

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070218\
 Data File : BN001933.D
 Acq On : 02 Jul 2018 14:27
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.483

Quant Time: Jul 03 01:52:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN061318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 03 01:51:51 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	6048	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	23846	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.34	164	13648	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	31662	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	24772	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	19707	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	15223	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	34351	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.53	128	24293	0.395	ng/ul	96
5) 2-Methylnaphthalene	12.15	142	17297	0.394	ng/ul	98
7) Acenaphthylene	14.06	152	22319	0.334	ng/ul	95
8) Acenaphthene	14.40	153	18920	0.386	ng/ul	97
9) Fluorene	15.39	166	21737	0.379	ng/ul#	94
11) Pentachlorophenol	16.74	266	1966	0.328	ng/ul	98
12) Phenanthrene	17.13	178	37046	0.392	ng/ul#	89
13) Anthracene	17.22	178	31325	0.350	ng/ul#	91
15) Fluoranthene	19.15	202	39960	0.356	ng/ul	96
17) Pyrene	19.51	202	39315	0.437	ng/ul	97
18) Benzo(a)anthracene	21.27	228	30426	0.355	ng/ul#	85
19) Chrysene	21.32	228	33025	0.388	ng/ul	97
21) Benzo(b)fluoranthene	22.85	252	29190	0.408	ng/ul	86
22) Benzo(k)fluoranthene	22.89	252	29815	0.418	ng/ul#	88
23) Benzo(a)pyrene	23.42	252	25344	0.382	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.75	276	28999	0.389	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.76	278	22817	0.384	ng/ul#	91
26) Benzo(g,h,i)perylene	26.43	276	24077	0.380	ng/ul	97

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

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