

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN123119\
 Data File : BN009285.D
 Acq On : 31 Dec 2019 17:57
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.408

Quant Time: Dec 31 18:28:55 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 : Tue Dec 31 11:02:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	2933	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	10264	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.14	164	5114	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	9944	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	7852	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8061	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.85	152	6556	0.36	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	11539	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.31	128	12668	0.407	ng/ul	99
5) 2-Methylnaphthalene	11.93	142	8258	0.351	ng/ul	99
7) Acenaphthylene	13.85	152	10396	0.387	ng/ul	99
8) Acenaphthene	14.20	153	8016	0.380	ng/ul	99
9) Fluorene	15.19	166	8776	0.352	ng/ul	98
11) Pentachlorophenol	16.55	266	1034	0.390	ng/ul	99
12) Phenanthrene	16.92	178	13921	0.403	ng/ul	99
13) Anthracene	17.02	178	12511	0.403	ng/ul	100
15) Fluoranthene	18.95	202	15339	0.388	ng/ul	99
17) Pyrene	19.31	202	15376	0.384	ng/ul	98
18) Benzo(a)anthracene	21.07	228	14240	0.408	ng/ul	99
19) Chrysene	21.12	228	14394	0.411	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	14169	0.376	ng/ul	99
22) Benzo(k)fluoranthene	22.63	252	15226	0.401	ng/ul	99
23) Benzo(a)pyrene	23.12	252	13774	0.418	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.30	276	16064	0.447	ng/ul	96
25) Dibenzo(a,h)anthracene	25.31	278	12900	0.451	ng/ul	99
26) Benzo(g,h,i)perylene	25.94	276	13759	0.455	ng/ul	98

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

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