

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM111020\
 Data File : BM027770.D
 Acq On : 10 Nov 2020 14:35
 Operator : JU/CG
 Sample : SSTD1.676
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.676

Quant Time: Nov 10 15:09:33 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	518	0.40	ng/ul	0.00
2) Naphthalene-d8	11.22	136	1876	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	1217	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	2622	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	2310	0.40	ng/ul	0.00
20) Perylene-d12	24.25	264	1950	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	5402	1.56	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	12421	1.59	ng/ul	0.00

Target Compounds					Ovalue	
3) Naphthalene	11.27	128	8892	1.483	ng/ul	96
5) 2-Methylnaphthalene	12.86	142	6332	1.531	ng/ul	100
7) Acenaphthylene	14.72	152	9159	1.513	ng/ul	96
8) Acenaphthene	15.06	153	6953	1.515	ng/ul	99
9) Fluorene	16.03	166	8484	1.565	ng/ul	99
11) Pentachlorophenol	17.35	266	1463	1.373	ng/ul	98
12) Phenanthrene	17.75	178	13100	1.540	ng/ul	97
13) Anthracene	17.84	178	13090	1.591	ng/ul	98
15) Fluoranthene	19.72	202	15524	1.602	ng/ul	98
17) Pyrene	20.08	202	15448	1.532	ng/ul	99
18) Benzo(a)anthracene	21.79	228	14054	1.537	ng/ul	98
19) Chrysene	21.85	228	14235	1.523	ng/ul	98
21) Benzo(b)fluoranthene	23.51	252	13696	1.597	ng/ul	94
22) Benzo(k)fluoranthene	23.55	252	13681	1.603	ng/ul#	95
23) Benzo(a)pyrene	24.14	252	12113	1.602	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	26.77	276	14006	1.564	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.78	278	11672	1.566	ng/ul	95
26) Benzo(g,h,i)perylene	27.55	276	11835	1.533	ng/ul	96

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

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