

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112019\
 Data File : BN008662.D
 Acq On : 20 Nov 2019 16:08
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.416

Quant Time: Nov 20 17:44:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 20 04:45:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	3445	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	13020	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	7228	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	14376	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	11739	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	11825	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	8177	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	16679	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.43	128	15140	0.386	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	10542	0.385	ng/ul	99
7) Acenaphthylene	13.95	152	14873	0.372	ng/ul	99
8) Acenaphthene	14.30	153	10791	0.364	ng/ul	99
9) Fluorene	15.29	166	12173	0.350	ng/ul	99
11) Pentachlorophenol	16.64	266	1755	0.323	ng/ul	99
12) Phenanthrene	17.03	178	19512	0.402	ng/ul	98
13) Anthracene	17.11	178	18349	0.401	ng/ul	99
15) Fluoranthene	19.04	202	22016	0.370	ng/ul	92
17) Pyrene	19.40	202	22429	0.476	ng/ul	96
18) Benzo(a)anthracene	21.16	228	21040	0.446	ng/ul	100
19) Chrysene	21.21	228	20094	0.433	ng/ul	99
21) Benzo(b)fluoranthene	22.70	252	19964	0.409	ng/ul	92
22) Benzo(k)fluoranthene	22.74	252	18969	0.392	ng/ul	96
23) Benzo(a)pyrene	23.25	252	18551	0.405	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.49	276	20939	0.391	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.50	278	16780	0.388	ng/ul	97
26) Benzo(g,h,i)perylene	26.13	276	17872	0.399	ng/ul	96

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

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