

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003950.D
 Acq On : 15 Dec 2018 09:48
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Dec 15 21:55:51 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 : Wed Dec 12 03:46:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	2264	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	10864	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6266	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	15201	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	19004	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	16479	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	6466	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	17703	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	11565	0.379	ng/ul	98
5) 2-Methylnaphthalene	12.29	142	8386	0.394	ng/ul	99
7) Acenaphthylene	14.20	152	12267	0.435	ng/ul	100
8) Acenaphthene	14.53	153	9353	0.378	ng/ul	99
9) Fluorene	15.52	166	10766	0.373	ng/ul	99
11) Pentachlorophenol	16.87	266	1199	0.438	ng/ul	99
12) Phenanthrene	17.26	178	18253	0.366	ng/ul	100
13) Anthracene	17.35	178	17084	0.417	ng/ul	99
15) Fluoranthene	19.27	202	23479	0.399	ng/ul	96
17) Pyrene	19.63	202	24050	0.353	ng/ul	96
18) Benzo(a)anthracene	21.38	228	25129	0.419	ng/ul	98
19) Chrysene	21.43	228	24358	0.355	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	24377	0.358	ng/ul	97
22) Benzo(k)fluoranthene	23.06	252	24178	0.364	ng/ul	98
23) Benzo(a)pyrene	23.61	252	22361	0.379	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.07	276	24208	0.358	ng/ul	98
25) Dibenzo(a,h)anthracene	26.08	278	19717	0.362	ng/ul	97
26) Benzo(g,h,i)perylene	26.80	276	19982	0.340	ng/ul	96

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

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