

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094620.D
 Acq On : 28 Oct 2017 13:27
 Operator : SJ/JU
 Sample : I5842-04
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA4

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:40:00 PM

Quant Time: Oct 30 03:48:00 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	1401	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	8212	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	4670	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9944	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	9534	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	8996m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3192	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	7694	0.26	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	533	0.027	ng/ul#	72
5) 2-Methylnaphthalene	12.35	142	482	0.034	ng/ul	97
12) Phenanthrene	17.29	178	9909	0.377	ng/ul#	91
13) Anthracene	17.38	178	1147	0.044	ng/ul#	83
15) Fluoranthene	19.31	202	23433	0.772	ng/ul	84
17) Pyrene	19.67	202	20983	0.712	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	9006	0.330	ng/ul#	89
19) Chrysene	21.46	228	11641	0.413	ng/ul	93
21) Benzo(b)fluoranthene	23.18	252	14694m	0.581	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	4587m	0.169	ng/ul	
23) Benzo(a)pyrene	23.85	252	8168m	0.323	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.67	276	6015m	0.230	ng/ul	
25) Dibenzo(a,h)anthracene	26.65	278	1447m	0.066	ng/ul	
26) Benzo(g,h,i)perylene	27.52	276	4587m	0.218	ng/ul	

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

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