

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122916\
 Data File : BM008714.D
 Acq On : 29 Dec 2016 11:13
 Operator : UM/SJ
 Sample : H6067-09DL 5X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA7G0DL

Manual Integrations
 APPROVED

Sohil
 12/30/2016 8:42:34 AM

Quant Time: Dec 30 01:37:48 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 30 00:54:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	327m	0.40	ng/ul	-0.15
2) Naphthalene-d8	10.49	136	993m	0.40	ng/ul	-0.06
6) Acenaphthene-d10	14.30	164	615m	0.40	ng/ul	-0.03
10) Phenanthrene-d10	17.06	188	1223m	0.40	ng/ul	-0.05
16) Chrysene-d12	21.26	240	1242m	0.40	ng/ul	-0.02
20) Perylene-d12	23.48	264	1197m	0.40	ng/ul	-0.03

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	72m	0.05	ng/ul	-0.03
14) Fluoranthene-d10	19.10	212	217m	0.06	ng/ul	-0.02

Target Compounds						Ovalue
7) Acenaphthylene	14.02	152	2561m	0.92	ng/ul	
12) Phenanthrene	17.11	178	509m	0.14	ng/ul	
13) Anthracene	17.19	178	2995m	0.83	ng/ul	
15) Fluoranthene	19.12	202	2098m	0.46	ng/ul	
17) Pyrene	19.48	202	3850m	0.80	ng/ul	
18) Benzo(a)anthracene	21.24	228	4901m	1.20	ng/ul	
19) Chrysene	21.29	228	11951m	2.68	ng/ul	
21) Benzo(b)fluoranthene	22.82	252	34185m	7.25	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	7203m	1.57	ng/ul	
23) Benzo(a)pyrene	23.39	252	22257m	5.09	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.72	276	22330m	4.29	ng/ul	
25) Dibenzo(a,h)anthracene	25.71	278	5939m	1.43	ng/ul	
26) Benzo(g,h,i)perylene	26.41	276	22952m	5.14	ng/ul	

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

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