

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094610.D
 Acq On : 27 Oct 2017 23:47
 Operator : SJ/JU
 Sample : I5842-14
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB2

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:37:12 PM

Quant Time: Oct 30 02:16:12 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 01:11:00 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1226	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.69	136	7171	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	4265	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9289m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	8536	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	9480m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3716	0.32	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	8825	0.32	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	2487m	0.101	ng/ul	
15) Fluoranthene	19.31	202	8440	0.298	ng/ul	84
17) Pyrene	19.67	202	8189	0.310	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	2662	0.109	ng/ul#	73
19) Chrysene	21.47	228	4988	0.198	ng/ul#	81
21) Benzo(b)fluoranthene	23.18	252	6832m	0.256	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	2069m	0.072	ng/ul	
23) Benzo(a)pyrene	23.85	252	3709m	0.139	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.67	276	2665m	0.097	ng/ul	
26) Benzo(g,h,i)perylene	27.52	276	2660m	0.120	ng/ul	

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

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