

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN006983.D
 Acq On : 07 Aug 2019 12:30
 Operator : HP/JU
 Sample : SSTD0.823
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.823

Quant Time: Aug 07 17:26:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:13:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	4869	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	18567	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	11862	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	30254	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	28146	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	30342	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	25327	0.77	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	74752	0.77	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.28	128	40862	0.765	ng/ul	98
5) 2-Methylnaphthalene	11.91	142	30764	0.767	ng/ul	100
7) Acenaphthylene	13.83	152	52598	0.765	ng/ul	99
8) Acenaphthene	14.18	153	35842	0.762	ng/ul	99
9) Fluorene	15.17	166	43794	0.768	ng/ul	99
11) Pentachlorophenol	16.52	266	6194	0.587	ng/ul#	82
12) Phenanthrene	16.90	178	73642	0.764	ng/ul	100
13) Anthracene	17.00	178	75498	0.765	ng/ul	99
15) Fluoranthene	18.93	202	92120	0.769	ng/ul	100
17) Pyrene	19.30	202	94069	0.749	ng/ul	100
18) Benzo(a)anthracene	21.07	228	94570	0.765	ng/ul	100
19) Chrysene	21.12	228	88288	0.763	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	93673	0.762	ng/ul	98
22) Benzo(k)fluoranthene	22.63	252	90737	0.761	ng/ul	97
23) Benzo(a)pyrene	23.12	252	88639	0.765	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.31	276	96920	0.776	ng/ul	100
25) Dibenzo(a,h)anthracene	25.33	278	77406	0.776	ng/ul	99
26) Benzo(g,h,i)perylene	25.94	276	81487	0.778	ng/ul	98

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

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