

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111020\
 Data File : BM027768.D
 Acq On : 10 Nov 2020 13:22
 Operator : JU/CG
 Sample : SSTD0.480
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.480

Quant Time: Nov 10 14:34:51 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 10 14:33:58 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	538	0.40	ng/ul	0.00
2) Naphthalene-d8	11.22	136	1880	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	1195	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	2633	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	2260	0.40	ng/ul	0.00
20) Perylene-d12	24.25	264	1913	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	1784	0.51	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	3966	0.51	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.27	128	2958	0.492	ng/ul	100
5) 2-Methylnaphthalene	12.86	142	2063	0.498	ng/ul	100
7) Acenaphthylene	14.72	152	2900	0.488	ng/ul	100
8) Acenaphthene	15.06	153	2246	0.498	ng/ul	100
9) Fluorene	16.03	166	2715	0.510	ng/ul	100
11) Pentachlorophenol	17.35	266	468	0.437	ng/ul	100
12) Phenanthrene	17.75	178	4261	0.499	ng/ul	100
13) Anthracene	17.84	178	4127	0.499	ng/ul	100
15) Fluoranthene	19.72	202	4966	0.510	ng/ul	100
17) Pyrene	20.08	202	4859	0.492	ng/ul	100
18) Benzo(a)anthracene	21.79	228	4448	0.497	ng/ul	100
19) Chrysene	21.85	228	4565	0.499	ng/ul	100
21) Benzo(b)fluoranthene	23.50	252	4281	0.509	ng/ul	100
22) Benzo(k)fluoranthene	23.55	252	4240	0.506	ng/ul	100
23) Benzo(a)pyrene	24.14	252	3769	0.508	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.77	276	4310	0.491	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.78	278	3632	0.497	ng/ul	100
26) Benzo(g,h,i)perylene	27.55	276	3695	0.488	ng/ul	100

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

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