

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA N\DATA\BN030218\
 Data File : BN000230.D
 Acq On : 02 Mar 2018 11:36
 Operator : JU/SJ
 Sample : SSTD0.855
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.855

Quant Time: Mar 05 09:37:10 2018
 Quant Method : Z:\HPCHEM1\BNA N\Methods\SOM-EPA-SIM-BN030218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 09:35:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	3643	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	11846	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	5170	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.78	188	11429	0.40	ng/ul	0.00
16) Chrysene-d12	21.90	240	7898	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6042	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	12772	0.82	ng/ul	0.00
14) Fluoranthene-d10	19.77	212	22147	0.84	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.34	128	30978	0.817	ng/ul#	88
5) 2-Methylnaphthalene	12.92	142	19786	0.822	ng/ul	99
7) Acenaphthylene	14.79	152	21217	0.825	ng/ul	95
8) Acenaphthene	15.12	153	20605	0.830	ng/ul	95
9) Fluorene	16.10	166	22774	0.845	ng/ul#	91
11) Pentachlorophenol	17.43	266	1789	0.779	ng/ul	97
12) Phenanthrene	17.82	178	39029	0.837	ng/ul#	88
13) Anthracene	17.91	178	29423	0.849	ng/ul#	90
15) Fluoranthene	19.80	202	38792	0.857	ng/ul	62
17) Pyrene	20.16	202	37465	0.806	ng/ul#	55
18) Benzo(a)anthracene	21.89	228	25403	0.820	ng/ul#	83
19) Chrysene	21.94	228	33763	0.800	ng/ul	95
21) Benzo(b)fluoranthene	23.63	252	30574	0.841	ng/ul	86
22) Benzo(k)fluoranthene	23.68	252	29443	0.859	ng/ul#	88
23) Benzo(a)pyrene	24.27	252	24217	0.851	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.94	276	27892	0.869	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.95	278	20666	0.892	ng/ul#	79
26) Benzo(g,h,i)perylene	27.72	276	26519	0.838	ng/ul#	83

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

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