

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121619\
 Data File : BN008997.D
 Acq On : 16 Dec 2019 13:58
 Operator : JU
 Sample : SSTD1.606
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.606

Quant Time: Dec 16 14:36:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 16 13:30:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	2472	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	10886	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	6787	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	16115	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	13226	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	10526	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	30850	1.73	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	79393	1.61	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.35	128	52036	1.586	ng/ul	99
5) 2-Methylnaphthalene	11.97	142	39559	1.726	ng/ul	100
7) Acenaphthylene	13.89	152	59711	1.591	ng/ul	99
8) Acenaphthene	14.24	153	45178	1.622	ng/ul	99
9) Fluorene	15.23	166	54103	1.655	ng/ul	99
11) Pentachlorophenol	16.58	266	7705	1.263	ng/ul	99
12) Phenanthrene	16.97	178	90299	1.661	ng/ul	100
13) Anthracene	17.06	178	84995	1.655	ng/ul	99
15) Fluoranthene	18.99	202	107871	1.616	ng/ul	98
17) Pyrene	19.35	202	105781	1.994	ng/ul	100
18) Benzo(a)anthracene	21.11	228	96033	1.806	ng/ul	100
19) Chrysene	21.17	228	93082	1.782	ng/ul	100
21) Benzo(b)fluoranthene	22.64	252	81418	1.876	ng/ul	95
22) Benzo(k)fluoranthene	22.68	252	80255	1.864	ng/ul	95
23) Benzo(a)pyrene	23.18	252	71043	1.744	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.37	276	76059	1.596	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.39	278	61174	1.590	ng/ul	98
26) Benzo(g,h,i)perylene	26.02	276	62945	1.579	ng/ul	98

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

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