

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN002989.D
 Acq On : 08 Oct 2018 16:26
 Operator : MJ/SJ
 Sample : SSTD0.890
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.890

Quant Time: Oct 08 17:03:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 16:26:19 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	1386	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	6062	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3593	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	8536	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	6852	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4987	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	7846	0.90	ng/ul	-0.01
14) Fluoranthene-d10	19.06	212	20029	0.91	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.46	128	12655	0.830	ng/ul	97
5) 2-Methylnaphthalene	12.08	142	9066	0.917	ng/ul	97
7) Acenaphthylene	13.99	152	13839	0.912	ng/ul	98
8) Acenaphthene	14.33	153	10455	0.844	ng/ul	99
9) Fluorene	15.33	166	12165	0.920	ng/ul	99
11) Pentachlorophenol	16.73	266	579	0.602	ng/ul	96
12) Phenanthrene	17.07	178	19808	0.841	ng/ul	99
13) Anthracene	17.17	178	18875	0.922	ng/ul	98
15) Fluoranthene	19.09	202	22798	0.898	ng/ul	100
17) Pyrene	19.46	202	22825	1.021	ng/ul	97
18) Benzo(a)anthracene	21.22	228	18143	0.958	ng/ul	99
19) Chrysene	21.27	228	18223	0.825	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	14292	0.867	ng/ul	94
22) Benzo(k)fluoranthene	22.83	252	14817	0.855	ng/ul#	93
23) Benzo(a)pyrene	23.36	252	13226	0.838	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.67	276	15322	0.805	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.67	278	12294	0.807	ng/ul	92
26) Benzo(g,h,i)perylene	26.35	276	12919	0.786	ng/ul	95

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

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