

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101317\
 Data File : BE094330.D
 Acq On : 13 Oct 2017 18:45
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.434

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:22:51 PM

Quant Time: Oct 14 03:54:13 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 14 02:15:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	1584	0.40	ng/ul	0.01
2) Naphthalene-d8	10.71	136	8026	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.54	164	4404	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	9668	0.40	ng/ul	0.02
16) Chrysene-d12	21.44	240	9460	0.40	ng/ul	0.01
20) Perylene-d12	23.98	264	9245	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	5047	0.41	ng/ul	0.01
14) Fluoranthene-d10	19.30	212	11699	0.36	ng/ul	0.01
Target Compounds						Qvalue
3) Naphthalene	10.77	128	7757	0.39	ng/ul#	93
5) 2-Methylnaphthalene	12.37	142	5359	0.42	ng/ul	99
7) Acenaphthylene	14.26	152	8406	0.49	ng/ul	97
8) Acenaphthene	14.60	153	5661	0.40	ng/ul	98
9) Fluorene	15.59	166	6576	0.40	ng/ul	94
11) Pentachlorophenol	16.96	266	332	0.29	ng/ul	91
12) Phenanthrene	17.31	178	10550m	0.42	ng/ul	
13) Anthracene	17.41	178	10564	0.43	ng/ul	97
15) Fluoranthene	19.33	202	12049	0.36	ng/ul	84
17) Pyrene	19.69	202	12732	0.48	ng/ul#	84
18) Benzo(a)anthracene	21.42	228	11413	0.45	ng/ul#	73
19) Chrysene	21.48	228	10562	0.36	ng/ul#	69
21) Benzo(b)fluoranthene	23.20	252	10955	0.44	ng/ul#	45
22) Benzo(k)fluoranthene	23.25	252	10684	0.37	ng/ul#	39
23) Benzo(a)pyrene	23.87	252	10570	0.41	ng/ul#	46
24) Indeno(1,2,3-cd)pyrene	26.68	276	11864	0.44	ng/ul#	89
25) Dibenzo(a,h)anthracene	26.69	278	10016	0.45	ng/ul#	46
26) Benzo(g,h,i)perylene	27.53	276	9977	0.43	ng/ul#	53

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

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