

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025157.D
 Acq On : 29 Feb 2020 06:08
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.496

Quant Time: Mar 02 00:40:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 02 00:36:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2461	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	9336	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	5640	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	13059	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	14788	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	16478	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	6720	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	18191	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	11505	0.405	ng/ul#	83
5) 2-Methylnaphthalene	12.12	142	7866	0.384	ng/ul	99
7) Acenaphthylene	14.03	152	11714	0.360	ng/ul	94
8) Acenaphthene	14.37	153	8606	0.394	ng/ul	98
9) Fluorene	15.36	166	10310	0.376	ng/ul	96
11) Pentachlorophenol	16.71	266	1613	0.404	ng/ul	98
12) Phenanthrene	17.09	178	17906	0.390	ng/ul	97
13) Anthracene	17.18	178	16165	0.363	ng/ul	96
15) Fluoranthene	19.11	202	24206	0.363	ng/ul	95
17) Pyrene	19.47	202	25009	0.375	ng/ul	96
18) Benzo(a)anthracene	21.22	228	23814	0.373	ng/ul	98
19) Chrysene	21.27	228	25271	0.397	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	26049	0.403	ng/ul	93
22) Benzo(k)fluoranthene	22.82	252	27223	0.425	ng/ul	100
23) Benzo(a)pyrene	23.34	252	23031	0.391	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.62	276	31559	0.427	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.64	278	26335	0.424	ng/ul#	94
26) Benzo(g,h,i)perylene	26.28	276	25772	0.414	ng/ul#	92

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

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