

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040521\
 Data File : BM029341.D
 Acq On : 05 Apr 2021 13:59
 Operator : CG/JU
 Sample : SSTD0.879
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.879

Quant Time: Apr 05 14:30:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 05 14:08:43 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	2041	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	6640	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	3404	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	7068	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	7882	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	8279	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	7064	0.75	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	14140	0.77	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.53	128	14005	0.847	ng/ul	99
5) 2-Methylnaphthalene	12.15	142	9066	0.819	ng/ul	100
7) Acenaphthylene	14.05	152	11322	0.815	ng/ul	99
8) Acenaphthene	14.39	153	9565	0.868	ng/ul	100
9) Fluorene	15.38	166	10488	0.858	ng/ul	99
11) Pentachlorophenol	16.73	266	1238	0.670	ng/ul	96
12) Phenanthrene	17.11	178	16362	0.828	ng/ul	100
13) Anthracene	17.20	178	14625	0.776	ng/ul	99
15) Fluoranthene	19.13	202	17795	0.791	ng/ul	98
17) Pyrene	19.49	202	18190	0.635	ng/ul	98
18) Benzo(a)anthracene	21.23	228	16145	0.600	ng/ul	99
19) Chrysene	21.28	228	17467	0.653	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	18736	0.661	ng/ul	95
22) Benzo(k)fluoranthene	22.84	252	19299	0.682	ng/ul	95
23) Benzo(a)pyrene	23.36	252	16313	0.646	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.68	276	23170	0.705	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.69	278	19608	0.711	ng/ul	96
26) Benzo(g,h,i)perylene	26.36	276	19615	0.717	ng/ul	98

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

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