

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080619\
 Data File : BM021882.D
 Acq On : 07 Aug 2019 05:37
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.437

Quant Time: Aug 07 06:09:57 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 04:02:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	651	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	3016	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	1694	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	3667	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	3350	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	3912	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1937	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	4398	0.36	ng/ul	0.00

Target Compounds						Ovalue
3) Naphthalene	10.47	128	3234	0.380	ng/ul	98
5) 2-Methylnaphthalene	12.11	142	2223	0.392	ng/ul	100
7) Acenaphthylene	14.01	152	3033	0.384	ng/ul	100
8) Acenaphthene	14.35	153	2517	0.389	ng/ul	97
9) Fluorene	15.35	166	2642	0.368	ng/ul	98
11) Pentachlorophenol	16.72	266	242	0.358	ng/ul	91
12) Phenanthrene	17.09	178	4139	0.388	ng/ul	99
13) Anthracene	17.19	178	3511	0.386	ng/ul	99
15) Fluoranthene	19.12	202	5068	0.355	ng/ul	97
17) Pyrene	19.48	202	5310	0.386	ng/ul#	94
18) Benzo(a)anthracene	21.23	228	4560	0.381	ng/ul	98
19) Chrysene	21.28	228	5864	0.381	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	5087	0.386	ng/ul	96
22) Benzo(k)fluoranthene	22.84	252	5669	0.380	ng/ul	97
23) Benzo(a)pyrene	23.37	252	5018	0.374	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.68	276	6011	0.375	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.69	278	4859	0.375	ng/ul	98
26) Benzo(g,h,i)perylene	26.36	276	5278	0.366	ng/ul	97

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

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