

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051219\
 Data File : BN005629.D
 Acq On : 11 May 2019 16:40
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.469

Quant Time: May 12 00:36:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 07 07:04:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	1777	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	7303	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.23	164	4238	0.40	ng/ul	-0.02
10) Phenanthrene-d10	16.99	188	9770	0.40	ng/ul	-0.01
16) Chrysene-d12	21.22	240	7438	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	6799	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	4801	0.42	ng/ul	-0.02
14) Fluoranthene-d10	19.03	212	10445	0.34	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.41	128	7824	0.405	ng/ul	100
5) 2-Methylnaphthalene	12.04	142	5618	0.426	ng/ul	100
7) Acenaphthylene	13.95	152	8925	0.418	ng/ul	99
8) Acenaphthene	14.29	153	6181	0.409	ng/ul	98
9) Fluorene	15.30	166	7067	0.401	ng/ul	98
11) Pentachlorophenol	16.69	266	667	0.264	ng/ul	94
12) Phenanthrene	17.03	178	11507	0.432	ng/ul	100
13) Anthracene	17.13	178	10649	0.401	ng/ul	99
15) Fluoranthene	19.06	202	13036	0.354	ng/ul	99
17) Pyrene	19.43	202	13278	0.483	ng/ul	98
18) Benzo(a)anthracene	21.20	228	10793	0.421	ng/ul	100
19) Chrysene	21.26	228	12024	0.382	ng/ul	98
21) Benzo(b)fluoranthene	22.77	252	9635	0.471	ng/ul	98
22) Benzo(k)fluoranthene	22.82	252	10907	0.426	ng/ul	100
23) Benzo(a)pyrene	23.33	252	9618	0.421	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.62	276	10387	0.406	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.63	278	8538	0.404	ng/ul	100
26) Benzo(g,h,i)perylene	26.29	276	8761	0.402	ng/ul	99

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

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