

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025135.D
 Acq On : 28 Feb 2020 16:20
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.494

Quant Time: Feb 28 16:53:37 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 : Fri Feb 28 12:40:55 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2468	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	8915	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.31	164	5272	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.06	188	12039	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	12898	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	14476	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	6321	0.38	ng/ul	-0.01
14) Fluoranthene-d10	19.08	212	15710	0.38	ng/ul	-0.01
Target Compounds					Ovalue	
3) Naphthalene	10.51	128	10989	0.405	ng/ul#	84
5) 2-Methylnaphthalene	12.12	142	7433	0.380	ng/ul	99
7) Acenaphthylene	14.03	152	10614	0.349	ng/ul	94
8) Acenaphthene	14.38	153	8092	0.396	ng/ul	99
9) Fluorene	15.37	166	9528	0.372	ng/ul	97
11) Pentachlorophenol	16.71	266	1298	0.353	ng/ul	96
12) Phenanthrene	17.10	178	16561	0.391	ng/ul	96
13) Anthracene	17.19	178	14239	0.347	ng/ul	97
15) Fluoranthene	19.12	202	20760	0.338	ng/ul	97
17) Pyrene	19.48	202	20786	0.358	ng/ul	96
18) Benzo(a)anthracene	21.23	228	19946	0.358	ng/ul	98
19) Chrysene	21.28	228	21791	0.393	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	22299	0.393	ng/ul	94
22) Benzo(k)fluoranthene	22.83	252	24563	0.437	ng/ul	98
23) Benzo(a)pyrene	23.35	252	19909	0.384	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.64	276	28421	0.438	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.65	278	23697	0.434	ng/ul	97
26) Benzo(g,h,i)perylene	26.29	276	23248	0.425	ng/ul#	91

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

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