

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121319\
 Data File : BN008962.D
 Acq On : 13 Dec 2019 11:38
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4111

Quant Time: Dec 13 14:08:09 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 : Thu Dec 12 00:49:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2324	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	9833	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	6389	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	14246	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	11333	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	9550	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	6811	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	16749	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	11833	0.399	ng/ul	100
5) 2-Methylnaphthalene	11.98	142	8855	0.428	ng/ul	100
7) Acenaphthylene	13.90	152	12877	0.365	ng/ul	98
8) Acenaphthene	14.25	153	9986	0.381	ng/ul	99
9) Fluorene	15.24	166	11706	0.380	ng/ul	99
11) Pentachlorophenol	16.59	266	1446	0.268	ng/ul	99
12) Phenanthrene	16.97	178	19700	0.410	ng/ul	99
13) Anthracene	17.06	178	17682	0.390	ng/ul	99
15) Fluoranthene	18.99	202	22524	0.382	ng/ul	99
17) Pyrene	19.36	202	22553	0.496	ng/ul	96
18) Benzo(a)anthracene	21.12	228	20533	0.451	ng/ul	100
19) Chrysene	21.17	228	20581	0.460	ng/ul	99
21) Benzo(b)fluoranthene	22.64	252	17762	0.451	ng/ul	100
22) Benzo(k)fluoranthene	22.68	252	17904	0.458	ng/ul	99
23) Benzo(a)pyrene	23.18	252	15673	0.424	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.38	276	15979	0.370	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.39	278	12917	0.370	ng/ul	99
26) Benzo(g,h,i)perylene	26.02	276	13569	0.375	ng/ul	98

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

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