

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110920\
 Data File : BM027753.D
 Acq On : 09 Nov 2020 12:28
 Operator : JU/CG
 Sample : SSTD1.681
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.681

Quant Time: Nov 10 06:25:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM02.4-EPA-SIM-BM110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 06:19:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.40	152	355	0.40	ng/ul	0.00
2) Naphthalene-d8	11.23	136	1201	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	781	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.72	188	1679	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	1527	0.40	ng/ul	0.00
20) Perylene-d12	24.26	264	1382	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	3454	1.58	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	8016	1.74	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.28	128	5790	1.648	ng/ul#	76
5) 2-Methylnaphthalene	12.87	142	4062	1.588	ng/ul	98
7) Acenaphthylene	14.73	152	5884	1.476	ng/ul#	78
8) Acenaphthene	15.07	153	4403	1.551	ng/ul	93
9) Fluorene	16.04	166	5313	1.512	ng/ul#	95
11) Pentachlorophenol	17.36	266	1059	2.317	ng/ul	91
12) Phenanthrene	17.76	178	8294	1.601	ng/ul#	85
13) Anthracene	17.85	178	8288	1.635	ng/ul#	84
15) Fluoranthene	19.73	202	9895	1.588	ng/ul	81
17) Pyrene	20.09	202	9705	1.393	ng/ul#	82
18) Benzo(a)anthracene	21.80	228	9100	1.520	ng/ul#	87
19) Chrysene	21.85	228	9217	1.486	ng/ul#	88
21) Benzo(b)fluoranthene	23.52	252	9081	1.631	ng/ul	69
22) Benzo(k)fluoranthene	23.56	252	9140	1.579	ng/ul#	69
23) Benzo(a)pyrene	24.15	252	8201	1.645	ng/ul#	65
24) Indeno(1,2,3-cd)pyrene	26.79	276	9959	1.551	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.81	278	8407	1.587	ng/ul#	62
26) Benzo(g,h,i)perylene	27.57	276	8567	1.581	ng/ul#	68

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

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