

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062419\
 Data File : BN006581.D
 Acq On : 25 Jun 2019 21:27
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.434

Quant Time: Jun 26 00:51:53 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 25 07:15:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	5317	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	20174	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	9439	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	16761	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	10302	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	12025	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	12160	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	16643	0.37	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.46	128	23849	0.428	ng/ul 100
5) 2-Methylnaphthalene	12.08	142	15350	0.396	ng/ul 99
7) Acenaphthylene	14.00	152	21249	0.405	ng/ul 100
8) Acenaphthene	14.34	153	15532	0.427	ng/ul 99
9) Fluorene	15.34	166	15886	0.373	ng/ul 97
11) Pentachlorophenol	16.71	266	2060	0.484	ng/ul 99
12) Phenanthrene	17.08	178	22095	0.411	ng/ul 99
13) Anthracene	17.17	178	20057	0.390	ng/ul 99
15) Fluoranthene	19.11	202	22023	0.371	ng/ul 97
17) Pyrene	19.48	202	22479	0.414	ng/ul 100
18) Benzo(a)anthracene	21.24	228	18282	0.401	ng/ul 99
19) Chrysene	21.29	228	19762	0.462	ng/ul 99
21) Benzo(b)fluoranthene	22.82	252	19341	0.386	ng/ul 95
22) Benzo(k)fluoranthene	22.87	252	20398	0.418	ng/ul 96
23) Benzo(a)pyrene	23.39	252	19572	0.419	ng/ul# 91
24) Indeno(1,2,3-cd)pyrene	25.73	276	22898	0.453	ng/ul# 97
25) Dibenzo(a,h)anthracene	25.74	278	18084	0.448	ng/ul 95
26) Benzo(g,h,i)perylene	26.41	276	19663	0.464	ng/ul 96

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

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