

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062417\
 Data File : BE093367.D
 Acq On : 24 Jun 2017 18:24
 Operator : SJ/JU
 Sample : I3627-12DL 10X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5B42DL

Manual Integrations
 APPROVED

mohammad
 6/29/2017 8:34:29 AM

Quant Time: Jun 26 02:00:07 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 26 01:58:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1326	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	7076	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	4099	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	10378	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	11979	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	10188	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	310	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	918	0.03	ng/ul	0.00
Target Compounds						
					Qvalue	
3) Naphthalene	10.56	128	1162	0.066	ng/ul#	90
5) 2-Methylnaphthalene	12.18	142	210	0.017	ng/ul	99
7) Acenaphthylene	14.08	152	11027	0.627	ng/ul	95
8) Acenaphthene	14.42	153	313	0.022	ng/ul#	87
9) Fluorene	15.42	166	418	0.025	ng/ul#	84
12) Phenanthrene	17.14	178	4777	0.167	ng/ul#	92
13) Anthracene	17.22	178	13438	0.541	ng/ul#	91
15) Fluoranthene	19.15	202	40772	1.164	ng/ul	83
17) Pyrene	19.51	202	98116	2.777	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	45090	1.411	ng/ul#	85
19) Chrysene	21.29	228	70391	2.124	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	168447m	5.448	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	55753m	1.698	ng/ul	
23) Benzo(a)pyrene	23.55	252	97282	3.328	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.18	276	63141	1.865	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	17332	0.608	ng/ul	92
26) Benzo(g,h,i)perylene	26.95	276	47366	1.632	ng/ul	96

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

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