

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
 Data File : BM012902.D
 Acq On : 11 Nov 2017 15:16
 Operator : SJ/JU
 Sample : I6319-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DAHL2

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:50:59 PM

Quant Time: Nov 13 01:37:23 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 01:07:29 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	70540	20.00	ng/ul	0.00
2) Naphthalene-d8	10.57	136	268625	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.42	164	158380	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.17	188	373566	20.00	ng/ul	0.00
17) Chrysene-d12	21.36	240	458269	20.00	ng/ul	0.00
22) Perylene-d12	23.63	264	509187	20.00	ng/ul	0.00
System Monitoring Compounds						
7) Acenaphthylene-d8	14.12	160	478006	29.02	ng/ul	0.00
10) Fluorene-d10	15.42	176	340568	30.23	ng/ul	0.00
14) Anthracene-d10	17.27	188	535815	29.67	ng/ul	0.00
18) Pyrene-d10	19.56	212	624835	30.47	ng/ul	0.00
25) Benzo(a)pyrene-d12	23.49	264	741203	30.83	ng/ul	0.00
Target Compounds						
23) Benzo(b)fluoranthene	22.94	252	42235m	1.395	ng/ul	Ovalue

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

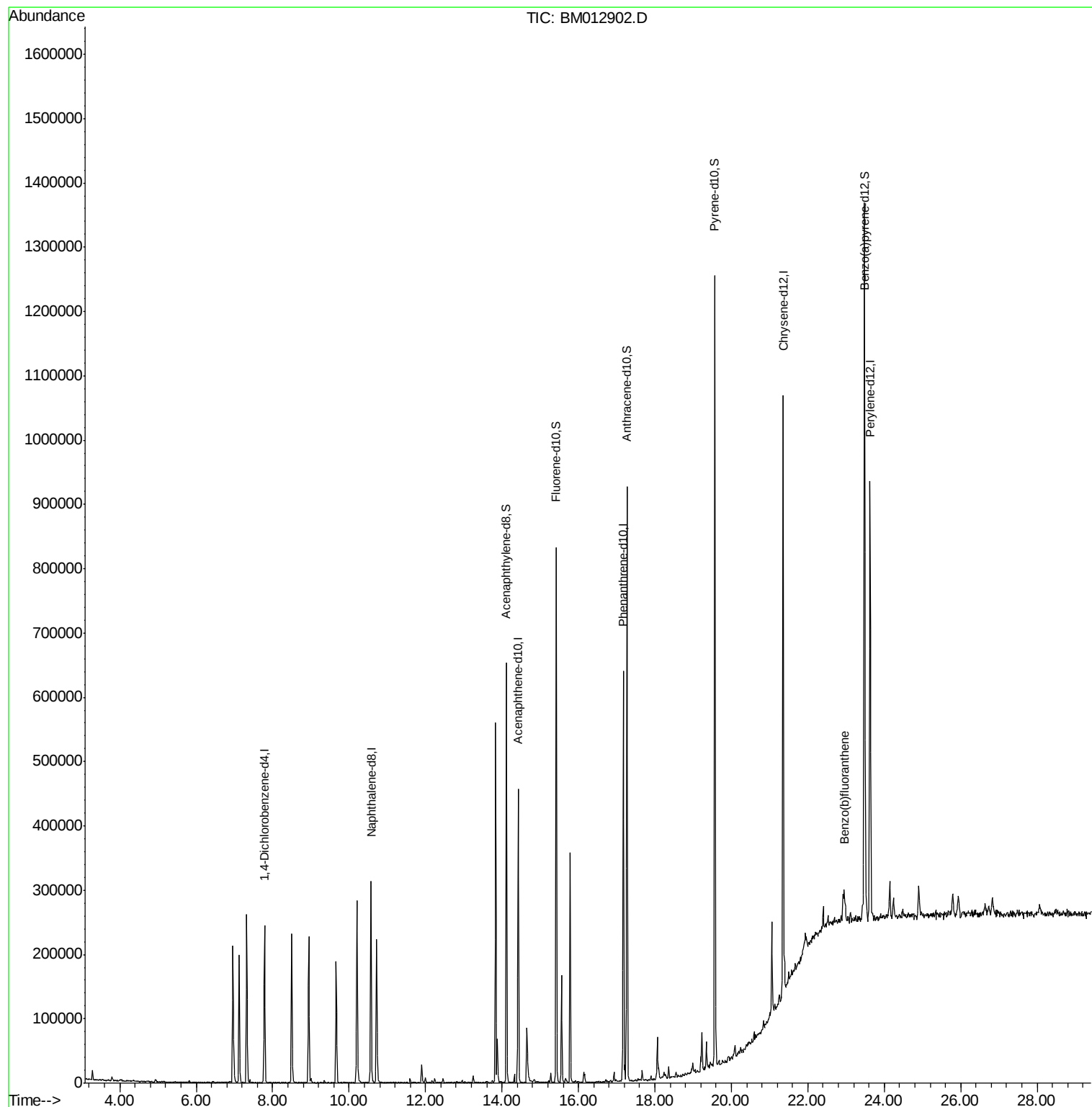
Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
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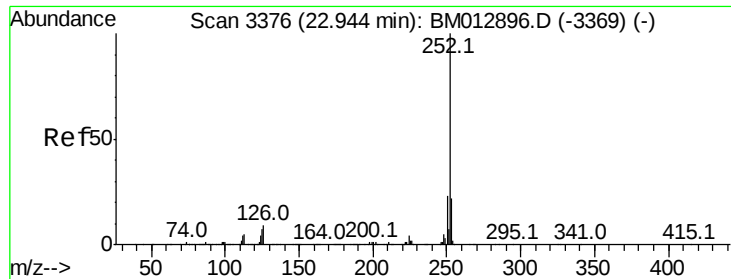
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#23

Benzo(b)fluoranthene

Concen: 1.395 ng/ul m

RT: 22.94 min Scan# 3376

Delta R.T. -0.00 min

Lab File: BM012902.D

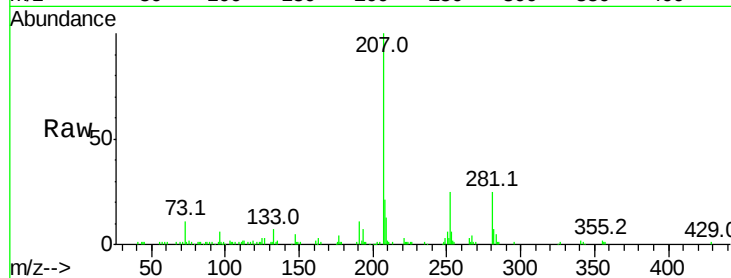
Acq: 11 Nov 2017 15:16

Instrument :

BNA_M

ClientSampleId :

DAHL2



Tot Ion:	252	Resp:	42235
Ion Ratio	Lower	Upper	
252	100		
253	23.2	18.0	27.0
125	10.4	8.0	12.0

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