

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112020\
 Data File : BN012680.D
 Acq On : 21 Nov 2020 11:13
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.410

Quant Time: Nov 23 01:00:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 00:58:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	886	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	3633	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2026	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	3920	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	3259	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	3372	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	2116	0.40	ng/ul	-0.01
14) Fluoranthene-d10	18.89	212	3924	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.26	128	4313	0.394	ng/ul	97
5) 2-Methylnaphthalene	11.89	142	2912	0.417	ng/ul	99
7) Acenaphthylene	13.82	152	4193	0.424	ng/ul	98
8) Acenaphthene	14.16	153	3197	0.384	ng/ul	98
9) Fluorene	15.16	166	3553	0.388	ng/ul	99
11) Pentachlorophenol	16.51	266	533	0.428	ng/ul	95
12) Phenanthrene	16.89	178	5440	0.399	ng/ul	99
13) Anthracene	16.99	178	4857	0.410	ng/ul	97
15) Fluoranthene	18.92	202	6128	0.389	ng/ul	98
17) Pyrene	19.29	202	6287	0.431	ng/ul	95
18) Benzo(a)anthracene	21.05	228	5362	0.459	ng/ul	99
19) Chrysene	21.10	228	6210	0.426	ng/ul	99
21) Benzo(b)fluoranthene	22.55	252	6291	0.406	ng/ul	98
22) Benzo(k)fluoranthene	22.59	252	6498	0.363	ng/ul	98
23) Benzo(a)pyrene	23.09	252	5790	0.397	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.26	276	7369	0.386	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.27	278	5866	0.382	ng/ul	97
26) Benzo(g,h,i)perylene	25.89	276	6486	0.377	ng/ul	96

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

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