

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007439.D
 Acq On : 27 Aug 2019 19:28
 Operator : HP/JU
 Sample : SSTD1.630
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.630

Quant Time: Aug 27 19:59:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 18:27:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	4217	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	17828	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	9167	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	21463	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	18889	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	19179	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	44629	1.41	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	110776	1.60	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.20	128	80370	1.567	ng/ul	99
5) 2-Methylnaphthalene	11.83	142	55411	1.439	ng/ul	100
7) Acenaphthylene	13.76	152	84060	1.582	ng/ul	100
8) Acenaphthene	14.11	153	59422	1.635	ng/ul	98
9) Fluorene	15.11	166	68061	1.545	ng/ul	99
11) Pentachlorophenol	16.46	266	9120	1.542	ng/ul	98
12) Phenanthrene	16.84	178	112234	1.641	ng/ul	98
13) Anthracene	16.93	178	110260	1.576	ng/ul	99
15) Fluoranthene	18.87	202	140161	1.649	ng/ul	99
17) Pyrene	19.24	202	142063	1.685	ng/ul	96
18) Benzo(a)anthracene	21.01	228	139910	1.686	ng/ul	100
19) Chrysene	21.06	228	138258	1.779	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	141797	1.824	ng/ul	96
22) Benzo(k)fluoranthene	22.55	252	137470	1.824	ng/ul#	93
23) Benzo(a)pyrene	23.04	252	134702	1.840	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.19	276	153659	1.947	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.21	278	122496	1.944	ng/ul	97
26) Benzo(g,h,i)perylene	25.82	276	129674	1.959	ng/ul	99

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

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