

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022520\
 Data File : BM025059.D
 Acq On : 26 Feb 2020 00:17
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.485

Quant Time: Feb 26 00:55:46 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	7.70	152	2510	0.40	ng/ul	0.00
2) Naphthalene-d8	10.47	136	9626	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	5919	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	13809	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	16313	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	18545	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	6850	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	18676	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.52	128	11559	0.394	ng/ul#	84
5) 2-Methylnaphthalene	12.14	142	8050	0.381	ng/ul	99
7) Acenaphthylene	14.04	152	12997	0.381	ng/ul	95
8) Acenaphthene	14.39	153	8931	0.390	ng/ul	99
9) Fluorene	15.37	166	10733	0.373	ng/ul	95
11) Pentachlorophenol	16.72	266	1562	0.370	ng/ul	99
12) Phenanthrene	17.11	178	18547	0.382	ng/ul	96
13) Anthracene	17.20	178	17414	0.370	ng/ul	96
15) Fluoranthene	19.12	202	23943	0.339	ng/ul	98
17) Pyrene	19.48	202	24337	0.331	ng/ul	99
18) Benzo(a)anthracene	21.23	228	26233	0.372	ng/ul	98
19) Chrysene	21.28	228	26980	0.385	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	28219	0.388	ng/ul	92
22) Benzo(k)fluoranthene	22.84	252	28430	0.395	ng/ul#	95
23) Benzo(a)pyrene	23.35	252	25324	0.382	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.64	276	32405	0.390	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.66	278	27461	0.393	ng/ul	96
26) Benzo(g,h,i)perylene	26.30	276	27259	0.389	ng/ul#	93

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

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