

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081617\
 Data File : BE093837.D
 Acq On : 16 Aug 2017 12:53
 Operator : SJ/JU
 Sample : I4672-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AG3

Manual Integrations
 APPROVED

Sohil
 8/17/2017 5:59:23 PM

Quant Time: Aug 17 01:41:30 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 17 00:58:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2397	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	12580	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	7602	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	15598	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	16198	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	15727m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4216	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	10569	0.21	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	44077	1.440	ng/ul#	93
5) 2-Methylnaphthalene	12.35	142	29080	1.303	ng/ul	99
7) Acenaphthylene	14.24	152	8642	0.272	ng/ul#	96
8) Acenaphthene	14.58	153	1524	0.061	ng/ul	95
9) Fluorene	15.57	166	1955	0.062	ng/ul#	87
12) Phenanthrene	17.29	178	59927	1.411	ng/ul#	92
13) Anthracene	17.38	178	12640	0.305	ng/ul#	93
15) Fluoranthene	19.30	202	127576	2.402	ng/ul	84
17) Pyrene	19.66	202	140659	2.975	ng/ul#	82
18) Benzo(a)anthracene	21.39	228	86983	1.852	ng/ul#	86
19) Chrysene	21.44	228	95166	2.113	ng/ul	99
21) Benzo(b)fluoranthene	23.14	252	121405	2.496	ng/ul	86
22) Benzo(k)fluoranthene	23.19	252	54549m	1.137	ng/ul	
23) Benzo(a)pyrene	23.80	252	96321m	2.081	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	59836m	1.171	ng/ul	
25) Dibenzo(a,h)anthracene	26.56	278	14742m	0.345	ng/ul	
26) Benzo(a,h,i)perylene	27.37	276	51449m	1.195	ng/ul	

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

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