

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071919\
 Data File : BN006927.D
 Acq On : 19 Jul 2019 16:53
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.439

Quant Time: Jul 19 23:33:20 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	2161	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	8396	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4477	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9292	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	7103	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	8006	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	4824	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	9697	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	9628	0.399	ng/ul	97
5) 2-Methylnaphthalene	12.07	142	5867	0.360	ng/ul	97
7) Acenaphthylene	13.97	152	8230	0.356	ng/ul	98
8) Acenaphthene	14.31	153	6793	0.381	ng/ul	98
9) Fluorene	15.31	166	6939	0.355	ng/ul	100
11) Pentachlorophenol	16.72	266	504	0.364	ng/ul	92
12) Phenanthrene	17.06	178	11096	0.371	ng/ul	98
13) Anthracene	17.16	178	8341	0.323	ng/ul	99
15) Fluoranthene	19.09	202	12663	0.343	ng/ul	98
17) Pyrene	19.46	202	13411	0.413	ng/ul	98
18) Benzo(a)anthracene	21.23	228	9634	0.330	ng/ul	100
19) Chrysene	21.28	228	14064	0.432	ng/ul	99
21) Benzo(b)fluoranthene	22.81	252	12245	0.373	ng/ul	98
22) Benzo(k)fluoranthene	22.86	252	13821	0.403	ng/ul	98
23) Benzo(a)pyrene	23.39	252	12303	0.383	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.71	276	14315	0.440	ng/ul	94
25) Dibenzo(a,h)anthracene	25.71	278	11103	0.429	ng/ul	97
26) Benzo(g,h,i)perylene	26.40	276	12722	0.468	ng/ul	97

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

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