

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094628.D
 Acq On : 28 Oct 2017 18:20
 Operator : SJ/JU
 Sample : I5842-20
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB9

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:40:11 PM

Quant Time: Oct 30 04:15:50 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	2276	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	13228	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	7548	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	16051	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	15361	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	17056m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	4982	0.23	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	11671	0.25	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.24	152	5330	0.156	ng/ul	99
8) Acenaphthene	14.57	153	859	0.035	ng/ul#	89
9) Fluorene	15.57	166	1088	0.039	ng/ul#	88
12) Phenanthrene	17.29	178	30256	0.714	ng/ul#	90
13) Anthracene	17.38	178	19420	0.464	ng/ul#	89
15) Fluoranthene	19.31	202	129212	2.636	ng/ul	84
17) Pyrene	19.67	202	231507	4.875	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	114691	2.612	ng/ul#	87
19) Chrysene	21.46	228	136678	3.011	ng/ul	97
21) Benzo(b)fluoranthene	23.19	252	165718	3.454	ng/ul	87
22) Benzo(k)fluoranthene	23.23	252	57071m	1.110	ng/ul	
23) Benzo(a)pyrene	23.85	252	74398m	1.551	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.66	276	27715m	0.558	ng/ul	
25) Dibenzo(a,h)anthracene	26.66	278	9624m	0.230	ng/ul	
26) Benzo(g,h,i)perylene	27.51	276	16335m	0.410	ng/ul	

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

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