

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022520\
 Data File : BM025076.D
 Acq On : 26 Feb 2020 11:42
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.486

Quant Time: Feb 26 12:43:37 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 : Tue Feb 25 11:03:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2354	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	8985	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	5513	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	12850	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	13958	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	15561	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	6487	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	17398	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.51	128	11052	0.404	ng/ul#	83
5) 2-Methylnaphthalene	12.13	142	7583	0.384	ng/ul	99
7) Acenaphthylene	14.04	152	11936	0.376	ng/ul	95
8) Acenaphthene	14.38	153	8466	0.396	ng/ul	99
9) Fluorene	15.37	166	10177	0.380	ng/ul	97
11) Pentachlorophenol	16.72	266	1454	0.370	ng/ul	98
12) Phenanthrene	17.10	178	17642	0.390	ng/ul	96
13) Anthracene	17.19	178	16345	0.373	ng/ul	96
15) Fluoranthene	19.12	202	22224	0.339	ng/ul	97
17) Pyrene	19.48	202	22359	0.356	ng/ul	97
18) Benzo(a)anthracene	21.23	228	22974	0.381	ng/ul	98
19) Chrysene	21.28	228	23676	0.394	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	24469	0.401	ng/ul	93
22) Benzo(k)fluoranthene	22.83	252	25247	0.418	ng/ul#	95
23) Benzo(a)pyrene	23.35	252	21948	0.394	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.64	276	28265	0.405	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.66	278	23723	0.405	ng/ul	96
26) Benzo(g,h,i)perylene	26.30	276	24137	0.410	ng/ul#	93

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

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