

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061417\
 Data File : BE093177.D
 Acq On : 14 Jun 2017 19:46
 Operator : SJ/MA
 Sample : I3520-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A39

Manual Integrations
 APPROVED

mohammad
 6/15/2017 3:53:21 PM

Quant Time: Jun 15 10:46:08 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 02:27:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3736	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	19043	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	11866	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	27648	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	25992	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	22434m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	5621	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	14218	0.19	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.14	178	6668	0.088	ng/ul#	90
15) Fluoranthene	19.15	202	16754	0.180	ng/ul	84
17) Pyrene	19.51	202	16898	0.220	ng/ul#	90
18) Benzo(a)anthracene	21.24	228	6793	0.098	ng/ul#	83
19) Chrysene	21.29	228	9178	0.128	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.18	276	5667m	0.076	ng/ul	
26) Benzo(g,h,i)perylene	26.95	276	5178	0.081	ng/ul#	63

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

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