

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102819\
 Data File : BN008436.D
 Acq On : 28 Oct 2019 08:39
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Oct 29 11:21:05 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 : Fri Oct 25 13:14:06 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1452	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	7602	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.20	164	5120	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	10889	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	10120	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	9825	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	5030	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	14163	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	8591	0.397	ng/ul	99
5) 2-Methylnaphthalene	12.01	142	5997	0.404	ng/ul	99
7) Acenaphthylene	13.92	152	8871	0.364	ng/ul	100
8) Acenaphthene	14.27	153	7207	0.380	ng/ul	99
9) Fluorene	15.26	166	8058	0.374	ng/ul	98
11) Pentachlorophenol	16.63	266	786	0.298	ng/ul	99
12) Phenanthrene	17.00	178	13269	0.416	ng/ul	100
13) Anthracene	17.09	178	11981	0.427	ng/ul	99
15) Fluoranthene	19.02	202	16472	0.391	ng/ul	97
17) Pyrene	19.39	202	17011	0.422	ng/ul	98
18) Benzo(a)anthracene	21.15	228	14751	0.401	ng/ul	99
19) Chrysene	21.20	228	16966	0.376	ng/ul	99
21) Benzo(b)fluoranthene	22.69	252	14768	0.462	ng/ul	94
22) Benzo(k)fluoranthene	22.73	252	15294	0.422	ng/ul	97
23) Benzo(a)pyrene	23.24	252	14293	0.425	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.48	276	15235	0.384	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.49	278	12386	0.395	ng/ul	99
26) Benzo(g,h,i)perylene	26.13	276	13049	0.359	ng/ul	98

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

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