

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
 Data File : BM012882.D
 Acq On : 11 Nov 2017 02:46
 Operator : SJ/JU
 Sample : I6322-11
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DAHM3

Quant Time: Nov 11 04:03:16 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 22:45:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	72195	20.00	ng/ul	0.00
2) Naphthalene-d8	10.57	136	277373	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.43	164	170136	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.17	188	410067	20.00	ng/ul	0.00
17) Chrysene-d12	21.36	240	510635	20.00	ng/ul	0.00
22) Perylene-d12	23.63	264	569298	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
7) Acenaphthylene-d8	14.12	160	441362	24.94	ng/ul	0.00
10) Fluorene-d10	15.42	176	326633	26.99	ng/ul	0.00
14) Anthracene-d10	17.27	188	554758	27.99	ng/ul	0.00
18) Pyrene-d10	19.57	212	643238	28.15	ng/ul	0.00
25) Benzo(a)pyrene-d12	23.49	264	790622	29.42	ng/ul	0.00

Target Compounds	Ovalue

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

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