

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102419\
 Data File : BN008422.D
 Acq On : 25 Oct 2019 12:16
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Oct 25 13:00:13 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 12:06:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1527	0.40	ng/ul	0.00
2) Naphthalene-d8	10.34	136	7721	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	5223	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	10640	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	8652	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	7766	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	5171	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	13862	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	8739	0.398	ng/ul	99
5) 2-Methylnaphthalene	12.01	142	6079	0.403	ng/ul	100
7) Acenaphthylene	13.93	152	9493	0.382	ng/ul	100
8) Acenaphthene	14.27	153	7305	0.378	ng/ul	97
9) Fluorene	15.27	166	8046	0.366	ng/ul	98
11) Pentachlorophenol	16.64	266	765	0.297	ng/ul	98
12) Phenanthrene	17.00	178	12922	0.415	ng/ul	99
13) Anthracene	17.10	178	11916	0.435	ng/ul	100
15) Fluoranthene	19.03	202	16005	0.389	ng/ul	96
17) Pyrene	19.39	202	16467	0.478	ng/ul	96
18) Benzo(a)anthracene	21.15	228	12679	0.403	ng/ul	99
19) Chrysene	21.20	228	14657	0.380	ng/ul	100
21) Benzo(b)fluoranthene	22.69	252	11141	0.441	ng/ul	93
22) Benzo(k)fluoranthene	22.74	252	12098	0.422	ng/ul	97
23) Benzo(a)pyrene	23.25	252	10985	0.413	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.49	276	11870	0.379	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.50	278	9590	0.387	ng/ul#	95
26) Benzo(g,h,i)perylene	26.14	276	10359	0.361	ng/ul	96

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

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