

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
 Data File : BM012864.D
 Acq On : 10 Nov 2017 16:04
 Operator : SJ/JU
 Sample : I6317-22
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DAHH3

Quant Time: Nov 10 16:57:43 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 14:59:43 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	45112	20.00	ng/ul	0.00
2) Naphthalene-d8	10.58	136	176581	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.43	164	106135	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.18	188	259113	20.00	ng/ul	0.00
17) Chrysene-d12	21.36	240	343964	20.00	ng/ul	0.00
22) Perylene-d12	23.64	264	388193	20.00	ng/ul	0.00
System Monitoring Compounds						
7) Acenaphthylene-d8	14.12	160	238452	21.60	ng/ul	0.00
10) Fluorene-d10	15.43	176	176213	23.34	ng/ul	0.00
14) Anthracene-d10	17.28	188	341112	27.23	ng/ul	0.00
18) Pyrene-d10	19.57	212	429476	27.91	ng/ul	0.00
25) Benzo(a)pyrene-d12	23.50	264	555736	30.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.63	128	51074	5.712	ng/ul	99
5) 1-Methylnaphthalene	12.46	142	71549	11.089	ng/ul	92
8) Acenaphthylene	14.15	152	38163	3.553	ng/ul	98
11) Fluorene	15.48	166	162660	18.955	ng/ul	92
15) Anthracene	17.32	178	138589	9.320	ng/ul	99
20) Benzo(a)anthracene	21.34	228	363110	17.157	ng/ul	100
23) Benzo(b)fluoranthene	22.96	252	473885	20.528	ng/ul#	96
24) Benzo(k)fluoranthene	23.00	252	195153	8.845	ng/ul#	96

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

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