

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091619\
 Data File : BM022696.D
 Acq On : 16 Sep 2019 18:13
 Operator : HP/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.467

Quant Time: Sep 17 01:34:49 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	7.84	152	8415	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	33987	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	18294	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	43467	0.40	ng/ul	0.00
16) Chrysene-d12	21.38	240	37801	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	33047	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	20982	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.23	212	46757	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.68	128	35223	0.364	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	24386	0.367	ng/ul	100
7) Acenaphthylene	14.19	152	31287	0.369	ng/ul	100
8) Acenaphthene	14.53	153	28146	0.397	ng/ul	100
9) Fluorene	15.52	166	29550	0.364	ng/ul	100
11) Pentachlorophenol	16.86	266	3362	0.381	ng/ul	97
12) Phenanthrene	17.25	178	49511	0.358	ng/ul	100
13) Anthracene	17.34	178	41546	0.350	ng/ul	99
15) Fluoranthene	19.26	202	56446	0.378	ng/ul	99
17) Pyrene	19.63	202	60977	0.356	ng/ul	99
18) Benzo(a)anthracene	21.37	228	49871	0.345	ng/ul	99
19) Chrysene	21.42	228	54429	0.358	ng/ul	100
21) Benzo(b)fluoranthene	22.98	252	52989	0.369	ng/ul	99
22) Benzo(k)fluoranthene	23.02	252	51121	0.349	ng/ul	99
23) Benzo(a)pyrene	23.56	252	44972	0.358	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.97	276	50284	0.328	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.99	278	41414	0.334	ng/ul	99
26) Benzo(g,h,i)perylene	26.68	276	42568	0.349	ng/ul	99

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

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