

Data Path : Z:\HPCHEM1\BNA M\Data\BM122016\Not Reviewed Data\NR\
 Data File : BM008494.D
 Acq On : 21 Dec 2016 00:05
 Operator : UM/SJ
 Sample : H6039-17
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Dec 21 09:33:26 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	7.728	152	16	0.40	ng/ul	0.03
2) Naphthalene-d8	10.476	136	27408	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.417	164	13504	0.40	ng/ul	0.10
10) Phenanthrene-d10	17.159	188	144978	0.40	ng/ul	0.09
16) Chrysene-d12	21.408	240	229	0.40	ng/ul	0.14
20) Perylene-d12	0.000	264	0	0.00	ng/ul	-23.50
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.070	152	36891	0.88	ng/ul	-0.01
14) Fluoranthene-d10	19.169	212	793	0.00	ng/ul	0.07
Target Compounds						
					Ovalue	
3) Naphthalene	10.500	128	50827	0.66	ng/ul#	94
5) 2-Methylnaphthalene	12.111	142	14503	0.28	ng/ul#	79
7) Acenaphthylene	14.091	152	45749	0.75	ng/ul	96
8) Acenaphthene	14.485	153	17958	0.39	ng/ul#	23
9) Fluorene	15.547	166	3242	0.06	ng/ul#	28
12) Phenanthrene	17.142	178	9222	0.02	ng/ul#	1
17) Pyrene	19.665	202	578	0.65	ng/ul#	1
18) Benzo(a)anthracene	21.377	228	473	0.63	ng/ul#	1
19) Chrysene	21.377	228	473	0.57	ng/ul#	1

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

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