

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN021419\
 Data File : BN004433.D
 Acq On : 14 Feb 2019 22:42
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 STD0.411

Quant Time: Feb 15 01:24:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 14:04:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	838	0.40	ng/ul	0.00
2) Naphthalene-d8	10.53	136	3395	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	2085	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.13	188	5752	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	7631	0.40	ng/ul	0.00
20) Perylene-d12	23.60	264	8161	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	2169	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.16	212	7007	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.58	128	3833	0.393	ng/ul	99
5) 2-Methylnaphthalene	12.20	142	2803	0.402	ng/ul	99
7) Acenaphthylene	14.11	152	3496	0.367	ng/ul	98
8) Acenaphthene	14.45	153	3205	0.388	ng/ul	99
9) Fluorene	15.43	166	3891	0.393	ng/ul	100
11) Pentachlorophenol	16.77	266	363	0.296	ng/ul	99
12) Phenanthrene	17.17	178	7321	0.392	ng/ul	99
13) Anthracene	17.26	178	5884	0.373	ng/ul	100
15) Fluoranthene	19.19	202	8850	0.386	ng/ul	99
17) Pyrene	19.56	202	9046	0.368	ng/ul	99
18) Benzo(a)anthracene	21.31	228	8988	0.365	ng/ul	99
19) Chrysene	21.36	228	11004	0.385	ng/ul	99
21) Benzo(b)fluoranthene	22.91	252	13006	0.399	ng/ul	98
22) Benzo(k)fluoranthene	22.95	252	12287	0.383	ng/ul	99
23) Benzo(a)pyrene	23.49	252	11710	0.401	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.88	276	16142	0.421	ng/ul	100
25) Dibenzo(a,h)anthracene	25.90	278	13216	0.425	ng/ul	99
26) Benzo(g,h,i)perylene	26.59	276	14352	0.413	ng/ul	99

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

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