

Data Path : Z:\HPCHEM1\BNA M\Data\BM122016\Not Reviewed Data\NR\
 Data File : BM008485.D
 Acq On : 20 Dec 2016 18:42
 Operator : UM/SJ
 Sample : H6039-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Dec 21 09:31:35 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 02:43:00 2016
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.666	152	398	0.40	ng/ul	-0.03
2) Naphthalene-d8	10.427	136	1278	0.40	ng/ul	-0.05
6) Acenaphthene-d10	14.297	164	732	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.142	188	97264	0.40	ng/ul	0.07
16) Chrysene-d12	21.273	240	1172	0.40	ng/ul	0.00
20) Perylene-d12	23.483	264	1478	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.059	152	460	0.23	ng/ul	-0.02
14) Fluoranthene-d10	19.096	212	992	0.00	ng/ul	0.00
Target Compounds						
					Ovalue	
3) Naphthalene	10.476	128	139	0.04	ng/ul#	15
9) Fluorene	15.290	166	207	0.07	ng/ul#	38
17) Pyrene	19.447	202	201	0.04	ng/ul#	1
21) Benzo(b)fluoranthene	22.815	252	137	0.02	ng/ul#	1
22) Benzo(k)fluoranthene	22.815	252	137	0.02	ng/ul#	1
23) Benzo(a)pyrene	23.326	252	414	0.08	ng/ul#	21

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

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