

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091219\
 Data File : BM022663.D
 Acq On : 12 Sep 2019 10:44
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.463

Quant Time: Sep 12 11:27:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 15:06:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	8259	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	33029	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	17183	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	40189	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	36328	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	34030	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	20220	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	42665	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	34324	0.365	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	23446	0.363	ng/ul	100
7) Acenaphthylene	14.19	152	29588	0.372	ng/ul	99
8) Acenaphthene	14.54	153	26517	0.398	ng/ul	99
9) Fluorene	15.52	166	27527	0.361	ng/ul	99
11) Pentachlorophenol	16.86	266	2713	0.333	ng/ul	99
12) Phenanthrene	17.25	178	45969	0.359	ng/ul	100
13) Anthracene	17.34	178	38749	0.353	ng/ul	100
15) Fluoranthene	19.26	202	51898	0.376	ng/ul	98
17) Pyrene	19.63	202	55825	0.339	ng/ul	99
18) Benzo(a)anthracene	21.37	228	47980	0.345	ng/ul	99
19) Chrysene	21.42	228	52881	0.362	ng/ul	99
21) Benzo(b)fluoranthene	22.98	252	52783	0.357	ng/ul	99
22) Benzo(k)fluoranthene	23.03	252	52489	0.348	ng/ul	99
23) Benzo(a)pyrene	23.57	252	45913	0.355	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.98	276	56426	0.358	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.00	278	46379	0.364	ng/ul	99
26) Benzo(g,h,i)perylene	26.69	276	48995	0.390	ng/ul	100

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

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