

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
 Data File : BM012915.D
 Acq On : 11 Nov 2017 22:58
 Operator : SJ/JU
 Sample : I6317-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DAHJ4

Quant Time: Nov 13 03:05:20 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 02:06:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	85953	20.00	ng/ul	0.00
2) Naphthalene-d8	10.57	136	322484	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.42	164	182917	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.17	188	424085	20.00	ng/ul	0.00
17) Chrysene-d12	21.35	240	524389	20.00	ng/ul	0.00
22) Perylene-d12	23.63	264	591622	20.00	ng/ul	0.00

System Monitoring Compounds						
7) Acenaphthylene-d8	14.12	160	303803	15.97	ng/ul	0.00
10) Fluorene-d10	15.42	176	226458	17.40	ng/ul	0.00
14) Anthracene-d10	17.27	188	342203	16.69	ng/ul	0.00
18) Pyrene-d10	19.56	212	401508	17.11	ng/ul	0.00
25) Benzo(a)pyrene-d12	23.48	264	477037	17.08	ng/ul	0.00

Target Compounds				Ovalue		
13) Phenanthrene	17.21	178	52970	2.261	ng/ul	99
16) Fluoranthene	19.23	202	116068	4.028	ng/ul#	92
19) Pyrene	19.59	202	92608	3.077	ng/ul	94
20) Benzo(a)anthracene	21.33	228	71331	2.211	ng/ul	97
21) Chrysene	21.39	228	62669	2.173	ng/ul	98
23) Benzo(b)fluoranthene	22.94	252	118875	3.379	ng/ul	96
26) Benzo(a)pyrene	23.53	252	67008	1.964	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.93	276	62059	1.410	ng/ul#	93
29) Benzo(g,h,i)perylene	26.63	276	53135	1.446	ng/ul	95

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

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