

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122416\
 Data File : BM008656.D
 Acq On : 25 Dec 2016 05:27
 Operator : UM/SJ
 Sample : H5995-12DL 10X
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA880DL

Manual Integrations
 APPROVED

sohil
 12/26/2016 6:28:21 PM

Quant Time: Dec 26 02:56:34 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 00:55:42 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	388	0.40	ng/ul	-0.03
2) Naphthalene-d8	10.48	136	1168	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	716	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.06	188	1333m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.26	240	1423	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	1402m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	34	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	137m	0.04	ng/ul	0.00
Target Compounds						
					Ovalue	
3) Naphthalene	10.52	128	326	0.10	ng/ul#	29
5) 2-Methylnaphthalene	12.15	142	507	0.23	ng/ul#	80
7) Acenaphthylene	14.02	152	424	0.13	ng/ul#	62
8) Acenaphthene	14.37	153	208	0.09	ng/ul#	85
9) Fluorene	15.38	166	253	0.09	ng/ul#	81
12) Phenanthrene	17.11	178	1760m	0.43	ng/ul	
13) Anthracene	17.19	178	788m	0.20	ng/ul	
15) Fluoranthene	19.12	202	4572	0.92	ng/ul	97
17) Pyrene	19.48	202	5116	0.93	ng/ul	98
18) Benzo(a)anthracene	21.25	228	3275	0.70	ng/ul#	93
19) Chrysene	21.29	228	2895	0.57	ng/ul#	91
21) Benzo(b)fluoranthene	22.83	252	6012m	1.09	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	2499m	0.46	ng/ul	
23) Benzo(a)pyrene	23.40	252	3550	0.69	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.74	276	2906	0.48	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.74	278	611	0.13	ng/ul#	52
26) Benzo(a,h,i)perylene	26.44	276	2368	0.45	ng/ul	92

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

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