

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120619\
 Data File : BN008863.D
 Acq On : 06 Dec 2019 17:14
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.402

Quant Time: Dec 06 17:59:27 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	7.57	152	2377	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	10322	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	6718	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	14717	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	13538	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	14894	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	6840	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	17931	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	12066	0.388	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	8952	0.412	ng/ul	100
7) Acenaphthylene	13.91	152	13390	0.360	ng/ul	99
8) Acenaphthene	14.25	153	10077	0.366	ng/ul	99
9) Fluorene	15.25	166	11847	0.366	ng/ul	98
11) Pentachlorophenol	16.60	266	1751	0.314	ng/ul	98
12) Phenanthrene	16.98	178	20146	0.406	ng/ul	99
13) Anthracene	17.07	178	18741	0.400	ng/ul	99
15) Fluoranthene	19.00	202	23842	0.391	ng/ul	100
17) Pyrene	19.37	202	24338	0.448	ng/ul	98
18) Benzo(a)anthracene	21.12	228	24278	0.446	ng/ul	99
19) Chrysene	21.17	228	23437	0.438	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	24317	0.396	ng/ul	100
22) Benzo(k)fluoranthene	22.70	252	24273	0.398	ng/ul	99
23) Benzo(a)pyrene	23.20	252	23475	0.407	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.41	276	26859	0.398	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.42	278	21877	0.402	ng/ul	99
26) Benzo(g,h,i)perylene	26.05	276	22362	0.396	ng/ul	98

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

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