

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071819\
 Data File : BN006918.D
 Acq On : 19 Jul 2019 07:45
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.437

Quant Time: Jul 19 08:18:28 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 19 03:57:16 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	3188	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	11983	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.24	164	6187	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	13084	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	10486	0.40	ng/ul	-0.01
20) Perylene-d12	23.46	264	11812	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	6822	0.37	ng/ul	-0.01
14) Fluoranthene-d10	19.05	212	13849	0.36	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	13552	0.394	ng/ul	98
5) 2-Methylnaphthalene	12.05	142	8328	0.358	ng/ul	98
7) Acenaphthylene	13.96	152	11659	0.365	ng/ul	100
8) Acenaphthene	14.30	153	9413	0.382	ng/ul	99
9) Fluorene	15.31	166	9808	0.364	ng/ul	100
11) Pentachlorophenol	16.71	266	638	0.327	ng/ul	98
12) Phenanthrene	17.05	178	15695	0.373	ng/ul	100
13) Anthracene	17.15	178	12871	0.354	ng/ul	100
15) Fluoranthene	19.08	202	18012	0.346	ng/ul	98
17) Pyrene	19.45	202	18784	0.392	ng/ul	98
18) Benzo(a)anthracene	21.22	228	15053	0.349	ng/ul	100
19) Chrysene	21.27	228	19266	0.401	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	17471	0.361	ng/ul	97
22) Benzo(k)fluoranthene	22.84	252	19854	0.392	ng/ul	98
23) Benzo(a)pyrene	23.36	252	17805	0.375	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.68	276	20713	0.432	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.68	278	16234	0.425	ng/ul	96
26) Benzo(g,h,i)perylene	26.37	276	17995	0.448	ng/ul	97

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

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