

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082819\
 Data File : BN007437.D
 Acq On : 27 Aug 2019 17:59
 Operator : HP/JU
 Sample : SSTD0.428
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.428

Quant Time: Aug 27 18:38:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 18:27:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	3631	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	15027	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	7728	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	17613	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	15365	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	16142	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	9025	0.34	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	21985	0.39	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.21	128	16548	0.383	ng/ul	100
5) 2-Methylnaphthalene	11.84	142	11086	0.341	ng/ul	100
7) Acenaphthylene	13.76	152	16529	0.369	ng/ul	100
8) Acenaphthene	14.11	153	12021	0.392	ng/ul	100
9) Fluorene	15.11	166	12921	0.348	ng/ul	100
11) Pentachlorophenol	16.48	266	1329	0.274	ng/ul	100
12) Phenanthrene	16.84	178	21128	0.376	ng/ul	100
13) Anthracene	16.94	178	19862	0.346	ng/ul	100
15) Fluoranthene	18.87	202	28053	0.402	ng/ul	100
17) Pyrene	19.24	202	29176	0.425	ng/ul	100
18) Benzo(a)anthracene	21.01	228	26369	0.391	ng/ul	100
19) Chrysene	21.06	228	29010	0.459	ng/ul	100
21) Benzo(b)fluoranthene	22.51	252	26791	0.409	ng/ul	100
22) Benzo(k)fluoranthene	22.55	252	29438	0.464	ng/ul	100
23) Benzo(a)pyrene	23.04	252	27161	0.441	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.20	276	30219	0.455	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.22	278	24147	0.455	ng/ul	100
26) Benzo(g,h,i)perylene	25.82	276	26118	0.469	ng/ul	100

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

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