

Data Path : Z:\HPCHEM1\BNA M\Data\BM122016\Not Reviewed Data\NR\
 Data File : BM008488.D
 Acq On : 20 Dec 2016 20:30
 Operator : UM/SJ
 Sample : H6039-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Dec 21 09:32:22 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	7.650	152	403	0.40	ng/ul	-0.05
2) Naphthalene-d8	10.258	136	15903	0.40	ng/ul	-0.22
6) Acenaphthene-d10	14.023	164	2972	0.40	ng/ul	-0.29
10) Phenanthrene-d10	16.696	188	120	0.40	ng/ul	-0.38
16) Chrysene-d12	20.762	240	1	0.40	ng/ul	-0.50
20) Perylene-d12	22.951	264	4	0.40	ng/ul	-0.55
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.882	152	503	0.02	ng/ul	-0.20
14) Fluoranthene-d10	0.000	212	0	0.00	ng/ul	
Target Compounds						
					Ovalue	
7) Acenaphthylene	13.725	152	315	0.02	ng/ul#	1
8) Acenaphthene	14.040	153	586	0.06	ng/ul#	1
9) Fluorene	15.136	166	253	0.02	ng/ul#	1
11) Pentachlorophenol	16.250	266	7	0.19	ng/ul#	38
12) Phenanthrene	16.713	178	293	0.80	ng/ul#	1
13) Anthracene	16.867	178	128	0.36	ng/ul#	1
17) Pyrene	19.120	202	1060	274.22	ng/ul#	1
18) Benzo(a)anthracene	20.731	228	10	3.05	ng/ul#	1
19) Chrysene	20.804	228	3	0.83	ng/ul#	1
21) Benzo(b)fluoranthene	22.346	252	9	0.57	ng/ul#	1
22) Benzo(k)fluoranthene	22.346	252	9	0.59	ng/ul#	1
23) Benzo(a)pyrene	22.815	252	111	7.60	ng/ul#	1
26) Benzo(g,h,i)perylene	25.849	276	4	0.27	ng/ul#	1

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

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