

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061917\
 Data File : BE093264.D
 Acq On : 19 Jun 2017 23:29
 Operator : SJ/MA
 Sample : I3600-09
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A17

Manual Integrations
 APPROVED

mohammad
 6/20/2017 3:12:58 PM

Quant Time: Jun 30 08:50:50 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 08:49:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2594	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	13303	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	8437	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	20337m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	20543	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	16614m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	6676	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	19374	0.35	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	812	0.025	ng/ul#	78
7) Acenaphthylene	14.09	152	1171	0.032	ng/ul#	87
12) Phenanthrene	17.14	178	989	0.018	ng/ul#	80
13) Anthracene	17.22	178	6105	0.125	ng/ul#	93
15) Fluoranthene	19.15	202	1872	0.027	ng/ul#	61
17) Pyrene	19.52	202	3566	0.059	ng/ul	99
19) Chrysene	21.29	228	2257	0.040	ng/ul#	66
21) Benzo(b)fluoranthene	22.93	252	4912	0.097	ng/ul#	75
22) Benzo(k)fluoranthene	22.98	252	1536m	0.029	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.18	276	1998m	0.036	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	1738	0.037	ng/ul#	62

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

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