

Data Path : Z:\HPCHEM1\BNA M\Data\BM122016\Not Reviewed Data\NR\
 Data File : BM008495.D
 Acq On : 21 Dec 2016 00:41
 Operator : UM/SJ
 Sample : H6039-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Dec 21 09:33:37 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	7.651	152	518	0.40	ng/ul	-0.05
2) Naphthalene-d8	10.258	136	4857	0.40	ng/ul	-0.22
6) Acenaphthene-d10	14.057	164	2179	0.40	ng/ul	-0.26
10) Phenanthrene-d10	16.765	188	451	0.40	ng/ul	-0.31
16) Chrysene-d12	20.898	240	16	0.40	ng/ul	-0.37
20) Perylene-d12	23.066	264	6	0.40	ng/ul	-0.43
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.841	152	74	0.01	ng/ul	-0.24
14) Fluoranthene-d10	0.000	212	0	0.00	ng/ul	
Target Compounds						
					Ovalue	
3) Naphthalene	10.258	128	3913	0.29	ng/ul#	1
7) Acenaphthylene	13.839	152	884	0.09	ng/ul#	1
9) Fluorene	15.171	166	204	0.02	ng/ul#	1
11) Pentachlorophenol	16.422	266	4	0.03	ng/ul#	20
12) Phenanthrene	16.782	178	149	0.11	ng/ul#	1
13) Anthracene	16.867	178	114	0.09	ng/ul#	1
17) Pyrene	19.108	202	3236	52.32	ng/ul#	1
18) Benzo(a)anthracene	20.846	228	31	0.59	ng/ul#	1
19) Chrysene	20.919	228	18	0.31	ng/ul#	1
21) Benzo(b)fluoranthene	22.388	252	12	0.51	ng/ul#	1
22) Benzo(k)fluoranthene	22.388	252	14	0.61	ng/ul#	1
23) Benzo(a)pyrene	22.961	252	26	1.19	ng/ul#	1
26) Benzo(g,h,i)perylene	25.953	276	4	0.18	ng/ul#	1

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

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