

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN003010.D
 Acq On : 09 Oct 2018 04:42
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.495

Quant Time: Oct 09 08:37:58 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 09 08:37:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	1091	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	5110	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3132	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	7722	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	8246	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	6778	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3138	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	9246	0.42	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.47	128	5228	0.399	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	3585	0.397	ng/ul	100
7) Acenaphthylene	14.00	152	5537	0.386	ng/ul	99
8) Acenaphthene	14.34	153	4295	0.395	ng/ul	98
9) Fluorene	15.34	166	5007	0.402	ng/ul	99
11) Pentachlorophenol	16.74	266	328	0.438	ng/ul	96
12) Phenanthrene	17.08	178	8430	0.392	ng/ul	99
13) Anthracene	17.18	178	7902	0.389	ng/ul	100
15) Fluoranthene	19.10	202	10496	0.416	ng/ul	99
17) Pyrene	19.47	202	10842	0.339	ng/ul	99
18) Benzo(a)anthracene	21.22	228	10228	0.385	ng/ul	100
19) Chrysene	21.27	228	10729	0.398	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	9420	0.386	ng/ul	96
22) Benzo(k)fluoranthene	22.83	252	9978	0.394	ng/ul	94
23) Benzo(a)pyrene	23.36	252	8514	0.389	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.67	276	8737	0.348	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.67	278	7070	0.351	ng/ul	96
26) Benzo(g,h,i)perylene	26.35	276	7345	0.346	ng/ul	97

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

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