

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102819\
 Data File : BN008439.D
 Acq On : 28 Oct 2019 13:58
 Operator : JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.420

Quant Time: Oct 29 11:21:20 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	1333	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	7583	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	5315	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	11054	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	8761	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	8169	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	5083	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	13552	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	8522	0.395	ng/ul	100
5) 2-Methylnaphthalene	12.01	142	6140	0.415	ng/ul	100
7) Acenaphthylene	13.92	152	9243	0.365	ng/ul	100
8) Acenaphthene	14.27	153	7439	0.378	ng/ul	99
9) Fluorene	15.26	166	8374	0.375	ng/ul	97
11) Pentachlorophenol	16.63	266	750	0.280	ng/ul	97
12) Phenanthrene	17.00	178	13466	0.416	ng/ul	100
13) Anthracene	17.09	178	12019	0.422	ng/ul	99
15) Fluoranthene	19.02	202	15732	0.368	ng/ul	95
17) Pyrene	19.39	202	16203	0.464	ng/ul	99
18) Benzo(a)anthracene	21.15	228	12524	0.393	ng/ul	99
19) Chrysene	21.20	228	14834	0.379	ng/ul	100
21) Benzo(b)fluoranthene	22.69	252	11992	0.451	ng/ul	89
22) Benzo(k)fluoranthene	22.73	252	13204	0.438	ng/ul	99
23) Benzo(a)pyrene	23.24	252	11922	0.426	ng/ul#	72
24) Indeno(1,2,3-cd)pyrene	25.49	276	12684	0.385	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.49	278	10242	0.393	ng/ul	99
26) Benzo(g,h,i)perylene	26.14	276	10995	0.364	ng/ul	97

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

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