

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094608.D
 Acq On : 27 Oct 2017 22:33
 Operator : SJ/JU
 Sample : I5842-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB1

Manual Integrations
 APPROVED

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

Quant Time: Oct 30 02:04:07 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	1062	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	6240	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	3753	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8195m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	7548	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	6801m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3488	0.34	ng/ul	-0.01
14) Fluoranthene-d10	19.29	212	8199	0.34	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.29	178	2294	0.106	ng/ul#	80
15) Fluoranthene	19.31	202	9947	0.397	ng/ul	85
17) Pyrene	19.68	202	10086	0.432	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	4507	0.209	ng/ul#	61
19) Chrysene	21.47	228	5470	0.245	ng/ul#	54
21) Benzo(b)fluoranthene	23.19	252	7235m	0.378	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	2646m	0.129	ng/ul	
23) Benzo(a)pyrene	23.85	252	4401m	0.230	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.67	276	2518m	0.127	ng/ul	
26) Benzo(g,h,i)perylene	27.52	276	1818m	0.114	ng/ul	

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

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