

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111819\
 Data File : BN008606.D
 Acq On : 18 Nov 2019 16:31
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.409

Quant Time: Nov 18 17:03:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 15 22:41:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2710	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	9727	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	5662	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	11542	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	9536	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	10243	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	6469	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	14375	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	11750	0.401	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	8230	0.402	ng/ul	100
7) Acenaphthylene	13.97	152	12575	0.402	ng/ul	100
8) Acenaphthene	14.32	153	8779	0.378	ng/ul	99
9) Fluorene	15.31	166	10082	0.370	ng/ul	99
11) Pentachlorophenol	16.65	266	1846	0.423	ng/ul	97
12) Phenanthrene	17.04	178	16060	0.413	ng/ul	98
13) Anthracene	17.13	178	16425	0.447	ng/ul	99
15) Fluoranthene	19.06	202	18787	0.393	ng/ul	96
17) Pyrene	19.42	202	19172	0.501	ng/ul	99
18) Benzo(a)anthracene	21.18	228	17499	0.456	ng/ul#	92
19) Chrysene	21.23	228	16380	0.435	ng/ul#	93
21) Benzo(b)fluoranthene	22.72	252	17033	0.403	ng/ul#	44
22) Benzo(k)fluoranthene	22.77	252	16452	0.393	ng/ul#	39
23) Benzo(a)pyrene	23.27	252	16436	0.415	ng/ul#	44
24) Indeno(1,2,3-cd)pyrene	25.52	276	19165	0.413	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.53	278	15451	0.413	ng/ul#	61
26) Benzo(g,h,i)perylene	26.17	276	16164	0.417	ng/ul#	77

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

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