

Data Path : Z:\HPCHEM1\BNA E\DATA\BE103117\
 Data File : BE094663.D
 Acq On : 31 Oct 2017 18:15
 Operator : SJ/JU
 Sample : I5786-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-38(1-3)_100917

Manual Integrations
 APPROVED

Sohil
 11/1/2017 9:31:58 AM

Quant Time: Nov 01 02:50:36 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 01:43:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	2456	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	8074	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4336	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8533m	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	7828	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	10186m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	6188	0.47	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	12981	0.51	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.34	142	8728	0.632	ng/ul	100
8) Acenaphthene	14.57	153	10717	0.760	ng/ul	97
9) Fluorene	15.56	166	7615	0.481	ng/ul	92
12) Phenanthrene	17.29	178	18544	0.823	ng/ul#	93
13) Anthracene	17.38	178	3988	0.179	ng/ul#	95
15) Fluoranthene	19.31	202	13257	0.509	ng/ul	84
17) Pyrene	19.67	202	13930	0.576	ng/ul#	84
18) Benzo(a)anthracene	21.41	228	4262	0.190	ng/ul#	88
19) Chrysene	21.46	228	4562	0.197	ng/ul#	87
21) Benzo(b)fluoranthene	23.18	252	4360m	0.152	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	1219m	0.040	ng/ul	
23) Benzo(a)pyrene	23.84	252	2545m	0.089	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.68	276	1394m	0.047	ng/ul	
26) Benzo(g,h,i)perylene	27.52	276	1073m	0.045	ng/ul	

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

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