

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008306.D
 Acq On : 22 Oct 2019 09:28
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.445

Quant Time: Oct 22 13:17:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 21 23:58:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1665	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	8202	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5504	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	11458	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	10213	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	9455	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	5463	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	14577	0.37	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.40	128	9355	0.401	ng/ul	100
5) 2-Methylnaphthalene	12.02	142	6480	0.405	ng/ul	100
7) Acenaphthylene	13.94	152	10001	0.381	ng/ul	100
8) Acenaphthene	14.28	153	7766	0.381	ng/ul	98
9) Fluorene	15.28	166	8654	0.374	ng/ul	97
11) Pentachlorophenol	16.65	266	828	0.299	ng/ul	98
12) Phenanthrene	17.01	178	13960	0.416	ng/ul	100
13) Anthracene	17.11	178	12854	0.435	ng/ul	100
15) Fluoranthene	19.04	202	16865	0.381	ng/ul	97
17) Pyrene	19.40	202	17627	0.433	ng/ul	98
18) Benzo(a)anthracene	21.17	228	15244	0.411	ng/ul	99
19) Chrysene	21.22	228	17474	0.383	ng/ul	100
21) Benzo(b)fluoranthene	22.71	252	13777	0.448	ng/ul	99
22) Benzo(k)fluoranthene	22.75	252	14894	0.427	ng/ul	100
23) Benzo(a)pyrene	23.26	252	13282	0.410	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.51	276	14106	0.370	ng/ul	93
25) Dibenzo(a,h)anthracene	25.52	278	11429	0.379	ng/ul	99
26) Benzo(g,h,i)perylene	26.16	276	12217	0.349	ng/ul	98

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

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