

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102617\
 Data File : BE094572.D
 Acq On : 26 Oct 2017 21:05
 Operator : SJ/JU
 Sample : I5786-14 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-66(2.5-4.5) 101017

Manual Integrations
 APPROVED

Sohil
 10/27/2017 5:35:40 PM

Quant Time: Oct 27 08:57:51 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 27 04:05:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1208	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	6916	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.51	164	3987	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8943	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	9222	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	9554	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	702	0.06	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	1787	0.07	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	4245	0.251	ng/ul#	96
5) 2-Methylnaphthalene	12.35	142	1036	0.088	ng/ul	97
7) Acenaphthylene	14.23	152	3371	0.187	ng/ul	99
8) Acenaphthene	14.57	153	1188	0.092	ng/ul#	95
9) Fluorene	15.57	166	3005	0.206	ng/ul	95
12) Phenanthrene	17.29	178	62559	2.648	ng/ul#	91
13) Anthracene	17.38	178	11997	0.515	ng/ul#	91
15) Fluoranthene	19.31	202	155852	5.707	ng/ul	83
17) Pyrene	19.67	202	149605	5.248	ng/ul#	80
18) Benzo(a)anthracene	21.40	228	71029	2.694	ng/ul#	86
19) Chrysene	21.46	228	82425	3.025	ng/ul	99
21) Benzo(b)fluoranthene	23.17	252	107500m	4.000	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	36051m	1.252	ng/ul	
23) Benzo(a)pyrene	23.83	252	74945	2.790	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.64	276	52475	1.888	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.64	278	12597	0.538	ng/ul#	93
26) Benzo(g,h,i)perylene	27.47	276	42556	1.907	ng/ul	97

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

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