

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA N\DATA\BN030218\
 Data File : BN000228.D
 Acq On : 02 Mar 2018 10:27
 Operator : JU/SJ
 Sample : SSTD0.253
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.253

Quant Time: Mar 05 09:36:54 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	4377	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	13958	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	5934	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.78	188	13285	0.40	ng/ul	0.00
16) Chrysene-d12	21.90	240	7687	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6186	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	3587	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.77	212	5797	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.34	128	8874	0.199	ng/ul#	88
5) 2-Methylnaphthalene	12.92	142	5517	0.195	ng/ul	99
7) Acenaphthylene	14.79	152	5717	0.194	ng/ul	97
8) Acenaphthene	15.13	153	5594	0.196	ng/ul	94
9) Fluorene	16.10	166	5905	0.191	ng/ul#	91
11) Pentachlorophenol	17.43	266	400	0.150	ng/ul	95
12) Phenanthrene	17.82	178	10546	0.195	ng/ul#	88
13) Anthracene	17.92	178	7397	0.184	ng/ul#	91
15) Fluoranthene	19.80	202	9536	0.181	ng/ul	62
17) Pyrene	20.16	202	9582	0.212	ng/ul#	55
18) Benzo(a)anthracene	21.88	228	5862	0.194	ng/ul#	84
19) Chrysene	21.94	228	8346	0.203	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	7300	0.196	ng/ul	89
22) Benzo(k)fluoranthene	23.67	252	6689	0.191	ng/ul#	90
23) Benzo(a)pyrene	24.26	252	5589	0.192	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.93	276	6060	0.184	ng/ul#	78
25) Dibenzo(a,h)anthracene	26.94	278	4246	0.179	ng/ul#	81
26) Benzo(g,h,i)perylene	27.72	276	6211	0.192	ng/ul#	82

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

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