

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN080620\
 Data File : BN011372.D
 Acq On : 06 Aug 2020 10:34
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.418

Quant Time: Aug 06 12:10:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 06 12:09:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	909	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	2825	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	1579	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	3387	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	3239	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	3140	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	2473	0.35	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	5711	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.26	128	3089	0.368	ng/ul	95
5) 2-Methylnaphthalene	11.89	142	1953	0.347	ng/ul	99
7) Acenaphthylene	13.81	152	2825	0.375	ng/ul	98
8) Acenaphthene	14.15	153	2234	0.373	ng/ul	99
9) Fluorene	15.15	166	2490	0.353	ng/ul	99
11) Pentachlorophenol	16.52	266	258	0.311	ng/ul	98
12) Phenanthrene	16.89	178	4066	0.362	ng/ul	98
13) Anthracene	16.98	178	3433	0.350	ng/ul	96
15) Fluoranthene	18.92	202	4856	0.351	ng/ul	97
17) Pyrene	19.28	202	4904	0.333	ng/ul	99
18) Benzo(a)anthracene	21.05	228	4084	0.328	ng/ul	99
19) Chrysene	21.10	228	5192	0.364	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	5022	0.383	ng/ul	91
22) Benzo(k)fluoranthene	22.60	252	5535	0.377	ng/ul	96
23) Benzo(a)pyrene	23.09	252	4291	0.377	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.25	276	5665	0.473	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.26	278	4582	0.480	ng/ul	97
26) Benzo(g,h,i)perylene	25.87	276	5213	0.482	ng/ul	99

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

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