

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094848.D
 Acq On : 23 Nov 2017 00:49
 Operator : SJ/JU
 Sample : I6496-08
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB8

Manual Integrations
 APPROVED

Sohil
 11/27/2017 3:26:03 PM

Quant Time: Nov 23 03:18:37 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 23 02:50:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	658	0.40	ng/ul	-0.04
2) Naphthalene-d8	10.68	136	4555	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.50	164	3244	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.24	188	7781	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	7856	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	8679	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1618	0.22	ng/ul	-0.01
14) Fluoranthene-d10	19.27	212	5712	0.21	ng/ul	-0.02
Target Compounds				Ovalue		
3) Naphthalene	10.72	128	426	0.039	ng/ul#	65
5) 2-Methylnaphthalene	12.36	142	303	0.040	ng/ul	100
7) Acenaphthylene	14.22	152	581	0.039	ng/ul#	73
12) Phenanthrene	17.27	178	4553	0.218	ng/ul#	90
13) Anthracene	17.36	178	1021	0.048	ng/ul#	77
15) Fluoranthene	19.30	202	15089	0.550	ng/ul	84
17) Pyrene	19.66	202	14328	0.583	ng/ul#	81
18) Benzo(a)anthracene	21.40	228	8916	0.373	ng/ul#	89
19) Chrysene	21.46	228	8803	0.369	ng/ul	97
21) Benzo(b)fluoranthene	23.17	252	10983m	0.425	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	4658m	0.169	ng/ul	
23) Benzo(a)pyrene	23.83	252	7628	0.297	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.66	276	4925	0.176	ng/ul#	73
25) Dibenzo(a,h)anthracene	26.64	278	1305	0.055	ng/ul#	68
26) Benzo(g,h,i)perylene	27.50	276	3962	0.169	ng/ul#	93

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

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