

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081220\
 Data File : BN011417.D
 Acq On : 12 Aug 2020 12:37
 Operator : CG/JU
 Sample : SSTD1.606
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.606

Quant Time: Aug 12 13:32:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 12 11:51:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	759	0.40	ng/ul	0.00
2) Naphthalene-d8	10.19	136	2081	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.08	164	1063	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.83	188	2344	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	2407	0.40	ng/ul	0.00
20) Perylene-d12	23.17	264	2135	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	4961	1.18	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	12983	1.26	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.24	128	9191	1.659	ng/ul	98
5) 2-Methylnaphthalene	11.86	142	6110	1.869	ng/ul	99
7) Acenaphthylene	13.79	152	8575	1.780	ng/ul	97
8) Acenaphthene	14.14	153	6727	1.757	ng/ul	98
9) Fluorene	15.14	166	8759	1.961	ng/ul	99
11) Pentachlorophenol	16.50	266	978	1.881	ng/ul	98
12) Phenanthrene	16.87	178	12725	1.768	ng/ul	97
13) Anthracene	16.97	178	11492	1.895	ng/ul	97
15) Fluoranthene	18.90	202	18162	2.039	ng/ul	98
17) Pyrene	19.27	202	14795	1.564	ng/ul	97
18) Benzo(a)anthracene	21.04	228	13708	1.823	ng/ul	97
19) Chrysene	21.09	228	16213	1.670	ng/ul	99
21) Benzo(b)fluoranthene	22.54	252	18011	2.234	ng/ul	93
22) Benzo(k)fluoranthene	22.58	252	19879	2.223	ng/ul	93
23) Benzo(a)pyrene	23.08	252	12905	1.788	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.22	276	17484	1.813	ng/ul	99
25) Dibenzo(a,h)anthracene	25.23	278	14535	1.852	ng/ul	96
26) Benzo(g,h,i)perylene	25.83	276	16726	1.799	ng/ul	97

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

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