

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112119\
 Data File : BN008691.D
 Acq On : 22 Nov 2019 00:39
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Nov 22 06:35:04 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 Nov 22 06:30:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	3920	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	13670	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	7858	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	15336	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	13392	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	13646	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	8497	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	18168	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	16002	0.388	ng/ul	99
5) 2-Methylnaphthalene	12.03	142	10844	0.377	ng/ul	100
7) Acenaphthylene	13.95	152	15482	0.356	ng/ul	98
8) Acenaphthene	14.29	153	11277	0.350	ng/ul	99
9) Fluorene	15.28	166	12840	0.339	ng/ul	99
11) Pentachlorophenol	16.63	266	2232	0.385	ng/ul	97
12) Phenanthrene	17.02	178	20960	0.405	ng/ul	99
13) Anthracene	17.11	178	19531	0.400	ng/ul	100
15) Fluoranthene	19.04	202	24307	0.383	ng/ul	94
17) Pyrene	19.40	202	24697	0.460	ng/ul	98
18) Benzo(a)anthracene	21.15	228	23663	0.439	ng/ul	100
19) Chrysene	21.20	228	23055	0.436	ng/ul	99
21) Benzo(b)fluoranthene	22.69	252	23281	0.414	ng/ul	97
22) Benzo(k)fluoranthene	22.73	252	22698	0.407	ng/ul	97
23) Benzo(a)pyrene	23.23	252	21593	0.409	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.47	276	25050	0.406	ng/ul	93
25) Dibenzo(a,h)anthracene	25.48	278	20157	0.404	ng/ul	99
26) Benzo(g,h,i)perylene	26.11	276	21293	0.412	ng/ul	97

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

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