

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071119\
 Data File : BN006812.D
 Acq On : 11 Jul 2019 23:27
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.424

Quant Time: Jul 12 08:12:16 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	7.60	152	2109	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	7929	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4226	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9333	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	7889	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	8869	0.40	ng/ul	-0.02

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	4554	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	10195	0.37	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.41	128	8832	0.388	ng/ul	96
5) 2-Methylnaphthalene	12.05	142	5489	0.357	ng/ul	98
7) Acenaphthylene	13.96	152	7919	0.363	ng/ul	99
8) Acenaphthene	14.31	153	6338	0.377	ng/ul	98
9) Fluorene	15.31	166	6743	0.366	ng/ul	99
11) Pentachlorophenol	16.71	266	491	0.353	ng/ul	93
12) Phenanthrene	17.05	178	11029	0.367	ng/ul	100
13) Anthracene	17.15	178	8903	0.343	ng/ul	99
15) Fluoranthene	19.08	202	13140	0.354	ng/ul	99
17) Pyrene	19.45	202	13907	0.386	ng/ul	99
18) Benzo(a)anthracene	21.21	228	11229	0.346	ng/ul	99
19) Chrysene	21.27	228	14298	0.395	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	13002	0.358	ng/ul	96
22) Benzo(k)fluoranthene	22.84	252	14147	0.372	ng/ul	98
23) Benzo(a)pyrene	23.36	252	12887	0.362	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.68	276	15099	0.419	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.68	278	11895	0.415	ng/ul	96
26) Benzo(g,h,i)perylene	26.36	276	13251	0.440	ng/ul	96

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

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