Quantitation Report (Qedit)

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061816\
Data File : BM005871.D
Acq On : 17 Jun 2016 21:11
Operator : UM/SJ
Sample : H3535-21
Misc :
ALS Vial : 24 Sample Multiplier: 1

Quant Time: Jun 17 23:19:08 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061716.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 17 10:48:37 2016
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
D9Y11

Manual Integrations APPROVED
umangi
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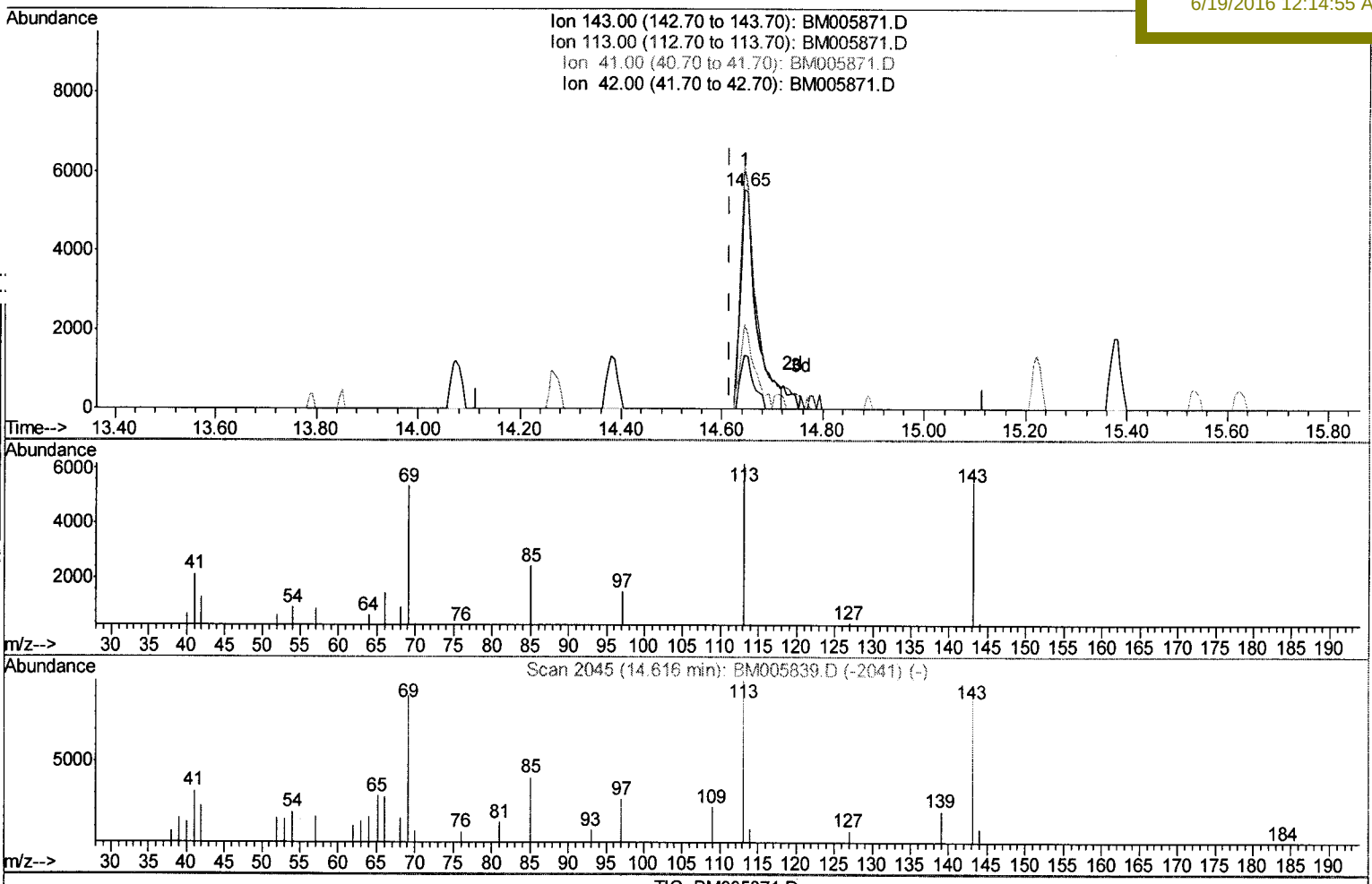

TIC: BM005871.D

(20) 4-Nitrophenol-d4 (S)

14.645min (+0.030) 15.82ng/ul

response 11963

Ion	Exp%	Act%
143.00	100	100
113.00	97.50	111.31
41.00	32.50	38.33
42.00	23.00	24.07

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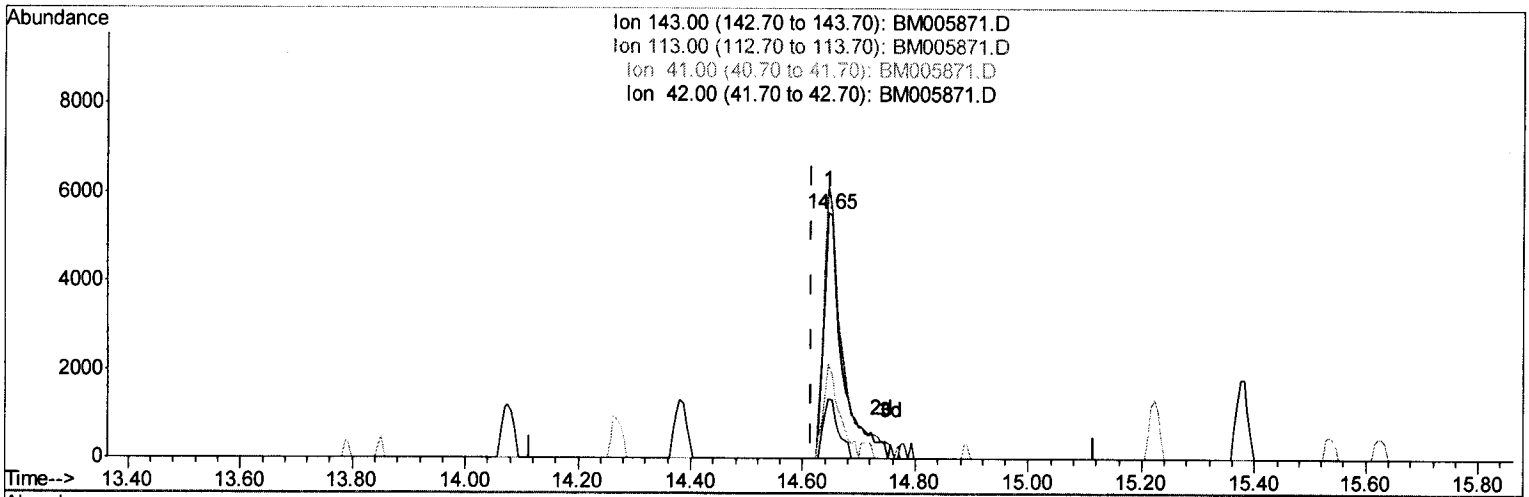

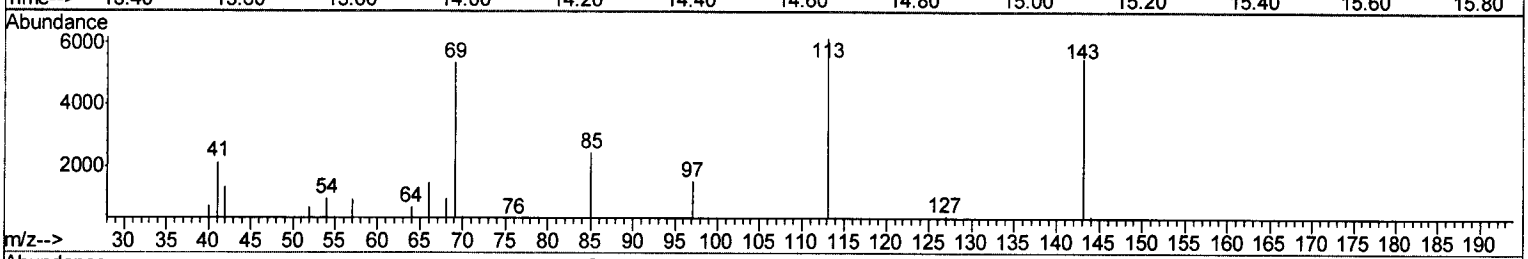

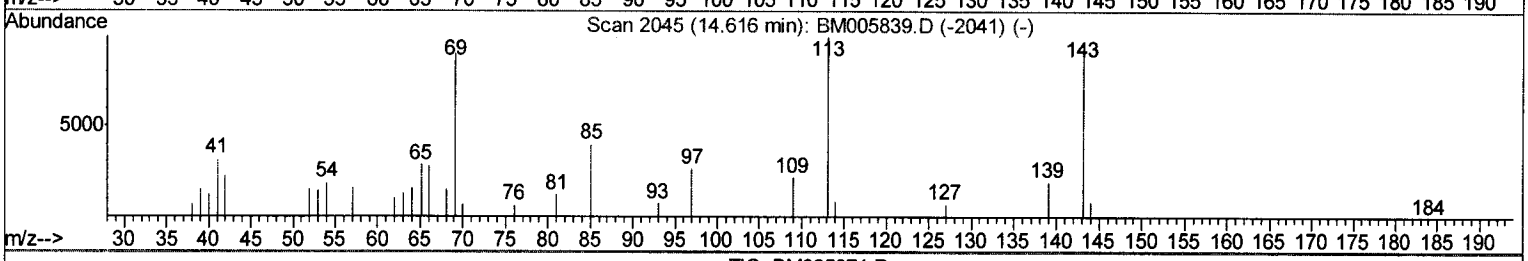

TIC: BM005871.D

(20) 4-Nitrophenol-d4 (S)

14.645min (+0.030) 16.92ng/ul m > U.M 06/20/16

response 12794

Ion	Exp%	Act%
143.00	100	100
113.00	97.50	111.31
41.00	32.50	38.33
42.00	23.00	24.07

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Instrument :
BNA_M
ClientSampleId :
D9Y11

Manual Integrations APPROVED
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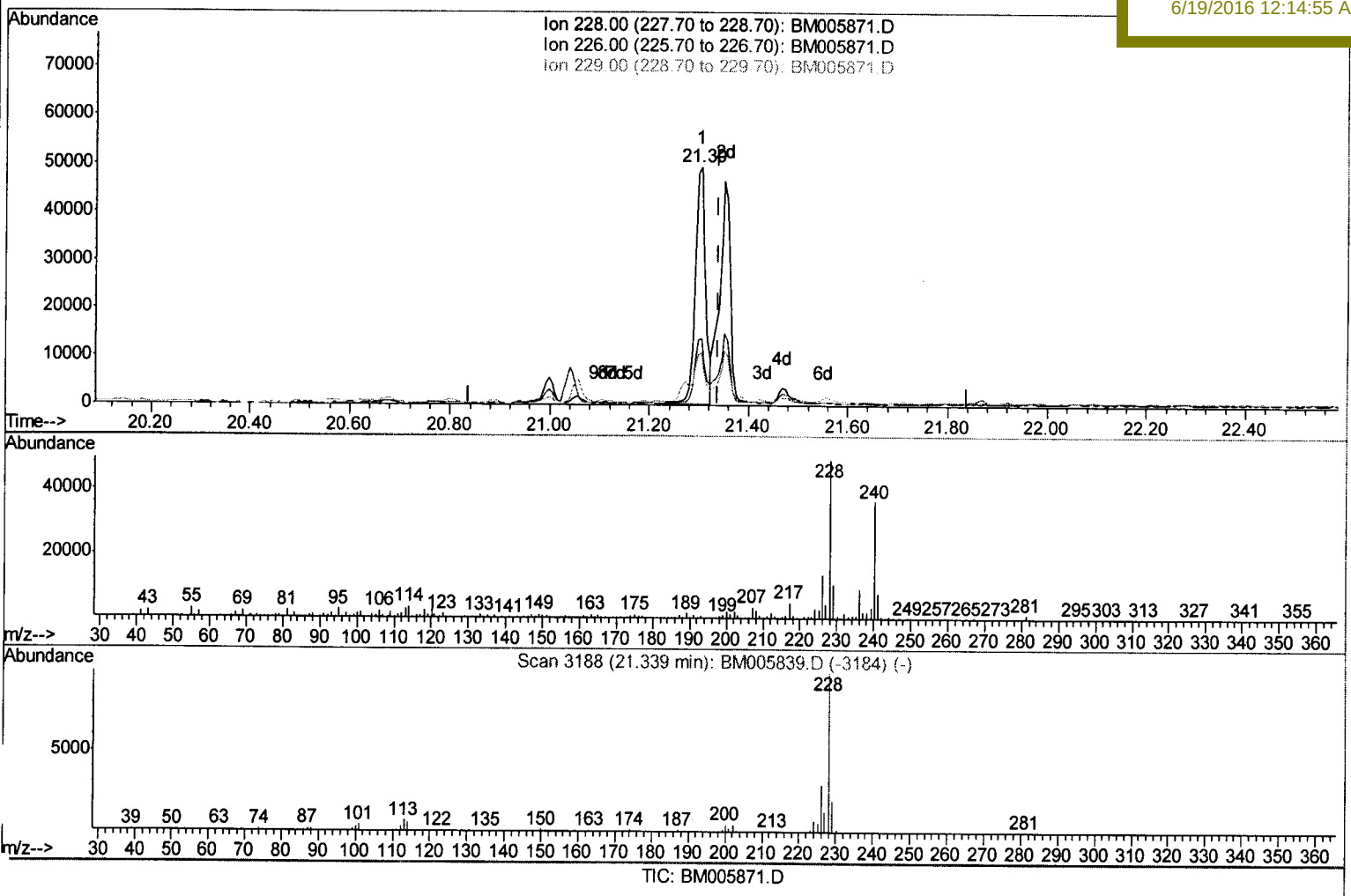

TIC: BM005871.D

(33) Chrysene

21.304min (-0.035) 8.05ng/ul

response 70794

Ion	Exp%	Act%
228.00	100	100
226.00	29.40	27.92
229.00	19.20	21.74
0.00	0.00	0.00

Quantitation Report (Qedit)

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ClientSampleId :
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Manual Integrations APPROVED
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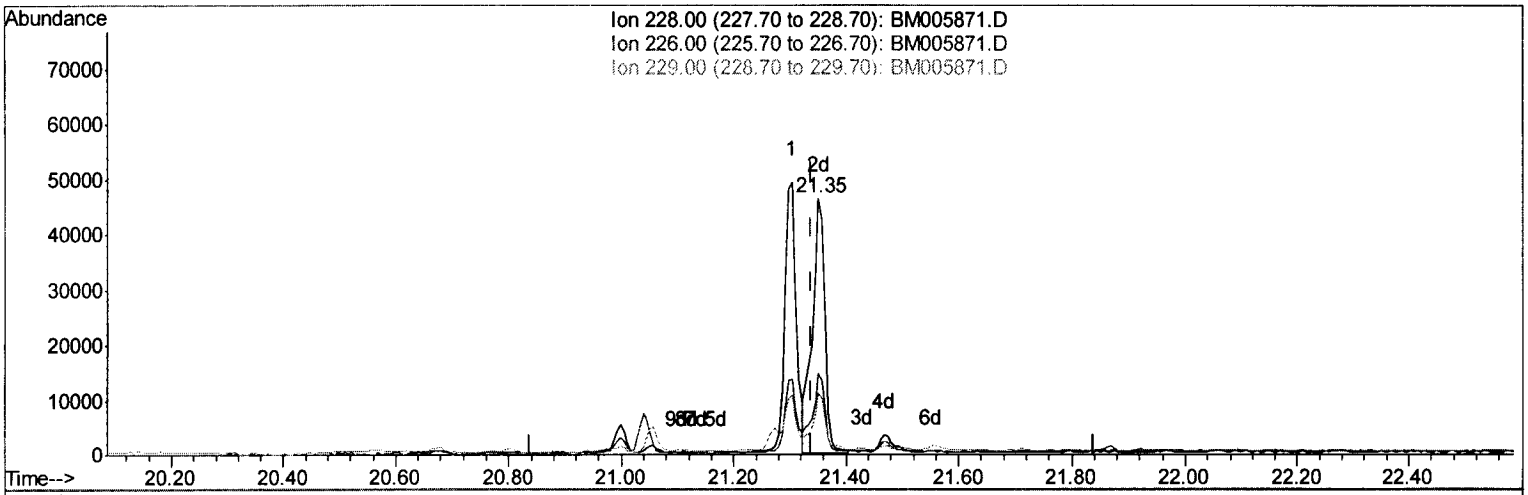

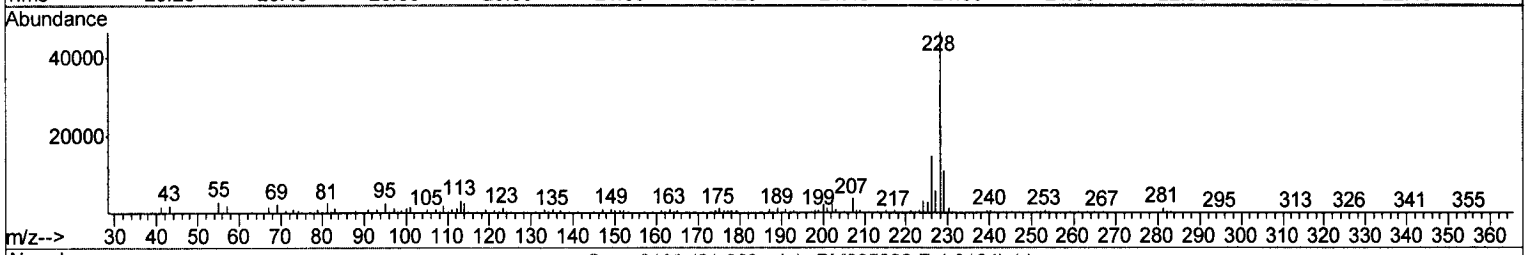

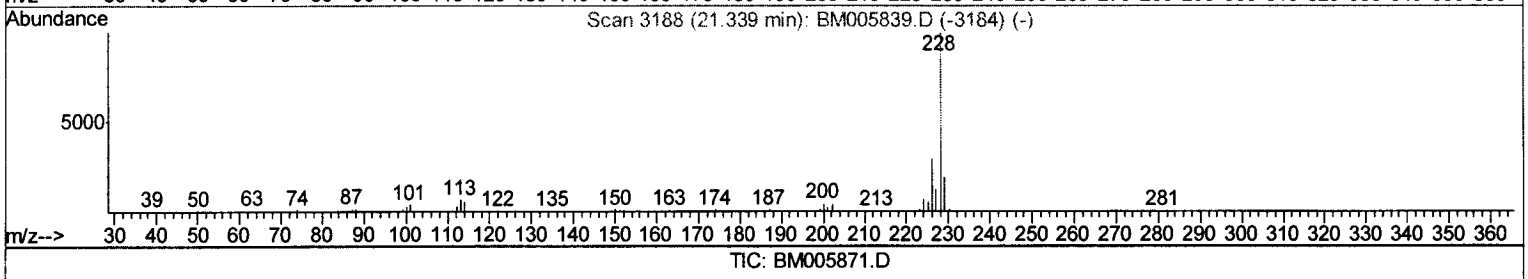

TIC: BM005871.D

(33) Chrysene

21.351min (+0.012) 8.32ng/ul m > U.M 06/20/16

response 73141

Ion	Exp%	Act%
228.00	100	100
226.00	29.40	31.96
229.00	19.20	23.97#
0.00	0.00	0.00

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Misc :
ALS Vial : 24 Sample Multiplier: 1

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Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
D9Y11

Manual Integrations APPROVED
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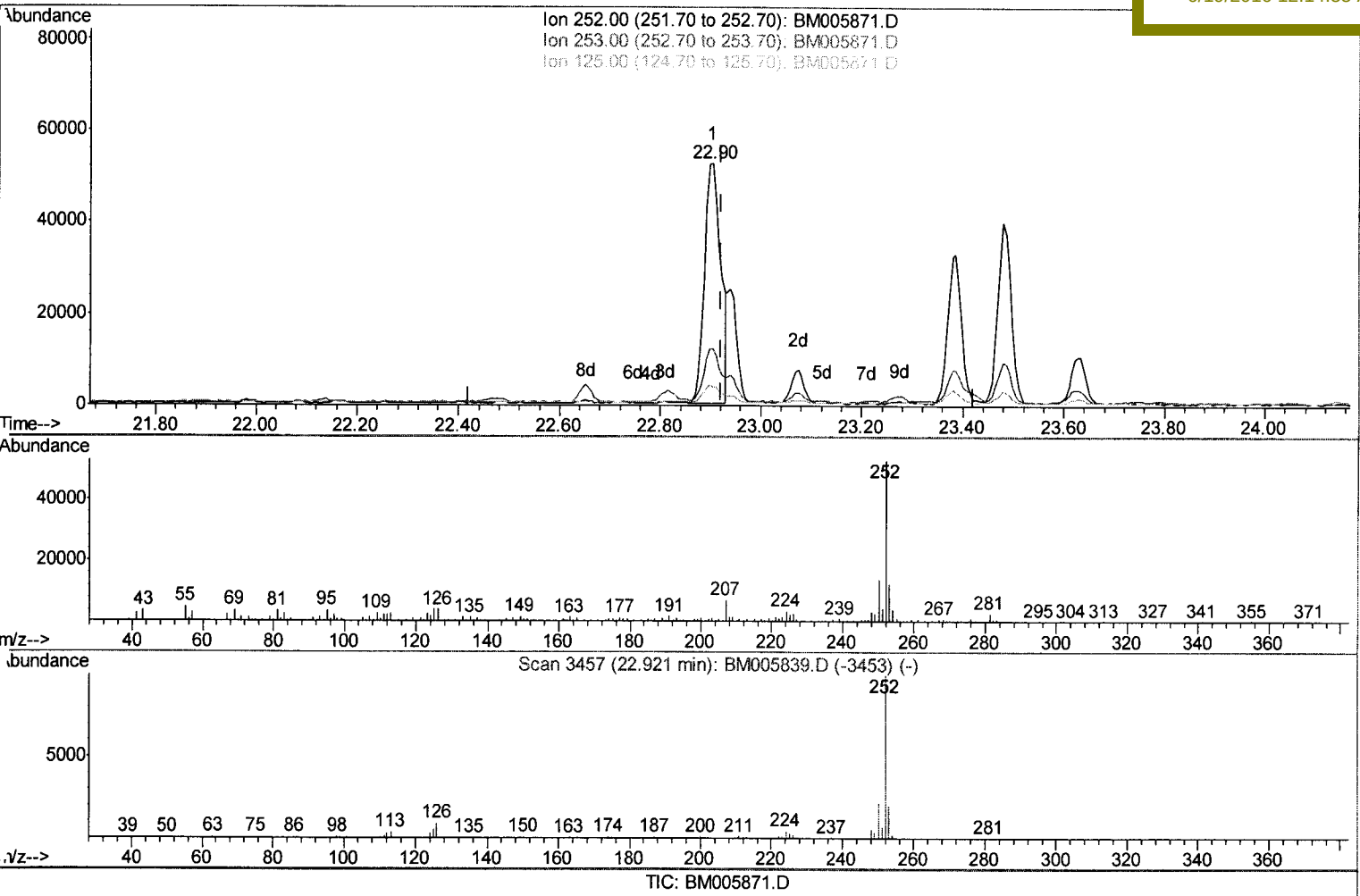

TIC: BM005871.D

(36) Benzo(k)fluoranthene

22.904min (-0.017) 13.90ng/ul

response 114671

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	23.65
125.00	6.00	7.85#
0.00	0.00	0.00

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Sample : H3535-21
Misc :
ALS Vial : 24 Sample Multiplier: 1

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BNA_M
ClientSampleId :
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Manual Integrations APPROVED
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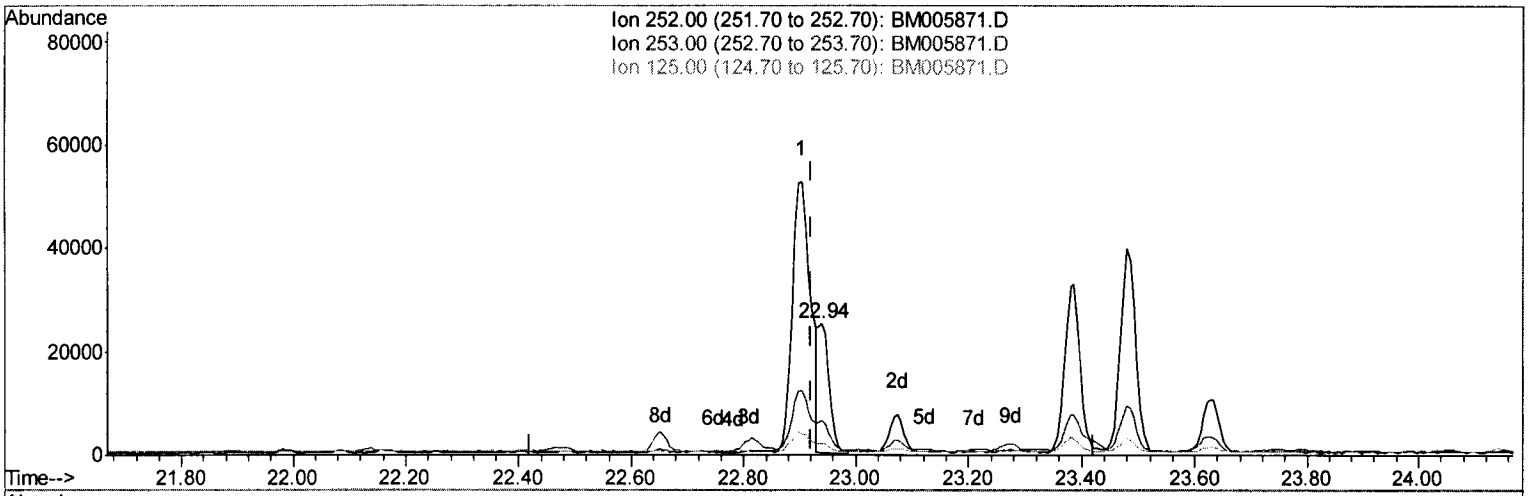

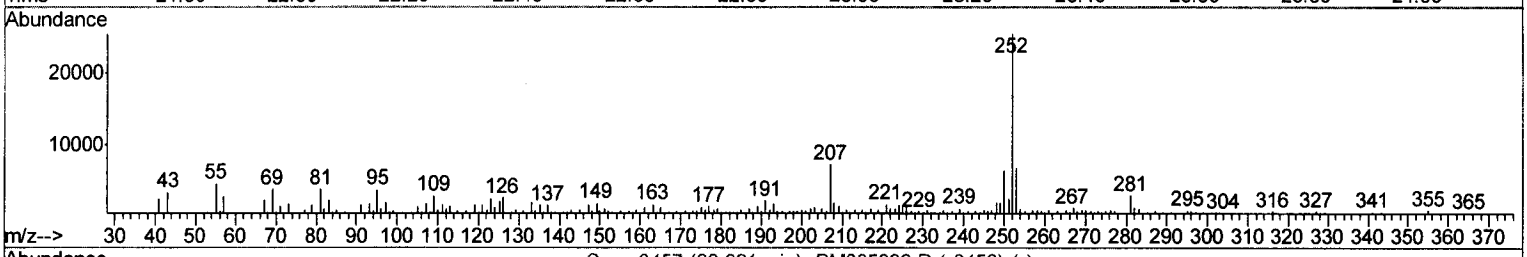

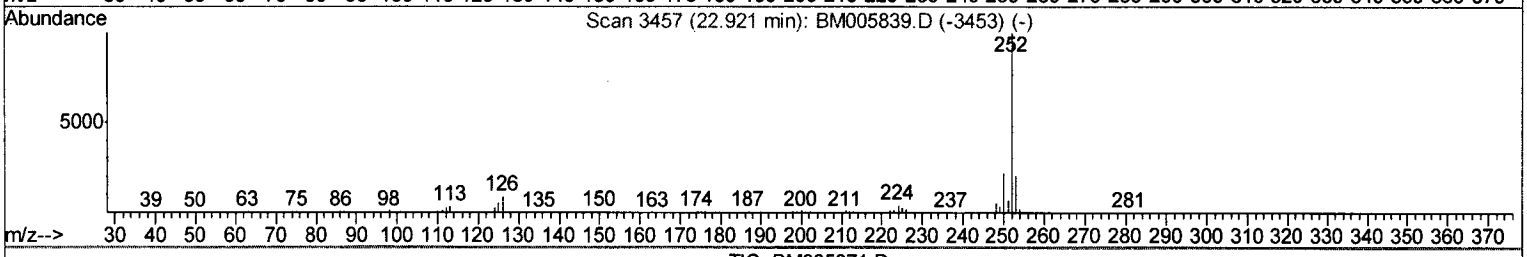

TIC: BM005871.D

(36) Benzo(k)fluoranthene

22.939min (+0.018) 4.42ng/ul m

U.M 06/20/16

response 36464

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	25.91
125.00	6.00	8.15#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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ata File : BM005871.D
:q On : 17 Jun 2016 21:11
perator : UM/SJ
mple : H3535-21
isc :
LS Vial : 24 Sample Multiplier: 1

uant Time: Jun 17 23:46:11 2016
uant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061716.M
uant Title : SVOA CALIBRATION
Last Update : Fri Jun 17 10:48:37 2016
esponse via : Initial Calibration

Instrument : BNA_M
ClientSampleId : D9Y11

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	24538	20.00	ng/ul	0.00
7) Naphthalene-d8	10.53	136	106626	20.00	ng/ul	0.01
15) Acenaphthene-d10	14.38	164	66333	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.13	188	157219	20.00	ng/ul	0.01
29) Chrysene-d12	21.32	240	146142	20.00	ng/ul	0.01
34) Perylene-d12	23.58	264	137229	20.00	ng/ul	0.02
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.18	96	597	3.15	ng/uL	0.00
3) Phenol-d5	6.93	99	44681	21.75	ng/ul	0.02
4) Bis-(2-Chloroethyl)ether-d	7.07	67	22560	21.19	ng/ul	0.00
5) 2-Chlorophenol-d4	7.28	132	40985	23.92	ng/ul	0.01
6) 4-Methylphenol-d8	8.46	113	38355	20.89	ng/ul	0.01
8) Nitrobenzene-d5	8.90	128	19468	25.87	ng/ul	0.00
9) 2-Nitrophenol-d4	9.62	143	22534	24.50	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	44405	24.45	ng/ul	0.01
12) 4-Chloroaniline-d4	10.68	131	18142	14.76	ng/ul	0.00
16) Dimethylphthalate-d6	13.79	166	145743	23.22	ng/ul	0.00
17) Acenaphthylene-d8	14.07	160	179358	24.81	ng/ul	0.00
20) 4-Nitrophenol-d4	14.65	143	12794m	16.92	ng/ul	0.03
21) Fluorene-d10	15.38	176	130378	24.38	ng/ul	0.01
24) 4,6-Dinitro-2-methylphenol	15.53	200	8946	8.43	ng/ul	0.02
26) Anthracene-d10	17.23	188	200036	24.37	ng/ul	0.01
30) Pyrene-d10	19.53	212	210732	30.42	ng/ul	0.01
37) Benzo(a)pyrene-d12	23.44	264	170298	24.82	ng/ul	0.02
Target Compounds						Qvalue
25) Phenanthrene	17.17	178	38281	4.09	ng/ul	99
27) Anthracene	17.26	178	9722	1.02	ng/ul	97
28) Fluoranthene	19.19	202	119810	10.19	ng/ul	98
31) Pyrene	19.56	202	110693	12.82	ng/ul	98
32) Benzo(a)anthracene	21.30	228	70794	7.76	ng/ul	97
33) Chrysene	21.35	228	73141m	8.32	ng/ul	
35) Benzo(b)fluoranthene	22.90	252	114671	13.42	ng/ul#	95
36) Benzo(k)fluoranthene	22.94	252	36464m	4.42	ng/ul	
38) Benzo(a)pyrene	23.48	252	73723	8.79	ng/ul#	96
39) Indeno(1,2,3-cd)pyrene	25.87	276	58126	5.67	ng/ul	97
40) Dibenzo(a,h)anthracene	25.87	278	16007	1.90	ng/ul#	69
41) Benzo(g,h,i)perylene	26.59	276	58997	6.76	ng/ul	97

U.M 06/20/16

U.M 06/20/16

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

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