

Data Path : Z:\HPCHEM1\BNA_M\Data\BM122416\
 Data File : BM008620.D
 Acq On : 24 Dec 2016 06:58
 Operator : UM/SJ
 Sample : H6069-08
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA8X8

Quant Time: Dec 25 01:06:49 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 24 04:56:03 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	502	0.40	ng/ul	-0.08
2) Naphthalene-d8	10.43	136	1651	0.40	ng/ul	-0.07
6) Acenaphthene-d10	14.28	164	966	0.40	ng/ul	-0.03
10) Phenanthrene-d10	17.14	188	95012	0.40	ng/ul	0.05
16) Chrysene-d12	21.25	240	1708	0.40	ng/ul	-0.02
20) Perylene-d12	23.33	264	83887	0.40	ng/ul	-0.18
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	561	0.22	ng/ul	-0.06
14) Fluoranthene-d10	19.08	212	991	0.00	ng/ul	-0.02
Target Compounds				Qvalue		
3) Naphthalene	10.46	128	5698	1.23	ng/ul	97
5) 2-Methylnaphthalene	12.11	142	1850	0.59	ng/ul	100
7) Acenaphthylene	14.01	152	639	0.15	ng/ul#	77
8) Acenaphthene	14.35	153	225	0.07	ng/ul#	86
9) Fluorene	15.34	166	175	0.05	ng/ul#	75
17) Pyrene	19.47	202	5686	0.86	ng/ul	98
18) Benzo(a)anthracene	21.29	228	4143	0.74	ng/ul	94
19) Chrysene	21.29	228	4143	0.67	ng/ul	95
21) Benzo(b)fluoranthene	22.80	252	8124	0.02	ng/ul	92
22) Benzo(k)fluoranthene	22.80	252	11294	0.04	ng/ul	93

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

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