

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN021119\
 Data File : BN004382.D
 Acq On : 11 Feb 2019 17:32
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.408

Quant Time: Feb 11 18:04:27 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	1121	0.40	ng/ul	0.00
2) Naphthalene-d8	10.54	136	4402	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.39	164	2591	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.14	188	6934	0.40	ng/ul	0.00
16) Chrysene-d12	21.33	240	8480	0.40	ng/ul	0.00
20) Perylene-d12	23.60	264	9092	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.13	152	2774	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.17	212	8515	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.59	128	4959	0.392	ng/ul	100
5) 2-Methylnaphthalene	12.20	142	3584	0.397	ng/ul	99
7) Acenaphthylene	14.11	152	4455	0.376	ng/ul	99
8) Acenaphthene	14.45	153	3976	0.387	ng/ul	97
9) Fluorene	15.44	166	4754	0.387	ng/ul	99
11) Pentachlorophenol	16.78	266	508	0.344	ng/ul	96
12) Phenanthrene	17.18	178	8716	0.387	ng/ul	100
13) Anthracene	17.27	178	7177	0.378	ng/ul	100
15) Fluoranthene	19.20	202	10448	0.378	ng/ul	100
17) Pyrene	19.56	202	10663	0.391	ng/ul	99
18) Benzo(a)anthracene	21.32	228	10353	0.379	ng/ul	100
19) Chrysene	21.37	228	12332	0.388	ng/ul	100
21) Benzo(b)fluoranthene	22.92	252	14167	0.390	ng/ul	99
22) Benzo(k)fluoranthene	22.96	252	13674	0.383	ng/ul	99
23) Benzo(a)pyrene	23.50	252	13046	0.401	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.89	276	17296	0.405	ng/ul	100
25) Dibenzo(a,h)anthracene	25.91	278	14174	0.409	ng/ul	100
26) Benzo(g,h,i)perylene	26.60	276	15189	0.392	ng/ul	100

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

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