

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM122416\
 Data File : BM008670.D
 Acq On : 25 Dec 2016 14:19
 Operator : UM/SJ
 Sample : H5995-09DL 5X
 Misc :
 ALS Vial : 84 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA874DL

Manual Integrations
 APPROVED

sohil
 12/26/2016 6:29:55 PM

Quant Time: Dec 30 10:14:40 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	7.65	152	488	0.40	ng/ul	-0.12
2) Naphthalene-d8	10.40	136	1877m	0.40	ng/ul	-0.11
6) Acenaphthene-d10	14.28	164	2234m	0.40	ng/ul	-0.05
10) Phenanthrene-d10	16.99	188	13705m	0.40	ng/ul	-0.10
16) Chrysene-d12	21.25	240	2099m	0.40	ng/ul	-0.03
20) Perylene-d12	23.48	264	1704m	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	266m	0.09	ng/ul	-0.17
14) Fluoranthene-d10	19.07	212	1455m	0.04	ng/ul	-0.05
Target Compounds				Qvalue		
3) Naphthalene	10.45	128	23044	4.36	ng/ul	98
5) 2-Methylnaphthalene	12.07	142	98138	27.61	ng/ul	98
7) Acenaphthylene	14.00	152	6160m	0.61	ng/ul	
8) Acenaphthene	14.35	153	7774m	1.02	ng/ul	
9) Fluorene	15.34	166	7864m	0.91	ng/ul	
12) Phenanthrene	17.07	178	126813m	3.03	ng/ul	
15) Fluoranthene	19.10	202	78857m	1.54	ng/ul	
17) Pyrene	19.48	202	71653m	8.83	ng/ul	
18) Benzo(a)anthracene	21.22	228	16587m	2.41	ng/ul	
19) Chrysene	21.28	228	19261m	2.55	ng/ul	
21) Benzo(b)fluoranthene	22.81	252	17074	2.54	ng/ul	96
22) Benzo(k)fluoranthene	22.83	252	5191m	0.79	ng/ul	
23) Benzo(a)pyrene	23.38	252	6782	1.09	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.72	276	4149m	0.56	ng/ul	
25) Dibenzo(a,h)anthracene	25.71	278	1492m	0.25	ng/ul	
26) Benzo(g,h,i)perylene	26.42	276	3730m	0.59	ng/ul	

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

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