

Data Path : Z:\HPCHEM1\BNA_M\Data\BM110416\
 Data File : BM007762.D
 Acq On : 04 Nov 2016 19:45
 Operator : UM/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.435

Quant Time: Nov 04 22:16:27 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM102516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 04 23:14:40 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	1142	0.40	ng/ul	0.00
2) Naphthalene-d8	10.54	136	3795	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	1923	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	3643	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	3304	0.40	ng/ul	0.00
20) Perylene-d12	23.58	264	3296	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	1980	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.16	212	3852	0.47	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.58	128	4132	0.42	ng/ul	97
5) 2-Methylnaphthalene	12.21	142	2672	0.43	ng/ul	100
7) Acenaphthylene	14.09	152	3846	0.43	ng/ul	96
8) Acenaphthene	14.43	153	2712	0.39	ng/ul	89
9) Fluorene	15.44	166	3002	0.39	ng/ul	98
11) Pentachlorophenol	16.82	266	449	0.53	ng/ul	98
12) Phenanthrene	17.18	178	4386	0.43	ng/ul	95
13) Anthracene	17.26	178	4180	0.45	ng/ul	98
15) Fluoranthene	19.20	202	6412	0.53	ng/ul	98
17) Pyrene	19.56	202	5402	0.48	ng/ul	97
18) Benzo(a)anthracene	21.31	228	4511	0.48	ng/ul	97
19) Chrysene	21.35	228	4702	0.44	ng/ul	92
21) Benzo(b)fluoranthene	22.89	252	4972	0.44	ng/ul	96
22) Benzo(k)fluoranthene	22.93	252	4985	0.42	ng/ul	99
23) Benzo(a)pyrene	23.48	252	4855	0.46	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.84	276	5841	0.46	ng/ul	99
25) Dibenzo(a,h)anthracene	25.85	278	4591	0.45	ng/ul	96
26) Benzo(g,h,i)perylene	26.54	276	5086	0.44	ng/ul	98

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

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