

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN122818\
 Data File : BN004266.D
 Acq On : 28 Dec 2018 17:21
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :

Quant Time: Dec 28 22:12:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN122818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 28 16:45:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	2235	0.40	ng/ul	0.00
2) Naphthalene-d8	10.62	136	9922	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5561	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	12985	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	13762	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	13639	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	6044	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	13366	0.35	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.67	128	10912	0.390	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	7727	0.398	ng/ul	98
7) Acenaphthylene	14.19	152	10768	0.387	ng/ul	99
8) Acenaphthene	14.53	153	8705	0.391	ng/ul	99
9) Fluorene	15.52	166	9798	0.387	ng/ul	100
11) Pentachlorophenol	16.90	266	592	0.315	ng/ul	97
12) Phenanthrene	17.26	178	15983	0.379	ng/ul	100
13) Anthracene	17.36	178	14492	0.378	ng/ul	100
15) Fluoranthene	19.28	202	17766	0.349	ng/ul	99
17) Pyrene	19.64	202	18227	0.387	ng/ul	99
18) Benzo(a)anthracene	21.39	228	18328	0.392	ng/ul	100
19) Chrysene	21.44	228	19628	0.391	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	20457	0.395	ng/ul	100
22) Benzo(k)fluoranthene	23.07	252	21959	0.400	ng/ul	99
23) Benzo(a)pyrene	23.63	252	19246	0.391	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.11	276	18528	0.363	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.12	278	15100	0.361	ng/ul	100
26) Benzo(g,h,i)perylene	26.84	276	15571	0.363	ng/ul	99

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

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