

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102518\
 Data File : BE098049.D
 Acq On : 25 Oct 2018 10:30
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4

Quant Time: Oct 25 14:52:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 25 14:48:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	965	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	4892	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3457	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	8497	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	10091	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	10089	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3269	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10373	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	4455	0.380	ng/ul	98
5) 2-Methylnaphthalene	12.35	142	3393	0.383	ng/ul	99
7) Acenaphthylene	14.25	152	5213	0.382	ng/ul	98
8) Acenaphthene	14.60	153	3794	0.372	ng/ul	99
9) Fluorene	15.59	166	4614	0.376	ng/ul	99
11) Pentachlorophenol	16.95	266	258	0.580	ng/ul	93
12) Phenanthrene	17.33	178	7367	0.376	ng/ul	100
13) Anthracene	17.42	178	7281	0.376	ng/ul	99
15) Fluoranthene	19.35	202	9390	0.376	ng/ul	99
17) Pyrene	19.71	202	11170	0.474	ng/ul	99
18) Benzo(a)anthracene	21.45	228	13629	0.459	ng/ul	100
19) Chrysene	21.51	228	10093	0.440	ng/ul	100
21) Benzo(b)fluoranthene	23.24	252	10349	0.370	ng/ul	99
22) Benzo(k)fluoranthene	23.29	252	10264	0.375	ng/ul	98
23) Benzo(a)pyrene	23.91	252	9931	0.369	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.71	276	11481	0.373	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.74	278	9586	0.373	ng/ul	99
26) Benzo(g,h,i)perylene	27.54	276	9723	0.375	ng/ul	99

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

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