

Data Path : S:\HPCHEM1\BNA_M\DATA\BM061115\
 Data File : BM001637.D
 Acq On : 10 Jun 2015 22:09
 Operator : TP/IZ
 Sample : SSTD0.286
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.286

Quant Time: Jun 11 07:49:21 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM061115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 11 07:43:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1680	0.40	ng/ul	0.00
4) Naphthalene-d8	10.48	136	5917	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.34	164	2712	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.09	188	5262	0.40	ng/ul	0.00
18) Chrysene-d12	21.27	240	3783	0.40	ng/ul	0.00
22) Perylene-d12	23.50	264	3346	0.40	ng/ul	0.00

System Monitoring Compounds						
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	12.08	152	1416	0.20	ng/ul	0.00
16) Fluoranthene-d10	19.12	212	2055	0.22	ng/ul	0.00

Target Compounds				Qvalue		
5) Naphthalene	10.53	128	3055	0.20	ng/ul	97
7) 2-Methylnaphthalene	12.14	142	1925	0.20	ng/ul	100
9) Acenaphthylene	14.05	152	2524	0.20	ng/ul	97
10) Acenaphthene	14.39	153	1927	0.19	ng/ul	100
11) Fluorene	15.38	166	2087	0.19	ng/ul	99
13) Pentachlorophenol	16.74	266	194	0.30	ng/ul	95
14) Phenanthrene	17.12	178	2975	0.20	ng/ul	97
15) Anthracene	17.22	178	2844	0.21	ng/ul	98
17) Fluoranthene	19.14	202	3118	0.21	ng/ul	96
19) Pyrene	19.50	202	3115	0.20	ng/ul	97
20) Benzo(a)anthracene	21.25	228	2439	0.22	ng/ul	98
21) Chrysene	21.31	228	2616	0.21	ng/ul	98
23) Benzo(b)fluoranthene	22.83	252	2507	0.21	ng/ul	93
24) Benzo(k)fluoranthene	22.87	252	2607	0.21	ng/ul#	93
25) Benzo(a)pyrene	23.40	252	2337	0.21	ng/ul#	92
26) Indeno(1,2,3-cd)pyrene	25.76	276	2729	0.23	ng/ul	99
27) Dibenzo(a,h)anthracene	25.78	278	2206	0.23	ng/ul	94
28) Benzo(g,h,i)perylene	26.44	276	2601	0.25	ng/ul	95

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

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