

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062317\
 Data File : BE093342.D
 Acq On : 23 Jun 2017 18:20
 Operator : SJ/JU
 Sample : I3653-10
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5C53

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:33:06 AM

Quant Time: Jun 24 04:34:39 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 24 04:29:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1530	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	8184	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5361	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	12459	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	12599	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	9894m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	2630	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	7583	0.22	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	465160	22.908	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	77770	5.601	ng/ul	99
7) Acenaphthylene	14.08	152	3704	0.161	ng/ul#	94
8) Acenaphthene	14.42	153	52604	2.834	ng/ul	97
9) Fluorene	15.42	166	59349	2.727	ng/ul#	96
12) Phenanthrene	17.14	178	413938	12.075	ng/ul#	90
13) Anthracene	17.22	178	31618	1.059	ng/ul#	91
15) Fluoranthene	19.15	202	191670	4.559	ng/ul	84
17) Pyrene	19.51	202	136519	3.674	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	37304	1.110	ng/ul#	86
19) Chrysene	21.29	228	35218	1.010	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	23750	0.791	ng/ul	88
22) Benzo(k)fluoranthene	22.97	252	9399m	0.295	ng/ul	
23) Benzo(a)pyrene	23.56	252	16991m	0.598	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.17	276	6578m	0.200	ng/ul	
25) Dibenzo(a,h)anthracene	26.19	278	1924m	0.070	ng/ul	
26) Benzo(g,h,i)perylene	26.95	276	5060	0.179	ng/ul#	93

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

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