

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080719\
 Data File : BM021900.D
 Acq On : 07 Aug 2019 17:40
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.440

Quant Time: Aug 07 18:36:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 08:28:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	757	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	3370	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	1925	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	4299	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	4203	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4857	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2220	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	5421	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	3642	0.383	ng/ul	98
5) 2-Methylnaphthalene	12.10	142	2532	0.399	ng/ul	100
7) Acenaphthylene	14.01	152	3537	0.394	ng/ul	98
8) Acenaphthene	14.35	153	2830	0.385	ng/ul	95
9) Fluorene	15.35	166	3096	0.379	ng/ul	100
11) Pentachlorophenol	16.71	266	349	0.440	ng/ul	89
12) Phenanthrene	17.09	178	4834	0.386	ng/ul	99
13) Anthracene	17.19	178	4436	0.416	ng/ul	98
15) Fluoranthene	19.11	202	6295	0.376	ng/ul	99
17) Pyrene	19.47	202	6673	0.386	ng/ul	96
18) Benzo(a)anthracene	21.23	228	6086	0.405	ng/ul	98
19) Chrysene	21.28	228	7071	0.366	ng/ul	97
21) Benzo(b)fluoranthene	22.79	252	6502	0.397	ng/ul	98
22) Benzo(k)fluoranthene	22.84	252	6823	0.368	ng/ul	99
23) Benzo(a)pyrene	23.36	252	6291	0.378	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.68	276	7747	0.390	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.69	278	6250	0.389	ng/ul	97
26) Benzo(g,h,i)perylene	26.36	276	6614	0.369	ng/ul	96

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

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