

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120719\
 Data File : BN008882.D
 Acq On : 07 Dec 2019 10:15
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.404

Quant Time: Dec 09 00:12:33 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 : Fri Dec 06 12:03:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2577	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	10697	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	7278	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	16255	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	14553	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	15101	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	7216	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	19181	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	12598	0.391	ng/ul	100
5) 2-Methylnaphthalene	11.99	142	9352	0.415	ng/ul	99
7) Acenaphthylene	13.90	152	14013	0.348	ng/ul	99
8) Acenaphthene	14.25	153	10894	0.365	ng/ul	99
9) Fluorene	15.25	166	13066	0.373	ng/ul	99
11) Pentachlorophenol	16.59	266	1808	0.294	ng/ul	99
12) Phenanthrene	16.97	178	22389	0.408	ng/ul	100
13) Anthracene	17.07	178	20335	0.393	ng/ul	100
15) Fluoranthene	19.00	202	26017	0.386	ng/ul	100
17) Pyrene	19.36	202	26241	0.449	ng/ul	99
18) Benzo(a)anthracene	21.12	228	25352	0.433	ng/ul	100
19) Chrysene	21.17	228	25482	0.443	ng/ul	100
21) Benzo(b)fluoranthene	22.65	252	25351	0.407	ng/ul	99
22) Benzo(k)fluoranthene	22.69	252	25873	0.419	ng/ul	99
23) Benzo(a)pyrene	23.19	252	23762	0.407	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.39	276	26328	0.385	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.41	278	21250	0.385	ng/ul	99
26) Benzo(g,h,i)perylene	26.03	276	21690	0.379	ng/ul	98

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

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