

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061417\
 Data File : BE093166.D
 Acq On : 14 Jun 2017 12:57
 Operator : SJ/MA
 Sample : I3520-05
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A37

Manual Integrations
 APPROVED

mohammad
 6/15/2017 3:52:53 PM

Quant Time: Jun 15 10:28:14 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 07:58:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3447	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	17320	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	11088	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	25404	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	25535	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	20517m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	5579	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	15001	0.22	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.13	178	4894	0.070	ng/ul#	90
15) Fluoranthene	19.15	202	11694	0.136	ng/ul	84
17) Pyrene	19.51	202	13627	0.181	ng/ul#	91
18) Benzo(a)anthracene	21.24	228	5131	0.075	ng/ul#	86
19) Chrysene	21.29	228	6940	0.098	ng/ul	96
21) Benzo(b)fluoranthene	22.93	252	9341	0.150	ng/ul	87
22) Benzo(k)fluoranthene	22.97	252	3551m	0.054	ng/ul	
23) Benzo(a)pyrene	23.55	252	5671m	0.096	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.17	276	4509m	0.066	ng/ul	
26) Benzo(g,h,i)perylene	26.95	276	4014	0.069	ng/ul#	84

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

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