

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082917\
 Data File : BE093897.D
 Acq On : 28 Aug 2017 15:00
 Operator : SJ/JU
 Sample : I4687-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AH2

Manual Integrations
 APPROVED

Sohil
 8/28/2017 5:51:52 PM

Quant Time: Aug 28 16:01:21 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 28 15:03:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	2211	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	11423	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	7461	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14173	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	14103	0.40	ng/ul	0.00
20) Perylene-d12	23.93	264	10707m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4906	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	11476	0.26	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.75	128	135198	4.866	ng/ul#	91
5) 2-Methylnaphthalene	12.36	142	127877	6.309	ng/ul	98
7) Acenaphthylene	14.24	152	22006	0.705	ng/ul#	92
8) Acenaphthene	14.58	153	14306	0.583	ng/ul	97
9) Fluorene	15.57	166	10774	0.347	ng/ul#	87
12) Phenanthrene	17.29	178	276368	7.160	ng/ul#	91
13) Anthracene	17.38	178	57601	1.532	ng/ul#	91
15) Fluoranthene	19.30	202	538144	11.149	ng/ul	84
17) Pyrene	19.66	202	495447	12.035	ng/ul#	84
18) Benzo(a)anthracene	21.40	228	295403	7.225	ng/ul#	88
19) Chrysene	21.45	228	308677	7.871	ng/ul	97
21) Benzo(b)fluoranthene	23.16	252	384404m	11.606	ng/ul	
22) Benzo(k)fluoranthene	23.20	252	150736m	4.615	ng/ul	
23) Benzo(a)pyrene	23.82	252	258200m	8.196	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.59	276	179997m	5.175	ng/ul	
25) Dibenzo(a,h)anthracene	26.59	278	47843m	1.646	ng/ul	
26) Benzo(g,h,i)perylene	27.41	276	182521m	6.227	ng/ul	

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

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