

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101419\
 Data File : BN008189.D
 Acq On : 16 Oct 2019 00:36
 Operator : JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Oct 16 03:43:57 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 : Tue Oct 15 15:13:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1427	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	7182	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	4914	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	10538	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	10128	0.40	ng/ul	-0.01
20) Perylene-d12	23.38	264	8353	0.40	ng/ul	-0.02

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	5152	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	14843	0.41	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.43	128	8796	0.430	ng/ul	100
5) 2-Methylnaphthalene	12.06	142	6106	0.435	ng/ul	99
7) Acenaphthylene	13.97	152	9418	0.402	ng/ul	100
8) Acenaphthene	14.31	153	7436	0.409	ng/ul	98
9) Fluorene	15.31	166	8250	0.399	ng/ul	97
11) Pentachlorophenol	16.67	266	826	0.324	ng/ul	100
12) Phenanthrene	17.04	178	13454	0.436	ng/ul	99
13) Anthracene	17.13	178	11995	0.442	ng/ul	100
15) Fluoranthene	19.06	202	17077	0.419	ng/ul	95
17) Pyrene	19.43	202	18127	0.450	ng/ul	100
18) Benzo(a)anthracene	21.18	228	15424	0.419	ng/ul	99
19) Chrysene	21.23	228	19003	0.421	ng/ul	100
21) Benzo(b)fluoranthene	22.73	252	14344	0.528	ng/ul	97
22) Benzo(k)fluoranthene	22.77	252	15743	0.511	ng/ul	99
23) Benzo(a)pyrene	23.28	252	13782	0.482	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.55	276	14114	0.419	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.56	278	11287	0.424	ng/ul	98
26) Benzo(g,h,i)perylene	26.20	276	12463	0.403	ng/ul	98

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

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