

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110317\
 Data File : BE094733.D
 Acq On : 4 Nov 2017 6:22
 Operator : SJ/JU
 Sample : I5933-02 5X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-9(5-6) 101617

Manual Integrations
 APPROVED

Sohil
 11/4/2017 2:17:11 PM

Quant Time: Nov 04 05:58:51 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	7.89	152	1814	0.40	ng/ul	-0.03
2) Naphthalene-d8	10.68	136	10883	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.51	164	5873	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.24	188	12678	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	12322	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	12710	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	1030	0.06	ng/ul	-0.02
14) Fluoranthene-d10	19.27	212	2282	0.06	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	4533	0.170	ng/ul#	89
5) 2-Methylnaphthalene	12.34	142	1339	0.072	ng/ul	95
7) Acenaphthylene	14.24	152	2525	0.095	ng/ul#	87
8) Acenaphthene	14.57	153	6222	0.326	ng/ul	95
9) Fluorene	15.56	166	6798	0.317	ng/ul	93
12) Phenanthrene	17.29	178	94552	2.824	ng/ul#	93
13) Anthracene	17.38	178	28941	0.876	ng/ul#	92
15) Fluoranthene	19.31	202	222314	5.742	ng/ul	84
17) Pyrene	19.67	202	304649	7.998	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	148548	4.217	ng/ul#	88
19) Chrysene	21.46	228	145421	3.994	ng/ul	99
21) Benzo(b)fluoranthene	23.18	252	143752m	4.021	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	54113m	1.412	ng/ul	
23) Benzo(a)pyrene	23.84	252	123567	3.458	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	26.65	276	57552	1.556	ng/ul#	74
25) Dibenzo(a,h)anthracene	26.64	278	15315	0.491	ng/ul	96
26) Benzo(g,h,i)perylene	27.49	276	52674	1.774	ng/ul	97

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

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