

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025218.D
 Acq On : 03 Mar 2020 18:47
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.483

Quant Time: Mar 03 19:18:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 03 14:30:08 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	2880	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	10529	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6626	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15119	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	14004	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	13622	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	6700	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	16379	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	11491	0.363	ng/ul#	83
5) 2-Methylnaphthalene	12.12	142	7843	0.340	ng/ul	100
7) Acenaphthylene	14.02	152	11115	0.358	ng/ul	94
8) Acenaphthene	14.37	153	8682	0.361	ng/ul	99
9) Fluorene	15.36	166	10309	0.342	ng/ul	95
11) Pentachlorophenol	16.71	266	1262	0.319	ng/ul	98
12) Phenanthrene	17.09	178	17239	0.354	ng/ul	97
13) Anthracene	17.18	178	15153	0.350	ng/ul	96
15) Fluoranthene	19.11	202	20217	0.331	ng/ul	95
17) Pyrene	19.47	202	19927	0.375	ng/ul	96
18) Benzo(a)anthracene	21.22	228	18284	0.355	ng/ul	97
19) Chrysene	21.27	228	20683	0.366	ng/ul	97
21) Benzo(b)fluoranthene	22.77	252	19917	0.365	ng/ul	94
22) Benzo(k)fluoranthene	22.82	252	20050	0.351	ng/ul#	96
23) Benzo(a)pyrene	23.33	252	16641	0.364	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.61	276	22223	0.390	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	18642	0.394	ng/ul#	94
26) Benzo(g,h,i)perylene	26.27	276	19253	0.397	ng/ul#	90

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

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