

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025127.D
 Acq On : 28 Feb 2020 09:37
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.493

Quant Time: Feb 28 10:52:25 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	2573	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.46	136	9568	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.31	164	5669	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.06	188	13031	0.40	ng/ul	-0.01
16) Chrysene-d12	21.24	240	13901	0.40	ng/ul	-0.02
20) Perylene-d12	23.44	264	15534	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	6764	0.38	ng/ul	-0.02
14) Fluoranthene-d10	19.09	212	17202	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.51	128	11691	0.401	ng/ul#	83
5) 2-Methylnaphthalene	12.13	142	8001	0.381	ng/ul	99
7) Acenaphthylene	14.03	152	11560	0.354	ng/ul	95
8) Acenaphthene	14.38	153	8707	0.396	ng/ul	99
9) Fluorene	15.37	166	10270	0.373	ng/ul	97
11) Pentachlorophenol	16.71	266	1403	0.352	ng/ul	98
12) Phenanthrene	17.10	178	17760	0.387	ng/ul	96
13) Anthracene	17.19	178	15809	0.356	ng/ul	96
15) Fluoranthene	19.11	202	22280	0.335	ng/ul	97
17) Pyrene	19.48	202	22318	0.356	ng/ul	97
18) Benzo(a)anthracene	21.22	228	22142	0.369	ng/ul	98
19) Chrysene	21.28	228	23497	0.393	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	24311	0.399	ng/ul	93
22) Benzo(k)fluoranthene	22.83	252	25676	0.426	ng/ul	97
23) Benzo(a)pyrene	23.34	252	21603	0.389	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.63	276	29616	0.425	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.64	278	24829	0.424	ng/ul	97
26) Benzo(g,h,i)perylene	26.29	276	24683	0.421	ng/ul#	92

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

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