

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094715.D
 Acq On : 3 Nov 2017 19:43
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.473

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:15:33 PM

Quant Time: Nov 03 23:51:30 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 04 00:03:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1141	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	6985	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.52	164	3873	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8347	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	8449	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	8366	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4140	0.37	ng/ul	-0.01
14) Fluoranthene-d10	19.29	212	9513	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	6174	0.362	ng/ul#	94
5) 2-Methylnaphthalene	12.36	142	4218	0.353	ng/ul	100
7) Acenaphthylene	14.24	152	6456	0.369	ng/ul	98
8) Acenaphthene	14.58	153	4513	0.358	ng/ul	98
9) Fluorene	15.57	166	5115	0.362	ng/ul	91
11) Pentachlorophenol	17.01	266	181m	0.434	ng/ul	
12) Phenanthrene	17.31	178	7985	0.362	ng/ul#	93
13) Anthracene	17.39	178	8786m	0.404	ng/ul	
15) Fluoranthene	19.32	202	9732	0.382	ng/ul	84
17) Pyrene	19.68	202	10456	0.400	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	9333	0.386	ng/ul#	89
19) Chrysene	21.47	228	9117	0.365	ng/ul	92
21) Benzo(b)fluoranthene	23.18	252	8690	0.369	ng/ul	91
22) Benzo(k)fluoranthene	23.23	252	9171	0.364	ng/ul#	91
23) Benzo(a)pyrene	23.86	252	8766	0.373	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	26.67	276	8232	0.338	ng/ul#	82
25) Dibenzo(a,h)anthracene	26.66	278	7133	0.348	ng/ul#	84
26) Benzo(g,h,i)perylene	27.52	276	6030	0.309	ng/ul#	79

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

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