

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122818\
 Data File : BN004261.D
 Acq On : 28 Dec 2018 14:26
 Operator : JU/SJ
 Sample : SSTD0.454
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.454

Quant Time: Dec 28 15:24:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN122818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 28 15:21:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	2164	0.40	ng/ul	0.00
2) Naphthalene-d8	10.62	136	9760	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5486	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	12655	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14273	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	13047	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.21	152	5957	0.46	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	14293	0.49	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.67	128	10745	0.344	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	7596	0.383	ng/ul	99
7) Acenaphthylene	14.19	152	10644	0.390	ng/ul	100
8) Acenaphthene	14.53	153	8540	0.324	ng/ul	100
9) Fluorene	15.52	166	9540	0.334	ng/ul	100
11) Pentachlorophenol	16.90	266	604	0.237	ng/ul	97
12) Phenanthrene	17.26	178	16290	0.315	ng/ul	100
13) Anthracene	17.36	178	14421	0.376	ng/ul	100
15) Fluoranthene	19.28	202	18683	0.373	ng/ul	100
17) Pyrene	19.64	202	19104	0.227	ng/ul	100
18) Benzo(a)anthracene	21.39	228	18974	0.339	ng/ul	100
19) Chrysene	21.44	228	20289	0.266	ng/ul	100
21) Benzo(b)fluoranthene	23.02	252	19862	0.253	ng/ul	100
22) Benzo(k)fluoranthene	23.07	252	20187	0.273	ng/ul	100
23) Benzo(a)pyrene	23.63	252	18317	0.298	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.10	276	18889	0.272	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.11	278	15486	0.310	ng/ul	100
26) Benzo(g,h,i)perylene	26.83	276	15503	0.227	ng/ul	100

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

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