

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122316\
 Data File : BM008597.D
 Acq On : 23 Dec 2016 16:08
 Operator : UM/SJ
 Sample : H5970-07DL 5X
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA846DL

Manual Integrations
 APPROVED

sohil
 12/26/2016 4:03:08 PM

Quant Time: Dec 26 10:57:19 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 03:21:10 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	309	0.40	ng/ul	-0.09
2) Naphthalene-d8	10.49	136	952m	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.31	164	552	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.07	188	993m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.28	240	1107	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	894m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	42	0.03	ng/ul	-0.01
14) Fluoranthene-d10	19.10	212	165m	0.06	ng/ul	-0.01
Target Compounds						
5) 2-Methylnaphthalene	12.19	142	55	0.03	ng/ul	92
17) Pyrene	19.49	202	414	0.10	ng/ul#	65
18) Benzo(a)anthracene	21.26	228	416	0.11	ng/ul#	65
19) Chrysene	21.31	228	308	0.08	ng/ul#	60
21) Benzo(b)fluoranthene	22.83	252	766m	0.22	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	247m	0.07	ng/ul	
23) Benzo(a)pyrene	23.39	252	431	0.13	ng/ul#	47
24) Indeno(1,2,3-cd)pyrene	25.77	276	353	0.09	ng/ul#	89
26) Benzo(g,h,i)perylene	26.46	276	297	0.09	ng/ul#	39

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

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