

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN123119\
 Data File : BN009273.D
 Acq On : 31 Dec 2019 10:09
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.407

Quant Time: Dec 31 11:04:18 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	2725	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	9507	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.13	164	4815	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	9349	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	7360	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	7526	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.85	152	6084	0.36	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	10871	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.31	128	11719	0.406	ng/ul	100
5) 2-Methylnaphthalene	11.93	142	7825	0.359	ng/ul	100
7) Acenaphthylene	13.85	152	9924	0.392	ng/ul	99
8) Acenaphthene	14.19	153	7572	0.381	ng/ul	98
9) Fluorene	15.19	166	8330	0.355	ng/ul	99
11) Pentachlorophenol	16.55	266	914	0.367	ng/ul	99
12) Phenanthrene	16.92	178	13046	0.401	ng/ul	99
13) Anthracene	17.02	178	11801	0.405	ng/ul	99
15) Fluoranthene	18.95	202	14444	0.388	ng/ul	99
17) Pyrene	19.31	202	14549	0.387	ng/ul	99
18) Benzo(a)anthracene	21.07	228	13332	0.408	ng/ul	99
19) Chrysene	21.12	228	13765	0.419	ng/ul	100
21) Benzo(b)fluoranthene	22.59	252	13576	0.386	ng/ul	98
22) Benzo(k)fluoranthene	22.63	252	14031	0.396	ng/ul	98
23) Benzo(a)pyrene	23.12	252	12778	0.415	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.30	276	15166	0.452	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.32	278	12149	0.455	ng/ul	98
26) Benzo(g,h,i)perylene	25.94	276	13023	0.461	ng/ul	98

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

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