

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011020\
 Data File : BN009334.D
 Acq On : 10 Jan 2020 14:52
 Operator : JU
 Sample : SSTD0.806
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.806

Quant Time: Jan 10 15:24:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN011020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 10 14:47:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	2317	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	7877	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	4476	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	10364	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	9129	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	9050	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	9872	0.71	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	24215	0.79	ng/ul	0.00

Target Compounds						Ovalue
3) Naphthalene	10.28	128	17914	0.750	ng/ul	98
5) 2-Methylnaphthalene	11.90	142	12588	0.697	ng/ul	99
7) Acenaphthylene	13.82	152	17916	0.762	ng/ul	99
8) Acenaphthene	14.17	153	14069	0.763	ng/ul	100
9) Fluorene	15.17	166	16430	0.753	ng/ul	99
11) Pentachlorophenol	16.53	266	2258	0.817	ng/ul	98
12) Phenanthrene	16.90	178	27344	0.759	ng/ul	97
13) Anthracene	17.00	178	24712	0.765	ng/ul	98
15) Fluoranthene	18.93	202	32990	0.800	ng/ul	99
17) Pyrene	19.29	202	32732	0.703	ng/ul	99
18) Benzo(a)anthracene	21.05	228	30890	0.762	ng/ul	99
19) Chrysene	21.10	228	31635	0.777	ng/ul	97
21) Benzo(b)fluoranthene	22.56	252	30660	0.726	ng/ul	92
22) Benzo(k)fluoranthene	22.60	252	33530	0.787	ng/ul#	91
23) Benzo(a)pyrene	23.09	252	29489	0.796	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.26	276	35536	0.881	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.27	278	28286	0.881	ng/ul	96
26) Benzo(g,h,i)perylene	25.89	276	30647	0.902	ng/ul	96

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

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