

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101919\
 Data File : BN008271.D
 Acq On : 19 Oct 2019 15:51
 Operator : JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.404

Quant Time: Oct 19 22:28:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 19 22:25:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1619	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	7594	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	5015	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	10610	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	10534	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	10663	0.40	ng/ul	-0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	5343	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	15133	0.41	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.41	128	9403	0.435	ng/ul	100
5) 2-Methylnaphthalene	12.03	142	6407	0.432	ng/ul	100
7) Acenaphthylene	13.95	152	9546	0.400	ng/ul	100
8) Acenaphthene	14.29	153	7705	0.415	ng/ul	98
9) Fluorene	15.28	166	8670	0.411	ng/ul	98
11) Pentachlorophenol	16.66	266	752	0.293	ng/ul	98
12) Phenanthrene	17.02	178	14084	0.453	ng/ul	100
13) Anthracene	17.11	178	12540	0.459	ng/ul	100
15) Fluoranthene	19.04	202	17623	0.430	ng/ul	98
17) Pyrene	19.40	202	18429	0.439	ng/ul	97
18) Benzo(a)anthracene	21.16	228	16324	0.426	ng/ul	99
19) Chrysene	21.22	228	19892	0.423	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	17076	0.492	ng/ul	96
22) Benzo(k)fluoranthene	22.75	252	18986	0.483	ng/ul	97
23) Benzo(a)pyrene	23.26	252	16836	0.461	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.51	276	17261	0.401	ng/ul	96
25) Dibenzo(a,h)anthracene	25.52	278	13876	0.408	ng/ul	98
26) Benzo(g,h,i)perylene	26.16	276	14852	0.376	ng/ul	98

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

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