

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM122816\
 Data File : BM008691.D
 Acq On : 28 Dec 2016 13:35
 Operator : UM/SJ
 Sample : H5995-03DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA864DL

Manual Integrations
 APPROVED

mohammad
 12/30/2016 8:53:04 AM

Quant Time: Dec 28 16:19:38 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 28 15:23:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	336m	0.40	ng/ul	-0.15
2) Naphthalene-d8	10.25	136	24726m	0.40	ng/ul	-0.29
6) Acenaphthene-d10	14.30	164	3883m	0.40	ng/ul	-0.03
10) Phenanthrene-d10	16.99	188	22556m	0.40	ng/ul	-0.10
16) Chrysene-d12	21.25	240	2594m	0.40	ng/ul	-0.03
20) Perylene-d12	23.49	264	1809	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	284m	0.01	ng/ul	-0.10
14) Fluoranthene-d10	19.07	212	4252m	0.07	ng/ul	-0.05
Target Compounds				Qvalue		
3) Naphthalene	10.45	128	29245m	0.42	ng/ul	
5) 2-Methylnaphthalene	12.07	142	53349m	1.14	ng/ul	
7) Acenaphthylene	14.00	152	3012m	0.17	ng/ul	
8) Acenaphthene	14.33	153	7026	0.53	ng/ul	93
9) Fluorene	15.34	166	28196m	1.87	ng/ul	
12) Phenanthrene	17.07	178	43419m	0.63	ng/ul	
15) Fluoranthene	19.11	202	10048	0.12	ng/ul	79
17) Pyrene	19.47	202	14911m	1.49	ng/ul	
23) Benzo(a)pyrene	23.39	252	1836	0.28	ng/ul#	31
24) Indeno(1,2,3-cd)pyrene	25.73	276	752	0.10	ng/ul#	78
26) Benzo(g,h,i)perylene	26.43	276	802	0.12	ng/ul#	18

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

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