

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081617\
 Data File : BE093842.D
 Acq On : 16 Aug 2017 15:58
 Operator : SJ/JU
 Sample : I4672-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AG1

Manual Integrations
 APPROVED

Sohil
 8/17/2017 5:59:32 PM

Quant Time: Aug 17 02:35:14 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	2210	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	11646	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	7139	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14739	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	15284	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	14085m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4309	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	10485	0.23	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	47742	1.685	ng/ul#	91
5) 2-Methylnaphthalene	12.36	142	34967	1.692	ng/ul	98
7) Acenaphthylene	14.24	152	28820	0.965	ng/ul	98
8) Acenaphthene	14.58	153	2728	0.116	ng/ul	92
9) Fluorene	15.57	166	3513	0.118	ng/ul#	83
12) Phenanthrene	17.29	178	109363	2.725	ng/ul#	90
13) Anthracene	17.38	178	27979	0.716	ng/ul#	90
15) Fluoranthene	19.30	202	226072	4.504	ng/ul	85
17) Pyrene	19.66	202	217860	4.883	ng/ul#	83
18) Benzo(a)anthracene	21.39	228	149196	3.367	ng/ul#	86
19) Chrysene	21.45	228	137939	3.245	ng/ul	98
21) Benzo(b)fluoranthene	23.14	252	200028	4.591	ng/ul	85
22) Benzo(k)fluoranthene	23.19	252	103157m	2.401	ng/ul	
23) Benzo(a)pyrene	23.81	252	155074m	3.742	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	114637m	2.505	ng/ul	
25) Dibenzo(a,h)anthracene	26.56	278	27821m	0.728	ng/ul	
26) Benzo(a,h,i)perylene	27.37	276	94895m	2.461	ng/ul	

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

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