

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM111616\
 Data File : BM007925.D
 Acq On : 16 Nov 2016 17:28
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.465

Quant Time: Nov 16 17:59:43 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM102516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 16 16:16:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	670	0.40	ng/ul	0.02
2) Naphthalene-d8	10.54	136	2206	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.37	164	1185	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	2381	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	2312	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	2214	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	1114	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.16	212	2523	0.47	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.58	128	2376	0.42	ng/ul	95
5) 2-Methylnaphthalene	12.21	142	1514	0.42	ng/ul	99
7) Acenaphthylene	14.09	152	2157	0.39	ng/ul	98
8) Acenaphthene	14.43	153	1615	0.38	ng/ul	99
9) Fluorene	15.44	166	1791	0.38	ng/ul	98
11) Pentachlorophenol	16.82	266	269	0.48	ng/ul	91
12) Phenanthrene	17.16	178	2714	0.41	ng/ul	98
13) Anthracene	17.26	178	2581	0.43	ng/ul	98
15) Fluoranthene	19.19	202	3531	0.44	ng/ul	98
17) Pyrene	19.55	202	3456	0.44	ng/ul	97
18) Benzo(a)anthracene	21.30	228	2914	0.44	ng/ul	95
19) Chrysene	21.35	228	3235	0.43	ng/ul	95
21) Benzo(b)fluoranthene	22.89	252	3336	0.44	ng/ul	98
22) Benzo(k)fluoranthene	22.93	252	3368	0.43	ng/ul	98
23) Benzo(a)pyrene	23.48	252	3100	0.44	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.84	276	3798	0.44	ng/ul	97
25) Dibenzo(a,h)anthracene	25.85	278	3037	0.44	ng/ul	97
26) Benzo(g,h,i)perylene	26.54	276	3404	0.44	ng/ul	96

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

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