

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122818\
 Data File : BE098566.D
 Acq On : 28 Dec 2018 10:01
 Operator : SJ/JU
 Sample : SSTD0.203
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.203

Quant Time: Dec 28 11:37:26 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	799	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	3726	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	2454	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	6232	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	7286	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	7384	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	1188	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	3867	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.66	128	2062	0.231	ng/ul	95
5) 2-Methylnaphthalene	12.29	142	1252	0.186	ng/ul	98
7) Acenaphthylene	14.19	152	2050	0.212	ng/ul	97
8) Acenaphthene	14.53	153	1627	0.225	ng/ul	98
9) Fluorene	15.53	166	1903	0.219	ng/ul	97
11) Pentachlorophenol	16.90	266	117	0.359	ng/ul	94
12) Phenanthrene	17.27	178	3308	0.230	ng/ul	97
13) Anthracene	17.37	178	2563	0.180	ng/ul	94
15) Fluoranthene	19.29	202	4165	0.227	ng/ul	91
17) Pyrene	19.66	202	4303	0.253	ng/ul#	91
18) Benzo(a)anthracene	21.39	228	4027	0.188	ng/ul	99
19) Chrysene	21.45	228	5498	0.332	ng/ul	98
21) Benzo(b)fluoranthene	23.16	252	4303	0.210	ng/ul	95
22) Benzo(k)fluoranthene	23.21	252	4983	0.249	ng/ul#	95
23) Benzo(a)pyrene	23.81	252	4238	0.215	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	26.57	276	4889	0.217	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.59	278	3947	0.210	ng/ul#	92
26) Benzo(g,h,i)perylene	27.38	276	4328	0.228	ng/ul#	95

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

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