

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071719\
 Data File : BN006883.D
 Acq On : 17 Jul 2019 21:16
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.433

Quant Time: Jul 18 01:08:48 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	2067	0.40	ng/ul	0.02
2) Naphthalene-d8	10.37	136	8165	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4260	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9021	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	7233	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	7933	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4567	0.36	ng/ul	0.01
14) Fluoranthene-d10	19.05	212	9571	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	9053	0.386	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	5486	0.346	ng/ul	99
7) Acenaphthylene	13.97	152	7891	0.359	ng/ul	99
8) Acenaphthene	14.31	153	6434	0.380	ng/ul	98
9) Fluorene	15.31	166	6594	0.355	ng/ul	99
11) Pentachlorophenol	16.73	266	333	0.248	ng/ul#	90
12) Phenanthrene	17.05	178	10546	0.363	ng/ul	99
13) Anthracene	17.16	178	8196	0.327	ng/ul	98
15) Fluoranthene	19.09	202	12370	0.345	ng/ul	99
17) Pyrene	19.45	202	13073	0.396	ng/ul	98
18) Benzo(a)anthracene	21.22	228	10050	0.338	ng/ul	100
19) Chrysene	21.27	228	13673	0.412	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	11764	0.362	ng/ul	97
22) Benzo(k)fluoranthene	22.83	252	13275	0.390	ng/ul	98
23) Benzo(a)pyrene	23.36	252	12014	0.377	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.68	276	13845	0.429	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.69	278	10790	0.421	ng/ul	97
26) Benzo(g,h,i)perylene	26.36	276	12322	0.457	ng/ul	96

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

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