

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073020\
 Data File : BN011304.D
 Acq On : 30 Jul 2020 18:04
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.408

Quant Time: Jul 30 18:40:00 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 Jul 30 16:02:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	1516	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	5601	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	3558	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	8154	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	7130	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	6608	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	6035	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	16213	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	6862	0.412	ng/ul	99
5) 2-Methylnaphthalene	11.89	142	4751	0.425	ng/ul	98
7) Acenaphthylene	13.82	152	7225	0.425	ng/ul	99
8) Acenaphthene	14.16	153	5725	0.424	ng/ul	100
9) Fluorene	15.16	166	6726	0.424	ng/ul	99
11) Pentachlorophenol	16.52	266	851	0.426	ng/ul	99
12) Phenanthrene	16.90	178	11241	0.416	ng/ul	99
13) Anthracene	16.99	178	9928	0.421	ng/ul	99
15) Fluoranthene	18.93	202	13638	0.409	ng/ul	97
17) Pyrene	19.29	202	13675	0.422	ng/ul	100
18) Benzo(a)anthracene	21.06	228	11611	0.424	ng/ul	100
19) Chrysene	21.11	228	12553	0.400	ng/ul	99
21) Benzo(b)fluoranthene	22.57	252	11288	0.409	ng/ul	98
22) Benzo(k)fluoranthene	22.62	252	12999	0.420	ng/ul	98
23) Benzo(a)pyrene	23.11	252	10126	0.422	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.27	276	10165	0.403	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.29	278	8019	0.400	ng/ul	97
26) Benzo(g,h,i)perylene	25.90	276	8819	0.387	ng/ul	99

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

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