

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122319\
 Data File : BN009084.D
 Acq On : 23 Dec 2019 11:51
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Dec 23 14:00:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 20 23:31:06 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	3326	0.40	ng/ul	0.00
2) Naphthalene-d8	10.29	136	11983	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.16	164	6187	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.91	188	10715	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	6229	0.40	ng/ul	0.00
20) Perylene-d12	23.26	264	6184	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.89	152	7788	0.37	ng/ul	0.00
14) Fluoranthene-d10	18.94	212	10715	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.34	128	14944	0.411	ng/ul	99
5) 2-Methylnaphthalene	11.96	142	10145	0.369	ng/ul	100
7) Acenaphthylene	13.88	152	12775	0.393	ng/ul	100
8) Acenaphthene	14.22	153	9915	0.389	ng/ul	99
9) Fluorene	15.22	166	10571	0.351	ng/ul	99
11) Pentachlorophenol	16.58	266	741	0.259	ng/ul	96
12) Phenanthrene	16.95	178	15256	0.409	ng/ul	99
13) Anthracene	17.04	178	13319	0.399	ng/ul	99
15) Fluoranthene	18.98	202	14483	0.340	ng/ul	99
17) Pyrene	19.34	202	14190	0.447	ng/ul	98
18) Benzo(a)anthracene	21.10	228	11570	0.418	ng/ul	100
19) Chrysene	21.15	228	12017	0.432	ng/ul	99
21) Benzo(b)fluoranthene	22.63	252	11281	0.391	ng/ul	97
22) Benzo(k)fluoranthene	22.66	252	11963	0.411	ng/ul	97
23) Benzo(a)pyrene	23.16	252	10581	0.418	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.35	276	11992	0.435	ng/ul	100
25) Dibenzo(a,h)anthracene	25.37	278	9325	0.425	ng/ul	98
26) Benzo(g,h,i)perylene	25.99	276	10579	0.456	ng/ul	98

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

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