

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA N\DATA\BN030218\
 Data File : BN000231.D
 Acq On : 02 Mar 2018 12:10
 Operator : JU/SJ
 Sample : SSTD1.656
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.656

Quant Time: Mar 05 09:37:16 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	3298	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	10912	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	4859	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.78	188	10495	0.40	ng/ul	0.00
16) Chrysene-d12	21.90	240	8304	0.40	ng/ul	0.00
20) Perylene-d12	24.37	264	6240	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	24558	1.71	ng/ul	0.00
14) Fluoranthene-d10	19.77	212	44533	1.83	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.34	128	58725	1.681	ng/ul#	87
5) 2-Methylnaphthalene	12.92	142	38248	1.726	ng/ul	99
7) Acenaphthylene	14.79	152	42094	1.741	ng/ul	95
8) Acenaphthene	15.13	153	39803	1.706	ng/ul	95
9) Fluorene	16.10	166	44769	1.767	ng/ul#	91
11) Pentachlorophenol	17.43	266	3866	1.832	ng/ul	98
12) Phenanthrene	17.82	178	74844	1.747	ng/ul#	88
13) Anthracene	17.92	178	59921	1.883	ng/ul#	90
15) Fluoranthene	19.80	202	80133	1.929	ng/ul	62
17) Pyrene	20.16	202	76397	1.563	ng/ul#	55
18) Benzo(a)anthracene	21.88	228	57586	1.767	ng/ul#	84
19) Chrysene	21.93	228	71687	1.616	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	65555	1.746	ng/ul	86
22) Benzo(k)fluoranthene	23.67	252	65216	1.843	ng/ul#	88
23) Benzo(a)pyrene	24.26	252	54446	1.853	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.93	276	64048	1.931	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.94	278	48276	2.018	ng/ul#	78
26) Benzo(g,h,i)perylene	27.71	276	58358	1.787	ng/ul#	82

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

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