

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112420\
 Data File : BN012719.D
 Acq On : 24 Nov 2020 16:51
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.413

Quant Time: Nov 24 17:33:06 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 : Tue Nov 24 11:59:48 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	782	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	3124	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	1760	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	3488	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	3207	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	3311	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.80	152	1866	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	3528	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.25	128	3836	0.408	ng/ul	98
5) 2-Methylnaphthalene	11.88	142	2628	0.438	ng/ul	99
7) Acenaphthylene	13.81	152	3797	0.442	ng/ul	99
8) Acenaphthene	14.16	153	2900	0.401	ng/ul	98
9) Fluorene	15.15	166	3227	0.406	ng/ul	99
11) Pentachlorophenol	16.51	266	539	0.487	ng/ul	95
12) Phenanthrene	16.89	178	4992	0.411	ng/ul	98
13) Anthracene	16.98	178	4515	0.428	ng/ul	99
15) Fluoranthene	18.92	202	6270	0.447	ng/ul	100
17) Pyrene	19.28	202	6373	0.444	ng/ul	95
18) Benzo(a)anthracene	21.05	228	6221	0.541	ng/ul	99
19) Chrysene	21.10	228	7214	0.503	ng/ul	100
21) Benzo(b)fluoranthene	22.56	252	7261	0.477	ng/ul	97
22) Benzo(k)fluoranthene	22.60	252	7887	0.449	ng/ul	97
23) Benzo(a)pyrene	23.09	252	6753	0.471	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.26	276	8748	0.467	ng/ul	98
25) Dibenzo(a,h)anthracene	25.28	278	6894	0.458	ng/ul	96
26) Benzo(g,h,i)perylene	25.90	276	7468	0.443	ng/ul	97

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

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