

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112316\
 Data File : BM008080.D
 Acq On : 24 Nov 2016 18:38
 Operator : UM/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.455

Quant Time: Nov 25 01:30:42 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 25 01:27:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	583	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	2070	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	1239	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.11	188	2408	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	2301	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	2243	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	1201	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	2601	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.55	128	2289	0.40	ng/ul	97
5) 2-Methylnaphthalene	12.17	142	1528	0.41	ng/ul	97
7) Acenaphthylene	14.06	152	2340	0.42	ng/ul	96
8) Acenaphthene	14.40	153	1671	0.39	ng/ul	92
9) Fluorene	15.41	166	1923	0.39	ng/ul	93
11) Pentachlorophenol	16.78	266	326	0.39	ng/ul	97
12) Phenanthrene	17.14	178	2966	0.40	ng/ul	96
13) Anthracene	17.23	178	2856	0.41	ng/ul	98
15) Fluoranthene	19.17	202	3615	0.37	ng/ul	94
17) Pyrene	19.53	202	3823	0.48	ng/ul	97
18) Benzo(a)anthracene	21.27	228	3344	0.47	ng/ul	97
19) Chrysene	21.32	228	3269	0.37	ng/ul	95
21) Benzo(b)fluoranthene	22.86	252	3419	0.38	ng/ul	95
22) Benzo(k)fluoranthene	22.90	252	3443	0.37	ng/ul	97
23) Benzo(a)pyrene	23.43	252	3304	0.39	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.79	276	3936	0.38	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.80	278	3085	0.37	ng/ul	96
26) Benzo(g,h,i)perylene	26.47	276	3381	0.37	ng/ul#	94

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

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