

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122416\
 Data File : BM008646.D
 Acq On : 24 Dec 2016 23:33
 Operator : UM/SJ
 Sample : H6069-06DL 2X
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA7J2DL

Manual Integrations
 APPROVED

sohil
 12/26/2016 6:28:03 PM

Quant Time: Dec 26 01:34:19 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 26 00:55:42 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	443	0.40	ng/ul	-0.08
2) Naphthalene-d8	10.43	136	1425	0.40	ng/ul	-0.05
6) Acenaphthene-d10	14.28	164	906	0.40	ng/ul	-0.03
10) Phenanthrene-d10	17.04	188	1869m	0.40	ng/ul	-0.03
16) Chrysene-d12	21.26	240	1620	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	1165m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	515	0.24	ng/ul	-0.04
14) Fluoranthene-d10	19.08	212	1435m	0.28	ng/ul	-0.01
Target Compounds						
3) Naphthalene	10.46	128	539	0.13	ng/ul#	66
5) 2-Methylnaphthalene	12.11	142	548	0.20	ng/ul	99
12) Phenanthrene	17.07	178	1264m	0.22	ng/ul	
15) Fluoranthene	19.12	202	1152	0.16	ng/ul	88
17) Pyrene	19.47	202	1210	0.19	ng/ul#	78
18) Benzo(a)anthracene	21.25	228	526	0.10	ng/ul#	62
19) Chrysene	21.30	228	602	0.10	ng/ul#	65
21) Benzo(b)fluoranthene	22.82	252	781m	0.17	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	317m	0.07	ng/ul	
23) Benzo(a)pyrene	23.38	252	397m	0.09	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.77	276	332m	0.07	ng/ul	
26) Benzo(g,h,i)perylene	26.44	276	310m	0.07	ng/ul	

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

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