

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022720\
 Data File : BM025104.D
 Acq On : 27 Feb 2020 09:34
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.490

Quant Time: Feb 27 16:39:05 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	2583	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	9719	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.32	164	6089	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	14460	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	14642	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	15467	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	7100	0.39	ng/ul	-0.01
14) Fluoranthene-d10	19.09	212	19145	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.51	128	12015	0.406	ng/ul#	84
5) 2-Methylnaphthalene	12.13	142	8289	0.388	ng/ul	99
7) Acenaphthylene	14.03	152	12614	0.359	ng/ul	95
8) Acenaphthene	14.38	153	9141	0.388	ng/ul	98
9) Fluorene	15.37	166	11177	0.378	ng/ul	98
11) Pentachlorophenol	16.71	266	1481	0.335	ng/ul	99
12) Phenanthrene	17.10	178	19570	0.385	ng/ul	95
13) Anthracene	17.19	178	17651	0.358	ng/ul	95
15) Fluoranthene	19.12	202	24212	0.328	ng/ul	95
17) Pyrene	19.48	202	24334	0.369	ng/ul	97
18) Benzo(a)anthracene	21.23	228	23586	0.373	ng/ul	98
19) Chrysene	21.28	228	24782	0.394	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	24532	0.405	ng/ul	94
22) Benzo(k)fluoranthene	22.83	252	25636	0.427	ng/ul	98
23) Benzo(a)pyrene	23.34	252	21613	0.390	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.63	276	28111	0.405	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.65	278	23667	0.406	ng/ul	96
26) Benzo(g,h,i)perylene	26.29	276	23567	0.403	ng/ul#	93

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

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