

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE121817\
 Data File : BE095004.D
 Acq On : 18 Dec 2017 10:55
 Operator : SJ/JU
 Sample : SSTD0.101
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.101

Quant Time: Dec 18 15:33:28 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE121817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:33:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4709	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	14834	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	8309	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	21280	0.40	ng/ul	0.00
16) Chrysene-d12	21.85	240	17958	0.40	ng/ul	0.00
20) Perylene-d12	24.51	264	15406	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	2462	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	5855	0.08	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.34	128	3953	0.112	ng/ul#	91
5) 2-Methylnaphthalene	12.92	142	2709	0.110	ng/ul	99
7) Acenaphthylene	14.77	152	3737	0.098	ng/ul	98
8) Acenaphthene	15.11	153	2864	0.107	ng/ul	93
9) Fluorene	16.07	166	3315	0.102	ng/ul	98
12) Phenanthrene	17.80	178	5350	0.094	ng/ul	94
13) Anthracene	17.89	178	4299	0.074	ng/ul	98
15) Fluoranthene	19.76	202	6242	0.083	ng/ul	84
17) Pyrene	20.11	202	6358	0.113	ng/ul#	81
18) Benzo(a)anthracene	21.83	228	5115	0.094	ng/ul#	86
19) Chrysene	21.89	228	5519	0.101	ng/ul	96
21) Benzo(b)fluoranthene	23.68	252	5105	0.111	ng/ul	94
22) Benzo(k)fluoranthene	23.73	252	4850	0.099	ng/ul	95
23) Benzo(a)pyrene	24.38	252	4443	0.097	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	27.35	276	4843	0.097	ng/ul#	84
25) Dibenzo(a,h)anthracene	27.37	278	3780	0.090	ng/ul	97
26) Benzo(g,h,i)perylene	28.23	276	4316	0.104	ng/ul	96

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

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