

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022720\
 Data File : BM025108.D
 Acq On : 27 Feb 2020 18:03
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.491

Quant Time: Feb 28 01:07:21 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	2377	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	9049	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.32	164	5598	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	12760	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	13756	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	15727	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	6586	0.39	ng/ul	-0.01
14) Fluoranthene-d10	19.09	212	17129	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.51	128	11083	0.402	ng/ul#	84
5) 2-Methylnaphthalene	12.13	142	7662	0.385	ng/ul	99
7) Acenaphthylene	14.03	152	11391	0.353	ng/ul	96
8) Acenaphthene	14.38	153	8394	0.387	ng/ul	97
9) Fluorene	15.37	166	10065	0.370	ng/ul	97
11) Pentachlorophenol	16.71	266	1394	0.358	ng/ul	95
12) Phenanthrene	17.10	178	17633	0.393	ng/ul	96
13) Anthracene	17.19	178	15643	0.360	ng/ul	96
15) Fluoranthene	19.12	202	21997	0.337	ng/ul	97
17) Pyrene	19.48	202	22166	0.358	ng/ul	96
18) Benzo(a)anthracene	21.23	228	22079	0.372	ng/ul	98
19) Chrysene	21.28	228	23186	0.392	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	24638	0.400	ng/ul	93
22) Benzo(k)fluoranthene	22.83	252	25429	0.416	ng/ul	97
23) Benzo(a)pyrene	23.35	252	21989	0.391	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.63	276	30329	0.430	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.65	278	25307	0.427	ng/ul	97
26) Benzo(g,h,i)perylene	26.29	276	25337	0.426	ng/ul#	92

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

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