

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093285.D
 Acq On : 20 Jun 2017 17:20
 Operator : SJ/JU
 Sample : I3625-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5E17

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:18:26 PM

Quant Time: Jun 21 03:10:04 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 21 01:49:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1862	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9562	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	6579	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	14288	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	14425	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11927m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	4419	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	12157	0.31	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.56	128	1096685	46.225	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	232739	14.346	ng/ul	99
7) Acenaphthylene	14.09	152	17514	0.620	ng/ul	97
8) Acenaphthene	14.42	153	238805	10.485	ng/ul	96
9) Fluorene	15.42	166	255758	9.578	ng/ul#	93
12) Phenanthrene	17.14	178	1055927	26.859	ng/ul#	86
13) Anthracene	17.22	178	151982	4.440	ng/ul#	88
15) Fluoranthene	19.15	202	668340	13.863	ng/ul	83
17) Pyrene	19.51	202	475801	11.185	ng/ul#	86
18) Benzo(a)anthracene	21.24	228	126102	3.277	ng/ul#	86
19) Chrysene	21.29	228	91938	2.303	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	54434	1.504	ng/ul	87
22) Benzo(k)fluoranthene	22.97	252	22593m	0.588	ng/ul	
23) Benzo(a)pyrene	23.56	252	36271m	1.060	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.18	276	9860m	0.249	ng/ul	
25) Dibenzo(a,h)anthracene	26.20	278	2877m	0.086	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	7340	0.216	ng/ul#	93

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

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