

Data Path : Z:\HPCHEM1\BNA_N\Data\BN030518\
 Data File : BN000253.D
 Acq On : 05 Mar 2018 19:44
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.461

Quant Time: Mar 06 04:39:17 2018
 Quant Method : Z:\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN030218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 04:37:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	4285	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	13680	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.06	164	5871	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.78	188	13245	0.40	ng/ul	0.00
16) Chrysene-d12	21.89	240	8122	0.40	ng/ul	0.00
20) Perylene-d12	24.37	264	6689	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.84	152	6526	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.77	212	10946	0.36	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.34	128	16203	0.370	ng/ul#	87
5) 2-Methylnaphthalene	12.92	142	10057	0.362	ng/ul	99
7) Acenaphthylene	14.79	152	10466	0.358	ng/ul	97
8) Acenaphthene	15.12	153	10410	0.369	ng/ul	94
9) Fluorene	16.10	166	11337	0.370	ng/ul#	91
11) Pentachlorophenol	17.42	266	847	0.318	ng/ul	100
12) Phenanthrene	17.82	178	19767	0.366	ng/ul#	87
13) Anthracene	17.91	178	14251	0.355	ng/ul#	91
15) Fluoranthene	19.80	202	18676	0.356	ng/ul	62
17) Pyrene	20.15	202	18460	0.386	ng/ul#	55
18) Benzo(a)anthracene	21.88	228	11564	0.363	ng/ul#	84
19) Chrysene	21.93	228	16214	0.374	ng/ul	95
21) Benzo(b)fluoranthene	23.61	252	15083	0.375	ng/ul	87
22) Benzo(k)fluoranthene	23.66	252	13830	0.365	ng/ul#	89
23) Benzo(a)pyrene	24.26	252	11740	0.373	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.91	276	12990	0.365	ng/ul#	76
25) Dibenzo(a,h)anthracene	26.93	278	9335	0.364	ng/ul#	80
26) Benzo(g,h,i)perylene	27.70	276	12625	0.361	ng/ul#	82

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

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