

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080619\
 Data File : BM021858.D
 Acq On : 06 Aug 2019 08:36
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.434

Quant Time: Aug 06 11:12:25 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	649	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	2907	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	1675	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	3771	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	3692	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4417	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1972	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	4985	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	3271	0.398	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	2302	0.421	ng/ul	99
7) Acenaphthylene	14.01	152	3208	0.411	ng/ul	98
8) Acenaphthene	14.35	153	2610	0.408	ng/ul	98
9) Fluorene	15.35	166	2794	0.394	ng/ul	99
11) Pentachlorophenol	16.72	266	281	0.404	ng/ul	92
12) Phenanthrene	17.09	178	4487	0.409	ng/ul	99
13) Anthracene	17.19	178	4039	0.432	ng/ul	98
15) Fluoranthene	19.11	202	5724	0.389	ng/ul	100
17) Pyrene	19.48	202	6065	0.400	ng/ul	97
18) Benzo(a)anthracene	21.23	228	5656	0.428	ng/ul	99
19) Chrysene	21.28	228	6574	0.387	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	6113	0.410	ng/ul	97
22) Benzo(k)fluoranthene	22.84	252	6685	0.397	ng/ul	98
23) Benzo(a)pyrene	23.37	252	6103	0.403	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.68	276	7366	0.407	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.69	278	5989	0.410	ng/ul	97
26) Benzo(g,h,i)perylene	26.36	276	6415	0.394	ng/ul	99

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

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