

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112816\
 Data File : BM008097.D
 Acq On : 28 Nov 2016 21:30
 Operator : UM/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.458

Quant Time: Nov 29 00:14:23 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	549	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	2058	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	1224	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.11	188	2238	0.40	ng/ul	0.00
16) Chrysene-d12	21.30	240	2293	0.40	ng/ul	0.00
20) Perylene-d12	23.54	264	2125	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	1183	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.14	212	2471	0.39	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.55	128	2250	0.39	ng/ul	98
5) 2-Methylnaphthalene	12.16	142	1493	0.40	ng/ul	99
7) Acenaphthylene	14.06	152	2324	0.42	ng/ul	96
8) Acenaphthene	14.40	153	1641	0.39	ng/ul	92
9) Fluorene	15.41	166	1817	0.38	ng/ul	93
11) Pentachlorophenol	16.78	266	317	0.41	ng/ul	97
12) Phenanthrene	17.14	178	2734	0.39	ng/ul	97
13) Anthracene	17.23	178	2675	0.41	ng/ul	98
15) Fluoranthene	19.16	202	4536	0.50	ng/ul	99
17) Pyrene	19.52	202	3512	0.44	ng/ul	98
18) Benzo(a)anthracene	21.28	228	3096	0.44	ng/ul	93
19) Chrysene	21.33	228	2888	0.33	ng/ul	91
21) Benzo(b)fluoranthene	22.86	252	3107	0.37	ng/ul	91
22) Benzo(k)fluoranthene	22.90	252	3050	0.35	ng/ul#	90
23) Benzo(a)pyrene	23.43	252	2994	0.38	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.79	276	3571	0.36	ng/ul	95
25) Dibenzo(a,h)anthracene	25.80	278	2865	0.36	ng/ul#	91
26) Benzo(g,h,i)perylene	26.48	276	3064	0.35	ng/ul#	92

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

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