

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062719\
 Data File : BN006600.D
 Acq On : 27 Jun 2019 13:29
 Operator : HP/JU
 Sample : SSTD1.617
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.617

Quant Time: Jun 27 14:21:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 14:04:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2410	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	9989	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	5542	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	12466	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	11425	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	11830	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	24960	1.63	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	59153	1.77	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.45	128	45393	1.647	ng/ul	99
5) 2-Methylnaphthalene	12.08	142	31723	1.651	ng/ul	98
7) Acenaphthylene	13.99	152	48024	1.559	ng/ul	99
8) Acenaphthene	14.34	153	35864	1.678	ng/ul	100
9) Fluorene	15.33	166	40715	1.628	ng/ul	99
11) Pentachlorophenol	16.71	266	3261	1.029	ng/ul	94
12) Phenanthrene	17.08	178	65620	1.640	ng/ul	99
13) Anthracene	17.17	178	60560	1.585	ng/ul	98
15) Fluoranthene	19.11	202	80135	1.814	ng/ul	99
17) Pyrene	19.48	202	81805	1.359	ng/ul	98
18) Benzo(a)anthracene	21.26	228	75901	1.502	ng/ul	99
19) Chrysene	21.32	228	80486	1.697	ng/ul	99
21) Benzo(b)fluoranthene	22.86	252	78479	1.591	ng/ul	96
22) Benzo(k)fluoranthene	22.90	252	82796	1.723	ng/ul	96
23) Benzo(a)pyrene	23.43	252	76393	1.663	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.77	276	79082	1.590	ng/ul	100
25) Dibenzo(a,h)anthracene	25.77	278	63073	1.587	ng/ul	97
26) Benzo(g,h,i)perylene	26.45	276	65159	1.564	ng/ul	99

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

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