

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112816\
 Data File : BM008101.D
 Acq On : 29 Nov 2016 10:27
 Operator : UM/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.460

Quant Time: Nov 29 10:59:37 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 29 00:11:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	496	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	1791	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	1071	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.11	188	2021	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	2028	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	2053	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	1016	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	2233	0.39	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.55	128	1959	0.39	ng/ul	95
5) 2-Methylnaphthalene	12.16	142	1297	0.40	ng/ul	98
7) Acenaphthylene	14.06	152	2051	0.42	ng/ul	97
8) Acenaphthene	14.40	153	1409	0.38	ng/ul	92
9) Fluorene	15.41	166	1597	0.38	ng/ul	98
11) Pentachlorophenol	16.78	266	279	0.40	ng/ul	95
12) Phenanthrene	17.14	178	2416	0.38	ng/ul	95
13) Anthracene	17.23	178	2290	0.39	ng/ul	97
15) Fluoranthene	19.16	202	3052	0.38	ng/ul	98
17) Pyrene	19.53	202	3168	0.45	ng/ul	96
18) Benzo(a)anthracene	21.27	228	2804	0.45	ng/ul	96
19) Chrysene	21.32	228	2754	0.36	ng/ul	94
21) Benzo(b)fluoranthene	22.86	252	2955	0.36	ng/ul	94
22) Benzo(k)fluoranthene	22.90	252	2887	0.34	ng/ul	94
23) Benzo(a)pyrene	23.43	252	2881	0.38	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.79	276	3450	0.36	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.80	278	2739	0.36	ng/ul	94
26) Benzo(g,h,i)perylene	26.47	276	2948	0.35	ng/ul	93

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

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