

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122316\
 Data File : BM008607.D
 Acq On : 23 Dec 2016 22:06
 Operator : UM/SJ
 Sample : H6069-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA8X9

Manual Integrations
 APPROVED

sohil
 12/26/2016 4:03:33 PM

Quant Time: Dec 24 00:26:24 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	7.67	152	430	0.40	ng/ul	-0.08
2) Naphthalene-d8	10.43	136	1407	0.40	ng/ul	-0.07
6) Acenaphthene-d10	14.28	164	844	0.40	ng/ul	-0.03
10) Phenanthrene-d10	17.04	188	1725m	0.40	ng/ul	-0.05
16) Chrysene-d12	21.26	240	1567	0.40	ng/ul	-0.02
20) Perylene-d12	23.48	264	1406m	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	395	0.18	ng/ul	-0.05
14) Fluoranthene-d10	19.09	212	988m	0.21	ng/ul	-0.02
Target Compounds						Ovalue
3) Naphthalene	10.48	128	4667	1.18	ng/ul	97
5) 2-Methylnaphthalene	12.11	142	1499	0.56	ng/ul	99
7) Acenaphthylene	14.01	152	482	0.13	ng/ul#	73
8) Acenaphthene	14.35	153	184	0.06	ng/ul#	86
9) Fluorene	15.34	166	152	0.05	ng/ul#	71
12) Phenanthrene	17.07	178	2789	0.53	ng/ul	97
13) Anthracene	17.18	178	776	0.15	ng/ul#	83
15) Fluoranthene	19.11	202	4993	0.77	ng/ul	96
17) Pyrene	19.48	202	4870	0.80	ng/ul	98
18) Benzo(a)anthracene	21.23	228	3831	0.74	ng/ul	95
19) Chrysene	21.29	228	3907m	0.69	ng/ul	
21) Benzo(b)fluoranthene	22.81	252	7751m	1.40	ng/ul	
22) Benzo(k)fluoranthene	22.85	252	2314m	0.43	ng/ul	
23) Benzo(a)pyrene	23.37	252	5073	0.99	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.72	276	4842	0.79	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.71	278	1009	0.21	ng/ul#	63
26) Benzo(a,h,i)perylene	26.40	276	4438	0.85	ng/ul#	94

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

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