

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094741.D
 Acq On : 4 Nov 2017 11:02
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.475

Manual Integrations
 APPROVED

Sohil
 11/6/2017 5:58:54 PM

Quant Time: Nov 06 00:29:11 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 06:57:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1074	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	6441	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	3563	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	7568	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	7971	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	7556	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3816	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	8653	0.39	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	5643	0.358	ng/ul#	95
5) 2-Methylnaphthalene	12.36	142	3925	0.356	ng/ul	98
7) Acenaphthylene	14.24	152	6152	0.382	ng/ul	98
8) Acenaphthene	14.58	153	4167	0.360	ng/ul	99
9) Fluorene	15.57	166	4763	0.366	ng/ul	94
11) Pentachlorophenol	17.00	266	221	0.585	ng/ul#	64
12) Phenanthrene	17.29	178	7520m	0.376	ng/ul	
13) Anthracene	17.40	178	7599	0.385	ng/ul	95
15) Fluoranthene	19.32	202	9000	0.389	ng/ul	83
17) Pyrene	19.68	202	9523	0.386	ng/ul#	85
18) Benzo(a)anthracene	21.41	228	8565	0.376	ng/ul#	77
19) Chrysene	21.47	228	8890	0.377	ng/ul#	66
21) Benzo(b)fluoranthene	23.19	252	8301	0.391	ng/ul#	62
22) Benzo(k)fluoranthene	23.24	252	8097	0.355	ng/ul#	51
23) Benzo(a)pyrene	23.86	252	8627	0.406	ng/ul#	64
24) Indeno(1,2,3-cd)pyrene	26.68	276	7961	0.362	ng/ul#	77
25) Dibenzo(a,h)anthracene	26.67	278	6963	0.376	ng/ul#	55
26) Benzo(g,h,i)perylene	27.53	276	5564	0.315	ng/ul#	52

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

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