

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN007004.D
 Acq On : 08 Aug 2019 01:45
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.428

Quant Time: Aug 08 02:41:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:30:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	6452	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	25142	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	16036	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	39859	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	31480	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	26292	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	17275	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	46863	0.36	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.28	128	26980	0.373	ng/ul	97
5) 2-Methylnaphthalene	11.91	142	20515	0.378	ng/ul	100
7) Acenaphthylene	13.82	152	34081	0.367	ng/ul	99
8) Acenaphthene	14.18	153	23656	0.372	ng/ul	99
9) Fluorene	15.17	166	28466	0.369	ng/ul	99
11) Pentachlorophenol	16.52	266	3288	0.299	ng/ul#	84
12) Phenanthrene	16.90	178	47465	0.374	ng/ul	100
13) Anthracene	16.99	178	46707	0.359	ng/ul	99
15) Fluoranthene	18.93	202	56164	0.356	ng/ul	100
17) Pyrene	19.29	202	57085	0.406	ng/ul	98
18) Benzo(a)anthracene	21.06	228	50829	0.368	ng/ul	99
19) Chrysene	21.11	228	47671	0.368	ng/ul	100
21) Benzo(b)fluoranthene	22.57	252	40368	0.379	ng/ul	99
22) Benzo(k)fluoranthene	22.62	252	39306	0.380	ng/ul	99
23) Benzo(a)pyrene	23.11	252	35958	0.358	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.30	276	37027	0.342	ng/ul	99
25) Dibenzo(a,h)anthracene	25.31	278	30239	0.350	ng/ul	99
26) Benzo(g,h,i)perylene	25.93	276	30950	0.341	ng/ul	98

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

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