

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060619\
 Data File : BN006148.D
 Acq On : 06 Jun 2019 22:19
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.497

Quant Time: Jun 07 00:29:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 07 00:25:28 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5597	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	21349	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10313	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	21222	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	18325	0.40	ng/ul	-0.03
20) Perylene-d12	23.50	264	19217	0.40	ng/ul	-0.06

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	11952	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	23304	0.41	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.48	128	22891	0.389	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	14962	0.364	ng/ul	98
7) Acenaphthylene	14.02	152	22285	0.389	ng/ul	100
8) Acenaphthene	14.36	153	15663	0.394	ng/ul	99
9) Fluorene	15.36	166	16605	0.357	ng/ul	98
11) Pentachlorophenol	16.71	266	1909	0.354	ng/ul	98
12) Phenanthrene	17.09	178	25359	0.372	ng/ul	99
13) Anthracene	17.19	178	23268	0.358	ng/ul	99
15) Fluoranthene	19.12	202	30817	0.410	ng/ul	97
17) Pyrene	19.49	202	32032	0.332	ng/ul	98
18) Benzo(a)anthracene	21.25	228	26947	0.332	ng/ul	99
19) Chrysene	21.30	228	31208	0.410	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	27916	0.348	ng/ul	95
22) Benzo(k)fluoranthene	22.87	252	30977	0.397	ng/ul	98
23) Benzo(a)pyrene	23.40	252	29237	0.392	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.73	276	33277	0.412	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.74	278	26335	0.408	ng/ul	94
26) Benzo(g,h,i)perylene	26.41	276	28090	0.415	ng/ul	97

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

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