

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072020\
 Data File : BN011205.D
 Acq On : 20 Jul 2020 11:45
 Operator : CG/JU
 Sample : SSTD1.603
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.603

Quant Time: Jul 20 12:17:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN072020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 20 11:25:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	1191	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	4153	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	2447	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.87	188	5183	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	3953	0.40	ng/ul	0.00
20) Perylene-d12	23.23	264	3145	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	16574	1.70	ng/ul	0.00
14) Fluoranthene-d10	18.91	212	37529	1.68	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.28	128	19350	1.579	ng/ul	97
5) 2-Methylnaphthalene	11.91	142	13176	1.715	ng/ul	100
7) Acenaphthylene	13.83	152	18979	1.740	ng/ul	98
8) Acenaphthene	14.18	153	14715	1.610	ng/ul	99
9) Fluorene	15.18	166	17207	1.712	ng/ul	99
11) Pentachlorophenol	16.54	266	1893	1.560	ng/ul	97
12) Phenanthrene	16.92	178	27483	1.629	ng/ul	99
13) Anthracene	17.01	178	24147	1.733	ng/ul	96
15) Fluoranthene	18.94	202	32680	1.727	ng/ul	98
17) Pyrene	19.31	202	31322	1.823	ng/ul	97
18) Benzo(a)anthracene	21.08	228	23402	1.667	ng/ul	97
19) Chrysene	21.13	228	26950	1.510	ng/ul	99
21) Benzo(b)fluoranthene	22.60	252	22597	1.613	ng/ul	87
22) Benzo(k)fluoranthene	22.64	252	24062	1.623	ng/ul#	88
23) Benzo(a)pyrene	23.14	252	18344	1.551	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	25.31	276	22310	1.376	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.33	278	17534	1.325	ng/ul#	86
26) Benzo(g,h,i)perylene	25.93	276	20820	1.388	ng/ul	96

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

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