

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003078.D
 Acq On : 10 Oct 2018 23:27
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.402

Quant Time: Oct 11 07:08:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 07:04:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	1273	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	5714	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3033	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	5916	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	4460	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4670	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3698	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	6386	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	6261	0.427	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	4209	0.417	ng/ul	100
7) Acenaphthylene	14.00	152	5917	0.426	ng/ul	100
8) Acenaphthene	14.33	153	4555	0.433	ng/ul	98
9) Fluorene	15.34	166	4926	0.408	ng/ul	99
11) Pentachlorophenol	16.75	266	189	0.329	ng/ul	98
12) Phenanthrene	17.08	178	7131	0.432	ng/ul	100
13) Anthracene	17.18	178	6541	0.420	ng/ul	99
15) Fluoranthene	19.10	202	7203	0.372	ng/ul	97
17) Pyrene	19.47	202	7246	0.418	ng/ul	99
18) Benzo(a)anthracene	21.22	228	6047	0.421	ng/ul	99
19) Chrysene	21.27	228	6496	0.446	ng/ul	100
21) Benzo(b)fluoranthene	22.79	252	6647	0.395	ng/ul	97
22) Benzo(k)fluoranthene	22.83	252	6947	0.398	ng/ul	96
23) Benzo(a)pyrene	23.36	252	6579	0.436	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.67	276	8060	0.465	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.67	278	6428	0.464	ng/ul	95
26) Benzo(g,h,i)perylene	26.35	276	6796	0.465	ng/ul	96

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

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