

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094839.D
 Acq On : 22 Nov 2017 19:30
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.494

Quant Time: Nov 23 00:59:00 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 23 00:56:52 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	653	0.40	ng/ul	0.03
2) Naphthalene-d8	10.71	136	4074	0.40	ng/ul	0.02
6) Acenaphthene-d10	14.51	164	2896	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	7768	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	8480	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	8150	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	2622	0.40	ng/ul	0.03
14) Fluoranthene-d10	19.28	212	9031	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	3540	0.365	ng/ul#	95
5) 2-Methylnaphthalene	12.38	142	2681	0.397	ng/ul	97
7) Acenaphthylene	14.24	152	4669	0.352	ng/ul	98
8) Acenaphthene	14.58	153	3376	0.362	ng/ul	98
9) Fluorene	15.57	166	4238	0.374	ng/ul	93
11) Pentachlorophenol	17.01	266	198	0.483	ng/ul	96
12) Phenanthrene	17.29	178	7494	0.359	ng/ul#	87
13) Anthracene	17.40	178	7515	0.354	ng/ul#	91
15) Fluoranthene	19.32	202	9305	0.340	ng/ul	84
17) Pyrene	19.68	202	9800	0.369	ng/ul#	80
18) Benzo(a)anthracene	21.41	228	9168	0.356	ng/ul#	86
19) Chrysene	21.47	228	9519	0.370	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	8739	0.360	ng/ul	87
22) Benzo(k)fluoranthene	23.24	252	9313	0.361	ng/ul#	89
23) Benzo(a)pyrene	23.86	252	8759	0.363	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.68	276	9677	0.368	ng/ul#	79
25) Dibenzo(a,h)anthracene	26.67	278	8141	0.367	ng/ul	92
26) Benzo(g,h,i)perylene	27.52	276	8150	0.370	ng/ul	95

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

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