

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061417\
 Data File : BE093190.D
 Acq On : 15 Jun 2017 3:46
 Operator : SJ/MA
 Sample : I3521-03
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5C30

Manual Integrations
 APPROVED

mohammad
 6/15/2017 3:54:05 PM

Quant Time: Jun 15 04:56:10 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 15 03:59:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3607	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	19064	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	11906	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	27808	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	27465	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	22563m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	6534	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	17979	0.24	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	8500	0.180	ng/ul#	93
5) 2-Methylnaphthalene	12.18	142	2133	0.066	ng/ul	99
7) Acenaphthylene	14.09	152	99812	1.954	ng/ul	96
8) Acenaphthene	14.43	153	3039	0.074	ng/ul	95
9) Fluorene	15.42	166	4384	0.091	ng/ul	90
12) Phenanthrene	17.14	178	443592	5.798	ng/ul#	84
13) Anthracene	17.22	178	150539	2.260	ng/ul#	86
15) Fluoranthene	19.16	202	1665659	17.752	ng/ul	83
17) Pyrene	19.52	202	1330673	16.429	ng/ul#	86
18) Benzo(a)anthracene	21.24	228	371419	5.070	ng/ul#	83
19) Chrysene	21.30	228	633336	8.334	ng/ul	99
21) Benzo(b)fluoranthene	22.94	252	754376	11.017	ng/ul	87
22) Benzo(k)fluoranthene	22.98	252	271792m	3.737	ng/ul	
23) Benzo(a)pyrene	23.56	252	289979m	4.479	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.19	276	252384m	3.367	ng/ul	
25) Dibenzo(a,h)anthracene	26.20	278	57301m	0.908	ng/ul	
26) Benzo(g,h,i)perylene	26.97	276	189830	2.953	ng/ul	96

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

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