

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102419\

Data File : BN008400.D

Acq On : 24 Oct 2019 22:36

Operator : JU

Sample : K5236-01DL 5X

Misc :

ALS Vial : 23 Sample Multiplier: 1

Quant Time: Oct 25 00:53:27 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Oct 23 14:07:55 2019

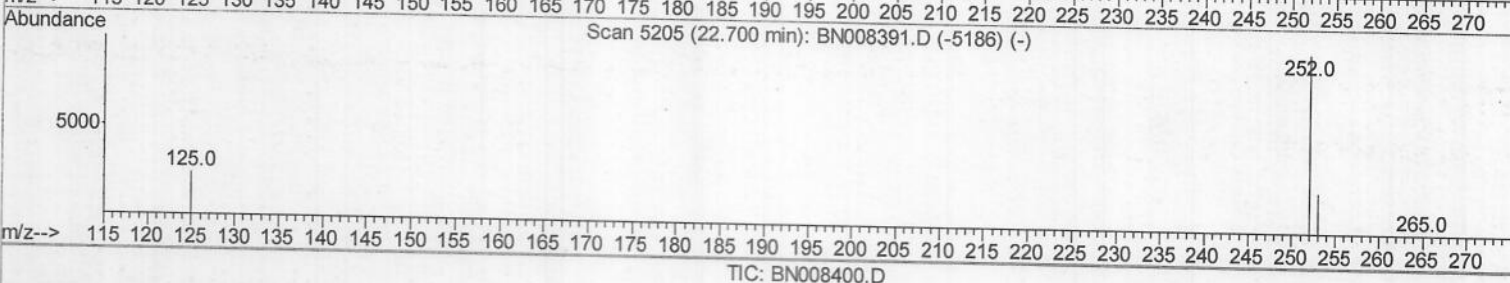
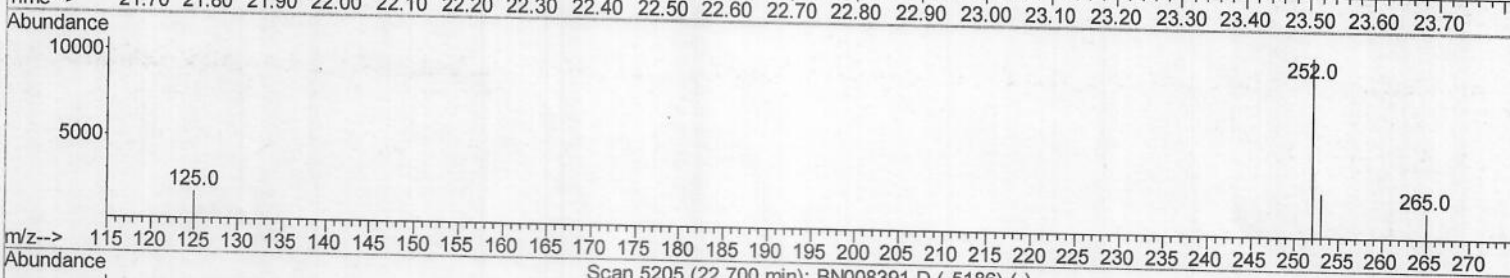
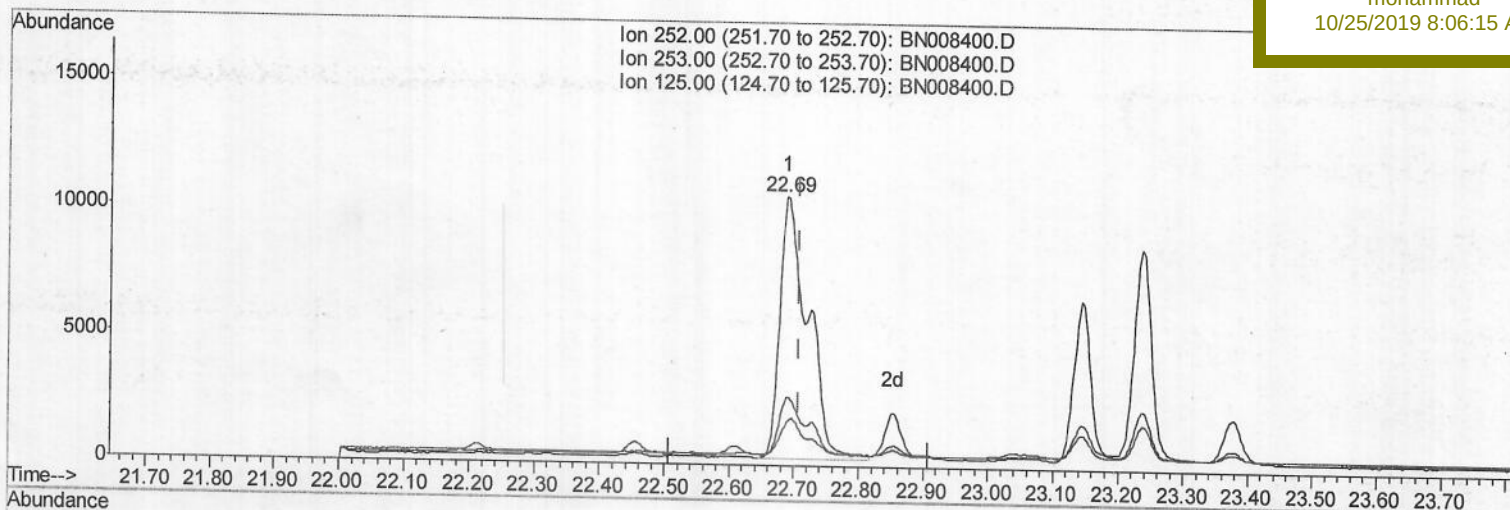
Response via : Initial Calibration

Instrument :

BNA_N

ClientSampleId :

BEPC2DL

Manual Integrations
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10/25/2019 8:06:15 AM

(21) Benzo(b)fluoranthene

22.688min (-0.020) 1.06ng/ul

response 31032

Ion	Exp%	Act%
252.00	100	100
253.00	26.70	25.39
125.00	16.20	16.05
0.00	0.00	0.00

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ClientSampleId :

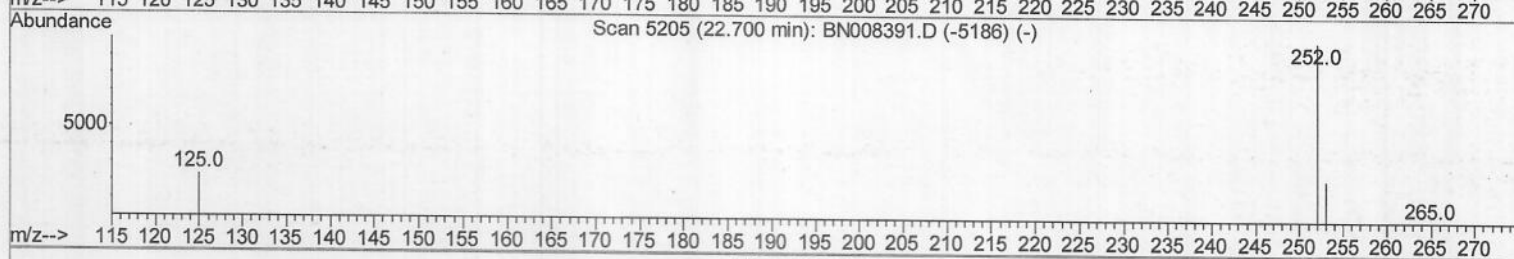
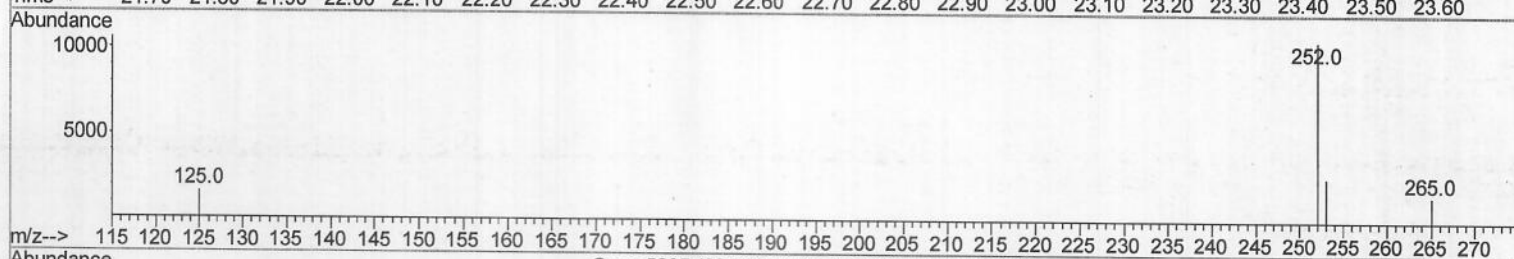
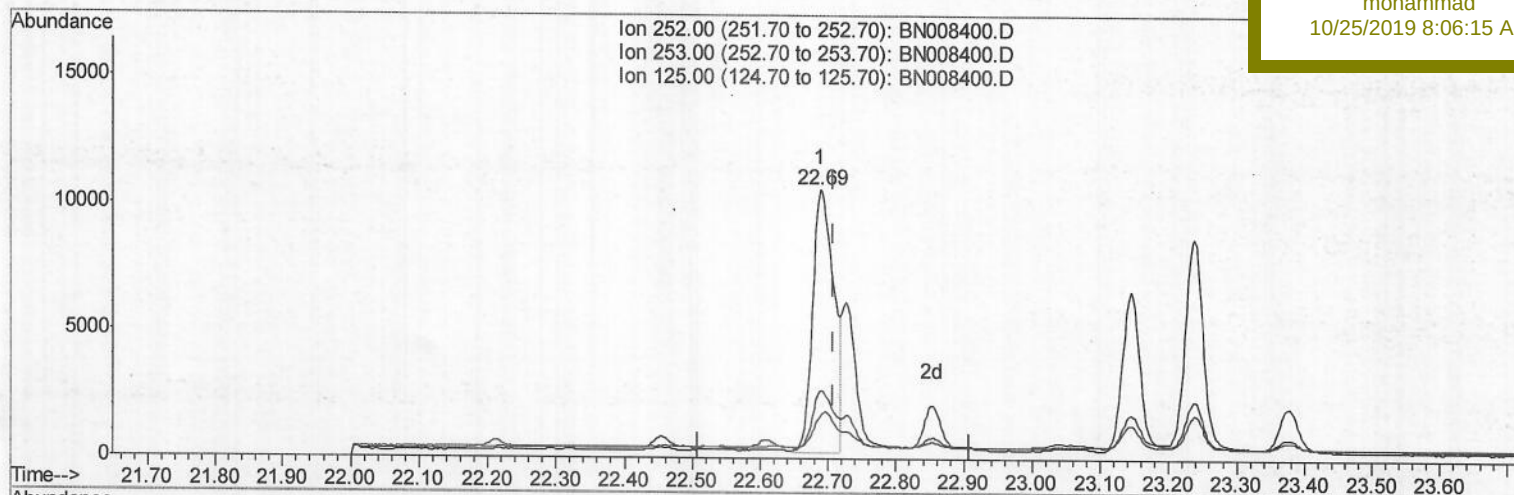
BEPC2DL

Manual Integrations

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(21) Benzo(b)fluoranthene

22.688min (-0.020) 0.77ng/ul m

response 22394

Ion	Exp%	Act%
252.00	100	100
253.00	26.70	25.39
125.00	16.20	16.05
0.00	0.00	0.00

J.U
10/25/19

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Operator : JU

Sample : K5236-01DL 5X

Misc :

ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :

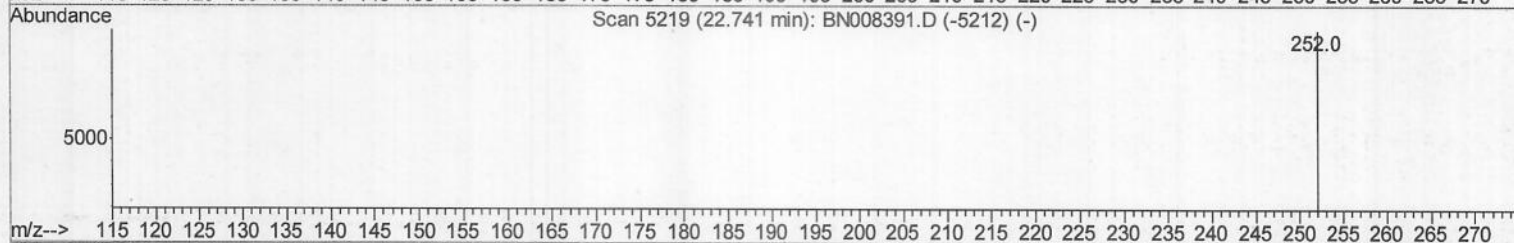
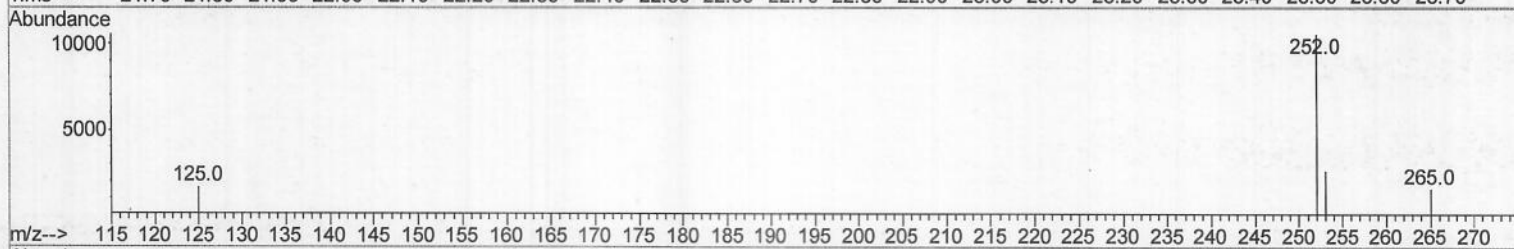
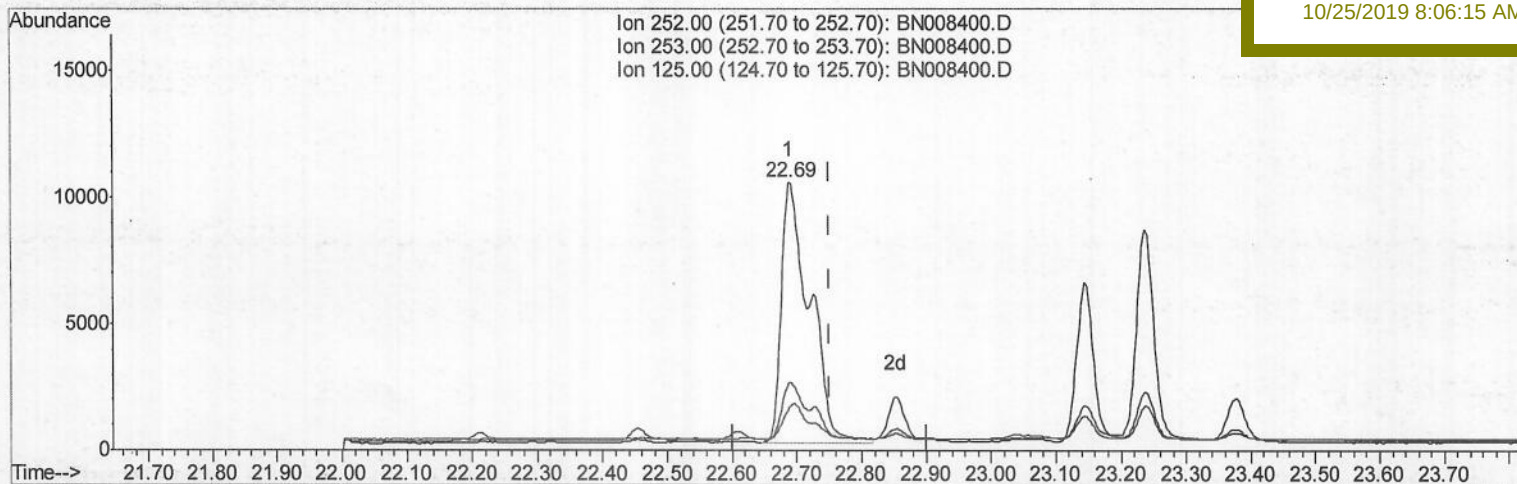
BEPC2DL

Manual Integrations

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TIC: BN008400.D

(22) Benzo(k)fluoranthene

22.688min (-0.064) 0.94ng/ul

response 31032

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	25.39
125.00	15.40	16.05
0.00	0.00	0.00

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Data File : BN008400.D

Acq On : 24 Oct 2019 22:36

Operator : JU

Sample : K5236-01DL 5X

Misc :

ALS Vial : 23 Sample Multiplier: 1

Quant Time: Oct 25 00:53:27 2019

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Instrument :

BNA_N

ClientSampled :

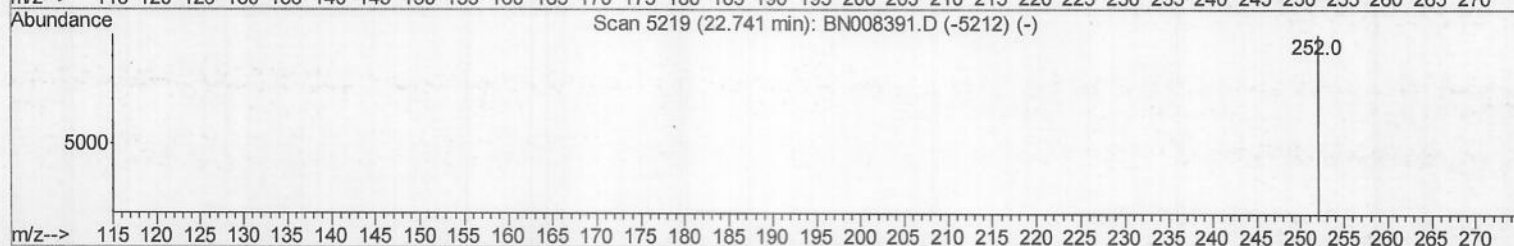
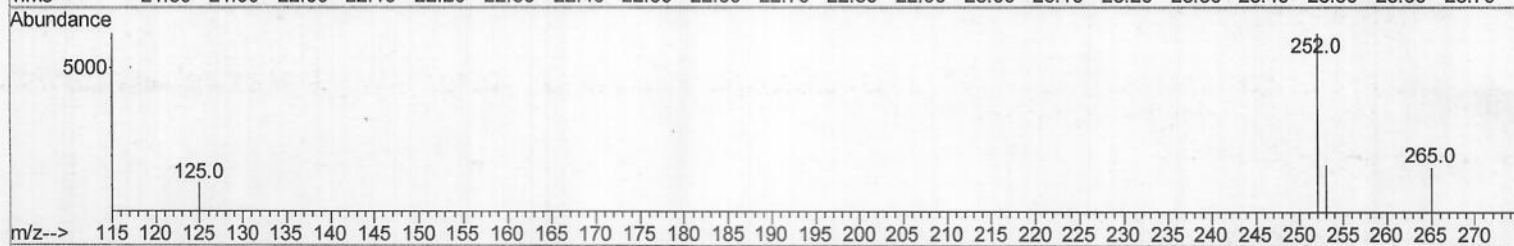
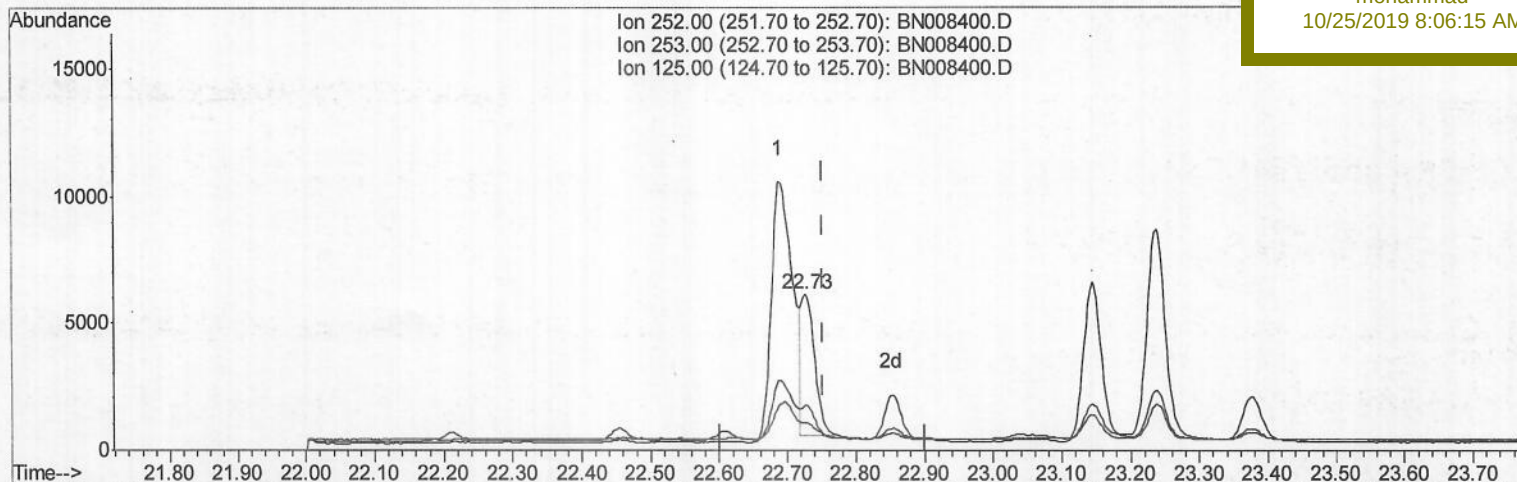
BEPC2DL

Manual Integrations

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TIC: BN008400.D

(22) Benzo(k)fluoranthene

22.726min (-0.026) 0.23ng/ul m →

response 7659

Ion	Exp%	Act%
252.00	100	100
253.00	26.60	27.91
125.00	15.40	17.40
0.00	0.00	0.00

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Instrument :
 BNA_N
 ClientSampleId :
 BEPC2DL

Quant Time: Oct 25 00:58:00 2019
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 14:07:55 2019
 Response via : Initial Calibration

Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1574	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.33	136	7731	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	5051	0.40	ng/ul	-0.01
10) Phenanthrene-d10	16.96	188	10975	0.40	ng/ul	-0.02
16) Chrysene-d12	21.16	240	10680	0.40	ng/ul	-0.02
20) Perylene-d12	23.33	264	8985	0.40	ng/ul	-0.03

System Monitoring Compounds

4) 2-Methylnaphthalene-d10	11.95	152	456	0.03	ng/ul	-0.01
14) Fluoranthene-d10	19.00	212	1516	0.04	ng/ul	-0.02

Target Compounds

						Ovalue
7) Acenaphthylene	13.93	152	1667	0.069	ng/ul#	90
8) Acenaphthene	14.27	153	405	0.022	ng/ul#	93
9) Fluorene	15.26	166	575	0.027	ng/ul#	95
12) Phenanthrene	17.00	178	11286	0.351	ng/ul	99
13) Anthracene	17.09	178	2823	0.100	ng/ul#	93
15) Fluoranthene	19.02	202	31943	0.753	ng/ul	96
17) Pyrene	19.39	202	34233	0.805	ng/ul	97
18) Benzo(a)anthracene	21.15	228	19689	0.507	ng/ul	96
19) Chrysene	21.20	228	19851	0.417	ng/ul	98
21) Benzo(b)fluoranthene	22.69	252	22394m	0.766	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	7659m	0.231	ng/ul	
23) Benzo(a)pyrene	23.24	252	15536	0.505	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.48	276	9692	0.267	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.48	278	2322	0.081	ng/ul#	75
26) Benzo(g,h,i)perylene	26.13	276	9576	0.288	ng/ul	96

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

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