

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062317\
 Data File : BE093346.D
 Acq On : 23 Jun 2017 20:46
 Operator : SJ/JU
 Sample : I3627-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A09

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:33:18 AM

Quant Time: Jun 24 04:56:05 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 24 04:29:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1648	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	8791	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5531	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	12894	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13785	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11698m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	2971	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	7969	0.23	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.56	128	29565	1.355	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	5239	0.351	ng/ul	98
7) Acenaphthylene	14.08	152	48641	2.049	ng/ul	96
8) Acenaphthene	14.42	153	7636	0.399	ng/ul	97
9) Fluorene	15.41	166	14118	0.629	ng/ul	91
12) Phenanthrene	17.13	178	106484	3.001	ng/ul#	90
13) Anthracene	17.22	178	57874	1.874	ng/ul#	89
15) Fluoranthene	19.15	202	255502	5.873	ng/ul	83
17) Pyrene	19.51	202	427118	10.507	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	257596	7.006	ng/ul#	83
19) Chrysene	21.30	228	356689	9.351	ng/ul	97
21) Benzo(b)fluoranthene	22.94	252	589834	16.615	ng/ul	86
22) Benzo(k)fluoranthene	22.98	252	212935m	5.648	ng/ul	
23) Benzo(a)pyrene	23.56	252	356784m	10.629	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.19	276	236609m	6.088	ng/ul	
25) Dibenzo(a,h)anthracene	26.19	278	62426m	1.907	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	171163	5.135	ng/ul	97

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

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