

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121618\
 Data File : BN003989.D
 Acq On : 16 Dec 2018 11:39
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.421

Quant Time: Dec 17 04:02:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 17 00:13:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	2159	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	10369	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5769	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	14088	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	15238	0.40	ng/ul	0.01
20) Perylene-d12	23.73	264	13246	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	6057	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	16664	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	11080	0.380	ng/ul	98
5) 2-Methylnaphthalene	12.30	142	7916	0.389	ng/ul	99
7) Acenaphthylene	14.20	152	11808	0.455	ng/ul	100
8) Acenaphthene	14.54	153	8728	0.383	ng/ul	100
9) Fluorene	15.52	166	9891	0.373	ng/ul	98
11) Pentachlorophenol	16.88	266	1139	0.449	ng/ul	93
12) Phenanthrene	17.26	178	16882	0.365	ng/ul	99
13) Anthracene	17.36	178	16619	0.437	ng/ul	99
15) Fluoranthene	19.28	202	21908	0.401	ng/ul	98
17) Pyrene	19.64	202	22582	0.413	ng/ul	97
18) Benzo(a)anthracene	21.39	228	21009	0.437	ng/ul	96
19) Chrysene	21.44	228	20051	0.364	ng/ul	96
21) Benzo(b)fluoranthene	23.02	252	19083	0.348	ng/ul	90
22) Benzo(k)fluoranthene	23.07	252	19025	0.356	ng/ul#	90
23) Benzo(a)pyrene	23.63	252	17828	0.376	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.09	276	20355	0.374	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.10	278	16471	0.376	ng/ul#	88
26) Benzo(g,h,i)perylene	26.82	276	16880	0.357	ng/ul#	89

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

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