

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122416\
 Data File : BM008666.D
 Acq On : 25 Dec 2016 11:57
 Operator : UM/SJ
 Sample : H6069-08DL 10X
 Misc :
 ALS Vial : 80 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA8X8DL

Manual Integrations
 APPROVED

sohil
 12/26/2016 6:29:41 PM

Quant Time: Dec 26 04:35:06 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	7.70	152	435	0.40	ng/ul	-0.08
2) Naphthalene-d8	10.49	136	1345m	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.31	164	772	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.07	188	1417m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.28	240	1544	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	1763	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	46	0.02	ng/ul	0.01
14) Fluoranthene-d10	19.11	212	178m	0.05	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.55	128	507	0.13	ng/ul#	35
5) 2-Methylnaphthalene	12.21	142	154	0.06	ng/ul	98
12) Phenanthrene	17.12	178	306m	0.07	ng/ul	
15) Fluoranthene	19.14	202	598m	0.11	ng/ul	
17) Pyrene	19.50	202	584	0.10	ng/ul#	69
18) Benzo(a)anthracene	21.26	228	425	0.08	ng/ul#	53
19) Chrysene	21.31	228	405	0.07	ng/ul#	57
21) Benzo(b)fluoranthene	22.84	252	933m	0.13	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	331m	0.05	ng/ul	
23) Benzo(a)pyrene	23.40	252	577	0.09	ng/ul#	42
24) Indeno(1,2,3-cd)pyrene	25.78	276	586	0.08	ng/ul#	87
26) Benzo(g,h,i)perylene	26.46	276	568	0.09	ng/ul#	56

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

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