

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082819\
 Data File : BN007440.D
 Acq On : 27 Aug 2019 20:05
 Operator : HP/JU
 Sample : SSTD3.231
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.231

Quant Time: Aug 27 20:36:08 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	3883	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	16398	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	8242	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	18868	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	15846	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	16078	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	79066	2.71	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	185839	3.06	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.20	128	141235	2.993	ng/ul	99
5) 2-Methylnaphthalene	11.83	142	96617	2.727	ng/ul	99
7) Acenaphthylene	13.76	152	147319	3.084	ng/ul	99
8) Acenaphthene	14.11	153	102909	3.149	ng/ul	97
9) Fluorene	15.11	166	118162	2.983	ng/ul	98
11) Pentachlorophenol	16.46	266	15636	3.007	ng/ul	98
12) Phenanthrene	16.84	178	189867	3.158	ng/ul	99
13) Anthracene	16.93	178	189537	3.081	ng/ul	99
15) Fluoranthene	18.87	202	230899	3.089	ng/ul	98
17) Pyrene	19.24	202	231195	3.268	ng/ul	99
18) Benzo(a)anthracene	21.01	228	230380	3.310	ng/ul	99
19) Chrysene	21.06	228	222720	3.417	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	225988	3.467	ng/ul	95
22) Benzo(k)fluoranthene	22.55	252	224543	3.554	ng/ul#	93
23) Benzo(a)pyrene	23.04	252	216732	3.531	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.19	276	246818	3.731	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.21	278	196543	3.720	ng/ul	96
26) Benzo(g,h,i)perylene	25.82	276	208174	3.751	ng/ul	98

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

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