

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111720\
 Data File : BM027810.D
 Acq On : 17 Nov 2020 08:37
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDD0.421

Quant Time: Nov 17 09:19:57 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	636	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	2308	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.98	164	1495	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.69	188	3366	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	2892	0.40	ng/ul	0.00
20) Perylene-d12	24.23	264	2521	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.77	152	1646	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	3945	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.26	128	2771	0.376	ng/ul	99
5) 2-Methylnaphthalene	12.84	142	1927	0.376	ng/ul	100
7) Acenaphthylene	14.71	152	2759	0.375	ng/ul	99
8) Acenaphthene	15.04	153	2156	0.382	ng/ul	98
9) Fluorene	16.01	166	2605	0.384	ng/ul	97
11) Pentachlorophenol	17.34	266	497	0.409	ng/ul	97
12) Phenanthrene	17.74	178	4278	0.390	ng/ul	99
13) Anthracene	17.83	178	4069	0.381	ng/ul	100
15) Fluoranthene	19.71	202	5035	0.401	ng/ul	99
17) Pyrene	20.07	202	5045	0.393	ng/ul	99
18) Benzo(a)anthracene	21.78	228	4600	0.398	ng/ul	99
19) Chrysene	21.83	228	4727	0.402	ng/ul	97
21) Benzo(b)fluoranthene	23.49	252	4575	0.397	ng/ul	98
22) Benzo(k)fluoranthene	23.53	252	4437	0.392	ng/ul	97
23) Benzo(a)pyrene	24.12	252	3907	0.385	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.74	276	4623	0.400	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.76	278	3803	0.394	ng/ul#	95
26) Benzo(g,h,i)perylene	27.52	276	3988	0.408	ng/ul	98

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

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