

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090420\
 Data File : BN011698.D
 Acq On : 04 Sep 2020 12:44
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.410

Quant Time: Sep 04 13:20:07 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 : Fri Sep 04 13:19:04 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1106	0.40	ng/ul	0.00
2) Naphthalene-d8	10.53	136	3864	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	2135	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.13	188	4725	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	4256	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	3951	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	2578	0.47	ng/ul	0.00
14) Fluoranthene-d10	19.16	212	5299	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	4655	0.413	ng/ul	99
5) 2-Methylnaphthalene	12.19	142	3146	0.442	ng/ul	98
7) Acenaphthylene	14.10	152	4291	0.399	ng/ul	98
8) Acenaphthene	14.44	153	3357	0.401	ng/ul	98
9) Fluorene	15.43	166	3946	0.400	ng/ul	98
11) Pentachlorophenol	16.78	266	626	0.558	ng/ul	96
12) Phenanthrene	17.16	178	6474	0.405	ng/ul	99
13) Anthracene	17.26	178	5898	0.449	ng/ul	97
15) Fluoranthene	19.19	202	7435	0.404	ng/ul	99
17) Pyrene	19.55	202	7490	0.450	ng/ul	98
18) Benzo(a)anthracene	21.30	228	6715	0.470	ng/ul	98
19) Chrysene	21.35	228	7533	0.391	ng/ul	100
21) Benzo(b)fluoranthene	22.90	252	7541	0.439	ng/ul	97
22) Benzo(k)fluoranthene	22.94	252	7929	0.418	ng/ul	97
23) Benzo(a)pyrene	23.47	252	6473	0.449	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.84	276	8588	0.440	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.85	278	7032	0.442	ng/ul	98
26) Benzo(g,h,i)perylene	26.53	276	7437	0.407	ng/ul	97

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

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