

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111720\
 Data File : BM027827.D
 Acq On : 17 Nov 2020 20:52
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.423

Quant Time: Nov 18 01:18:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM111020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 17 09:19:02 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1267	0.40	ng/ul	0.00
2) Naphthalene-d8	11.22	136	4871	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	2532	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	4129	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	2355	0.40	ng/ul	0.00
20) Perylene-d12	24.23	264	2335	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.78	152	3218	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	3783	0.31	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.26	128	6019	0.387	ng/ul	97
5) 2-Methylnaphthalene	12.85	142	3911	0.361	ng/ul	99
7) Acenaphthylene	14.71	152	5406	0.434	ng/ul	98
8) Acenaphthene	15.05	153	3819	0.400	ng/ul	97
9) Fluorene	16.02	166	4094	0.357	ng/ul	98
11) Pentachlorophenol	17.35	266	512	0.344	ng/ul	99
12) Phenanthrene	17.74	178	5332	0.397	ng/ul	99
13) Anthracene	17.83	178	5271	0.403	ng/ul	98
15) Fluoranthene	19.71	202	4957	0.322	ng/ul	93
17) Pyrene	20.07	202	5020	0.481	ng/ul#	90
18) Benzo(a)anthracene	21.78	228	3979	0.423	ng/ul#	83
19) Chrysene	21.84	228	3843	0.401	ng/ul#	84
21) Benzo(b)fluoranthene	23.49	252	3705	0.347	ng/ul#	70
22) Benzo(k)fluoranthene	23.54	252	3657	0.349	ng/ul#	67
23) Benzo(a)pyrene	24.13	252	3438	0.366	ng/ul#	65
24) Indeno(1,2,3-cd)pyrene	26.75	276	3981	0.372	ng/ul#	88
25) Dibenzo(a,h)anthracene	26.77	278	3343	0.374	ng/ul#	66
26) Benzo(g,h,i)perylene	27.53	276	3422	0.378	ng/ul#	79

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

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