

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102020\
 Data File : BN012378.D
 Acq On : 20 Oct 2020 14:43
 Operator : CG/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.428

Quant Time: Oct 20 15:39:31 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 20 10:51:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	997	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.36	136	4079	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	2228	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	4449	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	3935	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	3809	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	2599	0.43	ng/ul	-0.02
14) Fluoranthene-d10	19.02	212	4530	0.37	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.40	128	4711	0.400	ng/ul	99
5) 2-Methylnaphthalene	12.03	142	3181	0.425	ng/ul	99
7) Acenaphthylene	13.95	152	3986	0.412	ng/ul	98
8) Acenaphthene	14.29	153	3386	0.414	ng/ul	98
9) Fluorene	15.29	166	3787	0.416	ng/ul	97
11) Pentachlorophenol	16.64	266	516	0.401	ng/ul	97
12) Phenanthrene	17.02	178	5848	0.412	ng/ul	99
13) Anthracene	17.11	178	5131	0.422	ng/ul	98
15) Fluoranthene	19.05	202	6363	0.384	ng/ul	99
17) Pyrene	19.41	202	6551	0.369	ng/ul	99
18) Benzo(a)anthracene	21.17	228	5591	0.384	ng/ul	99
19) Chrysene	21.22	228	6335	0.357	ng/ul	98
21) Benzo(b)fluoranthene	22.71	252	7007	0.455	ng/ul	96
22) Benzo(k)fluoranthene	22.76	252	7089	0.416	ng/ul	97
23) Benzo(a)pyrene	23.27	252	6237	0.417	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.53	276	7552	0.393	ng/ul	100
25) Dibenzo(a,h)anthracene	25.54	278	6064	0.394	ng/ul	99
26) Benzo(g,h,i)perylene	26.19	276	6700	0.380	ng/ul	97

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

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