

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071119\
 Data File : BN006823.D
 Acq On : 12 Jul 2019 08:08
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.425

Quant Time: Jul 12 08:43:49 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	2193	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	8313	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4460	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9809	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	7942	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	8775	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4714	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	10422	0.36	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.42	128	9316	0.390	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	5651	0.350	ng/ul	98
7) Acenaphthylene	13.96	152	8127	0.353	ng/ul	99
8) Acenaphthene	14.31	153	6704	0.378	ng/ul	99
9) Fluorene	15.31	166	7044	0.362	ng/ul	99
11) Pentachlorophenol	16.72	266	389	0.266	ng/ul#	85
12) Phenanthrene	17.05	178	11562	0.366	ng/ul	99
13) Anthracene	17.15	178	8901	0.326	ng/ul	99
15) Fluoranthene	19.09	202	13527	0.347	ng/ul	99
17) Pyrene	19.45	202	14249	0.393	ng/ul	99
18) Benzo(a)anthracene	21.22	228	10829	0.332	ng/ul	99
19) Chrysene	21.28	228	14744	0.405	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	12620	0.351	ng/ul	97
22) Benzo(k)fluoranthene	22.85	252	14719	0.391	ng/ul	98
23) Benzo(a)pyrene	23.37	252	12852	0.365	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.70	276	15084	0.423	ng/ul	93
25) Dibenzo(a,h)anthracene	25.70	278	11757	0.415	ng/ul	96
26) Benzo(g,h,i)perylene	26.38	276	13372	0.449	ng/ul	97

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

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