

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094630.D
 Acq On : 28 Oct 2017 19:34
 Operator : SJ/JU
 Sample : I5842-06
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA6

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:40:15 PM

Quant Time: Oct 30 04:24:02 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 02:53:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1773	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	10276	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.51	164	5736	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12261	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12861	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	10635m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	3294	0.20	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	8600	0.24	ng/ul	0.00
Target Compounds				Ovalue		
5) 2-Methylnaphthalene	12.34	142	378	0.022	ng/ul	99
12) Phenanthrene	17.29	178	3130m	0.097	ng/ul	
13) Anthracene	17.29	178	3348	0.105	ng/ul	93
15) Fluoranthene	19.31	202	6906	0.184	ng/ul	83
17) Pyrene	19.67	202	7296	0.184	ng/ul#	88
18) Benzo(a)anthracene	21.41	228	2997	0.082	ng/ul#	77
19) Chrysene	21.46	228	4264	0.112	ng/ul#	80
21) Benzo(b)fluoranthene	23.19	252	5711m	0.191	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	1721m	0.054	ng/ul	
23) Benzo(a)pyrene	23.85	252	2912m	0.097	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.67	276	2485m	0.080	ng/ul	
26) Benzo(g,h,i)perylene	27.53	276	1851m	0.074	ng/ul	

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

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