

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE101018\
 Data File : BE097831.D
 Acq On : 10 Oct 2018 17:09
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.442

Quant Time: Oct 10 18:45:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 10 12:43:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2692	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	11742	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	7121	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	20835	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	36820	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	29594	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	8199	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	31032	0.46	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.74	128	12171	0.431	ng/ul	98
5) 2-Methylnaphthalene	12.36	142	8234	0.438	ng/ul	99
7) Acenaphthylene	14.26	152	11049	0.440	ng/ul	97
8) Acenaphthene	14.60	153	9169	0.439	ng/ul	98
9) Fluorene	15.59	166	11055	0.439	ng/ul	99
11) Pentachlorophenol	16.94	266	931	0.490	ng/ul	97
12) Phenanthrene	17.34	178	21496	0.432	ng/ul	99
13) Anthracene	17.43	178	18812	0.445	ng/ul	99
15) Fluoranthene	19.36	202	29809	0.454	ng/ul	100
17) Pyrene	19.72	202	38107	0.421	ng/ul	100
18) Benzo(a)anthracene	21.46	228	44028	0.431	ng/ul	99
19) Chrysene	21.52	228	40014	0.406	ng/ul	98
21) Benzo(b)fluoranthene	23.25	252	34991	0.458	ng/ul	99
22) Benzo(k)fluoranthene	23.30	252	36524	0.421	ng/ul	97
23) Benzo(a)pyrene	23.91	252	32493	0.464	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.71	276	34901	0.439	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.74	278	28864	0.435	ng/ul	99
26) Benzo(g,h,i)perylene	27.53	276	32279	0.430	ng/ul	99

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

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