

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122818\
 Data File : BE098588.D
 Acq On : 28 Dec 2018 23:36
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.410

Quant Time: Dec 29 00:18:28 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 : Sat Dec 29 00:05:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	798	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	3839	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	2554	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	6406	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	7646	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	7614	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.20	152	2555	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	8239	0.40	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.66	128	4040	0.386	ng/ul	98
5) 2-Methylnaphthalene	12.28	142	2745	0.413	ng/ul	99
7) Acenaphthylene	14.18	152	4338	0.396	ng/ul	99
8) Acenaphthene	14.53	153	3373	0.392	ng/ul	98
9) Fluorene	15.53	166	4053	0.396	ng/ul	96
11) Pentachlorophenol	16.89	266	495	0.521	ng/ul	98
12) Phenanthrene	17.27	178	6767	0.393	ng/ul	97
13) Anthracene	17.37	178	5974	0.414	ng/ul	99
15) Fluoranthene	19.29	202	8853	0.398	ng/ul	96
17) Pyrene	19.65	202	9183	0.376	ng/ul#	95
18) Benzo(a)anthracene	21.39	228	9258	0.415	ng/ul	99
19) Chrysene	21.45	228	10211	0.369	ng/ul	99
21) Benzo(b)fluoranthene	23.15	252	9397	0.403	ng/ul	98
22) Benzo(k)fluoranthene	23.21	252	9900	0.387	ng/ul	97
23) Benzo(a)pyrene	23.81	252	9251	0.407	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.57	276	10474	0.401	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.60	278	8566	0.403	ng/ul#	96
26) Benzo(g,h,i)perylene	27.38	276	8857	0.391	ng/ul	97

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

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