

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110118\
 Data File : BN003374.D
 Acq On : 01 Nov 2018 11:37
 Operator : MJ/SJ
 Sample : SSTD0.828
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.828

Quant Time: Nov 01 12:14:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 11:33:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	8252	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	31966	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	18374	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	44150	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	45567	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	34080	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	41978	0.85	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	103420	0.82	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.70	128	73141	0.892	ng/ul	100
5) 2-Methylnaphthalene	12.31	142	53176	0.942	ng/ul	100
7) Acenaphthylene	14.21	152	65372	0.778	ng/ul	100
8) Acenaphthene	14.56	153	58677	0.920	ng/ul	99
9) Fluorene	15.54	166	69708	0.954	ng/ul	100
11) Pentachlorophenol	16.88	266	6723	1.569	ng/ul	99
12) Phenanthrene	17.28	178	119518	0.971	ng/ul	100
13) Anthracene	17.36	178	98717	0.849	ng/ul	100
15) Fluoranthene	19.29	202	138362	0.958	ng/ul	100
17) Pyrene	19.65	202	132289	0.747	ng/ul	100
18) Benzo(a)anthracene	21.39	228	117298	0.800	ng/ul	100
19) Chrysene	21.45	228	131773	0.885	ng/ul	100
21) Benzo(b)fluoranthene	23.02	252	121623	0.991	ng/ul	99
22) Benzo(k)fluoranthene	23.07	252	121695	0.956	ng/ul	99
23) Benzo(a)pyrene	23.62	252	101973	0.927	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.08	276	109308	0.865	ng/ul	100
25) Dibenzo(a,h)anthracene	26.10	278	89989	0.889	ng/ul	99
26) Benzo(g,h,i)perylene	26.81	276	92764	0.870	ng/ul	100

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

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