

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101919\
 Data File : BN008261.D
 Acq On : 19 Oct 2019 09:53
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.449

Quant Time: Oct 19 22:27:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 23:47:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1794	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	8309	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	5322	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	10790	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	10464	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	10995	0.40	ng/ul	-0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	5800	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	14840	0.40	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.41	128	10343	0.437	ng/ul	100
5) 2-Methylnaphthalene	12.03	142	6949	0.428	ng/ul	100
7) Acenaphthylene	13.95	152	10120	0.399	ng/ul	100
8) Acenaphthene	14.29	153	8174	0.415	ng/ul	98
9) Fluorene	15.28	166	9085	0.406	ng/ul	97
11) Pentachlorophenol	16.67	266	743	0.285	ng/ul	98
12) Phenanthrene	17.02	178	14433	0.457	ng/ul	100
13) Anthracene	17.11	178	12710	0.457	ng/ul	100
15) Fluoranthene	19.04	202	17426	0.418	ng/ul	96
17) Pyrene	19.41	202	18140	0.435	ng/ul	99
18) Benzo(a)anthracene	21.17	228	16376	0.430	ng/ul	100
19) Chrysene	21.22	228	19602	0.420	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	17764	0.496	ng/ul	98
22) Benzo(k)fluoranthene	22.75	252	19304	0.476	ng/ul	97
23) Benzo(a)pyrene	23.26	252	17385	0.462	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.51	276	18071	0.407	ng/ul	96
25) Dibenzo(a,h)anthracene	25.53	278	14588	0.416	ng/ul	98
26) Benzo(g,h,i)perylene	26.17	276	15602	0.384	ng/ul	98

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

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