

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM033020\
 Data File : BM025857.D
 Acq On : 30 Mar 2020 13:12
 Operator : CG/JU
 Sample : SSTD1.678
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.678

Quant Time: Mar 30 13:47:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM033020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 30 12:31:41 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1450	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	5448	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	3375	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	7600	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	8042	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	7641	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	15987	1.57	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	39291	1.59	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.38	128	27260	1.663	ng/ul	97
5) 2-Methylnaphthalene	12.00	142	18633	1.560	ng/ul	99
7) Acenaphthylene	13.92	152	24343	1.540	ng/ul	97
8) Acenaphthene	14.27	153	20291	1.655	ng/ul	99
9) Fluorene	15.26	166	24551	1.598	ng/ul	99
11) Pentachlorophenol	16.61	266	2866	1.441	ng/ul	97
12) Phenanthrene	16.99	178	41099	1.677	ng/ul	99
13) Anthracene	17.08	178	36378	1.673	ng/ul	98
15) Fluoranthene	19.01	202	50360	1.642	ng/ul	99
17) Pyrene	19.38	202	48928	1.604	ng/ul	100
18) Benzo(a)anthracene	21.13	228	45937	1.554	ng/ul	99
19) Chrysene	21.18	228	51583	1.588	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	50495	1.652	ng/ul	95
22) Benzo(k)fluoranthene	22.70	252	52108	1.625	ng/ul#	94
23) Benzo(a)pyrene	23.20	252	41578	1.620	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.41	276	54316	1.698	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.42	278	43095	1.626	ng/ul	92
26) Benzo(g,h,i)perylene	26.05	276	50351	1.853	ng/ul	97

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

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