

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122818\
 Data File : BE098569.D
 Acq On : 28 Dec 2018 11:52
 Operator : SJ/JU
 Sample : SSTD1.606
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.606

Quant Time: Dec 28 12:28:28 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE122818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 28 11:34:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	893	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	4173	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	2669	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	6539	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	7892	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	7423	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.20	152	11308	1.52	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	34751	1.64	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.65	128	18205	1.823	ng/ul	96
5) 2-Methylnaphthalene	12.27	142	12608	1.670	ng/ul	99
7) Acenaphthylene	14.18	152	19230	1.826	ng/ul	96
8) Acenaphthene	14.53	153	14781	1.879	ng/ul	99
9) Fluorene	15.52	166	17851	1.885	ng/ul	100
11) Pentachlorophenol	16.89	266	1776	5.191	ng/ul	96
12) Phenanthrene	17.27	178	29073	1.926	ng/ul	99
13) Anthracene	17.36	178	27000	1.812	ng/ul	99
15) Fluoranthene	19.29	202	38015	1.976	ng/ul	97
17) Pyrene	19.65	202	42551	2.309	ng/ul	98
18) Benzo(a)anthracene	21.39	228	39455	1.700	ng/ul	99
19) Chrysene	21.45	228	43884	2.445	ng/ul	99
21) Benzo(b)fluoranthene	23.15	252	39560	1.924	ng/ul	97
22) Benzo(k)fluoranthene	23.20	252	39933	1.984	ng/ul	97
23) Benzo(a)pyrene	23.81	252	37371	1.889	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	26.56	276	43297	1.910	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.59	278	35699	1.890	ng/ul	97
26) Benzo(g,h,i)perylene	27.38	276	37086	1.944	ng/ul	95

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

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