

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122316\
 Data File : BM008604.D
 Acq On : 23 Dec 2016 20:20
 Operator : UM/SJ
 Sample : H6069-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA7J1

Manual Integrations
 APPROVED

Sohil
 12/27/2016 5:20:42 PM

Quant Time: Dec 24 00:15:21 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 23 23:37:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	446	0.40	ng/ul	-0.09
2) Naphthalene-d8	10.42	136	1503	0.40	ng/ul	-0.08
6) Acenaphthene-d10	14.28	164	996	0.40	ng/ul	-0.03
10) Phenanthrene-d10	17.04	188	1910m	0.40	ng/ul	-0.05
16) Chrysene-d12	21.26	240	1562	0.40	ng/ul	-0.02
20) Perylene-d12	23.48	264	1342m	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	987	0.43	ng/ul	-0.07
14) Fluoranthene-d10	19.09	212	2089m	0.40	ng/ul	-0.02
Target Compounds						Ovalue
3) Naphthalene	10.46	128	2101	0.50	ng/ul	98
5) 2-Methylnaphthalene	12.10	142	2847	1.00	ng/ul	98
12) Phenanthrene	17.07	178	3211	0.55	ng/ul	100
15) Fluoranthene	19.11	202	666	0.09	ng/ul	71
17) Pyrene	19.48	202	1213	0.20	ng/ul#	83
18) Benzo(a)anthracene	21.23	228	478	0.09	ng/ul#	52
19) Chrysene	21.29	228	629	0.11	ng/ul#	66
21) Benzo(b)fluoranthene	22.81	252	631m	0.12	ng/ul	
22) Benzo(k)fluoranthene	22.84	252	188m	0.04	ng/ul	
23) Benzo(a)pyrene	23.37	252	453m	0.09	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.74	276	262m	0.04	ng/ul	
26) Benzo(g,h,i)perylene	26.43	276	330m	0.07	ng/ul	

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

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