

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011751.D
 Acq On : 08 Sep 2020 23:44
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.498

Quant Time: Sep 09 06:37:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN090820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 09 06:29:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2272	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8625	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4601	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	9750	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	8082	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	7662	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	5776	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	11034	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	10943	0.429	ng/ul	99
5) 2-Methylnaphthalene	12.18	142	7232	0.425	ng/ul	100
7) Acenaphthylene	14.09	152	9490	0.428	ng/ul	100
8) Acenaphthene	14.44	153	7574	0.425	ng/ul	100
9) Fluorene	15.43	166	8668	0.421	ng/ul	98
11) Pentachlorophenol	16.77	266	1189	0.404	ng/ul	98
12) Phenanthrene	17.16	178	13888	0.424	ng/ul	98
13) Anthracene	17.25	178	12569	0.430	ng/ul	99
15) Fluoranthene	19.18	202	15620	0.429	ng/ul	99
17) Pyrene	19.54	202	15584	0.434	ng/ul	97
18) Benzo(a)anthracene	21.29	228	13886	0.442	ng/ul	100
19) Chrysene	21.35	228	14997	0.427	ng/ul	98
21) Benzo(b)fluoranthene	22.88	252	15070	0.420	ng/ul	98
22) Benzo(k)fluoranthene	22.92	252	15394	0.418	ng/ul	99
23) Benzo(a)pyrene	23.46	252	13464	0.434	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.82	276	17493	0.429	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.83	278	14130	0.425	ng/ul	98
26) Benzo(g,h,i)perylene	26.50	276	14624	0.411	ng/ul	98

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

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