

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080619\
 Data File : BM021867.D
 Acq On : 06 Aug 2019 20:31
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.436

Quant Time: Aug 06 21:03:03 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 : Tue Aug 06 01:29:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	673	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	3242	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	1798	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	4000	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3716	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4313	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	2189	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	5082	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	3632	0.397	ng/ul	98
5) 2-Methylnaphthalene	12.10	142	2551	0.418	ng/ul	99
7) Acenaphthylene	14.01	152	3398	0.405	ng/ul	99
8) Acenaphthene	14.35	153	2872	0.419	ng/ul	99
9) Fluorene	15.35	166	3049	0.400	ng/ul	98
11) Pentachlorophenol	16.72	266	303	0.411	ng/ul	93
12) Phenanthrene	17.09	178	4731	0.406	ng/ul	99
13) Anthracene	17.19	178	4190	0.422	ng/ul	98
15) Fluoranthene	19.11	202	5803	0.372	ng/ul	98
17) Pyrene	19.48	202	6120	0.401	ng/ul	96
18) Benzo(a)anthracene	21.23	228	5376	0.404	ng/ul	99
19) Chrysene	21.28	228	6645	0.389	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	5981	0.411	ng/ul	95
22) Benzo(k)fluoranthene	22.84	252	6522	0.396	ng/ul	97
23) Benzo(a)pyrene	23.37	252	5846	0.395	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.69	276	7050	0.399	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.70	278	5702	0.399	ng/ul	97
26) Benzo(g,h,i)perylene	26.36	276	6241	0.393	ng/ul	98

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

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