

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027871.D
 Acq On : 21 Nov 2020 02:27
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.434

Quant Time: Nov 21 03:27:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 21 02:40:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	909	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	3622	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2215	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4560	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2956	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2209	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	2402	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	4687	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	4342	0.400	ng/ul	98
5) 2-Methylnaphthalene	12.80	142	3095	0.419	ng/ul	99
7) Acenaphthylene	14.67	152	4412	0.408	ng/ul	99
8) Acenaphthene	15.01	153	3205	0.396	ng/ul	99
9) Fluorene	15.98	166	3727	0.397	ng/ul	98
11) Pentachlorophenol	17.31	266	586	0.377	ng/ul	98
12) Phenanthrene	17.70	178	5707	0.396	ng/ul	99
13) Anthracene	17.79	178	5470	0.394	ng/ul	99
15) Fluoranthene	19.68	202	6216	0.389	ng/ul	99
17) Pyrene	20.04	202	6194	0.413	ng/ul	99
18) Benzo(a)anthracene	21.74	228	4749	0.398	ng/ul	99
19) Chrysene	21.80	228	4652	0.400	ng/ul	99
21) Benzo(b)fluoranthene	23.44	252	3746	0.404	ng/ul	98
22) Benzo(k)fluoranthene	23.49	252	3679	0.405	ng/ul	99
23) Benzo(a)pyrene	24.08	252	3224	0.397	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.68	276	3594	0.409	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.69	278	3008	0.411	ng/ul	98
26) Benzo(g,h,i)perylene	27.45	276	3052	0.414	ng/ul	99

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

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