

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121619\
 Data File : BN008996.D
 Acq On : 16 Dec 2019 13:22
 Operator : JU
 Sample : SSTD0.805
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.805

Quant Time: Dec 16 14:35:52 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	2284	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	10008	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	6269	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	14970	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	12093	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	9944	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	13537	0.82	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	34100	0.75	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.36	128	22784	0.755	ng/ul	99
5) 2-Methylnaphthalene	11.97	142	17487	0.830	ng/ul	99
7) Acenaphthylene	13.89	152	25337	0.731	ng/ul	99
8) Acenaphthene	14.24	153	19887	0.773	ng/ul	99
9) Fluorene	15.23	166	23555	0.780	ng/ul	99
11) Pentachlorophenol	16.59	266	3011	0.531	ng/ul	99
12) Phenanthrene	16.97	178	39671	0.786	ng/ul	100
13) Anthracene	17.06	178	36092	0.757	ng/ul	99
15) Fluoranthene	18.99	202	46495	0.750	ng/ul	98
17) Pyrene	19.35	202	45692	0.942	ng/ul	99
18) Benzo(a)anthracene	21.11	228	40777	0.838	ng/ul	100
19) Chrysene	21.17	228	41190	0.862	ng/ul	99
21) Benzo(b)fluoranthene	22.64	252	35559	0.867	ng/ul	97
22) Benzo(k)fluoranthene	22.68	252	36035	0.886	ng/ul	99
23) Benzo(a)pyrene	23.18	252	31030	0.806	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.38	276	34180	0.759	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.39	278	27509	0.757	ng/ul	99
26) Benzo(g,h,i)perylene	26.01	276	28532	0.758	ng/ul	99

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

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