

Data Path : Z:\HPCHEM1\BNA_E\Data\BE110317\
 Data File : BE094701.D
 Acq On : 3 Nov 2017 10:25
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.472

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:14:56 PM

Quant Time: Nov 03 22:56:09 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 03 23:54:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1250	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	7647	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	4117	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	8803	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	9191	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	8527	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	4519	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	9772	0.37	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.77	128	6803	0.364	ng/ul#	94
5) 2-Methylnaphthalene	12.37	142	4640	0.355	ng/ul	99
7) Acenaphthylene	14.25	152	6816	0.366	ng/ul	98
8) Acenaphthene	14.59	153	4877	0.364	ng/ul	98
9) Fluorene	15.58	166	5323	0.354	ng/ul#	94
11) Pentachlorophenol	17.10	266	123m	0.280	ng/ul	
12) Phenanthrene	17.31	178	8370	0.360	ng/ul#	91
13) Anthracene	17.41	178	8591	0.374	ng/ul#	94
15) Fluoranthene	19.32	202	9984	0.371	ng/ul	85
17) Pyrene	19.68	202	10559	0.372	ng/ul#	83
18) Benzo(a)anthracene	21.42	228	9724	0.370	ng/ul#	87
19) Chrysene	21.47	228	10199	0.376	ng/ul	98
21) Benzo(b)fluoranthene	23.19	252	8677	0.362	ng/ul	89
22) Benzo(k)fluoranthene	23.24	252	10415m	0.405	ng/ul	
23) Benzo(a)pyrene	23.86	252	8854	0.369	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.69	276	7381	0.298	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.67	278	6396	0.306	ng/ul	93
26) Benzo(g,h,i)perylene	27.52	276	5273	0.265	ng/ul	97

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

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