

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040220\
 Data File : BM025890.D
 Acq On : 01 Apr 2020 16:45
 Operator : CG/JU
 Sample : SSTD1.679
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.679

Quant Time: Apr 01 17:19:01 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 16:04:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2047	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	8200	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	5658	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	13417	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	15067	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	15698	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	25224	1.80	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	73561	1.89	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.38	128	40824	1.675	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	29430	1.795	ng/ul	99
7) Acenaphthylene	13.92	152	42179	1.757	ng/ul	98
8) Acenaphthene	14.27	153	34360	1.713	ng/ul	99
9) Fluorene	15.26	166	43204	1.820	ng/ul	99
11) Pentachlorophenol	16.61	266	5987	2.156	ng/ul	96
12) Phenanthrene	16.99	178	73128	1.738	ng/ul	98
13) Anthracene	17.08	178	67639	1.922	ng/ul	98
15) Fluoranthene	19.01	202	93373	1.914	ng/ul	98
17) Pyrene	19.37	202	91694	1.577	ng/ul	99
18) Benzo(a)anthracene	21.13	228	94723	1.912	ng/ul	99
19) Chrysene	21.18	228	98679	1.651	ng/ul	100
21) Benzo(b)fluoranthene	22.65	252	102016	1.679	ng/ul	97
22) Benzo(k)fluoranthene	22.69	252	106445	1.711	ng/ul	97
23) Benzo(a)pyrene	23.20	252	90365	1.808	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.41	276	123561	1.944	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.42	278	103510	2.105	ng/ul	98
26) Benzo(g,h,i)perylene	26.05	276	102845	1.663	ng/ul	99

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

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