

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122316\
 Data File : BM008596.D
 Acq On : 23 Dec 2016 15:32
 Operator : UM/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.462

Manual Integrations
 APPROVED

sohil
 12/26/2016 4:04:32 PM

12/26/2016 4:04:32 PM

Quant Time: Dec 23 16:36:50 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 23 15:11:17 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	342	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	1295	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.31	164	752	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	1368	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	1404	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	1389	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	771	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.11	212	1533	0.41	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.55	128	1434	0.39	ng/ul#	78
5) 2-Methylnaphthalene	12.18	142	932	0.38	ng/ul	98
7) Acenaphthylene	14.04	152	1404	0.41	ng/ul#	90
8) Acenaphthene	14.38	153	1031	0.40	ng/ul	97
9) Fluorene	15.41	166	1101	0.38	ng/ul#	94
11) Pentachlorophenol	16.78	266	163	0.38	ng/ul	99
12) Phenanthrene	17.12	178	1637	0.39	ng/ul#	94
13) Anthracene	17.23	178	1724m	0.43	ng/ul	
15) Fluoranthene	19.14	202	2126	0.42	ng/ul	95
17) Pyrene	19.51	202	2257	0.42	ng/ul	95
18) Benzo(a)anthracene	21.26	228	1807	0.39	ng/ul	94
19) Chrysene	21.30	228	2044m	0.40	ng/ul	
21) Benzo(b)fluoranthene	22.83	252	2021	0.37	ng/ul	96
22) Benzo(k)fluoranthene	22.88	252	2036	0.38	ng/ul	93
23) Benzo(a)pyrene	23.41	252	2009	0.40	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.76	276	2362	0.39	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.76	278	1876	0.39	ng/ul#	88
26) Benzo(g,h,i)perylene	26.43	276	2058	0.40	ng/ul#	94

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

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