

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN121718\
 Data File : BN004034.D
 Acq On : 17 Dec 2018 14:15
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.425

Quant Time: Dec 18 00:11:19 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 06:22:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1280	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	6780	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4113	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	10381	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	12543	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	11459	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	4301	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	13847	0.46	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.68	128	7644	0.401	ng/ul 98
5) 2-Methylnaphthalene	12.29	142	5656	0.425	ng/ul 99
7) Acenaphthylene	14.19	152	8812	0.476	ng/ul 99
8) Acenaphthene	14.53	153	6621	0.408	ng/ul 99
9) Fluorene	15.52	166	7657	0.405	ng/ul 100
11) Pentachlorophenol	16.87	266	845	0.452	ng/ul 97
12) Phenanthrene	17.26	178	13526	0.397	ng/ul 98
13) Anthracene	17.35	178	13097	0.468	ng/ul 98
15) Fluoranthene	19.28	202	18225	0.453	ng/ul 96
17) Pyrene	19.64	202	18710	0.416	ng/ul 97
18) Benzo(a)anthracene	21.39	228	18292	0.462	ng/ul 97
19) Chrysene	21.44	228	17600	0.388	ng/ul 99
21) Benzo(b)fluoranthene	23.02	252	17290	0.365	ng/ul 94
22) Benzo(k)fluoranthene	23.07	252	17351	0.376	ng/ul# 94
23) Benzo(a)pyrene	23.63	252	16406	0.400	ng/ul# 94
24) Indeno(1,2,3-cd)pyrene	26.09	276	18725	0.398	ng/ul# 97
25) Dibenzo(a,h)anthracene	26.10	278	15317	0.404	ng/ul 94
26) Benzo(g,h,i)perylene	26.82	276	15822	0.387	ng/ul 94

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

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