

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE101417\
 Data File : BE094374.D
 Acq On : 15 Oct 2017 2:44
 Operator : SJ/JU
 Sample : I5548-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB3

Manual Integrations
 APPROVED

Sohil
 10/16/2017 8:25:36 PM

Quant Time: Oct 16 05:40:26 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 05:09:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1835	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	10209	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	6069	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12971	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12060	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	13098m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3144	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	7870	0.18	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	1132	0.045	ng/ul#	77
5) 2-Methylnaphthalene	12.35	142	923	0.057	ng/ul	98
7) Acenaphthylene	14.25	152	1068	0.045	ng/ul#	81
8) Acenaphthene	14.59	153	703	0.036	ng/ul#	92
9) Fluorene	15.57	166	668	0.029	ng/ul#	77
12) Phenanthrene	17.29	178	20280	0.598	ng/ul#	87
13) Anthracene	17.39	178	2997	0.091	ng/ul#	90
15) Fluoranthene	19.31	202	60042	1.348	ng/ul	84
17) Pyrene	19.67	202	53527	1.595	ng/ul#	84
18) Benzo(a)anthracene	21.41	228	26032	0.814	ng/ul#	92
19) Chrysene	21.46	228	31763	0.860	ng/ul#	93
21) Benzo(b)fluoranthene	23.18	252	39049	1.097	ng/ul	91
22) Benzo(k)fluoranthene	23.22	252	15783m	0.384	ng/ul	
23) Benzo(a)pyrene	23.84	252	26727m	0.723	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.64	276	20564m	0.544	ng/ul	
25) Dibenzo(a,h)anthracene	26.64	278	5206m	0.164	ng/ul	
26) Benzo(a,h,i)perylene	27.47	276	16797m	0.515	ng/ul	

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

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