

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080417\
 Data File : BE093665.D
 Acq On : 4 Aug 2017 11:15
 Operator : SJ/JU
 Sample : I4373-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA4

Manual Integrations
 APPROVED

Sohil
 8/4/2017 5:28:07 PM

Quant Time: Aug 04 16:37:09 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 04 16:09:39 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	3224	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	16324	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	9878	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	22010	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	23817	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	20616m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4328	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	12751	0.18	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	1118	0.028	ng/ul#	88
5) 2-Methylnaphthalene	12.36	142	924	0.032	ng/ul	97
8) Acenaphthene	14.59	153	931	0.029	ng/ul	97
9) Fluorene	15.57	166	1070	0.026	ng/ul#	96
12) Phenanthrene	17.29	178	24981	0.417	ng/ul#	89
13) Anthracene	17.38	178	5056	0.087	ng/ul#	92
15) Fluoranthene	19.30	202	40805	0.544	ng/ul	84
17) Pyrene	19.66	202	37181	0.535	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	17513	0.254	ng/ul#	88
19) Chrysene	21.44	228	19033	0.287	ng/ul	96
21) Benzo(b)fluoranthene	23.14	252	25306m	0.397	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	9922m	0.158	ng/ul	
23) Benzo(a)pyrene	23.80	252	18958m	0.313	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.54	276	12598m	0.188	ng/ul	
25) Dibenzo(a,h)anthracene	26.56	278	3809m	0.068	ng/ul	
26) Benzo(g,h,i)perylene	27.36	276	13772m	0.244	ng/ul	

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

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