

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071819\
 Data File : BN006885.D
 Acq On : 18 Jul 2019 11:29
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.434

Quant Time: Jul 18 14:16:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 06:39:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	2384	0.40	ng/ul	0.02
2) Naphthalene-d8	10.37	136	9075	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.25	164	4651	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9739	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	7392	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	7769	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	5091	0.36	ng/ul	0.01
14) Fluoranthene-d10	19.06	212	10063	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	10119	0.388	ng/ul	97
5) 2-Methylnaphthalene	12.06	142	6204	0.352	ng/ul	98
7) Acenaphthylene	13.97	152	8467	0.353	ng/ul	98
8) Acenaphthene	14.31	153	7098	0.384	ng/ul	98
9) Fluorene	15.31	166	7229	0.357	ng/ul	98
11) Pentachlorophenol	16.73	266	399	0.275	ng/ul	95
12) Phenanthrene	17.06	178	11610	0.371	ng/ul	99
13) Anthracene	17.16	178	8752	0.323	ng/ul	99
15) Fluoranthene	19.09	202	13172	0.340	ng/ul	99
17) Pyrene	19.46	202	13952	0.413	ng/ul	100
18) Benzo(a)anthracene	21.23	228	10039	0.331	ng/ul	100
19) Chrysene	21.28	228	14186	0.419	ng/ul	99
21) Benzo(b)fluoranthene	22.81	252	12233	0.384	ng/ul	98
22) Benzo(k)fluoranthene	22.86	252	13504	0.405	ng/ul	98
23) Benzo(a)pyrene	23.39	252	11708	0.375	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.71	276	13785	0.437	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.71	278	10483	0.418	ng/ul	96
26) Benzo(g,h,i)perylene	26.40	276	12499	0.474	ng/ul	97

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

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