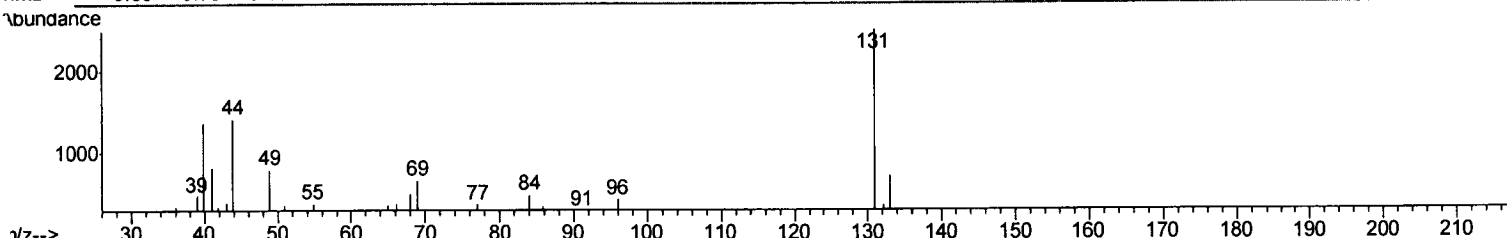
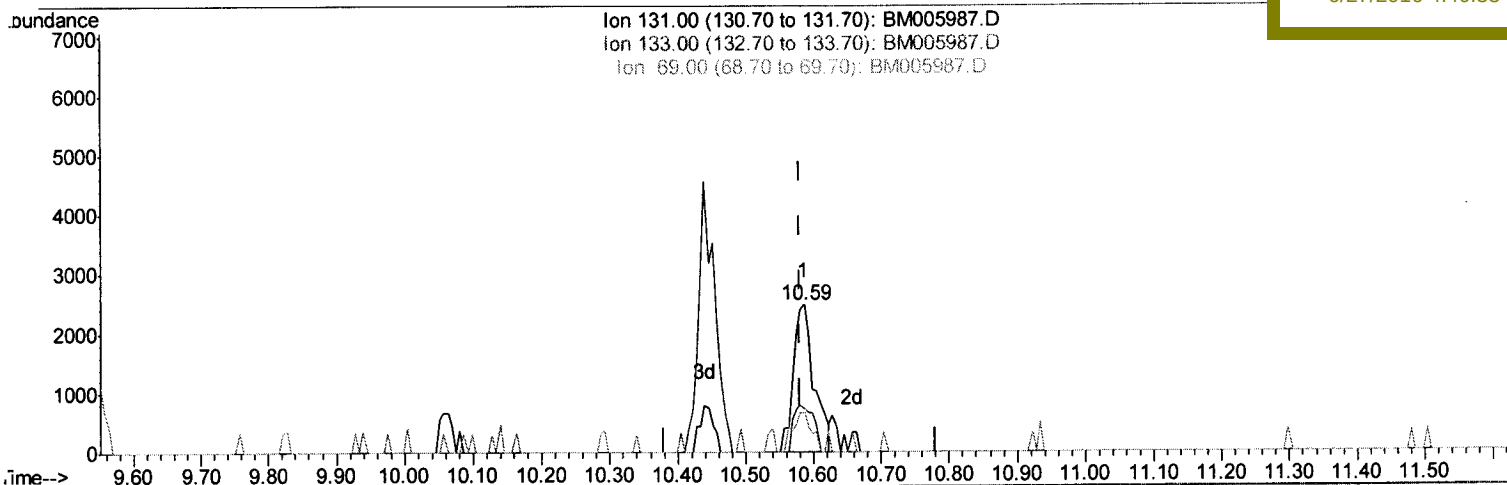
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TIC: BM005987.D

(29) 4-Chloroaniline-d4 (S)

10.586min (+0.006) 0.30ng/ul

response 5240

Ion	Exp%	Act%
131.00	100	100
133.00	32.00	28.91
69.00	19.30	26.70#
0.00	0.00	0.00

Quantitation Report (Qedit)

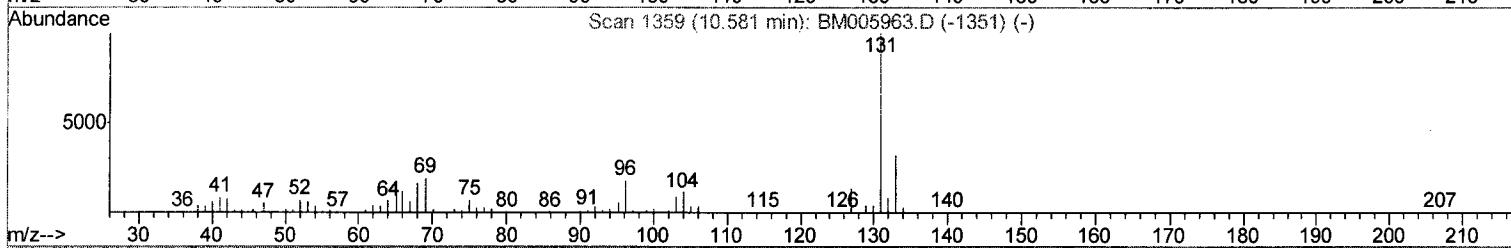
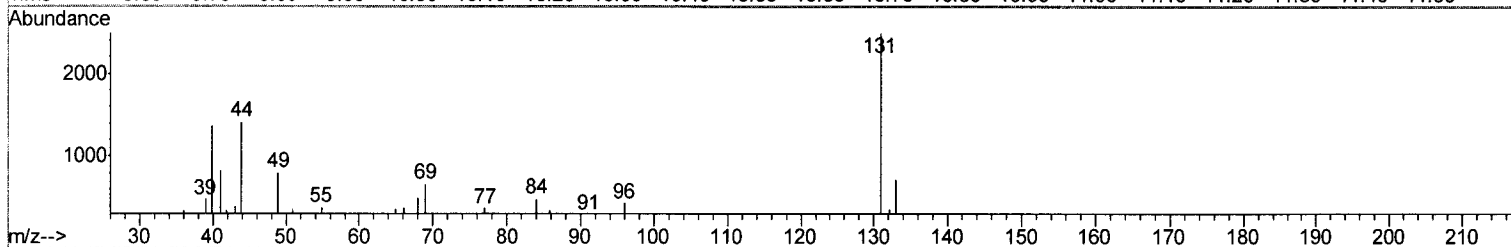
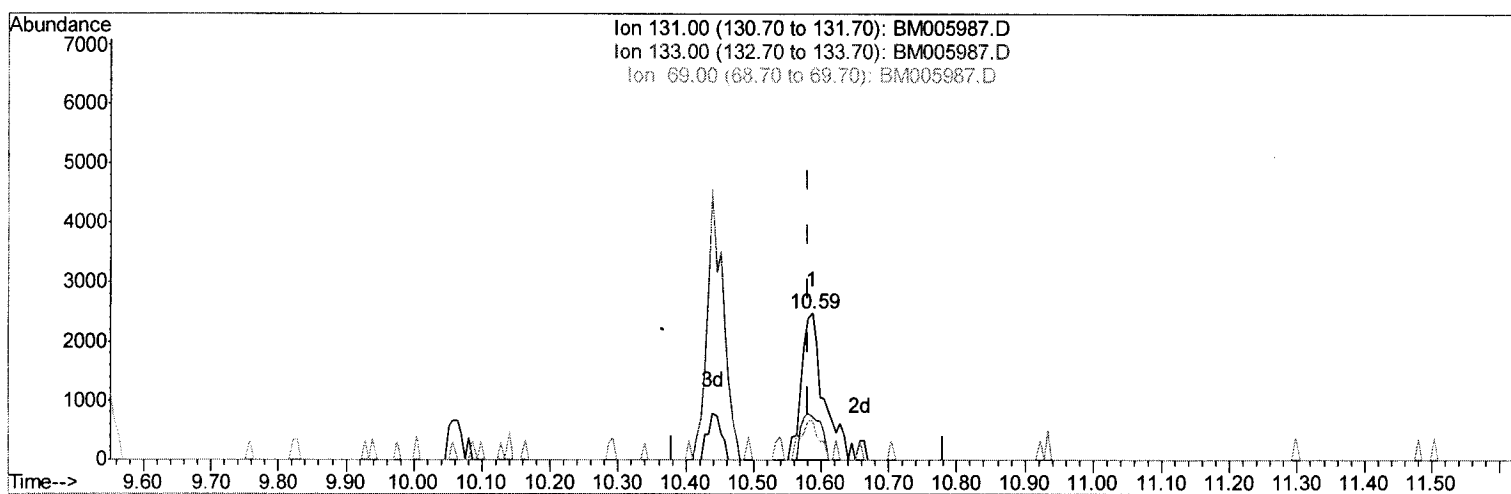
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Manual Integration

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TIC: BM005987.D

(29) 4-Chloroaniline-d4 (S)

10.586min (+0.006) 0.35ng/ul m

response 5952

Ion Exp% Act%

131.00 100 100

133.00 32.00 28.91

69.00 19.30 26.70#

0.00 0.00 0.00

UM
 06/28/16

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062416\
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	247424	20.00	ng/ul	0.00
18) Naphthalene-d8	10.44	136	1067796	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.31	164	561187	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	1237080	20.00	ng/ul	0.00
75) Chrysene-d12	21.24	240	1284268	20.00	ng/ul	0.00
83) Perylene-d12	23.46	264	1208648	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	766	0.12	ng/uL	0.00
5) Phenol-d5	6.83	99	46796	2.20	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.00	67	31786	2.44	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	36769	2.19	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	34478	2.07	ng/ul	0.00
19) Nitrobenzene-d5	8.81	128	17338	2.21	ng/ul	0.00
22) 2-Nitrophenol-d4	9.53	143	16632	2.02	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	34719	2.20	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	5952m	0.35	ng/ul	0.00
43) Dimethylphthalate-d6	13.72	166	111208	2.49	ng/ul	0.00
46) Acenaphthylene-d8	14.00	160	139388	2.50	ng/ul	0.00
51) 4-Nitrophenol-d4	14.50	143	3880	0.45	ng/ul	0.00
57) Fluorene-d10	15.30	176	109170	2.83	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.42	200	2386	0.39	ng/ul	0.00
70) Anthracene-d10	17.15	188	166265	2.95	ng/ul	0.00
76) Pyrene-d10	19.45	212	188806	3.17	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.32	264	164708	3.02	ng/ul	0.00

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Target Compounds					Qvalue
68) Pentachlorophenol	16.70	266	11663	1.38 ng/ul	93
74) Fluoranthene	19.12	202	109429	1.49 ng/ul	99
77) Pyrene	19.48	202	97865	1.25 ng/ul	98
85) Benzo(b)fluoranthene	22.80	252	130652	1.78 ng/ul#	92
91) Benzo(g,h,i)perylene	26.36	276	68660	1.05 ng/ul	96

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