

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071919\
 Data File : BN006939.D
 Acq On : 19 Jul 2019 23:54
 Operator : HP/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.440

Quant Time: Jul 20 00:27:53 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 23:29:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	2404	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	9397	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4862	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	10292	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	7977	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	8943	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	5343	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	10854	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	10666	0.395	ng/ul	98
5) 2-Methylnaphthalene	12.07	142	6517	0.358	ng/ul	99
7) Acenaphthylene	13.97	152	9146	0.364	ng/ul	98
8) Acenaphthene	14.31	153	7492	0.387	ng/ul	98
9) Fluorene	15.31	166	7684	0.362	ng/ul	99
11) Pentachlorophenol	16.72	266	699	0.456	ng/ul	95
12) Phenanthrene	17.05	178	12239	0.370	ng/ul	98
13) Anthracene	17.16	178	9553	0.334	ng/ul	98
15) Fluoranthene	19.09	202	14083	0.344	ng/ul	98
17) Pyrene	19.45	202	14935	0.410	ng/ul	98
18) Benzo(a)anthracene	21.22	228	11233	0.343	ng/ul	100
19) Chrysene	21.27	228	15418	0.422	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	13636	0.372	ng/ul	97
22) Benzo(k)fluoranthene	22.83	252	15222	0.397	ng/ul	98
23) Benzo(a)pyrene	23.36	252	13894	0.387	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.68	276	15796	0.435	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.68	278	12358	0.428	ng/ul	98
26) Benzo(g,h,i)perylene	26.36	276	13895	0.457	ng/ul	98

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

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