

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094367.D
 Acq On : 14 Oct 2017 22:27
 Operator : SJ/JU
 Sample : I5548-10
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC8

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:22 PM

Quant Time: Oct 15 02:42:22 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 14 05:22:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1521	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.70	136	8033	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	5157	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11694	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	11934	0.40	ng/ul	-0.02
20) Perylene-d12	23.95	264	10696m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	2289	0.18	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	6359	0.16	ng/ul	-0.02
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	4842	0.246	ng/ul#	94
5) 2-Methylnaphthalene	12.35	142	3210	0.251	ng/ul	97
7) Acenaphthylene	14.24	152	4113	0.205	ng/ul	97
8) Acenaphthene	14.58	153	3185	0.192	ng/ul	98
9) Fluorene	15.57	166	3525	0.182	ng/ul	92
12) Phenanthrene	17.29	178	105330	3.447	ng/ul#	88
13) Anthracene	17.38	178	16121	0.542	ng/ul#	90
15) Fluoranthene	19.31	202	222047	5.530	ng/ul	84
17) Pyrene	19.67	202	194985	5.872	ng/ul#	79
18) Benzo(a)anthracene	21.40	228	93559	2.955	ng/ul#	86
19) Chrysene	21.46	228	111890	3.060	ng/ul	99
21) Benzo(b)fluoranthene	23.17	252	137856m	4.745	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	47375m	1.411	ng/ul	
23) Benzo(a)pyrene	23.83	252	84374m	2.795	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.62	276	62714m	2.030	ng/ul	
25) Dibenzo(a,h)anthracene	26.62	278	16780m	0.648	ng/ul	
26) Benzo(a,h,i)perylene	27.46	276	52540m	1.972	ng/ul	

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

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