

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093305.D
 Acq On : 21 Jun 2017 5:37
 Operator : SJ/JU
 Sample : I3599-15DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5C91DL

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:19:32 PM

Quant Time: Jun 21 07:07:06 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 21 03:23:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1818	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9755	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5838	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14524	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	14883	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	12749	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	504	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	1474	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	8521	0.352	ng/ul#	93
5) 2-Methylnaphthalene	12.18	142	1954	0.118	ng/ul	100
7) Acenaphthylene	14.08	152	31918	1.274	ng/ul	96
8) Acenaphthene	14.42	153	1520	0.075	ng/ul	96
9) Fluorene	15.41	166	2095	0.088	ng/ul	90
12) Phenanthrene	17.13	178	26853	0.672	ng/ul#	89
13) Anthracene	17.22	178	39549	1.137	ng/ul#	89
15) Fluoranthene	19.15	202	120539	2.460	ng/ul	83
17) Pyrene	19.51	202	129916	2.960	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	70234	1.769	ng/ul#	84
19) Chrysene	21.29	228	96974	2.355	ng/ul	99
21) Benzo(b)fluoranthene	22.93	252	232117m	5.999	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	64663m	1.574	ng/ul	
23) Benzo(a)pyrene	23.55	252	104986	2.870	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	26.18	276	106571	2.516	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.19	278	27611	0.774	ng/ul	92
26) Benzo(g,h,i)perylene	26.96	276	65494	1.803	ng/ul	97

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

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