

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051319\
 Data File : BN005668.D
 Acq On : 14 May 2019 00:52
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.474

Quant Time: May 14 05:28:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 14 05:24:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1205	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	5126	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	3330	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	8459	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	8307	0.40	ng/ul	-0.01
20) Perylene-d12	23.43	264	8119	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	3471	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	10609	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.41	128	5504	0.406	ng/ul	98
5) 2-Methylnaphthalene	12.04	142	4036	0.436	ng/ul	100
7) Acenaphthylene	13.95	152	6987	0.417	ng/ul	99
8) Acenaphthene	14.29	153	4896	0.413	ng/ul	99
9) Fluorene	15.30	166	5806	0.419	ng/ul	99
11) Pentachlorophenol	16.70	266	561	0.256	ng/ul#	82
12) Phenanthrene	17.03	178	9733	0.422	ng/ul	100
13) Anthracene	17.13	178	9485	0.413	ng/ul	100
15) Fluoranthene	19.06	202	12692	0.398	ng/ul	97
17) Pyrene	19.43	202	13307	0.434	ng/ul	98
18) Benzo(a)anthracene	21.21	228	12440	0.434	ng/ul	99
19) Chrysene	21.26	228	13642	0.388	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	11589	0.475	ng/ul	93
22) Benzo(k)fluoranthene	22.82	252	12591	0.412	ng/ul#	95
23) Benzo(a)pyrene	23.34	252	11407	0.418	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.63	276	11435	0.375	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.64	278	9372	0.372	ng/ul#	94
26) Benzo(g,h,i)perylene	26.30	276	9327	0.358	ng/ul	98

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

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