

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003375.D
 Acq On : 01 Nov 2018 12:12
 Operator : MJ/SJ
 Sample : SSTD1.629
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.629

Quant Time: Nov 01 12:45:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 11:33:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	8394	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	33029	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	19189	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	44432	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	47991	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	34891	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	86326	1.69	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	213206	1.68	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.70	128	149829	1.768	ng/ul	99
5) 2-Methylnaphthalene	12.31	142	109996	1.885	ng/ul	100
7) Acenaphthylene	14.21	152	140939	1.605	ng/ul	100
8) Acenaphthene	14.55	153	121345	1.821	ng/ul	100
9) Fluorene	15.54	166	143642	1.882	ng/ul	99
11) Pentachlorophenol	16.88	266	15025	3.485	ng/ul	99
12) Phenanthrene	17.28	178	240220	1.939	ng/ul	99
13) Anthracene	17.36	178	208647	1.783	ng/ul	100
15) Fluoranthene	19.29	202	284958	1.961	ng/ul	99
17) Pyrene	19.65	202	272324	1.461	ng/ul	99
18) Benzo(a)anthracene	21.39	228	253892	1.644	ng/ul	100
19) Chrysene	21.45	228	271857	1.734	ng/ul	100
21) Benzo(b)fluoranthene	23.02	252	251807	2.004	ng/ul	98
22) Benzo(k)fluoranthene	23.07	252	253649	1.947	ng/ul	99
23) Benzo(a)pyrene	23.62	252	214107	1.901	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.08	276	226711	1.752	ng/ul	100
25) Dibenzo(a,h)anthracene	26.10	278	187419	1.809	ng/ul	99
26) Benzo(g,h,i)perylene	26.81	276	188971	1.732	ng/ul	99

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

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