

Data Path : Z:\HPCHEM1\BNA E\DATA\BE050318\
 Data File : BE096322.D
 Acq On : 3 May 2018 11:56
 Operator : SJ/JU
 Sample : SSTD1.674
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.674

Quant Time: May 03 12:42:23 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE050318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 03 12:01:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	562	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2180	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	1313	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	3575	0.40	ng/ul	0.00
16) Chrysene-d12	21.74	240	4391	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	3640	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	6535	1.67	ng/ul	0.00
14) Fluoranthene-d10	19.63	212	23697	1.56	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.22	128	9504	1.583	ng/ul#	89
5) 2-Methylnaphthalene	12.80	142	6657	1.689	ng/ul	99
7) Acenaphthylene	14.67	152	11157	1.645	ng/ul	97
8) Acenaphthene	15.00	153	7925	1.614	ng/ul	93
9) Fluorene	15.97	166	11087	1.680	ng/ul#	93
11) Pentachlorophenol	17.31	266	820	1.731	ng/ul	96
12) Phenanthrene	17.69	178	17086	1.585	ng/ul#	88
13) Anthracene	17.78	178	17022	1.647	ng/ul#	91
15) Fluoranthene	19.66	202	23837	1.546	ng/ul	82
17) Pyrene	20.02	202	7354	1.463	ng/ul#	75
18) Benzo(a)anthracene	21.73	228	9787	1.389	ng/ul#	86
19) Chrysene	21.78	228	21434	1.644	ng/ul	96
21) Benzo(b)fluoranthene	23.54	252	20412	1.649	ng/ul	83
22) Benzo(k)fluoranthene	23.59	252	19276	1.601	ng/ul#	86
23) Benzo(a)pyrene	24.23	252	18827	1.617	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	27.12	276	21145	1.646	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.14	278	17282	1.645	ng/ul#	86
26) Benzo(g,h,i)perylene	27.98	276	17749	1.640	ng/ul#	92

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

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