

Data Path : S:\HPCHEM1\BNA_M\DATA\BM061115\
 Data File : BM001642.D
 Acq On : 11 Jun 2015 10:10
 Operator : TP/IZ
 Sample : SSTD0.490
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.490

Quant Time: Jun 11 14:53:06 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 07:52:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1714	0.40	ng/ul	0.00
4) Naphthalene-d8	10.48	136	6112	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.34	164	2925	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.09	188	6124	0.40	ng/ul	0.00
18) Chrysene-d12	21.28	240	4472	0.40	ng/ul	0.00
22) Perylene-d12	23.52	264	3421	0.40	ng/ul	0.01

System Monitoring Compounds						
2) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
6) 2-Methylnaphthalene-d10	12.08	152	3049	0.41	ng/ul	0.00
16) Fluoranthene-d10	19.12	212	5127	0.41	ng/ul	0.00

Target Compounds				Qvalue		
5) Naphthalene	10.53	128	6509	0.40	ng/ul	99
7) 2-Methylnaphthalene	12.14	142	4145	0.41	ng/ul	99
9) Acenaphthylene	14.05	152	5640	0.40	ng/ul	99
10) Acenaphthene	14.39	153	4305	0.41	ng/ul	98
11) Fluorene	15.38	166	4753	0.42	ng/ul	100
13) Pentachlorophenol	16.73	266	425	0.35	ng/ul	94
14) Phenanthrene	17.12	178	7300	0.41	ng/ul	99
15) Anthracene	17.21	178	6690	0.39	ng/ul	98
17) Fluoranthene	19.14	202	7792	0.41	ng/ul	99
19) Pyrene	19.51	202	7876	0.44	ng/ul	98
20) Benzo(a)anthracene	21.25	228	5757	0.40	ng/ul	99
21) Chrysene	21.31	228	6369	0.41	ng/ul	99
23) Benzo(b)fluoranthene	22.84	252	5845	0.44	ng/ul	96
24) Benzo(k)fluoranthene	22.88	252	4933	0.37	ng/ul	99
25) Benzo(a)pyrene	23.41	252	4856	0.40	ng/ul	99
26) Indeno(1,2,3-cd)pyrene	25.77	276	5022	0.36	ng/ul	99
27) Dibenzo(a,h)anthracene	25.78	278	3954	0.35	ng/ul	99
28) Benzo(g,h,i)perylene	26.46	276	4530	0.35	ng/ul	99

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

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