

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061319\
 Data File : BN006324.D
 Acq On : 13 Jun 2019 22:52
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Jun 14 01:06:40 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 : Thu Jun 13 11:20:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	4730	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	18678	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9230	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	18155	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	13512	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	14952	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	10910	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	19322	0.40	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.47	128	21074	0.409	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	13591	0.378	ng/ul	97
7) Acenaphthylene	14.01	152	19402	0.378	ng/ul	99
8) Acenaphthene	14.36	153	14592	0.410	ng/ul	99
9) Fluorene	15.35	166	15049	0.361	ng/ul	98
11) Pentachlorophenol	16.72	266	1269	0.275	ng/ul	98
12) Phenanthrene	17.09	178	22337	0.383	ng/ul	99
13) Anthracene	17.19	178	19423	0.349	ng/ul	99
15) Fluoranthene	19.12	202	25507	0.396	ng/ul	96
17) Pyrene	19.49	202	26457	0.372	ng/ul	100
18) Benzo(a)anthracene	21.25	228	21561	0.361	ng/ul	98
19) Chrysene	21.30	228	25064	0.447	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	22778	0.365	ng/ul	94
22) Benzo(k)fluoranthene	22.87	252	25245	0.416	ng/ul	96
23) Benzo(a)pyrene	23.40	252	23465	0.404	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.73	276	26944	0.429	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.74	278	21192	0.422	ng/ul	94
26) Benzo(g,h,i)perylene	26.42	276	23144	0.440	ng/ul	95

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

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