

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA M\Data\BM070816\
 Data File : BM006311.D
 Acq On : 07 Jul 2016 09:49
 Operator : UM/SJ
 Sample : PB91709BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SBLK09

Quant Time: Jul 07 13:46:34 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM070716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 07 05:51:09 2016
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.651	152	3908	0.40	ng/ul	0.00
4) Naphthalene-d8	10.428	136	15940	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.297	164	9566	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.142	188	1479145	0.40	ng/ul	0.10
18) Chrysene-d12	21.245	240	15575	0.40	ng/ul	0.00
22) Perylene-d12	23.319	264	929222	0.40	ng/ul	-0.13
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.166	96	9797	5.93	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.028	152	7440	0.40	ng/ul	0.01
16) Fluoranthene-d10	19.077	212	19258	0.01	ng/ul	0.00
Target Compounds						
					Ovalue	
11) Fluorene	15.291	166	2714	0.08	ng/ul#	21
19) Pyrene	19.452	202	2599	0.04	ng/ul#	1

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

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