

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080719\
 Data File : BM021951.D
 Acq On : 09 Aug 2019 02:01
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.444

Quant Time: Aug 09 02:42:56 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 : Fri Aug 09 01:29:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	794	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	3679	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	1981	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	4317	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	3896	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4407	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2262	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	4962	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	3752	0.361	ng/ul	98
5) 2-Methylnaphthalene	12.09	142	2561	0.370	ng/ul	100
7) Acenaphthylene	14.01	152	3156	0.342	ng/ul	98
8) Acenaphthene	14.34	153	2850	0.377	ng/ul	97
9) Fluorene	15.34	166	3022	0.360	ng/ul	99
11) Pentachlorophenol	16.71	266	282	0.354	ng/ul	94
12) Phenanthrene	17.08	178	4572	0.364	ng/ul	99
13) Anthracene	17.18	178	4043	0.377	ng/ul	97
15) Fluoranthene	19.11	202	5525	0.328	ng/ul	97
17) Pyrene	19.47	202	5813	0.363	ng/ul#	95
18) Benzo(a)anthracene	21.22	228	5023	0.360	ng/ul	99
19) Chrysene	21.27	228	6183	0.345	ng/ul	100
21) Benzo(b)fluoranthene	22.79	252	5582	0.376	ng/ul	94
22) Benzo(k)fluoranthene	22.83	252	6039	0.359	ng/ul	96
23) Benzo(a)pyrene	23.36	252	5412	0.358	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.67	276	6535	0.362	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.68	278	5244	0.360	ng/ul	96
26) Benzo(g,h,i)perylene	26.35	276	5721	0.352	ng/ul	96

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

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