

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080919\
 Data File : BN007039.D
 Acq On : 09 Aug 2019 12:26
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.431

Quant Time: Aug 09 14:22:27 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	8102	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	28551	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.11	164	16328	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	38064	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	30198	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	32928	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	18478	0.36	ng/ul	-0.01
14) Fluoranthene-d10	18.90	212	41997	0.34	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.27	128	30605	0.372	ng/ul	97
5) 2-Methylnaphthalene	11.90	142	22082	0.358	ng/ul	100
7) Acenaphthylene	13.82	152	33311	0.352	ng/ul	99
8) Acenaphthene	14.17	153	23920	0.369	ng/ul	99
9) Fluorene	15.17	166	27749	0.354	ng/ul	100
11) Pentachlorophenol	16.52	266	3714	0.354	ng/ul#	82
12) Phenanthrene	16.90	178	45669	0.376	ng/ul	100
13) Anthracene	16.99	178	43038	0.347	ng/ul	99
15) Fluoranthene	18.93	202	51613	0.342	ng/ul	98
17) Pyrene	19.29	202	51585	0.383	ng/ul	99
18) Benzo(a)anthracene	21.07	228	47433	0.358	ng/ul	99
19) Chrysene	21.12	228	46622	0.375	ng/ul	98
21) Benzo(b)fluoranthene	22.58	252	45038	0.337	ng/ul	99
22) Benzo(k)fluoranthene	22.62	252	45564	0.352	ng/ul	98
23) Benzo(a)pyrene	23.11	252	43306	0.345	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.31	276	51732	0.382	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.33	278	42558	0.393	ng/ul	98
26) Benzo(g,h,i)perylene	25.94	276	43644	0.384	ng/ul	97

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

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