

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101118\
 Data File : BE097849.D
 Acq On : 11 Oct 2018 9:46
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4

Quant Time: Oct 11 16:16:13 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 : Thu Oct 11 14:16:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2393	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10705	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	7659	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	23250	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	33968	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	28347	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	7782	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	33538	0.45	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	11231	0.431	ng/ul	99
5) 2-Methylnaphthalene	12.36	142	8150	0.432	ng/ul	98
7) Acenaphthylene	14.26	152	13044	0.450	ng/ul	99
8) Acenaphthene	14.60	153	9925	0.441	ng/ul	98
9) Fluorene	15.60	166	12871	0.439	ng/ul	98
11) Pentachlorophenol	16.95	266	1368	0.417	ng/ul	98
12) Phenanthrene	17.34	178	24088	0.438	ng/ul	99
13) Anthracene	17.43	178	22670	0.444	ng/ul	99
15) Fluoranthene	19.36	202	31648	0.448	ng/ul	98
17) Pyrene	19.72	202	39608	0.391	ng/ul	98
18) Benzo(a)anthracene	21.46	228	41763	0.407	ng/ul	98
19) Chrysene	21.51	228	35892	0.389	ng/ul	98
21) Benzo(b)fluoranthene	23.24	252	33505	0.442	ng/ul	99
22) Benzo(k)fluoranthene	23.30	252	35225	0.434	ng/ul	98
23) Benzo(a)pyrene	23.91	252	32709	0.440	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.71	276	37156	0.440	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.75	278	30995	0.444	ng/ul	97
26) Benzo(g,h,i)perylene	27.54	276	31919	0.434	ng/ul	98

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

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