

Quantitation Report (Qedit)

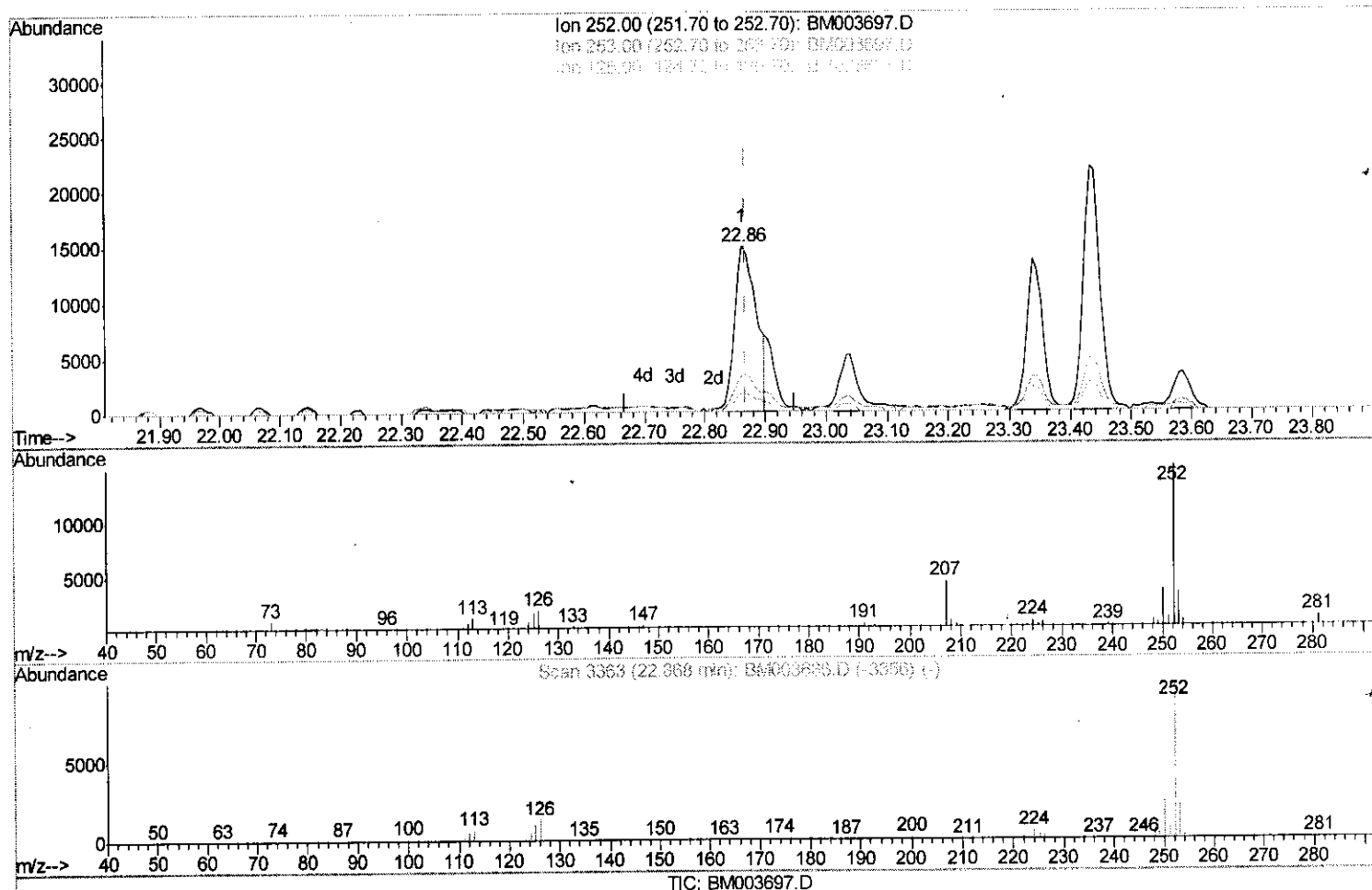
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM122115\
 Data File : BM003697.D
 Acq On : 22 Dec 2015 03:41
 Operator : UM/SJ
 Sample : G4801-17DL 30X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 Client Sampled :
 G9C65DL

Manual Integrations
 APPROVED

apatel
 12/22/2015 4:50:07 PM

Quant Time: Dec 22 04:39:41 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM121915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 22 02:40:18 2015
 Response via : Initial Calibration



(88) Benzo(b)fluoranthene

22.862min (-0.006) 4.13ng/ul m

response 37551

Ion Exp% Act%

252.00 100 100

253.00 21.90 22.05

125.00 10.60 11.68

0.00 0.00 0.00

> U.M
 12/24/15

Quantitation Report (Qedit)

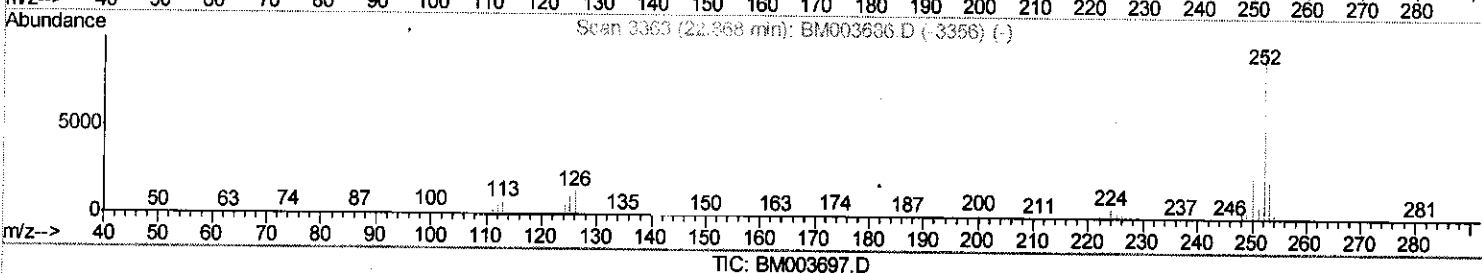
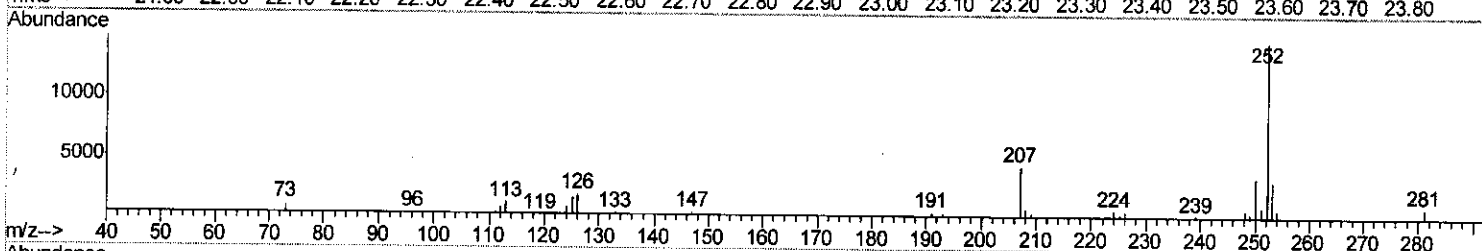
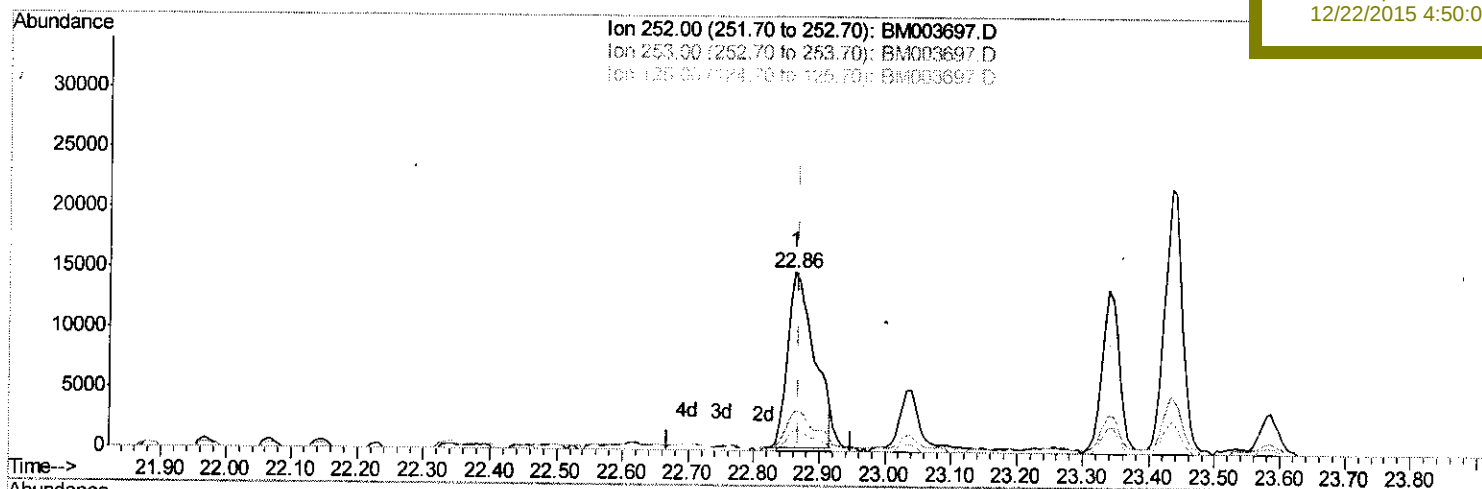
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(88) Benzo(b)fluoranthene

22.862min (-0.006) 4.56ng/ui

response 41406

Ion	Exp%	Act%
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252.00	100	100
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253.00	21.90	22.05
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125.00	10.60	11.68
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0.00	0.00	0.00
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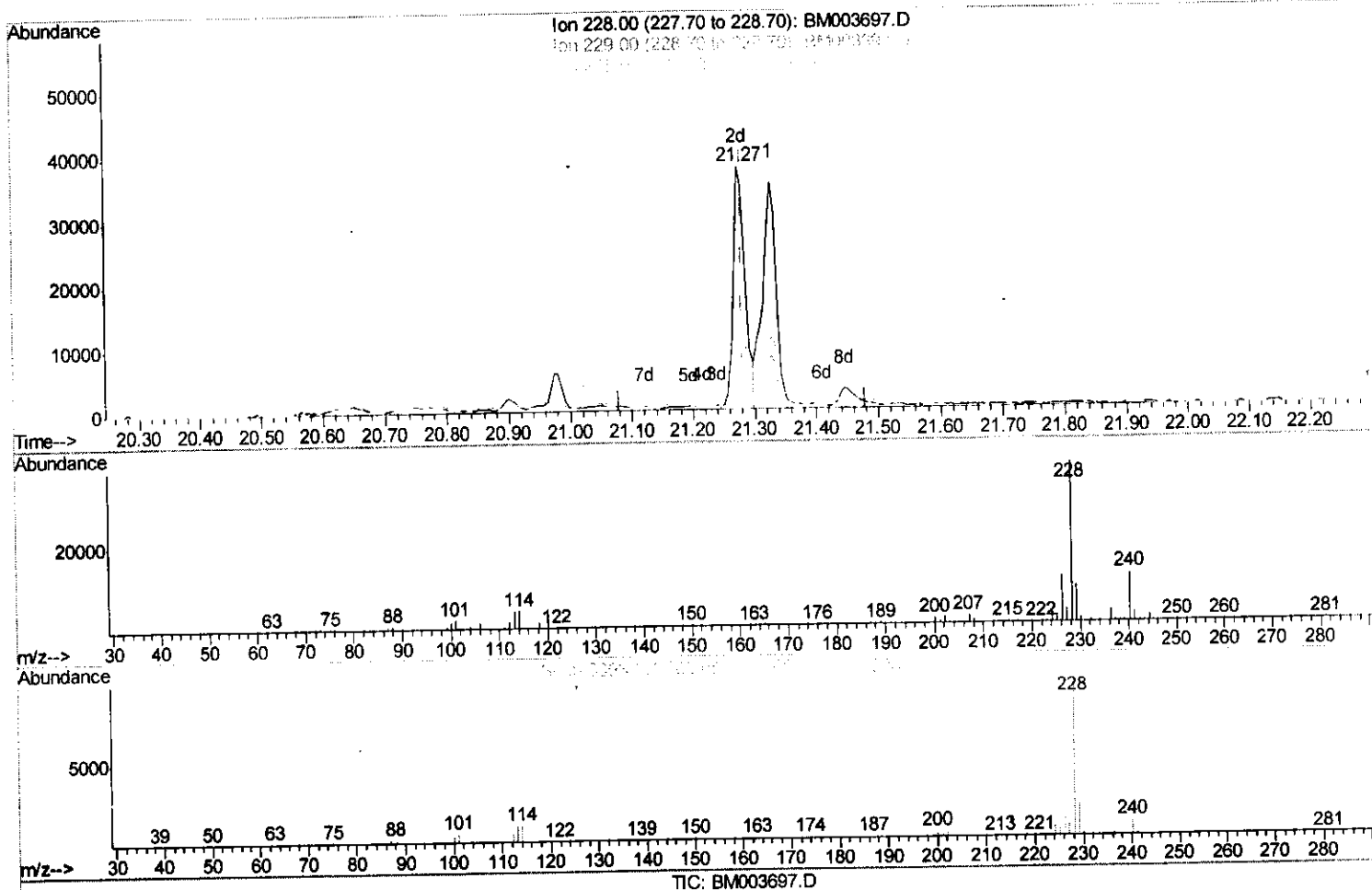
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(83) Benzo(a)anthracene

21.274min (-0.006) 5.53ng/ul m

response 50969

Ion	Exp%	Act%
228.00	100	100
229.00	19.60	23.92#
226.00	27.20	30.11
0.00	0.00	0.00

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 12/24/15

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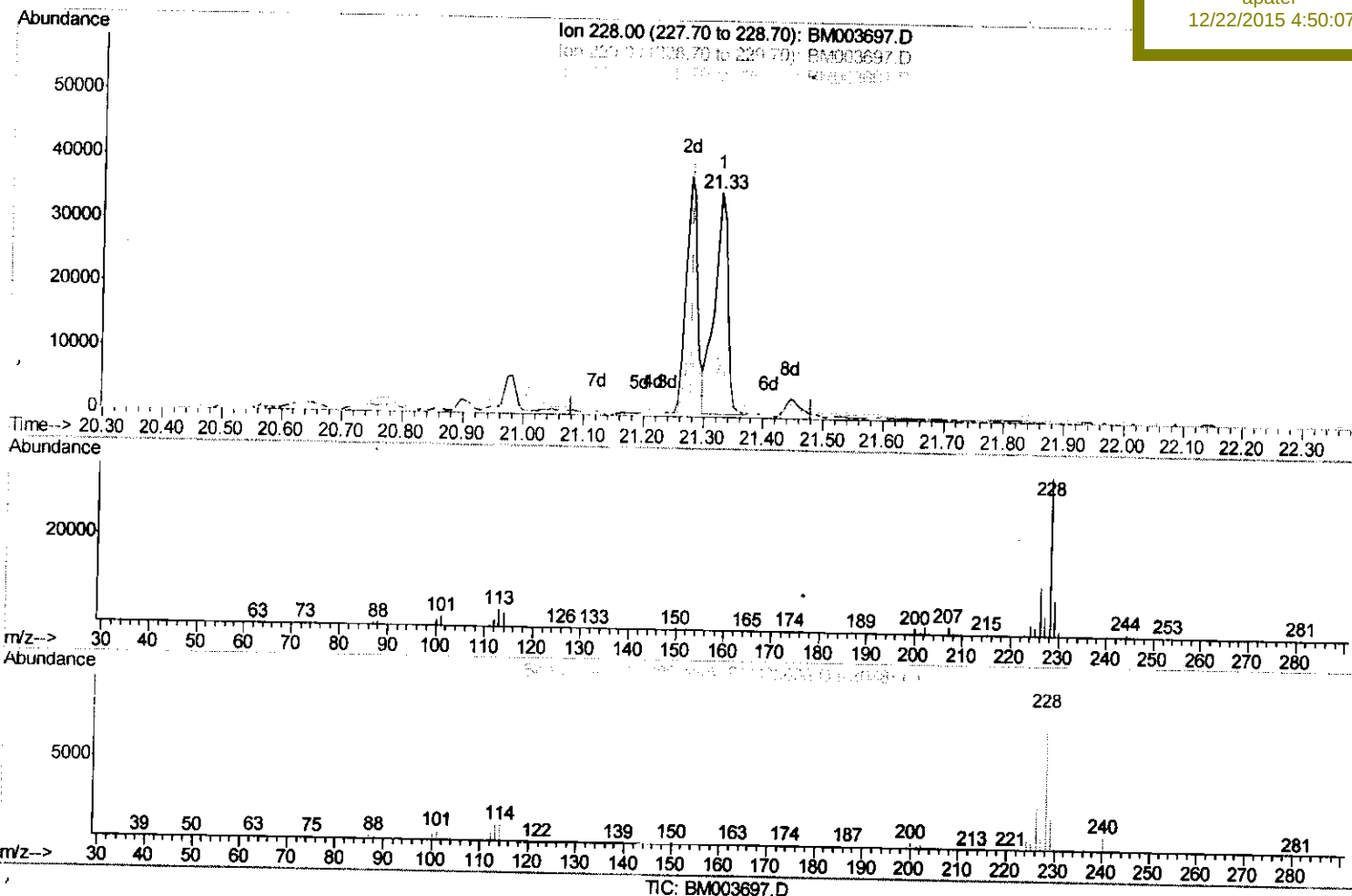
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(83) Benzo(a)anthracene

21.327min (+0.047) 5.81ng/ul

response 53541

Ion	Exp%	Act%
228.00	100	100
229.00	19.60	23.12
226.00	27.20	31.95
0.00	0.00	0.00

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Quant Time: Dec 24 01:56:37 2015
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	29247	20.00	ng/ul	0.00
18) Naphthalene-d8	10.48	136	121383	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.35	164	65297	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.09	188	139630	20.00	ng/ul	0.00
78) Chrysene-d12	21.29	240	152959	20.00	ng/ul	0.00
86) Perylene-d12	23.53	264	150680	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	6.87	99	504	0.20	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.03	67	249	0.19	ng/ul	0.00
9) 2-Chlorophenol-d4	7.23	132	415	0.21	ng/ul	0.00
13) 4-Methylphenol-d8	8.23	113	522	0.24	ng/ul	-0.18
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	10.60	131	706	0.29	ng/ul	-0.02
44) Dimethylphthalate-d6	13.75	166	2587	0.48	ng/ul	0.00
47) Acenaphthylene-d8	14.04	160	3034	0.49	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.34	176	2868	0.64	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.19	188	4055	0.64	ng/ul	0.00
79) Pyrene-d10	19.49	212	4537	0.68	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.39	264	4269	0.57	ng/ul	0.00

Target Compounds

					Qvalue
28) Naphthalene	10.53	128	211767	33.86	ng/ul 100
34) 2-Methylnaphthalene	12.15	142	167507	35.87	ng/ul 99
48) Acenaphthylene	14.06	152	36244	5.39	ng/ul# 91
50) Acenaphthene	14.41	153	54417	12.13	ng/ul 100
54) Dibenzofuran	14.75	168	10954	1.71	ng/ul# 76
59) Fluorene	15.40	166	49360	9.60	ng/ul 98
70) Phenanthrene	17.14	178	241376	31.25	ng/ul 100
72) Anthracene	17.23	178	61760	7.84	ng/ul 99
77) Fluoranthene	19.16	202	77557	8.74	ng/ul 99
80) Pyrene	19.52	202	172411	19.59	ng/ul 100
83) Benzo(a)anthracene	21.27	228	50969m	5.53	ng/ul
85) Chrysene	21.33	228	54754	6.16	ng/ul 95
88) Benzo(b)fluoranthene	22.86	252	37551m	4.13	ng/ul
91) Benzo(a)pyrene	23.43	252	41926	4.83	ng/ul 98
92) Indeno(1,2,3-cd)pyrene	25.79	276	14378	1.54	ng/ul 95
94) Benzo(g,h,i)perylene	26.49	276	15670	2.04	ng/ul 99

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

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