

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071919\
 Data File : BN006920.D
 Acq On : 19 Jul 2019 12:05
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.438

Quant Time: Jul 19 23:32:52 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	2367	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	9087	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4705	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9732	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	7560	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	8810	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	5189	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	10155	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	10388	0.398	ng/ul	97
5) 2-Methylnaphthalene	12.06	142	6297	0.357	ng/ul	97
7) Acenaphthylene	13.96	152	8697	0.358	ng/ul	99
8) Acenaphthene	14.30	153	7220	0.386	ng/ul	99
9) Fluorene	15.31	166	7341	0.358	ng/ul	98
11) Pentachlorophenol	16.71	266	526	0.363	ng/ul	97
12) Phenanthrene	17.05	178	11699	0.374	ng/ul	99
13) Anthracene	17.16	178	9009	0.333	ng/ul	98
15) Fluoranthene	19.09	202	13247	0.342	ng/ul	100
17) Pyrene	19.45	202	14044	0.407	ng/ul	99
18) Benzo(a)anthracene	21.22	228	10505	0.338	ng/ul	100
19) Chrysene	21.28	228	14952	0.431	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	12861	0.356	ng/ul	98
22) Benzo(k)fluoranthene	22.85	252	14628	0.387	ng/ul	98
23) Benzo(a)pyrene	23.38	252	13456	0.380	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.70	276	15159	0.423	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.70	278	11770	0.413	ng/ul	97
26) Benzo(g,h,i)perylene	26.38	276	13375	0.447	ng/ul	97

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

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