

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080319\
 Data File : BM021786.D
 Acq On : 03 Aug 2019 22:58
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.426

Quant Time: Aug 04 04:11:07 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 : Sat Aug 03 22:53:27 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	643	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	2828	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	1558	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	3415	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	3220	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	3769	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	1884	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	4457	0.39	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.50	128	3288	0.412	ng/ul	99
5) 2-Methylnaphthalene	12.13	142	2153	0.405	ng/ul	100
7) Acenaphthylene	14.03	152	2925	0.403	ng/ul	99
8) Acenaphthene	14.37	153	2455	0.413	ng/ul	96
9) Fluorene	15.37	166	2591	0.392	ng/ul	97
11) Pentachlorophenol	16.75	266	242	0.384	ng/ul#	77
12) Phenanthrene	17.11	178	4010	0.404	ng/ul	100
13) Anthracene	17.21	178	3361	0.397	ng/ul	97
15) Fluoranthene	19.13	202	5114	0.384	ng/ul	97
17) Pyrene	19.49	202	5499	0.416	ng/ul#	93
18) Benzo(a)anthracene	21.25	228	4304	0.374	ng/ul	99
19) Chrysene	21.29	228	6168	0.416	ng/ul	99
21) Benzo(b)fluoranthene	22.81	252	4956	0.390	ng/ul	98
22) Benzo(k)fluoranthene	22.86	252	5868	0.408	ng/ul	98
23) Benzo(a)pyrene	23.39	252	4989	0.386	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.71	276	6023	0.390	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.73	278	4814	0.386	ng/ul	94
26) Benzo(g,h,i)perylene	26.39	276	5630	0.405	ng/ul	96

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

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