

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025244.D
 Acq On : 04 Mar 2020 11:04
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.485

Quant Time: Mar 04 11:43:50 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 : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	3891	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	15647	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	11133	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	27432	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	23491	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	20598	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	10578	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	30100	0.34	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.50	128	17174	0.365	ng/ul#	83
5) 2-Methylnaphthalene	12.12	142	12367	0.361	ng/ul	99
7) Acenaphthylene	14.02	152	18726	0.359	ng/ul	93
8) Acenaphthene	14.37	153	14636	0.362	ng/ul	99
9) Fluorene	15.35	166	18249	0.360	ng/ul	96
11) Pentachlorophenol	16.70	266	2716	0.378	ng/ul	100
12) Phenanthrene	17.09	178	31672	0.358	ng/ul	96
13) Anthracene	17.18	178	28257	0.360	ng/ul	95
15) Fluoranthene	19.10	202	37463	0.338	ng/ul	96
17) Pyrene	19.47	202	36576	0.411	ng/ul	96
18) Benzo(a)anthracene	21.22	228	32046	0.371	ng/ul	97
19) Chrysene	21.27	228	34662	0.365	ng/ul	97
21) Benzo(b)fluoranthene	22.77	252	29711	0.361	ng/ul	93
22) Benzo(k)fluoranthene	22.81	252	32106	0.371	ng/ul#	96
23) Benzo(a)pyrene	23.33	252	25435	0.368	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.61	276	31446	0.365	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	26326	0.368	ng/ul#	94
26) Benzo(g,h,i)perylene	26.27	276	25599	0.350	ng/ul#	91

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

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