

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN100118\
 Data File : BN002887.D
 Acq On : 01 Oct 2018 19:43
 Operator : MJ/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.476

Quant Time: Oct 02 02:29:28 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 02 02:29:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	1162	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	5278	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	3034	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	6985	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	6333	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4948	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3493	0.46	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	8656	0.48	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	5803	0.437	ng/ul#	96
5) 2-Methylnaphthalene	12.11	142	4042	0.470	ng/ul	99
7) Acenaphthylene	14.00	152	5769	0.450	ng/ul	97
8) Acenaphthene	14.34	153	4679	0.448	ng/ul	97
9) Fluorene	15.34	166	5227	0.468	ng/ul	94
11) Pentachlorophenol	16.73	266	407	0.517	ng/ul	97
12) Phenanthrene	17.08	178	8528	0.442	ng/ul#	89
13) Anthracene	17.18	178	7660	0.457	ng/ul#	93
15) Fluoranthene	19.10	202	9978	0.480	ng/ul	96
17) Pyrene	19.47	202	10079	0.488	ng/ul	97
18) Benzo(a)anthracene	21.22	228	8378	0.479	ng/ul#	85
19) Chrysene	21.27	228	9288	0.455	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	7811	0.478	ng/ul	89
22) Benzo(k)fluoranthene	22.83	252	8362	0.486	ng/ul#	91
23) Benzo(a)pyrene	23.36	252	7050	0.450	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.67	276	6985	0.370	ng/ul#	88
25) Dibenzo(a,h)anthracene	25.67	278	5553	0.367	ng/ul	95
26) Benzo(g,h,i)perylene	26.34	276	5922	0.363	ng/ul	99

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

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