

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110218\
 Data File : BN003405.D
 Acq On : 02 Nov 2018 16:19
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.435

Quant Time: Nov 02 22:17:54 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 : Fri Nov 02 23:13:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	9640	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	39707	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	21741	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	48945	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	45723	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	37340	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	25157	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	57261	0.42	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.70	128	45764	0.411	ng/ul	100
5) 2-Methylnaphthalene	12.31	142	32565	0.407	ng/ul	100
7) Acenaphthylene	14.21	152	44979	0.476	ng/ul	99
8) Acenaphthene	14.55	153	34279	0.406	ng/ul	99
9) Fluorene	15.54	166	38892	0.391	ng/ul	100
11) Pentachlorophenol	16.88	266	5375	0.548	ng/ul	99
12) Phenanthrene	17.28	178	63661	0.397	ng/ul	100
13) Anthracene	17.36	178	58782	0.449	ng/ul	100
15) Fluoranthene	19.29	202	74241	0.412	ng/ul	100
17) Pyrene	19.65	202	74846	0.450	ng/ul	99
18) Benzo(a)anthracene	21.40	228	68414	0.477	ng/ul	99
19) Chrysene	21.45	228	64099	0.393	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	59368	0.369	ng/ul	97
22) Benzo(k)fluoranthene	23.07	252	56997	0.366	ng/ul	96
23) Benzo(a)pyrene	23.62	252	53954	0.402	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.09	276	59100	0.411	ng/ul	100
25) Dibenzo(a,h)anthracene	26.10	278	47878	0.408	ng/ul	96
26) Benzo(g,h,i)perylene	26.81	276	49398	0.400	ng/ul	97

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

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