

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101118\
 Data File : BE097877.D
 Acq On : 12 Oct 2018 3:19
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.453

Quant Time: Oct 12 05:00:03 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 12 04:56:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2327	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	12295	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	8061	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.31	188	19501	0.40	ng/ul	0.00
16) Chrysene-d12	21.49	240	20697	0.40	ng/ul	0.00
20) Perylene-d12	24.06	264	20939	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	9037	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.34	212	26856	0.43	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	12870	0.430	ng/ul	99
5) 2-Methylnaphthalene	12.36	142	9541	0.441	ng/ul	99
7) Acenaphthylene	14.27	152	14295	0.469	ng/ul	97
8) Acenaphthene	14.61	153	10674	0.451	ng/ul	98
9) Fluorene	15.61	166	12739	0.413	ng/ul	98
11) Pentachlorophenol	16.97	266	1019	0.371	ng/ul	97
12) Phenanthrene	17.35	178	20246	0.439	ng/ul	98
13) Anthracene	17.44	178	20361	0.476	ng/ul	100
15) Fluoranthene	19.37	202	25290	0.427	ng/ul	96
17) Pyrene	19.73	202	30966	0.502	ng/ul	99
18) Benzo(a)anthracene	21.47	228	34091	0.545	ng/ul	94
19) Chrysene	21.53	228	26066	0.464	ng/ul#	91
21) Benzo(b)fluoranthene	23.27	252	25153	0.449	ng/ul#	79
22) Benzo(k)fluoranthene	23.32	252	25468	0.425	ng/ul#	76
23) Benzo(a)pyrene	23.94	252	25248	0.460	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	26.77	276	28205	0.452	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.79	278	23968	0.465	ng/ul#	80
26) Benzo(g,h,i)perylene	27.60	276	23854	0.439	ng/ul#	86

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

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