

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061419\
 Data File : BN006339.D
 Acq On : 14 Jun 2019 16:49
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Jun 14 18:05:57 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 : Thu Jun 13 11:20:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	3599	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	14166	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	7067	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	13817	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	10359	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	11531	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	8180	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	14922	0.40	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.47	128	15972	0.409	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	10270	0.377	ng/ul	98
7) Acenaphthylene	14.01	152	14733	0.375	ng/ul	100
8) Acenaphthene	14.36	153	11095	0.407	ng/ul	99
9) Fluorene	15.35	166	11287	0.354	ng/ul	98
11) Pentachlorophenol	16.72	266	1085	0.309	ng/ul	95
12) Phenanthrene	17.09	178	16887	0.381	ng/ul	99
13) Anthracene	17.19	178	14639	0.346	ng/ul	98
15) Fluoranthene	19.12	202	19487	0.398	ng/ul	98
17) Pyrene	19.49	202	20321	0.372	ng/ul	98
18) Benzo(a)anthracene	21.25	228	16494	0.360	ng/ul	99
19) Chrysene	21.31	228	19239	0.447	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	17656	0.367	ng/ul	96
22) Benzo(k)fluoranthene	22.88	252	19753	0.422	ng/ul	96
23) Benzo(a)pyrene	23.41	252	18233	0.407	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.75	276	20721	0.427	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.76	278	16326	0.421	ng/ul	94
26) Benzo(g,h,i)perylene	26.43	276	17857	0.440	ng/ul	96

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

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