

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061318\
 Data File : BN001860.D
 Acq On : 14 Jun 2018 01:11
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.479

Quant Time: Jun 14 05:26:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN061318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 13 12:02:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	6383	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	25110	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	14950	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	35964	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	30309	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	23800	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.09	152	16378	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	41854	0.38	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.54	128	25769	0.398	ng/ul	96
5) 2-Methylnaphthalene	12.16	142	18775	0.406	ng/ul	98
7) Acenaphthylene	14.07	152	28549	0.391	ng/ul	95
8) Acenaphthene	14.41	153	21370	0.398	ng/ul	97
9) Fluorene	15.40	166	24894	0.397	ng/ul#	93
11) Pentachlorophenol	16.74	266	2868	0.421	ng/ul	97
12) Phenanthrene	17.14	178	42053	0.392	ng/ul#	89
13) Anthracene	17.23	178	40029	0.394	ng/ul#	92
15) Fluoranthene	19.16	202	48254	0.379	ng/ul	96
17) Pyrene	19.52	202	48269	0.438	ng/ul	97
18) Benzo(a)anthracene	21.28	228	42661	0.407	ng/ul#	85
19) Chrysene	21.33	228	40904	0.392	ng/ul	97
21) Benzo(b)fluoranthene	22.86	252	36348	0.421	ng/ul	86
22) Benzo(k)fluoranthene	22.90	252	36671	0.426	ng/ul#	91
23) Benzo(a)pyrene	23.43	252	32973	0.411	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.77	276	32975	0.366	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.78	278	26298	0.366	ng/ul#	89
26) Benzo(g,h,i)perylene	26.44	276	26954	0.352	ng/ul	98

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

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