

Quantitation Report (QT Reviewed)

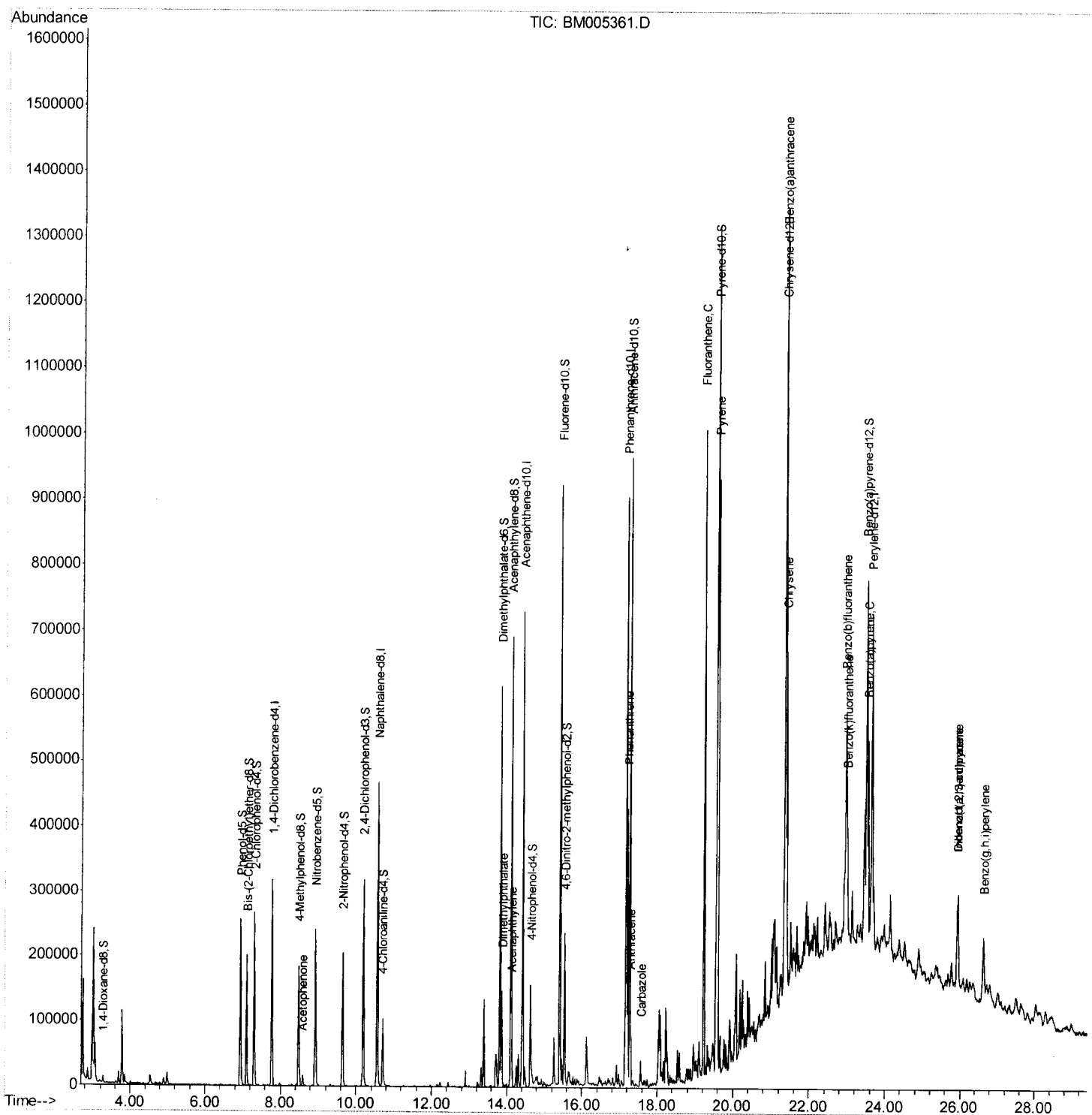
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM050916\
 Data File : BM005361.D
 Acq On : 09 May 2016 15:37
 Operator : UM/SJ
 Sample : H2888-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 D9WW2

Manual Integrations
 APPROVED

sohil
 5/10/2016 6:45:33 PM

Quant Time: May 10 05:06:04 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM050516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 10 04:23:52 2016
 Response via : Initial Calibration



Quantitation Report (Qedit)

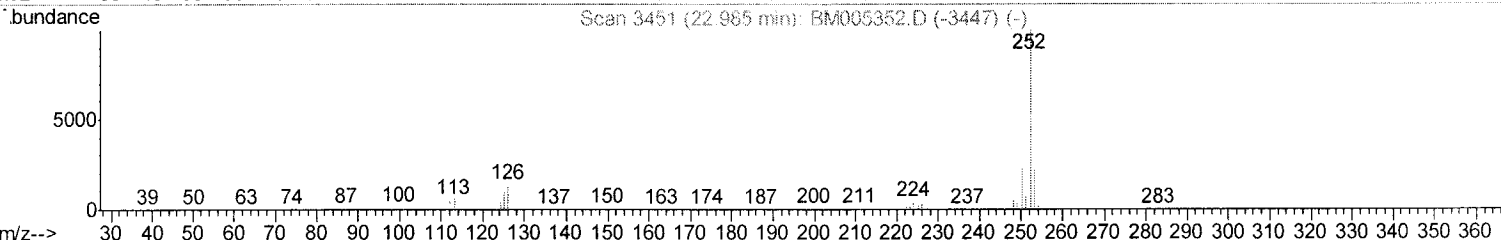
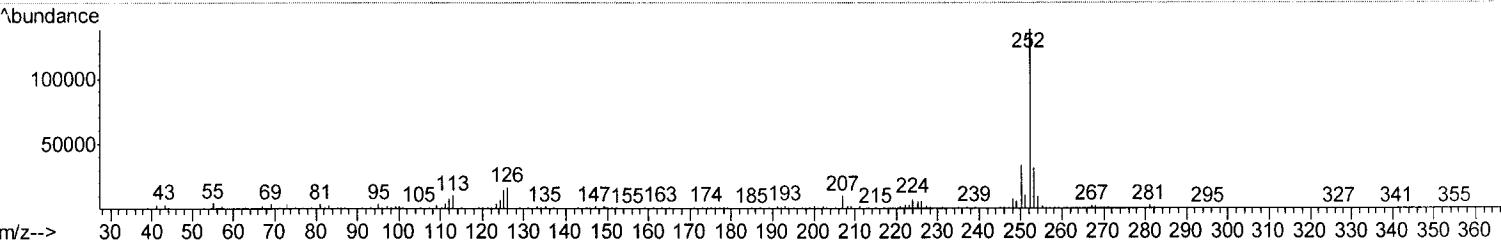
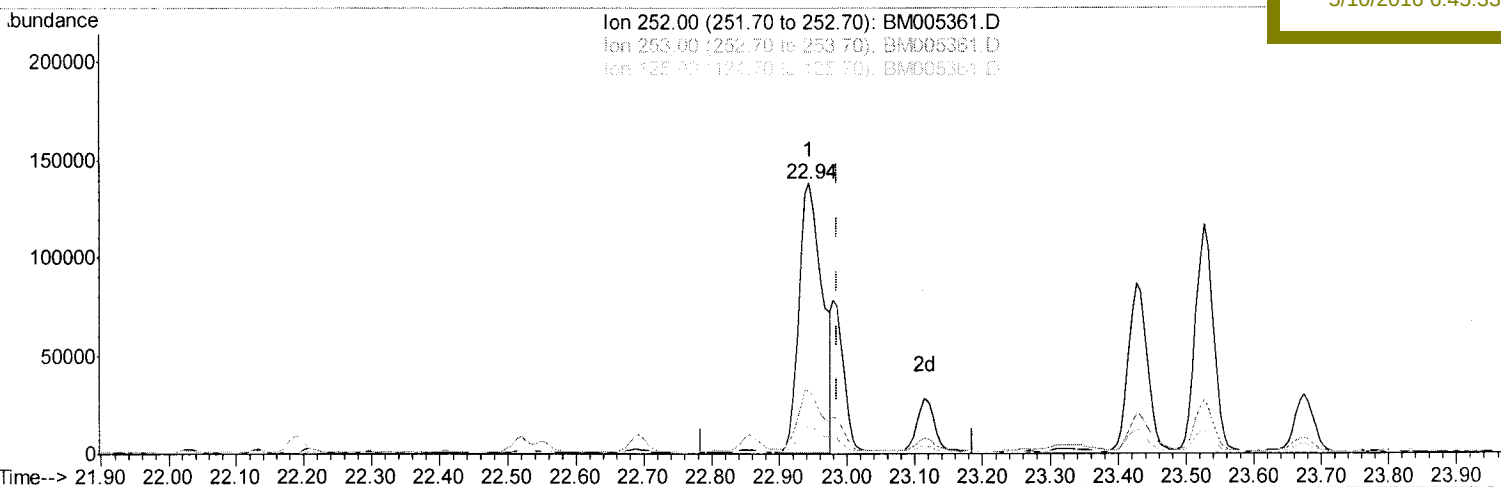
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Quant Time: May 10 05:04:51 2016
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TIC: BM005361.D

(86) Benzo(k)fluoranthene

22.944min (-0.041) 15.97ng/ul

response 327702

Ion	Exp%	Act%
252.00	100	100
253.00	21.60	23.06
125.00	10.00	10.43
0.00	0.00	0.00

Quantitation Report (Qedit)

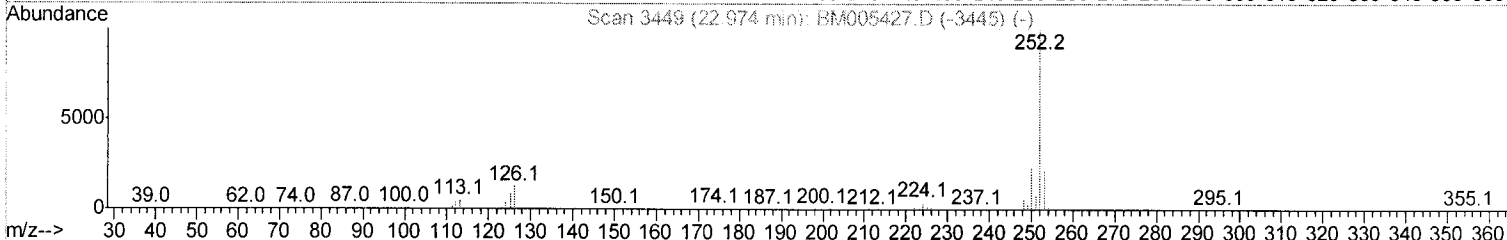
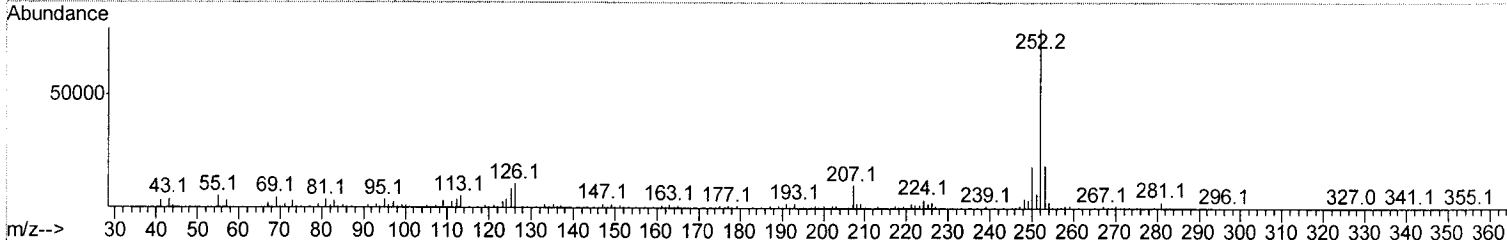
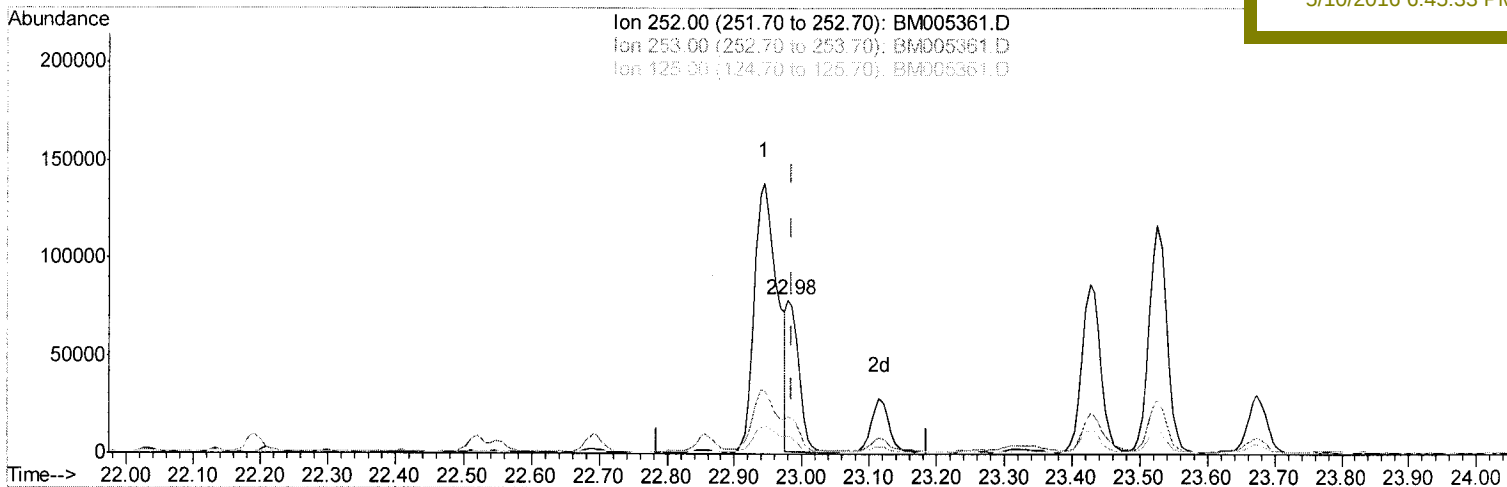
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TIC: BM005361.D

(86) Benzo(k)fluoranthene

22.980min (-0.006) 4.78ng/ul m U.M

response 98085

Ion	Exp%	Act%
252.00	100	100
253.00	21.60	24.32
125.00	10.00	11.19
0.00	0.00	0.00

05/18/18

Quantitation Report (Qedit)

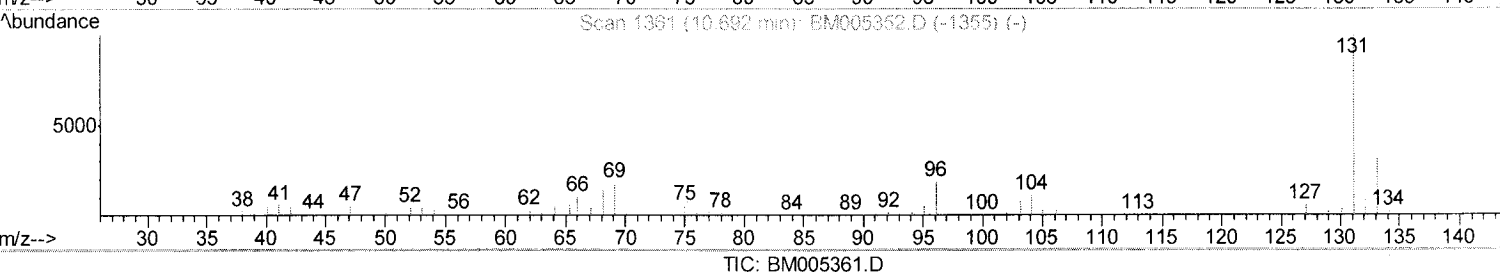
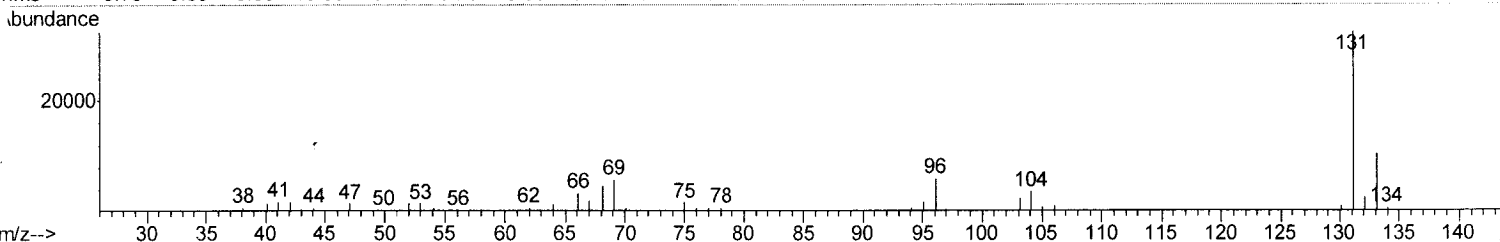
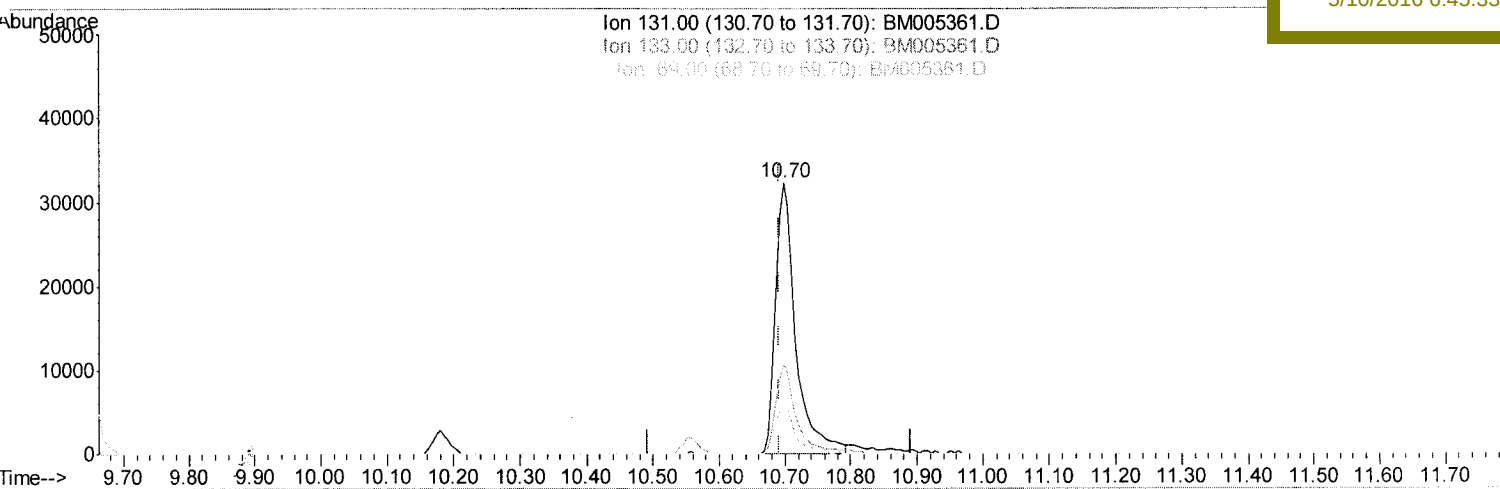
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Instrument :
BNA_M
ClientSampleId :
D9WW2

Manual Integrations
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5/10/2016 6:45:33 PM

Quant Time: May 10 04:25:32 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM050516.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 10 04:23:52 2016
Response via : Initial Calibration



(29) 4-Chloroaniline-d4 (S)

10.698min (+0.006) 9.48ng/ul

response 68430

Ion	Exp%	Act%
131.00	100	100
133.00	32.30	32.98
69.00	19.20	18.17
0.00	0.00	0.00

Quantitation Report (Qedit)

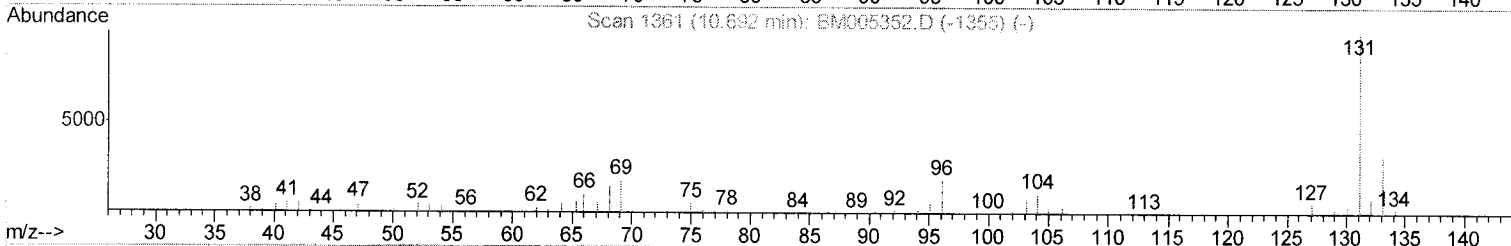
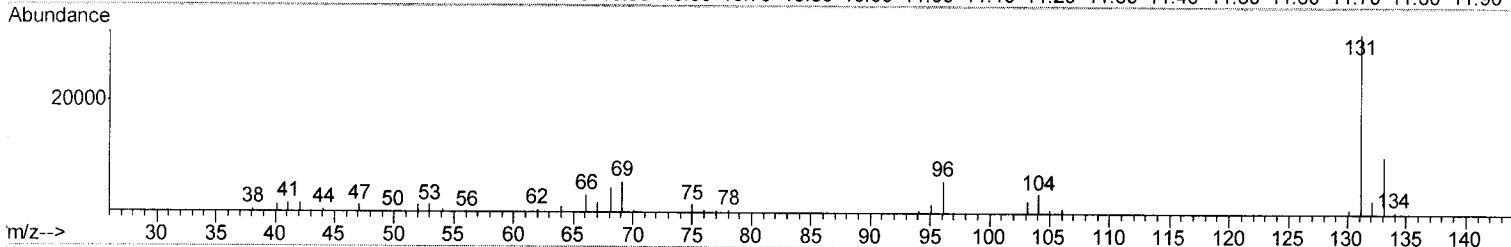
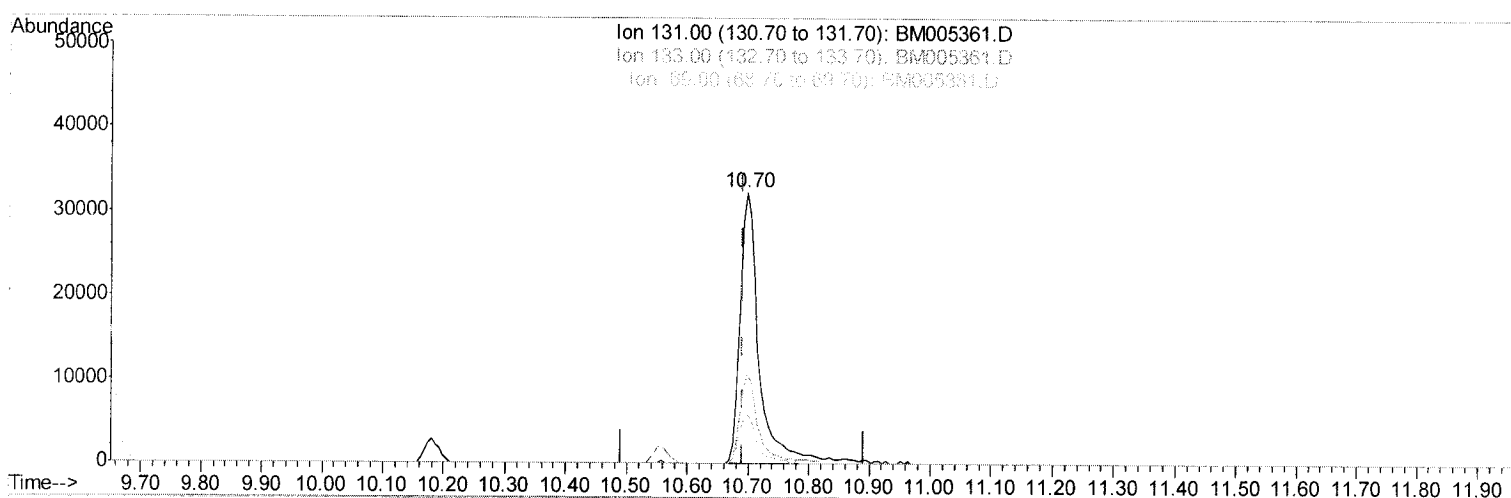
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 Operator : UM/SJ
 Sample : H2888-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9WW2

Manual Integrations
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TIC: BM005361.D

(29) 4-Chloroaniline-d4 (S)

10.698min (+0.006) 10.07ng/ul m *U.M*

response 72691

Ion	Exp%	Act%
131.00	100	100
133.00	32.30	32.98
69.00	19.20	18.17
0.00	0.00	0.00

05/18/16

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Instrument :
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 D9VW2

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Manual Integrations
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 5/10/2016 6:45:33 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	87282	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	399121	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.41	164	237610	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.16	188	528332	20.00	ng/ul	0.00
75) Chrysene-d12	21.35	240	474176	20.00	ng/ul	0.00
83) Perylene-d12	23.63	264	372834	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.27	96	5922	3.19	ng/uL	0.00
5) Phenol-d5	6.94	99	152438	19.26	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.10	67	89115	19.73	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	123683	20.69	ng/ul	0.00
13) 4-Methylphenol-d8	8.47	113	72844	11.13	ng/ul	0.00
19) Nitrobenzene-d5	8.93	128	60506	21.23	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	69485	21.54	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.18	165	125419	20.92	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	72691m	10.07	ng/ul	0.00
43) Dimethylphthalate-d6	13.82	166	416036	21.84	ng/ul	0.00
46) Acenaphthylene-d8	14.10	160	503071	22.52	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	48454	13.94	ng/ul	0.00
57) Fluorene-d10	15.40	176	367397	22.34	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.53	200	52531	17.67	ng/ul	0.00
70) Anthracene-d10	17.26	188	542005	23.21	ng/ul	0.00
76) Pyrene-d10	19.56	212	592547	27.07	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.48	264	382189	23.16	ng/ul	0.00

V.M
 05/18/16

Target Compounds

					Qvalue
14) Acetophenone	8.59	105	10255	1.06	ng/ul 96
44) Dimethylphthalate	13.86	163	86817	4.56	ng/ul 99
47) Acenaphthylene	14.13	152	29966	1.27	ng/ul 99
69) Phenanthrene	17.20	178	257343	9.26	ng/ul 99
71) Anthracene	17.29	178	68457	2.47	ng/ul 99
72) Carbazole	17.56	167	25899	1.06	ng/ul 96
74) Fluoranthene	19.22	202	606552	19.71	ng/ul 98
77) Pyrene	19.59	202	567459	20.63	ng/ul 97
80) Benzo(a)anthracene	21.34	228	290304	10.64	ng/ul 97
82) Chrysene	21.39	228	283016	10.94	ng/ul 97
85) Benzo(b)fluoranthene	22.94	252	327702	15.04	ng/ul 98
86) Benzo(k)fluoranthene	22.98	252	98085m	4.78	ng/ul 97
88) Benzo(a)pyrene	23.53	252	211036	10.24	ng/ul 97
89) Indeno(1,2,3-cd)pyrene	25.92	276	133639	5.94	ng/ul# 93
90) Dibenzo(a,h)anthracene	25.92	278	35866	1.91	ng/ul# 79
91) Benzo(g,h,i)perylene	26.62	276	119726	6.29	ng/ul 97

V.M
 05/18/16

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